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# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

CIRCULAR OF THE NATIONAL BUREAU OF STANDARDS 500

U. S. DEPARTMENT OF COMMERCE  
NATIONAL BUREAU OF STANDARDS









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# SELECTED VALUES of CHEMICAL THERMODYNAMIC PROPERTIES

CIRCULAR OF THE NATIONAL BUREAU OF STANDARDS 500

by  
Frederick D. Rossini, Donald D. Wagman,  
William H. Evans, Samuel Levine, and Irving Jaffe

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National Bureau of Standards

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This project was conducted, in part, under the sponsorship of the Office of Naval Research, Department of the Navy. The interest and assistance of that Office in recent years, as explained in the Preface, insured the continuation of the work.

## PREFACE

The assembly of the table of values of heats of formation in the section on Thermochemistry in volume V of the International Critical Tables, compiled by F. R. Bichowsky under the editorship of Edward W. Washburn and published in 1929 (McGraw-Hill Book Co., New York, N. Y.), was the first attempt ever made to collate all the published data involving heats of reaction and to prepare therefrom a self-consistent table of selected values of the heats of formation of the chemical substances. These tables were revised and extended by F. R. Bichowsky and F. D. Rossini and published in book form in 1936 (Reinhold Publishing Corp., New York, N. Y.) under the title "Thermochemistry of the Chemical Substances," and constituted the "assembly of a self-consistent table of 'best' values for the heats of formation of the chemical substances (except carbon compounds containing more than two carbon atoms), including heats of transition, fusion, and vaporization." However, it became obvious that the continued maintenance of a complete thermochemical table is a work too great to be carried on individual shoulders through part-time work as an "extracurricular" activity.

In 1940, the National Bureau of Standards assigned one investigator, D. D. Wagman, to work with F. D. Rossini on the maintenance of tables of thermochemical and chemical thermodynamic data. World War II interrupted the work. In 1946, an additional investigator, E. J. Blau, was assigned. In 1947, additional financial support for the work was received from the Office of Naval Research of the United States Department of the Navy, and two additional investigators, W. H. Evans and S. Levine, joined in the work. In 1948, E. J. Blau left to resume graduate school work and his place was taken by I. Jaffe. After the completion of these tables in June 1950, F. D. Rossini resigned from the National Bureau of Standards to accept a position at the Carnegie Institute of Technology, Pittsburgh, Pa. The project is now being carried on at the Bureau under the direction of D. D. Wagman.

Beginning as of March 31, 1947, and ending as of March 31, 1950, the tables were issued quarterly as prepared, in loose-leaf form, without references, one set to each laboratory in industry, educational institutions, and other organizations having need of and requesting the data. The number of sets so distributed in loose-leaf form as of March 31, 1950, was about 1,700.

The present volume includes all the tables of Series I and II issued in loose-leaf form plus all the references, which were not previously issued.

Grateful acknowledgment is made to Edmund J. Blau for his early work on the tables, and to Alberta L. Kelvie, Marie T. Lynch, Helen Marion Flanagan, Rita M. Braun, Corinne R. DeCelles, Agnes C. Fahl, and others of the staff of the American Petroleum Institute Research Project 44, for assistance in the editorial work on the book.

The authors will greatly appreciate receiving comments regarding errors that have escaped attention.

A. V. ASTIN, *Acting Director.*

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# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

By Frederick D. Rossini, Donald D. Wagman, William H. Evans, Samuel Levine, and Irving Jaffe

## I. INTRODUCTION

### Unit of Energy and Fundamental Constants

All the values of energy given in these tables are expressed, insofar as possible, in terms of the thermochemical calorie, now defined in terms of the absolute joule, but formerly defined in terms of the international electrical joule, the relations being such as to keep unchanged, in the transition from the international to the absolute joule, the amount of energy represented by the calorie. These relations, and the values of the fundamental constants used in this work, follow:

1 atm = 1,013,250 dynes/cm<sup>2</sup>; definition; atm = standard atmosphere.

1 mm Hg (pressure) = (1/760) atm = 1333.2237 dynes/cm<sup>2</sup>; definition; mm Hg (pressure) = standard millimeter mercury.

1 abs joule = 0.999835 ± 0.000052 int. joule; abs = absolute; int. = international.

1 cal = 4.1840 abs joule = 4.1833 int. joule; definition; cal = thermochemical calorie.

$T_0$ °C = 273.160 ± 0.010 °K; absolute temperature of the ice point, 0°C.

$R$  = 8.31439 ± 0.00034 abs joule/deg mole = 1.98719 ± 0.00008 cal/deg mole = 82.0567 ± 0.0034 cm<sup>3</sup> atm/deg mole;  $R$  = gas constant per mole.

$F$  = 96485.30 ± 10.0 abs coul/g-equiv, or abs joule/abs volt g-equiv, = 96501.2 ± 10.0 int. coul/g-equiv, or int. joule/int. volt g-equiv, = 23060.5 ± 2.4 cal/abs volt g-equiv = 23068.1 ± 2.4 cal/int. volt g-equiv;  $F$  = Faraday constant.

$Z = Nhc = 11.9600 ± 0.0036$  abs joule cm/mole = 2.85851 ± 0.0009 cal cm/mole;  $Z$  = constant relating wave number,  $\nu$ , and energy per mole,  $E$ , in the relation  $(\Delta E) = Nhc (\Delta \nu) = Z (\Delta \nu)$ .

$(Z/R) = (\hbar c/k) = c_2 = 1.43847 ± 0.00045$  cm deg;  $c_2$  = second radiation constant.

The reader is referred to Mueller and Rossini-1<sup>1</sup> for a discussion of the calorie and the joule, and to Rossini-29 for a discussion of the relations among the fundamental constants. In connection with the Faraday constant, the value given above is that which was generally accepted as "best"

until last year, when new calculations and data indicated that a higher value was better. However, it was necessary, for preservation of the internal consistency of the present tables, to use the same value throughout.

### Scale of Temperature

The values of temperature given in these tables conform to the International Temperature Scale of 1948 (see Stinson-1).

### Definition of Symbols

The symbols used in these tables are defined as follows:  $P$  = pressure;  $V$  = volume;  $T$  = absolute temperature;  $E$  = energy, internal or intrinsic;  $S$  = entropy;  $H = E + PV$  = heat content (or enthalpy);  $F = E + PV - TS = H - TS$  = free energy;  $K$  = equilibrium constant;  $C_p = (dH/dT)_p$  = heat capacity at constant pressure.

Circular superscript, °, denotes the standard state (thermodynamic standard reference state).

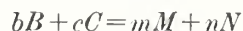
Numerical subscript, as <sub>298.16</sub> or <sub>0</sub>, denotes the absolute temperature in degrees Kelvin.

$S^\circ$  = the entropy of the given substance in the standard state at the indicated temperature, omitting nuclear contributions, as isotopic mixing, nuclear spins, etc., which contributions cancel out in all reactions of practical importance.

$\Delta$  denotes the increment of a given property for a given process or reaction, taken as the value for the final state (sum of products) less that for the initial state (sum of reactants).

$\Delta H$ ,  $\Delta F$ ,  $\Delta S$ ,  $\Delta C_p$  = the increment in heat content, free energy, entropy, and heat capacity, respectively, for the given process or reaction, with each substance in a specified thermodynamic state.

$K$  = the equilibrium constant for a given process or reaction, defined as the proper quotient of activities ( $a$ ), of the substances involved in the reaction, at thermodynamic equilibrium at the given temperature, pressure, etc. For example, for the reaction



with each of the reactants and products at equilibrium,

<sup>1</sup>The numbers following a dash after a name in the text (Part I) and in the Tables of Specific References (Parts IV and V) refer to the General List of References, Part VI.

$$K = \frac{(a_M)^m (a_N)^n}{(a_B)^b (a_C)^c}.$$

The equilibrium constant for a given process or reaction is related to the standard change in free energy for that process or reaction by the relation:

$$\log_{10} K = -\Delta F^\circ / (2.302585RT).$$

$\Delta Hf^\circ$  = the standard heat of formation, which is the increment in heat content associated with the reaction of forming the given compound from its elements, with each substance in its standard state at the given temperature.

$\Delta Ff^\circ$  = the standard free energy of formation, which is the increment in free energy associated with the reaction of forming the given compound from its elements, with each substance in its standard state at the given temperature.

$\log_{10} Kf$  = the logarithm (to the base 10) of the equilibrium constant of formation, for the reaction of forming 1 mole of the given compound from its elements, with each substance in its standard state at the given temperature. The equilibrium constant of formation is related to the standard free energy of formation, per mole, by the relation:

$$\log_{10} Kf = -\Delta Ff^\circ / (2.302585RT).$$

### Abbreviations Regarding Physical States

The physical state of each substance is indicated as gaseous (g), liquid (liq), crystalline (c), glass (gls), amorphous (amorp), colloidal (coll), or in aqueous (aq) or other solution. The composition of a solution is indicated by the number of moles of solvent associated with 1 mole of the substance (solute). When the concentration is not specified, the solution is understood to be "dilute".

### Standard States

The standard states used in these tables follow the conventions of Lewis and Randall and are defined as follows:

The standard state for a pure solid or liquid at 25° C is normally taken as the pure substance in its thermodynamically stable modification at a pressure of one atmosphere at 25° C (exceptions will be clearly indicated);

The standard state for a gas at any temperature is taken as the hypothetical reference state of unit fugacity (1 atmosphere) at the given temperature, in which state the heat content (enthalpy) is the same as that of the real gas at zero pressure at the same temperature;

The standard state for a solute in aqueous solution is taken as the hypothetical ideal state of unit molality (indicated as "hyp. m=1"), in which state the partial molal heat content (enthalpy) and heat capacity (also the partial molal volume) of the solute are the same as those in the infinitely dilute solution.

For further description of these standard states,

the reader is referred to Lewis and Randall-7 and Rossini-29.

### Conventions Regarding Reference States

The values for the heats of formation of substances in aqueous or other solution represent values of the apparent molal heat content, rather than the partial molal heat content at the given concentration. At infinite dilution, the apparent molal heat content and the apparent molal heat capacity are the same as the partial molal heat content and the partial molal heat capacity, respectively.

The values of the chemical thermodynamic properties tabulated for individual ions in aqueous solution are based on the convention that the values of  $\Delta Hf^\circ$ ,  $\Delta Ff^\circ$ ,  $S^\circ$  and  $C_p^\circ$  for  $H^+$  (aq, hyp. m=1) are zero. The values of the quantities  $\Delta Hf^\circ$ ,  $\Delta Ff^\circ$ ,  $S^\circ$ , and  $C_p^\circ$  for a strong electrolyte in aqueous solution, in the standard state, are equal to the sums of these values for the appropriate kinds and number of individual ions assumed to constitute the undissociated molecule of the given electrolyte.

In connection with the reference states for aqueous ions, it is to be noted that this book follows Latimer-1 in assigning for  $H^+$  (aq) the value zero for  $S^\circ$  rather than for  $\Delta Sf^\circ$ . For a discussion of the possible advantages of assigning the value zero to  $\Delta Sf^\circ$  for  $H^+$  (aq), the reader is referred to Rossini-29.

### Properties Given in the Tables of Series I

The tables of Series I give the following information, as known, in the order of columns, left to right:

1. The molecular formula of the given substance.
2. Information, as necessary, to complete the description of the given substance, such as the name of a particular crystalline form, or the name of the solvent if the substance is in solution and the solvent is other than water, or the isomeric name if the substance is one of a number of isomers, etc.
3. The physical state of the given substance.
4. The value of  $\Delta Hf^\circ$ , the standard heat of formation of the given substance from its elements, at 0°K.
5. The value of  $\Delta Hf^\circ$ , the standard heat of formation of the given substance from its elements, at the reference temperature, 298.16°K (25°C).
6. The value of  $\Delta Ff^\circ$ , the standard free energy of formation of the given substance from its elements, at the reference temperature, 298.16°K (25°C).
7. The value of  $\log_{10} Kf$ , the logarithm of the equilibrium constant for the reaction of forming the given substance from its elements, at the reference temperature, 298.16°K (25°C).



8. The value of  $S^\circ$ , the entropy of the given substance in its thermodynamic standard state, at the reference temperature, 298.16°K (25°C).

9. The value of  $C_p^\circ$ , the heat capacity of the given substance in its thermodynamic standard state, at the reference temperature, 298.16°K (25°C).

### Properties Given in the Tables of Series II

The tables of Series II give the following information, as known, in the order of columns, left to right:

1. The molecular formula of the given substance.

2. Information, as necessary, to complete the description of the given substance.

3. The name of the process, as transition, fusion, vaporization, or sublimation, to which the tabulated thermodynamic properties apply.

4. The initial and final states for the given process.

5. The pressure.

6. The temperature, °K and °C.

7. The value of  $\Delta H$ , the increment in heat content associated with the given process.

8. The value of  $\Delta S$ , the increment in entropy associated with the given process.

9. The value of  $\Delta C_p$ , the heat capacity of the given substance in the final state less that in the initial state.

It is to be noted that when the pressure is 1 atm the values of temperature given for the process of vaporization correspond to values of the normal boiling point. Values of the temperature of the liquid-vapor equilibrium at other pressures may also be given.

The values of the temperature of the solid-liquid equilibrium correspond to values of the freezing (or melting) point of the given substance. When the same values of both temperature and pressure are given for the solid-liquid and the liquid-gas or solid-gas equilibrium, the given values of temperature and pressure are those of the so-called triple point, at which the solid, liquid, and gaseous phases of the given substance are in thermodynamic equilibrium.

### Internal Consistency

The tables of values in this volume are internally consistent in the sense that all the known physical and thermodynamic relations existing between different properties for a given substance, or the same property for different substances, are satisfied by the tabulated values of the properties. For example, the value of any thermodynamic property for a reaction must come out the same regardless of the path chosen for the evaluation. In particular, the values of  $\Delta H$ ,  $\Delta F$ , and  $\Delta S$  for the same process or reaction must satisfy the well-known thermodynamic relation involving these properties and the temperature.

It is realized that new data will have become available on certain compounds after the selection of values for them was made for these tables. In such cases, it was not always possible to change the selected value to conform to the new information because of the necessity of maintaining the internal consistency of the tables. However, readers encountering such cases can of course rearrange their own particular calculations to take advantage of the new information.

### Uncertainties

Following Bichowsky and Rossini-1, all the values given in these tables have been calculated from original experimental data using consistent values for all subsidiary quantities. Heats of reaction, of dilution, of solution, of neutralization, etc., may be calculated from the appropriate tabulated values with an accuracy as good as is known. The number of significant figures in any one value in the table does not indicate the absolute accuracy of that value. The uncertainty in any value will be greater than that of the least accurate determination in the total chain of reactions used to calculate the value. More definite information regarding the uncertainty to be attached to any given tabulated value may be obtained by examination of the material in the cited references.

### Methods of Calculation

The methods of calculation which may be employed in the utilization of the values of thermodynamic properties given in these tables are quite varied. For information regarding the calculation of values of  $\Delta H$  for any process or reaction, the reader is referred to pages 12 and 13 of Bichowsky and Rossini-1, to Lewis and Randall-7, and to Rossini-29. For discussion of the other thermodynamic properties and their interrelations, and for more general methods of thermodynamic calculations, including free energies and entropies, the reader is referred to the latter two references.

### References

Part VI of this book gives a complete general list of references in alphabetical order of the first-named author. The references to a given author or authors will follow in numerical suborder in the list, as Lewis and Randall-1, Lewis and Randall-2, etc.

The reader interested in the origin of the value of a given property of a given compound, in either Series I or Series II, will find in the tables of specific references, in Part IV for Series I and in Part V for Series II, reference to the appropriate investigators, as Lewis and Randall-3. Then the complete reference to the literature is obtained from the general list of references in Part VI.



## Order of Arrangement of the Chemical Substances

The substances listed in the tables are arranged in a manner similar to that used by Bichowsky and Rossini-1, with some modification. The chemical elements are arranged in the order shown in figure 1 and in table 1, "List of tables and cor-

responding elements." If the elements are taken to be arranged in the order, a, b, c, d, e, etc., the chemical substances are in the order: a; b, ba; c, ca, cb, cba; d, da, db, dba, dc, dca, dcba, dcba; etc. Table 2, "List of elements and corresponding tables," gives the number of the table corresponding to a given element, listed alphabetically.

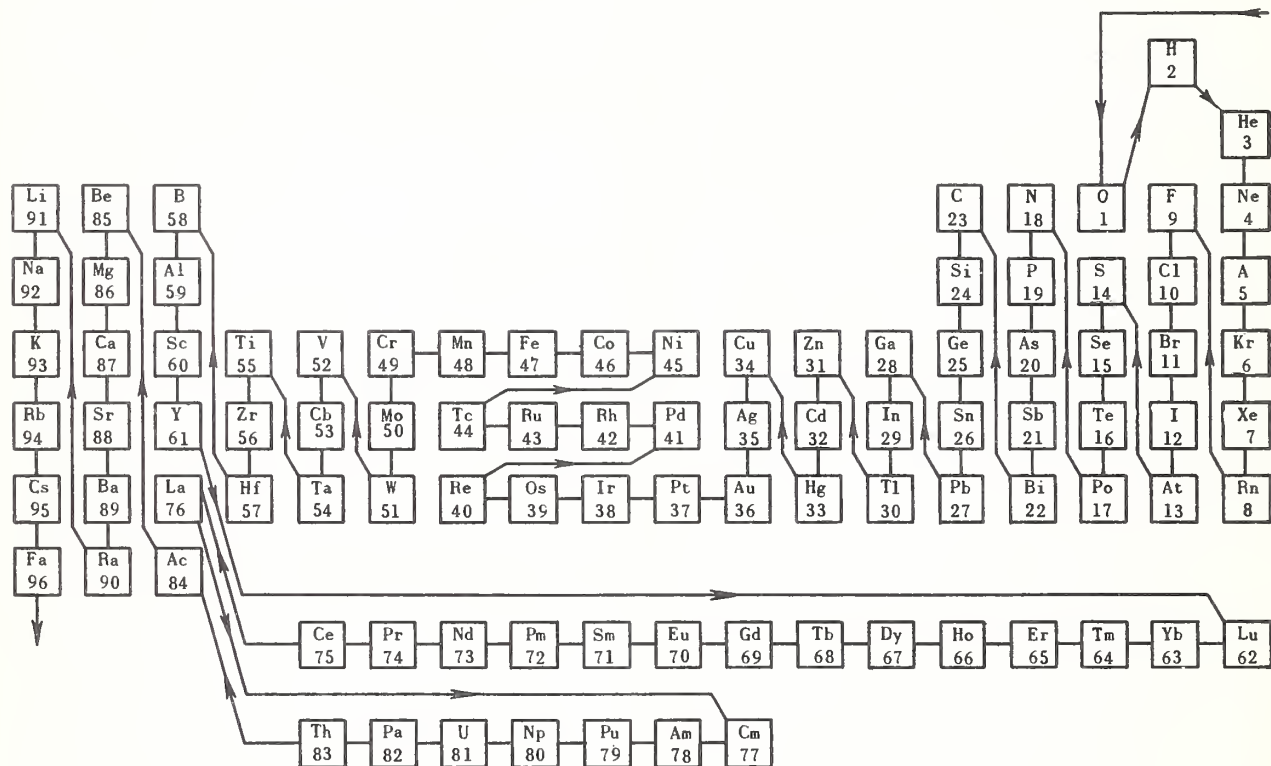


FIGURE 1. Standard order of arrangement of the chemical elements.

TABLE 1.—List of tables and corresponding elements

Table	Element	Symbol	Atomic number	Table	Element	Symbol	Atomic number
1	Oxygen	O	8	49	Chromium	Cr	24
2	Hydrogen	H	1	50	Molybdenum	Mo	42
3	Helium	He	2	51	Tungsten	W	74
4	Neon	Ne	10	52	Vanadium	V	23
5	Argon	A	18	53	Columbium (Niobium)	Cb	41
6	Krypton	Kr	36	54	Tantalum	Ta	73
7	Xenon	Xe	54	55	Titanium	Ti	22
8	Radon	Rn	86	56	Zirconium	Zr	40
9	Fluorine	F	9	57	Hafnium	Hf	72
10	Chlorine	Cl	17	58	Boron	B	5
11	Bromine	Br	35	59	Aluminum	Al	13
12	Iodine	I	53	60	Scandium	Sc	21
13	Astatine	At	85	61	Yttrium	Y	39
14	Sulfur	S	16	62	Lutetium	Lu	71
15	Selenium	Se	34	63	Ytterbium	Yb	70
16	Tellurium	Te	52	64	Thulium	Tm	69
17	Polonium	Po	84	65	Erbium	Er	68
18	Nitrogen	N	7	66	Holmium	Ho	67
19	Phosphorus	P	15	67	Dysprosium	Dy	66
20	Arsenic	As	33	68	Terbium	Tb	65
21	Antimony	Sb	51	69	Gadolinium	Gd	64
22	Bismuth	Bi	83	70	Europium	Eu	63
23	Carbon	C	6	71	Samarium	Sm	62
24	Silicon	Si	14	72	Promethium	Pm	61
25	Germanium	Ge	32	73	Neodymium	Nd	60
26	Tin	Sn	50	74	Praseodymium	Pr	59
27	Lead	Pb	82	75	Cerium	Ce	58
28	Gallium	Ga	31	76	Lanthanum	La	57
29	Indium	In	49	77	Curium	Cm	96
30	Thallium	Tl	81	78	Americium	Am	95
31	Zinc	Zn	30	79	Plutonium	Pu	94
32	Cadmium	Cd	48	80	Neptunium	Np	93
33	Mercury	Hg	80	81	Uranium	U	92
34	Copper	Cu	29	82	Protactinium	Pa	91
35	Silver	Ag	47	83	Thorium	Th	90
36	Gold	Au	79	84	Actinium	Ac	89
37	Platinum	Pt	78	85	Beryllium	Be	4
38	Iridium	Ir	77	86	Magnesium	Mg	12
39	Osmium	Os	76	87	Calcium	Ca	20
40	Rhenium	Re	75	88	Strontium	Sr	38
41	Palladium	Pd	46	89	Barium	Ba	56
42	Rhodium	Rh	45	90	Radium	Ra	88
43	Ruthenium	Ru	44	91	Lithium	Li	3
44	Technetium	Tc	43	92	Sodium	Na	11
45	Nickel	Ni	28	93	Potassium	K	19
46	Cobalt	Co	27	94	Rubidium	Rb	37
47	Iron	Fe	26	95	Cesium	Cs	55
48	Manganese	Mn	25	96	Francium	Fa	87

TABLE 2.—List of elements and corresponding tables

Element	Table	Symbol	Atomic number	Element	Table	Symbol	Atomic number
Actinium	84	Ac	89	Molybdenum	50	Mo	42
Aluminum	59	Al	13	Neodymium	73	Nd	60
Americium	78	Am	95	Neon	4	Ne	10
Antimony	21	Sb	51	Neptunium	80	Np	93
Argon	5	A	18	Nickel	45	Ni	28
Arsenic	20	As	33	Nitrogen	18	N	7
Astatine	13	At	85	Osmium	39	Os	76
Barium	89	Ba	56	Oxygen	1	O	8
Beryllium	85	Be	4	Palladium	41	Pd	46
Bismuth	22	Bi	83	Phosphorus	19	P	15
Boron	58	B	5	Platinum	37	Pt	78
Bromine	11	Br	35	Plutonium	79	Pu	94
Cadmium	32	Cd	48	Polonium	17	Po	84
Calcium	87	Ca	20	Potassium	93	K	19
Carbon	23	C	6	Praseodymium	74	Pr	59
Cerium	75	Ce	58	Promethium	72	Pm	61
Cesium	95	Cs	55	Protactinium	82	Pa	91
Chlorine	10	Cl	17	Radium	90	Ra	88
Chromium	49	Cr	24	Radon	8	Rn	86
Cobalt	46	Co	27	Rhenium	40	Re	75
Columbium (Niobium)	53	Cb	41	Rhodium	42	Rh	45
Copper	34	Cu	29	Rubidium	94	Rb	37
Curium	77	Cm	96	Ruthenium	43	Ru	44
Dysprosium	67	Dy	66	Samarium	71	Sm	62
Erbium	65	Er	68	Scandium	60	Sc	21
Europium	70	Eu	63	Selenium	15	Se	34
Fluorine	9	F	9	Silicon	24	Si	14
Francium	96	Fa	87	Silver	35	Ag	47
Gadolinium	69	Gd	64	Sodium	92	Na	11
Gallium	28	Ga	31	Strontium	88	Sr	38
Germanium	25	Ge	32	Sulfur	14	S	16
Gold	36	Au	79	Tantalum	54	Ta	73
Hafnium	57	Hf	72	Technetium	44	Tc	43
Helium	3	He	2	Tellurium	16	Te	52
Holmium	66	Ho	67	Terbium	68	Tb	65
Hydrogen	2	H	1	Thallium	30	Tl	81
Indium	29	In	49	Thorium	83	Th	90
Iodine	12	I	53	Thulium	64	Tm	69
Iridium	38	Ir	77	Tin	26	Sn	50
Iron	47	Fe	26	Titanium	55	Ti	22
Krypton	6	Kr	36	Tungsten	51	W	74
Lanthanum	76	La	57	Uranium	81	U	92
Lead	27	Pb	82	Vanadium	52	V	23
Lithium	91	Li	3	Xenon	7	Xe	54
Lutetium	62	Lu	71	Ytterbium	63	Yb	70
Magnesium	86	Mg	12	Yttrium	61	Y	39
Manganese	48	Mn	25	Zinc	31	Zn	30
Mercury	33	Hg	80	Zirconium	56	Zr	40



## II. TABLES OF SELECTED VALUES OF PROPERTIES OF SERIES I:

Heat of Formation at 25°C

Free Energy of Formation at 25°C

Entropy at 25°C

Heat Capacity at 25°C

Logarithm of the Equilibrium Constant of Formation at 25°C

Heat of Formation at 0°K

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 1-1. Oxygen (at. no., 8; at. wt., 16.0000)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1947; March 31, 1950

Substance		At 0°K		At 298.16°K (25°C)		log <sub>10</sub> K <sub>f</sub>	S°	C <sub>p</sub> °
Formula	Description	State	ΔH <sub>f</sub> ° kcal/mole	ΔH <sub>f</sub> ° kcal/mole	ΔF <sub>f</sub> ° kcal/mole			
O		g	58.586	59.159	54.994	-40.31022	38.4689	5.2364
O*		g	372.555	374.609				
O**		g	1183.09	1186.624				
O***		g	2450.0	2454.98				
O****		g	4234.8	4241.30				
O*****		g	6660.9	6668.9				
O*****		g	10045.	10054.8				
O*****		g	27091.	27102.				
O <sup>-</sup>		g	-12.21	-13.12				
O <sup>--</sup>		g						
O <sub>2</sub>		g	0.000	0.000	0.000	0.0000	49.003	7.017
O <sub>2</sub> <sup>+</sup>		aq	281.6	-3.8				
		g		283.1				
O <sub>3</sub>	ozone	g		34.0	39.06	-28.631	56.8	9.12
		aq		32.2				
O <sub>4</sub>		g		-0.16				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 2-1. Hydrogen (at. no., 1; at. wt., 1.0080)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1947; June 30, 1947; March 31, 1950

Substance		State	Description	At 298.16°K (25°C)				
Formula	ΔHf° kcal/mole			ΔFf°		log10Kf	S° cal/deg mole	Cp° cal/deg mole
				kcal/mole	kcal/mole			
H	51.620	g	-  std. state hyp. m = 1	52.089	48.575	-35.60481	27.3927	4.9680
H +	365.138	g		367.088				
H +		aq		0.000	0.000	0.00000	0.000	0.000
H2	0.000	g	aq	0.000	0.000	0.00000	31.211	6.892
H2								
OH	10.0	g	std. state, hyp. m = 1	10.06	8.93	-6.546	43.888	7.141
OH -	-75.	g		-76.4				
OH -		aq		-54.957	-37.595	27.5566	-2.519	-32.0
H2O	-57.107	g	water  liq	-57.7979	-54.6357	40.04724	45.106	8.025
				-68.3174	-56.6902	41.55313	16.716	17.996

SERIES I

Washington, D.C.

National Bureau of Standards

Table 2-2. Hydrogen (at. no., 1; at. wt., 1.0080)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1947; June 30, 1947; March 31, 1950

Substance								
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
			At 0°K kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
H <sub>2</sub> O <sub>2</sub>	std. state, hyp. m = 1 in 1 H <sub>2</sub> O 2 H <sub>2</sub> O 3 H <sub>2</sub> O 5 H <sub>2</sub> O 10 H <sub>2</sub> O ∞ H <sub>2</sub> O	g		-31.83				
		liq		-44.84				
		aq		-45.68				
		aq		-45.43				
		aq		-45.55				
		aq		-45.60				
		aq		-45.64				
		aq		-45.66				
		aq		-45.68				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 2a-1. Hydrogen isotope of mass 1 (at. no., 1; at. wt., 1.0078) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1947; March 31, 1948; March 31, 1950											
Substance			State	Description	At 0°K kcal/mole	ΔH <sup>o</sup> kcal/mole	ΔF <sup>o</sup> kcal/mole	ΔF <sup>o</sup> kcal/mole	log <sub>10</sub> K <sub>f</sub>	S <sup>o</sup> cal/deg mole	C <sub>p</sub> <sup>o</sup> cal/deg mole
Formula		At 0°K kcal/mole									
1H		g			51.620	52.069	48.575	-35.6048		27.392	4.9660
1H <sup>+</sup>		g			365.138	367.088					
1H <sup>-</sup>		g			35.066	34.054					
1H <sub>2</sub>		g			0.000	0.000	0.000	0.0000		31.210	6.892
1H <sub>2</sub> <sup>+</sup>		g			355.67	357.15					
1H <sub>2</sub> O		g			-57.1040	-57.7976	-54.6352	40.04685		45.105	8.025
		liq				-68.3171	-56.6902	41.55314		16.716	17.996

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 2b-1. Hydrogen isotope of mass 2 (at. no., 1; at. wt., 2.0142)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1947; March 31, 1948; March 31, 1950

Substance		State	Description	At 298.16°K (25°C)					
Formula	$\Delta H_f^\circ$			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
	kcal/mole			kcal/mole	kcal/mole		cal/deg mole	cal/deg mole	
$2H$		g		52.525	52.982	49.358	-36.1787	29.456	4.9680
$2H^+$		g		366.128	368.066				
$2H^-$		g		35.967	34.943				
$2H_2$		g		0.000	0.000	0.000	0.0000	34.602	6.98
$1H^2H$		g		0.037	0.037	-0.391	0.2866	34.341	6.98
$2H_2O$		g			-59.5628	-56.0670	41.09634	47.379	8.19
		liq			-70.4133	-58.2062	42.66436	18.162	19.70
$1H^2HO$		g		-58.035	-58.735	-55.828	40.9212	47.66	8.06
		liq			-69.393	-57.926	42.4590	18.95	18.85

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 3-1. Helium (at. no., 2; at. wt., 4.003)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1947; March 31, 1950

Substance		State	Description	Formula	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f^\dagger$	$S^\circ$	$C_p^\circ$			
At 0°K					At 298.16°K (25°C)							
kcal/mole					kcal/mole	kcal/mole		cal/deg mole	cal/deg mole			
He	-	g			0.000	0.000	0.000	30.126	4.9660			
He <sup>+</sup>		g			566.86	568.34						
He <sup>++</sup>		g			1821.48	1824.44						
He <sub>2</sub> <sup>+</sup>		g										



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 4-1. Neon (at. no., 10; at. wt., 20.183) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1947; March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	At 0°K kcal/mole			kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole	cal/deg mole		
Ne			g	0.000	0.000	0.0000	34.948	4.9680	
Ne <sup>+</sup>			g	437.19	498.67				
Ne <sup>++</sup>			g	1444.35	1447.32				
Ne <sup>+++</sup>			g	2914.0	2918.5				
Ne <sup>++++</sup>			g	5154.8	5160.7				
Ne <sup>+++++</sup>			g	8070.3	8077.7				
Ne <sup>+++++</sup>			g	11712.	11721.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 5-1. Argon (at. no., 18; at. wt., 39.944)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1947; March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
A			g	0.000	0.000	0.000	0.000	36.983.	4.9680
A-5H <sub>2</sub> O			c		-357.2				
A +			g	363.345	364.826				
A ++			g	1000.28	1003.24				
A +++			g	1943.5	1947.9				
A ++++			g	3322.	3328.				
A +++++			g	5052.	5056.				
A ++++++			g	7158.	7167.				
A +++++++			g	10017.	10027.				
A ++++++++			g	13326.	13338.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 6-1. Krypton (at. no., 36; at. wt., 83.7)

HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1947; March 31, 1950

Substance		State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description							
Kr		g	0.000	0.000	0.000	0.0000	39.19	4.9680
Kr+5H <sub>2</sub> O		c		-357.1				
Kr +		g	322.7	324.25				
Kr ++		g	889.27	892.23				
Kr +++		g	1741.2	1745.6				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I


National Bureau of Standards

Washington, D.C.

Table 7-1. Xenon (at. no., 54; at. wt., 131.3)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1947; March 31, 1950

Substance								
Formula	Description	State	$\Delta H_f^{\circ}$	$\Delta H_f^{\circ}$	$\Delta F_f^{\circ}$	$\log_{10} K_f$	$S^{\circ}$	$C_p^{\circ}$
			At 0°K kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
Xe		g	0.000	0.000	0.000	0.000	40.53	4.9680
Xe·6H <sub>2</sub> O		c		-4.28.				
Xe <sup>+</sup>		g	279.66	281.14				
Xe <sup>++</sup>		g	768.66	771.62				
Xe <sup>+++</sup>		g	1509.3	1513.7				
Xe <sup>++++</sup>		g	2584.	2590.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 8-1. Radon (at. no., 86; at. wt., 222) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1947; March 31, 1950										
Substance		Description	State	$\Delta H_f^{\circ}$		$\Delta F_f^{\circ}$		$\log_{10} K_f$	$S^{\circ}$	$C_p^{\circ}$
Formula				At 0°K		At 298.16°K (25°C)				
				kcal/mole	kcal/mole	kcal/mole	kcal/mole			
Rn			g	0.000	0.000	0.000	0.000	0.0000	42.10	4.9680
Rn +			g	247.8	249.3					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 9-1. Fluorine (at. no., 9; at. wt., 19.00)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1947; June 30, 1947; March 31, 1950

Substance		State	At 298.16°K (25°C)			
Formula	Description		$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole
F		g	17.8	18.3	-10.41	37.917
F <sup>+</sup>		g	419.57	421.55		
F <sup>++</sup>		g	1226.62	1229.68		
F <sup>+++</sup>		g	2670.9	2675.9		
F <sup>++++</sup>		g	4682.7	4689.1		
F <sup>+++++</sup>		g	7316.6	7324.5		
F <sup>++++++</sup>		g	10940.	10949.		
F <sup>+++++++</sup>		g	15210.	15221.		
F <sup>+++++++</sup>		g	37201.	37214.		
F <sup>-</sup>		g	-78.5	-79.5		
F <sup>-</sup>	std. state, hyp. m=1	aq		-78.66	48.435	-2.3
F <sub>2</sub>		g	0.000	0.000	0.000	48.6
F <sub>2</sub> O		g		5.5	-7.126	58.95
HF		g		-64.2	47.402	41.47
HF	std. state, hyp. m=1 in 2 H <sub>2</sub> O	aq		-66.08	48.435	-2.3
		aq		-75.167		6.95
						-29.5
						7.52
						-29.5



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 9-2. Fluorine (at. no., 9; at. wt., 19.00)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1947; June 30, 1947; March 31, 1950

Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description		At 0°K	At 298.16°K (25°C)				
			kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
HF	in	aq						
	3 H <sub>2</sub> O	aq			-75.385			
	4 H <sub>2</sub> O	aq			-75.478			
	5 H <sub>2</sub> O	aq			-75.525			
	10 H <sub>2</sub> O	aq			-75.605			
	20 H <sub>2</sub> O	aq			-75.634			
	30 H <sub>2</sub> O	aq			-75.646			
	40 H <sub>2</sub> O	aq			-75.654			
	50 H <sub>2</sub> O	aq			-75.660			
	100 H <sub>2</sub> O	aq			-75.686			
	200 H <sub>2</sub> O	aq			-75.704			
	300 H <sub>2</sub> O	aq			-75.722			
	400 H <sub>2</sub> O	aq			-75.74			
	500 H <sub>2</sub> O	aq			-75.760			
	1000 H <sub>2</sub> O	aq			-75.872			
	∞ H <sub>2</sub> O	aq			-78.66			
		aq			-153.6			
		g			-426.0			
HF <sub>2</sub> <sup>-</sup>								
(HF) <sub>6</sub>								
								-29.5

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 10-1. Chlorine (at. no., 17; at. wt., 36.457)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1947; June 30, 1947; June 30, 1948

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 298.16°K (25°C)				
			At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
Cl		g	28.61	29.012	25.192	-18.4651	5.2203
Cl <sup>+</sup>		g	327.466	329.349			
Cl <sup>++</sup>		g	876.30	879.66			
Cl <sup>+++</sup>		g	1796.55	1801.39			
Cl <sup>++++</sup>		g	3029.21	3035.53			
Cl <sup>+++++</sup>		g	4592.8	4600.6			
Cl <sup>+++++</sup>		g					
Cl <sup>+++++</sup>		g					
Cl <sup>+++++</sup>		g					
Cl <sup>-</sup>		g	-57.2	-58.3			
Cl <sub>2</sub>	std. state, hyp. m=1	aq	0.0000	-40.023	-31.350	22.9792	-30.0
	saturated solution			0.0000	0.0000	0.0000	8.11
	in CCl <sub>4</sub> (tetrachloromethane)			-6.0			
				-4.6			
				-564.4			
Cl <sub>2</sub> ·8H <sub>2</sub> O		c		25.			
Cl <sub>3</sub>		g					

## SERIES I

Washington, D.C.

Table 10-2. Chlorine (at. no., 17; at. wt., 35.457)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

Substance		State	Description	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K kcal/mole			At 298.16°K (25°C) kcal/mole	At 298.16°K (25°C) cal/deg mole	At 298.16°K (25°C) cal/deg mole			
ClO		g				33.			
ClO <sup>-</sup>	std. state, hyp. m=1	aq						10.3	
ClO <sub>2</sub>		g		25.5	29.5	-21.62		59.6	
		liq							
		aq			18.1				
		aq			-16.5		-2.56	24.0	
ClO <sub>2</sub> <sup>-</sup>	std. state, hyp. m=1	aq			37.0				-18.
ClO <sub>3</sub>		g			-23.50	-0.62	0.454	39.0	
ClO <sub>3</sub> <sup>-</sup>	std. state, hyp. m=1	aq			-31.41	-2.57	1.884	43.5	
ClO <sub>4</sub> <sup>-</sup>	std. state, hyp. m=1	aq			18.20	22.40	-16.419	63.70	
Cl <sub>2</sub> O		g			9.2				
	in 400 H <sub>2</sub> O	aq			63.4				
Cl <sub>2</sub> O <sub>7</sub>		g							
		o		-22.019	-22.063	16.6696		44.617	6.96
	std. state, hyp. m=1	aq			-40.023	-31.350	22.9792	13.16	-30.0
	in 1 H <sub>2</sub> O	aq			-28.331				
HCl									

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 10-3. Chlorine (at. no., 17; at. wt., 35.457)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1947; March 31, 1950

Substance									
Formula	Description		State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
				At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole	cal/deg mole	
HCl	in	2 H <sub>2</sub> O	aq		-33.731				
		3 H <sub>2</sub> O	aq		-35.651				
		4 H <sub>2</sub> O	aq		-36.691				
		5 H <sub>2</sub> O	aq		-37.371				
		6 H <sub>2</sub> O	aq		-37.811				
		8 H <sub>2</sub> O	aq		-38.371				
		10 H <sub>2</sub> O	aq		-38.671				
		12 H <sub>2</sub> O	aq		-38.847				
		15 H <sub>2</sub> O	aq		-39.030				
		20 H <sub>2</sub> O	aq		-39.218				
		25 H <sub>2</sub> O	aq		-39.335				
		30 H <sub>2</sub> O	aq		-39.413				
		40 H <sub>2</sub> O	aq		-39.516				
		50 H <sub>2</sub> O	aq		-39.577				
		75 H <sub>2</sub> O	aq		-39.665				
		100 H <sub>2</sub> O	aq		-39.713				
		200 H <sub>2</sub> O	aq		-39.798				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 10-4. Chlorine (at. no., 17; at. wt., 35.457) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1947; June 30, 1947; March 31, 1950											
Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
Formula	in			At 0°K	At 298.16°K (25°C)						
				kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole		
HCl	in	300 H <sub>2</sub> O	aq		-39.837						
		400 H <sub>2</sub> O	aq		-39.859						
		500 H <sub>2</sub> O	aq		-39.874						
		700 H <sub>2</sub> O	aq		-39.895						
		1000 H <sub>2</sub> O	aq		-39.913						
		2000 H <sub>2</sub> O	aq		-39.946						
		3000 H <sub>2</sub> O	aq		-39.960						
		4000 H <sub>2</sub> O	aq		-39.968						
		5000 H <sub>2</sub> O	aq		-39.972						
		7000 H <sub>2</sub> O	aq		-39.982						
		10000 H <sub>2</sub> O	aq		-39.987						
		20000 H <sub>2</sub> O	aq		-39.998						
		50000 H <sub>2</sub> O	aq		-40.007						
		100000 H <sub>2</sub> O	aq		-40.012						
		$\infty$ H <sub>2</sub> O	aq		-40.023						
			in CCl <sub>4</sub>		-25.6						
			in C <sub>2</sub> H <sub>5</sub> OH		-32.6						
									-30.0		

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 10-5. Chlorine (at. no., 17; at. wt., 35.457)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1947; June 30, 1948; March 31, 1950

Formula	Substance	Description	State	At 0°K					At 298.16°K (25°C)		
				kcal/mole					kcal/mole		
									cal/deg mole		
HC10	in 400 H <sub>2</sub> O		aq								
HC10 <sub>2</sub>			aq								
HC10 <sub>3</sub>			aq								



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 10-6. Chlorine (at. no., 17; at. wt., 35.457)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1948; March 31, 1950

Substance		State	Description	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	At 298.16°K (25°C)									
	kcal/mole			kcal/mole	kcal/mole		kcal/mole	cal/deg mole	cal/deg mole	
HClO <sub>4</sub>		liq								
	std. state, hyp. m = 1	aq			-11.1					
	in 1 H <sub>2</sub> O	aq			-31.41					
	3 H <sub>2</sub> O	aq			-19.7					
	6 H <sub>2</sub> O	aq			-30.2					
	10 H <sub>2</sub> O	aq			-31.08					
	50 H <sub>2</sub> O	aq			-31.38					
	100 H <sub>2</sub> O	aq			-31.43					
	200 H <sub>2</sub> O	aq			-31.47					
	600 H <sub>2</sub> O	aq			-31.45					
HClO <sub>4</sub> ·H <sub>2</sub> O HClO <sub>4</sub> ·2H <sub>2</sub> O	∞ H <sub>2</sub> O	aq			-31.45					
		aq			-31.41					
		c			-92.1					
		liq			-162.8					
ClF ClF <sub>3</sub>		g			-13.3	-13.6	9.97	52.05	7.67	
		g			-37.0	-27.2	19.94	66.61	15.33	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 11-1. Bromine (at. no., 35; at. wt., 79.916)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1947; June 30, 1947; June 30, 1948; March 31, 1950

Formula	Substance	Description	State	At 0°K					At 298.16°K (25°C)		
				kcal/mole					kcal/mole		
				$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$
Br			g								
Br <sup>+</sup>			g								
Br <sup>++</sup>			g								
Br <sup>+++</sup>			g								
Br <sup>++++</sup>			g								
Br <sup>-</sup>			g								
Br <sub>2</sub>			aq								
		std. state, hyp. m = 1	g								
			liq								
			aq								
		in CCl <sub>4</sub> (tetrachloromethane)									
		in CHCl <sub>3</sub> (trichloromethane)									
		in CS <sub>2</sub> (carbon disulfide)									
Br <sub>2</sub> ·10H <sub>2</sub> O			c								
Br <sub>3</sub> <sup>-</sup>			aq								
Br <sub>5</sub> <sup>-</sup>			aq								
BrO <sup>-</sup>			aq								
BrO <sub>3</sub> <sup>-</sup>			aq								
		std. state, hyp. m = 1									

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 11-2. Bromine (at. no., 35; at. wt., 79.916)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1947; June 30, 1947; June 30, 1948

Substance								
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 0°K	At 298.16°K (25°C)				
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
HBr	g							
	std. state, hyp. m = 1	aq		-8.66	9.3269	47.437	6.96	
	in 1 H <sub>2</sub> O	aq		-28.90	18.0124	19.29	-30.7	
	2 H <sub>2</sub> O	aq		-18.56				
	3 H <sub>2</sub> O	aq		-22.18				
	4 H <sub>2</sub> O	aq		-24.36				
	5 H <sub>2</sub> O	aq		-25.54				
	6 H <sub>2</sub> O	aq		-26.27				
	8 H <sub>2</sub> O	aq		-26.78				
	10 H <sub>2</sub> O	aq		-27.39				
	12 H <sub>2</sub> O	aq		-27.72				
	15 H <sub>2</sub> O	aq		-27.90				
	20 H <sub>2</sub> O	aq		-28.08				
	25 H <sub>2</sub> O	aq		-28.25				
	30 H <sub>2</sub> O	aq		-28.34				
	40 H <sub>2</sub> O	aq		-28.41				
		aq		-28.48				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 11-3. Bromine (at. no., 35; at. wt., 79.916)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1947; March 31, 1950

Substance			State	Description	At 298.16°K (25°C)					
Formula	Formula	State			Description	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
						kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
HBr	100 H <sub>2</sub> O	aq			-28.54					
	75 H <sub>2</sub> O	aq			-28.61					
	100 H <sub>2</sub> O	aq			-28.65					
	200 H <sub>2</sub> O	aq			-28.71					
	300 H <sub>2</sub> O	aq			-28.74					
	400 H <sub>2</sub> O	aq			-28.76					
	500 H <sub>2</sub> O	aq			-28.77					
	700 H <sub>2</sub> O	aq			-28.78					
	1000 H <sub>2</sub> O	aq			-28.80					
	2000 H <sub>2</sub> O	aq			-28.83					
	3000 H <sub>2</sub> O	aq			-28.84					
	4000 H <sub>2</sub> O	aq			-28.85					
	5000 H <sub>2</sub> O	aq			-28.85					
	7000 H <sub>2</sub> O	aq			-28.86					
	10000 H <sub>2</sub> O	aq			-28.86					
	20000 H <sub>2</sub> O	aq			-28.87					
	50000 H <sub>2</sub> O	aq			-28.88					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 11-4. Bromine (at. no., 35; at. wt., 79.916)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1947; June 30, 1947; March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	At 298.16°K (25°C)				
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
HBr		in 100000 H <sub>2</sub> O	aq		-28.89				
		∞ H <sub>2</sub> O	aq		-28.90				
HBrO			aq		-9.5				
HBrO <sub>3</sub>			aq		3.51	-0.210	0.1539	57.34	
BrCl			g		-2.30				
			in CCl <sub>4</sub>						-30.7



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 12-1. Iodine (at. no., 53; at. wt., 126.92)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1947; June 30, 1947; June 30, 1948; March 31, 1950

Formula	Substance	Description	State	At 0°K					At 298.16°K (25°C)				
				kcal/mole					kcal/mole				
I			g	25.592	25.482	16.765	-12.2833	43.1641	4.9680				
I <sup>+</sup>			g	266.366	267.737								
I <sup>++</sup>			g	704.75	707.60								
I <sup>-</sup>			g	-48.6	-50.2								
I <sub>2</sub>		std. state, hyp. m = 1	aq	15.640	-13.37	-12.35	9.052	26.14	-21.7				
			g	0.000	14.876	4.33	-3.3937	64.260	8.81				
			c	0.000	0.000	0.000	11.181	27.9	13.14				
			aq		5.0								
		in C <sub>6</sub> H <sub>6</sub> (benzene)			6.0								
		in C <sub>2</sub> H <sub>5</sub> O (ethanol)			1.7								
		in C <sub>4</sub> H <sub>10</sub> O (diethyl ether)			1.5								
		in CCl <sub>4</sub> (tetrachloromethane)			5.99								
		in CHCl <sub>3</sub> (trichloromethane)			6.0								
		in CS <sub>2</sub> (carbon disulfide)			5.0								
I <sub>3</sub> <sup>-</sup>		std. state, hyp. m = 1	aq		-12.4	-12.31	31.62	41.7					
IO <sup>-</sup>			aq		-34.								
IO <sub>3</sub> <sup>-</sup>			aq		-55.0	-32.4	-3.75						
IO <sub>4</sub> <sup>-</sup>		std. state, hyp. m = 1	aq										
IO <sub>6</sub> <sup>-----</sup>			aq										
I <sub>2</sub> O <sub>5</sub>			c		-42.34								
3I <sub>2</sub> O <sub>5</sub> ·H <sub>2</sub> O			c		-201.12								

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

Washington, D.C.

National Bureau of Standards

Table 12-2. Iodine (at. no., 53; at. wt., 126.92)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1947; June 30, 1947; June 30, 1948

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	At 298.16°K (25°C)				
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
HI			g	6.73	6.20	0.31	-0.227	49.314	6.97
			liq						
			aq		-13.37				
			aq		-6.4	-12.35	9.052	26.14	-31.0
			aq		-8.72				
			aq		-11.29				
			aq		-12.48				
			aq		-12.90				
			aq		-12.97				
			aq		-13.04				
			aq		-13.10				
			aq		-13.18				
			aq		-13.22				
			aq		-13.24				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

<p>Table 12-3. Iodine (at. no., 53; at. wt., 126.92) HEAT OF FORMATION AT 0°K, HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1947; June 30, 1947; March 31, 1950</p>									
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
HI	in 400 H <sub>2</sub> O		aq		-13.25				
	500 H <sub>2</sub> O		aq		-13.26				
	1000 H <sub>2</sub> O		aq		-13.28				
	2000 H <sub>2</sub> O		aq		-13.30				
	$\infty$ H <sub>2</sub> O		aq		-13.37				-31.0
HIO HIO <sub>3</sub> HIO <sub>6</sub> H <sub>2</sub> IO <sub>6</sub> H <sub>3</sub> IO <sub>6</sub> H <sub>4</sub> IO <sub>6</sub> H <sub>5</sub> IO <sub>6</sub>			aq		-38.				
			c		-57.03				
			aq		-54.9				
			aq		-158.9				
			aq		-171.8				
			aq		-182.5				
			aq		-174.4				
			c		-184.4				
			aq		-183.0				
			c		-99.7				
I <sub>2</sub> O <sub>5</sub> ·HI O <sub>3</sub> ICl			g		4.20				
			liq			-1.32	0.965	59.12	8.46

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 12-4. Iodine (at. no., 53; at. wt., 126.92) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1947; June 30, 1947; March 31, 1950									
Substance		$\Delta H_f^\circ$	$\Delta E_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
Formula	Description	State	At 0°K			At 298.16°K (25°C)			cal/deg mole
			kcal/mole	kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
ICl	$\beta$	c							
	$\alpha$	c							
ICl <sub>3</sub>		in CCl <sub>4</sub>		-3.24					
		c		-21.1	-5.36	3.929	41.1		
IBr		g		9.75	0.91	-0.667	61.80	8.69	
		liq							
		in CCl <sub>4</sub>		1.72	-0.314	0.2302	39.0		

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 14-1. Sulfur (at. no., 16; at. wt., 32.066)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

September 30, 1947; March 31, 1950

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 298.16°K (25°C)				
			At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	cal/deg mole	cal/deg mole
S	$\lambda$	g		53.25	43.57	-31.9362	40.085
	$\mu$	liq		0.253			6.9
	rhombic	liq		0.67			6.9
	monoclinic	c, II	0.000	0.000	0.000	0.0000	7.62
	in 2 CS <sub>2</sub>	c, I		0.071	0.023	-0.0169	7.78
	6 CS <sub>2</sub>			0.402			5.65
S <sup>+</sup> S <sup>++</sup> S <sup>+++</sup> S <sup>++++</sup> S <sup>+++++</sup> S <sup>+++++</sup> S <sup>+++++</sup> S <sup>+++++</sup> S <sup>+++++</sup> S <sup>----</sup>		g		293.586			
		g		834.82			
		g		1644.56			
		g		2736.7			
		g		4409.5			
		g		6441.1			
		g		12923.			
		g		20507.			
		g		125.2			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 14-2. Sulfur (at. no., 16; at. wt., 32.066)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
September 30, 1947; March 31, 1950

Formula	Substance	State	At 298.16°K (25°C)					$C_p^0$
			$\Delta H_f^0$ kcal/mole	$\Delta H_f^0$ kcal/mole	$\Delta F_f^0$ kcal/mole	$\log_{10} K_f$	$S^0$ cal/deg mole	
S <sup>---</sup>	std. state, hyp. m = 1	aq						
S <sub>2</sub>		g						
S <sub>2</sub> <sup>---</sup>		aq		10.0	20.0	-14.660	5.3	
S <sub>3</sub> <sup>---</sup>		aq		29.86				
S <sub>4</sub> <sup>---</sup>		aq		9.3				
S <sub>4</sub> <sup>---</sup>		aq		7.8				
S <sub>6</sub>		g		5.3				
S <sub>8</sub>		g		25.3				
S <sub>0</sub>		g		24.1	12.78	-9.368	53.04	
				19.02				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

<p>Table 14-3. Sulfur (at. no., 16; at. wt., 32.066) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1947; March 31, 1950</p>									
Substance		Description	State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula									
SO <sub>2</sub>	in	200 H <sub>2</sub> O	g						
		300 H <sub>2</sub> O	aq						
		400 H <sub>2</sub> O	aq						
		500 H <sub>2</sub> O	aq						
		1000 H <sub>2</sub> O	aq						
		2000 H <sub>2</sub> O	aq						
		3000 H <sub>2</sub> O	aq						
		4000 H <sub>2</sub> O	aq						
		5000 H <sub>2</sub> O	aq						
		10000 H <sub>2</sub> O	aq						
SO <sub>2</sub> ·7H <sub>2</sub> O			c						
SO <sub>3</sub>			g						
		low-melting, wool-like	c, II						
		high-melting, wool-like	c, I						
			liq						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 14-4. Sulfur (at. no., 16; at. wt., 32.066) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1947; March 31, 1950										
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	At 0°K			At 298.16°K (25°C)						
	kcal/mole			kcal/mole	kcal/mole				cal/deg mole	cal/deg mole
$SO_3^{--}$		std. state, hyp. m = 1	aq		-149.2	-118.8	87.08	10.4		
$SO_4^{--}$		std. state, hyp. m = 1	aq		-216.90	-177.34	129.988	4.1		4.0
$S_2O_3^{--}$		std. state, hyp. m = 1	aq		-154.	-127.2	93.24	29.		
$S_2O_4^{--}$		std. state, hyp. m = 1	aq		-164.	-138.	101.2	57.		
$S_2O_5^{--}$			aq		-231.					
$S_2O_6^{--}$			aq		-280.4					
$S_2O_7$			c		-194.3					
$S_2O_8^{--}$			aq		-325.3					
$S_3O_6^{--}$			aq		-295.					
$S_4O_6^{--}$		std. state, hyp. m = 1	aq		-290.	-246.3	180.53	62.		
$S_5O_6^{--}$			aq		-281.					



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 14-5. Sul fur (at. no., 16; at. wt., 32.066) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1947; March 31, 1950									
Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	Description		At 0°K						At 298.16°K (25°C)
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole	
HS <sup>-</sup>	std. state, hyp. m=1	aq		-4.22	3.01	-2.206	14.6		
H <sub>2</sub> S		g		-4.815	-7.892	5.7847	49.15	8.12	
H <sub>2</sub> S·6H <sub>2</sub> O	std. state, hyp. m=1	aq		-9.4	-6.54	4.797	29.2		
		c		-431.2					
H <sub>2</sub> S <sub>2</sub>		liq		-5.5					
H <sub>2</sub> S <sub>3</sub>		liq							
H <sub>2</sub> S <sub>5</sub>		liq		+0.7					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 14-6. Sulfur (at. no., 16; at. wt., 32.066)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

September 30, 1947; March 31, 1950

Formula	Substance	State	Description	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	At 298.16°K (25°C)				
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
$\text{HSO}_3^-$	std. state, hyp. m=1	aq			-150.09	-126.03	92.378	31.64	
$\text{HSO}_4^-$	std. state, hyp. m=1	aq			-211.70	-179.94	131.893	30.32	
$\text{H}_2\text{SO}_3$	in 200 H <sub>2</sub> O	aq			-146.82				
	300 H <sub>2</sub> O	aq			-146.92				
	400 H <sub>2</sub> O	aq			-147.06				
	500 H <sub>2</sub> O	aq			-147.17				
	1000 H <sub>2</sub> O	aq			-147.59				
	2000 H <sub>2</sub> O	aq			-148.06				
	3000 H <sub>2</sub> O	aq			-148.37				
	4000 H <sub>2</sub> O	aq			-148.58				
	5000 H <sub>2</sub> O	aq			-148.73				
	10000 H <sub>2</sub> O	aq			-149.18				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 14-7. Sulfur (at. no. 16; at. wt., 32.066)

HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
September 30, 1947; March 31, 1950

Formula	Substance	Description	State					
				$\Delta H_f^\circ$		$\Delta F_f^\circ$		$\log_{10} K_f$
				At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole	
H <sub>2</sub> SO <sub>4</sub>	std. state, hyp. m = 1		liq					
	in 0.5 H <sub>2</sub> O		aq		-193.91			32.88
	1.0 H <sub>2</sub> O		aq		-216.90			4.0
	1.5 H <sub>2</sub> O		aq		-197.67			
	2 H <sub>2</sub> O		aq		-200.62			
	3 H <sub>2</sub> O		aq		-202.73			
	4 H <sub>2</sub> O		aq		-203.93			
	5 H <sub>2</sub> O		aq		-205.62			
	6 H <sub>2</sub> O		aq		-206.83			
	7 H <sub>2</sub> O		aq		-207.78			
	8 H <sub>2</sub> O		aq		-208.43			
	10 H <sub>2</sub> O		aq		-208.95			
	12 H <sub>2</sub> O		aq		-209.33			
	15 H <sub>2</sub> O		aq		-210.32			
	20 H <sub>2</sub> O		aq		-210.68			
	25 H <sub>2</sub> O		aq		-211.00			
						-177.34	129.968	4.1

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 14-8. Sulfur (at. no., 16; at. wt., 32.066) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1947; March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K					
				kcal/mole	kcal/mole			kcal/mole	cal/deg mole
$H_2SO_4$	30 $H_2O$		aq		-211.28				
	40 $H_2O$		aq		-211.38				
	50 $H_2O$		aq		-211.44				
	75 $H_2O$		aq		-211.52				
	100 $H_2O$		aq		-211.59				
	200 $H_2O$		aq		-211.82				
	300 $H_2O$		aq		-212.00				
	400 $H_2O$		aq		-212.13				
	500 $H_2O$		aq		-212.25				
	600 $H_2O$		aq		-212.35				
	700 $H_2O$		aq		-212.45				
	800 $H_2O$		aq		-212.54				
	900 $H_2O$		aq		-212.62				
	1000 $H_2O$		aq		-212.69				
	2000 $H_2O$		aq		-213.24				
	3000 $H_2O$		aq		-213.63				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 14-9. Sulfur (at. no., 16; at. wt., 32.066) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1947; March 31, 1950									
Substance				State	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	At 0°K kcal/mole							
$H_2S^{(14)}$	in	4000 $H_2O$	aq		-213.90				
		5000 $H_2O$	aq		-214.09				
		6000 $H_2O$	aq		-214.25				
		7000 $H_2O$	aq		-214.38				
		8000 $H_2O$	aq		-214.51				
		9000 $H_2O$	aq		-214.62				
		10000 $H_2O$	aq		-214.72				
		15000 $H_2O$	aq		-215.09				
		20000 $H_2O$	aq		-215.33				
		30000 $H_2O$	aq		-215.64				
		40000 $H_2O$	aq		-215.84				
		50000 $H_2O$	aq		-215.98				
		70000 $H_2O$	aq		-216.15				
		100000 $H_2O$	aq		-216.23				
		200000 $H_2O$	aq		-216.50				
		500000 $H_2O$	aq		-216.69				
		$\infty$ $H_2O$	aq		-216.90				4.10

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 14-10. Sulfur (at. no., 16; at. wt., 32.066) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1947; March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	At 298.16°K (25°C)				
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
$H_2SO_4$		in 5 $(C_2H_5)_2O$ 10 $(C_2H_5)_2O$ 20 $(C_2H_5)_2O$	liq liq liq		-205.29 -205.51 -208.27				
$H_2SO_4 \cdot H_2O$ $H_2SO_4 \cdot 2H_2O$ $H_2SO_4 \cdot 4H_2O$					-269.32				
$H_2SO_5$ $H_2S_2O_4$ $H_2S_2O_6$ $H_2S_2O_7$ $H_2S_2O_8$			c aq aq liq c		-164. -280.0 -302.7				
$H_2S_4O_6$		in 200 $H_2O$	aq aq		-319.3 -290.				
$SF_4$ $SF_6$ $S_2F_2$			g g g		-262. -237.		173.7	69.5	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 14-11. Sulfur (at. no., 16; at. wt., 32.066) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1947; March 31, 1950									
Formula	Substance	Description	State	At 298.16°K (25°C)					
				$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$G_p^\circ$	
SCl <sub>4</sub>			liq						
S <sub>2</sub> Cl <sub>2</sub>			liq	-13.6					30.
S <sub>2</sub> Cl <sub>4</sub>			liq	-14.4					
			liq	-24.0					
SOCl <sub>2</sub>			liq	-49.2					28.8
SO <sub>2</sub> Cl <sub>2</sub>			liq	-93.0					31.5
S <sub>2</sub> O <sub>5</sub> Cl <sub>2</sub>			liq	-166.6					55.5
HSO <sub>3</sub> Cl			liq	-142.7					
S <sub>2</sub> 8r <sub>2</sub>			liq	-3.6					
S08r <sub>2</sub>			liq						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 15-1. Selenium (at. no., 34; at. wt., 78.96)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
September 30, 1947; March 31, 1950

Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 298.16°K (25°C)								
	kcal/mole			kcal/mole	kcal/mole		cal/deg mole	cal/deg mole	
Se		gray, hexagonal	g	48.23	48.37	38.77	-28.418	42.21	4.968
		black, monoclinic	c, I	0.000	0.000	0.000	0.0000	10.0	5.95
		red, monoclinic	c, II						
		vitreous	c, III		0.2				
		precipitated	gls		1.05				6.13
			amorp		1.05				6.13
Se <sup>+</sup>			g	273.08	274.70				
Se <sup>++</sup>			g	769.19	772.29				
Se <sup>+++</sup>			g	1555.06	1559.64				
Se <sup>++++</sup>			g	2544.37	2550.43				
Se <sup>+++++</sup>			g	4230.26	4237.80				
Se <sup>+++++</sup>			g	6113.99	6123.01				
Se <sup>++++++</sup>			g	9949.81	9960.31				
Se <sup>-----</sup>		std. state, hyp. m=1	aq		31.6	37.2	-27.27	20.0	



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 15-2. Selenium (at. no., 34; at. wt., 78.96)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
September 30, 1947; March 31, 1950

Formula	Substance	State	Description	At 0°K					At 298.16°K (25°C)		
				$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole			
Se		g		33.61	33.14	-15.503	61.72	8.445			
Se <sub>6</sub>		g									
Se <sub>8</sub>		g		10.46	3.48						
Se <sub>16</sub>		c			-55.01						
SeO <sub>2</sub>		aq			-54.07						
SeO <sub>3</sub>		aq	std. state, hyp. m=1		-122.33	-89.34	51.3				
SeO <sub>4</sub>		aq	std. state, hyp. m=1		-145.3	-105.42	77.271				
SeO <sub>2</sub>		aq	std. state, hyp. m=1		24.6	23.57	42.3				
SeO <sub>3</sub>		aq	std. state, hyp. m=1		24.5	17.0	51.9				
SeO <sub>4</sub>		aq	std. state, hyp. m=1		14.1	18.4	39.4				
H <sub>2</sub> SeO <sub>3</sub>		aq	std. state, hyp. m=1		-123.5	-98.3	72.05				
H <sub>2</sub> SeO <sub>4</sub>		aq	std. state, hyp. m=1		-143.1	-108.2	79.31				
H <sub>2</sub> SeO <sub>3</sub>		aq	std. state, hyp. m=1		-126.5						
H <sub>2</sub> SeO <sub>4</sub>		aq	std. state, hyp. m=1		-122.39	-101.8	74.6				
H <sub>2</sub> SeO <sub>4</sub>		aq	std. state, hyp. m=1		-138.6						
H <sub>2</sub> SeO <sub>4</sub>		aq	std. state, hyp. m=1		-145.9	-105.42	77.271				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards SERIES I Washington, D.C.

Table 15-3. Selenium (at. no., 34; at. wt., 78.96) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1947; March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K					
				kcal/mole	kcal/mole			kcal/mole	cal/deg mole
$H_2SeO_4$	in	100 $H_2O$	aq		-141.5				
		200 $H_2O$	aq		-141.7				
		500 $H_2O$	aq		-142.2				
		1000 $H_2O$	aq		-142.6				
		1200 $H_2O$	aq		-142.7				
		2000 $H_2O$	aq		-142.9				
		5000 $H_2O$	aq		-143.8				
		$\infty$ $H_2O$	aq		-145.3				
			c		-203.0				
			c						
$SeF_6$			g		-246.	-222.	162.7	75.10	
$SeCl_2$			g		-9.7				
$SeCl_4$			c		-45.0				
$Se_2Cl_2$			liq		-20.0				
$Se(OH)_3ClO_4$			c		-149.0				
$SeO_2 \cdot 3SO_3$			c		-167.1				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 16-1. Tellurium (at. no., 52; at. wt., 127.61) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1947; September 30, 1948									
Substance		Description	State	$\Delta H_f^o$	$\Delta H_f^o$	$\Delta F_f^o$	$\log_{10} K_f$	$S^o$	$C_p^o$
Formula	At 298.16°K (25°C)								
	At 0°K kcal/mole			kcal/mole	kcal/mole		cal/deg mole	cal/deg mole	
Te	precipitated		g		47.6	38.1	-27.93	43.64	4.968
Te <sup>+</sup>			c,II	0.000	0.000	0.000	0.0000	11.88	6.15
Te <sup>++</sup>			amorp		2.7				
Te <sup>+++</sup>			g		256.80				
Te <sup>++++</sup>			g		755.09				
Te <sup>+++++</sup>			g		1462.50				
Te <sup>+++++</sup>			g		2336.08				
Te <sup>+++++</sup>			g		3727.50				
Te <sup>+++++</sup>			g		5396.9				
Te <sup>+++++</sup>			g		8562.3				
Te <sub>2</sub>			g		41.0	29.0	-21.26	64.07	8.72
TeO			g		43.0				
TeO <sub>2</sub>			c		-77.69	-64.60	47.353	16.99	15.89
TeO <sub>3</sub>			aq		-142.1				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 16-2. Tellurium (at. no., 52; at. wt., 127.61)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
September 30, 1947; March 31, 1950

Formula	Substance	State	Description	Substance				
				$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole
$\text{TeO}_4$		aq			-170.2			
$\text{H}_2\text{Te}$		g			36.9	33.1	-24.26	56.
$\text{H}_2\text{TeO}_3$		c			-144.7	-115.7	84.81	47.7
		aq			-144.7			
$\text{H}_2\text{TeO}_4$		aq			-166.7			
$\text{H}_2\text{TeO}_4 \cdot 2\text{H}_2\text{O}$		c			-306.6	-245.3	179.80	47.
$\text{TeF}_6$		g			-315.	-292.	214.0	80.67
$\text{TeCl}_4$		c			-77.2			
$\text{TeBr}_4$		c			-43.8			
$(\text{TeO}_2)_2\text{SO}_3$		c			-292.8			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 17-1. Polonium (at. no., 84; at. wt., 210)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
September 30, 1947; March 31, 1950

Formula	Substance	Description	State	At 0°K					At 298.16°K (25°C)		
				$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\Delta F^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole		
Po			c	0.000	0.000	0.000	0.0000				
Po <sup>++</sup>			aq			30.	-22.6				
PoO <sub>2</sub>			c			-46.	33.7				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 18-1. Nitrogen (at. no., 7; at. wt., 14.006)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$			
Formula	At 298.16°K (25°C)											
	At 0°K kcal/mole			kcal/mole	kcal/mole		cal/deg mole	cal/deg mole				
N			g	85.120	85.565	81.471	-59.71717	36.6147	4.9680			
N <sup>+</sup>				420.552	422.478							
N <sup>++</sup>				1103.30	1106.71							
N <sup>+++</sup>				2197.04	2201.93							
N <sup>++++</sup>				3983.18	3989.55							
N <sup>+++++</sup>				6240.07	6247.92							
N <sup>++++++</sup>				18968.5	18977.8							
N <sup>+++++++</sup>				34347.	34358.							
N <sub>2</sub>				0.000	0.000	0.000				0.0000	45.767	6.960
N <sub>2</sub> <sup>+</sup>				434.69	436.17							
N <sub>2</sub> <sup>++</sup>				1130.	1133.							
N <sub>3</sub> <sup>-</sup>					58.6							

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 18-2. Nitrogen (at. no., 7; at. wt., 14.008) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1950									
Formula	Substance	Description	State	ΔH <sup>fo</sup>		ΔF <sup>fo</sup>	log <sub>10</sub> K <sup>fo</sup>	S <sup>o</sup>	C <sub>p</sub> <sup>o</sup>
				At 0°K					
				kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
NO			g	21.477	21.600	20.719	-15.18688	50.339	7.137
NO <sub>2</sub>			g	8.682	8.091	12.390	-9.0817	57.47	9.06
NO <sub>2</sub> <sup>-</sup>		std. state, hyp. m = 1	aq		-25.4				
NO <sub>3</sub>			g		13.				
NO <sub>3</sub> <sup>-</sup>		std. state, hyp. m = 1	aq		-49.372	-26.41	19.358	35.0	
N <sub>2</sub> O			g	20.31	19.49	24.76	-18.149	52.58	9.251
N <sub>2</sub> O·6H <sub>2</sub> O			aq		-2.59	33.0	-24.19	6.6	
N <sub>2</sub> O <sub>2</sub> <sup>==</sup>			g		20.0				
N <sub>2</sub> O <sub>3</sub>			g	4.489	2.309	23.491	-17.2186	72.73	18.90
N <sub>2</sub> O <sub>4</sub>			q		3.6				
N <sub>2</sub> O <sub>5</sub>			c		-10.0				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 18-3. Nitrogen (at. no. 7; at. wt., 14.008)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance		$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K		At 298.16°K (25°C)		
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
NH		g	59.	59.			6.96
NH <sub>3</sub>		g	-9.368	-11.04	-3.976	46.01	8.523
	std. state, hyp. m = 1	aq		-19.32	-6.37	26.3	
	in 1 H <sub>2</sub> O	aq		-18.1			
	2 H <sub>2</sub> O	aq		-18.7			
	3 H <sub>2</sub> O	aq		-18.87			
	4 H <sub>2</sub> O	aq		-18.99			
	5 H <sub>2</sub> O	aq		-19.07			
	10 H <sub>2</sub> O	aq		-19.23			
	20 H <sub>2</sub> O	aq		-19.27			
	30 H <sub>2</sub> O	aq		-19.28			
	40 H <sub>2</sub> O	aq		-19.28			
	50 H <sub>2</sub> O	aq		-19.27			
	100 H <sub>2</sub> O	aq		-19.30			
	200 H <sub>2</sub> O	aq		-19.32			
	$\infty$ H <sub>2</sub> O	aq		-19.32			



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

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Table 18-4. Nitrogen (at. no., 7; at. wt., 14.008) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1950										
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula				At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole		
$\text{NH}_4^+$			g		150.					
$\text{N}_2\text{H}_4$		std. state, hyp. m=1	aq		-31.74	-19.00	13.927	26.97		
		hydrazine	liq		12.05					
		in 300 H <sub>2</sub> O	aq		8.16					
$\text{N}_2\text{H}_4 \cdot \text{H}_2\text{O}$			liq		-57.95					
$\text{N}_2\text{H}_4\text{H}^+$			aq		-1.7					
$\text{N}_2\text{H}_4\text{H}_2^{++}$			aq		-4.0					
$\text{HN}_3$		hydrogen azide	g		70.3	78.5	-57.54	56.74	10.12	
		in 100 H <sub>2</sub> O	aq		61.51					
		200 H <sub>2</sub> O	aq		60.84					
		300 H <sub>2</sub> O	aq		60.61					
		400 H <sub>2</sub> O	aq		60.50					
$\text{NH}_4\text{N}_3$			c		20.4					
$\text{N}_2\text{H}_4\text{N}_3\text{H}$		hydrazine hydrazide	aq		26.90					
			c							

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 13-5. Nitrogen (at. no., 7; at. wt., 14.008)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1947; March 31, 1950

Substance		State	At 298.16°K (25°C)				S°	C <sub>p</sub> <sup>o</sup>
Formula	Description		ΔH <sup>fo</sup>	ΔH <sup>ro</sup>	ΔF <sup>ro</sup>	log <sub>10</sub> K <sub>f</sub>		
			At 0°K kcal/mole	kcal/mole	kcal/mole			
HNO <sub>2</sub>	std. state, hyp. m=1	aq			-28.4			
		g			-41.404		63.62	
HNO <sub>3</sub>	std. state, hyp. m=1	liq			-49.372	14.0000	37.19	26.26
	in	aq			-44.538	19.358	35.0	
	1 H <sub>2</sub> O	aq			-46.204			
	2 H <sub>2</sub> O	aq			-47.212			
	3 H <sub>2</sub> O	aq			-47.852			
	4 H <sub>2</sub> O	aq			-48.270			
	5 H <sub>2</sub> O	aq			-48.535			
	6 H <sub>2</sub> O	aq			-48.843			
	8 H <sub>2</sub> O	aq			-49.014			
	10 H <sub>2</sub> O	aq			-49.162			
	15 H <sub>2</sub> O	aq			-49.212			
	20 H <sub>2</sub> O	aq			-49.230			
	25 H <sub>2</sub> O	aq			-49.234			
	30 H <sub>2</sub> O	aq			-49.232			
	40 H <sub>2</sub> O	aq						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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<p>Table 18-6. Nitrogen (at. no., 7; at. wt., 14.008) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1950</p>									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K	At 298.16°K (25°C)				
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
HNO <sub>3</sub>	in 50 H <sub>2</sub> O		aq		-49.230				
	75 H <sub>2</sub> O		aq		-49.229				
	100 H <sub>2</sub> O		aq		-49.231				
	200 H <sub>2</sub> O		aq		-49.244				
	300 H <sub>2</sub> O		aq		-49.255				
	500 H <sub>2</sub> O		aq		-49.267				
	700 H <sub>2</sub> O		aq		-49.277				
	1000 H <sub>2</sub> O		aq		-49.286				
	2000 H <sub>2</sub> O		aq		-49.303				
	5000 H <sub>2</sub> O		aq		-49.323				
	10000 H <sub>2</sub> O		aq		-49.336				
	50000 H <sub>2</sub> O		aq		-49.355				
	$\infty$ H <sub>2</sub> O		aq		-49.372				
	in (C <sub>2</sub> H <sub>5</sub> ) <sub>2</sub> O								

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 18-7. Nitrogen (at. no., 7; at. wt., 14.008) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1950									
Formula	Substance	Description	State	$\Delta H^{\circ}$		$\Delta F^{\circ}$	$\log_{10} f^{\circ}$	$S^{\circ}$	$C_p^{\circ}$
				At 0°K	At 298.16°K (25°C)				
$\text{HNO}_3 \cdot \text{H}_2\text{O}$			liq						
$\text{HNO}_3 \cdot 3\text{H}_2\text{O}$			liq						
$\text{NH}_2\text{OH}$			c						
$\text{NH}_2\text{OH} \cdot \text{H}^+$			aq						
$\text{NH}_2\text{OH} \cdot \text{H}_2\text{O}$			aq						
$\text{NH}_4\text{OH}$			aq						
	std. state, hyp. m = 1								
	in								
	2 $\text{H}_2\text{O}$								
	3 $\text{H}_2\text{O}$								
	4 $\text{H}_2\text{O}$								
	5 $\text{H}_2\text{O}$								
	10 $\text{H}_2\text{O}$								
	20 $\text{H}_2\text{O}$								
	30 $\text{H}_2\text{O}$								
	40 $\text{H}_2\text{O}$								
	50 $\text{H}_2\text{O}$								
	100 $\text{H}_2\text{O}$								

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

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Table 18-8. Nitrogen (at. no., 7; at. wt., 14.006)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance						
Formula	Description	State	$\Delta H_f^\circ$	$\Delta G_f^\circ$	$\log_{10} K_f$	$S^\circ$
			At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole
NH <sub>4</sub> OH	in 200 H <sub>2</sub> O	aq		-87.64		
	$\infty$ H <sub>2</sub> O	aq		-87.64		
HN <sub>2</sub> O <sub>2</sub> <sup>-</sup> H <sub>2</sub> N <sub>2</sub> O <sub>2</sub>	std. state, hyp. m = 1	aq		-9.4	18.0	34.
	std. state, hyp. m = 1	aq		-13.7	8.4	52.
NH <sub>4</sub> NO <sub>2</sub>		c		-63.1		
	std. state, hyp. m = 1	aq		-57.1		
NH <sub>4</sub> NO <sub>3</sub>		c, IV		-87.27		43.5
	std. state, hyp. m = 1	aq		-81.11		
	in 3 H <sub>2</sub> O	aq		-83.30		
	4 H <sub>2</sub> O	aq		-83.20		
	5 H <sub>2</sub> O	aq		-83.03		
	10 H <sub>2</sub> O	aq		-82.46		
	20 H <sub>2</sub> O	aq		-81.95		
	30 H <sub>2</sub> O	aq		-81.72		
	40 H <sub>2</sub> O	aq		-81.58		
	50 H <sub>2</sub> O	aq		-81.49		
	100 H <sub>2</sub> O	aq		-81.30		

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 18-9. Nitrogen (at. no., 7, at. wt., 14.003)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} f^*$	$S^\circ$	$C_p^\circ$
Formula	At 0°K			At 298.16°K (25°C)					
	kcal/mole			kcal/mole	kcal/mole		kcal/mole	cal/deg mole	
$\text{NH}_4\text{NO}_3$	in 100 $\text{H}_2\text{O}$		aq		-81.19				
	500 $\text{H}_2\text{O}$		aq		-81.12				
	1000 $\text{H}_2\text{O}$		aq		-81.10				
	$\infty$ $\text{H}_2\text{O}$		aq		-81.11				
$\text{NH}_4\text{NO}_3 \cdot 1\frac{1}{2}\text{NH}_3$			liq		-93.6				
$\text{NH}_2\text{OH} \cdot \text{HNO}_3$			c		-86.3				
$\text{N}_2\text{H}_4 \cdot \text{HNO}_3$	stable		aq		-80.5				
$\text{N}_2\text{H}_4 \cdot 2\text{HNO}_3$			c, l		-51.1				
			aq		-102.7				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 19-10. Nitrogen (at. no., 7; at. wt., 14.008)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
NF <sub>3</sub>			g		-27.2				
NOF			g						
N <sub>2</sub> O			g						
NO <sub>2</sub> F			g						
NH <sub>4</sub> F			g						
NH <sub>4</sub> F			g						
NH <sub>4</sub> F			c		-114.6				
	in 200 H <sub>2</sub> O		aq		-110.14				
	in H <sub>2</sub> O		aq		-110.40				
NH <sub>4</sub> F·HF			c						
N <sub>2</sub> H <sub>4</sub> ·2HF			c						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 18-11. Nitrogen (at. no., 7; at. wt., 14.005)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1949

Substance						
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$
			At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole		cal/deg mole
NCI <sub>3</sub>	in CCl <sub>4</sub>	liq				
NOCl		g	13.25	54.7		
NO <sub>2</sub> Cl		g		12.57		
NH <sub>2</sub> Cl		g			-11.625	63.0
NH <sub>4</sub> Cl	std. state, hyp. m = 1 in 10 H <sub>2</sub> O 25 H <sub>2</sub> O 50 H <sub>2</sub> O 100 H <sub>2</sub> O 200 H <sub>2</sub> O 300 H <sub>2</sub> O 400 H <sub>2</sub> O 500 H <sub>2</sub> O 1000 H <sub>2</sub> O 2000 H <sub>2</sub> O	c, II				
		aq		-75.38		
		aq		-71.76		
		aq		-71.57		
		aq		-71.597		
		aq		-71.624		
		aq		-71.628		
		aq		-71.646		
		aq		-71.656		
		aq		-71.664		
		aq		-71.671		
		aq		-71.691		
		aq		-71.710		
				-48.73	35.718	22.6
						20.1



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 18-12. Nitrogen (at. no., 7; at. wt., 14.008)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F^\circ$	$\log_{10} f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)			
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
NH <sub>4</sub> Cl	in 3000 H <sub>2</sub> O	aq		-71.718			
	4000 H <sub>2</sub> O	aq		-71.724			
	5000 H <sub>2</sub> O	aq		-71.728			
	10000 H <sub>2</sub> O	aq		-71.737			
	$\infty$ H <sub>2</sub> O	aq		-71.76			
NH <sub>4</sub> Cl · 3NH <sub>3</sub> NH <sub>4</sub> Cl · 6NH <sub>3</sub>		liq		-133.9			
		liq		-183.9			
N <sub>2</sub> H <sub>4</sub> · HCl		c		-46.93			
	in 50 H <sub>2</sub> O	aq		-41.775			
	100 H <sub>2</sub> O	aq		-41.73			
	200 H <sub>2</sub> O	aq		-41.712			
	300 H <sub>2</sub> O	aq		-41.711			
	400 H <sub>2</sub> O	aq		-41.710			
	500 H <sub>2</sub> O	aq		-41.710			
	1000 H <sub>2</sub> O	aq		-41.708			

## SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

Table 13-13. Nitrogen (at. no., 7; at. wt., 14.008)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	At 298.16°K (25°C) kcal/mole		cal/deg mole	cal/deg mole
$N_2H_4 \cdot 2HCl$			c		-90.0				
			aq		-84.0				
$NH_2OH \cdot HCl$			c		-74.0				
			aq		-70.7				22
$NH_4ClO_4$			c, II		-69.42				
			aq		-63.2				
			aq		-63.15				
		in 500 $H_2O$ $\infty$ $H_2O$							

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Table 18-14. Nitrogen (at. no., 7; at. wt., 14.008)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance									
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 0°K	At 298.16°K (25°C)					
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole			cal/deg mole
NBr <sub>3</sub>		g		19.56	19.70	-14.440	65.16		
NOBr		c, II		-64.61					
NH <sub>4</sub> Br	in 200 H <sub>2</sub> O	aq		-60.51					
	∞ H <sub>2</sub> O	aq		-60.74					
NH <sub>4</sub> Br·NH <sub>3</sub>		liq		-63.6					
NH <sub>4</sub> Br·3NH <sub>3</sub>		liq		-122.3					
NH <sub>4</sub> Br·6NH <sub>3</sub>		liq		-172.0					
N <sub>2</sub> H <sub>4</sub> Br		c							
N <sub>2</sub> H <sub>4</sub> ·2HBr		c							

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Table 18-15. Nitrogen (at. no., 7; at. wt., 14.008)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance		Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K			At 298.16°K (25°C)		cal/deg mole	cal/deg mole	
	kcal/mole			kcal/mole	kcal/mole			kcal/mole
$\text{NL}_3$			c					
$\text{NH}_4\text{I}$			c, l					
		in 200 $\text{H}_2\text{O}$	aq	-48.30				
		$\infty$ $\text{H}_2\text{O}$	aq	-44.98				
			liq	-45.11				
$\text{NH}_4\text{I} \cdot \text{NH}_3$			liq	-69.1				
$\text{NH}_4\text{I} \cdot 3\text{NH}_3$			liq	-104.5				
$\text{NH}_4\text{I} \cdot 4\text{NH}_3$			liq					
$\text{NH}_4\text{I} \cdot 6\text{NH}_3$			liq	-154.7				
$\text{N}_2\text{H}_4 \cdot \text{HI}$			c					
$\text{N}_2\text{H}_4 \cdot 2\text{HI}$			c					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 18-16. Nitrogen (at. no., 7; at. wt., 14.008) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1950						
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$
				At 0°K	At 298.16°K (25°C)	
				kcal/mole	kcal/mole	cal/deg mole
N <sub>2</sub>			g			
N <sub>2</sub> S <sub>5</sub>			liq			
N <sub>4</sub> S <sub>4</sub>			c		127.6	
N <sub>2</sub> O <sub>3</sub> ·2SO <sub>3</sub>			c			
NH <sub>4</sub> HS			c		-33.10	
(NH <sub>4</sub> ) <sub>2</sub> S			aq		-35.1	
(NH <sub>4</sub> ) <sub>2</sub> S <sub>4</sub>			aq		-54.5	
(NH <sub>4</sub> ) <sub>2</sub> S <sub>5</sub>			c		-67.4	
			aq		-60.0	
			c		-66.8	
			aq		-61.2	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 19-17. Nitrogen (at. no., 7; at. wt., 14.008)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance		$\Delta H_f^\circ$	$\Delta G_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 298.16°K (25°C)			
			kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
NO <sub>2</sub> SO <sub>2</sub> OH NH <sub>4</sub> HSO <sub>3</sub>	std. state, hyp. m = 1	c				
		c		-183.8		
		aq		-181.83		
		aq		-181.5		
NH <sub>4</sub> HSO <sub>4</sub>	in 10 H <sub>2</sub> O 20 H <sub>2</sub> O 50 H <sub>2</sub> O 100 H <sub>2</sub> O 200 H <sub>2</sub> O 400 H <sub>2</sub> O 800 H <sub>2</sub> O	c		-244.83		34.2
		aq		-244.34		
		aq		-244.76		
		aq		-244.93		
		aq		-245.05		
		aq		-245.26		
		aq		-245.53		
		aq		-245.86		

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 18-18. Nitrogen (at. no., 7; at. wt., 14.008) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1950									
Formula	Substance	Description	State	At 0°K					
				At 298.16°K (25°C)					
				$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
$\text{NH}_4\text{OH} \cdot \text{H}_2\text{SO}_4$			c		-245.1				
$\text{N}_2\text{H}_4 \cdot \text{H}_2\text{SO}_4$			aq		-244.4				
			c		-231.6				
		in 1000 $\text{H}_2\text{O}$	aq		-223.44				
$(\text{NH}_4)_2\text{SO}_3$			c		-212.0				
			aq		-211.3				
$(\text{NH}_4)_2\text{SO}_3 \cdot \text{H}_2\text{O}$			c		-284.22				
$(\text{NH}_4)_2\text{SO}_4$			c		-281.86	-215.19	157.732	52.65	44.81
		in 20 $\text{H}_2\text{O}$	aq		-280.47				
		30 $\text{H}_2\text{O}$	aq		-280.40				
		40 $\text{H}_2\text{O}$	aq		-280.35				
		50 $\text{H}_2\text{O}$	aq		-280.31				
		100 $\text{H}_2\text{O}$	aq		-280.212				
		200 $\text{H}_2\text{O}$	aq		-280.150				
		300 $\text{H}_2\text{O}$	aq		-280.135				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 18-19. Nitrogen (at. no., 7; at. wt., 14.008) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole		cal/deg mole	cal/deg mole	
(NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub>	in	400 H <sub>2</sub> O	aq		-280.124				
		500 H <sub>2</sub> O	aq		-280.116				
		1000 H <sub>2</sub> O	aq		-280.122				
		5000 H <sub>2</sub> O	aq		-280.196				
		∞ H <sub>2</sub> O	aq		-280.38				
(NH <sub>2</sub> OH) <sub>2</sub> *H <sub>2</sub> SO <sub>4</sub>			c		-282.5				
			aq		-276.7				



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 18-20. Nitrogen (at. no., 7; at. wt., 14.008)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1947; March 31, 1950

Formula	Substance	Description	State	$\Delta f^{\circ}$	$\Delta f^{\circ}$	$\Delta f^{\circ}$	$\log_{10} K^{\circ}$	$S^{\circ}$	$C_p^{\circ}$
				At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
$(\text{NH}_4)_2\text{S}_2\text{O}_5$	std. state, hyp. m=1		c			-295.3			
$(\text{NH}_4)_2\text{S}_2\text{O}_8$			aq			-396.4			
$\text{NH}(\text{SO}_2\text{NH}_2)_2$			c			-387.8			
$(\text{NH}_4)_2\text{H}_2\text{SO}_4$	std. state, hyp. m=1		aq			-217.8			
$\text{NOSO}_3\text{F}$			c						
$\text{NSe}$			g						
$\text{N}_4\text{Se}_4$	std. state, hyp. m=1		c			169.2			
$\text{NH}_4\text{HSe}$			aq			-5.3			
$(\text{NH}_4)_2\text{Se}$			aq			-5.6			
	std. state, hyp. m=1		aq			-26.2			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 19-1. Phosphorus (at. no., 15; at. wt., 30.98) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1949									
Substance		Description	State	At 298.16°K (25°C)					
				$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
P	white red black		g		75.13	66.71	-48.897	38.98	4.9680
			c, III		0.000	0.000	0.0000	10.6	5.55
			c, II		-4.4				
			c, I		-10.3				
P <sup>+</sup> P <sup>++</sup> P <sup>+++</sup> P <sup>++++</sup> P <sup>+++++</sup>			g		329.81				
			g		784.51				
			g		1481.43				
			g		2667.2				
			g		4167.9				
P <sub>2</sub>			g		33.82	24.60	-18.031	52.13	7.63
P <sub>4</sub>			g		13.12	5.82	-4.266	66.90	16.0
			liq		0.601				24.2

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 19-2. Phosphorus (at. no., 15; at. wt. 30.98) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°K) December 31, 1947; March 31, 1949									
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	At 298.16°K (25°K)				
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
P <sub>0</sub>			g						
P <sub>03</sub>		std. state, hyp. m=1	aq		-9.7				
P <sub>04</sub>		std. state, hyp. m=1	aq		-235.6				
P <sub>2</sub> O <sub>7</sub>		std. state, hyp. m=1	aq		-306.9	-245.1	179.65	-52.	
P <sub>2</sub> O <sub>5</sub>		std. state, hyp. m=1	aq		-543.9				
P <sub>4</sub> O <sub>6</sub>			liq						
P <sub>4</sub> O <sub>10</sub>			c, II		-720.0				
			amorp		-734.0				
PH <sub>3</sub>			g		2.21	4.36	-3.196	50.2	
PH <sub>3</sub> ·6H <sub>2</sub> O			c		-422.7				
P <sub>2</sub> H <sub>4</sub>			c		-14.5				
P <sub>2</sub> H <sub>6</sub>			liq						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 19-3. Phosphorus (at. no., 15; at. wt., 30.96) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1950									
Substance		Description	State	$\Delta H_f^{\circ}$	$\Delta F_f^{\circ}$	$\log_{10} K_f$	$S^{\circ}$	$C_p^{\circ}$	
Formula				At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole	cal/deg mole	
HP0 <sub>3</sub>			c		-228.2				
HP0 <sub>3</sub> <sup>---</sup>			aq		-234.9				
HP0 <sub>4</sub> <sup>---</sup>			aq		-233.0				
H <sub>2</sub> P0 <sub>3</sub> <sup>-</sup>		std. state, hyp. m = 1	aq		-310.4	191.68	-8.6		
H <sub>2</sub> P0 <sub>4</sub> <sup>-</sup>			aq		-233.3				
		std. state, hyp. m = 1	aq		-311.3	198.86	21.3		
H <sub>3</sub> P0 <sub>2</sub>			c		-145.5				
			liq		-143.2				
			aq		-145.6				
H <sub>3</sub> P0 <sub>3</sub>			c		-232.2				
			liq		-229.1				
			aq		-232.2				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 19-4. Phosphorus (at. no., 15; at. wt., 30.98) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1950									
Substance		State	Description	ΔH <sup>o</sup> kcal/mole	ΔF <sup>o</sup> kcal/mole	log <sub>10</sub> K <sub>f</sub>	S <sup>o</sup> cal/deg mole	C <sub>p</sub> <sup>o</sup> cal/deg mole	
Formula									
H <sub>3</sub> PO <sub>4</sub>	in	c							
	1 H <sub>2</sub> O	aq		-306.2					
	2 H <sub>2</sub> O	aq		-305.6					
	3 H <sub>2</sub> O	aq		-306.4					
	4 H <sub>2</sub> O	aq		-307.2					
	5 H <sub>2</sub> O	aq		-307.7					
	10 H <sub>2</sub> O	aq		-308.0					
	20 H <sub>2</sub> O	aq		-308.56					
	30 H <sub>2</sub> O	aq		-308.82					
	40 H <sub>2</sub> O	aq		-308.97					
	50 H <sub>2</sub> O	aq		-309.04					
	100 H <sub>2</sub> O	aq		-309.06					
	200 H <sub>2</sub> O	aq		-309.18					
	300 H <sub>2</sub> O	aq		-309.28					
400 H <sub>2</sub> O	aq		-309.30						
				-309.32					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 19-5. Phosphorus (at. no., 15; at. wt., 30.96)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance		State	Description	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula									
$H_3PO_4$	in 500 $H_2O$	aq							
	1000 $H_2O$	aq			-309.34				
	2000 $H_2O$	aq			-309.36				
	3000 $H_2O$	aq			-309.39				
$H_3PO_4 \cdot 3H_2O$		c			-309.4				
		liq			-342.9				
$HP_2O_7^{---}$		aq			-339.3				
		aq			-546.7				
		aq			-547.0				
$H_4P_2O_7^{--}$		aq			-546.7				
		aq			-391.2				
		c							
$H_4P_2O_6$		c							
		c							

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 19-6. Phosphorus (at. no., 15; at. wt., 30.98) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1950							
Formula	Substance	Description	State	At 298.16°K (25°C)			
				$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole
$H_4P_2O_7$			c				
$H_4P_2O_7 \cdot \frac{1}{2}H_2O$			aq	-538.0			
			c	-545.9			
			liq	-643.9			
				-640.8			
$H_6P_4O_{13}$			c				
$PF_3$			g				64.12
$PF_5$			g				
$POF_3$			g				
$HPO_2F_2$			liq				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 19-7. Phosphorus (at. no., 15; at. wt., 30.98)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1947; March 31, 1950

Formula	Substance	Description	State	At 0°K					At 298.16°K (25°C)		
				kcal/mole					kcal/mole		
				$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
PCl <sub>3</sub>			liq		-81.0	-68.42	50.151	74.49			
			g		-73.22						
PCl <sub>5</sub>			c		-110.7	-77.59	56.872	84.3			
			g		-95.35						
P <sub>2</sub> Cl <sub>4</sub>			liq								
POCl <sub>3</sub>			liq		-151.0	-130.3	95.508	77.59			
			g		-141.5						
PH <sub>4</sub> Cl			g		-36.2						
PF <sub>2</sub> Cl			g								
PFC1 <sub>2</sub>			g								
POF <sub>2</sub> Cl			g								
POFC1 <sub>2</sub>			liq								



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 19-8. Phosphorus (at. no., 15; at. wt., 30.98)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°K)  
December 31, 1947; March 31, 1950

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°K)				
			kcal/mole	kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
PBr <sub>3</sub>		g liq in CS <sub>2</sub>		-35.9 -47.5 -46.9	-41.2	30.20	83.11	
PBr <sub>5</sub>		c		-66.				
POBr <sub>3</sub>		c		-114.6				
PH <sub>4</sub> Br		c		-29.5				
PF <sub>2</sub> Br		g						
PFBr <sub>2</sub>		liq						
PF <sub>3</sub> Br <sub>2</sub>		liq						
POF <sub>2</sub> Br		liq						
POFBr <sub>2</sub>		liq						
POCl <sub>2</sub> Br		liq						
POClBr <sub>2</sub>		c						
POFClBr		liq						

## National Bureau of Standards

## Washington, D.C.

Table 19-9. Phosphorus (at. no., 15; at. wt., 30.98)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance							
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K^*$	$S^\circ$	$C_p^\circ$
			At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	
At 298.16°K (25°C)							
			kcal/mole	kcal/mole			
Pt <sub>3</sub>		c		-10.9			
P <sub>2</sub> I <sub>4</sub>		c		-19.76			
PH <sub>4</sub> I		c		-15.8			26.5
P <sub>2</sub> S <sub>3</sub>		c					
P <sub>4</sub> S <sub>3</sub>		c					
P <sub>4</sub> S <sub>7</sub>		c					
P <sub>4</sub> S <sub>10</sub>		c					
P <sub>4</sub> S <sub>6</sub> O <sub>4</sub>		c					
PSF <sub>3</sub>		g				79.34	
PSCl <sub>3</sub>		g					
PSF <sub>2</sub> Cl		g					
PSFCl <sub>2</sub>		liq					
PSBr <sub>3</sub>		c					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 19-10. Phosphorus (at. no., 15; at. wt., 30.98) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1949							
Substance							
Formula	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$
			At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole			
PSF <sub>2</sub> Br		liq					
PSFBr <sub>2</sub>		liq					
PSCl <sub>2</sub> Br		liq					
PSClBr <sub>2</sub>		liq					
PN		g					
1/n (PN) <sub>n</sub>		c					
P <sub>3</sub> N <sub>5</sub>		c					
NH <sub>2</sub> OH·H <sub>3</sub> PO <sub>2</sub>		c	-20.2	-25.3	18.54	50.45	7.097
NH <sub>4</sub> H <sub>2</sub> PO <sub>3</sub>		c	-16.8				35.6
			-75.7				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 19-11. Phosphorus (at. no., 15; at. wt., 30.98) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	kcal/mole			At 298.16°K (25°C) cal/deg mole	cal/deg mole
$\text{NH}_4\text{H}_2\text{PO}_4$	in 20 H <sub>2</sub> O		c						
	25 H <sub>2</sub> O		aq		-346.75		212.903	36.32	34.0
	30 H <sub>2</sub> O		aq		-343.07				
	35 H <sub>2</sub> O		aq		-343.06				
	40 H <sub>2</sub> O		aq		-343.05				
	45 H <sub>2</sub> O		aq		-343.03				
	50 H <sub>2</sub> O		aq		-343.01				
	100 H <sub>2</sub> O		aq		-342.97				
	200 H <sub>2</sub> O		aq		-342.95				
	300 H <sub>2</sub> O		aq		-342.93				
	400 H <sub>2</sub> O		aq		-342.92				
	500 H <sub>2</sub> O		aq		-342.91				
	600 H <sub>2</sub> O		aq		-342.9				
	700 H <sub>2</sub> O		aq		-342.89				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 19-12. Phosphorus (at. no., 15; at. wt., 30.98) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula				At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
$N_2H_4 \cdot H_3PO_3$ $NH_4NH_2HPO_2$ $N_2H_4 \cdot H_3PO_4$  $(NH_4)_2HPO_4$			c					43.5	
			c						
			c						
			c						
	in 11 H <sub>2</sub> O		aq	-376.12					
	12 H <sub>2</sub> O		aq	-374.07					
	13 H <sub>2</sub> O		aq	-374.06					
	14 H <sub>2</sub> O		aq	-374.05					
	15 H <sub>2</sub> O		aq	-374.04					
	20 H <sub>2</sub> O		aq	-374.03					
	30 H <sub>2</sub> O		aq	-373.96					
	40 H <sub>2</sub> O		aq	-373.82					
50 H <sub>2</sub> O		aq	-373.71						
				-373.63					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 19-13. Phosphorus (at. no., 15; at. wt., 30.98)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$				
Formula	At 0°K			At 298.16°K (25°C)									
				kcal/mole	kcal/mole					kcal/mole	cal/deg mole	cal/deg mole	
(NH <sub>4</sub> ) <sub>2</sub> HPO <sub>4</sub>	in 100 H <sub>2</sub> O		aq		-373.39								
	20 H <sub>2</sub> O		aq		-373.21								
	300 H <sub>2</sub> O		aq		-373.12								
	400 H <sub>2</sub> O		aq		-373.07								
	500 H <sub>2</sub> O		aq		-373.04								
	600 H <sub>2</sub> O		aq		-373.0								
	700 H <sub>2</sub> O		aq		-372.99								
(NH <sub>4</sub> ) <sub>3</sub> PO <sub>4</sub>			c		-401.8								
	in 66 H <sub>2</sub> O		aq		-384.0								
(NH <sub>4</sub> ) <sub>3</sub> PO <sub>4</sub> · 3H <sub>2</sub> O			c		-612.8								

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 20-1. Arsenic (at. no., 33; at. wt., 74.91)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	At 298.16°K (25°C) cal/deg mole	At 298.16°K (25°C) cal/deg mole	At 298.16°K (25°C) cal/deg mole
As		$\alpha$ , gray, metallic	g		60.64	-37.192	41.62	4.968
		$\beta$	c	0.000	0.000	0.0000	8.4	5.97
		$\gamma$ , yellow, cubic	amorp		1.0			
			c		3.53			
As <sup>+</sup>			g		304.24			
As <sup>++</sup>			g		771.03			
As <sup>+++</sup>			g		1402.0			
As <sup>++++</sup>			g		2559.4			
As <sup>+++++</sup>			g		4004.8			
As <sup>++++++</sup>			g		6948.			
As <sub>2</sub>			g		29.6	-12.83	57.3	8.366
As <sub>4</sub>			g		35.7	-18.47	69.	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards Washington, D.C.

SERIES I

Table 20-2. Arsenic (at. no., 33; at. wt., 74.91)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1949

Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
			At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				
AsO		g						
AsO <sup>+</sup>	std. state, hyp. m=1	aq		4.79	-39.1	28.69		
AsO <sub>2</sub> <sup>-</sup>	std. state, hyp. m=1	aq			-83.7	61.33		
AsO <sub>4</sub> <sup>3-</sup>	std. state, hyp. m=1	aq		-208.	-152.	111.4	-34.6	
As <sub>2</sub> O <sub>5</sub>		c			-184.6	135.41	25.2	27.85
As <sub>2</sub> O <sub>5</sub> ·4H <sub>2</sub> O		aq		-218.6				
3As <sub>2</sub> O <sub>5</sub> ·5H <sub>2</sub> O		c		-224.6				
		c		-500.3				
		c		-1007.5				
As <sub>4</sub> O <sub>6</sub>	octahedral	c, II		-313.94	-275.36	201.835	51.2	45.72
	monoclinic	c, I		-312.8				
		aq		-299.4				
As <sub>2</sub> O <sub>3</sub> ·As <sub>2</sub> O <sub>5</sub>		c		-351.1				



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 20-3. Arsenic (at. no., 33; at. wt., 74.91)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance		State	Description	$\Delta F^{\circ}$					$S^{\circ}$	$C_p^{\circ}$
Formula	$\Delta F^{\circ}$									
	At 0°K			$\Delta F^{\circ}$	$\log_{10} K_f$	$\Delta F^{\circ}$				
		kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole	cal/deg mole	cal/deg mole	cal/deg mole
AsH <sub>3</sub>										
AsH <sub>3</sub> · 6H <sub>2</sub> O		g		41.0						
		c		-386.7						
HAso <sub>2</sub>		aq	std. state, hyp. m = 1	-109.0	-96.25	70.550	30.3			
HAso <sub>3</sub> <sup>--</sup>		aq		-157.1						
HAso <sub>4</sub> <sup>--</sup>		aq	std. state, hyp. m = 1	-214.8	-169.	123.9	0.9			
H <sub>2</sub> AsO <sub>3</sub> <sup>-</sup>		aq		-170.3						
H <sub>2</sub> AsO <sub>4</sub> <sup>-</sup>		aq	std. state, hyp. m = 1	-216.2	-178.9	131.13	28.			
H <sub>3</sub> AsO <sub>3</sub>		aq	std. state, hyp. m = 1	-177.3	-152.9	112.07	47.0			
H <sub>3</sub> AsO <sub>4</sub>		c		-215.2						
H <sub>3</sub> AsO <sub>4</sub> · 2H <sub>2</sub> O		aq	std. state, hyp. m = 1	-214.8	-183.8	134.72	49.3			
		c								

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 20-4. Arsenic [at. no., 33; at. wt., 74.91]

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} f_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)			cal/deg mole
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	
AsF <sub>3</sub>		g		-218.3	-214.7	157.37	69.08
AsF <sub>5</sub>		liq		-226.8	-215.5	157.96	43.31
AsCl <sub>2</sub>		g		7.35			
AsCl <sub>3</sub>		g		-71.5	-68.5	50.21	78.2
AsBr <sub>3</sub>		liq		-80.2	-70.5	51.68	55.8
AsI <sub>2</sub>		c		-46.61			
AsI <sub>3</sub>		c		-13.7			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 20-5. Arsenic (at. no., 33; at. wt., 74.91)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance		State	At 298.16°K (25°C)				
Formula	Description		$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
$As_2S_2$		g	-4.22				
$As_2S_3$		c	-31.9				
$As_2O_3 \cdot SO_3$		c	-35.				
		c	-285.26		*		
ASN		g	7.				7.274
$NH_4H_2AsO_4$		c	-251.47	-197.24	144.574	41.12	
$(NH_4)_2HAsO_4$		aq	-247.9				
		c	-280.24				
$(NH_4)_3AsO_4$		aq	-278.3				
		c	-306.11				
		aq	-303.				
$(NH_4)_3AsO_4 \cdot 3H_2O$		c	-516.6				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 21-1. Antimony (at. no., 51; at. wt., 121.76) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1950									
Formula	Substance	Description	State	At 298.16°K (25°C)					$C_p^0$ cal/deg mole
				$\Delta H_f^0$ kcal/mole	$\Delta H_f^0$ kcal/mole	$\Delta F_f^0$ kcal/mole	$\log_{10} K_f$	$S^0$ cal/deg mole	
				At 0°K kcal/mole	kcal/mole	kcal/mole		cal/deg mole	
Sb			g		60.8	51.1	-37.46	43.05	4.968
			c, III	0.000	0.000	0.000	0.0000	10.5	6.08
		explosive	c, IV		2.44				
Sb <sup>+</sup>			g		261.52				
Sb <sup>++</sup>			g		678.1				
Sb <sup>+++</sup>			g		1249.				
Sb <sup>++++</sup>			g		2263.				
Sb <sup>+++++</sup>			g		3542.				
Sb <sup>+++++</sup>			g		6013.				
Sb <sub>2</sub>			g		52.	40.	-29.3	60.9	8.684
Sb <sub>4</sub>			g		48.8				
SbO			g		45.				
SbO <sup>+</sup>		std. state, hyp. m=1	aq			-42.0	30.79		

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 21-2. Antimony (at. no., 51; at. wt., 121.76)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance							
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
			At 0°K	At 298.16°K (25°C)			
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
$Sb_2O_3$		aq		-166.5			
$Sb_2O_4$		c		-214.	137.8	30.3	27.4
$Sb_2O_5$		c		-234.4	146.96	29.9	28.1
		aq		-226.4			
$Sb_4O_6$		c, II		-336.8	218.43	58.3	48.46
$Sb_6O_{13}$		c		-670.			
$SbH_3$		g					
$H_3SbO_4$		aq		-215.7			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 21-3. Antimony (at. no., 51; at. wt., 121.76)

HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance									
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f^\circ$	$S^\circ$	$C_p^\circ$	
			At 0°K			At 298.16°K (25°C)			
			kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole	
SbF		g		0.					
SbF <sub>3</sub>		c		-217.2					
		aq		-216.1					
SbF <sub>5</sub>		liq							
2SbF <sub>3</sub> •SbF <sub>5</sub>		liq							
H <sub>3</sub> SbF <sub>6</sub>		aq		-444.3					
SbCl		g		-4.					
SbCl <sub>3</sub>		g		-75.2	-72.3	52.99	80.8	18.5	
		c		-91.34	-77.62	56.894	44.5		
SbCl <sub>5</sub>		g		-93.9					
		liq		-104.8					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

Washington, D.C.

National Bureau of Standards

Table 21-4. Antimony (at. no., 51; at. wt., 121.76) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1950											
	Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
					At 0°K	At 298.16°K (25°C)					
					kcal/mole	kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
SbOCl				c		-90.8					
Sb <sub>4</sub> O <sub>5</sub> Cl <sub>2</sub>				c		-350.					
SbCl <sub>3</sub> F <sub>2</sub>				c							
SbBr <sub>3</sub>				c		-62.1					
		in CS <sub>2</sub>				-58.4					
SbI <sub>3</sub>				c		-23.0					
				aq		-22.6					
SbS <sub>3</sub>				aq		-4.					
Sb <sub>2</sub> S <sub>3</sub>		black		c		-43.5					
		orange		amorp		-36.0					
Sb <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>				c		-575.3					
Sb <sub>2</sub> Se <sub>3</sub>											
Sb <sub>2</sub> Te <sub>3</sub>											

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 21-5. Antimony (at. no., 51; at. wt., 121.76) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1950									
	Substance			State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
	Formula	Description							
SbN			g		74.4				
SbF <sub>3</sub> ·NH <sub>3</sub>			c		-242.3				
SbF <sub>3</sub> ·2NH <sub>3</sub>			c		-265.5				
SbF <sub>3</sub> ·3NH <sub>3</sub>			c		-285.7				
SbF <sub>3</sub> ·4NH <sub>3</sub>			c		-304.4				
SbF <sub>3</sub> ·6NH <sub>3</sub>			c		-341.4				
3NH <sub>4</sub> Cl·2SbCl <sub>3</sub>			c						



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 22-1. Bismuth (at. no., 83; at. wt., 209.00) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1947; March 31, 1950									
Substance		State	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole	
Formula	Description								
Bi		g	0.000	49.7	40.4	-29.61	44.67	4.964	
Bi <sup>+</sup>		c		0.000	0.000	0.0000	13.6	6.1	
Bi <sup>++</sup>		g		219.1					
Bi <sup>+++</sup>		g		605.3					
Bi <sup>++++</sup>		g		1196.					
Bi <sup>+++++</sup>		g		2242.					
Bi <sup>+++++</sup>		g		3535.					
Bi <sup>+++++</sup>		g		5715.					
Bi <sub>2</sub>		g		59.4	48.0	-35.18	65.4	8.836	
BiO		g		16.					
BiO <sup>+</sup>		c		-49.85					
Bi <sub>2</sub> O <sub>3</sub>	std. state, hyp. m = 1	aq		-137.9	-34.54	25.317			
		c			-118.7	87.1	46.4	27.2	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 22-2. Bismuth (at. no., 83; at. wt., 209.00)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1949

Substance		State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description							
BiH		g						
Bi(OH) <sub>3</sub>		c		43. -169.6				
BiF		g		7.				
BiF <sub>3</sub>		c						
BiF <sub>5</sub>		c						
BiCl		g		10.7	5.2	-3.81	58.9	
BiCl <sub>3</sub>		g		-64.7	-62.2	45.59	85.3	
		c		-90.61	-76.23	55.876	45.3	19.02
	in aq HCl	aq		-101.6				
BiCl <sub>4</sub>		c						
BiOCl		c		-87.3	-77.0	56.44	20.6	
BiBr		g		12.7	3.8	-2.79	61.6	
BiBr <sub>3</sub>		c, II						
BiI		g		16.	11.	-8.06	63.4	

Table 22-3. Bismuth (at. no., 83; at. wt., 209.00)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1947; March 31, 1950

Substance						
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$
			At 0°K	At 298.16°K (25°C)		
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole
BiS		c				
Bi <sub>2</sub> S <sub>3</sub>		c		-43.8	28.88	35.3
Bi <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c		-607.		
BiSe		c, II				
Bi <sub>2</sub> Se <sub>3</sub>		c				
Bi <sub>2</sub> Te <sub>3</sub>		c				
Bi <sub>2</sub> Te <sub>3</sub> ·Bi <sub>2</sub> S <sub>3</sub>		c				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 23-1. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 21, 1948; March 19-20

Table 23-1. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 21, 1948; March 1949									
Substance		Description	State	$\Delta H_f^0$	$\Delta H_f^0$	$\Delta F_f^0$	$\log_{10} K_f^*$	$S^0$	$C_p^0$
Formula	At 0°K			At 298.16°K (25°C)	cal/mole	kcal/mole	cal/deg mole	cal/deg mole	
C			g	170.39	171.698	160.845	-117.8971	37.7611	4.985
		diamond	c	0.5766	0.4532	0.6850	-0.50210	0.5829	1.449
		graphite	c	0.000	0.000	0.000	0.0000	1.3609	2.066
C <sup>+</sup>			g	430.167	431.648				
C <sup>++</sup>			g	992.318	995.281				
C <sup>+++</sup>			g	2096.16	2100.60				
C <sup>++++</sup>			g	3583.09	3589.02				
C <sup>+++++</sup>			g	12623.0	12360.4				
C <sup>+++++</sup>			g	23920.	23930.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 23-2. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1948; March 31, 1950

Substance		At 0°K		At 298.16°K (25°C)		S°	C <sub>p</sub> <sup>o</sup>
Formula	Description	State	ΔH <sup>o</sup> kcal/mole	ΔH <sup>o</sup> kcal/mole	ΔF <sup>o</sup> kcal/mole		
CO		g	-27.2019	-26.4157	-32.6079	47.401	6.447
CO <sup>+</sup>		g	291.	292.5			
CO <sub>2</sub>		g	-93.9686	-94.0518	-94.2598	51.061	8.474
	std. state, hyp. m = 1	aq		-98.69	-92.31	29.3	
CO <sub>2</sub> ·6H <sub>2</sub> O		c		-520.			
CO <sub>2</sub> <sup>+</sup>		g	222.6	224.1			
CO <sub>3</sub> <sup>++</sup>		g					
	std. state, hyp. m = 1	aq		-161.63	-126.22	-12.7	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 23-3. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1948; March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	At 0°K kcal/mole			kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole	cal/deg mole		
CH			g	142.0	142.1				
CH <sup>+</sup>			g						
CH <sub>2</sub>			g						
CH <sub>2</sub> <sup>+</sup>			g						
CH <sub>3</sub>			g	33.4	32.0				
CH <sub>3</sub> <sup>+</sup>			g						
CH <sub>4</sub>			g	-15.987	-17.889 -445.	-12.140	44.50	8.536	
CH <sub>4</sub> · 6H <sub>2</sub> O			c						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 23-4. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1948; March 31, 1950

Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 298.16°K (25°C)								
	At 0°K kcal/mole			kcal/mole	kcal/mole		cal/deg mole	cal/deg mole	
$\text{HCOO}^-$		formate ion, std. state, hyp. m=1	aq		-98.0	-80.0	58.64	21.9	
$\text{HCO}_3^-$		bicarbonate ion std. state, hyp. m=1	aq		-165.18	-140.31	102.645	22.7	
$\text{CH}_2\text{O}$		formaldehyde	g		-27.7	-26.3	19.28	52.26	8.45
		in 60 $\text{H}_2\text{O}$	aq		-42.5				
		in 40 $\text{H}_3\text{OH}$			-42.7				
$\text{CH}_2\text{O}_2$		formic acid, monomer	g		-86.67	-80.24	58.815	60.0	
		formic acid	liq		-97.8	-82.7	60.62	30.82	23.67
		std. state, hyp. m=1	aq		-98.0	-85.1	62.36	39.1	
		in 0.2 $\text{H}_2\text{O}$	aq		-97.86				
		0.5 $\text{H}_2\text{O}$	aq		-97.93				
		1.0 $\text{H}_2\text{O}$	aq		-98.00				
		1.5 $\text{H}_2\text{O}$	aq		-98.01				
		2.0 $\text{H}_2\text{O}$	aq		-98.01				
		3.0 $\text{H}_2\text{O}$	aq		-97.99				
		4 $\text{H}_2\text{O}$	aq		-97.96				
		5 $\text{H}_2\text{O}$	aq		-97.96				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards Washington, D.C.

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Table 23-5. Carbon (at. no., 6; at. wt., 12.010)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1948; March 31, 1950

Substance		Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K			At 298.16°K (25°C)					
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
CH <sub>2</sub> O <sub>2</sub>	formic acid, in 10 H <sub>2</sub> O 15 H <sub>2</sub> O 25 H <sub>2</sub> O 50 H <sub>2</sub> O 100 H <sub>2</sub> O 200 H <sub>2</sub> O ∞ H <sub>2</sub> O	aq		-97.94					
				-97.93					
				-97.94					
				-97.95					
				-97.96					
				-97.97					
H <sub>2</sub> CO <sub>3</sub> CH <sub>3</sub> O <sup>-</sup>	carbonic acid, std. state, hyp. m=1 methylate ion	aq		-167.0	-149.00	109.215	45.7		
		aq							



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 23-6. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1948; March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	At 298.16°K (25°C)				
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
CH <sub>3</sub> OH	methanol		g						
			liq						
	std. state, hyp. m = 1		aq						
	in 0.25 H <sub>2</sub> O		aq						
	0.5 H <sub>2</sub> O		aq						
	1.0 H <sub>2</sub> O		aq						
	1.5 H <sub>2</sub> O		aq						
	2.0 H <sub>2</sub> O		aq						
	3 H <sub>2</sub> O		aq						
	4 H <sub>2</sub> O		aq						
CH <sub>3</sub> O <sub>2</sub>	5 H <sub>2</sub> O		aq						
	10 H <sub>2</sub> O		aq						
	25 H <sub>2</sub> O		aq						
	∞ H <sub>2</sub> O		aq						
	methanol hydrogen peroxide		liq						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

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Table 23-7. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1948; March 31, 1950										
Substance		State	Description	ΔH <sup>o</sup>		ΔF <sup>o</sup>	log <sub>10</sub> K <sup>o</sup>	S <sup>o</sup>	C <sub>p</sub> <sup>o</sup>	
Formula				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole					cal/deg mole
CF <sub>4</sub>	tetrafluoromethane	g			-162.5	-151.8	111.27	62.7		
COF <sub>2</sub>	carbonyl fluoride	g							8.95	
CH <sub>3</sub> F	fluoromethane	g						53.30	10.24	
CH <sub>2</sub> F <sub>2</sub>	difluoromethane	g							12.68	
CHF <sub>3</sub>	trifluoromethane	g								
CCl <sub>4</sub>	tetrachloromethane	g			-25.5	-15.3	11.21	73.95	19.96	
COCl <sub>2</sub>	carbonyl chloride	liq			-33.3	-16.4	12.02	51.25	31.49	
		g			-53.30	-50.31	36.877	69.13	14.51	
CH <sub>3</sub> Cl	chloromethane	g			-19.6	-14.0	10.26	55.97	9.75	
CH <sub>3</sub> Cl·9H <sub>2</sub> O		c			-650.					
CH <sub>2</sub> Cl <sub>2</sub>	dichloromethane	g			-21.	-14.	10.3	64.68	12.28	
		liq			-28.0	-15.1	11.07	42.7	23.9	
CHCl <sub>3</sub>	trichloromethane	g			-24.	-16.	11.7	70.86	15.73	
		liq			-31.5	-17.1	12.53	48.5	27.8	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 23-8. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1948; March 31, 1950

Substance		State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description							
CF <sub>3</sub> Cl	trifluorochloromethane	g						15.24
CF <sub>2</sub> Cl <sub>2</sub>	difluorodichloromethane	g						16.56
CFCl <sub>3</sub>	fluorotrichloromethane	g					74.06	18.73
		liq						29.00
COFCl	carbonyl fluorochloride	g						
CH <sub>2</sub> FCl	fluorochloromethane	g						11.37
CHF <sub>2</sub> Cl	difluorochloromethane	g						13.79
CHFCl <sub>2</sub>	fluorodichloromethane	g						14.64

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 23-9. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1948; March 31, 1950

Substance										
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
			At 298.16°K (25°C)							
			kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole		
CBr <sub>4</sub>	tetrabromomethane	g		12.	8.6	-6.30	85.6	21.84		
COBr <sub>2</sub>	carbonyl bromide	c, li		-23.				35.0		
		liq		-30.						
CH <sub>3</sub> Br	bromomethane	g		-8.5	-6.2	4.54	58.74	10.18		
CH <sub>2</sub> Br <sub>2</sub>	dibromomethane	g		-1.	-1.4	1.03	70.16	13.15		
CHBr <sub>3</sub>	tribromomethane	g		6.	3.8	-2.78	79.18	17.08		
		liq		-4.8	0.7	-0.51	53.0			
CF <sub>3</sub> Br	trifluorobromomethane	g						16.8		
CF <sub>2</sub> Br <sub>2</sub>	difluorodibromomethane	g						18.5		
CFBr <sub>3</sub>	fluorotribromomethane	g						19.4		

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-10. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1948; March 31, 1950

Substance		State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description		At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
$\text{CH}_2\text{FBr}$	fluorobromomethane	g					11.8
$\text{CHF}_2\text{Br}$	difluorobromomethane	g					13.9
$\text{CHBrBr}_2$	fluorodibromomethane	g					15.41
$\text{CCl}_3\text{Br}$	trichlorobromomethane	g					20.60
$\text{CCl}_2\text{Br}_2$	dichlorodibromomethane	g					20.92
$\text{CClBr}_3$	chlorotribromomethane	g					21.24
$\text{CH}_2\text{ClBr}$	chlorobromomethane	g					12.8
$\text{CHCl}_2\text{Br}$	dichlorobromomethane	g					16.08
$\text{CHClBr}_2$	chlorodibromomethane	g					16.52
$\text{CHFClBr}$	fluorochlorobromomethane	g					15.1

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-11. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1948; March 31, 1950

Substance		State	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description		At 0°K					
$Cl_4$	tetraiodomethane	c						
$CH_3I$	iodomethane	g		4.9	5.3	-3.88	60.85	10.55
$CH_2I_2$	diiodomethane	liq		-2.0	4.9	-3.59	38.9	
		g						13.81
		liq		15.9				
$CHI_3$	triiodomethane	c		33.6				
$CHF_2I$	difluoroiodomethane	g						
$CHF_2I_2$	fluorodiiodomethane	liq						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 23-12. Carbon (at. no., 6; at. wt., 12.010)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1948; March 31, 1950

Substance									
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole	
CS		g					50.4	7.11	
$\frac{1}{n}(\text{CS})_n$		c		13.					
CS <sub>2</sub>	carbon disulfide	g		27.55	15.55	-11.398	56.84	10.91	
		liq		21.0	15.2	-11.14	36.10	18.1	
COS	carbon oxysulfide	g		-32.80	-40.45	29.649	55.34	9.92	
CH <sub>3</sub> S	methanethiol	g		-2.97	0.21	-0.154	60.90	12.1	
CSCl <sub>2</sub>	carbon sulfochloride	g					70.17	15.45	
CSe <sub>2</sub>	carbon diselenide	liq		37.1					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 23-13. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1948; March 31, 1950

Table 23-13. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1948; March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\Delta F^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	kcal/mole			kcal/mole	kcal/mole	At 298.16°K (25°C)		cal/deg mole	cal/deg mole
						At 0°K	kcal/mole		
CN			g					48.40	6.96
CN <sup>-</sup>			g						
		cyanide ion, std. state, hyp. m = 1	aq		36.1	39.6	-29.03	28.2	
CN <sub>4</sub>			c		92.6				
$\frac{2}{3}(\text{CN}_4)_n$		cyanogen azide	c		82.2				
CNO <sup>-</sup>			aq		-33.5	-23.6	17.30	31.1	
C(NO <sub>2</sub> ) <sub>4</sub>		cyanate ion, std. state, hyp. m = 1 tetranitromethane	liq		8.8				
HCN		hydrogen cyanide	g		31.2	28.7	-21.04	48.23	8.58
			liq		25.2	29.0	-21.26	26.97	16.88
		hydrocyanic acid, std. state, hyp. m = 1	aq		25.2	26.8	-19.64	30.8	



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-14. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1948; March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
CH <sub>5</sub> N	methyl amine in 400 H <sub>2</sub> O 600 H <sub>2</sub> O 800 H <sub>2</sub> O 1000 H <sub>2</sub> O		g aq aq aq aq		-6.7 -17.84 -17.88 -17.93 -17.99		-4.84	57.73	12.9
CH <sub>6</sub> N <sup>+</sup>	methyl ammonium ion, std. state, hyp. m=1		aq		-23.653				
CH <sub>2</sub> N <sub>2</sub>	cyanamide in 800 H <sub>2</sub> O		c aq		9.2 12.9				
NH <sub>4</sub> CN	ammonium cyanide		c aq		0.0 4.6				32.0

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-15. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1948; March 31, 1950									
Substance									
Formula	Description	State	$\Delta H^{\circ}$		$\Delta F^{\circ}$	$\log_{10} K^{\circ}$	$S^{\circ}$	$C_p^{\circ}$	
			At 0°K kcal/mole	kcal/mole					
$\text{CH}_5\text{N}_3$	guanidine	c							
		aq		-17.0					
		aq		-18.3					
$\text{CH}_6\text{N}_3^{+}$	guanidine ion	aq		-32.2					
$\text{HCNO}$	cyanoic acid, std. state, hyp. m=1	aq		-35.1	-28.9	21.18	43.6		
$\text{HNCO}$	hydrogen isocyanate	g							
$\text{CH}_2\text{O}_2\text{N}^{-}$	nitromethane ion of nitro form	aq		-14.5					
$\text{CH}_3\text{ON}$	formamide	liq		-61.6					
		aq		-56.0					
$\text{CH}_3\text{O}_2\text{N}$	nitroethane	liq		-21.28	2.26	-1.656	41.1	25.3	
		aq		-20.7					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-16. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1945; March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta G_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole		cal/deg mole	cal/deg mole	
$\text{CH}_3\text{O}_2\text{N}$	methyl nitrite		g						
$\text{CH}_3\text{O}_2\text{N}$	methyl nitrate		liq						
$\text{CH}_5\text{O}_2\text{N}$	ammonium formate		c	-132.8					
			aq	-129.83					
$\text{NH}_4\text{HCO}_3$	ammonium bicarbonate		c	-203.7					
	std. state, hyp. m = 1		aq	-196.92	-159.31	116.77?	49.7		
$\text{CH}_4\text{ON}_2$	urea		c	-79.634	-47.12?	34.5383	25.0?		
	std. state, hyp. m = 1		aq	-76.30	-48.72	35.711	41.7?		
	in 5 $\text{H}_2\text{O}$		aq	-75.790					
	7 $\text{H}_2\text{O}$		aq	-75.879					
	10 $\text{H}_2\text{O}$		aq	-75.970					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)		Table 23-17. Carbon (at. no., 6; at. wt., 12.010) March 31, 1948; March 31, 1950				
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$
				At 0°K	At 298.16°K (25°C)	$S^\circ$
				kcal/mole	kcal/mole	cal/deg mole
CH <sub>4</sub> ON <sub>2</sub>	in 15 H <sub>2</sub> O		aq		-76.056	
	20 H <sub>2</sub> O		aq		-76.107	
	25 H <sub>2</sub> O		aq		-76.142	
	50 H <sub>2</sub> O		aq		-76.219	
	100 H <sub>2</sub> O		aq		-76.259	
	200 H <sub>2</sub> O		aq		-76.281	
	400 H <sub>2</sub> O		aq		-76.29	
	∞ H <sub>2</sub> O		aq		-76.301	
			c		-74.7	
			aq		-68.5	
NH <sub>4</sub> CNO	ammonium cyanate					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-18. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1948; March 31, 1950

Substance		$\Delta H^{\circ}$		$\Delta F^{\circ}$	$\log_{10} K$	$S^{\circ}$	$C_p^{\circ}$
Formula	Description	State	At 0°K		At 298.16°K (25°C)		cal/deg mole
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	
$\text{CH}_3\text{O}_2\text{N}_2$	ammonium carbamate	c		-154.21	-109.47	80.240	39.70
$(\text{NH}_4)_2\text{CO}_3$	ammonium carbonate, std. state, hyp. m = 1	aq		-150.4			
		aq		-225.11	-164.22	120.371	41.2
$\text{CH}_3\text{N}_3$	trinitromethane	liq		-18.6			
$\text{CH}_5\text{O}_4\text{N}_3$	urea nitrate	c		-114.8			
$\text{CH}_4\text{O}_2\text{N}_4$	nitroguanidine	c		-22.14			
$\text{CH}_6\text{O}_3\text{N}_4$	guanidine nitrate	c		-91.1			
		aq		-81.6			
CNF	cyanogen fluoride	g					
$\text{COF}_3\text{N}$	N,N-difluorofluoroformamide	g					
$\text{COF}_3\text{N}$	trifluoronitrosomethane	g					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 23-19. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1948; March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole				
				At 298.16°K (25°C)					
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
CNCl	cyanogen chloride		g		34.5	32.9	-24.12	56.31	10.70
CH <sub>3</sub> NCI	methyllamine hydrochloride		c		-68.31	-35.09	25.720	33.13	21.75
	std. state, hyp. m = 1		aq		-69.876				
	in 1 H <sub>2</sub> O		aq		-69.737				
	2 H <sub>2</sub> O		aq		-69.756				
	3 H <sub>2</sub> O		aq		-69.765				
	5 H <sub>2</sub> O		aq		-69.78				
	10 H <sub>2</sub> O		aq		-69.798				
	20 H <sub>2</sub> O		aq		-69.818				
	50 H <sub>2</sub> O		aq		-69.837				
	100 H <sub>2</sub> O		aq		-69.848				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 23-20. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1948; March 31, 1950

Substance		$\Delta H_f^\circ$	$\Delta G_f^\circ$	$\log_{10} f_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)		
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole
CH <sub>3</sub> NCI	methylaniline hydrochloride	aq				
	in 200 H <sub>2</sub> O			-69.855		
	500 H <sub>2</sub> O	aq		-63.862		
	1000 H <sub>2</sub> O	aq		-69.856		
	$\infty$ H <sub>2</sub> O	aq		-69.876		
CH <sub>3</sub> ON <sub>3</sub> Cl	semicarbazide hydrochloride	c				34.
CNBr	cyanogen bromide	g			59.05	11.11
CNI	cyanogen iodide	g			61.26	11.54
		c		-34.96	47.7	
				-31.23	42.6	
		aq		-31.74	43.3	
	std. state, hyp. m = 1					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-21. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1948; March 31, 1950

Substance		State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description							
CNS <sup>-</sup>	thiocyanate ion	aq		17.2				
HCNS	hydrogen thiocyanate	liq						
	thiocyanic acid	aq		17.7				
CH <sub>4</sub> N <sub>2</sub> S	thiourea	c		-22.1				
		aq		-16.8				
NH <sub>4</sub> CNS	ammonium thiocyanate	c		-20.0				
		aq		-14.5				
CH <sub>5</sub> O <sub>3</sub> N <sub>3</sub> S	thiourea nitrate	c		-74.5				



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

Washington, D.C.

Table 23-22. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1948; March 31, 1950

Substance		State	Description	$\Delta H_f^\circ$			$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K			At 298.16°K (25°C)					
	kcal/mole			kcal/mole	kcal/mole	cal/deg mole			
CP		g						51.92	7.14
CONPF <sub>2</sub>	phosphorus difluoroisocyanate	g							
CNPSF <sub>2</sub>	phosphorus difluoroisothiocyanate	liq							
CNPSCl <sub>2</sub>	phosphorus dichlorothiocyante	liq							
CH <sub>3</sub> As	methyl arsine	g							
CH <sub>3</sub> F <sub>2</sub> As	methyl difluoroarsine	liq							
CH <sub>3</sub> Cl <sub>2</sub> As	methyl dichloroarsine	liq							

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

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Table 23-23. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1948; March 31, 1950

Formula	Substance	Description	State	At 298.16°K (25°C)				
				$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
C <sub>2</sub>			g	233.1				
C <sub>2</sub> O <sub>4</sub> <sup>2-</sup>	oxalate ion; std. state, hyp. m = 1		aq		-161.3	118.23	12.2	
C <sub>2</sub> H <sub>2</sub>	ethyne (acetylene)		g	54.329	50.000	-36.6490	47.997	10.499
			aq		50.6			
C <sub>2</sub> H <sub>2</sub> ·6H <sub>2</sub> O			c		-371.1			
C <sub>2</sub> H <sub>4</sub>	ethene (ethylene)		g	14.522	16.282	-11.9345	52.45	10.41
C <sub>2</sub> H <sub>4</sub> ·6H <sub>2</sub> O			c		-412.8			
C <sub>2</sub> H <sub>6</sub>	ethane		g	-16.517	-20.236	5.7613	54.85	12.585
HC <sub>2</sub> O <sub>4</sub> <sup>-</sup>	bioxalate ion; std. state, hyp. m = 1		aq		-167.1	122.48	36.7	
C <sub>2</sub> H <sub>2</sub> O	ketene		g		-14.6			
C <sub>2</sub> H <sub>2</sub> O <sub>2</sub>	glyoxal		c		-83.7			
			aq		-82.3			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 23-24. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1948									
Substance		State		At 0°K		At 298.16°K (25°C)		S°	C <sub>p</sub> <sup>o</sup>
Formula	Description			ΔH <sup>fo</sup> kcal/mole	ΔF <sup>fo</sup> kcal/mole	log <sub>10</sub> K <sup>o</sup>	cal/deg mole		
C <sub>2</sub> H <sub>2</sub> O <sub>4</sub>	oxalic acid	c							
	in 300 H <sub>2</sub> O	aq							
	2100 H <sub>2</sub> O	aq							
	in CH <sub>4</sub> O (methanol)								
	in C <sub>2</sub> H <sub>5</sub> O (ethanol)								
C <sub>2</sub> H <sub>2</sub> O <sub>4</sub> · 2H <sub>2</sub> O	in C <sub>3</sub> H <sub>8</sub> O (1-propanol)								
	oxalic acid dihydrate	c							

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-25. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1948 (corrected)									
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 298.16°K (25°C)					
				At 0°K kcal/mole	kcal/mole	kcal/mole		kcal/deg mole	cal/deg mole
$C_2H_3O_2^-$	acetate ion, std. state, hyp. m = 1		aq		-116.843				
$C_2H_3O_3^-$	glycollate ion		aq		-155.2				
$C_2H_4O$	acetaldehyde		g	-37.4	-39.76	-31.96	23.426	63.5	15.0
	in 50 H <sub>2</sub> O		aq		-49.88				
$C_2H_4O \cdot H_2O$	acetaldehyde monohydrate		c						
$C_2H_4O$	ethylene oxide		g		-12.19	-2.79	2.045	58.1	11.5
$C_2H_4O_2$	acetic acid		liq		-116.4	-93.8	68.75	38.2	29.5
	std. state, hyp. m = 1		aq		-116.27				
	in 0.5 H <sub>2</sub> O		aq		-116.25				
	1 H <sub>2</sub> O		aq		-116.23				
	1.5 H <sub>2</sub> O		aq		-116.24				
	2.0 H <sub>2</sub> O		aq		-116.26				
	3 H <sub>2</sub> O		aq		-116.31				
	5 H <sub>2</sub> O		aq		-116.40				
	8 H <sub>2</sub> O		aq		-116.40				
	10 H <sub>2</sub> O		aq		-116.450				

Table 23-26. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1948									
Substance		State	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	Description								
$C_2H_4O_2$	acetic acid, in 25 $H_2O$	aq							
	50 $H_2O$	aq							
	100 $H_2O$	aq							
	200 $H_2O$	aq							
	300 $H_2O$	aq							
	400 $H_2O$	aq							
	500 $H_2O$	aq							
	800 $H_2O$	aq							
	1000 $H_2O$	aq							
	3000 $H_2O$	aq							
	5000 $H_2O$	aq							
	$\infty H_2O$ (undissociated)	aq							
	in 5 $H_2SO_4$								
	10 $H_2SO_4$								
	15 $H_2SO_4$								
	20 $H_2SO_4$								
	25 $H_2SO_4$								
	30 $H_2SO_4$								
	50 $H_2SO_4$								

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-27. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1948

Substance		State	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description						
		At 0°K kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
C <sub>2</sub> H <sub>4</sub> O <sub>2</sub>	acetic acid, in 5 C <sub>3</sub> H <sub>6</sub> O (acetone)		-117.05				
	10 C <sub>3</sub> H <sub>6</sub> O		-117.14				
	15 C <sub>3</sub> H <sub>6</sub> O		-117.21				
	25 C <sub>3</sub> H <sub>6</sub> O		-117.30				
	50 C <sub>3</sub> H <sub>6</sub> O		-117.38				
	in 5 C <sub>4</sub> H <sub>10</sub> O (diethyl ether)		-116.45				
	10 C <sub>4</sub> H <sub>10</sub> O		-116.43				
	15 C <sub>4</sub> H <sub>10</sub> O		-116.42				
	in 5 C <sub>5</sub> H <sub>12</sub> (n-pentane)		-115.40				
	10 C <sub>5</sub> H <sub>12</sub>		-115.18				
	15 C <sub>5</sub> H <sub>12</sub>		-115.04				
	25 C <sub>5</sub> H <sub>12</sub>		-114.89				
	50 C <sub>5</sub> H <sub>12</sub>		-114.74				
	in 5 C <sub>6</sub> H <sub>6</sub> (benzene)		-115.9				
	10 C <sub>6</sub> H <sub>6</sub>		-115.7				
	15 C <sub>6</sub> H <sub>6</sub>		-115.6				
	25 C <sub>6</sub> H <sub>6</sub>		-115.4				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-28. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1948

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C)		cal/deg mole	
$C_2H_4O_2$	methyl formate		g		-83.6			29.
	in 33 H <sub>2</sub> O		liq		-90.4			
	hydroxyacetaldehyde		aq		-91.6			
$C_2H_4O_2$	hydroxyacetic acid (glycollic acid)		c, I		-158.2			83.1
$C_2H_4O_3$			c, II		-158.1			
	in 400 H <sub>2</sub> O		aq		-155.3			
$C_2H_4O_4$	dihydroxyacetic acid (glyoxylic acid)		c		-199.6			
	formic acid dimer		aq		-197.0			
$C_2H_4O_4$			g		-187.7	120.09	-163.8	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-29. Carbon (at. no., 6; at. wt., 12.010)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1948

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
$C_2H_5O^-$	ethylate ion		g						
$C_2H_5O$	ethanol		liq						
	in 0.05 $H_2O$		aq						
	0.1 $H_2O$		aq						
	0.2 $H_2O$		aq						
	0.5 $H_2O$		aq						
	1 $H_2O$		aq						
	1.5 $H_2O$		aq						
	2 $H_2O$		aq						
	3 $H_2O$		aq						
	4 $H_2O$		aq						
	5 $H_2O$		aq						
	6 $H_2O$		aq						
	7 $H_2O$		aq						
	8 $H_2O$		aq						
	9 $H_2O$		aq						
	10 $H_2O$		aq						



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-30. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1948

Substance		$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description						
$C_2H_6O$	ethanol, in 15 $H_2O$						
	25 $H_2O$						
	50 $H_2O$						
	100 $H_2O$						
	200 $H_2O$						
	in 5 $H_2SO_4$						
	10 $H_2SO_4$						
	15 $H_2SO_4$						
	20 $H_2SO_4$						
	25 $H_2SO_4$						
	30 $H_2SO_4$						
	50 $H_2SO_4$						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-31. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1948									
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole		cal/deg mole	cal/deg mole	
$C_2H_6O$	ethanol, in 0.2 $C_6H_6$ (benzene) 0.5 $C_6H_6$ 1 $C_6H_6$ 3 $C_6H_6$ 5 $C_6H_6$ 10 $C_6H_6$ 25 $C_6H_6$ 50 $C_6H_6$ 100 $C_6H_6$				-66.50				
					-66.65				
					-66.90				
					-67.50				
					-67.90				
					-68.50				
					-69.15				
					-69.75				
					-70.00				
					-44.3	19.98	63.72	15.76	
$C_2H_6O$	dimethyl ether in 200 $H_2O$		g aq		-27.3				
					-52.5				

Table 23-32. Carbon (at. no., 6; at. wt., 12.010)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1948

Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 298.16°K (25°C)								
	kcal/mole			kcal/mole	kcal/mole		cal/deg mole	cal/deg mole	
C <sub>2</sub> H <sub>6</sub> O <sub>2</sub>	1,2-ethanediol (ethylene glycol)	liq							
	in 0.1 H <sub>2</sub> O	aq							
	0.2 H <sub>2</sub> O	aq							
	0.3 H <sub>2</sub> O	aq							
	0.5 H <sub>2</sub> O	aq							
	0.7 H <sub>2</sub> O	aq							
	1 H <sub>2</sub> O	aq							
	1.5 H <sub>2</sub> O	aq							
	2 H <sub>2</sub> O	aq							
	3 H <sub>2</sub> O	aq							
	4 H <sub>2</sub> O	aq							
	5 H <sub>2</sub> O	aq							
	7 H <sub>2</sub> O	aq							
	10 H <sub>2</sub> O	aq							
	15 H <sub>2</sub> O	aq							
25 H <sub>2</sub> O	aq								
50 H <sub>2</sub> O	aq								
C <sub>2</sub> H <sub>6</sub> O <sub>2</sub>	ethyl hydrogen peroxide	liq							

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-33. Carbon (at. no. 6; at. wt., 12.010)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1948

Formula	Substance	State	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
$C_2F_4$	tetrafluoroethene	g						
$C_2F_6$	hexafluoroethane	g		-24.0				
$C_2OF_4$	trifluoroacetyl fluoride	g						
$C_2H_3F$	fluoroethene	g						
$C_2H_5F$	fluoroethane	g						
$C_2H_4F_2$	1,1-difluoroethane	g						
$C_2HF_3$	trifluoroethene	g						
$C_2H_3F_3$	1,1,1-trifluoroethane	g					66.87	
$C_2H_3OF$	acetyl fluoride	g						
$C_2H_3O_2F$	fluoroacetic acid	c		-180.9				
$C_2H_5OF$	2-fluoroethanol	liq		-109.7				
$C_2H_2O_2F_2$	difluoroacetic acid	liq		-134.4				
$C_2H_4OF_2$	2,2-difluoroethanol	liq		-161.4				
$C_2HO_2F_3$	trifluoroacetic acid	liq						
$C_2H_3OF_3$	2,2,2-trifluoroethanol	liq						

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Table 23-34. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)									
June 30, 1948; March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula									
$C_2Cl_2$	dichloroethyne	liq							
$C_2Cl_4$	tetrachloroethene	liq			-3.			35.8	
$C_2Cl_6$	hexachloroethane	c			-48.			41.	
$C_2O_2Cl_3^-$	trichloroacetate ion	aq			-12 <sup>2,10</sup>				
$C_2OCl_4$	trichloroacetyl chloride	liq			-89.3				
$C_2O_2Cl_4$	trichloromethyl chloroformate								

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-35. Carbon (at. no., 6; at. wt., 12.010)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1948

Substance		State	At 298.16°K (25°C)					C <sub>p</sub> <sup>o</sup>
Formula	Description		ΔBf <sup>o</sup>	ΔBf <sup>o</sup>	ΔBf <sup>o</sup>	log <sub>10</sub> K <sub>f</sub>	S <sup>o</sup>	
			kcal/mole	kcal/mole	kcal/mole		cal/deg mole	
C <sub>2</sub> H <sub>3</sub> Cl	chloroethene	g						
C <sub>2</sub> H <sub>5</sub> Cl	chloroethane	g						
C <sub>2</sub> H <sub>2</sub> Cl <sub>2</sub>	1,1-dichloroethene	liq		7.5				15.
C <sub>2</sub> H <sub>2</sub> Cl <sub>2</sub>	cis-1,2-dichloroethene	liq		-25.1	-12.7	9.331	65.90	27.
C <sub>2</sub> H <sub>2</sub> Cl <sub>2</sub>	trans-1,2-dichloroethene	liq						27.
C <sub>2</sub> H <sub>4</sub> Cl <sub>2</sub>	1,1-dichloroethane	liq		-36.4				
C <sub>2</sub> H <sub>4</sub> Cl <sub>2</sub>	1,2-dichloroethane	liq		-39.7	-19.2	14.11	49.84	30.83
C <sub>2</sub> HCl <sub>3</sub>	tri-chloroethene	liq					54.36	29.
C <sub>2</sub> H <sub>3</sub> Cl <sub>3</sub>	1,1,1-trichloroethane	liq						34.5
C <sub>2</sub> H <sub>3</sub> Cl <sub>3</sub>	1,1,2-trichloroethane	liq						
C <sub>2</sub> H <sub>2</sub> Cl <sub>4</sub>	1,1,1,2-tetrachloroethane	liq						45.
C <sub>2</sub> H <sub>2</sub> Cl <sub>4</sub>	1,1,2,2-tetrachloroethane	liq						54.
C <sub>2</sub> HCl <sub>5</sub>	pentachloroethane	liq						

Table 23-36. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1948; March 31, 1950

Formula	Substance	State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
$C_2H_2O_2Cl^-$	chloroacetate ion	aq		-121.2				27.
$C_2H_3OCl$	acetyl chloride	liq		-65.8				
$C_2H_3OCl$	chloroacetaldehyde	liq		-61.3				
$C_2H_3O_2Cl$	chloroacetic acid	c, I		-124.2				34.4
		c, II		-124.1				
		c, III		-123.4				
$C_2H_5OCl$	2-chloroethanol	aq		-120.6				
$C_2H_5OCl$	chloromethyl methyl ether	liq		-70.3				
$C_2H_7OCl$	dimethyl ether hydrochloride	g		-71.1	-49.1	36.00	89.3	
$C_2HO_2Cl_2^-$	dichloroacetate ion	aq		-123.4				
$C_2H_2OCl_2$	chloroacetyl chloride	liq		-68.9				
$C_2H_2OCl_2$	dichloroacetaldehyde							

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 23-37. Carbon (at. no., 6; at. wt., 12.010)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1948

Substance									
Formula	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 0°K					At 298.16°K (25°C)	
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole		
$C_2H_2O_2Cl_2$	dichloroacetic acid	liq							47.
	in 200 $H_2O$	aq		-120.2					
	300 $H_2O$	aq		-122.33					
	500 $H_2O$	aq		-122.50					
	800 $H_2O$	aq		-122.68					
$C_2H_4O_2Cl_2$ $C_2HOC l_3$	dichloroacetaldehyde monohydrate	c		-122.80					
	trichloroacetaldehyde (chloral)	liq		-51.1					36.0
	in 80 $H_2O$	aq		-63.0					
$C_2HO_2Cl_3$	trichloroacetic acid	c		-122.8					75.0
	in 200 $H_2O$	aq		-124.80					
	500 $H_2O$	aq		-125.11					



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 23-38. Carbon (at. no., 6; at. wt., 12.010)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1948

Substance		State	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description						
$C_2H_2O_2Cl_3$	trichloroacetic acid,						
	in 10 $C_2H_6O$ (ethanol)						
	20 $C_2H_6O$			-122.67			
	30 $C_2H_6O$			-122.85			
	50 $C_2H_6O$			-122.97			
	100 $C_2H_6O$			-123.16			
	in 5 $C_3H_6O$ (acetone)			-123.58			
	10 $C_3H_6O$			-126.45			
	20 $C_3H_6O$			-126.57			
	30 $C_3H_6O$			-126.72			
	50 $C_3H_6O$			-126.81			
	100 $C_3H_6O$			-126.92			
	in 20 $C_4H_{10}O$ (diethyl ether)			-122.71			
	50 $C_4H_{10}O$			-123.18			
	100 $C_4H_{10}O$			-123.40			
	in 10 $C_5H_{12}$ (n-pentane)			-115.15			
	20 $C_5H_{12}$			-115.40			
	30 $C_5H_{12}$			-115.55			
	50 $C_5H_{12}$			-115.73			
	100 $C_5H_{12}$			-116.08			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 23-39. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1948

Substance		State	Description	At 298.16°K (25°C)					C <sub>p</sub> <sup>o</sup>
Formula	ΔH <sub>f</sub> <sup>o</sup>			ΔH <sub>f</sub> <sup>o</sup>	ΔF <sub>f</sub> <sup>o</sup>	log <sub>10</sub> K <sub>f</sub>	S <sup>o</sup>		
	kcal/mole			kcal/mole	kcal/mole		cal/deg mole	cal/deg mole	
C <sub>2</sub> H <sub>3</sub> OC <sub>2</sub> Cl <sub>3</sub> C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> Cl <sub>3</sub>	2,2,2-trichloroethanol trichloroacetaldehyde monohydrate in 45 CHCl <sub>3</sub> (trichloromethane)	liq c aq		-132.2 -131.3 -126.2				35.	
C <sub>2</sub> F <sub>3</sub> Cl C <sub>2</sub> F <sub>4</sub> Cl <sub>2</sub> C <sub>2</sub> F <sub>3</sub> Cl <sub>3</sub> C <sub>2</sub> F <sub>2</sub> Cl <sub>4</sub>	trifluorochloroethene 1,1,2,2-tetrafluoro-1,2-dichloroethane 1,1,2-trifluoro-1,2,2-trichloroethane 1,2-difluoro-1,1,2,2-tetrachloroethane	g g liq c						42.	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 23-40. Carbon (at. no., 6; at. wt., 12.010)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1948

Substance		State	$\Delta H_f^\circ$		$\Delta F^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description		At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				
$C_2H_5Br$	bromoethane	g						17.
$C_2H_4Br_2$	1,2-dibromoethane	liq						24.
		liq			-4.94	3.621	53.37	32.51
$C_2H_3OBr$	acetyl bromide	liq						
$C_2H_3O_2Br$	bromoacetic acid	c						
$C_2H_2O_2Br_2$	dibromoacetic acid	c						
$C_2HO_2Br_3$	tribromoacetic acid	c						
$C_2H_3O_2Br_3$	tribromoacetaldehyde monohydrate	c						77.
		aq						
$C_2H_4ClBr$	1-chloro-2-bromoethane	liq						31.

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Substance		Table 23-41. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1948				
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F^\circ$	$\log_{10} K_f$
			At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C) cal/deg mole
$C_2I_4$	tetraiodoethene	c				
$C_2H_5I$	iodoethane	liq		73.		
$C_2H_4I_2$	1,2-diiodoethane	c		-7.4		
$C_2H_3OI$	acetyl iodide	liq		0.1		
$C_2H_3O_2I$	iodoacetic acid	c		-38.3		
$C_2H_4BrI$	1-bromo-2-iodoethane	c		-11.3		
						26.



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 23-43. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)									
June 30, 1948									
Substance									
Formula	Description	State	ΔH <sup>o</sup> <sub>f</sub>		ΔF <sup>o</sup> <sub>f</sub>		log <sub>10</sub> K <sub>f</sub>		C <sub>p</sub> <sup>o</sup>
			At 0°K kcal/mole	kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
C <sub>2</sub> N <sub>2</sub>	cyanogen	g		73.60	70.81	-51.903	57.86	13.60	
C <sub>2</sub> H <sub>3</sub> N	acetonitrile	g		21.0	25.2	-18.47	58.18	12.48	
C <sub>2</sub> H <sub>3</sub> N	methy isocyanide	liq		12.7	24.0	-17.59	34.5		
C <sub>2</sub> H <sub>3</sub> N		g		35.9	40.0	-29.32	58.78	12.76	
C <sub>2</sub> H <sub>7</sub> N	ethylamine	liq		28.1					
C <sub>2</sub> H <sub>7</sub> N	in 100 H <sub>2</sub> O	g		-11.6				16.7	
C <sub>2</sub> H <sub>7</sub> N	dimethylamine	aq		-24.5					
C <sub>2</sub> H <sub>7</sub> N	in 500 H <sub>2</sub> O	g		-6.6	14.1	-10.34	65.30		
C <sub>2</sub> H <sub>8</sub> N <sup>+</sup>	1000 H <sub>2</sub> O	aq		-19.3					
C <sub>2</sub> H <sub>8</sub> N <sup>+</sup>	ethylammonium ion	aq		-19.5					
C <sub>2</sub> H <sub>8</sub> N <sub>2</sub>	dimethylammonium ion	aq		-37.4					
C <sub>2</sub> H <sub>8</sub> N <sub>2</sub>	1,2-ethanediamine (ethylenediamine)	liq		-30.81					
C <sub>2</sub> H <sub>8</sub> N <sub>2</sub> ·H <sub>2</sub> O		aq		-6.36					
C <sub>2</sub> H <sub>8</sub> N <sub>2</sub> ·H <sub>2</sub> O	1,2-ethanediamine monohydrate	c		-13.86					
C <sub>2</sub> H <sub>4</sub> N <sub>4</sub>	dicyandiamide	c		-77.1				38.	
		aq		4.4					
				10.2					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 23-44. Carbon (at. no., 6; at. wt., 12.010)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1948

Substance		State	At 298.16°K (25°C)			
Formula	Description		$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole
$C_2H_2O_3N^-$	oxamate ion	aq	-154.5			
$C_2H_3ON$	methyl isocyanate	liq	-21.5			
$C_2H_3ON$	glycollonitrile	liq	-34.1			
$C_2H_3O_3N$	oxamic acid	c	-161.4			
	in 800 H <sub>2</sub> O	aq	-154.9			
$C_2H_4O_2N^-$	aminoacetate ion	aq	-112.13			
$C_2H_4O_2N^-$	nitroethane ion	aq	-27.			
$C_2H_5ON$	acetamide	c	-76.60			
	in 20 H <sub>2</sub> O	aq	-74.36			
	200 H <sub>2</sub> O	aq	-74.40			
	in C <sub>2</sub> H <sub>6</sub> O (ethanol)		-79.13			
$C_2H_5ON$	acetaldehyde oxime	c	-18.2			
						15.9

## Washington, D.C.

## SERIES I

National Bureau of Standards

Table 23-45. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1948

Substance			At 298.16°K (25°C)					
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} f_f$	$S^\circ$	$C_p^\circ$
			kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
$C_2H_5O_2N$	aminoacetic acid (glycine)	c						
	in 50 H <sub>2</sub> O	aq		-126.33	-88.61	64.950	26.1	24.0
	100 H <sub>2</sub> O	aq		-122.817				
	200 H <sub>2</sub> O	aq		-122.775				
	500 H <sub>2</sub> O	aq		-122.75				
	$\infty$ H <sub>2</sub> O (undissociated)	aq		-122.733				
		aq		-122.721				
	nitroethane	liq		-30.				33.
	nitro form	aq		-29.				
	aci form	aq		-28.				
$C_2H_5O_2N$	ethyl nitrite	g		-24.8				
$C_2H_5O_3N$	ethyl nitrate	liq		-44.3				



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 23-46. Carbon (at. no., 6; at. wt., 12.010)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1948; March 31, 1950

Formula	Substance	State	At 0°K				At 298.16°K (25°C)			
			kcal/mole				kcal/mole			
			$\Delta H_f^0$	$\Delta G_f^0$	$\log_{10} K_f$	$S^0$	$\Delta H_f^0$	$\log_{10} K_f$	$S^0$	$C_p^0$
$C_2H_3O_2N$	ammonium acetate	c								
	in 5 H <sub>2</sub> O	aq		-147.6						
	10 H <sub>2</sub> O	aq		-146.97						
	15 H <sub>2</sub> O	aq		-147.32						
	25 H <sub>2</sub> O	aq		-147.51						
	50 H <sub>2</sub> O	aq		-147.70						
	100 H <sub>2</sub> O	aq		-147.90						
	200 H <sub>2</sub> O	aq		-148.00						
$C_2H_3O_3N$	400 H <sub>2</sub> O	aq		-148.06						
	methylammonium bicarbonate	aq		-148.1						
	ammonium glycolate	c		-195.						
	ammonium bioxalate	aq		-189.8						
$C_2H_5O_4N$	ammonium glyoxylate	aq		-186.8						
	glycinium ion	c		-196.8						
	ammonium glyoxylate	aq		-227.02						
$C_2H_6O_2N^+$		aq		-153.74						
$C_2H_7O_4N$		aq		-226.0						
										36.

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 23-47. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1948							
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole		cal/deg mole
$C_2H_4O_2N_2$	glyoxime		c		-21.1		
$C_2H_4O_2N_2$	oxamide		c		-121.2		
$C_2H_4O_2N_2$	formylurea		c		-117.5		
			aq		-110.8		
$C_2H_4O_3N_2$	nitroacetamide		c		58.		
$C_2H_4O_6N_2$	glycol dinitrate		liq		0.7		
$C_2H_6ON_2$	dimethylnitrosamine		liq		-20.2		
$C_2H_6O_2N_2$	ethylnitramine		liq		-268.72		
$C_2H_6O_4N_2$	ammonium oxalate		c		-260.6		
	in 2100 H <sub>2</sub> O		aq		-340.62		
$C_2H_8O_4N_2 \cdot H_2O$	ammonium oxalate monohydrate		c		-99.8		
$C_2H_7O_4N_5$	guanylurea nitrate		c				
							54.

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 23-48. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1948									
Substance		State	Description	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	kcal/mole	kcal/mole			
$C_2H_5F_2N$	2,2-difluoroethylamine	liq			-115.5				
$C_2H_4OFN$	fluoroacetamide	c			-116.3				
$C_2H_3OF_2N$	difluoroacetamide	c			-165.0				
$C_2H_2OF_3N$	trifluoroacetamide	c							
$C_2H_4O_2F_2N_2$	2,2-difluoroethylnitramine	liq			-116.1				
$C_2H_5ClN$	ethylamine hydrochloride	aq			-77.4				
$C_2H_5ClN$	dimethylamine hydrochloride, in 100 H <sub>2</sub> O	aq			-70.73				
	200 H <sub>2</sub> O	aq			-70.74				
	500 H <sub>2</sub> O	aq			-70.751				
	1000 H <sub>2</sub> O	aq			-70.764				
	5000 H <sub>2</sub> O	aq			-70.794				
	$\infty$ H <sub>2</sub> O	aq			-70.833				
$C_2H_{10}Cl_2N_2$	1,2-ethanediamine dihydrochloride	c			-123.14				
		aq			-116.49				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 23-49. Carbon (at. no., 6; at. wt., 12.010) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)									
June 30, 1948									
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole		cal/deg mole	cal/deg mole	
$C_2H_4OClN$	chloroacetamide		c						
	in 2500 $H_2O$		aq		-81.7				
$C_2H_6O_2ClN$	glycine hydrochloride		aq		-76.4				
$C_2H_5O_2ClN$	ammonium chloroacetate, in 200 $H_2O$		aq		-163.76				
$C_2H_3OCl_2N$	dichloroacetamide		c		-152.5				
	in 5000 $H_2O$		aq		-77.7				
$C_2H_2OCl_3N$	trichloroacetamide		c		-87.				
			aq		-78.7				
$C_2H_4O_2Cl_3N$	ammonium trichloroacetate, in 200 $H_2O$		aq		-157.2				
$C_2H_3SN$	methyl thiocyanate		liq		25.				
$C_2H_3SN$	methyl isothiocyanate		c		18.				
$C_2H_7O_3SN$	2-aminoethylsulfonic acid (taurine)		c		-187.8	100.20	36.8		33.6
	in 50 $H_2O$		aq		-181.66				
	100 $H_2O$		aq		-181.72				
	500 $H_2O$		aq		-181.80				
	$\infty H_2O$		aq		-181.82				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 24-1. Silicon (at. no., 14; at. wt., 28.06)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

September 30, 1948

Substance								
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
			At 298.16°K (25°C)					
			kcal/mole	kcal/mole	kcal/mole			cal/deg mole
Si		q	87.	88.04	77.41	-56.740	40.120	5.318
		c	0.000	0.000	0.000	0.0000	4.47	4.75
		amorp		1.0				
Si +		q	274.93	277.45				
Si ++		q	650.04	654.04				
Si +++		q	1421.6	1427.1				
Si ++++		q	2462.4	2469.4				
Si +++++		g	6307.	6316.				
Si ++++++		g	11037.	11047.				
Si ++++++		g	16718.	16730.				
Si ++++++		q	23708.	23721.				
Si ++++++		g	31802.	31817.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 24-2. Silicon (at. no., 14; at. wt., 28.06)

HEAT OF FORMATION, AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
September 30, 1948 (Corrected)

Substance		State	$\Delta H_f^0$	$\Delta H_f^0$	$\Delta F_f^0$	$\log_{10} K_f$	$S^0$	$C_p^0$
Formula	Description		At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
SiO		g	-27.	-26.72	-32.77	24.020	49.26	7.14
SiO <sub>2</sub>	quartz, II	c		-205.4	-192.4	141.03	10.00	10.62
	cristobalite, II	c		-205.0	-192.1	140.81	10.19	10.56
	tridymite, IV	c		-204.8	-191.9	140.66	10.36	10.66
	vitreous	gls		-202.5	-190.9	139.92	11.2	10.60
		aq		-201.0				
SiH <sub>4</sub>		g		-14.8	-9.4	6.89	48.7	10.24
Si <sub>2</sub> H <sub>6</sub>		g						
Si <sub>3</sub> H <sub>8</sub>		liq						
H <sub>2</sub> SiO <sub>3</sub>		c		-270.7				
H <sub>4</sub> SiO <sub>4</sub>		c		-340.6				

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Table 24-4. Silicon (at. no., 14; at. wt., 28.06) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1948										
Substance		Formula	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
					At 0°K kcal/mole	kcal/mole				
SiI <sub>4</sub>				c						
SiH <sub>3</sub> I				liq						
SiH <sub>2</sub> I <sub>2</sub>				liq						
SiHI <sub>3</sub>				liq						
SiS <sub>2</sub>			white	c			-34.7			
Si <sub>3</sub> N <sub>4</sub>				c			-179.3	113.39	22.4	
(NH <sub>4</sub> ) <sub>2</sub> SiF <sub>6</sub>				c			-629.7			
				aq			-622.0			
SiC				c		-26.46		19.13	3.935	6.37
(CH <sub>3</sub> )SiH <sub>3</sub>			methyl silane	g			-26.7			
(CH <sub>3</sub> ) <sub>2</sub> SiH <sub>2</sub>			dimethyl silane	g						
Si(C <sub>2</sub> H <sub>5</sub> O) <sub>4</sub>			tetraethoxy silane	liq			-318.			



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 25-1. Germanium (at. no., 32; at. wt., 72.60)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
September 30, 1948

Substance		State	Description	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K			At 298.16°K (25°C)					
				kcal/mole	kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
Ge		g		78.0	78.44	69.50	-50.943	40.106	7.346
Ge *		c		0.000	0.000	0.000	0.0000	10.14	6.24
Ge **		q		265.40	267.32				
Ge ***		g		633.10	636.50				
Ge ****		g		1422.2	1427.0				
Ge *****		q		2476.1	2482.4				
Ge *****		q		4631.	4639.				
GeO		g		-22.5	-22.8	-28.2	20.63	52.56	7.39
GeO <sub>2</sub>	tetragonal	c,II							
	glassy	amorp			-128.3				
		aq			-131.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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<p>Table 25-2. Germanium (at. no., 32; at. wt., 72.60) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1948 (Corrected)</p>									
Substance									
Formula	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$		$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
			At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole			
GeH <sub>4</sub>		g							
Ge <sub>2</sub> H <sub>6</sub>		liq							
Ge <sub>3</sub> H <sub>8</sub>		liq							
HGeO <sub>3</sub> <sup>-</sup>		aq							
H <sub>2</sub> GeO <sub>3</sub>		aq		-199.3					
GeF <sub>4</sub>		g							
GeCl		g	32.6	32.45					8.41
GeCl <sub>4</sub>		liq		-130.					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 25-3. Germanium (at. no., 32; at. wt., 72.60) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1948									
Substance				$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 298.16°K (25°C)						
			At 0°K kcal/mole	kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole
GeBr		g		34.8					8.64
GeBr <sub>4</sub>		c							
GeI <sub>4</sub>		c							
GeS		g		1.35					8.05
GeSe		g							9.41
GeTe		g							9.58
Ge <sub>3</sub> N <sub>4</sub>		c		-14.8					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 26-1. Tin (at. no., 50; at. wt., 118.70) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1948 (Corrected)									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				
Sn		gray, cubic	g		72.		-47.5	40.245	5.081
		white, tetragonal	c, III		0.6		-0.806	10.7	6.16
Sn *			c, II		0.000		0.0000	12.3	6.30
Sn **			g		242.57				
Sn ***			g		581.4				
Sn ****			g		1289.8				
Sn *****			g		2231.				
			g		4094.				
SnO			g					54.98	21.05
			c		-68.4	-61.5	45.08	13.5	10.6
SnO <sub>2</sub>			c		-138.8	-124.2	91.037	12.5	12.57

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 26-2. Tin (at. no., 50; at. wt., 118.70)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

September 30, 1948

Substance		State	Description	At 298.16°K (25°C)				S°	C <sub>p</sub> <sup>o</sup>
Formula	ΔH <sup>o</sup>			ΔH <sup>o</sup>	ΔF <sup>o</sup>	log <sub>10</sub> f <sup>o</sup>			
							kcal/mole		
SnH <sub>4</sub>		g							
Sn(OH) <sup>+</sup>	std. state, hyp. m = 1	aq			-61.1	44.79			
HSnO <sub>2</sub> <sup>-</sup>	std. state, hyp. m = 1	aq			-97.9	71.76			
Sn(OH) <sub>2</sub>		c		-138.3	-117.6	86.199	23.1		
Sn(OH) <sub>4</sub>		c		-270.5					
SnF <sub>4</sub>		c							
SnF <sub>6</sub> <sup>--</sup>	std. state, hyp. m = 1	aq		-474.7					
H <sub>2</sub> SnF <sub>6</sub>		aq		-473.1					
SnCl <sub>2</sub>		c		-83.6					
	in aq HCl	aq		-84.0					
		c		-225.9					
SnCl <sub>2</sub> ·2H <sub>2</sub> O		liq		-130.3	-113.3	89.047	61.8		39.5
SnCl <sub>4</sub>		aq		-160.2					
SnOCl <sub>2</sub>	in aq HCl	aq		-147.5					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards Washington, D.C.

Table 26-3. $T_{in}$ (at. no., 50; at. wt., 118.70) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1948 (Corrected)									
Substance									
Formula	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$		$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
			At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole			
SnBr <sub>2</sub>		c		-63.6					
SnBr <sub>4</sub>		aq		-62.0					
		c, l		-97.1					
SnBr <sub>4</sub> ·8H <sub>2</sub> O		aq		-113.0					
SnI <sub>2</sub>		liq		-656.6					
		c		-34.4					
		aq		-28.6					
SnS		c		-18.6	-19.7	14.44	23.6		
Sn(SO <sub>4</sub> ) <sub>2</sub>		c		-393.4					
SnCl <sub>2</sub> ·2½NH <sub>3</sub>		c		-149.6					
SnCl <sub>2</sub> ·4NH <sub>3</sub>		c		-180.1					
SnCl <sub>2</sub> ·9NH <sub>3</sub>		c		-273.3					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 26-4. Tin (at. no., 50; at. wt., 118.70)

HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
September 30, 1948; March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole		cal/deg mole	cal/deg mole
$\text{SnBr}_2 \cdot \text{NH}_3$			c		-93.8			
$\text{SnBr}_2 \cdot 2\text{NH}_3$			c		-118.4			
$\text{SnBr}_2 \cdot 3\text{NH}_3$			c		-142.3			
$\text{SnBr}_2 \cdot 5\text{NH}_3$			c		-184.0			
$\text{SnBr}_2 \cdot 9\text{NH}_3$			c		-258.3			
$\text{SnI}_2 \cdot \text{NH}_3$			c		-61.4			
$\text{SnI}_2 \cdot 2\text{NH}_3$			c		-85.7			
$\text{SnI}_2 \cdot 3\text{NH}_3$			c		-109.0			
$\text{SnI}_2 \cdot 5\text{NH}_3$			c		-151.5			
$\text{SnI}_2 \cdot 9\text{NH}_3$			c		-227.2			
$\text{SnCl}_4 \cdot \frac{1}{2}\text{PH}_3$			c		-153.6			
$\text{SnH}_{12}\text{C}_4$	tetramethyltin		liq					
$\text{SnH}_{20}\text{C}_8$	tetraethyltin		liq		-49.			
$\text{SnH}_{28}\text{C}_{12}$	tetra-n-propyltin		liq		-56.			
$\text{SnH}_{36}\text{C}_{16}$	tetra-n-butyltin		liq		-94.			
$\text{SnH}_{44}\text{C}_{20}$	tetra-n-amyltin		liq		-132.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 27-1. Lead (at. no., 82; at. wt., 207.21)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
September 30, 1948

Substance			Description	State	$\Delta H_f^o$	$\Delta F_f^o$	$\log_{10} K_f$	$S^o$	$C_p^o$
Formula		At 298.16°K (25°c)							
		kcal/mole			kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
Pb		g	46.5	46.34	38.47	-28.198	41.890	4.968	
Pb +		c	0.000	0.000	0.000	0.0000	15.51	6.41	
Pb ++		q	217.50	218.82					
		g	564.09	566.89					
Pb +++		aq		0.39	-5.81	4.259	5.1		
Pb ++++		g	1303.8	1308.1					
Pb +++++		g	2278.2	2284.0					
Pb <sub>2</sub>		g	3886.	3894.					
PbO		g	79.	78.			57.4		
	red	c, II		-52.40	-45.25	33.168	16.2		
	yellow	c, I		-52.07	-45.05	33.021	16.6	11.60	
PbO <sub>2</sub>		c		-66.12	-52.34	38.364	18.3	15.4	
Pb <sub>2</sub> O		c		-51.2					
Pb <sub>3</sub> O <sub>4</sub>		c		-175.6	-147.6	108.19	50.5	35.14	



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 27-2. Lead (at. no., 82; at. wt., 207.21)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

September 30, 1948

Substance		State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description		At 0°K	At 298.16°K (25°C)				
			kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
HPbO <sub>2</sub> <sup>-</sup>	std. state, hyp. m = 1	aq		-81.00				
Pb(OH) <sub>2</sub>		c		-123.0	-100.6	73.738	21.	
PbF <sub>2</sub>		c		-158.5	-148.1	108.55	29.	
PbF <sub>4</sub>		aq		-155.7				
		c		-222.3				
PbCl <sub>2</sub>		c		-85.85	-75.04	55.003	32.6	18.4
PbCl <sub>2</sub> ·PbO	std. state, hyp. m = 1	aq		-79.65	-68.51	50.217	31.4	
PbCl <sub>2</sub> ·2PbO		c		-144.5				
PbCl <sub>2</sub> ·3PbO		c		-199.2				
		c		-252.7				

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Table 27-3. Lead (at. no., 82; at. wt., 207.21) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1948									
Substance		State		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	At 0°K		At 298.16°K (25°C)					
		kcal/mole	kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole	cal/deg mole
PbBr <sub>2</sub>	std. state, hyp. m=1			-66.21	-62.24	45.621	38.6	19.15	
Pb(BrO <sub>3</sub> ) <sub>2</sub>				-57.41	-54.95	40.278	43.7		
PbBr <sub>2</sub> ·PbO	std. state, hyp. m=1			-122.3					
PbBr <sub>2</sub> ·2PbO				-176.4					
PbBr <sub>2</sub> ·3PbO				-230.8					
PbI <sub>2</sub>				-41.85	-41.53	30.441	42.3		
PbI <sub>3</sub> <sup>-</sup>	std. state, hyp. m=1			-26.35	-30.51	22.363	57.8		
PbI <sub>4</sub> <sup>--</sup>	std. state, hyp. m=1				-48.55				
Pb(IO <sub>3</sub> ) <sub>2</sub>					-63.02				
PbI <sub>2</sub> ·HI·5H <sub>2</sub> O	std. state, hyp. m=1			-400.1					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 27-4. Lead (at. no., 82; at. wt., 207.21) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1948									
Substance		State	$\Delta H_f^0$ kcal/mole	$\Delta H_f^0$ kcal/mole	$\Delta F_f^0$ kcal/mole	$\log_{10} K_f$	$S^0$	$C_p^0$	
Formula	Description								
PbS		c		-22.54	-22.15	16.236	21.8	11.83	
PbSO <sub>4</sub>		c, II		-219.50	-193.89	142.119	35.2	24.9	
PbS <sub>2</sub> O <sub>3</sub>		c		-158.					
PbS <sub>3</sub> O <sub>6</sub>		c		-289.					
		aq		-295.					
PbSO <sub>4</sub> ·PbO		c		-282.5					
PbSO <sub>4</sub> ·2PbO		c		-344.					
PbSO <sub>4</sub> ·3PbO		c		-403.					
PbSO <sub>4</sub> ·2HCl		c							
PbSe		c		-18.					
PbSeO <sub>4</sub>		c		-148.					
PbTe		c		-17.5					

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Table 27-5. Lead (at. no., 82; at. wt., 207.21)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
September 30, 1948

Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				
Pb(N <sub>3</sub> ) <sub>2</sub>			c		104.3				
Pb(NO <sub>3</sub> ) <sub>2</sub>			c		-107.35				
	std. state, hyp. m = 1		aq		-98.35		42.982	75.1	
	in 400 H <sub>2</sub> O		aq		-100.10				
	∞ H <sub>2</sub> O		aq		-98.35				
Pb(OH)NO <sub>3</sub>			c		-117.9				
PbCl <sub>2</sub> ·NH <sub>3</sub>			c		-109.9				
PbCl <sub>2</sub> ·1½NH <sub>3</sub>			c		-121.1				
PbCl <sub>2</sub> ·2NH <sub>3</sub>			c		-132.1				
PbCl <sub>2</sub> ·8NH <sub>3</sub>			c		-249.0				
2PbCl <sub>2</sub> ·NH <sub>4</sub> Cl			c		-247.4				
PbBr <sub>2</sub> ·NH <sub>3</sub>			c		-92.3				
PbBr <sub>2</sub> ·2NH <sub>3</sub>			c		-114.7				
PbBr <sub>2</sub> ·3NH <sub>3</sub>			c		-135.3				
PbBr <sub>2</sub> ·5½NH <sub>3</sub>			c		-186.0				
PbBr <sub>2</sub> ·8NH <sub>3</sub>			c		-234.7				

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Table 27-6. Lead (at. no., 82; at. wt., 207.21)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

September 30, 1948

Substance		State	Description	$\Delta H_f^\circ$		$\Delta F^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K kcal/mole			At 298.16°K (25°C) kcal/mole					
					kcal/mole				
$PbI_2 \cdot \frac{1}{2}NH_3$		c		-54.8					
$PbI_2 \cdot NH_3$		c		-66.8					
$PbI_2 \cdot 2NH_3$		c		-89.1					
$PbI_2 \cdot 5NH_3$		c		-149.8					
$PbI_2 \cdot 8NH_3$		c		-205.1					
$3PbI_2 \cdot 4NH_4I$		c		-314.3					
$3PbI_2 \cdot 4NH_4I \cdot 6H_2O$		c		-734.3					
$PbSO_4 \cdot (NH_4)_2SO_4$		c		-503.34					
$Pb_3(PO_4)_2$		c		-620.3	-581.4	426.16	84.45	61.2	
$PbHPO_3$		c		-234.5					
$3PbI_2 \cdot PI_3$		c		-131.8					
$3PbI_2 \cdot PI_3 \cdot 12H_2O$		c		-977.7					

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Table 27-7. Lead (at. no., 82; at. wt., 207.21)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

September 30, 1948

Substance			State	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	Description	At 298.16°K (25°C)							
		At 0°K kcal/mole							kcal/mole
$3PbI_2 \cdot AsI_3$		c	-111.3						
$3PbI_2 \cdot AsI_3 \cdot 12H_2O$		c	-952.8						
$3PbI_2 \cdot SbI_3$		c	-127.7						
$3PbI_2 \cdot SbI_3 \cdot 12H_2O$		c	-977.6						
$3PbI_2 \cdot BiI_3$		c							
$3PbI_2 \cdot BiI_3 \cdot 12H_2O$		c							
$PbCO_3$		c	-167.3	-149.7	109.73	31.3	20.9		
$PbC_2O_4$	lead oxalate	c	-205.1						
$PbO \cdot PbCO_3$		c	-220.0	-195.6	143.37	48.5	32.2		
$2PbO \cdot PbCO_3$		c	-273.	-242.	177.4	65.			
$Pb(C_2H_5)_4$	tetraethyllead	liq	52.						
$Pb(CHO_2)_2$	lead formate	c	-200.1						
		aq	-194.0						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 27-8. Lead (at. no., 82; at. wt., 207.21) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1948									
Substance		Description	State	$\Delta F^{\circ}$	$\Delta H^{\circ}$	$\Delta P^{\circ}$	$\log_{10} P^{\circ}$	$S^{\circ}$	$C_p^{\circ}$
Formula	At 0°K kcal/mole			kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole	
$Pb(C_2H_3O_2)_2$		lead acetate	c		-230.5				
$Pb(C_2H_3O_2)_2 \cdot 3H_2O$			aq		-232.6				
$Pb(C_2H_3O_3)_2$		lead glycolate	c		-443.1				
			c		-316.3				
			aq		-311.5				
$Pb(CN)_2 \cdot 2PbO \cdot H_2O$			c		-123.0				
$Pb(CNS)_2$			c		27.5				
$PbSiO_3$			c		-258.8	-239.0	175.18	27.	
$Pb_2SiO_4$			amorp		-256.9				
			c		-312.7	-285.7	209.41	43.	
$PbI_2 \cdot SnI_2$			amorp		-309.0				
$PbI_2 \cdot SnI_2 \cdot 8H_2O$			c		-64.2				
			c		-631.6				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 28-1. Gallium (at. no., 31; at. wt., 69.72)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1948

Substance		$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$G_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)		
			kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
Ga		g				
Ga <sup>+</sup>		c, l	0.000	66.0	-41.78	6.053
Ga <sup>++</sup>		g		0.000	0.0000	6.35
Ga <sup>+++</sup>		g		205.77		
		g		680.21		
		g		1388.5		
	std. state, hyp. m=1	aq		-50.4	26.83	-83.
Ga <sup>++++</sup>		g		2868.		
GaO		g		66.		
GaO <sub>3</sub> <sup>---</sup>		aq		-148.	108.5	
Ga <sub>2</sub> O	std. state, hyp. m=1	c		-82.		
Ga <sub>2</sub> O <sub>3</sub>		c		-258.		20.



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 28-2. Gallium (at. no., 31; at. wt., 69.72)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1948

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
Ga(OH) <sup>++</sup>	std. state, hyp. m=1		aq			-69.8	65.82		
HGaO <sub>3</sub> <sup>--</sup>	std. state, hyp. m=1		aq			-164.	120.2		
Ga(OH) <sub>2</sub> <sup>+</sup>	std. state, hyp. m=1		aq			-142.1	104.16		
H <sub>2</sub> GaO <sub>3</sub> <sup>-</sup>	std. state, hyp. m=1		aq			-178.	130.5		
Ga(OH) <sub>3</sub>			c			-199.	145.9		
GaCl			g		9.				
GaCl <sub>3</sub>			c		-125.4				
	std. state, hyp. m=1		aq		-170.5	-130.7	95.801	-43.	
	in aq HCl		aq		-164.4				
GaBr			g		11.				
GaBr <sub>3</sub>			c		-92.4				
	in aq HCl		aq		-130.8				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 28-3. Gallium (at. no., 31; at. wt., 69.72)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1948

Substance		$\Delta H^{\circ}$			$\Delta F^{\circ}$	$\log_{10} K^{\circ}$	$S^{\circ}$	$C_p^{\circ}$
Formula	Description	State	At 0°K		At 298.16°K (25°C)		cal/deg mole	cal/deg mole
			kcal/mole	kcal/mole	kcal/mole	kcal/mole		
GaI	in aq HCl	a		27.				62.
GaI <sub>3</sub>		c		-51.2				
		aq		-84.2				
Ga <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c						
GaN		c		-25.				
GaCl <sub>3</sub> ·NH <sub>3</sub>		c		-170.8				
GaCl <sub>3</sub> ·3NH <sub>3</sub>		c		-231.3				
GaCl <sub>3</sub> ·5NH <sub>3</sub>		c		-285.6				
GaCl <sub>3</sub> ·6NH <sub>3</sub>		c		-308.0				
GaCl <sub>3</sub> ·7NH <sub>3</sub>		c		-328.5				
GaCl <sub>3</sub> ·14NH <sub>3</sub>		c		-459.7				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 28-4. Gallium (at. no., 31; at. wt., 69.72)  
HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1948

Substance		State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	At 298.16°K (25°C)		$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description					$\log_{10} K_f$			
$\text{GaBr}_3 \cdot \text{NH}_3$		c		-135.2					
$\text{GaBr}_3 \cdot 5\text{NH}_3$		c		-256.1					
$\text{GaBr}_3 \cdot 6\text{NH}_3$		c		-280.4					
$\text{GaBr}_3 \cdot 7\text{NH}_3$		c		-301.8					
$\text{GaBr}_3 \cdot 9\text{NH}_3$		c		-340.5					
$\text{GaBr}_3 \cdot 14\text{NH}_3$		c		-433.7					
$\text{GaI}_3 \cdot \text{NH}_3$		c		-90.6					
$\text{GaI}_3 \cdot 5\text{NH}_3$		c		-211.5					
$\text{GaI}_3 \cdot 6\text{NH}_3$		c		-237.7					
$\text{GaI}_3 \cdot 7\text{NH}_3$		c		-259.6					
$\text{GaI}_3 \cdot 9\text{NH}_3$		c		-298.7					
$\text{GaI}_3 \cdot 13\text{NH}_3$		c		-374.8					
$\text{GaI}_3 \cdot 20\text{NH}_3$		c		-506.0					
$\text{Ga}_2(\text{C}_2\text{O}_4)_3$	gallium oxalate	c		-515.					
$\text{Ga}_2(\text{C}_2\text{O}_4)_3 \cdot 2\text{H}_2\text{O}$		c		-689.					
$\text{Ga}_2(\text{C}_2\text{O}_4)_3 \cdot 4\text{H}_2\text{O}$		c		-847.					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 29-1. Indium (at. no., 49; at. wt., 114.76)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1948

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	At 298.16°K (25°C) kcal/mole	At 298.16°K (25°C) cal/deg mole	At 298.16°K (25°C) cal/deg mole	At 298.16°K (25°C) cal/deg mole
In			g	0.000	58.2	49.6	-36.36	41.51	6.015
In <sup>+</sup>			c		0.000	0.000	0.0000	12.5	6.55
In <sup>++</sup>			g		193.09				
In <sup>+++</sup>			g		629.62				
			g		1277.5				
		std. state, hyp. m=1	aq		-23.7	-32.0	23.46	-62.	
In <sup>++++</sup>			g		2616.				
In <sup>+++++</sup>			g						
InO			g		91.				22.
In <sub>2</sub> O <sub>3</sub>			c		-222.5				
InH			g		51.	45.	-33.0	49.6	7.06
In(OH) <sup>++</sup>			aq			-55.5	40.68		
In(OH) <sub>3</sub>		std. state, hyp. m=1	c		-214.	-182.	133.4	25.	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Substance		State	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} f_f$	$S^\circ$	$C_p^\circ$
Formula	Description		At 0°K	At 298.16°K (25°C)				
InCl		g		-18.	-23.	16.9	59.3	8.60
InCl <sub>2</sub>		c, II		-44.5				
InCl <sub>3</sub>		c		-86.8				
		c		-128.4				
		aq		-152.0				
	in aq HCl	aq		-147.2				
InBr		g		10.	1.	-0.73	62.0	8.77
InBr <sub>3</sub>		c		-96.5				
	in aq HCl	aq		-113.9				
InI		g		20.	9.	-6.6	63.8	8.83
InI <sub>3</sub>		c		-55.0				
	in aq HCl	aq		-67.3				
In <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c		-695.				67.
InN		c		-4.8				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards Washington, D.C.

Table 29-3. Indium (at. no., 49; at. wt., 114.76) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1948											
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
				At 0°K						At 298.16°K (25°C)	
				kcal/mole	kcal/mole					kcal/mole	cal/deg mole
InCl <sub>3</sub> ·NH <sub>3</sub>			c		-163.1						
InCl <sub>3</sub> ·2NH <sub>3</sub>			c		-195.5						
InCl <sub>3</sub> ·3NH <sub>3</sub>			c		-227.9						
InCl <sub>3</sub> ·5NH <sub>3</sub>			c		-276.6						
InCl <sub>3</sub> ·7NH <sub>3</sub>			c		-315.9						
InCl <sub>3</sub> ·15NH <sub>3</sub>			c		-465.4						
InBr <sub>3</sub> ·3NH <sub>3</sub>			c		-193.3						
InBr <sub>3</sub> ·5NH <sub>3</sub>			c		-246.9						
InBr <sub>3</sub> ·7NH <sub>3</sub>			c		-289.4						
InBr <sub>3</sub> ·15NH <sub>3</sub>			c		-444.9						
InI <sub>3</sub> ·NH <sub>3</sub>			c		-87.6						
InI <sub>3</sub> ·2NH <sub>3</sub>			c		-117.7						
InI <sub>3</sub> ·5NH <sub>3</sub>			c		-201.9						
InI <sub>3</sub> ·7NH <sub>3</sub>			c		-251.0						
InI <sub>3</sub> ·9NH <sub>3</sub>			c		-290.9						
InI <sub>3</sub> ·13NH <sub>3</sub>			c		-365.6						
InI <sub>3</sub> ·21NH <sub>3</sub>			c		-513.2						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 30-1. Thallium (at. no., 81; at. wt., 204.39)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1948; March 31, 1950

	Formula	Substance	State	Description	At 298.16°K (25°C)				
					$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$
					At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
Tl			g			43.34	35.05	-25.691	43.23
Tl <sup>+</sup>			c, II		0.000	0.000	0.000	0.0000	15.4
			g			185.64			
Tl <sup>++</sup>		std. state, hyp. m = 1	aq			1.38	-7.755	5.6843	30.4
Tl <sup>+++</sup>			g			657.6			
			g			1346.9			
Tl <sup>++++</sup>		std. state, hyp. m = 1	aq			27.7	50.0	-36.65	-106.
			g			2514.			
Tl <sub>2</sub> O			c			-41.9	-32.5	23.82	23.8
TlH			g			48.	42.	-30.8	51.39
TlOH			c			-56.9	-45.5	33.35	17.3
		in 500 H <sub>2</sub> O	aq			-53.7			
		2000 H <sub>2</sub> O	aq			-54.0			
Tl(OH) <sub>3</sub>			c			-122.6			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 30-2. Thallium (at. no., 81; at. wt., 204.39) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1948; March 31, 1950											
Substance			State	Description	At 0°K kcal/mole	ΔHf° kcal/mole	ΔFf° kcal/mole	log10Kf	S° cal/deg mole	Cp° cal/deg mole	
Formula											
TlF		g				-33.					
HTlF2		aq				-77.3					
		aq				-152.5					
TlCl		g				-16.	-22.	16.1	61.1	8.66	
		c				-48.99	-44.19	32.391	25.9		
TlCl3	std. state, hyp. m=1	aq				-38.64	-39.105	28.6634	43.6		
		c				-83.9					
		aq				-92.3					
TlCl3·4H2O		c				-367.7					
TlBr		g				-5.	-14.	10.3	63.8	8.81	
		c				-41.2	-39.7	29.10	26.6		
	std. state, hyp. m=1	aq				-27.5	-32.3	23.68	49.7		
TlBr3		c				-59.0					
TlBr3·4H2O		c				-334.6					
TlBrO3		c				-24.0					
	std. state, hyp. m=1	aq				-10.2	1.3	-0.953	68.6		



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 30-3. Thallium (at. no., 81; at. wt., 204.39) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1948; March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole				
TlClBr <sub>2</sub>			aq		-70.1				
TlClBr <sub>2</sub> ·4H <sub>2</sub> O			c		-346.3				
TlI		yellow	g		7.	-4.	2.9	65.6	8.66
			c, II		-29.7	-29.7	21.77	29.4	
		std. state, hyp. m = 1	aq		-12.0	-19.9	14.59	56.5	
TlIO <sub>3</sub>			c			-47.6	34.89		
		std. state, hyp. m = 1	aq		-53.5	-40.1	29.33	58.1	
Tl <sub>2</sub> S			c		-20.8				
Tl <sub>2</sub> SO <sub>4</sub>			c		-221.7				
		std. state, hyp. m = 1	aq		-214.14	-192.85	141.356	65.1	
TlHSO <sub>4</sub>			aq		-210.3				
Tl <sub>2</sub> Se			c		-18.				
Tl <sub>2</sub> Te			c		-7.				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 30-4. Thallium (at. no., 81; at. wt., 204.39) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)										
December 31, 1948										
Substance		Formula	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
					At 0°K					
					kcal/mole	kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
TlN <sub>3</sub>	thallous azide	c	56.3							
TlNO <sub>3</sub>	std. state, hyp. m = 1	c, III aq	-58.01 -47.99	-36.07 -34.22	26.439 25.063	38.2 65.6				
TlCl·3NH <sub>3</sub>		c	-103.4							
TlCl <sub>3</sub> ·3NH <sub>3</sub>		c	-145.							
TlBr·3NH <sub>3</sub>		c	-95.6							
TlI·3NH <sub>3</sub>		c	-84.1							
Tl(C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> )	thallous acetate	aq	-115.5							
TlOC <sub>2</sub> H <sub>5</sub>	thallous ethoxide in C <sub>2</sub> H <sub>6</sub> O (ethanol)	liq	-56.5 -56.0							
TlONC	thallous fulminate	c	27.6							

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 31-1. Zinc (at. no., 30; at. wt. 65.38) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1948; March 31, 1950										
Substance			State	Description	At 0°K kcal/mole	ΔH <sup>o</sup> kcal/mole	ΔF <sup>o</sup> kcal/mole	log <sub>10</sub> K <sup>o</sup> (25°C)	S <sup>o</sup> cal/deg mole	C <sub>p</sub> <sup>o</sup> cal/deg mole
Formula										
Zn			g				31.19	-16.631	38.45	4.968
			c		0.000		0.000	0.0000	9.95	5.99
		in Hg	g				0.00	0.0000		
Zn <sup>+</sup>			g				249.251			
Zn <sup>++</sup>			g				664.902			
		std. state, hyp. m = 1	aq				-36.43	25.7894	-25.45	
Zn <sup>+++</sup>			g				1581.			
Zn <sup>++++</sup>			g							
ZnO			c				-83.17	-76.05	10.5	9.62
ZnO <sub>2</sub> ·2H <sub>2</sub> O			c				-207.9			
Zn <sub>3</sub> O <sub>5</sub> ·2H <sub>2</sub> O			c				-371.1			
Zn <sub>3</sub> O <sub>5</sub> ·3H <sub>2</sub> O			c				-437.7			
ZnH			g				54.4	47.5	48.70	7.71
Zn(OH) <sub>2</sub>		stable	c, I				-153.5	-34.82		17.3
		unstable	c, II				-153.1			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

<p>Table 31-2. Zinc (at. no., 30; at. wt., 65.38) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1948</p>									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	kcal/mole				
ZnF <sub>2</sub>			aq		-187.9				
ZnCl			g		1.		3.7	58.3	8.45
ZnCl <sub>2</sub>			c		-99.40		64.6897	25.9	18.3
	std. state, hyp. m = 1		aq		-116.48		71.7476	0.89	
	in 4 H <sub>2</sub> O		aq		-106.30				
	5 H <sub>2</sub> O		aq		-107.21				
	7 H <sub>2</sub> O		aq		-108.13				
	10 H <sub>2</sub> O		aq		-108.98				
	15 H <sub>2</sub> O		aq		-109.88				
	25 H <sub>2</sub> O		aq		-111.00				
	50 H <sub>2</sub> O		aq		-112.55				
	100 H <sub>2</sub> O		aq		-114.00				
	200 H <sub>2</sub> O		aq		-114.86				
	400 H <sub>2</sub> O		aq		-115.23				
	800 H <sub>2</sub> O		aq		-115.43				
	$\infty$ H <sub>2</sub> O		aq		-116.48				
	in C <sub>2</sub> H <sub>5</sub> O (ethanol)		aq		-109.2				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 31-3. Zinc (at. no. 30; at. wt., 65.38)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1948

Substance		State	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\log_{10} K_f$ (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description							
$Zn(ClO_3)_2$		c						
$Zn(ClO_4)_2$		c						
$ZnCl_2 \cdot 3ZnO \cdot 5H_2O$		c		-708.1				
$ZnCl_2 \cdot 4ZnO \cdot 11H_2O$		c		-1206.6				
$ZnCl_2 \cdot 5ZnO \cdot 8H_2O$		c		-1086.4				
$ZnCl_2 \cdot 8ZnO \cdot 10H_2O$		c		-1483.4				
$ZnBr$		g					60.8	8.57
$ZnBr_2$		c		-78.17	-74.142	54.3450	32.84	
	std. state, hyp. m = 1	aq		-94.23	-84.332	61.8142	13.13	
$ZnBr_2 \cdot 2H_2O$		c		-220.9	-190.6	139.71	56.2	
$ZnBr_2 \cdot 4ZnO \cdot 13H_2O$		c		-1324.7				
$ZnI$		g		15.	3.	-2.2	63.0	8.76
$ZnI_2$		c		-49.98	-50.01	36.657	38.0	
	std. state, hyp. m = 1	aq		-63.17	-59.884	43.8941	26.83	
$Zn(IO_3)_2$		c						
$ZnI_2 \cdot 5ZnO \cdot 11H_2O$		c		-1218.5				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 31-4. Zinc (at. no., 30; at. wt., 65.38) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1948									
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
				At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole			
ZnS	sphalerite		g		-14.				
	wurtzite		c, II		-48.5				
ZnSO <sub>3</sub>			c, I		-45.3				
			c						
ZnSO <sub>4</sub>			c		-233.88				
	std. state, hyp. m=1				-253.33				
	in 18 H <sub>2</sub> O		aq		-251.58				
	20 H <sub>2</sub> O		aq		-251.71				
	25 H <sub>2</sub> O		aq		-251.68				
	30 H <sub>2</sub> O		aq		-251.95				
	40 H <sub>2</sub> O		aq		-252.01				
	50 H <sub>2</sub> O		aq		-252.057				
	100 H <sub>2</sub> O		aq		-252.139				
	200 H <sub>2</sub> O		aq		-252.233				
	300 H <sub>2</sub> O		aq		-252.283				
	500 H <sub>2</sub> O		aq		-252.349				
	1000 H <sub>2</sub> O		aq		-252.462				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 31-5. Zinc (at. no., 30; at. wt., 65.38)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1943; March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
ZnSO <sub>4</sub>	in	1500 H <sub>2</sub> O	aq		-252.532				
		2000 H <sub>2</sub> O	aq		-252.585				
		3000 H <sub>2</sub> O	aq		-252.661				
		5000 H <sub>2</sub> O	aq		-252.750				
		10,000 H <sub>2</sub> O	aq		-252.866				
		20,000 H <sub>2</sub> O	aq		-252.987				
		30,000 H <sub>2</sub> O	aq		-253.050				
		50,000 H <sub>2</sub> O	aq		-253.120				
		100,000 H <sub>2</sub> O	aq		-253.198				
		200,000 H <sub>2</sub> O	aq		-253.254				
		500,000 H <sub>2</sub> O	aq		-253.290				
		1,000,000 H <sub>2</sub> O	aq		-253.302				
		$\infty$ H <sub>2</sub> O	aq		-253.33				
	ZnSO <sub>4</sub> ·H <sub>2</sub> O		c		-310.6	-269.9	197.83	34.9	34.7
	ZnSO <sub>4</sub> ·6H <sub>2</sub> O		c		-663.3	-555.0	406.81	86.8	80.8
	ZnSO <sub>4</sub> ·7H <sub>2</sub> O		c		-735.1	-611.9	448.51	92.4	93.7

## National Bureau of Standards

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Table 31-6. Zinc (at. no., 30; at. wt. 65.38)

HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1948

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} f_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)			
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
$ZnS_2O_3$		c					
$ZnS_2O_4$		aq		-200.4			
$ZnS_2O_6$		c					
		aq		-316.8			
$ZnS_2O_6 \cdot 6H_2O$		c		-728.9			
$ZnS_4O_6$		c					
$Zn(HSO_3)_2$		aq		-336.6			
$ZnSe$		c		-34.			
$ZnTe$		g		30.			
		c		-30.	-29.	21.3	19.
$Zn(N_3)_2$	zinc azide	c		50.8			
$Zn_3N_2$		c		-6.9			25.7



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 31-7. Zinc (at. no., 30; at. wt., 65.38)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1948

Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} f_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K kcal/mole			kcal/mole	kcal/mole	At 298.16°K (25°C)			
						cal/deg mole	cal/deg mole	cal/deg mole	
Zn(NO <sub>2</sub> ) <sub>2</sub> Zn(NO <sub>3</sub> ) <sub>2</sub>	in	1.5 H <sub>2</sub> O	c						
	2	H <sub>2</sub> O	c		-115.12				
	2	H <sub>2</sub> O	aq		-120.9				
	3	H <sub>2</sub> O	aq		-122.8				
	3	H <sub>2</sub> O	aq		-125.6				
	4	H <sub>2</sub> O	aq		-127.9				
	5	H <sub>2</sub> O	aq		-129.63				
	6	H <sub>2</sub> O	aq		-131.05				
	7	H <sub>2</sub> O	aq		-132.13				
	8	H <sub>2</sub> O	aq		-132.93				
	10	H <sub>2</sub> O	aq		-133.88				
	15	H <sub>2</sub> O	aq		-134.79				
	20	H <sub>2</sub> O	aq		-135.03				
	30	H <sub>2</sub> O	aq		-135.13				
	50	H <sub>2</sub> O	aq		-135.08				
	100	H <sub>2</sub> O	aq		-134.99				
200	H <sub>2</sub> O	aq		-135.03					
400	H <sub>2</sub> O	aq		-135.18					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 31-8. Zinc (at. no., 30; at. wt., 65.38) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1948									
Substance		State		At 0°K		At 298.16°K (25°C)		log <sub>10</sub> K <sub>f</sub>	
Formula	Description	State		ΔH <sup>o</sup> kcal/mole	ΔH <sup>o</sup> kcal/mole	ΔF <sup>o</sup> kcal/mole	S <sup>o</sup> cal/deg mole	C <sub>p</sub> <sup>o</sup> cal/deg mole	
Zn(NO <sub>3</sub> ) <sub>2</sub> ·H <sub>2</sub> O		c			-191.99				
Zn(NO <sub>3</sub> ) <sub>2</sub> ·2H <sub>2</sub> O		c			-264.94				
Zn(NO <sub>3</sub> ) <sub>2</sub> ·4H <sub>2</sub> O		c			-405.76				
Zn(NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O		c			-550.92				
Zn(NH <sub>2</sub> ) <sub>2</sub>		c			-38.8				
ZnCl <sub>2</sub> ·NH <sub>3</sub>		c			-135.5				
ZnCl <sub>2</sub> ·2NH <sub>3</sub>		c			-165.7				
ZnCl <sub>2</sub> ·2NH <sub>4</sub> Cl		c							
ZnCl <sub>2</sub> ·3NH <sub>4</sub> Cl		c							
ZnCl <sub>2</sub> ·4NH <sub>3</sub>		c			-211.5				
ZnCl <sub>2</sub> ·2N <sub>2</sub> H <sub>4</sub>		c			-117.8				
ZnCl <sub>2</sub> ·5NH <sub>3</sub> ·H <sub>2</sub> O		c			-295.3				
ZnCl <sub>2</sub> ·6NH <sub>3</sub>		c			-255.2				
ZnCl <sub>2</sub> ·10NH <sub>3</sub>		c			-327.6				
2ZnCl <sub>2</sub> ·4NH <sub>3</sub> ·H <sub>2</sub> O		c			-362.4				
3ZnCl <sub>2</sub> ·6NH <sub>4</sub> Cl·H <sub>2</sub> O		c			-831.3				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 31-9. Zinc (at. no., 30; at. wt., 65.38)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1948

Formula	Substance	Description	State	$\Delta H^{\circ}$		$\Delta F^{\circ}$	$\log_{10} K^{\circ}$	$S^{\circ}$	$C_p^{\circ}$
				At 0°K	At 298.16°K (25°C)				
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
$2ZnCl_2 \cdot 6NH_4Cl \cdot ZnO$			c		-891.4				
$3ZnCl_2 \cdot 10NH_4Cl \cdot ZnO$			c		-1140.8				
$6ZnCl_2 \cdot 12NH_3 \cdot ZnO$			c		-1349.1				
$ZnBr_2 \cdot NH_3$			c		-113.3				
$ZnBr_2 \cdot 2NH_3$			c		-144.3				
$ZnBr_2 \cdot 2N_2H_4$			c		-94.2				
$ZnBr_2 \cdot 4NH_3$			c		-193.3				
$ZnBr_2 \cdot 6NH_3$			c		-237.4				
$ZnI_2 \cdot NH_3$			c		-80.9				
$ZnI_2 \cdot 2NH_3$			c		-114.1				
$ZnI_2 \cdot 2N_2H_4$			c		-65.7				
$ZnI_2 \cdot 4NH_3$			c		-166.1				
$ZnI_2 \cdot 6NH_3$			c		-210.1				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 31-10. Zinc (at. no., 30; at. wt., 65.38)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1948

Formula	Substance	Description	State	At 0°K					At 298.16°K (25°C)		
				kcal/mole					kcal/mole		
$\text{ZnSO}_4 \cdot \frac{1}{2}\text{NH}_3$			c						-251.9		
$\text{ZnSO}_4 \cdot \text{NH}_3$			c						-268.6		
$\text{ZnSO}_4 \cdot 2\text{NH}_3$			c						-298.6		
$\text{Zn}(\text{NH}_4)_2(\text{SO}_4)_2 \cdot 2\text{H}_2\text{O}$			c						-662.8		
$\text{Zn}(\text{NH}_4)_2(\text{SO}_4)_2 \cdot 6\text{H}_2\text{O}$			c						-955.6		
$\text{ZnSO}_4 \cdot 3\text{NH}_3$			c						-325.1		
$\text{ZnSO}_4 \cdot 4\text{NH}_3$			c						-349.8		
$\text{ZnSO}_4 \cdot 5\text{NH}_3$			c						-371.7		
$\text{Zn}(\text{NH}_4)_2(\text{SeO}_4)_2$			c								
ZnSb			c						-36.		12.4
$\text{ZnCO}_3$			c						-194.2	-174.8	19.7
$\text{Zn}(\text{CH}_3)_2$	dimethylzinc		liq						6.0		
$\text{Zn}(\text{C}_2\text{H}_5)_2$	diethylzinc		liq						5.0		

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 31-11. Zinc (at. no., 30; at. wt., 65.38)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1948 (Corrected)

Formula	Substance	Description	State	At 0°K					At 298.16°K (25°C)		
				$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$\Delta H_f^\circ$ kcal/mole	$\log_{10} K_f$	$C_p^\circ$ cal/deg mole
$ZnC_2O_4 \cdot 2H_2O$	zinc oxalate dihydrate		c			-373.8					
$Zn(CHO_2)_2$	zinc formate		c			-225.0					
			aq			-229.0					
$Zn(CHO_2)_2 \cdot 2H_2O$	zinc formate dihydrate		c			-367.9					
$Zn(C_2H_3O_2)_2$	zinc acetate		c			-258.1					
	in 50 $H_2O$		aq			-264.30					
	100 $H_2O$		ao			-265.49					
	200 $H_2O$		aq			-266.55					
	400 $H_2O$		aq			-267.43					
	800 $H_2O$		aq			-267.9					
$Zn(C_2H_3O_2)_2 \cdot H_2O$	zinc acetate hydrate		c			-329.2					
$Zn(C_2H_3O_2)_2 \cdot 2H_2O$	zinc acetate dihydrate		c			-398.8					
$Zn(C_2H_3O_2)_3$	zinc glycollate		c			-350.7					
			aq			-348.7					
$Zn(C_2H_3O_2)_2 \cdot 2H_2O$	zinc glycollate dihydrate		c			-490.4					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 31-12. Zinc (at. no., 30; at. wt., 65.38)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1948

Formula	Substance	State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
$Zn(CN)_2$	zinc chloride ethylenediamine	c		18.4				
$Zn(CN)_2 \cdot ZnO$		c		-33.1				
$Zn(CN)_4^{--}$		aq		82.0				
$ZnCl_2 \cdot C_2H_4(NH_2)_2$		c		-138.0				
$ZnCl_2 \cdot 3C_2H_4(NH_2)_2$	zinc chloride trisethylenediamine	c		-179.4				
$ZnBr_2 \cdot C_2H_4(NH_2)_2$		c		-112.5				
$ZnBr_2 \cdot 3C_2H_4(NH_2)_2$	zinc bromide trisethylenediamine	c		-157.1				
$ZnI_2 \cdot C_2H_4(NH_2)_2$		c		-86.9				
$ZnI_2 \cdot 3C_2H_4(NH_2)_2$	zinc iodide trisethylenediamine	c		-133.3				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Substance		Table 31-13. Zinc (at. no., 30; at. wt., 65.38) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1948				
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K$	$C_p^\circ$
			At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole
$Zn(CHO_2)_2 \cdot 2NH_3$		c		-273.1		
$Zn(CHO_2)_2 \cdot 2\frac{1}{2}NH_3$		c		-283.9		
$Zn(CHO_2)_2 \cdot 4NH_3$		c		-316.2		
$Zn(CHO_2)_2 \cdot 5NH_3$		c		-337.8		
$Zn(CHO_2)_2 \cdot 6NH_3$		c		-354.9		
$Zn(CNS)_2$		c				
$ZnSiO_3$		c		-294.6		
$Zn_2SiO_4$		c		-360.8		
		g's		-351.8		
$2ZnI_2 \cdot PbI_2$		c		-142.9		

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 32-1. Cadmium (at. no., 48; at. wt., 112.41)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1948

Formula	Substance	Description	State	At 298.16°K (25°C)				
				$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
Cd			g	26.97	26.97	-13.700	40.067	4.968
			c, $\alpha$		0.000	0.0000	12.3	6.19
			c, $\gamma$		0.140	-0.1026		
					-2.188	1.6038		
Cd <sup>+</sup>		in Hg (saturated)	g	234.321	235.802			
Cd <sup>++</sup>			g	624.155	627.117			
		std. state, hyp. m=1	aq		-17.30		-14.6	
Cd <sup>+++</sup>			g	1431.	1436.			
Cd <sup>++++</sup>			g					
Cd <sub>2</sub>			g	51.93	51.04			
CdO			c		-60.86	39.427	13.1	10.38
					-53.79			
CdH			g	62.96	62.54	-40.849	50.76	
(CdH) <sup>+</sup>			g	238.	239.1			
Cd(OH) <sub>2</sub>			c		-112.46	82.432	22.8	



## Washington, D.C.

## National Bureau of Standards

Table 32-2. Cadmium (at. no., 48; at. wt., 112.41)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1948

Substance							
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} f_f$	$S^\circ$	$C_p^\circ$
			At 0°K	At 298.16°K (25°C)			
			kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
CdF <sub>2</sub>		c		-164.9		113.47	27.
		aq		-173.6			
CdCl		g	5.	4.6	-1.8	1.32	60.36
CdCl <sub>2</sub>		c		-93.00	-81.88	60.017	28.3
	std. state, hyp m = 1	aq		-97.39	-81.26	59.577	11.6
	in 200 H <sub>2</sub> O	aq		-96.290			
	300 H <sub>2</sub> O	aq		-96.341			
	400 H <sub>2</sub> O	aq		-96.38			
	500 H <sub>2</sub> O	aq		-96.410			
	1000 H <sub>2</sub> O	aq		-96.520			
	2000 H <sub>2</sub> O	aq		-96.650			
	3000 H <sub>2</sub> O	aq		-96.725			
	4000 H <sub>2</sub> O	aq		-96.780			
	5000 H <sub>2</sub> O	aq		-96.830			
	10000 H <sub>2</sub> O	aq		-96.965			
	50000 H <sub>2</sub> O	aq		-97.225			
	100000 H <sub>2</sub> O	aq		-97.270			
	$\infty$ H <sub>2</sub> O	aq		-97.39			

SERIES I

Washington, D.C.

Table 32-3. Cadmium (at. no., 48; at. wt., 112.41)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1948

Substance		State	Description	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K			At 298.16°K (25°C)					
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
$\text{CdCl}_2 \cdot \text{H}_2\text{O}$	c			-164.13	-140.13	102.713	40.8		
$\text{CdCl}_2 \cdot 2\frac{1}{2}\text{H}_2\text{O}$	c			-269.97	-225.47	165.266	55.6		
$\text{CdCl}_2 \cdot 4\text{H}_2\text{O}$	c								
$\text{CdCl}_2 \cdot 2\text{HCl} \cdot 7\text{H}_2\text{O}$	c			-656.1					
$\text{CdCl}_2 \cdot \text{CdO} \cdot \text{H}_2\text{O}$	c			-236.46					
$\text{CdBr}$	g			-12.	-22.	16.1	63.09		
$\text{CdBr}_2$	c			-75.15	-70.14	51.412	31.9		
	aq			-75.822					
std. state, hyp. m = 1	aq			-75.827					
in 100 $\text{H}_2\text{O}$	aq			-75.874					
200 $\text{H}_2\text{O}$	aq			-75.900					
300 $\text{H}_2\text{O}$	aq			-75.915					
400 $\text{H}_2\text{O}$	aq			-75.926					
500 $\text{H}_2\text{O}$	aq			-75.958					
1000 $\text{H}_2\text{O}$	aq			-75.955					
2000 $\text{H}_2\text{O}$	aq			-75.941					
3000 $\text{H}_2\text{O}$	aq								

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 32-4. Cadmium (at. no., 48; at. wt., 112.41) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1948									
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_r^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
CdBr <sub>2</sub>	in 4000 H <sub>2</sub> O 5000 H <sub>2</sub> O 10000 H <sub>2</sub> O 50000 H <sub>2</sub> O 100000 H <sub>2</sub> O $\infty$ H <sub>2</sub> O		aq		-75.930				
					-75.919				
					-75.890				
					-75.830				
					-75.820				
					-75.822				
CdBr <sub>2</sub> ·4H <sub>2</sub> O CdBr <sub>2</sub> ·CdO·H <sub>2</sub> O			c		-356.32	-297.64	218.166	74.7	
					-216.10				
CdI CdI <sub>2</sub>			g	21.	19.6	8.1	-5.94	64.97	
					-48.0	-48.00	35.183	40.2	
std. state, hyp m = 1 in 400 H <sub>2</sub> O 500 H <sub>2</sub> O 1000 H <sub>2</sub> O 2000 H <sub>2</sub> O			aq		-47.46				
					-47.345				
					-47.000				
					-46.680				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

Washington, D.C.

Table 32-5. Cadmium (at. no., 48; at. wt., 112.41) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1948									
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
CdI <sub>2</sub>	in	3000 H <sub>2</sub> O	aq			-46.470			
		4000 H <sub>2</sub> O	aq			-46.320			
		5000 H <sub>2</sub> O	aq			-46.190			
		10000 H <sub>2</sub> O	aq			-45.840			
		50000 H <sub>2</sub> O	aq			-45.390			
CdI <sub>2</sub> ·CdO·H <sub>2</sub> O		100000 H <sub>2</sub> O	aq			-44.990			
		∞ H <sub>2</sub> O	aq			-44.660			
			c			-178.7			
			c			-34.5	24.63	17.	
			c			-221.36	143.657	32.8	
CdS	std. state, hyp. m=1		aq			-234.20	-195.92	-10.47	
	in	50 H <sub>2</sub> O	aq			-232.190			
		100 H <sub>2</sub> O	aq			-232.413			
		200 H <sub>2</sub> O	aq			-232.578			
		300 H <sub>2</sub> O	aq			-232.670			
CdSO <sub>4</sub>		400 H <sub>2</sub> O	aq			-232.735			
		500 H <sub>2</sub> O	aq			-232.785			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 32-6. Cadmium (at. no., 48; at. wt., 112.41)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1948

Substance		State	Description	At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula									
CdSO <sub>4</sub>	in	700 H <sub>2</sub> O	aq		-232.855				
		1000 H <sub>2</sub> O	aq		-232.935				
		2000 H <sub>2</sub> O	aq		-233.085				
		3000 H <sub>2</sub> O	aq		-233.180				
		4000 H <sub>2</sub> O	aq		-233.250				
		5000 H <sub>2</sub> O	aq		-233.315				
		10000 H <sub>2</sub> O	aq		-233.500				
		20000 H <sub>2</sub> O	aq		-233.675				
		30000 H <sub>2</sub> O	aq		-233.760				
		40000 H <sub>2</sub> O	aq		-233.865				
		50000 H <sub>2</sub> O	aq		-233.905				
		100000 H <sub>2</sub> O	aq		-233.950				
		500000 H <sub>2</sub> O	aq		-234.080				
		$\infty$ H <sub>2</sub> O	aq		-234.20				
CdSO <sub>4</sub> · H <sub>2</sub> O CdSO <sub>4</sub> · $\frac{8}{9}$ H <sub>2</sub> O		c			-234.37	-254.84	166.794	41.1	
		c			-411.82	-349.63	256.274	57.9	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 32-7. Cadmium (at. no., 48; at. wt., 112.41)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1948

Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description		At 0°K kcal/mole	kcal/mole				
$\text{CdSO}_4 \cdot 2\frac{1}{2}\text{H}_2\text{SO}_4$	in 80% $\text{H}_2\text{O}$ or $\text{H}_2\text{O}$	c		-766.4				
$\text{CdSO}_4$		aq		-297.2				
		aq		-297.7				
CdSe	cadmium azide	g	2.2	1.6				
CdTe		c		-24.30	-23.82	17.460	22.6	
$\text{Cd}(\text{N}_3)_2$	std. state, hyp. m=1	c		107.8				
$\text{Cd}_3\text{N}_2$		c		38.6				
$\text{Cd}(\text{NO}_3)_2$		c		-107.98	-71.41	52.342	20.4	
		aq		-116.04				
		aq						
$\text{Cd}(\text{NO}_3)_2 \cdot 2\text{H}_2\text{O}$		c		-251.19				
$\text{Cd}(\text{NO}_3)_2 \cdot 4\text{H}_2\text{O}$		c		-394.02				
$\text{Cd}(\text{NH}_2)_2$		c		-14.71				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 32-8. Cadmium (at. no., 48; at. wt., 112.41) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1948												
Substance			State	Description	At 0°K kcal/mole	ΔH <sup>o</sup> kcal/mole	ΔF <sup>o</sup> kcal/mole	log <sub>10</sub> K <sup>o</sup>	S <sup>o</sup>	C <sub>p</sub> <sup>o</sup>		
Formula		At 298.16°K (25°C)										
		kcal/mole									kcal/mole	cal/deg mole
CdCl <sub>2</sub> ·NH <sub>3</sub>		c				-122.5						
CdCl <sub>2</sub> ·NH <sub>4</sub> Cl·½H <sub>2</sub> O		c				-206.7						
CdCl <sub>2</sub> ·2NH <sub>3</sub>		c				-150.5						
CdCl <sub>2</sub> ·4NH <sub>3</sub>		c				-194.0						
CdCl <sub>2</sub> ·4NH <sub>4</sub> Cl		c				-395.0						
CdCl <sub>2</sub> ·6NH <sub>3</sub>		c				-236.5						
CdCl <sub>2</sub> ·10NH <sub>3</sub>		c				-310.3						
CdBr <sub>2</sub> ·NH <sub>3</sub>		c				-103.5						
CdBr <sub>2</sub> ·NH <sub>4</sub> Br·½H <sub>2</sub> O		c				-177.1						
CdBr <sub>2</sub> ·2NH <sub>3</sub>		c				-131.1						
CdBr <sub>2</sub> ·6NH <sub>3</sub>		c				-218.1						
CdBr <sub>2</sub> ·12NH <sub>3</sub>		c				-329.9						
CdI <sub>2</sub> ·NH <sub>4</sub> I·½H <sub>2</sub> O		c				-133.9						
CdI <sub>2</sub> ·2NH <sub>3</sub>		c				-98.5						
CdI <sub>2</sub> ·6NH <sub>3</sub>		c				-189.4						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 32-9. Cadmium (at. no., 48; at. wt., 112.41)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1948

Substance		$\Delta H_f^\circ$		$\Delta G_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K		At 298.16°K (25°C)		cal/deg mole
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	
$\text{CdSO}_4 \cdot 4\text{NH}_3$		c		-316.8			
$\text{CdSO}_4 \cdot 6\text{NH}_3$		c		-362.6			
$\text{CdSb}$		c		-3.29			
$\text{Cd}_3\text{Sb}_2$		c		7.83	1.59	78.8	
$\text{CdCO}_3$		c		-178.7	-160.2	117.42	
		amorp		-177.5			
$\text{Cd}(\text{CH}_3)_2$		liq		16.5			
$\text{Cd}(\text{C}_2\text{H}_5)_2$		liq		14.5			
$\text{Cd}(\text{CN})_2$		c		39.0			
		aq		30.3			
$\text{Cd}(\text{ONC})_2$		c		37.7			
$2\text{Cd}(\text{CN})_2 \cdot \text{CdC} \cdot 3\text{H}_2\text{O}$	cadmium fulminate	c		-358.0			
$2\text{CdI}_2 \cdot \text{PbI}_2$		c		-128.9			



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

SERIES I

Washington, D.C.

Table 33-1. Mercury (at. no., 80; at. wt., 200.61) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1948									
Substance		State	At 0°K				At 298.16°K (25°C)		
Formula	Description		$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$c_p^\circ$ cal/deg mole	
Hg		g	15.02	14.54	7.59	-5.563	41.60	4.968	
Hg <sup>+</sup>		liq	0.000	0.000	0.000	0.0000	18.5	6.65	
Hg <sup>++</sup>		g	255.645	257.126					
Hg <sup>+++</sup>		g	688.08	689.56					
	std. state, hyp. m = 1	aq			39.38	-28.865			
Hg <sub>2</sub>		g	28.6	27.1					
Hg <sub>2</sub> <sup>++</sup>		aq			36.79	-26.967			
HgO	red	c, II		-21.68	-13.990	10.2545	17.2	10.93	
	yellow	c, I		-21.56	-13.959	10.2317	17.5		
Hg <sub>2</sub> O		c		-21.8					
HgH		g	58.07	58.06	52.60	-38.555	52.42	7.147	
HgO <sub>2</sub> <sup>-</sup>	std. state, hyp. m = 1	aq			-45.42	33.292			
Hg(OH) <sub>2</sub>	std. state, hyp. m = 1	aq			-66.0	48.38			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 33-2. Mercury (at. no., 80; at. wt., 200.61)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1946 (Corrected)

Substance								
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
			At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)		
			kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
HgF		g	15.	14.				
HgF <sub>2</sub>		c						
Hg <sub>2</sub> F <sub>2</sub>		c						
HgCl		g	19.	19.	14.	-10.3	62.2	8.65
HgCl <sub>2</sub>		c		-55.0				18.3
		aq		-51.8				
HgCl <sub>3</sub> <sup>-</sup>		aq		-93.7				
Hg <sub>2</sub> Cl <sub>2</sub>		c		-63.32	-50.350	36.9058	46.8	24.3
HgCl <sub>2</sub> ·HgO		c		-80.9				
HgCl <sub>2</sub> ·2HgO		c		-105.3				
HgCl <sub>2</sub> ·3HgO		c		-128.4				
HgCl <sub>2</sub> ·4HgO		c		-151.3				
HgBr		g		23.	18.	-13.2	65.0	8.82
HgBr <sub>2</sub>		c		-40.5				
		aq		-37.1				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

<p>Table 33-3. Mercury (at. no., 80; at. wt., 200.61) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1948</p>									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	kcal/mole				
HgBr <sub>4</sub> <sup>---</sup>			aq		-99.9				
Hg <sub>2</sub> Br <sub>2</sub>			c		-49.42		31.3088	50.9	
HgBr <sub>2</sub> ·HgO			c		-59.6				
HgBr <sub>2</sub> ·2HgO			c		-78.5				
HgBr <sub>2</sub> ·3HgO			c		-98.3				
HgBr <sub>2</sub> ·4HgO			c		-117.6				
HgI			g	34.	33.	23.	-16.9	67.1	8.89
HgI <sub>2</sub>	red		c, II		-25.2				
	yellow		c, I		-24.55				
HgI <sub>3</sub> <sup>-</sup>			aq		-55.4				
HgI <sub>4</sub> <sup>---</sup>			aq		-28.91		19.497	57.2	25.3
Hg <sub>2</sub> I <sub>2</sub>	yellow		c			-26.60			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

Washington, D.C.

National Bureau of Standards

Table 33-4. Mercury (at. no., 80; at. wt., 200.61) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1948									
Formula	Substance	Description	State	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
$\text{HgS}$  $\text{HgSO}_4$ $\text{Hg}_2\text{SO}_4$ $\text{HgSO}_4 \cdot 2\text{HgO}$ $\text{HgSe}$		red, cinnabar  black, metacinnabarite	g c, II c, I c c c			3. -13.90 -12.90 -168.3 -177.34 -222.3	8.5539 8.0995	18.6 19.9	31.55
$\text{Hg}_2(\text{N}_3)_2$ $\text{Hg}(\text{NO}_3)_2$ $\text{Hg}(\text{NO}_3)_2 \cdot \frac{1}{2}\text{H}_2\text{O}$ $\text{Hg}_2(\text{NO}_3)_2$ $\text{Hg}_2(\text{NO}_3)_2 \cdot 2\text{H}_2\text{O}$ $\text{Hg}(\text{NO}_3)_2 \cdot 2\text{HgO} \cdot \text{H}_2\text{O}$		mercurous azide	c aq c aq c c			133. -58.0 -93.0 -58.4 -206.9 -178.2	109.303	47.98	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 33-5. Mercury (at. no., 80; at. wt., 200.61) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1948									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula									
						At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole
(Hg <sub>2</sub> N) <sub>2</sub> O	Millon's oxide	c			76.3				
Hg <sub>2</sub> NOH	Millon's hydroxide	c			2.5				
Hg <sub>2</sub> NOH·1½H <sub>2</sub> O		c			-103.9				
Hg <sub>2</sub> NOH·2H <sub>2</sub> O	Millon's base	c			-139.2				
Hg <sub>2</sub> NCI	Millon's chloride	c							
Hg <sub>2</sub> NCI·½H <sub>2</sub> O		c			-28.0				
Hg <sub>2</sub> NCI·H <sub>2</sub> O		c			-63.5				
Hg <sub>2</sub> NCI·½HgCl <sub>2</sub>		c			-24.6				
HgCl <sub>2</sub> ·2NH <sub>3</sub>		c			-112.0				
HgCl <sub>2</sub> ·8NH <sub>3</sub>		c			-226.8				
HgCl <sub>2</sub> ·9½NH <sub>3</sub>		c			-254.2				
Hg <sub>2</sub> NCI·½NH <sub>3</sub>		c			-5.5				
Hg <sub>2</sub> NCI·NH <sub>3</sub>		c			-13.6				
Hg <sub>2</sub> NCI·NH <sub>4</sub> Cl		c			-73.8				
Hg <sub>2</sub> NCI·3NH <sub>4</sub> Cl		c			-222.9				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 33-6. Mercury (at. no., 80; at. wt., 200.61)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1948

Formula	Substance	State	$\Delta H^{\circ}$		$\Delta F^{\circ}$	$\log_{10} K_f$	$S^{\circ}$	$C_p^{\circ}$
			At 0°K kcal/mole	kcal/mole		At 298.16°K (25°C)		
Hg <sub>2</sub> NBr	Millon's bromide	c		15.0				
HgBr <sub>2</sub> ·2NH <sub>3</sub>		c		-94.7				
HgBr <sub>2</sub> ·8NH <sub>3</sub>		c		-207.7				
Hg <sub>2</sub> NBr·NH <sub>4</sub> Br		c		-55.9				
Hg <sub>2</sub> NBr·3NH <sub>4</sub> Br		c		-180.5				
Hg <sub>2</sub> NBr·4HgBr <sub>2</sub>		c		1.0				
Hg <sub>2</sub> NBr·2HgBr <sub>2</sub>	dimethylmercury diethylmercury mercuric oxalate	c		-11.8				
HgI <sub>2</sub> ·18NH <sub>3</sub>		c		-56.6				
HgI <sub>2</sub> ·2NH <sub>3</sub>		c		-71.6				
HgI <sub>2</sub> ·6NH <sub>3</sub>		c		-146.0				
Hg(CH <sub>3</sub> ) <sub>2</sub>		liq		18.				
Hg(C <sub>2</sub> H <sub>5</sub> ) <sub>2</sub>		liq		15.				
HgC <sub>2</sub> O <sub>4</sub>		c		-161.8				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 33-7. Mercury (at. no., 80; at. wt., 200.61)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1948

Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description		At 0°K kcal/mole	kcal/mole				
$\text{Hg}(\text{C}_2\text{H}_3\text{O}_2)_2$	mercuric acetate	c		-199.4				
$\text{Hg}_2(\text{C}_2\text{H}_3\text{O}_2)_2$	mercurous acetate	aq		-195.2				
$\text{HgCl}_2 \cdot \text{CH}_3\text{OH}$		c		-201.1				
		c		-113.3				
$\text{Hg}(\text{CN})_2$		c		62.5				
$\text{Hg}(\text{CN})_4$		aq		66.1				
$\text{Hg}(\text{ONC})_2$		aq		126.0				
$\text{Hg}(\text{CN})_2 \cdot \text{H}_2\text{O}$	mercuric fulminate	c		64.				
$3\text{Hg}(\text{CN})_2 \cdot \text{H}_2\text{O}$		c		41.2				
		c		173.4				
$\text{HgCl}_2 \cdot \text{C}_2\text{H}_4(\text{NH}_2)_2$		c		-88.5				
$\text{HgBr}_2 \cdot \text{C}_2\text{H}_4(\text{NH}_2)_2$		c		-65.7				
$\text{HgI}_2 \cdot \text{C}_2\text{H}_4(\text{NH}_2)_2$		c		-47.8				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 33-8. Mercury (at. no., 80; at. wt., 200.61)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K. (25°C)

December 31, 1948

Formula	Substance	Description	State	At 298.16°K. (25°C)				
				$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
$\text{Hg}(\text{CN})_2 \cdot \text{NH}_4\text{Cl} \cdot 2\text{H}_2\text{O}$			c		-64.5			
$\text{Hg}(\text{CN})_2 \cdot \text{NH}_4\text{Br} \cdot 2\text{H}_2\text{O}$			c		-77.8			
$\text{Hg}(\text{CN})_2 \cdot \text{NH}_4\text{I} \cdot 2\text{H}_2\text{O}$			c		-10.0			
$\text{Hg}(\text{CNS})_2$			c		48.0			
$\text{Hg}_5\text{I}_2$			c		-2.5			
$\text{Hg}(\text{CN})_2 \cdot \text{ZnCl}_2 \cdot 7\text{H}_2\text{O}$			c		-475.2			
$2\text{Hg}(\text{CN})_2 \cdot \text{ZnBr}_2 \cdot 8\text{H}_2\text{O}$			c		-530.7			
$\text{Hg}(\text{CN})_2 \cdot \text{CdCl}_2 \cdot 2\text{H}_2\text{O}$			c		-176.6			
$\text{Hg}(\text{CN})_2 \cdot \text{CdBr}_2 \cdot 3\text{H}_2\text{O}$			c		-228.0			
$2\text{Hg}(\text{CN})_2 \cdot \text{CdI}_2 \cdot 8\text{H}_2\text{O}$			c		-486.1			



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 34-1. Copper (at. no., 29; at. wt., 63.54)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 298.16°K (25°C)					
			At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole
Cu	in Hg (two phase)	g	81.24	81.52	72.04	-52.804	39.744	4.968
		c	0.000	0.000	0.000	0.0000	7.96	5.848
Cu <sup>+</sup>		g	259.627	250.828				
Cu <sup>++</sup>		aq		12.4	12.0	-8.796	-6.3	
Cu <sup>+++</sup>	std. state, hyp. m=1	g	727.436	730.118				
Cu <sup>++++</sup>	std. state, hyp. m=1	aq		15.39	15.53	-11.383	-23.6	
		g						
		g						
CuO		g	36.	35.				
		c		-37.1	-30.4	22.28	10.4	10.6
CuO <sub>2</sub> <sup>---</sup>		aq			-43.3	31.74		
Cu <sub>2</sub> O	std. state, hyp. m=1	c		-39.84	-34.98	25.640	24.1	16.7

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 34-2. Copper [at. no., 29; at. wt., 63.54]

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1949; March 31, 1950

Substance											
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$			
			At 298.16°K (25°C)								
			At 0°K								
			kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole			
CuH	std. state, hyp. m = 1	g	71.	71.	64.	-46.9	46.89	6.98			
HCuO <sub>2</sub> <sup>-</sup>		aq			-61.2	44.86					
Cu(OH) <sub>2</sub>		c		-107.2							
CuF		g	44.	44.							
CuF <sub>2</sub>		c		-126.9							
CuF <sub>2</sub> · 2H <sub>2</sub> O		c		-274.5	-235.2	172.40	36.2				
CuCl	in aq HCl	g	32.	32.	25.	-18.3	56.50	8.40			
		c		-32.2	-26.4	20.82	21.9				
		aq		-25.8							
		c		-49.2							
		aq		-55.43							
CuCl <sub>2</sub>	in 10 H <sub>2</sub> O	aq		-57.48							
		aq		-58.36							
		aq		-59.18							
		aq		-59.98							
		aq		-60.52							
		aq		-60.94							
		aq		-61.24							
		aq		-61.4							

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 34-3. Copper (at. no., 29; at. wt., 63.54)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1949; March 31, 1950

Formula	Substance	Description	State	At 298.16°K (25°C)				
				$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole
CuCl <sub>2</sub>	in 10 C <sub>2</sub> H <sub>6</sub> O (ethanol)	20 C <sub>2</sub> H <sub>6</sub> O			-53.19			
		50 C <sub>2</sub> H <sub>6</sub> O			-53.55			
		100 C <sub>2</sub> H <sub>6</sub> O			-53.89			
		200 C <sub>2</sub> H <sub>6</sub> O			-54.06			
		in HCl·8H <sub>2</sub> O	aq		-54.20			
CuCl <sub>2</sub> ·2H <sub>2</sub> O	std. state, hyp. m = 1		c		-55.5			
			aq		-193.0	-58.1	42.53	51.1
Cu(ClO <sub>3</sub> ) <sub>2</sub>			aq		-31.6			
			aq		-19.0			
CuCl <sub>2</sub> ·CuO			c		-85.3			
			c		-154.9			
CuCl <sub>2</sub> ·3CuO			c		-450.0			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 34-4. Copper (at. no., 29; at. wt., 63.54)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1949; March 31, 1950

Formula	Substance	Description	State	At 298.16°K (25°C)				
				$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
CuBr			g		38.	28.	59.22	8.61
CuBr <sub>2</sub>			c		-25.1	-25.81	21.9	
CuBr <sub>2</sub> ·4H <sub>2</sub> O		in 400 H <sub>2</sub> O	c		-33.2			
			aq		-42.0			
CuBr <sub>2</sub> ·3Cu(OH) <sub>2</sub>		std. state, hyp. m=1	c		-316.4			
CuBr <sub>2</sub> <sup>-</sup>			aq		-373.6	-45.2	31.31	
CuI			c					
			d	63.	62.	50.	51.05	8.70
			c		-16.2	-16.62	23.1	12.92
CuI <sub>2</sub>			c, II		-1.7			20.1
			aq		-7.9			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 34-5. Copper (at. no., 29; at. wt., 63.54)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1943

Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K			At 298.16°K (25°C)					
				kcal/mole	kcal/mole	kcal/mole	cal/deg mole		
CuS			c	-11.6	-11.7	0.176	15.9	11.43	
Cu <sub>2</sub> S			c, II	-19.0	-20.6	15.10	28.9	18.24	
CuSO <sub>4</sub>			c	-184.00	-158.2	115.95	27.1	24.1	
	std. state, hyp m = 1		aq	-201.51	-161.81	118.04	-19.5		
	in 50 H <sub>2</sub> O		aq	-199.998					
	100 H <sub>2</sub> O		aq	-200.110					
	200 H <sub>2</sub> O		aq	-200.201					
	500 H <sub>2</sub> O		aq	-200.342					
	1000 H <sub>2</sub> O		aq	-200.491					
	2000 H <sub>2</sub> O		a	-200.641					
	3000 H <sub>2</sub> O		aq	-200.723					
	5000 H <sub>2</sub> O		aq	-200.825					
	10,000 H <sub>2</sub> O		aq	-200.964					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 34-6. Copper (at. no., 29; at. wt., 63.54)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1949

Formula	Substance	Description	State	At 298.16°K (25°C)				
				$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
CuSO <sub>4</sub>	in 20,000 H <sub>2</sub> O		aq		-201.110			
	30,000 H <sub>2</sub> O		aq		-201.192			
	50,000 H <sub>2</sub> O		aq		-201.281			
	100,000 H <sub>2</sub> O		aq		-201.364			
	200,000 H <sub>2</sub> O		aq		-201.427			
	500,000 H <sub>2</sub> O		aq		-201.469			
	$\infty$ H <sub>2</sub> O		aq		-201.51			
CuSO <sub>4</sub> ·H <sub>2</sub> O			c		-259.00	-219.2	35.8	31.3
CuSO <sub>4</sub> ·3H <sub>2</sub> O			c		-402.27	-334.6	53.8	49.0
CuSO <sub>4</sub> ·5H <sub>2</sub> O			c, II		-544.45	-449.3	73.0	67.2
Cu <sub>2</sub> O								
Cu <sub>2</sub> O <sub>6</sub>			aq		-265.0			
Cu <sub>2</sub> O <sub>6</sub> ·5H <sub>2</sub> O			c		-611.5			
Cu <sub>2</sub> SO <sub>4</sub>			c		-179.2			
			aq		-130.8			
CuSO <sub>4</sub> ·3CuO·4H <sub>2</sub> O			c		-592.6			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 34-7. Copper (at. no., 29; at. wt., 63.54)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1949; March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole		cal/deg mole	cal/deg mole
CuSe			c		-15.1			21.2
Cu <sub>2</sub> Se			c, II		-5.5			
CuSeO <sub>4</sub>			aq		-129.3			
CuSeO <sub>4</sub> ·5H <sub>2</sub> O			c		-473.8			
CuN <sub>3</sub>			c		60.5			
Cu <sub>3</sub> N		cuprous azide	c		17.8			21.7
Cu(NO <sub>3</sub> ) <sub>2</sub>			c		-73.4			
	in 10 H <sub>2</sub> O		aq		-83.11			
	12 H <sub>2</sub> O		aq		-83.57			
	15 H <sub>2</sub> O		aq		-83.65			
	20 H <sub>2</sub> O		aq		-84.05			
	50 H <sub>2</sub> O		aq		-84.01			
	100 H <sub>2</sub> O		aq		-83.68			
	200 H <sub>2</sub> O		aq		-83.64			
	400 H <sub>2</sub> O		aq		-83.60			
	800 H <sub>2</sub> O		aq		-83.6			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 34-8. Copper (at. no., 29; at. wt., 63.54)  
HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1949

Substance		Description	State	$\Delta H_f^\circ$		$\Delta F^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K			At 298.16°K (25°C)					
				kcal/mole	kcal/mole				
$\text{Cu}(\text{NO}_3)_2 \cdot 3\text{H}_2\text{O}$			c		-291.3				
$\text{Cu}(\text{NO}_3)_2 \cdot 6\text{H}_2\text{O}$			c		-504.3				
$\text{Cu}(\text{NO}_3)_2 \cdot 3\text{CuO} \cdot 3\text{H}_2\text{O}$			c		-415.7				
$4\text{CuO} \cdot 3\text{N}_2\text{O}_5$			c		-236.				
$\text{Cu}(\text{NH}_3)^+$		std. state, hyp. m=1	aq			-26.5	19.42		
$\text{Cu}(\text{NH}_3)^{++}$		std. state, hyp. m=1	aq			-29.8	21.84		
$\text{Cu}(\text{NH}_3)_2^+$		std. state, hyp. m=1	aq			-39.2	28.73		
$\text{Cu}(\text{NH}_3)_2^{++}$		std. state, hyp. m=1	aq			-41.2	30.20	133.3	
$\text{Cu}(\text{NH}_3)_3^{++}$		std. state, hyp. m=1	aq			-51.7	37.90		
$\text{Cu}(\text{NH}_3)_4^{++}$		std. state, hyp. m=1	aq			-61.2	44.86	192.8	
$\text{Cu}(\text{NH}_3)_5^{++}$		std. state, hyp. m=1	aq			-67.0	49.11		



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 34-9. Copper (at. no., 29; at. wt., 63.54)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949; March 31, 1950

Substance		State	Description	$\Delta H^\circ$		$\log_{10} f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K			At 298.16°K (25°C)				
				kcal/mole	kcal/mole	kcal/mole	cal/deg mole	
$\text{Cu}(\text{NO}_3)_2 \cdot 2\text{NH}_3$	c	-139.7						
$\text{Cu}(\text{NO}_3)_2 \cdot 4\text{NH}_3$	c	-196.6						
$\text{Cu}(\text{NO}_3)_2 \cdot 6\text{NH}_3$	c	-238.3						
$\text{CuCl} \cdot \text{NH}_3$	c	-60.3						
$\text{CuCl} \cdot 1\frac{1}{2}\text{NH}_3$	c	-72.1						
$\text{CuCl}_2 \cdot 2\text{NH}_3$	c	-114.1						
$\text{CuCl}_2 \cdot 2\text{NH}_3 \cdot \frac{1}{2}\text{H}_2\text{O}$	c	-130.5						
$\text{CuCl}_2 \cdot 2\text{NH}_4\text{Cl}$	c	-199.7						
$\text{CuCl}_2 \cdot 2\text{NH}_4\text{Cl} \cdot 2\text{H}_2\text{O}$	c	-334.4						
$\text{CuCl}_2 \cdot 3\text{NH}_3$	c	-102.9						
$\text{CuCl}_2 \cdot 3\frac{1}{2}\text{NH}_3$	c	-147.6						
$\text{CuCl}_2 \cdot 4\text{NH}_3 \cdot 2\text{H}_2\text{O}$	c	-307.0						
$\text{CuCl}_2 \cdot 5\text{NH}_3$	c	-186.4						
$\text{CuCl}_2 \cdot 5\text{NH}_3 \cdot \frac{1}{2}\text{H}_2\text{O}$	c	-200.9						
$\text{CuCl}_2 \cdot 5\text{NH}_3 \cdot 1\frac{1}{2}\text{H}_2\text{O}$	c	-269.4						
$\text{CuCl}_2 \cdot 6\text{NH}_3$	c	-212.4						
$\text{CuCl}_2 \cdot 10\text{NH}_3$	c	-266.9						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 34-10. Copper (at. no., 29; at. wt., 63.54)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1949; March 31, 1950

Substance		State	Description	Formula	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} f_f$	$S^\circ$	$C_p^\circ$			
					At 0°K	At 298.16°K (25°C)						
					kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole			
	CuBr·NH <sub>3</sub>	c				-50.8						
	CuBr·1½NH <sub>3</sub>	c				-62.9						
	CuBr <sub>2</sub> ·2NH <sub>3</sub>	c				-96.5						
	CuBr·3NH <sub>3</sub>	c				-93.6						
	CuBr <sub>2</sub> ·3½NH <sub>3</sub>	c				-132.0						
	CuBr <sub>2</sub> ·5NH <sub>3</sub>	c				-171.9						
	CuBr <sub>2</sub> ·10NH <sub>3</sub>	c				-266.1						
	CuI·½NH <sub>3</sub>	c				-29.5						
	CuI·NH <sub>3</sub>	c				-42.4						
	CuI·2NH <sub>3</sub>	c				-64.7						
	CuI <sub>2</sub> ·2NH <sub>3</sub>	c				-61.8						
	CuI·3NH <sub>3</sub>	c				-86.1						
	CuI <sub>2</sub> ·3½NH <sub>3</sub>	c				-96.8						
	CuI <sub>2</sub> ·5NH <sub>3</sub>	c				-137.2						
	CuI <sub>2</sub> ·10NH <sub>3</sub>	c				-227.4						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 34-11. Copper (at. no., 29; at. wt., 63.54) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1949									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole			At 298.16°K (25°C) kcal/mole	cal/deg mole
$\text{CuSO}_4 \cdot \text{NH}_3$ $\text{CuSO}_4 \cdot 2\text{NH}_3$ $\text{CuSO}_4 \cdot (\text{NH}_4)_2\text{SO}_4$ $\text{CuSO}_4 \cdot (\text{NH}_4)_2\text{SO}_4 \cdot 2\text{H}_2\text{O}$ $\text{CuSO}_4 \cdot (\text{NH}_4)_2\text{SO}_4 \cdot 6\text{H}_2\text{O}$			c		-218.4				100.
					-248.8				
					-470.6				
					-606.3				
					-900.9				
$\text{CuSO}_4 \cdot 4\text{NH}_3$ $\text{CuSO}_4 \cdot 4\text{NH}_3 \cdot 1\frac{1}{2}\text{H}_2\text{O}$ $\text{CuSO}_4 \cdot 5\text{NH}_3$ $\text{CuSO}_4 \cdot 6\text{NH}_3$			c		-300.2				
					-406.9				
					-325.6				
					-352.7				
$\text{CuP}_2$ $\text{Cu}_3\text{P}$			c		-28.9				
					-36.4				
$\text{CuCl} \cdot \text{PH}_3$ $\text{CuCl} \cdot 2\text{PH}_3$			c		-42.7				
					-49.1				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 34-12. Copper (at. no., 29; at. wt., 63.54) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1949; March 31, 1950									
Substance		Description	State	At 298.16°K (25°C)					cal/deg mole
Formula	ΔH <sup>o</sup>			log <sub>10</sub> K <sup>o</sup>	S <sup>o</sup>	C <sub>p</sub> <sup>o</sup>			
	kcal/mole						kcal/mole		
CuBr·PH <sub>3</sub>			c		-34.6				
CuBr·2PH <sub>3</sub>			c		-42.3				
CuI·PH <sub>3</sub>			c		-24.5				
CuI·2PH <sub>3</sub>			c		-31.8				
Cu <sub>2</sub> Sb			c		-2				18.3
Cu <sub>3</sub> Sb			c						24.5
CuCO <sub>3</sub>			c		-142.2				
Cu(CHO <sub>2</sub> ) <sub>2</sub>	cupric formate		c		-178.6				
	in 200 H <sub>2</sub> O		aq		-179.7				
Cu(CHO <sub>2</sub> ) <sub>2</sub> ·4H <sub>2</sub> O			c		-460.8				
Cu(C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>2</sub>	cupric acetate		c		-214.3				
	in 400 H <sub>2</sub> O		aq		-217.0				
Cu(C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>2</sub> ·2H <sub>2</sub> O			c		-285.3				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 34-13. Copper (at. no., 29; at. wt., 63.54)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1949; March 31, 1950

Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description		At 0°K kcal/mole	kcal/mole				
$\text{Cu}(\text{C}_2\text{H}_3\text{O}_2)_2$	cupric glycolate	c		-297.4				
		aq		-295.6				
$2\text{CuCl} \cdot \text{CO}$		aq		-89.4				
$2\text{CuCl} \cdot \text{CO} \cdot 2\text{H}_2\text{O}$		c		-229.4				
$2\text{CuCl} \cdot \text{C}_2\text{H}_2$	cupric ethylsulfate	c		2.3				
$\text{CuCl}_2 \cdot 2\text{CH}_3\text{OH}$		c		-166.4				
$\text{Cu}(\text{C}_2\text{H}_5\text{SO}_4)_2$		aq		-402.0				
$\text{CuCN}$	cuprous fulminate	aq		23.1				
$\text{CuONC}$		c		26.3				
$\text{Cu}(\text{CHO}_2)_2 \cdot 2\text{NH}_3$	cupric formate diammine	c		-226.8				
$\text{Cu}(\text{CHO}_2)_2 \cdot 4\text{NH}_3$		c		-275.7				
$\text{Cu}(\text{CHO}_2)_2 \cdot 6\text{NH}_3$		c		-314.8				
$\text{Cu}(\text{CHO}_2)_2 \cdot 8\text{NH}_3$		c		-352.5				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 34-14. Copper (at. no., 29; at. wt., 63.54)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1949; March 31, 1950

Formula	Substance	State	Description	At 0°K				At 298.16°K (25°C)			
				$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\Delta P_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
$\text{Cu}(\text{C}_2\text{H}_3\text{O}_2)_2 \cdot 2\text{NH}_3$	cupric acetate diammine	c			-269.5						
$\text{Cu}(\text{C}_2\text{H}_3\text{O}_2)_2 \cdot 4\text{NH}_3$		c			-314.3						
$\text{Cu}(\text{C}_2\text{H}_3\text{O}_2)_2 \cdot 5\text{NH}_3$		c			-335.3						
$\text{Cu}(\text{C}_2\text{H}_3\text{O}_2)_2 \cdot 8\text{NH}_3$		c			-394.8						
$\text{Cu}(\text{C}_2\text{H}_3\text{O}_3)_2 \cdot 2\text{NH}_3$	cupric glycolate diammine	c			-351.5						
$\text{Cu}(\text{C}_2\text{H}_3\text{O}_3)_2 \cdot 4\text{NH}_3$		c			-397.2						
$\text{Cu}(\text{C}_2\text{H}_3\text{O}_3)_2 \cdot 6\text{NH}_3$		c			-440.0						
$\text{Cu}(\text{C}_2\text{H}_3\text{O}_3)_2 \cdot 8\text{NH}_3$		c			-480.3						
$\text{CuCl}_2 \cdot \text{C}_2\text{H}_4(\text{NH}_2)_2$	cupric chloride ethylenediamine	c			-69.3						
$\text{CuCl}_2 \cdot 2\text{C}_2\text{H}_4(\text{NH}_2)_2$	cupric chloride bisethylenediamine	c			-114.7						
$\text{CuBr}_2 \cdot \text{C}_2\text{H}_4(\text{NH}_2)_2$	cupric bromide ethylenediamine	c			-71.8						
$\text{CuBr}_2 \cdot 2\text{C}_2\text{H}_4(\text{NH}_2)_2$	cupric bromide bisethylenediamine	c			-102.4						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 34-15. Copper (at. no., 29; at. wt., 63.54)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		State	At 298.16°K (25°C)				
Formula	Description		$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$
$\text{Cu}_3\text{Si}$		c					23.0
$\text{CuSiO}_3 \cdot \text{H}_2\text{O}$		c					29.
$\text{Cu}_3\text{Sn}$		c	-8.				
$\text{Cu}_2\text{Zn}_3$		c	-16.				
$\text{Cu}_2\text{Cd}_3$		c	-4.7				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 35-1. Silver (at. no., 47; at. wt., 107.860)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 298.16°K (25°C)				
			At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
Ag		g	69.02	69.12	59.84	-43.862	4.9680
Ag <sup>+</sup>		c	0.000	0.000	0.000	10.206	6.092
Ag <sup>++</sup>		g	243.693	245.274			
Ag <sup>+++</sup>	std. state, hyp. m=1	aq		25.31	18.430	-13.5089	9.
Ag <sub>2</sub> O		g	738.99	742.05			
Ag <sub>2</sub> O <sub>2</sub>		g	1571.5	1576.0			
AgH		c	-6.940	-7.306	-2.586	1.8955	15.67
		c		-6.3			
		g	68.	67.7	60.8	-44.56	6.99



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 35-2. Silver (at. no., 47; at. wt., 107.880)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance							
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$C_p^\circ$	
AgF	std. state, hyp. m = 1 in 3 H <sub>2</sub> O 4 H <sub>2</sub> O 5 H <sub>2</sub> O 7 H <sub>2</sub> O 10 H <sub>2</sub> O 15 H <sub>2</sub> O 20 H <sub>2</sub> O 30 H <sub>2</sub> O 40 H <sub>2</sub> O 50 H <sub>2</sub> O 100 H <sub>2</sub> O 400 H <sub>2</sub> O $\infty$ H <sub>2</sub> O	c	At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole
		aq		-48.5	-44.2	32.40	20.
		aq		-53.35	-47.65	34.927	15.4
		aq		-52.50			
		aq		-52.70			
		aq		-52.83			
		aq		-52.955			
		aq		-53.015			
		aq		-53.053			
		aq		-53.072			
		aq		-53.080			
		aq		-53.085			
		aq		-53.088			
		aq		-53.091			
		aq		-53.1			
aq		-53.35					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 35-3. Silver (at. no., 47; at. wt., 107.860)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		State	Description	At 0°K					At 298.16°K (25°C)		
Formula				$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$G_p^\circ$ cal/deg mole		
AgF·H <sub>2</sub> O		c			-120.4						
AgF·2H <sub>2</sub> O		c			-191.2						
AgF·4H <sub>2</sub> O		c			-331.5						
Ag <sub>2</sub> F		c			-50.4						
AgF <sub>2</sub>		c			-88.5						
AgHF <sub>2</sub>		aq			-130.8						
AgCl		g			23.23	16.79	-12.307	58.5	8.54		
Ag <sub>2</sub> Cl		c			-30.362	-26.224	19.2218	22.97	12.14		
AgClO <sub>2</sub>		c			-30.9						
AgClO <sub>3</sub>		c			0.0	16.0	-11.73	32.16	20.87		
AgClO <sub>4</sub>		c			-5.73						
AgClO <sub>4</sub>		aq			1.81						
AgClO <sub>4</sub>		c			-7.75						
AgClO <sub>4</sub>		aq			-6.10						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 35-4. Silver (at. no., 47; at. wt., 107.880)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949 (Corrected)

Formula	Substance	Description	State	At 298.16°K (25°C)				
				$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
AgBr			c, II					
AgBrO <sub>3</sub>			c					
AgI			c, II					
AgIO <sub>3</sub>			c					
Ag <sub>2</sub> H <sub>3</sub> IO <sub>6</sub>			c, I					
3AgI·HI·7H <sub>2</sub> O			c					
Ag <sub>2</sub> S		rhombic, α	c, II					
		β	c, I					
Ag <sub>2</sub> SO <sub>4</sub>			c, II					
Ag <sub>2</sub> S <sub>2</sub> O <sub>6</sub>			aq					
Ag <sub>2</sub> S <sub>2</sub> O <sub>6</sub> ·2H <sub>2</sub> O			aq					
Ag(S <sub>2</sub> O <sub>3</sub> ) <sub>2</sub> <sup>2-</sup>			c					
Ag <sub>2</sub> SO <sub>4</sub> ·2HCl			aq					
			c					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 35-5. Silver (at. no., 47; at. wt., 107.880)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)		cal/deg mole	cal/deg mole
			kcal/mole	kcal/mole	kcal/mole		
Ag <sub>2</sub> Se		c, II					
Ag <sub>2</sub> SeO <sub>4</sub>		c		-66.5	50.21	48.3	
Ag <sub>2</sub> Te		c, II					
AgN <sub>3</sub>	silver azide	c		66.8			
AgNO <sub>2</sub>		c		-10.605			18.8
	std. state, hyp. m=1	aq		-0.09	-3.4773	30.62	
AgNO <sub>3</sub>		c, II		-29.43	-7.176	49.0	
	std. state, hyp. m=1	aq		-24.06	5.637	33.68	22.24
Ag <sub>2</sub> N <sub>2</sub> O <sub>2</sub>		c		14.	5.849	52.67	
Ag(NH <sub>3</sub> ) <sub>2</sub> <sup>+</sup>		aq		-26.724			
AgNO <sub>3</sub> ·2NH <sub>3</sub>		c		-84.5			
		aq		-76.3			
AgNO <sub>3</sub> ·3NH <sub>3</sub>		c		-105.0			
		aq		-95.6			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 35-6. Silver (at. no., 47; at. wt., 107.880)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		State	Description	At 298.16°K (25°C)					
Formula				$\Delta F^{o}$	$\Delta F^{o}$	$\log_{10} f$	$S^o$		
	$\Delta H^{o}$ kcal/mole			kcal/mole	kcal/mole	cal/deg mole	cal/deg mole		
AgCl · NH <sub>3</sub>		c			-52.5				
AgCl · 1½NH <sub>3</sub>		c			-64.3				
AgCl · 3NH <sub>3</sub>		c			-95.1				
AgClO <sub>4</sub> · 2NH <sub>3</sub>		c			-67.5				
		aq			-57.4				
AgClO <sub>4</sub> · 3NH <sub>3</sub>		c			-89.2				
		aq			-78.7				
AgBr · NH <sub>3</sub>		c			-45.4				
AgBr · 1½NH <sub>3</sub>		c			-55.8				
AgBr · 3NH <sub>3</sub>		c			-85.6				
AgI · ½NH <sub>3</sub>		c			-26.2				
AgI · NH <sub>3</sub>		c			-36.1				
AgI · 1½NH <sub>3</sub>		c			-45.5				
AgI · 2NH <sub>3</sub>		c			-54.9				
AgI · 3NH <sub>3</sub>		c			-73.2				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 35-7. Silver (at. no., 47; at. wt., 107.880)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance									
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 298.16°K (25°C)						
			At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole		
AgP <sub>2</sub>	silver acetylide	c		-9.8					
AgP <sub>3</sub>		c		-15.0					
Ag <sub>2</sub> C <sub>2</sub>		c		81.9					
Ag <sub>2</sub> CO <sub>3</sub>		c		-120.97	-104.48	76.5824	40.0	26.8	
Ag <sub>2</sub> C <sub>2</sub> O <sub>4</sub>	silver oxalate	c		-159.1					
Ag <sub>2</sub> C <sub>2</sub> H <sub>3</sub> O <sub>2</sub>	silver acetate	c		-93.41					
		aq		-89.0					
Ag <sub>2</sub> C <sub>2</sub> ·AgCl	(Ag <sub>2</sub> C <sub>2</sub> ) <sub>2</sub> ·AgCl	c		50.4					
(Ag <sub>2</sub> C <sub>2</sub> ) <sub>2</sub> ·AgCl		c		130.8					
Ag <sub>2</sub> C <sub>2</sub> ·AgI		c		67.9					
Ag <sub>2</sub> C <sub>2</sub> ·2AgI		c		51.8					
Ag <sub>2</sub> C <sub>2</sub> ·Ag <sub>2</sub> SO <sub>4</sub>		c		-90.1					
(Ag <sub>2</sub> C <sub>2</sub> ) <sub>2</sub> ·Ag <sub>2</sub> SO <sub>4</sub>		c		-15.4					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 35-8. Silver. (at. no., 47; at. wt., 107.880)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K			At 298.16°K (25°C)					cal/deg mole
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	
AgCN			c		34.94	39.20	-28.733	20.0	
$\text{Ag}(\text{CN})_2^-$		std. state, hyp. m = 1	aq		64.5	72.05	-52.812	49.0	
$\text{Ag}_2\text{CN}_2$		silver cyanamide	c		51.5				
AgCNO		silver cyanate	c		-21.1				
AgONC		silver fulminate	c		43.2				
$\text{Ag}_2\text{C}_2\cdot\text{AgNO}_3$			c		51.				
$\text{AgCl}\cdot\text{CH}_3\text{NH}_2$			c		-49.6				
$\text{AgBr}\cdot\text{CH}_3\text{NH}_2$			c		-42.8				
$\text{AgI}\cdot\frac{1}{2}\text{CH}_3\text{NH}_2$			c		-22.5				
$\text{AgI}\cdot\text{CH}_3\text{NH}_2$			c		-32.1				
AgSCN			c		21.0				
			aq		42.5				
$\text{Ag}_2\text{Zn}_3$			c		-9.5				
$\text{Ag}_3\text{Hg}_4$			c		0.7				
$\text{Ag}_2\text{HgI}_4$			c						47.5

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 36-1. Gold (at. no., 79; at. wt., 197.2)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		State	Description	At 298.16°K (25°C)				
Formula	$\Delta H_f^\circ$			$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
	kcal/mole			kcal/mole	kcal/mole	cal/deg mole		cal/deg mole
Au		g		82.29	72.83	-53.383	43.120	4.9680
Au <sup>+</sup>		c		0.000	0.000	0.0000	11.4	6.03
Au <sup>++</sup>		g			296.62			
		g						
AuO <sub>3</sub> <sup>---</sup>		aq				4.25		
Au <sub>2</sub> O <sub>3</sub>	std. state, hyp. m = 1	c		19.3	39.0	-28.6	30.	
HAuO <sub>3</sub> <sup>---</sup>	std. state, hyp. m = 1	aq			-27.6	20.23		
H <sub>2</sub> AuO <sub>3</sub> <sup>-</sup>	std. state, hyp. m = 1	aq			-45.8	33.57		
Au(OH) <sub>3</sub>	H <sub>3</sub> AuO <sub>3</sub>	c			-100.0	50.80	29.	
	std. state, hyp. m = 1	aq			-61.8	45.30		



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 36-2. Gold (at. no., 79; at. wt., 197.2) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1949									
Substance		Description	State	$\Delta H^{\circ}$	$\Delta F^{\circ}$	$\Delta P^{\circ}$	$\log_{10} K_f$	$S^{\circ}$	$C_p^{\circ}$
Formula	At 0°K			At 298.16°K (25°C)					
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
AuCl			c		-8.4				
AuCl <sub>3</sub>			c		-28.3				
			aq		-32.8				
AuCl <sub>3</sub> ·2H <sub>2</sub> O			c		-167.7				
AuCl <sub>4</sub> <sup>-</sup>		std. state, hyp. m = 1	aq		-77.8	-56.2	41.19	61.	
HAuCl <sub>4</sub>		std. state, hyp. m = 1	aq		-77.8	-56.2	41.19	61.	
HAuCl <sub>4</sub> ·3H <sub>2</sub> O			c		-279.2				
HAuCl <sub>4</sub> ·4H <sub>2</sub> O			c		-356.9				
AuBr			c		-4.4				
AuBr <sub>2</sub> <sup>-</sup>		std. state, hyp. m = 1	aq		-27.1		19.86		
AuBr <sub>3</sub>			c		-13.0				
			aq		-9.2				
AuBr <sub>4</sub> <sup>-</sup>		std. state, hyp. m = 1	aq		-45.5	-38.1	27.93	75.	
HAuBr <sub>4</sub>		std. state, hyp. m = 1	aq		-45.5	-38.1	27.93	75.	
HAuBr <sub>4</sub> ·5H <sub>2</sub> O			c		-398.5				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 36-3. Gold (at. no., 79; at. wt., 197.2)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	$\Delta H_f^\circ$		At 298.16°K (25°C)		cal/deg mole
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	
AuI		c		0.2			
AuCl·NH <sub>3</sub>		c		-43.9			
AuCl·2NH <sub>3</sub>		c		-69.0			
AuCl·6NH <sub>3</sub>		c		-142.0			
AuBr·NH <sub>3</sub>		c		-38.0			
AuBr·2NH <sub>3</sub>		c		-61.1			
AuBr·3NH <sub>3</sub>		c		-80.4			
AuBr·4NH <sub>3</sub>		c		-99.6			
AuBr·6NH <sub>3</sub>		c		-137.6			
AuI·NH <sub>3</sub>		c		-26.3			
AuI·2NH <sub>3</sub>		c		-46.4			
AuI·3NH <sub>3</sub>		c		-66.4			
AuI·6NH <sub>3</sub>		c		-123.5			
AuI·8NH <sub>3</sub>		c		-158.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 36-4. Gold (at. no., 79; at. wt., 197.2)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		State	$\Delta H_f^\circ$		$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description		At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	At 298.16°K (25°C) kcal/mole	At 298.16°K (25°C) kcal/mole				
Au <sub>2</sub> P <sub>3</sub>		c			-24.1					
AuI·PH <sub>3</sub>		c			-8.4					
AuSb <sub>2</sub>		c, III			-4.8					
Au(CN) <sub>2</sub> <sup>-</sup>	std. state, hyp. m = 1	aq			58.4	51.5	-37.75	99.	19.0	
AuSn		c							12.5	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 37-1. Platinum (at. no., 78; at. wt., 195.23)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		State	At 298.16°K (25°C)				
Formula	Description		$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	
			kcal/mole	kcal/mole		cal/deg mole	
Pt		g		121.6	-81.288	45.96	6.102
Pt <sup>+</sup>		c	0.000	0.000	0.0000	10.0	6.35
Pt(OH) <sub>2</sub>		g		328.			
PtCl		c		-87.2	49.99	26.5	
PtCl <sub>2</sub>		c		-17.7			
PtCl <sub>3</sub>		c		-35.5			
PtCl <sub>4</sub>		c		-49.9			
PtCl <sub>4</sub> ·5H <sub>2</sub> O		c		-62.9			
PtCl <sub>4</sub> ---	std. state, hyp. m = 1	aq		-82.4			
PtCl <sub>6</sub> ---	std. state, hyp. m = 1	aq		-425.8			
				-123.4	67.36	42.	
				-167.4	90.231	52.6	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES I

Washington, D.C.

Table 37-2. Platinum (at. no., 78; at. wt., 195.23) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1949									
Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	Description		At 0°K kcal/mole	kcal/mole					
$\text{HPtCl}_5 \cdot 2\text{H}_2\text{O}$		c		-249.8					
$\text{H}_2\text{PtCl}_6$		aq		-167.3					
$\text{H}_2\text{PtCl}_6 \cdot 6\text{H}_2\text{O}$		c		-572.9					
$\text{PtBr}_4$		c		-41.3					
$\text{PtBr}_4^{--}$		aq		-51.1					
$\text{PtBr}_6^{--}$		aq		-91.1					
		aq		-117.1					
$\text{H}_2\text{PtBr}_6$		aq		-117.0					
$\text{H}_2\text{PtBr}_6 \cdot 9\text{H}_2\text{O}$		c		-734.7					
$\text{PtI}_2$		c		-21.6					
$\text{PtI}_4$		c		-55.7					
$\text{PtI}_6^{--}$		aq							
$\text{PtS}$		c		-20.8					
$\text{PtS}_2$		c		-27.8					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 37-3. Platinum (at. no., 78; at. wt., 195.23)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1949

Formula	Substance	Description	State	$\Delta H_f^\circ$		$\log_{10} f$	$S^\circ$	$C_p^\circ$
				At 298.16°K (25°C)				
				At 0°K kcal/mole	kcal/mole		kcal/mole	cal/deg mole
$(\text{NH}_4)_2\text{PtCl}_4$			c	-195.3				
$\text{Pt}(\text{OH})_2 \cdot 4\text{NH}_3$			aq	-186.9				
$\text{PtCl}_2 \cdot 2\text{NH}_3$			c	-202.9				
$\text{PtCl}_2 \cdot 4\text{NH}_3$			c	-120.3				
			c	-184.0				
			aq	-172.9				
$\text{PtCl}_2 \cdot 4\text{NH}_3 \cdot \text{H}_2\text{O}$			c	-250.				
$\text{PtCl}_2 \cdot 5\text{NH}_3$			c	-205.3				
$\text{PtI}_2 \cdot 2\text{NH}_3$			c					
$\text{PtI}_2 \cdot 4\text{NH}_3$			c					
$\text{PtI}_2 \cdot 6\text{NH}_3$			c					
$\text{PtSO}_4 \cdot 4\text{NH}_3$			c	-309.0				17.
$\text{PtSb}_2$			c, II					12.
$\text{PtSn}$			c					
$\text{Ag}_2\text{PtCl}_6$			c	-131.6				
$\text{Ag}_2\text{PtBr}_6$			c	-98.8				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 38-1. Iridium (at. no., 77; at. wt., 193.1) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1949; March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\Delta P_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K kcal/mole			kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole	
Ir			g		165.	154.	-112.9	46.25	4.968
Ir <sup>+</sup>			c	0.000	0.000	0.000	0.0000	8.7	5.9
IrO <sub>2</sub>			g		378.				
			c		-40.1				13.7
IrF <sub>6</sub>			liq						
IrCl			c		-130.				
IrCl <sub>2</sub>			c		-22.3				
IrCl <sub>3</sub>			c		-42.8				
IrCl <sub>6</sub> <sup>---</sup>			c		-61.5				
IrCl <sub>6</sub> <sup>---</sup>			aq		-154.				
IrCl <sub>6</sub> <sup>---</sup>			aq		-189.				
IrS <sub>2</sub>			c		-30.				
Ir <sub>2</sub> S <sub>3</sub>			c		-51.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 39-1. Osmium (at. no. 76; at. wt., 190.2)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1949

Substance										
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} f_f$	$S^\circ$	$C_p^\circ$		
			At 0°K kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole		
Os		g		174.	163.	-119.5	45.97	4.968		
Os <sup>+</sup>		c	0.000	0.000	0.000	0.0000	7.8	5.9		
		g		376.						
OsO <sub>4</sub>		g		-79.9	-67.9	49.77	65.6			
	white	c, I		-91.7	-70.5	51.68	34.7			
	yellow	c, II		-93.4	-70.7	51.82	29.7			
	std. state, hyp. m = 1	aq			-68.59	50.276				
HOsO <sub>5</sub> <sup>-</sup>		aq			-108.78	79.7342				
H <sub>2</sub> OsO <sub>5</sub>		aq			-125.28	91.8285				
OsS <sub>2</sub>		c		-35.						
OsP <sub>2</sub>		c		-40.						



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 40-1. Rhenium (at. no., 75; at. wt., 186.31) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1949									
Substance		State	Description	At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula									
Re		g		0.000	189.	179.	-131.2	45.13	4.968
ReO <sub>3</sub>		c			-83.				
ReO <sub>4</sub> <sup>-</sup>		aq			-190.3				
Re <sub>2</sub> O <sub>7</sub>		c			-297.5				
Re <sub>2</sub> O <sub>8</sub>		c			-148.3				
HReO <sub>4</sub>	in 5000 H <sub>2</sub> O	aq			-188.6				
	7000 H <sub>2</sub> O	aq			-188.9				
	10,000 H <sub>2</sub> O	aq			-189.2				
	20,000 H <sub>2</sub> O	aq			-189.4				
ReF <sub>6</sub>		g			-273.				
ReS <sub>2</sub>		c			-44.3				
ReAs <sub>2</sub>		c			1.				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 41-1. Palladium (at. no., 46; at. wt., 106.7)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949; March 31, 1950

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta P_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)		cal/deg mole	cal/deg mole
			kcal/mole	kcal/mole	kcal/mole		
Pd		g		93.	-61.6	39.91	4.968
Pd <sup>+</sup>		c	0.000	0.000	0.0000	8.9	6.3
Pd <sup>++</sup>		g		276.			
		g		734.			
PdO		c		-20.4			7.5
Pd <sub>2</sub> H		c		-8.9			
Pd(OH) <sub>2</sub>		c		-92.1			
Pd(OH) <sub>4</sub>		c		-169.4			
PdCl <sub>2</sub>		c		-45.4			
PdCl <sub>4</sub> <sup>---</sup>		aq		-128.3	70.88	41.	
PdCl <sub>6</sub> <sup>---</sup>	std. state, hyp. m = 1	aq		-152.			
H <sub>2</sub> PdCl <sub>4</sub>		aq		-129.3			
H <sub>2</sub> PdCl <sub>6</sub>		aq		-152.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 41-2. Palladium (at. no., 46; at. wt., 106.7)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949; March 31, 1950

Formula	Substance	Description	State	At 0°K				At 298.16°K (25°C)			
				kcal/mole				kcal/mole			
PdBr <sub>2</sub>			c								
PdBr <sub>4</sub>			aq								
PdI <sub>2</sub>			c								
PdI <sub>2</sub> ·H <sub>2</sub> O			c								
PdCl <sub>2</sub> ·2NH <sub>3</sub>			c								
PdCl <sub>2</sub> ·4NH <sub>3</sub>			c								
PdI <sub>2</sub> ·2NH <sub>3</sub>			c								
PdI <sub>2</sub> ·4NH <sub>3</sub>			c								
PdSb			c								11.9
PdSb <sub>2</sub>			c								17.3
PdSb <sub>3</sub>			c								24.8
Pd(CN) <sub>2</sub>			c								
PdCu			c								10.6
PdCu <sub>3</sub>			c								22.2

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 42-1. Rhodium (at. no., 45; at. wt., 102.91)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)			cal/deg mole
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	
Rh		g		138.	127.	44.39	5.001
Rh <sup>+</sup>		c	0.000	0.000	0.0000	7.6	6.1
Rh <sup>++</sup>		g		306.			
RhO		g					
Rh <sub>2</sub> O		c		-21.7			11.5
Rh <sub>2</sub> O <sub>3</sub>		c		-22.7			17.5
		c		-68.3			24.9
RhCl		c		-16.			
RhCl <sub>2</sub>		c		-36.			
RhCl <sub>3</sub>		c		-56.			
RhCl <sub>6</sub> ---		aq		-207.8			
Rh <sub>2</sub> S <sub>3</sub>		c					
Rh <sub>2</sub> S <sub>5</sub>		c					
Rh <sub>3</sub> S <sub>4</sub>		c					
Rh <sub>3</sub> S <sub>8</sub>		c					
RhP <sub>2</sub>		c					
RhP <sub>3</sub>		c					

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## National Bureau of Standards

Table 43-1. Ruthenium (at. no., 44; at. wt., 101.7)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		State	Description	$\Delta H^{\circ}$		$\log_{10} f$	$S^{\circ}$	$C_p^{\circ}$
Formula	At 0°K			$\Delta H^{\circ}$				
					kcal/mole			
				At 298.16°K (25°C)				
				kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
Ru		g		0.000	160.	-109.2	44.57	
Ru <sup>+</sup>		c, IV						
Ru <sup>++</sup>		g				0.0000	6.9	5.6
RuO <sub>2</sub>		c			-52.5			
RuCl <sub>3</sub>		c			-63.			
RuS <sub>2</sub>		c			-48.1			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards  
Washington, D.C.

Table 44-1. Technetium (at. no., 43; at. wt., 99) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1949 (Corrected)									
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K					
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
Tc			c	0.000	0.000	0.000	0.0000	9.	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 45-1. Nickel [at. no., 28; at. wt., 58.69]

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		State	Description	At 298.16°K (25°C)				S°	C <sub>p</sub> <sup>o</sup>
Formula				ΔH <sup>fo</sup>	ΔH <sup>fo</sup>	ΔF <sup>fo</sup>	log <sub>10</sub> f <sup>o</sup>		
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
Ni		g		101.14	101.61	90.77	-66.533	43.592	5.5986
		c, II		0.000	0.000	0.000	0.0000	7.20	6.21
Ni <sup>+</sup>		g		277.16	279.11				
Ni <sup>++</sup>		g		697.4	700.8				
		aq	std. state, hyp. m=1		-15.3	-11.1	8.136	-38.1	
Ni <sup>+++</sup>		g							
NiO		g		59.3	59.3	51.8	-37.97	57.	
		c		-57.8	-58.4	-51.7	37.90	9.22	10.60
NiH		g		93.	93.				
		c			-2.7				
NiH <sub>2</sub>		c			-6.2				
Ni(OH) <sub>2</sub>		c			-128.6	-108.3	79.382	19.	
Ni(OH) <sub>3</sub>		c			-162.1				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 45-2. Nickel (at. no., 28; at. wt., 58.69) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)											
March 31, 1949											
Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
Formula				At 0°K	At 298.16°K (25°C)						
				kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole		
NiF <sub>2</sub>			c		-159.5						
NiF <sub>2</sub> ·4H <sub>2</sub> O			aq		-171.5						
			c			-379.9	278.46				
NiCl			g	15.	15.						
NiCl <sub>2</sub>			c		-75.5	-65.1	47.72	25.6	18.6		
	in		aq		-93.20						
	25 H <sub>2</sub> O		aq		-93.88						
	50 H <sub>2</sub> O		aq		-94.20						
	100 H <sub>2</sub> O		aq		-94.44						
	200 H <sub>2</sub> O		aq		-94.70						
	500 H <sub>2</sub> O		aq		-94.8						
	800 H <sub>2</sub> O		aq		-94.84						
	1000 H <sub>2</sub> O		aq		-94.96						
	2000 H <sub>2</sub> O		aq		-95.08						
	5000 H <sub>2</sub> O		aq		-95.13						
	10000 H <sub>2</sub> O		aq								
NiCl <sub>2</sub> ·2H <sub>2</sub> O			c		-220.8						
NiCl <sub>2</sub> ·4H <sub>2</sub> O			c		-364.7						
NiCl <sub>2</sub> ·6H <sub>2</sub> O			c		-505.8	-410.5	300.89	75.2			



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 45-3. Nickel (at. no., 28; at. wt., 58.69) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1949									
Substance			State	Description	At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	log <sub>10</sub> K <sub>f</sub>	S° cal/deg mole	C <sub>p</sub> ° cal/deg mole
Formula		At 0°K kcal/mole							
NiBr <sub>2</sub>		c				-54.2			
NiBr <sub>2</sub> ·3H <sub>2</sub> O		aq				-73.1			
		c				-277.8			
NiI <sub>2</sub>		c				-20.5			
		aq				-42.0			
Ni(IO <sub>3</sub> ) <sub>2</sub>		c				-124.5			
		aq				-125.1			
Ni(IO <sub>3</sub> ) <sub>2</sub> ·2H <sub>2</sub> O		c, I				-264.9			
		c, II				-263.9			
Ni(IO <sub>3</sub> ) <sub>2</sub> ·4H <sub>2</sub> O		c				-401.7			
NiS		c, I				-17.5			
		c, II				-18.6			
Ni <sub>3</sub> S <sub>2</sub>		c				-43.4			
				precipitated					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 45-4. Nickel (at. no., 28; at. wt., 58.69)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K		At 298.16°K (25°C)		cal/deg mole
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	
$\text{NiSO}_4$	std. state, hyp. m=1	c					
	in 200 $\text{H}_2\text{O}$	aq		-213.0	-184.9	135.53	33.4
	500 $\text{H}_2\text{O}$	aq		-232.2	-188.4	138.09	
	1000 $\text{H}_2\text{O}$	aq		-230.3			
	2000 $\text{H}_2\text{O}$	aq		-230.41			
	5000 $\text{H}_2\text{O}$	aq		-230.52			
	10000 $\text{H}_2\text{O}$	aq		-230.63			
	20000 $\text{H}_2\text{O}$	aq		-230.78			
	50000 $\text{H}_2\text{O}$	aq		-230.90			
	$\infty \text{H}_2\text{O}$	aq		-231.03			
$\text{NiSO}_4 \cdot 6\text{H}_2\text{O}$	green	c, I		-231.14			
	blue	c, II		-232.2			
$\text{Ni}_2\text{S}_2\text{O}_6$				-644.98			
				-642.5	-531.0	389.22	82.
$\text{Ni}_2\text{S}_2\text{O}_6 \cdot 6\text{H}_2\text{O}$				-712.9			
				-295.6			
		c		-707.9			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Substance		HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)				
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$C_p^\circ$
			At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole	cal/deg mole
NiSe		c				
NiTe		c				
Ni(N <sub>3</sub> ) <sub>2</sub> ·H <sub>2</sub> O	nickel azide hydrate	c				
Ni(NO <sub>3</sub> ) <sub>2</sub>		c				
Ni(NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O		aq				
Ni(NH <sub>3</sub> ) <sub>4</sub> <sup>++</sup>		c				
Ni(NH <sub>3</sub> ) <sub>6</sub> <sup>++</sup>	std. state, hyp. m = 1	aq				
Ni(NH <sub>3</sub> ) <sub>6</sub> <sup>++</sup>	std. state, hyp. m = 1	aq				
Ni(NO <sub>3</sub> ) <sub>2</sub> ·6NH <sub>3</sub>		c, l				
NiCl <sub>2</sub> ·NH <sub>3</sub>		c				
NiCl <sub>2</sub> ·2NH <sub>3</sub>		c				
NiCl <sub>2</sub> ·6NH <sub>3</sub>		c				
NiBr <sub>2</sub> ·NH <sub>3</sub>		c				
NiBr <sub>2</sub> ·2NH <sub>3</sub>		c				
NiBr <sub>2</sub> ·6NH <sub>3</sub>		c				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

Washington, D.C.

National Bureau of Standards

Table 45-6. Nickel (at. no., 28; at. wt., 58.69)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance									
Formula	Description	State	$\Delta F^\circ$	$\Delta F^\circ$	$\Delta F^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 0°K kcal/mole	kcal/mole	kcal/mole				
						At 298.16°K (25°K)	cal/deg mole	cal/deg mole	
NiI <sub>2</sub> ·2NH <sub>3</sub>		c		-83.1					
NiI <sub>2</sub> ·6NH <sub>3</sub>		c		-192.4					
NiSO <sub>4</sub> ·½NH <sub>3</sub>		c		-231.0					
NiSO <sub>4</sub> ·2NH <sub>3</sub>		c		-275.8					
NiSO <sub>4</sub> ·4NH <sub>3</sub>		c		-330.6					
NiSO <sub>4</sub> ·6NH <sub>3</sub>		c		-382.1					
Ni <sub>2</sub> P		c		-44.					
Ni <sub>3</sub> P		c		-53.					
Ni <sub>5</sub> P <sub>2</sub>		c		-105.					
NiSb		c		-15.6					
Ni <sub>5</sub> Sb <sub>2</sub>		c		-36.4					
Ni <sub>3</sub> C		c		11.0					
NiCO <sub>3</sub>		c			-146.7	107.53		97.	
Ni(CO) <sub>4</sub>		g liq						74.	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 45-7. Nickel (at. no., 28; at. wt., 58.69)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Table 45-7. Nickel (at. no., 28; at. wt., 58.69)											
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)											
March 31, 1949											
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
				At 298.16°K (25°C)						cal/deg mole	cal/deg mole
				kcal/mole	kcal/mole	kcal/mole					
NiBr <sub>2</sub> ·6CH <sub>3</sub> OH			c		-453.4						
Ni(CN) <sub>2</sub>			c		27.1						
Ni(CN) <sub>4</sub> <sup>2-</sup>			aq		86.9						
NiSi			c		-20.6				10.9		
Ni <sub>2</sub> Si			c		-33.6				16.8		
NiSn			c		-14.8				23.8		
Ni <sub>3</sub> Sn			c		-37.5						
Ni <sub>3</sub> Sn <sub>2</sub>			c		-85.7						
2NiI <sub>2</sub> ·PbI <sub>2</sub>			c		-341.9						
2NiI <sub>2</sub> ·FeI <sub>2</sub> ·3H <sub>2</sub> O			c								

## SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

# SERIES I

Washington, D.C.

Table 46-1. Cobalt (at. no., 27; at. wt., 58.94)

Table 46-1. Cobalt (at. no., 27; at. wt., 58.94)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} f$	$S^\circ$	$C_p^\circ$
At 0°K					At 298.16°K (25°C)					
kcal/mole					kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
Co				g c, III		105. 0.000	94. 0.000	-68.9 0.0000	42.881 6.8	5.5043 6.11
Co <sup>+</sup>				g		287.27				
Co <sup>++</sup>				g		687.7				
			std. state, hyp. m=1	aq		-16.1	-12.3	9.016	-37.1	
Co <sup>+++</sup>				g						
			std. state, hyp. m=1	aq			29.6	-21.70		
CoO				c		-57.2	-51.0	37.38	10.5	
Co <sub>3</sub> O <sub>4</sub>				c		-210.				
CoH				c		-4.1				
CoH <sub>2</sub>				c		-10.2				
Co(OH) <sub>2</sub>				c		-131.2				
Co(OH) <sub>3</sub>				c		-176.6				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 46-2. Cobalt (at. no., 27; at. wt., 58.94)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K		At 298.16°K (25°C)		cal/deg mole
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	
CoF <sub>2</sub>		c		-159.			
CoF <sub>2</sub> ·4H <sub>2</sub> O		aq		-173.6			
CoF <sub>3</sub>		c			279.19		
		c		-187.			
CoCl <sub>2</sub>		c		-77.8		49.48	18.8
		aq		-96.1		25.4	
CoCl <sub>2</sub> ·2H <sub>2</sub> O		c		-222.9			
CoCl <sub>2</sub> ·4H <sub>2</sub> O		c		-367.2			
CoCl <sub>2</sub> ·6H <sub>2</sub> O		c		-508.9			
CoBr <sub>2</sub>		c		-55.5			
		aq		-73.9			
CoBr <sub>2</sub> ·6H <sub>2</sub> O		c		-485.1			
CoI <sub>2</sub>		c		-24.4			
		aq		-42.8			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 46-3. Cobalt (at. no., 27; at. wt., 58.94) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1949									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				
Formula									
$\text{Co}(\text{IO}_3)_2$			c		-124.3				
$\text{Co}(\text{IO}_3)_2 \cdot 2\text{H}_2\text{O}$			aq		-125.9				
$\text{Co}(\text{IO}_3)_2 \cdot 4\text{H}_2\text{O}$			c		-264.9				
			c		-401.9				
$\text{CoS}$			c, I		-20.2				11.4
		precipitated	c, II		-21.4				
$\text{Co}_2\text{S}_3$			c		-51.				
$\text{CoSO}_4$			c		-207.5	-182.1	133.48	27.1	
$\text{CoSO}_4 \cdot 6\text{H}_2\text{O}$			aq		-232.0				
$\text{CoSO}_4 \cdot 7\text{H}_2\text{O}$			c		-643.2				
			c		-713.8				96.
$\text{CoSe}$			c		-10.				
$\text{CoTe}$			c		-9.				



## SERIES I

Washington, D.C.

Table 46-4. Cobalt (at. no., 27; at. wt., 58.94)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance			State	Description	At 0°K	ΔH° kcal/mole	ΔF° kcal/mole	ΔP° kcal/mole	log <sub>10</sub> f <sub>r</sub> (25°C)	S° cal/deg mole	C <sub>p</sub> <sup>o</sup> cal/deg mole			
Formula		At 0°K										kcal/mole	kcal/mole	kcal/mole
Co(NO <sub>3</sub> ) <sub>2</sub>		c				-102.9								
Co(NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O		aq				-114.8								
Co(NH <sub>3</sub> ) <sub>6</sub> <sup>++</sup>	std. state, hyp. m = 1	c				-529.7								
Co(NH <sub>3</sub> ) <sub>6</sub> <sup>+++</sup>	std. state, hyp. m = 1	aq						-57.2	41.93					
		aq						-54.5	39.95					
[Co(NH <sub>3</sub> ) <sub>5</sub> H <sub>2</sub> O] <sup>+++</sup>	std. state, hyp. m = 1	aq				-192.9		-106.2	77.843	73.4				
[Co(NH <sub>3</sub> ) <sub>5</sub> NO <sub>3</sub> ] <sup>++</sup>		aq				-173.0								
[Co(NH <sub>3</sub> ) <sub>5</sub> NO <sub>3</sub> ](NO <sub>3</sub> ) <sub>2</sub>		c				-288.0								
		aq				-271.7								
		c				-357.4								
		aq				-340.7								
[Co(NH <sub>3</sub> ) <sub>5</sub> H <sub>2</sub> O]F <sub>3</sub>		c				-435.7								

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 46-5. Cobalt (at. no., 27; at. wt., 58.94)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		State	Description	At 298.16°K (25°C)					log <sub>10</sub> <i>f</i> <sub>f</sub>	S°	C <sub>p</sub> <sup>o</sup>
Formula	Δ <i>H</i> <sub>f</sub> <sup>o</sup>			Δ <i>H</i> <sub>f</sub> <sup>o</sup>	Δ <i>F</i> <sub>f</sub> <sup>o</sup>		cal/deg mole				
					kcal/mole	kcal/mole		cal/deg mole			
CoCl <sub>2</sub> ·NH <sub>3</sub>		c		-110.3							
CoCl <sub>2</sub> ·2NH <sub>3</sub>	rose	c, I		-140.2							
	blue	c, II									
CoCl <sub>2</sub> ·2N <sub>2</sub> H <sub>4</sub>		c		-100.4							
[Co(NH <sub>3</sub> ) <sub>4</sub> Cl <sub>2</sub> ] <sup>+</sup>	<i>cis</i>	aq		-171.4							
	<i>trans</i>	aq		-173.3							
[Co(NH <sub>3</sub> ) <sub>4</sub> Cl <sub>2</sub> ]Cl	<i>cis</i>	c		-220.9							
		aq		-211.4							
	<i>trans</i>	c		-221.6							
		aq		-213.3							
[Co(NH <sub>3</sub> ) <sub>5</sub> Cl] <sup>++</sup>	std. state, hyp. m = 1	aq		-162.1	-86.2			63.18	96.1		
[Co(NH <sub>3</sub> ) <sub>5</sub> Cl]Cl <sub>2</sub>		c		-254.5							
		aq		-242.1							
CoCl <sub>2</sub> ·6NH <sub>3</sub>		c		-240.2							
CoCl <sub>3</sub> ·6NH <sub>3</sub>		c									
[Co(NH <sub>3</sub> ) <sub>5</sub> H <sub>2</sub> O]Cl <sub>3</sub>		c		-319.0							
		aq		-312.5							
											76.7

76.7

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 46-6. Cobalt (at. no., 27; at. wt., 58.94)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance		State	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description							
$\text{CoBr}_2 \cdot \text{NH}_3$		c		-88.2				
$\text{CoBr}_2 \cdot 2\text{NH}_3$	rose	c, I		-120.1				
	blue	c, II						
$\text{CoBr}_2 \cdot 2\text{N}_2\text{H}_4$		c		-80.2				
$[\text{Co}(\text{NH}_3)_5\text{Br}]^{++}$		aq		-152.7				
$[\text{Co}(\text{NH}_3)_5\text{Br}]\text{Br}_2$		c		-223.2				
		aq		-210.5				
$\text{CoBr}_2 \cdot 6\text{NH}_3$		c		-222.8				
$[\text{Co}(\text{NH}_3)_5\text{H}_2\text{O}]\text{Br}_3$		c		-289.7				
		aq		-280.5				
$\text{CoI}_2 \cdot 2\text{NH}_3$	blue	c, I		-87.6				
	green	c, II						
$\text{CoI}_2 \cdot 6\text{NH}_3$		c		-193.1				69.2
$\text{CoI}_3 \cdot 6\text{NH}_3$		c						74.3

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 46-7. Cobalt (at. no., 27; at. wt., 58.94)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1949

Substance								
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
			At 0°K					
			kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
$\text{CoSO}_4 \cdot \frac{1}{2}\text{NH}_3$		c		-292.5				
$\text{CoSO}_4 \cdot 2\text{NH}_3$		c		-338.2				
$\text{CoSO}_4 \cdot 3\text{NH}_3$		c		-366.0				
$\text{CoSO}_4 \cdot 4\text{NH}_3$		c		-391.9				
$[\text{Co}(\text{NH}_3)_5\text{SO}_4]^+$		aq			-231.9	169.98		
$\text{CoSO}_4 \cdot 6\text{NH}_3$	std. state, hyp. m = 1	c		-441.7				
CoP		c		-35.				32.9
CoP <sub>3</sub>		c		-65.				12.2
Co <sub>2</sub> P		c		-47.3				
CoAs <sub>2</sub> ·CoS <sub>2</sub>		c						
CoSb		c		-10.0				
CoSb <sub>2</sub>		c		-13.2				
Co <sub>3</sub> C		c		9.5	7.1	-5.20	29.8	
CoCO <sub>3</sub>		c		-172.7				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES I

Washington, D.C.

Table 46-8. Cobalt (at. no., 27; at. wt., 58.94) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1949 (Corrected)														
Substance			State	Description	Formula	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$			
						At 298.16°K (25°C)								
						kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole			
			c					-221.9						
			c					-295.1						
			c					-422.1						
			c					-183.3						
			c					-199.8						
			c					-290.2						
			c					-255.8						
			c					-401.8						
			c					-114.2						
	cis		aq					-127.1						
	trans		aq					-128.9						
	cis		c					-175.1						
			aq					-167.1						
			c					-174.2						
	trans		aq					-168.9						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 46-9. Cobalt (at. no., 27; at. wt., 58.94)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1949 (Corrected)

Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description		At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				
$[\text{Co}(\text{C}_2\text{H}_5(\text{NH}_2)_2)_2\text{Cl}]_2\text{Cl} \cdot \text{NH}_3$	<i>cis</i>	c		-202.				
	<i>trans</i>	c		-198.				
$[\text{Co}(\text{C}_2\text{H}_5(\text{NH}_2)_2)_2\text{Cl}]_2\text{Cl} \cdot 2\text{NH}_3$	<i>cis</i>	c		-226.				
	<i>trans</i>	c		-219.				
$[\text{Co}(\text{C}_2\text{H}_5(\text{NH}_2)_2)_2\text{Cl}]_2\text{Cl} \cdot 4\text{NH}_3$	<i>cis</i>	c		-268.				
	<i>trans</i>	c		-259.				
$[\text{Co}(\text{C}_2\text{H}_5(\text{NH}_2)_2)_2\text{Cl}]_2\text{Cl} \cdot 6\text{NH}_3$	<i>cis</i>	c		-308.				
$\text{CoCl}_2 \cdot 3\text{C}_2\text{H}_5(\text{NH}_2)_2$		c		-160.5				
$\text{CoBr}_2 \cdot 4\frac{1}{2}\text{C}_2\text{H}_5(\text{NH}_2)_2$		c		-109.2				
$\text{CoBr}_2 \cdot 3\text{C}_2\text{H}_5(\text{NH}_2)_2$		c		-145.3				
$\text{CoI}_2 \cdot 3\text{C}_2\text{H}_5(\text{NH}_2)_2$		c		-116.9				
$\text{Co}(\text{C}_2\text{H}_5\text{SO}_4)_2$	cobalt ethylsulfate	aq		-433.7				
CoSi		c		-24.0				
CoSi <sub>2</sub>		c		-24.6				
CoSi <sub>3</sub>		c		-25.6				
Co <sub>2</sub> Si		c		-27.6				
Co <sub>2</sub> Sn		c						19.

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

		Table 47-1. Iron (at. no., 26; at. wt., 55.85) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1949						
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	At 298.16°K (25°C)		cal/deg mole	cal/deg mole
				kcal/mole	kcal/mole	kcal/mole		
Fe			g		96.68	85.76	43.11	6.13
Fe <sup>+</sup>			c	0.000	0.000	0.0000	6.49	6.03
Fe <sup>++</sup>			g		277.9			
			g		650.7			
Fe <sup>+++</sup>		std. state, hyp. m=1	aq		-21.0	-20.30	-27.1	
			g		1354.			
Fe <sup>++++</sup>		std. state, hyp. m=1	aq		-11.4	-2.52	-70.1	
			g					
Fe <sub>0.95</sub> O		"FeO" wustite	c		-63.7	-58.4	12.9	
Fe <sub>2</sub> O <sub>3</sub>		hematite	c		-196.5	-177.1	21.5	25.0
Fe <sub>3</sub> O <sub>4</sub>		magnetite	c		-267.0	-242.4	35.0	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 47-2. Iron (at. no., 26; at. wt., 55.85) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1949								
Substance			State	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description	At 0°K kcal/mole						
FeH <sub>2</sub> FeH <sub>3</sub>		c c						
Fe(OH) <sup>++</sup> FeO(OH) Fe(OH) <sub>2</sub> Fe(OH) <sub>2</sub> <sup>+</sup> Fe(OH) <sub>3</sub>	std. state, hyp. m = 1	aq c c aq c		-67.4 -135.8 -197.0	-55.91 -115.57 -106.2	40.981 84.7112 77.843	-23.2 19.	
FeF <sub>2</sub> FeF <sub>3</sub>		aq aq		-177.8 -243.1				



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 47-3. Iron (at. no., 26; at. wt., 55.85) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1949										
Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula				At 0°K	At 298.16°K (25°C)					
						kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
FeCl <sup>++</sup>		std. state, hyp. m=1	aq			-42.9	-35.9	-22.	18.25	
FeCl <sub>2</sub>		std. state, hyp. m=1	c			-81.5	-72.2	28.6		
		in 200 H <sub>2</sub> O	aq			-101.0				
		500 H <sub>2</sub> O	aq			-100.40				
		1000 H <sub>2</sub> O	aq			-100.6				
		5000 H <sub>2</sub> O	aq			-100.70				
FeCl <sub>2</sub> ·2H <sub>2</sub> O			aq			-100.95				
FeCl <sub>2</sub> ·4H <sub>2</sub> O			c			-228.2				
FeCl <sub>3</sub>			c			-370.7				
			c			-96.8				
			aq			-127.9				
FeCl <sub>3</sub> ·6H <sub>2</sub> O			c			-532.0				

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Washington, D.C.

Table 47-4. Iron (at. no., 26; at. wt., 55.85)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1949

Substance								
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} f_f$	$S^\circ$	$C_p^\circ$	
			At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole		cal/deg mole		
FeBr <sup>++</sup>	std. state, hyp. m =1	aq		-34.2	20.45	-28.		
FeBr <sub>2</sub>		c		-60.02				
		aq		-79.1				
FeBr <sub>3</sub>		aq		-98.1				
FeI <sub>2</sub>		c		-29.98				
		aq		-49.03				
FeS		$\alpha$	c		-22.72	17.093	16.1	13.1
		$\beta$	c		-21.35			
	pyrites	c		-42.52	29.202	12.7	14.8	
FeS <sub>2</sub>	markasite	c		-36.88				

## Washington, D.C.

## National Bureau of Standards

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Substance						
Formula	Description	State	$\Delta E_f^\circ$	$\Delta E_f^\circ$	$\log_{10} K_f$	$S^\circ$
			At 0°K	At 298.16°K (25°C)		
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole
FeSO <sub>4</sub>		c		-220.5		
FeSO <sub>4</sub> ·H <sub>2</sub> O	in 200 H <sub>2</sub> O	aq		-236.0		
FeSO <sub>4</sub> ·4H <sub>2</sub> O		c		-296.3		
FeSO <sub>4</sub> ·7H <sub>2</sub> O		c		-507.5		
		c		-718.7		
Fe <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c				
	in 400 H <sub>2</sub> O	aq		-653.0		
	1000 H <sub>2</sub> O	aq		-653.20		
	2000 H <sub>2</sub> O	aq		-653.41		
	3000 H <sub>2</sub> O	aq		-653.62		

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 47-6. Iron (at. no., 26; at. wt., 55.85)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1949

Formula	Substance	Description	State	At 0°K					At 298.16°K (25°C)		
				$\Delta H_f^\circ$ kcal/mole	$\Delta G_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole		
$\text{Fe}(\text{HSO}_4)_3$			aq		-645.7						
$\text{FeSe}$			c		-16.5						
$\text{FeTe}$			amorp		-13.9						
		precipitated	c		-18.6						
$\text{Fe}_2\text{N}$			c		-0.9	2.6	-1.91	24.2	16.8		
$\text{Fe}_4\text{N}$			c		-2.55	0.89	-0.652	37.3	29.3		
$\text{Fe}(\text{NO}_3)_3$			aq		-160.4						
$\text{Fe}(\text{NO}_3)_3 \cdot 9\text{H}_2\text{O}$			c		-784.4						
$\text{Fe}(\text{NO})\text{Cl}_2$			aq		-89.7						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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## SERIES I

Washington, D.C.

Table 47-7. Iron (at. no., 26; at. wt., 55.85)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)			
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
$\text{FeCl}_2 \cdot \text{NH}_3$		c		-115.0			
$\text{FeCl}_2 \cdot 2\text{NH}_3$		c		-144.3			
$\text{FeCl}_2 \cdot 6\text{NH}_3$		c		-237.8			
$\text{FeCl}_3 \cdot 6\text{NH}_3$		c		-214.7			
$\text{FeCl}_2 \cdot 10\text{NH}_3$		c		-312.4			
$\text{FeBr}_2 \cdot \text{NH}_3$		c		-91.84			
$\text{FeBr}_2 \cdot 2\text{NH}_3$		c		-122.75			
$\text{FeBr}_2 \cdot 6\text{NH}_3$		c		-220.3			
$\text{FeBr}_3 \cdot 6\text{NH}_3$		c					
$\text{FeI}_2 \cdot 2\text{NH}_3$		c		-92.98			
$\text{FeI}_2 \cdot 6\text{NH}_3$		c		-195.1			
$\text{Fe}(\text{NO})\text{SO}_4$		aq		-224.5			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 47-8. Iron (at. no., 26; at. wt., 55.85)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Formula	Substance	Description	State	At 298.16°K (25°C)				
				$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
$\text{FeSO}_4 \cdot \text{NH}_3$			c		-252.0			
$\text{FeSO}_4 \cdot 2\text{NH}_3$			c		-280.9			
$\text{FeSO}_4 \cdot 3\text{NH}_3$			c		-306.0			
$\text{FeSO}_4 \cdot 4\text{NH}_3$			c		-330.6			
$\text{FeSO}_4 \cdot 6\text{NH}_3$			c		-378.2			
$\text{Fe}_2(\text{SO}_4)_3 \cdot 12\text{NH}_3$			c					
FeP			c		-28.			
$\text{FeP}_2$			c		-42.			
$\text{Fe}_2\text{P}$			c		-36.			
$\text{Fe}_3\text{P}$			c		-40.			
$\text{FePO}_4$			c		-299.6			
$\text{FePO}_4 \cdot 2\text{H}_2\text{O}$			c		-440.8			
$\text{FePO}_4 \cdot 4\text{H}_2\text{O}$			c		-578.8			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Washington, D.C.

Table 47-9. Iron (at. no., 26; at. wt. 55.85) HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1949									
Substance									
Formula	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$		$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
			At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole			
FeSb		c		2.4					
FeSb <sub>2</sub>		c		3.6					
Fe <sub>3</sub> C	cementite	c		5.0	3.5		-2.56	25.7	25.3
FeCO <sub>3</sub>	siderite	c		-178.70	-161.06		118.055	22.2	19.63
Fe(CO) <sub>5</sub>		liq		-187.8					
Fe <sub>2</sub> (C <sub>2</sub> O <sub>4</sub> ) <sub>3</sub>	ferric oxalate in 100 H <sub>2</sub> O	aq		-611.7					
	400 H <sub>2</sub> O	aq		-611.4					
Fe(HC <sub>2</sub> O <sub>4</sub> ) <sub>3</sub>	ferric bioxalate	aq		-600.1					
Fe(C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>3</sub>	ferric acetate in 300 H <sub>2</sub> O	aq		-355.5					
	600 H <sub>2</sub> O	aq		-355.2					
	1800 H <sub>2</sub> O	aq		-353.8					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards Washington, D.C.

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Table 47-10. Iron (at. no., 26; at. wt., 55.85)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1949

Substance		Description	State	$\Delta H_f^\circ$		$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	$\Delta H_f^\circ$			At 298.16°K (25°C)				
				At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
$\text{Fe}(\text{CO})_4\text{Cl}_2$	in 1700 $\text{H}_2\text{O}$		c		-195.7			
$\text{Fe}(\text{CO})_4\text{Br}_2$			c		-188.0			
$\text{Fe}(\text{CO})_4\text{I}_2$			aq		-178.5			
$\text{Fe}(\text{CN})_6^{4-}$			c		126.7			
$\text{Fe}_4[\text{Fe}(\text{CN})_6]_3$			aq		338.6			
$\text{FeCO}(\text{CN})_5^{--}$			aq		47.2			
$\text{Fe}_2\text{CO}(\text{CN})_5$			c		98.5			



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 47-11. Iron (at. no., 26; at. wt., 55.85) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1949									
Formula	Substance	Description	State	At 298.16°K (25°C)					
				$\Delta H_f^\circ$		$\Delta F_f^\circ$		$\log_{10} K_f$	
				kcal/mole	kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
HFe(CN) <sub>6</sub> <sup>---</sup>			aq		126.9				
H <sub>2</sub> Fe(CN) <sub>6</sub> <sup>--</sup>			aq		127.1				
H <sub>3</sub> Fe(CN) <sub>6</sub>			aq		153.				
H <sub>3</sub> Fe(CN) <sub>6</sub> <sup>-</sup>			aq		127.1				
H <sub>4</sub> Fe(CN) <sub>6</sub>			c		127.8				
(NH <sub>4</sub> ) <sub>4</sub> Fe(CN) <sub>6</sub>			aq		127.4				
(NH <sub>4</sub> ) <sub>4</sub> Fe(CN) <sub>6</sub> ·6H <sub>2</sub> O			aq		1.4				
HFecO(CN) <sub>5</sub> <sup>---</sup>			c		-415.3				
H <sub>2</sub> FeCO(CN) <sub>5</sub> <sup>-</sup>			aq		47.3				
H <sub>3</sub> FeCO(CN) <sub>5</sub>			aq		47.4				
H <sub>3</sub> FeCO(CN) <sub>5</sub> ·H <sub>2</sub> O			aq		47.6				
			c		-15.7				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 47-12. Iron (at. no., 26; at. wt., 55.85) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1949							
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole	cal/deg mole
Fe(CNS) <sup>++</sup>			aq				
FeSi			c	-19.2			11.8
Fe <sub>3</sub> Si			c	-20.			
FeSiO <sub>3</sub>			c	-276.			
Fe <sub>2</sub> SiO <sub>4</sub>			c	-343.7	-319.8	234.41	31.75
2FeI <sub>2</sub> ·PbI <sub>2</sub>			c	-101.5			
Zn <sub>2</sub> Fe(CN) <sub>6</sub>			c	49.4			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 49-1. Manganese (at. no., 25; at. wt., 54.93)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949 (Corrected)

Substance			$\Delta H_f^\circ$		$\Delta F_f^\circ$	At 298.16°K (25°C)		$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K		$\Delta H_f^\circ$	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole
			kcal/mole	kcal/mole						
Mn	$\alpha$  $\gamma$	g  c, IV  c, II	68.05	68.34	58.23	-42.682	41.493	4.9680		
			0.000	0.000	0.000	0.0000	7.59	6.29		
			0.342	0.37	0.33	-0.242	7.72	6.59		
Mn <sup>+</sup> Mn <sup>++</sup> Mn <sup>+++</sup>	std. state, hyp. m = 1	g  aq  g  aq	239.382	240.863						
			599.97	601.45						
			-52.3	-53.4	39.14	-20.				
MnO		g  c, I  c, I	34.4	34.6	-86.8	63.62	14.4	10.27		
			-92.0	-124.5	-111.4	81.655	12.7	12.91		
MnO <sub>2</sub>		aq		-123.9	-101.6	74.471	45.4			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 48-2. Manganese (at. no., 25; at. wt., 54.93) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1949									
Formula	Substance	Description	State	At 298.16°K (25°C)					
				$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Mn <sub>2</sub> O <sub>3</sub>			c, II						
Mn <sub>3</sub> O <sub>4</sub>			c, I						
HMnO <sub>4</sub>			aq						
Mn(OH) <sub>2</sub>		precipitated	amorp						
Mn(OH) <sub>3</sub>		precipitated	amorp						
MnF <sub>2</sub>			c, I						
MnF <sub>3</sub>			aq						
			aq						
MnCl <sub>2</sub>			c						
MnCl <sub>2</sub> ·H <sub>2</sub> O		in 400 H <sub>2</sub> O	aq						
MnCl <sub>2</sub> ·2H <sub>2</sub> O			c						
MnCl <sub>2</sub> ·4H <sub>2</sub> O			c						
H <sub>2</sub> MnCl <sub>6</sub>			c						
			aq						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Substance		HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)				
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$C_p^\circ$
			At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	At 298.16°K (25°C) cal/deg mole	At 298.16°K (25°C) cal/deg mole
MnBr <sub>2</sub>		c		-90.7		
MnBr <sub>2</sub> ·H <sub>2</sub> O		aq		-110.1		
MnBr <sub>2</sub> ·4H <sub>2</sub> O		c		-164.4		
MnBr <sub>3</sub>		c		-367.5		
		aq		-111.		
MnI <sub>2</sub>		c		-59.3		
		aq		-79.0		
MnI <sub>2</sub> ·H <sub>2</sub> O		c		-127.9		
MnI <sub>2</sub> ·2H <sub>2</sub> O		c		-194.5		
MnI <sub>2</sub> ·4H <sub>2</sub> O		c		-327.5		
MnI <sub>2</sub> ·6H <sub>2</sub> O		c		-451.		

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 48-4. Manganese (at. no., 25; at. wt., 54.93)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Formula	Substance	Description	State	At 298.16°K (25°C)				
				$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
MnS	green red		c, I c, II		-48.8 -47.6	-49.9	36.58	18.7 11.94
MnSO <sub>4</sub>	std. state, hyp. m=1 in 30 H <sub>2</sub> O 50 H <sub>2</sub> O 100 H <sub>2</sub> O 200 H <sub>2</sub> O 400 H <sub>2</sub> O 1000 H <sub>2</sub> O 2000 H <sub>2</sub> O 3000 H <sub>2</sub> O 5000 H <sub>2</sub> O		c aq aq aq aq aq aq aq aq aq		-254.24 -269.2 -266.52 -267.08 -267.47 -267.76 -268.0 -268.28 -268.41 -268.48 -268.58	-228.48 -230.7	167.473 169.10	26.8 -16. 23.94

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Substance		HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)				
Formula	Description	State	At 298.16°K (25°C)			
			$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$C_p^\circ$ cal/deg mole
MnSO <sub>4</sub>	in 10000 H <sub>2</sub> O	aq				
	20000 H <sub>2</sub> O	aq	-268.74			
	50000 H <sub>2</sub> O	aq	-268.88			
	100000 H <sub>2</sub> O	aq	-269.02			
	∞ H <sub>2</sub> O	aq	-269.10			
		aq	-269.2			
MnSO <sub>4</sub> · H <sub>2</sub> O		c, I	-328.5			
MnSO <sub>4</sub> · 4H <sub>2</sub> O		c, II	-322.4			
MnSO <sub>4</sub> · 5H <sub>2</sub> O		c	-539.3			
MnSO <sub>4</sub> · 7H <sub>2</sub> O		c	-609.6			
		c	-750.0			78.
Mn <sub>2</sub> O <sub>6</sub>						
Mn <sub>2</sub> O <sub>6</sub> · 2H <sub>2</sub> O		aq	-332.7			
Mn <sub>2</sub> O <sub>6</sub> · 6H <sub>2</sub> O		c				66.7
		c	-744.5			57.68

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 48-6. Manganese (at. no., 25; at. wt., 54.93)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} f_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
$Mn_2(SO_4)_3$			c		-666.9				
MnSe			aq		-699.				
MnTe			c, I		-28.0	-29.2	21.40	21.7	12.20
			c, I					22.4	17.3
$Mn(N_3)_2$			c		92.2				
$Mn_5N_2$		manganese azide	c		-57.8				43.8
$Mn_8N_2$			c		-81.				61.0
$Mn(NO_3)_2$			c		-166.32				
	in 2.5 H <sub>2</sub> O		aq		-142.58				
	3 H <sub>2</sub> O		aq		-143.59				
	4 H <sub>2</sub> O		aq		-145.16				
	5 H <sub>2</sub> O		aq		-146.21				
	6 H <sub>2</sub> O		aq		-146.99				



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 48-7. Manganese (at. no., 25; at. wt., 54.93)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1949 (Corrected)

Substance		$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)		cal/deg mole
			kcal/mole	kcal/mole	kcal/mole	
$Mn(NO_3)_2$	in 8 H <sub>2</sub> O	aq		-148.03		
	10 H <sub>2</sub> O	aq		-148.69		
	15 H <sub>2</sub> O	aq		-149.66		
	25 H <sub>2</sub> O	aq		-150.47		
	50 H <sub>2</sub> O	aq		-151.08		
	100 H <sub>2</sub> O	aq		-151.41		
	200 H <sub>2</sub> O	aq		-151.53		
	400 H <sub>2</sub> O	aq		-151.59		
	1000 H <sub>2</sub> O	aq		-151.65		
$Mn(NO_3)_2 \cdot 3H_2O$		c		-355.1		
$Mn(NO_3)_2 \cdot 6H_2O$		c		-566.50		
		liq		-556.89		146.7

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

Washington, D.C.

National Bureau of Standards

Table 48-8. Manganese (at. no., 25; at. wt., 54.93)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Substance		State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description							
$MnCl_2 \cdot NH_3$		c		-148.1				
$MnCl_2 \cdot 2NH_3$		c		-175.9				
$MnCl_2 \cdot 2NH_4Cl \cdot 2H_2O$		c		-428.5				
$MnCl_2 \cdot 6NH_3$		c		-267.8				
$MnBr_2 \cdot NH_3$		c		-123.7				
$MnBr_2 \cdot 2NH_3$		c		-155.2				
$MnBr_2 \cdot 6NH_3$		c		-256.5				
$MnI_2 \cdot 2NH_3$		c		-124.7				
$MnI_2 \cdot 6NH_3$		c		-233.4				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 48-9. Manganese (at. no., 25; at. wt., 54.93) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1949									
Substance									
Formula	Description	State	$\Delta F_f^\circ$		$\Delta F_f^\circ$		$\log_{10} K_f$		$S^\circ$
			At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
MnSO <sub>4</sub> ·2NH <sub>3</sub>		c		-269.6					
MnSO <sub>4</sub> ·NH <sub>3</sub>		c		-280.6					
MnSO <sub>4</sub> ·2NH <sub>3</sub>		c		-308.9					
MnSO <sub>4</sub> ·(NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub> ·2H <sub>2</sub> O		c		-692.					
MnSO <sub>4</sub> ·(NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub> ·6H <sub>2</sub> O		c		-968.					
MnSO <sub>4</sub> ·5NH <sub>3</sub>		c		-379.2					
MnSO <sub>4</sub> ·6NH <sub>3</sub>		c		-402.0					
Mn <sub>2</sub> (PO <sub>4</sub> ) <sub>3</sub>	precipitated	c		-771.					
Mn <sub>3</sub> C		c, II		-1.		0.73		23.6	22.34
MnCO <sub>3</sub>	precipitated	c		-213.9		143.23		20.5	19.49
	std. state, hyp. m = 1	aq		-211.0					
				-213.9		131.64		-32.7	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 48-10. Manganese (at. no., 25; at. wt., 54.93)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Substance		State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description							
$Mn_2O_4$	manganese oxalate	c		-258.2				
$MnC_2O_4 \cdot 2H_2O$		c		-388.6				
$MnC_2O_4 \cdot 3H_2O$		c		-455.4				
$Mn(CHO_2)_2$	manganese formate	c		-242.2				
$Mn(CHO_2)_2 \cdot 2H_2O$		aq		-246.5				
		c		-386.0				
$Mn(C_2H_3O_2)_2$	manganese acetate	c		-273.0				
		aq		-285.2				
$Mn(C_2H_3O_2)_2 \cdot 4H_2O$		c		-556.9				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 48-11. Manganese (at. no., 25; at. wt., 54.93) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1949										
Substance		Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\Delta P_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	At 0°K						At 298.16°K (25°C)			
	kcal/mole			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole		
MnSiO <sub>3</sub>			c							
MnI <sub>2</sub> ·PbI <sub>2</sub>			g†							
MnI <sub>2</sub> ·PbI <sub>2</sub> ·3H <sub>2</sub> O			c							
			c							
ZnMnO <sub>4</sub>		std. state, hyp. m = 1	aq							

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 49-1. Chromium (at. no., 24; at. wt., 52.01)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Substance									
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 0°K kcal/mole			At 298.16°K (25°C) kcal/mole	cal/deg mole	cal/deg mole	
Cr		g	80.	80.5	69.8	-51.16	41.637	4.9680	
Cr <sup>+</sup>		c	0.000	0.000	0.000	0.0000	5.68	5.58	
Cr <sup>++</sup>		g	235.91	237.39					
		g	618.7	621.7					
Cr <sup>+++</sup>		aq		-33.2					
		g	1306.	1310.					
[Cr(6H <sub>2</sub> O)] <sup>+++</sup>	violet	aq		-471.0					
Cr <sup>++++</sup>		g	2468.	2474.					
Cr <sup>+++++</sup>		g	4151.	4158.					
Cr <sup>+++++</sup>		g	6231.	6240.					
Cr <sup>+++++</sup>		g	9946.	9956.					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 49-2. Chromium (at. no., 24; at. wt., 52.01)

HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Substance		State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description							
CrO	in 2 H <sub>2</sub> O 3 H <sub>2</sub> O 4 H <sub>2</sub> O 5 H <sub>2</sub> O 6 H <sub>2</sub> O 8 H <sub>2</sub> O 10 H <sub>2</sub> O 15 H <sub>2</sub> O 25 H <sub>2</sub> O 50 H <sub>2</sub> O 80 H <sub>2</sub> O	g	65.	65.				
CrO <sub>3</sub>		c		-138.4				
		aq		-139.37				
		aq		-139.65				
		aq		-139.87				
		aq		-140.03				
		aq		-140.14				
		aq		-140.32				
		aq		-140.43				
		aq		-140.60				
		aq		-140.78				
		aq		-140.88				
		aq		-140.9				
CrO <sub>4</sub> <sup>2-</sup>	std. state, hyp. m = 1	aq		-206.3	-168.8	123.73	9.2	

# SERIES I

Washington, D.C.

Table 49-3. Chromium (at. no., 24; at. wt., 52.01)

Table 49-3. Chromium (at. no., 24; at. wt., 52.01)

June 30, 1949

Substance							
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	
			At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C) cal/deg mole	cal/deg mole
$\text{Cr}_2\text{O}_3$		c		-269.7	-250.2	183.39	28.38
$\text{Cr}_2\text{O}_3 \cdot \text{H}_2\text{O}$		c		-358.			
$\text{Cr}_2\text{O}_3 \cdot 2\text{H}_2\text{O}$		c		-439.			
$\text{Cr}_2\text{O}_3 \cdot 3\text{H}_2\text{O}$		c		-517.			
$\text{Cr}_2\text{O}_7^{--}$	std. state, hyp. m=1	aq		-349.1	-300.5	220.26	51.1
$\text{Cr}_2\text{O}_9^{--}$		aq		-305.			
$\text{Cr}_7\text{H}_2$		c		-3.7			
$[\text{Cr}(5\text{H}_2\text{O})(\text{OH})]^{++}$		aq		-456.7			
$\text{HCrO}_4^-$	std. state, hyp. m=1	aq		-212.8	-177.5	130.10	16.5
$[\text{Cr}(4\text{H}_2\text{O})(\text{OH})_2]^+$		aq		-448.9			



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

Washington, D.C.

National Bureau of Standards

Table 49-4. Chromium (at. no., 24; at. wt., 52.01) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)									
June 30, 1949									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole				
$H_2CrO_4$	std. state, hyp. m=1 in 80 $H_2O$ $\infty H_2O$		aq		-213.5	-178.5	130.84	17.5	
			aq		-209.2				
			aq		-213.5				
$Cr(OH)_3$			c		-247.1				
			c		-507.5				
			c		-579.8				
$HCr_2O_9^{--}$			aq		-322.				
$CrF_2$			c		-181.0				
$CrF_3$			c		-265.2				
$[Cr(6H_2O)]F_3$			aq		-704.3				
$H_3[Cr(6H_2O)]F_6$			aq		-931.9				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 49-5. Chromium (at. no., 24; at. wt., 52.01)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Substance		State	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description							
$[\text{Cr}(\text{SH}_2\text{O})\text{Cl}]^{++}$	blue green	aq		-477.7				
$\text{CrCl}_2$		c		-94.56	-85.15	62.414	27.4	16.87
$\text{CrCl}_2 \cdot 2\text{H}_2\text{O}$	light green	aq		-113.2				
$\text{CrCl}_2 \cdot 3\text{H}_2\text{O}$	pale blue	c		-237.				
$\text{CrCl}_2 \cdot 4\text{H}_2\text{O}$	dark blue	c		-309.0				
		c		-384.5				
$[\text{Cr}(\text{4H}_2\text{O})\text{Cl}_2]^+$	green	aq		-404.4				
$\text{CrCl}_3$		c		-134.6	-118.0	86.492	30.0	21.53
$[\text{Cr}(\text{4H}_2\text{O})\text{Cl}_2]\text{Cl}$	green	c		-436.0				
	green	aq		-444.4				
$[\text{Cr}(\text{4H}_2\text{O})\text{Cl}_2]\text{Cl} \cdot 2\text{H}_2\text{O}$	green	c		-581.1				
$[\text{Cr}(\text{4H}_2\text{O})\text{Cl}_2]\text{Cl} \cdot 6\text{H}_2\text{O}$	green	c		-854.3				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 49-6. Chromium (at. no., 24; at. wt., 52.01)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Substance		Description	State	$\Delta H^{\circ}$	$\Delta F^{\circ}$	$\log_{10} f$	$S^{\circ}$	$C_p^{\circ}$
Formula	At 298.16°K (25°C)							
	At 0°K kcal/mole			kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
$[\text{Cr}(\text{5H}_2\text{O})\text{Cl}]\text{Cl}_2$ $[\text{Cr}(\text{6H}_2\text{O})]\text{Cl}_3$	blue green		aq		-517.7			
	violet		c		-579.0			
	violet		aq		-591.0			
$\text{CrCl}_4$ $\text{CrO}_2\text{Cl}_2$			g		-104.			
			liq		-135.7			
$[\text{Cr}(\text{4H}_2\text{O})(\text{OH})_2]\text{Cl}$ $[\text{Cr}(\text{5H}_2\text{O})(\text{OH})]\text{Cl}_2$			aq		-488.9			
			aq		-536.7			
$[\text{Cr}(\text{4H}_2\text{O})\text{Br}_2]\text{Br}$ $[\text{Cr}(\text{4H}_2\text{O})\text{Br}_2]\text{Br}\cdot 2\text{H}_2\text{O}$ $[\text{Cr}(\text{6H}_2\text{O})]\text{Br}_3$	green		aq		-409.5			
	green		c		-545.5			
	violet		c		-543.3			
	violet		aq		-557.7			
$\text{CrI}_2$			c		-54.2			
			aq		-59.9			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 49-7. Chromium (at. no., 24; at. wt., 52.01)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description		At 0°K kcal/mole	kcal/mole			At 298.16°K (25°C) kcal/mole	cal/deg mole
$[\text{Cr}_2(6\text{H}_2\text{O})(\text{SO}_4)_3]$	green	aq		-1164.7				
$[\text{Cr}_2(6\text{H}_2\text{O})(\text{SO}_4)_3 \cdot 2\text{H}_2\text{O}]$	green	c		-1287.7				
$[\text{Cr}_2(8\text{H}_2\text{O})(\text{SO}_4)_2](\text{SO}_4)$		aq		-1304.9				
$[\text{Cr}_2(10\text{H}_2\text{O})(\text{SO}_4)](\text{SO}_4)_2$		aq		-1447.4				
$[\text{Cr}_2(12\text{H}_2\text{O})](\text{SO}_4)_3$	violet	aq		-1589.5				
$[\text{Cr}_2(12\text{H}_2\text{O})](\text{SO}_4)_3 \cdot 2\text{H}_2\text{O}$	violet	c		-1716.0				
$[\text{Cr}_2(12\text{H}_2\text{O})](\text{SO}_4)_3 \cdot 3\text{H}_2\text{O}$	violet	c		-1786.2				
$[\text{Cr}_2(12\text{H}_2\text{O})](\text{SO}_4)_3 \cdot 4\text{H}_2\text{O}$		c		-1855.3				
$[\text{Cr}_2(12\text{H}_2\text{O})](\text{SO}_4)_3 \cdot 5\text{H}_2\text{O}$		c		-1924.8				
$[\text{Cr}_2(12\text{H}_2\text{O})](\text{SO}_4)_3 \cdot 6\text{H}_2\text{O}$		c		-1993.2				
$\text{CrN}$		c		-29.8				
$\text{Cr}_2\text{N}$		c		-23.4				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 49-3. Chromium (at. no., 24; at. wt. 52.01) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1949									
Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
$(\text{NH}_4)_2\text{CrO}_4$		in 300 $\text{H}_2\text{O}$	c		-275.2				
$(\text{NH}_4)_2\text{Cr}_2\text{O}_7$			aq		-269.4				
			c		-425.5				
			aq		-412.6				
CrP			c						12.4
CrP <sub>2</sub>			c						19.8
CrSb			c						
CrSb <sub>2</sub>			c						
Cr <sub>3</sub> C <sub>2</sub>			c		-21.0	-21.2	15.54	20.4	23.38
Cr <sub>4</sub> C			c		-16.4	-16.8	12.31	25.3	25.89
Cr <sub>7</sub> C <sub>3</sub>			c		-42.5	-43.8	32.10	48.0	49.92

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 49-9. Chromium (at. no. 24; at. wt., 52.01)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)			cal/deg mole
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	
PbCrO <sub>4</sub>		c		-217.7			
2CrI <sub>2</sub> ·PbI <sub>2</sub>		c		-140.2			
2CrI <sub>2</sub> ·PbI <sub>2</sub> ·3H <sub>2</sub> O		c		-368.9			
Ag <sub>2</sub> CrO <sub>4</sub>		c		-170.15	108.900	51.8	34.00
FeCr <sub>2</sub> O <sub>4</sub>		c		-341.9	232.87	34.9	31.9

## Washington, D.C.

## National Bureau of Standards

Table 50-1. Molybdenum (at. no., 42; at. wt., 95.95)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Substance		State	Description	At 298.16°K (25°C)				
Formula	$\Delta H^{\circ}$			$\Delta F^{\circ}$	$\log_{10} f^{\circ}$	$S^{\circ}$	$C_p^{\circ}$	
	kcal/mole			kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
Mo		g		155.5	144.2	-105.70	43.462	4.9680
Mo <sup>+</sup>		c		0.000	0.000	0.0000	6.83	5.61
Mo <sup>++</sup>		g		327.23				
		g						
MoO <sub>2</sub>		c		-130.				
MoO <sub>3</sub>		c		-180.33	-161.95	118.707	18.68	17.59
		aq		-188.1				
MoO <sub>4</sub>		aq		-173.5				
MoO <sub>4</sub> <sup>--</sup>		aq		-254.3				
MoO <sub>5</sub>		aq		-155.1				
H <sub>2</sub> MoO <sub>4</sub>	MoO <sub>3</sub> ·H <sub>2</sub> O white	c		-256.9				
		aq		-256.4				
H <sub>2</sub> MoO <sub>4</sub> ·H <sub>2</sub> O	MoO <sub>3</sub> ·2H <sub>2</sub> O yellow	c		-331.4				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 50-2. Molybdenum (at. no., 42; at. wt., 95.95) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1949									
Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\Delta F_f^\circ$	$\log_{10} f$	$S^\circ$	$G_p^\circ$	
Formula	Description	State	At 0°K	At 298.16°K (25°C)					
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole	
MoCl <sub>2</sub>		c		-44.					
MoCl <sub>3</sub>		c		-65.					
MoCl <sub>4</sub>		c		-79.					
MoCl <sub>5</sub>		c		-90.8					
MoCl <sub>6</sub>		c		-90.					
MoBr <sub>2</sub>		c		-29.					
MoBr <sub>3</sub>		c		-41.					
MoBr <sub>4</sub>		c		-45.					
MoBr <sub>5</sub>		c		-51.					
MoI <sub>2</sub>		c		-12.					
MoI <sub>3</sub>		c		-15.					
MoI <sub>4</sub>		c		-18.					
MoI <sub>5</sub>		c		-18.					



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 50-3. Molybdenum (at. no., 42; at. wt., 95.95) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1949									
Substance		State	Description	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				
MoS <sub>2</sub>		c			-55.5		39.43	15.1	15.17
MoS <sub>3</sub>		c			-61.2				
MoP		c							
MoP <sub>2</sub>		c							
Mo <sub>2</sub> C		c			4.3	2.9	-2.13	19.7	
PbMoO <sub>4</sub>		c			-265.8				
CuMoO <sub>4</sub>		c			-226.6				
FeMoO <sub>4</sub>		c			-257.5				
Fe <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>		c			-708.2				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 51-1. Tungsten (at. no., 74; at. wt., 183.92)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1949

Substance		$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)		
			kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
W		g				
W <sup>+</sup>		c	0.000	201.6	-140.44	5.0903
W <sup>++</sup>		g		0.000	0.0000	5.97
		g		387.2		
WO <sub>2</sub>		c		-136.3		
WO <sub>3</sub>	yellow	c		-200.84	-182.47	19.90
WO <sub>4</sub> <sup>--</sup>		aq		-266.6		
W <sub>2</sub> O <sub>5</sub>		c		-337.9		
H <sub>2</sub> WO <sub>4</sub>	WO <sub>3</sub> · H <sub>2</sub> O	c		-279.6		

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 51-2. Tungsten (at. no., 74; at. wt., 183.92)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Substance		State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$
Formula	Description							
WC1 <sub>2</sub>		c		-38.				
WC1 <sub>4</sub>		c		-71.				
WC1 <sub>5</sub>		c		-84.				
WC1 <sub>6</sub>		c		-98.7				
WBr <sub>2</sub>		c		-19.				
WBr <sub>4</sub>		c		-35.				
WBr <sub>5</sub>		c		-42.				
WBr <sub>6</sub>		c		-44.				
WI <sub>2</sub>		c		-1.				
WI <sub>4</sub>		c		0.				
WI <sub>5</sub>		c		27.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 51-3. Tungsten (at. no., 74; at. wt., 183.92)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
WS <sub>2</sub>			c		-46.3	-46.2	33.86	23.	
WP			c						
WP <sub>2</sub>			c						
WC			c		-4.09				
CuWO <sub>4</sub>			c		-250.0				
CuWO <sub>4</sub> ·2H <sub>2</sub> O			c		-393.6				
FeWO <sub>4</sub>			c		-274.1				
FeWO <sub>4</sub> ·3H <sub>2</sub> O			c		-486.7				
Fe <sub>2</sub> (WO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O			c		-1352.8				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES I

Washington, D.C.

Table 52-1. Vanadium (at. no., 23; at. wt., 50.95) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1949											
Substance		State	Description	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
Formula	Description			At 0°K						At 298.16°K (25°C)	
				kcal/mole	kcal/mole					kcal/mole	cal/deg mole
V		g		119.2	120.	109.	-79.90	43.546	6.2166		
		c		0.000	0.000	0.000	0.0000	7.05	5.85		
V +		g		274.17	275.65						
V ++		g		601.6	604.8						
V +++		g		1213.5	1217.9						
V ++++		g		2331.2	2337.1						
V ++++		g		3817.6	3825.0						
V ++++		g		6776.3	6785.2						
V ++++		g		10235.	10245.						
V ++++		g		14220.	14232.						
V ++++		g									
VO		g		51.	52.						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 52-2. Vanadium (at. no., 23; at. wt., 50.95)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949 (Corrected)

Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$ (25°C)	$S^\circ$	$C_p^\circ$
Formula	At 0°K								
	kcal/mole			kcal/mole	kcal/mole	cal/deg mole	cal/deg mole		
$\text{VO}_4^-$			aq		-210.9				
$\text{VO}_5^-$			aq		-194.9				
$\text{V}_2\text{O}_2$			c		-200.				
$\text{V}_2\text{O}_3$			c	-288.86	-290.	-271.	198.6	23.58	24.83
$\text{V}_2\text{O}_4$			c, II	-341.94	-344.	-318.	233.1	24.65	28.30
$\text{V}_2\text{O}_4 \cdot 2\text{H}_2\text{O}$	red		c						
$\text{V}_2\text{O}_5$			c	-370.71	-373.	-344.	252.1	31.3	31.00
$\text{V}_3\text{O}_5$			aq		-673.5				
$\text{HV}_6\text{O}_{17}$		std. state, hyp. m = 1	aq			-1127.3	826.295		
$\text{H}_2\text{V}_6\text{O}_{17}$		std. state, hyp. m = 1	aq			-1130.	828.27		
$\text{VF}_3$			aq						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES Y

National Bureau of Standards

Washington, D.C.

Table 52-3. Vanadium (at. no., 23; at. wt., 50.95)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Substance			State	Description	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	At 0°K				At 298.16°K (25°C)					
	kcal/mole	kcal/mole			kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole	
VCl <sub>2</sub>	c			-106.	-97.	71.1	23.2	17.26		
VCl <sub>3</sub>	c			-137.	-120.	67.96	31.3	22.27		
VCl <sub>4</sub>	liq			-138.						
VOCl <sub>3</sub>	c			-172.						
VOSO <sub>4</sub>	c			-312.5						
VN	c			-41.	-35.	25.7	8.91	9.08		
NH <sub>4</sub> VO <sub>3</sub>	c			-251.2	-211.8	155.25	33.6	30.91		
VC	c						6.77	7.97		

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

Washington, D.C.

National Bureau of Standards

Table 53-1. Columbium (at. no., 41; at. wt., 92.91) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1949									
Formula	Substance	Description	State	At 298.16°K (25°C)					
				$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Cb			g	0.000	184.5	173.7	-127.32	44.492	7.2082
Cb <sub>2</sub> O <sub>4</sub>			c	0.000	0.000	0.000	0.0000	8.3	30.3
Cb <sub>2</sub> O <sub>5</sub>			c	0.000	-387.8	0.000	0.0000	8.3	30.3



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 54-1. Tantalum (at. no., 73; at. wt., 180.88)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1949 (Corrected)

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	At 298.16°K (25°C)			cal/deg mole
				kcal/mole	kcal/mole	kcal/mole	cal/deg mole	
Ta			g	184.9	185.	-128.3	44.244	5.0054
				0.000	0.000	0.0000	9.9	6.05
Ta <sub>2</sub> O <sub>5</sub>			c		-499.9	344.94	34.2	32.30
TaN			c					9.7
TaP			c		-58.2			
TaP <sub>2</sub>			c					
TaC			c				10.1	8.79

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 55-1. Titanium (at. no., 22; at. wt., 47.90)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1949; March 31, 1950

Substance		State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description							
Ti		g	0.000	112.	101.	-74.0	43.069	5.8385
		c, II		0.000	0.000	0.0000	7.24	6.010
Ti *		g		271.09				
Ti **		g		587.0				
Ti ***		g		1225.0				
Ti ****		g		2223.5				
Ti *****		g		4527.4				
Ti +++++		g		7290.				
Ti ++++++		g		10538.				
Ti +++++++		g						
TiO		g		43.			8.31	9.55
		c, II						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 55-2. Titanium (at. no., 22; at. wt., 47.90)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1949; March 31, 1950

Substance		State	Description	At 298.16°K (25°C)				
Formula	$\Delta H_f^\circ$			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
	At 0°K kcal/mole			kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
TiO <sub>2</sub>	rutile, III anatase, III hydrated precipitate	c		-218.0	-203.8	149.38	12.01	13.16
		c					11.93	13.22
Ti <sub>2</sub> O <sub>3</sub>		amorp		-207.				
Ti <sub>3</sub> O <sub>5</sub>		c, II		-367.	-346.	253.6	18.83	23.27
	c, II		-584.	-550.	403.1	30.92	37.00	
TiF <sub>2</sub>		c		-198.				
TiF <sub>3</sub>		c		-315.				
TiF <sub>4</sub>		c		-370.				
H <sub>2</sub> TiF <sub>6</sub>		aq		-555.1				
TiCl		g		122.				
TiCl <sub>2</sub>		c		-114.				
TiCl <sub>3</sub>		c		-165.				
TiCl <sub>4</sub>		g						22.88
	hydrolysed	liq		-179.3	-161.2	118.16	84.4	37.5
		aq		-236.5			60.4	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 55-3. Titanium (at. no., 22; at. wt., 47.90)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta P_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)				
			kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
TiBr <sub>2</sub>		c		-95.				
TiBr <sub>3</sub>		c		-132.				
TiBr <sub>4</sub>		c		-155.				
TiI <sub>2</sub>		c		-61.				
TiI <sub>3</sub>		c		-80.				
TiI <sub>4</sub>		c		-102.				
TiCl <sub>4</sub> ·H <sub>2</sub> S		c		-193.0				
TiCl <sub>4</sub> ·2H <sub>2</sub> S		c		-205.3				
TiBr <sub>4</sub> ·H <sub>2</sub> S		c		-168.4				
TiBr <sub>4</sub> ·2H <sub>2</sub> S		c		-180.4				
TiN		c		-73.0	-66.1	48.45	7.20	8.86

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 55-4. Titanium (at. no., 22; at. wt., 47.90) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1949									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	kcal/mole				
TiCl <sub>4</sub> ·PH <sub>3</sub>			c		-195.7				
TiCl <sub>4</sub> ·2PH <sub>3</sub>			c		-204.3				
TiBr <sub>4</sub> ·PH <sub>3</sub>			c		-168.8				
TiBr <sub>4</sub> ·2PH <sub>3</sub>			c		-178.4				
TiC			c		-54.		38.8	5.8	8.04
FeTiO <sub>3</sub>			c		-288.5		197.10	25.3	23.78

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

Washington, D.C.

Table 56-1. Zirconium (at. no., 40; at. wt., 91.22)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1949 (Corrected)

Substance		State	Description	At 298.16°K (25°C)				
Formula	$\Delta H_f^\circ$			$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
	kcal/mole			kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
Zr		g c, II		125. 0.000	115. 0.000	-84.29 0.0000	43.313 9.18	6.3624
Zr <sup>+</sup>		g		286.78				
Zr <sup>++</sup>		g		611.84				
Zr <sup>+++</sup>		g		1169.0				
Zr <sup>++++</sup>		g		1953.9				
Zr <sup>++++</sup>		g						
ZrO <sup>++</sup>		aq		-223.1				
ZrO <sub>2</sub>		c, III		-258.2	-244.4	179.14	12.03	
ZrO <sub>2</sub> ·5H <sub>2</sub> O		аморф		-625.3				
ZrO(OH) <sub>2</sub>		c		-338.0				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards Washington, D.C.

SERIES I

Table 56-2. Zirconium (at. no., 40; at. wt., 91.22)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
June 30, 1949

Substance		$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)		
			kcal/mole	$\Delta H_f^\circ$	$\Delta F_f^\circ$	cal/deg mole
				kcal/mole	kcal/mole	
Zr(OH) <sub>4</sub>		c		-411.2		
Zr(OH) <sub>4</sub> ·H <sub>2</sub> O		c		-482.9		
Zr(OH) <sub>4</sub> ·2H <sub>2</sub> O		c		-554.2		
ZrF <sub>2</sub>		c		-230.		
ZrF <sub>3</sub>		c		-350.		
ZrF <sub>4</sub>		c		-445.		
ZrCl <sub>2</sub>		c		-145.		
ZrCl <sub>3</sub>		c		-208.		
ZrCl <sub>4</sub>		c		-230.	-209.	44.5
ZrOCl <sub>2</sub>		aq		-303.1		
ZrOCl <sub>2</sub> ·2H <sub>2</sub> O		c		-423.7		
ZrOCl <sub>2</sub> ·3H <sub>2</sub> O		c		-533.2		
ZrOCl <sub>2</sub> ·6H <sub>2</sub> O		c		-712.3		
ZrOCl <sub>2</sub> ·8H <sub>2</sub> O		c		-852.7		

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

Washington, D.C.

Table 56-3. Zirconium (at. no., 40; at. wt., 91.22) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) June 30, 1949 (Corrected)									
Substance		$\Delta H_f^\circ$		$\Delta F_f^\circ$		$\log_{10} K_f$		$S^\circ$	
Formula	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$		$\log_{10} K_f$		$C_p^\circ$
			At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole	
ZrBr <sub>2</sub>		c		-120.					
ZrBr <sub>3</sub>		c		-174.					
ZrBr <sub>4</sub>		c		-192.					
ZrOBr <sub>2</sub>		aq		-280.9					
ZrOBr <sub>2</sub> ·3H <sub>2</sub> O		c		-511.0					
ZrOBr <sub>2</sub> ·8H <sub>2</sub> O		c		-829.5					
ZrI <sub>2</sub>		c		-90.					
ZrI <sub>3</sub>		c		-128.					
ZrI <sub>4</sub>		c		-130.					
ZrOSO <sub>4</sub>		aq		-486.3					
Zr(SO <sub>4</sub> ) <sub>2</sub>		c		-597.4					
Zr(SO <sub>4</sub> ) <sub>2</sub> ·H <sub>2</sub> O		c		-677.9					
Zr(SO <sub>4</sub> ) <sub>2</sub> ·4H <sub>2</sub> O		c		-693.1					26.7



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

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Table 56-4. Zirconium (at. no., 40; at. wt., 91.22) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°K) June 30, 1949									
Substance			State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°K)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description								
ZrN			c						
ZrO(NO <sub>3</sub> ) <sub>2</sub>			aq						
ZrO(NO <sub>3</sub> ) <sub>2</sub> ·2H <sub>2</sub> O			c		-82.2	-75.4	55.27	9.23	
ZrO(NO <sub>3</sub> ) <sub>2</sub> ·3H <sub>2</sub> O			c		-321.8				
ZrO(NO <sub>3</sub> ) <sub>2</sub> ·3H <sub>2</sub> O			c		-456.3				
ZrO(NO <sub>3</sub> ) <sub>2</sub> ·3H <sub>2</sub> O			c		-527.3				
ZrO(NO <sub>3</sub> ) <sub>2</sub> ·3H <sub>2</sub> O			c		-562.8				
ZrO(NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O			c		-737.6				
ZrC			c		-45.				
ZrSiO <sub>4</sub>			c					20.1	23.53

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards Washington, D.C.

Table 57-1. Hafnium (at. no., 72; at. wt., 178.6)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

June 30, 1949

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)			
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
Hf		g	0.000	0.000	0.0000	44.65	6.15
HfO <sub>2</sub>		c				13.1	
		c		-271.5			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 58-1. Boron (at. no., 5; at. wt., 10.82) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1949									
Substance		Description	State	$\Delta H_f^0$	$\Delta H_f^0$	$\Delta F_f^0$	$\log_{10} K_f$	$S^0$	$C_p^0$
Formula	At 298.16°K (25°C)								
	At 0°K kcal/mole			kcal/mole	kcal/mole		kcal/mole	cal/deg mole	
B			g	96.	97.2	86.7	-63.55	36.649	4.971
			c	0.000	0.000	0.000	0.0000	1.56	2.66
			amorp		0.4			1.56 + x <sup>a</sup>	2.859
B <sup>+</sup>			g	287.32	290.00				
B <sup>++</sup>			g	867.30	871.46				
B <sup>+++</sup>			g	1741.81	1747.45				
B <sup>++++</sup>			g	7721.70	7728.82				
B <sup>+++++</sup>			g	15565.6	15574.2				
B <sub>2</sub>			g	123.	124.5				

a x = entropy of randomness of arrangement of B [amorp] at 0°K

a. x = entropy of randomness of arrangement of B (amorp) at 0°K

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 58-2. Boron (at. no., 5; at. wt., 10.82)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

September 30, 1949

Substance		State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description							
B0		g	-6.	-5.3	-11.6	8.50	47.22	6.95
B0 <sub>2</sub> <sup>-</sup>		aq		-183.5				
B <sub>2</sub> O <sub>3</sub>		c		-302.0	-283.0	207.44	12.91	14.88
B <sub>4</sub> O <sub>7</sub> <sup>---</sup>		gls		-297.6	-280.4	205.53	18.8	14.6
BH		aq						
BH		g	73.	73.8	67.1	-49.18	39.62	6.95
B <sub>2</sub> H <sub>6</sub>	diborane	g		7.5	19.8	-14.51	55.66	13.48
B <sub>4</sub> H <sub>10</sub>	tetraborane	g						
B <sub>5</sub> H <sub>9</sub>	pentaborane	g		15.0	39.6	-29.03	65.88	19.
B <sub>5</sub> H <sub>9</sub>	pentaborane	liq		7.8	38.8	-28.44	44.16	35.8
B <sub>5</sub> H <sub>11</sub>		liq						
B <sub>10</sub> H <sub>14</sub>	decaborane	c		8.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES I

Washington, D.C.

Table 58-3. Boron (at. no., 5; at. wt., 10.82)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

September 30, 1949

Substance									
Formula	Description	State	$\Delta H_f^\circ$		$\Delta P_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 0°K						
			kcal/mole	kcal/mole					
HBO <sub>2</sub>		c		-186.9	-170.5	124.97	11.		
H <sub>2</sub> BO <sub>3</sub> <sup>-</sup>	std. state, hyp. m = 1	aq		-186.9					
H <sub>3</sub> BO <sub>3</sub>		aq		-251.8	-217.63	159.520	7.3		
	std. state, hyp. m = 1	c		-260.2	-230.2	168.73	21.41	19.61	
	in 50 H <sub>2</sub> O	aq		-255.2	-230.24	168.762	38.2		
	700 H <sub>2</sub> O	aq		-255.28					
	∞ H <sub>2</sub> O	aq		-255.21					
		aq		-255.2					
H <sub>2</sub> B <sub>4</sub> O <sub>7</sub>		c		-676.5					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 58-4. Boron (at. no., 5; at. wt., 10.82)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
September 30, 1949

Substance		State	Description	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K			At 298.16°K (25°C)					
				kcal/mole	kcal/mole			kcal/mole	cal/deg mole
BF	std. state, hyp. m = 1	g		-18.	-17.4				
BF <sub>3</sub>		g			-265.4		191.53	60.70	12.06
BF <sub>4</sub> <sup>-</sup>		aq				-289.8			
BH <sub>4</sub> <sup>-</sup>		aq				-365.	251.4	40.	
BH <sub>4</sub>		aq				-365.			
BCl		g		25.	25.6				5.59
BCl <sub>3</sub>		g			-94.5	-90.9	66.63	69.29	14.97
		liq				-100.0	66.41	50.0	
BBr		g			25.				
BBr <sub>3</sub>		g			-44.6	-51.0	37.38	77.49	16.25
		liq				-52.8	38.41	54.7	
B <sub>2</sub> H <sub>5</sub> Br		g							

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 58-5. Boron (at. no., 5; at. wt., 10.82)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°K)  
September 30, 1949

Substance									
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 0°K						
			kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole	
B <sub>2</sub> S <sub>3</sub>		c		-57.0					
BN		g	90.	90.6					
B <sub>3</sub> N <sub>3</sub> H <sub>6</sub>	hexahydro-s-triazatriborine, borazole	c		-32.1	-27.2	19.94	8.	5.9	
		g					73.7	23.3	
NH <sub>4</sub> BO <sub>2</sub>		aq		-215.2					
NH <sub>4</sub> BO <sub>3</sub>		aq		-193.7					
NH <sub>4</sub> BO <sub>3</sub> ·H <sub>2</sub> O		c		-270.8					
(NH <sub>4</sub> ) <sub>2</sub> HBO <sub>3</sub>		aq		-305.5					
B <sub>4</sub> C		c					6.47	12.55	
B(CH <sub>3</sub> ) <sub>3</sub>		liq		-31.4					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 59-1. Aluminum (at. no., 13; at. wt., 26.97)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
September 30, 1949

Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 298.16°K (25°C)								
	kcal/mole			kcal/mole	kcal/mole			cal/deg mole	cal/deg mole
Al			g	74.4	75.0	65.3	-47.86	39.303	5.112
Al +			c	0.000	0.000	0.000	0.0000	6.769	5.817
Al ++			g	212.406	214.487				
Al ***			q	646.500	650.062				
			q	1302.40	1307.44				
		std. state, hyp. m=1	aq		-125.4	-115.0	84.29	-74.9	
Al ++++			g	4068.82	4075.34				
Al +++++			q	7615.09	7623.09				
Al ++++++			q	12006.6	12016.1				
Al ++++++			g	17585.9	17596.9				
Al ++++++			g	24162.	24174.				
Al ++++++			g	31775.	31789.				
Al ++++++			q	40966.	40982.				
Al ++++++			g	51156.	51173.				



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 59-2. Aluminum (at. no., 13; at. wt., 26.97) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1949									
Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\Delta P_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	Description	State	At 298.16°K (25°C)						
			At 0°K kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole	
AlO		g	45.	45.3					
AlO <sub>2</sub> <sup>-</sup>		aq		-218.6					
Al <sub>2</sub> O <sub>3</sub>	α, corundum	c	-396.19	-399.09	-376.77	276.167	12.186	18.88	
	γ	c		-384.84					
Al <sub>2</sub> O <sub>3</sub> ·H <sub>2</sub> O		c		-471.	-435.	318.8	23.15	31.37	
Al <sub>2</sub> O <sub>3</sub> ·3H <sub>2</sub> O	hydrargillite	c		-613.7	-547.9	401.60	33.51	44.49	
AlH		g	59.	59.2					
Al(OH) <sub>3</sub>		amorp		-304.2					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 59-3. Aluminum [at. no., 13; at. wt., 26.97]  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
September 30, 1949

Formula	Substance	State	Description	At 0°K					At 298.16°K (25°C)		
				$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	$\Delta H_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$
AlF		g		51.	51.4						
AlF <sub>3</sub>		c			-311.	215.5	23.		-294.		
AlF <sub>3</sub> ·½H <sub>2</sub> O		c			-361.4						
AlF <sub>3</sub> ·3H <sub>2</sub> O		c			-357.4	244.52	28.		-333.6		
H <sub>3</sub> AlF <sub>6</sub>		c			-549.1	359.46	50.		-490.4		
		aq			-597.2						
AlCl		g		-5.	-4.9						
AlCl <sub>3</sub>		c			-166.2	111.56	40.		-152.2		21.3
		aq			-245.5						
AlCl <sub>3</sub> ·6H <sub>2</sub> O		c			-641.1	397.57	90.		-542.4		
Al <sub>2</sub> Cl <sub>6</sub>		g			-303.6						
			in 600 H <sub>2</sub> O								

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 59-4. Aluminum (at. no., 13; at. wt., 26.97) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1949										
Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	At 0°K			At 298.16°K (25°C)			cal/deg mole	cal/deg mole		
	kcal/mole			kcal/mole	kcal/mole	kcal/mole			kcal/mole	cal/deg mole
AlBr			g		1.8					
AlBr <sub>3</sub>			c		-125.8	-120.7	88.47	44.	24.5	
			aq		-211.9					
AlI			g	31.	30.6					
AlI <sub>3</sub>			c		-75.2	-75.0	54.97	48.		
			aq		-165.8					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Table 59-5. Aluminum (at. no., 13; at. wt., 26.97)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

September 30, 1949

Substance		State	Description	$\Delta F^{\circ}$					$\log_{10} K_f$	$S^{\circ}$	$C_p^{\circ}$
Formula	$\Delta F^{\circ}$			$\Delta F^{\circ}$							
	At 0°K kcal/mole			kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole				
$Al_2S_3$		c			-121.6	-117.7	86.27	23.			
$Al_2(SO_4)_3$		c			-820.98	-738.99	541.670	57.2		62.00	
$Al_2(SO_4)_3 \cdot 6H_2O$		aq			-897.1						
$Al_2(SO_4)_3 \cdot 14H_2O$		c			-1268.14	-1105.14	810.054	112.1		117.8	
		c			-2118.5						
$AlCl_3 \cdot \frac{1}{2}SO_2$		c			-210.8						
$AlCl_3 \cdot SO_2$		c			-251.6						
$AlCl_3 \cdot \frac{1}{2}H_2S$		c			-173.3						
$AlCl_3 \cdot H_2S$		c			-180.2						
$AlBr_3 \cdot H_2S$		c			-140.3						
$AlI_3 \cdot 2H_2S$		c			-103.1						
$AlI_3 \cdot 4H_2S$		c			-125.7						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Washington, D.C.

Table 59-6. Aluminum (at. no., 13; at. wt., 26.97) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1949									
Substance		Description	State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula									
AlN			c		-57.7	-50.1	36.72	5.	7.67
Al(NO <sub>3</sub> ) <sub>3</sub>			aq		-273.65				
Al(NO <sub>3</sub> ) <sub>3</sub> ·6H <sub>2</sub> O			c		-680.65	-525.82	365.419	111.8	103.5
Al(NO <sub>3</sub> ) <sub>3</sub> ·9H <sub>2</sub> O			c		-897.34	-700.2	513.24	136.	
AlF <sub>3</sub> ·2NH <sub>4</sub> F· $\frac{3}{2}$ H <sub>2</sub> O			c		-673.7				
AlCl <sub>3</sub> ·NH <sub>3</sub>			c		-211.3				
AlCl <sub>3</sub> ·NH <sub>4</sub> Cl			c		-255.0				
AlCl <sub>3</sub> ·3NH <sub>3</sub>			c		-283.1				
AlCl <sub>3</sub> ·5NH <sub>3</sub>			c		-341.5				
AlCl <sub>3</sub> ·6NH <sub>3</sub>			c		-366.5				94.2
AlCl <sub>3</sub> ·7NH <sub>3</sub>			c		-367.7				
AlCl <sub>3</sub> ·NH <sub>4</sub> Cl·6NH <sub>3</sub>			c		-444.2				
AlCl <sub>3</sub> ·14NH <sub>3</sub>			c		-519.6				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 59-7. Aluminum (at. no., 13; at. wt., 26.97) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1949									
Formula	Substance	Description	State	At 298.16°K (25°C)					$C_p^0$
				$\Delta H_f^0$ kcal/mole	$\Delta H_f^0$ kcal/mole	$\Delta F_f^0$ kcal/mole	$\log_{10} K_f$	$S^0$ cal/deg mole	
$AlBr_3 \cdot NH_3$			c		-178.1				
$AlBr_3 \cdot 3NH_3$			c		-254.8				
$AlBr_3 \cdot 5NH_3$			c		-319.8				
$AlBr_3 \cdot 6NH_3$			c		-347.2				
$AlBr_3 \cdot 7NH_3$			c		-368.9				
$AlBr_3 \cdot 9NH_3$			c		-407.8				
$AlBr_3 \cdot 14NH_3$			c		-501.6				
$AlI_3 \cdot NH_3$			c		-119.9				
$AlI_3 \cdot 3NH_3$			c		-208.3				
$AlI_3 \cdot 5NH_3$			c		-291.8				
$AlI_3 \cdot 6NH_3$			c		-317.2				
$AlI_3 \cdot 7NH_3$			c		-339.9				
$AlI_3 \cdot 9NH_3$			c		-379.2				
$AlI_3 \cdot 13NH_3$			c		-454.4				
$AlI_3 \cdot 20NH_3$			c		-585.4				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 59-8. Aluminum (at. no., 13; at. wt., 26.97)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
September 30, 1949

Substance		State	Description	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K		kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
				kcal/mole	kcal/mole				
$(\text{NH}_4)_2\text{Al}(\text{SO}_4)_2$		c							
	in 1500 H <sub>2</sub> O	aq			-561.24	-485.95	356.195	51.7	54.12
	2000 H <sub>2</sub> O	aq			-591.00				
	2500 H <sub>2</sub> O	aq			-591.23				
	3000 H <sub>2</sub> O	aq			-591.46				
	3500 H <sub>2</sub> O	aq			-591.60				
$(\text{NH}_4)_2\text{Al}(\text{SO}_4)_2 \cdot 12\text{H}_2\text{O}$		aq			-591.74				
	ammonium alunite	c			-1419.40	-1179.02	864.207	166.6	163.3
	ammonium basic alum	c			-2406.45				
		c			-2758.65				
$\text{AlCl}_3 \cdot \text{PH}_3$		c			-173.9				
$\text{AlBr}_3 \cdot \text{PH}_3$		c			-137.9				
$\text{AlI}_3 \cdot \text{PH}_3$		c			-91.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 59-9. Aluminum (at. no., 13; at. wt., 26.97)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
September 30, 1949

Formula	Substance	State	Description	$\Delta H_f^\circ$			$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole			
$Al_4C_3$		c			-30.9	-29.0	21.26	25.	33.6
$Al(CH_3)_3$		liq			-26.9				
$Al_2SiO_5$	andalusite	c			-642.2	-607.8	445.51	25.0	32.14
	disthene	c			-642.7	-607.0	444.92	20.7	29.0
	sillimanite	c			-648.9	-615.0	450.79	27.0	30.45
$2AlI_3 \cdot 3PbI_2$		c			-234.				
$2AlI_3 \cdot 3PbI_2 \cdot 10H_2O$		c			-1122.				
$AlCu$		c			-10.				
$AlCu_2$		c			-16.				
$Al_2Cu$		c			-10.				
$AlCl_3 \cdot AgCl$		c			-198.9				



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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## SERIES I

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Table 59-10. Aluminum (at. no., 13; at. wt., 26.97) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) September 30, 1949									
Substance			State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	At 0°K							
		kcal/mole		kcal/mole	kcal/mole	cal/deg mole	cal/deg mole		
AlNi			c		-34.				
AlNi <sub>3</sub>			c		-38.				
Al <sub>2</sub> Ni			c		-38.				
Al <sub>3</sub> Ni			c		-38.				
AlCo			c		-26.				
Al <sub>4</sub> Co			c		-38.				
Al <sub>5</sub> Co <sub>2</sub>			c		-70.				
AlFe			c		-12.				
Al <sub>2</sub> Fe			c		-20.				
Al <sub>3</sub> Fe			c		-27.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 60-1. Scandium (at. no., 21; at. wt., 45.10)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1949

Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description		At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				
Sc		g	0.000	93.	0.000	0.0000	41.76	5.282
Sc +		g		245.75				
Sc ++		g		544.51				
Sc +++		g		1116.81				
Sc ++++		aq		-146.8	-143.7	105.33		
Sc +++++		g		2822.8				
Sc +++++		g		4942.				
Sc ++++++		g		7505.				
Sc2O3	std. state, hyp. m=1	c						21.

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 60-2. Scandium (at. no., 21; at. wt., 45.10)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1949

Substance		State	At 298.16°K (25°C)			
Formula	Description		$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole
ScCl <sub>3</sub>		c				
ScBr <sub>3</sub>	std. state, hyp. m=1	aq	-220.8	-237.7	174.23	
		c	-268.9			
	std. state, hyp. m=1	aq	-179.4	-217.4	159.35	
			-235.5			
Sc <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c				62.
ScF <sub>3</sub> ·3NH <sub>4</sub> F		c				90.0
Sc <sub>2</sub> (C <sub>2</sub> O <sub>4</sub> ) <sub>3</sub>	scandium oxalate	c				114.1
Sc(CH <sub>3</sub> O <sub>2</sub> ) <sub>3</sub>	scandium formate	c				53.8

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Substance		Table 61-1. Yttrium (at. no., 39; at. wt., 88.92) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1949				
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$C_p^\circ$
			At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole	cal/deg mole
Y		g	0.000	0.000	0.0000	6.181
Y <sup>+</sup>		c	103.			
Y <sup>++</sup>		g	255.04			
Y <sup>+++</sup>		g	542.37			
Y <sup>++++</sup>		g	1016.33			
Y <sup>++++</sup>	std. state, hyp. m = 1	aq	-168.0	-164.1	120.28	
Y <sub>2</sub> O <sub>3</sub>		g				23.
Y(OH) <sub>3</sub>		c				
Y(OH) <sub>3</sub>		c	-337.6	-308.3	225.98	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 61-2. Yttrium (at. no., 39; at. wt., 88.92)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1949

Substance		State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description							
YCl <sub>3</sub>	Y	c						
	std. state, hyp. m = 1	aq		-234.8 -288.1	-258.1	189.18		
YI <sub>3</sub>	β	c		-143.2				
	std. state, hyp. m = 1	aq		-208.1	-201.1	147.40		
Y <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c						61.
	std. state, hyp. m = 1	aq		-986.7	-860.2	630.51		
Y <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O		c		-1512.6	-1327.3	972.89		138.
Y(NO <sub>3</sub> ) <sub>3</sub>		c						75.7
Y <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>		c						105.

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SERIES I

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Washington, D.C.

Table 62-1. Lutetium (at. no., 71; at. wt., 174.99)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1949

Substance		Formula	Description	State	$\Delta F^\circ$		$\Delta F^\circ$	$\log_{10} f^\circ$	$S^\circ$	$C_p^\circ$
At 0°K					At 298.16°K (25°C)		kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole
kcal/mole	kcal/mole				kcal/mole	kcal/mole				
Lu				g	0.000	87.	0.000	0.0000	44.14	4.988
				c						
Lu <sup>+</sup>				g		203.8				
Lu <sup>++</sup>				g						
Lu <sup>+++</sup>			std. state, hyp. m = 1	aq		-160.1	-155.0	113.61		
LuCl <sub>3</sub>			γ	c		-227.9				
			std. state, hyp. m = 1	aq		-280.2	-249.0	182.51		
LuI <sub>3</sub>			β	c		-133.2				
			std. state, hyp. m = 1	aq		-200.2	-192.0	140.73		
Lu <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>			std. state, hyp. m = 1	aq		-970.9	-842.0	617.17		
Lu <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O				c			-1308.1	958.82		

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 63-1. Ytterbium (at. no., 70; at. wt., 173.04)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1949

Substance		State	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description							
Yb		g	0.000	87.	0.000	0.0000	41.30	4.968
Yb <sup>+</sup>		c		0.000				
Yb <sup>++</sup>		g		231.5				
Yb <sup>+++</sup>		g		511.				
	std. state, hyp. m = 1	aq			-129.0	94.56		
	std. state, hyp. m = 1	g			-155.5	113.98		
YbCl <sub>3</sub>		aq		-160.6				
	Y	c		-228.7				
	std. state, hyp. m = 1	aq		-280.7	-249.5	182.88		
Yb <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		aq		-971.9	-843.0	617.91		
Yb <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O		c			-1308.8	959.33		

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

SERIES I

Washington, D.C.

Table 64-1. Thulium (at. no., 69; at. wt., 169.4) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1949									
Substance		$\Delta H_f^\circ$		$\Delta F_f^\circ$		$\log_{10} K_f$		$S^\circ$	
Formula	Description	State	At 0°K		At 298.16°K (25°C)		cal/deg mole	cal/deg mole	$C_p^\circ$
			kcal/mole	kcal/mole	kcal/mole				
$Tm$		g	0.000	0.000	0.000		45.2		
$Tm^{++}$		c							
$Tm^{+++}$		g							
	std. state, hyp. m=1	aq		-161.3	-156.5	114.71			
$TmCl_3$	$\gamma$	c		-229.5					
	std. state, hyp. m=1	aq		-281.4	-250.5	163.61			
$TmI_3$	$\beta$	c		-137.8					
	std. state, hyp. m=1	aq		-201.4	-193.5	141.83			



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Washington, D.C.

Table 65-1. Erbium (at. no., 68; at. wt., 167.2) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1949 (Corrected)									
Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 298.16°K (25°C)					
				At 0°K kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
Er			g						
Er <sup>+</sup>			c	0.000	0.000	0.000	0.000		
Er <sup>+++</sup>			g						
		std. state, hyp. m = 1	aq		-162.3	-157.5	115.45		
Er <sub>2</sub> O <sub>3</sub>			c						25.
Er(OH) <sub>3</sub>			c		-340.5				
ErCl <sub>3</sub>			c		-231.8				
	γ		aq		-282.4	-251.5	184.35		
ErI <sub>3</sub>	β		c		-140.0				
		std. state, hyp. m = 1	aq		-202.4	-194.5	142.57		
Er <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>			c						65.
		std. state, hyp. m = 1	aq		-975.3	-847.0	620.84		
Er <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O			c			-1313.9	963.07		139.
Er(C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>3</sub>			aq		-512.8				
Er(C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>3</sub> ·4H <sub>2</sub> O		erbium acetate	c		-785.6				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 66-1. Holmium (at. no., 67; at. wt., 164.94)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1949

Formula	Substance	Description	State	At 0°K					At 298.16°K (25°C)				
				kcal/mole					kcal/mole				
Ho			g	0.000	0.000	0.000	0.000	0.0000					
Ho <sup>+</sup>			g										
Ho <sup>+++</sup>			aq										
HoCl <sub>3</sub>		std. state, hyp. m=1	c										
HoI <sub>3</sub>		std. state, hyp. m=1	aq										
Ho <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		std. state, hyp. m=1	aq										
Ho <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O			c										

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 67-1. Dysprosium (at. no., 66; at. wt., 162.46)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1949

Substance		$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K		At 298.16°K (25°C)		
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
Dy		g	0.000	87.	0.000		
Dy <sup>+</sup>		c		0.000			
Dy <sup>++</sup>		g		245.8			
Dy <sup>+++</sup>		aq		-166.0			
	std. state, hyp. m=1						
Dy(DH) <sub>3</sub>		c					
DyCl <sub>3</sub>		c					
	β			-237.8			
	γ			-234.8			
	std. state, hyp. m=1						
	β	aq		-286.1			
DyI <sub>3</sub>		c		-144.5			
	std. state, hyp. m=1						
		aq		-206.1			
	std. state, hyp. m=1						
Dy <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		aq		-982.7			
Dy <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O		c					
				-854.4			
				-1322.0			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 68-1. Terbium (at. no., 65; at. wt., 159.2)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1949

Formula	Substance	State	At 0°K					At 298.16°K (25°C)		
			$\Delta H_f^\circ$ kcal/mole		$\Delta H_f^\circ$ kcal/mole		$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole	
Tb		g			87.					
		c	0.000		0.000	0.000	0.0000			
Tb <sup>+</sup>		g			244.0					
Tb <sup>++</sup>		g								
Tb <sup>+++</sup>	std. state, hyp. m=1	aq			-168.4	-163.9	120.14			
TbCl <sub>3</sub>	$\beta$	c			-241.6					
	std. state, hyp. m=1	aq			-288.5	-257.9	189.04			
Tb <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>	std. state, hyp. m=1	aq			-987.5	-859.8	630.22			
Tb <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O		c				-1328.2	973.55			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES I

Washington, D.C.

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)		Table 69-1. Gadolinium (at. no., 64; at. wt., 156.9)					
		December 31, 1949					
Formula	Substance	Description	State	At 298.16°K (25°C)			
				$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$
				kcal/mole	kcal/mole		cal/deg mole
Gd			g				
			c	0.000	87.	-56.4	46.41
Gd <sup>+</sup>			g		0.000	0.000	14.
Gd <sup>++</sup>			g		230.55		
Gd <sup>+++</sup>			aq		-168.8	120.65	-47.1
Gd(OH) <sub>3</sub>		std. state, hyp. m=1	c		-308.1	225.83	
GdCl <sub>3</sub>		$\alpha$	c		-245.5		
		std. state, hyp. m=1	aq		-288.9	189.55	-7.6
GdI <sub>3</sub>		$\beta$	c		-147.6		
		std. state, hyp. m=1	aq		-208.9	147.77	-31.3
Gd <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		std. state, hyp. m=1	aq		-968.3	631.25	-81.9
Gd <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O			c		-1518.9	974.72	155.8
							140.5

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 70-1. Europium (at. no., 63; at. wt., 152.0)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1949

Substance									
Formula	Description	State	$\Delta F_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 298.16°K (25°C)						
			At 0°K						
			kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole	
Eu		g		87.					4.968
		c	0.000	0.000	0.000	0.0000	45.10		
Eu <sup>+</sup>		g		219.23					
		g		480.0					
Eu <sup>++</sup>	std. state, hyp. m = 1	aq			-155.2	113.76			
		g							
Eu <sup>+++</sup>	std. state, hyp. m = 1	aq		-169.3	-165.1	121.02			
	$\alpha$	c		-247.1					
EuCl <sub>3</sub>	std. state, hyp. m = 1	aq		-289.4	-259.2	189.99			
Eu <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>	std. state, hyp. m = 1	aq		-989.3	-862.2	631.98			
		c			-1331.0	975.60			146.3

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 71-1. Samarium (at. no., 62; at. wt., 150.43)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1949

Formula	Substance	Description	State	At 298.16°K (25°C)				
				$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
Sm			g	0.000	87.	0.000	43.74	7.254
Sm <sup>+</sup>			c		0.000			
Sm <sup>++</sup>			g	217.7				
Sm <sup>+++</sup>			g	482.				
		std. state, hyp. m = 1	aq	-169.8	-165.9	121.60		
Sm(OH) <sub>3</sub>			c		-308.8	226.35		
SmCl <sub>3</sub>		$\alpha$	c	-249.8				
		std. state, hyp. m = 1	aq	-269.9	-259.9	190.50		
SmI <sub>3</sub>		$\beta$	c	-153.4				
		std. state, hyp. m = 1	aq	-209.9	-202.9	148.72		

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

Washington, D.C.

## SERIES I

National Bureau of Standards

Table 71-2. Samarium (at. no., 62; at. wt., 150.43)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1949

Formula	Substance	Description	State	At 0°K					At 298.16°K (25°C)		
				kcal/mole	ΔH <sup>o</sup>	kcal/mole	ΔF <sup>o</sup>	log <sub>10</sub> K <sup>o</sup>	S <sup>o</sup>	C <sub>p</sub> <sup>o</sup>	
Sm <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>			aq			-990.3	-863.8	633.15			
Sm <sub>2</sub> (SO <sub>4</sub> )·8H <sub>2</sub> O		std. state, hyp. m = 1	c				-1332.6	976.78			
SmCl <sub>3</sub> ·NH <sub>3</sub>			c			-282.3					
SmCl <sub>3</sub> ·2NH <sub>3</sub>			c			-310.3					
SmCl <sub>3</sub> ·3NH <sub>3</sub>			c			-336.8					
SmCl <sub>3</sub> ·4NH <sub>3</sub>			c			-361.9					
SmCl <sub>3</sub> ·5NH <sub>3</sub>			c			-385.2					
SmCl <sub>3</sub> ·8NH <sub>3</sub>			c			-452.2					
SmCl <sub>3</sub> ·9½NH <sub>3</sub>			c			-483.8					
SmCl <sub>3</sub> ·11½NH <sub>3</sub>			c			-524.3					



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 72-1. Promethium (at. no., 61; at. wt., 147.1) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1949									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole					
Pm			g c	0.400	0.300	0.0000			
Pm <sup>+</sup> Pm <sup>+++</sup>		std. state, hyp. m = 1	g aq		-170.4				
PmCl <sub>3</sub>		std. state, hyp. m = 1	c aq		-251.9 -290.5				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 73-1. Neodymium (at. no., 60; at. wt., 144.27)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1949

Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole		cal/deg mole	cal/deg mole
Nd			g	0.000	87.	0.000	0.000		7.2
			c, III						
Nd <sup>+</sup>			g		234.1				
Nd <sup>++</sup>			g						
Nd <sup>+++</sup>		std. state, hyp. m = 1	aq		-171.2	-167.6	122.85		
Nd <sub>2</sub> O <sub>3</sub>			c		-442.0				
Nd(OH) <sub>3</sub>			c			-309.6	226.93		
NdCl <sub>3</sub>			c		-254.3				
	$\alpha$		aq		-291.3	-261.6	191.75		
	std. state, hyp. m = 1		c		-692.3				
NdCl <sub>3</sub> ·6H <sub>2</sub> O			c		-158.9				
NdI <sub>3</sub>			c		-211.3				
	$\alpha$		aq			-204.6	149.97		
	std. state, hyp. m = 1								

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 73-2. Neodymium (at. no., 60; at. wt., 144.27)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1949

Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K								
	kcal/mole			kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole	
Nd <sub>2</sub> S <sub>3</sub>			c		-281.8				
Nd <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>			c		-948.1				
		std. state, hyp. m=1	aq		-993.1	-867.2	635.65		
Nd <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·5H <sub>2</sub> O			c		-1318.4				
Nd <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O			c		-1524.7	-1334.5	978.17		
NdCl <sub>3</sub> ·NH <sub>3</sub>			c		-286.0				
NdCl <sub>3</sub> ·2NH <sub>3</sub>			c		-314.5				
NdCl <sub>3</sub> ·4NH <sub>3</sub>			c		-364.6				
NdCl <sub>3</sub> ·5NH <sub>3</sub>			c		-388.3				
NdCl <sub>3</sub> ·8NH <sub>3</sub>			c		-444.2				
NdCl <sub>3</sub> ·11NH <sub>3</sub>			c		-505.8				
NdCl <sub>3</sub> ·12NH <sub>3</sub>			c		-525.1				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards SERIES I Washington, D.C.

Table 74-1. Praseodymium (at. no., 59; at. wt., 140.92) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1949									
Formula	Substance	Description	State	At 298.16°K (25°C)					
				$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole	
Pr			g						
			c	0.000		0.0000			6.8
Pr <sup>+</sup>			g						
Pr <sup>++</sup>			g	221.4					
Pr <sup>+++</sup>			aq	-172.7	-169.1	123.95			
PrO <sub>2</sub>		std. state, hyp. m=1	c						
Pr <sub>2</sub> O <sub>3</sub>			c	-234.0					
Pr <sub>6</sub> O <sub>11</sub>			c	-444.5					
			c	-1391.					
Pr(OH) <sub>3</sub>			c		-310.7	227.74			
PrCl <sub>3</sub>			c	-257.8					
	$\alpha$		aq	-292.8	-263.1	192.85			
PrCl <sub>3</sub> ·H <sub>2</sub> O		std. state, hyp. m=1	c	-330.8					
PrCl <sub>3</sub> ·7H <sub>2</sub> O			c	-764.5					



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 75-1. Cerium (at. no., 58; at. wt., 140.13)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1949

Formula	Substance	State	Description	ΔH <sup>o</sup> At 0°K kcal/mole	ΔH <sup>o</sup> kcal/mole	ΔH <sup>o</sup> kcal/mole	log <sub>10</sub> K <sub>f</sub> At 298.16°K (25°C)	S <sup>o</sup> cal/deg mole	C <sub>p</sub> <sup>o</sup> cal/deg mole
Ce		g		0.000	85.	0.000	0.000	13.8	6.19
Ce <sup>+</sup>		c, III			0.000				
Ce <sup>++</sup>		g			245.9				
Ce <sup>+++</sup>		g			531.1				
Ce <sup>++++</sup>		g			994.				
Ce <sup>++++</sup>		aq	std. state, hyp. m = 1		-173.8		124.97	-44.	
Ce <sup>++++</sup>		g			1842.				
Ce <sup>++++</sup>		aq	in HClO <sub>4</sub> (aq)		-134.4				
Ce <sup>++++</sup>		g			-170.5				
CeO <sub>2</sub>		c			-233.				15.1
CeO <sub>3</sub> ·2H <sub>2</sub> O		c			-389.				
Ce <sub>3</sub> Mg		c			-170.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 75-2. Cerium (at. no., 58; at. wt., 140.13) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)									
Substance		State	Description	At 298.16°K (25°C)					
Formula				$\Delta H_f^0$ kcal/mole	$\Delta H_f^0$ kcal/mole	$\Delta F_f^0$ kcal/mole	$\log_{10} K_f$	$S^0$ cal/deg mole	$C_p^0$ cal/deg mole
$\text{Ce}(\text{OH})^{+++}$		aq	std. state, hyp. $m=1$			-187.7	137.58		
$\text{Ce}(\text{OH})_2^{++}$		aq	std. state, hyp. $m=1$			-244.0	178.85		
$\text{CeCl}_3$		c	$\alpha$			-260.3			
		aq	std. state, hyp. $m=1$			-293.9	193.87	-5.	
$\text{CeI}_3$		c	$\alpha$			-164.4			
		aq	std. state, hyp. $m=1$			-213.9	152.09	34.	
$\text{CeS}_2$		c				-153.9			
$\text{Ce}_2\text{S}_3$		c				-298.7			
$\text{Ce}(\text{SO}_4)_2$		c				-560.			66.
$\text{Ce}_2(\text{SO}_4)_3$		c							
		aq	std. state, hyp. $m=1$			-873.0	639.90	-76.	
$\text{Ce}_2(\text{SO}_4)_3 \cdot 4\text{H}_2\text{O}$		c				-1308.			132.
$\text{Ce}_2(\text{SO}_4)_3 \cdot 6\text{H}_2\text{O}$		c				-1340.2	982.35		
$\text{Ce}_2(\text{SO}_4)_3 \cdot 9\text{H}_2\text{O}$		c				-1366.8	1023.83		

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 75-3. Cerium (at. no., 58; at. wt., 140.13)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1949

Formula	Substance	State	Description	At 298.16°K (25°C)				
				$\Delta H_f^0$ kcal/mole	$\Delta F_f^0$ kcal/mole	$\log_{10} K_f$	$S^0$ cal/deg mole	$C_p^0$ cal/deg mole
CeN		c						11.2
CeCl <sub>3</sub> ·2NH <sub>3</sub>		c		-78.3				
CeCl <sub>3</sub> ·4NH <sub>3</sub>		c		-322.0				
CeCl <sub>3</sub> ·8NH <sub>3</sub>		c		-375.1				
CeCl <sub>3</sub> ·8NH <sub>3</sub>		c		-455.2				
CeCl <sub>3</sub> ·12NH <sub>3</sub>		c		-525.4				
CeCl <sub>3</sub> ·20NH <sub>3</sub>		c		-657.7				
CeHg <sub>4</sub>		c		-21.5				96.
Ce <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>		c						
CeAl <sub>4</sub>		c, II		-39.				
Ce <sub>3</sub> Al		c		-22.				



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 76-1. Lanthanum (at. no., 57; at. wt., 138.92)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1949

Substance									
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole	
La		g		88.	79.	-57.9	43.57	5.438	
		c, III	0.000	0.000	0.000	0.0000	13.7	6.6	
La <sup>+</sup>		g		218.95					
La <sup>++</sup>		g		484.10					
La <sup>+++</sup>		g		927.59					
	std. state, hyp. m=1	aq		-176.2	-172.9	126.73	-44.		
La <sup>++++</sup>		g							
La <sub>2</sub> O <sub>3</sub>		c		-458.				24.2	
La <sub>3</sub> H <sub>8</sub>		c		-382.				37.	
La(OH) <sub>3</sub>		c			-312.8	229.28			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Substance		Description		State		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula						At 0°K	At 298.16°K (25°C)				
						kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
LaCl <sub>3</sub>	$\alpha$				c						
	std. state, hyp. m = 1				aq						
	in 400 H <sub>2</sub> O				aq		-263.6	-266.9	195.63	-5.	
	500 H <sub>2</sub> O				aq		-296.3				
	600 H <sub>2</sub> O				aq		-294.984				
	700 H <sub>2</sub> O				aq		-295.062				
	800 H <sub>2</sub> O				aq		-295.112				
	1000 H <sub>2</sub> O				aq		-295.158				
	1500 H <sub>2</sub> O				aq		-295.194				
	2000 H <sub>2</sub> O				aq		-295.254				
	3000 H <sub>2</sub> O				aq		-295.358				
	5000 H <sub>2</sub> O				aq		-295.430				
	10000 H <sub>2</sub> O				aq		-295.522				
	20000 H <sub>2</sub> O				aq		-295.638				
	50000 H <sub>2</sub> O				aq		-295.784				
	100000 H <sub>2</sub> O				aq		-295.914				
	200000 H <sub>2</sub> O				aq		-296.044				
	500000 H <sub>2</sub> O				aq		-296.114				
	1000000 H <sub>2</sub> O				aq		-296.168				
	$\infty$ H <sub>2</sub> O				aq		-296.216				
					aq		-296.234				
					aq		-296.3				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 76-3. Lanthanum (at. no., 57; at. wt., 138.92)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1949

Substance						
Formula	Description	State	$\Delta H_f^0$	$\Delta H_f^0$	$\Delta F_f^0$	$\log_{10} K_f$
			At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	At 298.16°K (25°C) kcal/mole	At 298.16°K (25°C) cal/deg mole
LaI <sub>3</sub>	$\alpha$	c		-167.4		
	std. state, hyp. m = 1	aq		-216.3	-209.9	153.85
LaS <sub>2</sub>		c		-156.7		
		c		-306.8		
La <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c				
	std. state, hyp. m = 1	aq		-1003.1	-877.8	643.41
	in 1800 H <sub>2</sub> O	aq		-994.06		
	2000 H <sub>2</sub> O	aq		-994.14		
	3000 H <sub>2</sub> O	aq		-994.41		
	4000 H <sub>2</sub> O	aq		-994.58		
	5000 H <sub>2</sub> O	aq		-994.68		
	7500 H <sub>2</sub> O	aq		-994.87		
	10000 H <sub>2</sub> O	aq		-995.02		
	20000 H <sub>2</sub> O	aq		-995.54		
	50000 H <sub>2</sub> O	aq		-996.60		

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards Washington, D.C.

## SERIES I

Table 76-4. Lanthanum (at. no., 57; at. wt., 138.92)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1949

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
$\text{La}_2(\text{SO}_4)_3$	in 100000 $\text{H}_2\text{O}$ 200000 $\text{H}_2\text{O}$ 500000 $\text{H}_2\text{O}$ 1000000 $\text{H}_2\text{O}$ 2000000 $\text{H}_2\text{O}$ $\infty$ $\text{H}_2\text{O}$		aq		-997.61				
					-998.60				
					-999.73				
					-1000.72				
					-1001.58				
					-1003.1				
$\text{La}_2(\text{SO}_4)_3 \cdot 9\text{H}_2\text{O}$			c			-1403.1	1028.45		152.
LaN			c		-72.1				11.1
$\text{La}(\text{NO}_3)_3 \cdot 6\text{H}_2\text{O}$			c, II						100.7
$\text{La}(\text{CHO}_2)_3$	lanthanum formate		c		-229.				79.7
$\text{La}_2(\text{CN}_2)_3$	lanthanum cyanamide		c						
$\text{La}_2(\text{MoO}_4)_3$			c						86.
$\text{LaAl}_2$			c		-36.1				
$\text{LaAl}_4$			c, II		-42.2				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards Washington, D.C.

Table 79-1. Plutonium (at. no., 94; at. wt., 239.) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1949									
Formula	Substance	Description	State	$\Delta H_f^{\circ}$	$\Delta H_f^{\circ}$	$\Delta F^{\circ}$	$\log_{10} K_f$	$S^{\circ}$	$C_p^{\circ}$
				At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole		cal/deg mole	cal/deg mole
Pu			g						
Pu <sup>+</sup>			c	0.000	0.000	0.000	0.0000		
Pu <sup>+++</sup>			g						
Pu <sup>++++</sup>			aq						
			aq						
		in HClO <sub>4</sub> ·55H <sub>2</sub> O							
PuO <sub>2</sub>			c						
PuO <sub>2</sub> <sup>+</sup>		in HClO <sub>4</sub> ·55H <sub>2</sub> O	aq						
PuO <sub>2</sub> <sup>++</sup>		in HClO <sub>4</sub> ·55H <sub>2</sub> O	aq						
PuH <sub>2</sub>			c						
Pu(OH) <sub>4</sub>			c						
PuF <sub>3</sub>			aq						
PuCl <sub>3</sub>			c						
			aq						
PuOCl <sub>3</sub>			c						
PuBr <sub>3</sub>			c						
PuI <sub>3</sub>			aq						
			c						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 80-1. Neptunium (at. no., 93; at. wt., 237)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1949

Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	At 0°K			At 298.16°K (25°C)						
	kcal/mole			kcal/mole	kcal/mole					kcal/mole
Np			g	0.000	0.000	0.000	0.0000			
Np <sup>+</sup>			c							
Np <sup>+++</sup>		in HCl•55H <sub>2</sub> O	g							
Np <sup>++++</sup>		in HCl•55H <sub>2</sub> O	aq		-127.3					
			aq		-133.2					
NpO <sub>2</sub> <sup>+</sup>		in HCl•55H <sub>2</sub> O	aq		-230.8					
NpO <sub>2</sub> <sup>++</sup>		in HCl•55H <sub>2</sub> O	aq		-208.5					
NpF <sub>3</sub>			c		-360.					
NpF <sub>4</sub>			c		-428.					
NpCl <sub>3</sub>			c		-216.					
		in HCl•55H <sub>2</sub> O	aq		-291.2					
NpCl <sub>4</sub>			c		-237.					
NpCl <sub>5</sub>			c		-246.					
NpBr <sub>3</sub>			c		-174.					
NpBr <sub>4</sub>			c		-183.					
NpI <sub>3</sub>			c		-120.					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 81-1. Uranium (at. no., 92; at. wt., 238.07) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1949									
Substance			$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	Description	State	At 0°K	At 298.16°K (25°C)					
			kcal/mole	kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole
U		g		125.					
U <sup>+</sup>		c, III	0.000	0.000	0.000	0.0000	12.03	6.57	
U <sup>++</sup>		g		218.					
U <sup>+++</sup>		g							
	std. state, hyp. m = 1.	aq		-123.0	-124.4	91.18	-30.		
U <sup>++++</sup>	std. state, hyp. m = 1	aq		-146.7	-138.4	101.44	-78.		

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 81-2. Uranium (at. no., 92; at. wt., 238.07)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1949

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	At 298.16°K (25°C)				
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
UO <sub>2</sub>			c		-270.	-257.	188.4	18.6	
UO <sub>2</sub> <sup>+</sup>		std. state, hyp. m=1	aq		-247.4	-237.6	174.16	12.	
UO <sub>2</sub> <sup>++</sup>		std. state, hyp. m=1	aq		-250.4	-236.4	173.28	-17.	
UO <sub>3</sub>			c		-302.	-283.	207.4	23.57	
UO <sub>3</sub> ·H <sub>2</sub> O			c		-375.4				
UO <sub>3</sub> ·2H <sub>2</sub> O			c		-446.2				
UO <sub>4</sub> ·2H <sub>2</sub> O			c		-436.				
U <sub>3</sub> O <sub>8</sub>			c		-898.				
UH <sub>3</sub>			c		-30.4				
U(OH) ***		std. state, hyp. m=1	aq		-204.1	-193.5	141.83	-30.	



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 81-3. Uranium (at. no., 92; at. wt., 238.07) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1949									
Substance		Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	At 0°K			At 298.16°K (25°C)					
	kcal/mole			kcal/mole	kcal/mole				cal/deg mole
UF <sub>3</sub>			c	-357.	-339.	248.5	26.		
UF <sub>4</sub>			c	-443.	-421.	308.6	36.1	28.12	
UF <sub>5</sub>			c	-488.	-461.	337.9	43.		
UF <sub>6</sub>			g	-505.	-485.	355.5	90.76		
			c	-517.	-486.	356.2	54.45		
UO <sub>2</sub> F <sub>2</sub>			c				32.40	24.68	
UCl <sub>3</sub>			c	-213.0	-196.9	144.32	37.99		
UCl <sub>4</sub>			c	-251.2	-230.0	168.59	47.4		
UCl <sub>5</sub>			c	-262.1	-237.4	174.01	62.		
UCl <sub>6</sub>			c	-272.4	-241.5	177.02	68.3		
UO <sub>2</sub> Cl <sub>2</sub>			aq	-331.					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 81-4. Uranium (at. no., 92; at. wt., 238.07)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1949

Substance										
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f^\circ$	$S^\circ$	$C_p^\circ$		
			At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole		
UBr <sub>3</sub>		c		-170.1	-164.7	120.72	49.			
UBr <sub>4</sub>		c		-196.6	-188.5	138.17	58.			
UO <sub>2</sub> Br <sub>2</sub>		aq		-308.2						
UI <sub>3</sub>		c		-114.7	-115.3	84.51	56.			
UI <sub>4</sub>		c		-127.0	-126.1	92.43	65.			
UCl <sub>3</sub> I		c		-219.9	-204.4	149.82	54.			
UBr <sub>3</sub> I		c		-177.1						
UO <sub>2</sub> SO <sub>4</sub>	std. state, hyp. m = 1	aq		-467.3	-413.7	303.24	-13.			
UO <sub>2</sub> SO <sub>4</sub> ·3H <sub>2</sub> O		c		-666.8	-586.0	429.53	63.			
U(SO <sub>4</sub> ) <sub>2</sub>		c		-563.						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 81-5. Uranium (at. no., 92; at. wt., 238.07)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1949

Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 298.16°K (25°C)					cal/deg mole	cal/deg mole		
	At 0°K kcal/mole			kcal/mole	kcal/mole			kcal/mole	
UN			c		-80.	-75.	55.0	18.	
U <sub>2</sub> N <sub>3</sub>			c		-213.	-194.	142.2	29.	
UO <sub>2</sub> (NO <sub>3</sub> ) <sub>2</sub>			c		-329.2	-273.1	200.18	66.	
		std. state, hyp. m=1	aq		-349.1	-289.2	211.98	53.	
UO <sub>2</sub> (NO <sub>3</sub> ) <sub>2</sub> ·H <sub>2</sub> O			c		-404.8	-335.3	245.77	76.	
UO <sub>2</sub> (NO <sub>3</sub> ) <sub>2</sub> ·2H <sub>2</sub> O			c		-480.0	-396.6	290.70	85.	
UO <sub>2</sub> (NO <sub>3</sub> ) <sub>2</sub> ·3H <sub>2</sub> O			c		-552.2	-454.7	333.29	94.	
UO <sub>2</sub> (NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O			c		-764.3	-625.0	458.12	120.85	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 81-6. Uranium (at. no., 92; at. wt., 238.07)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1949

Substance		State	Description	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} f_f^\circ$	$S^\circ$	$C_p^\circ$	
Formula	At 0°K			At 298.16°K (25°C)			cal/deg mole	cal/deg mole		
	kcal/mole			kcal/mole	kcal/mole	kcal/mole				
UC		c								
UC <sub>2</sub>		c								
UO <sub>2</sub> (C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>2</sub>		aq	uranyl acetate		-42.	-42.	30.8	14.		
UO <sub>2</sub> (C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>2</sub> ·2H <sub>2</sub> O		c			-484.0					
UO <sub>2</sub> (C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>2</sub> ·NH <sub>4</sub> C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ·6H <sub>2</sub> O		c			-624.9					
					-1045.8					
UO <sub>2</sub> CrO <sub>4</sub>		aq			-456.7					
UO <sub>2</sub> CrO <sub>4</sub> ·5½H <sub>2</sub> O		c			-838.8					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 83-1. Thorium (at. no., 90; at. wt., 232.12) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1949									
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole				
Th			g	0.000	0.000	0.000	0.0000	45.43	7.7
Th <sup>+</sup>			c					13.6	
Th <sup>++++</sup>			g		-183.0				
			aq						
ThO <sub>2</sub>			c		-292.				20.38
ThH <sub>4</sub>			c		-43.				
Th(OH) <sub>4</sub>		"soluble"	c		-421.5				
ThF <sub>4</sub>			c		-477.				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards  
 SERIES I  
 Washington, D.C.

Table 83-2. Thorium (at. no., 90; at. wt., 232.12)  
 HEAT OF FORMATION AT 0°K, HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
 December 31, 1949

Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
			At 0°K	At 298.16°K (25°C)				
Formula	Description		kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole
$\text{ThCl}_4$		c		-285.				
$\text{ThCl}_4 \cdot 2\text{H}_2\text{O}$		aq		-343.0				
$\text{ThCl}_4 \cdot 4\text{H}_2\text{O}$		c		-437.4				
$\text{ThCl}_4 \cdot 7\text{H}_2\text{O}$		c		-589.2				
$\text{ThCl}_4 \cdot 8\text{H}_2\text{O}$		c		-805.9				
$\text{ThCl}_4 \cdot 8\text{H}_2\text{O}$		c		-877.6				
$\text{ThOCl}_2$		c		-274.8				
$\text{Th}(\text{OH})\text{Cl}_3 \cdot \text{H}_2\text{O}$		c		-377.4				
$\text{ThBr}_4$		c		-227.1				
$\text{ThBr}_4 \cdot 7\text{H}_2\text{O}$		aq		-298.6				
$\text{ThBr}_4 \cdot 10\text{H}_2\text{O}$		c		-753.7				
$\text{ThBr}_4 \cdot 12\text{H}_2\text{O}$		c		-971.6				
$\text{ThBr}_4 \cdot 12\text{H}_2\text{O}$		c		-1116.0				
$\text{ThOBr}_2$		c		-252.8				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Substance		HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1949				
Formula	Description	State	$\Delta H_f^\circ$	$\Delta G_f^\circ$	$\log_{10} K_f$	$S^\circ$
			At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole
ThI <sub>4</sub>		c		-131.		
ThOI <sub>2</sub>		c		-228.2		
ThOI <sub>2</sub> ·3H <sub>2</sub> O		c		-479.3		
Th(OH)I <sub>3</sub> ·10H <sub>2</sub> O		c		-952.4		
Th <sub>2</sub> S <sub>3</sub>		c		-262.0		
ThOSO <sub>4</sub>		c		-487.		
Th(SO <sub>4</sub> ) <sub>2</sub>		c		-602.		
Th(SO <sub>4</sub> ) <sub>2</sub> ·4H <sub>2</sub> O		aq		-616.8		
Th(SO <sub>4</sub> ) <sub>2</sub> ·8H <sub>2</sub> O		c		-882.7		
		c		-1168.6		
						41.

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 83-4. Thorium (at. no., 90; at. wt., 232.12)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1949

Formula	Substance	State	Description	At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	log <sub>10</sub> K <sub>f</sub>	S° cal/deg mole	C <sub>p</sub> <sup>o</sup> cal/deg mole
Th <sub>3</sub> N <sub>4</sub>		c						
Th(NO <sub>3</sub> ) <sub>4</sub>		aq						
	in 20 H <sub>2</sub> O	aq						
	25 H <sub>2</sub> O	aq						
	50 H <sub>2</sub> O	aq						
	100 H <sub>2</sub> O	aq						
	500 H <sub>2</sub> O	aq						
Th(NO <sub>3</sub> ) <sub>4</sub> ·NH <sub>4</sub> NO <sub>3</sub> ·H <sub>2</sub> O		c						
Th(NO <sub>3</sub> ) <sub>4</sub> ·NH <sub>4</sub> NO <sub>3</sub> ·8H <sub>2</sub> O		c						
5Th(NO <sub>3</sub> ) <sub>4</sub> ·4NH <sub>4</sub> NO <sub>3</sub> ·11H <sub>2</sub> O		c						
5Th(NO <sub>3</sub> ) <sub>4</sub> ·4NH <sub>4</sub> NO <sub>3</sub> ·25H <sub>2</sub> O		c						



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 83-5. Thorium (at. no., 90; at. wt., 232.12)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1949

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)		cal/deg mole	cal/deg mole
			kcal/mole	kcal/mole	kcal/mole		
$\text{ThCl}_4 \cdot \text{NH}_4\text{Cl}$		c					
$\text{ThCl}_4 \cdot 2\text{NH}_4\text{Cl} \cdot 10\text{H}_2\text{O}$		c					
$\text{ThCl}_4 \cdot 4\text{NH}_3$	$\alpha$	c					
$\text{ThCl}_4 \cdot 6\text{NH}_3$	$\alpha$	c					
$[\text{Th}(\text{NH}_3)_6]\text{Cl}_4$	$\beta$	c					
$\text{ThCl}_4 \cdot 7\text{NH}_3$	$\alpha$	c					
$[\text{Th}(\text{NH}_3)_6]\text{Cl}_4 \cdot \text{NH}_3$	$\beta$	c					
$\text{ThCl}_4 \cdot 12\text{NH}_3$	$\alpha$	c					
$[\text{Th}(\text{NH}_3)_6]\text{Cl}_4 \cdot 6\text{NH}_3$	$\beta$	c					
$\text{ThCl}_4 \cdot 18\text{NH}_3$	$\alpha$	c					
$[\text{Th}(\text{NH}_3)_6]\text{Cl}_4 \cdot 12\text{NH}_3$	$\beta$	c					
$\text{ThC}_2$		c					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 85-1. Beryllium (at. no., 4; at. wt. 9.013)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1949

Substance		At 298.16°K (25°C)				
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$
			At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole
Be		g				
		c				
Be <sup>+</sup>		g				
Be <sup>++</sup>		g				
	in acid solution	aq				
Be <sup>+++</sup>		g				
Be <sup>++++</sup>		g				
BeO		g				
		c				
BeO <sub>2</sub> <sup>---</sup>		aq				
BeH		g				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 85-2. Beryllium (at. no., 4; at. wt., 9.013) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1949									
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K					
				kcal/mole	kcal/mole			kcal/mole	cal/deg mole
Be(OH) <sub>2</sub>		$\alpha$	c		-216.8				
		$\beta$	c		-216.1				
BeO•Be(OH) <sub>2</sub>		precipitated	c		-366.2				
BeF <sub>2</sub>			aq		-251.4				
BeCl <sub>2</sub>			c		-122.3				
		in HCl (aq)	aq		-173.4				
		in C <sub>2</sub> H <sub>5</sub> O (ethanol)	c		-160.				
BeCl <sub>2</sub> •4H <sub>2</sub> O					-436.8				
BeBr <sub>2</sub>			c		-88.4				
		in HCl (aq)	aq		-151.				
BeI <sub>2</sub>			c		-50.6				
		in HCl (aq)	aq		-120.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

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Table 85-3. Beryllium (at. no., 4; at. wt., 9.013) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) December 31, 1949									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} f_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				
BeS			c		-55.9				21.
BeSO <sub>4</sub>	in	15 H <sub>2</sub> O	c		-286.0				
		18 H <sub>2</sub> O	aq		-301.50				
		20 H <sub>2</sub> O	aq		-301.89				
		25 H <sub>2</sub> O	aq		-302.08				
		30 H <sub>2</sub> O	aq		-302.42				
		40 H <sub>2</sub> O	aq		-302.65				
		50 H <sub>2</sub> O	aq		-302.92				
		100 H <sub>2</sub> O	aq		-303.09				
		200 H <sub>2</sub> O	aq		-303.50				
		400 H <sub>2</sub> O	aq		-303.83				
		1000 H <sub>2</sub> O	aq		-304.1				
		2000 H <sub>2</sub> O	aq		-304.46				
		5000 H <sub>2</sub> O	aq		-304.79				
			aq		-305.10				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

Washington, D.C.

Table 85-4. Beryllium (at. no. 4; at. wt., 9.013)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
December 31, 1949

Substance								
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole	cal/deg mole	
$\text{BeSO}_4 \cdot \text{H}_2\text{O}$		c		-361.				
$\text{BeSO}_4 \cdot 2\text{H}_2\text{O}$		c		-433.2				
$\text{BeSO}_4 \cdot 4\text{H}_2\text{O}$		c		-576.3				
$\text{BeSO}_4 \cdot 4\text{BeO}$		c		-871.4				
$\text{Be}_2\text{r}_2 \cdot 2\text{H}_2\text{S}$		c		-115.7				
$\text{BeI}_2 \cdot 2\text{H}_2\text{S}$		c		-74.2				
$\text{Be}_3\text{N}_2$		c		-135.7				
$\text{Be}(\text{NO}_3)_2$		aq		-186.3				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 85-5. Beryllium (at. no., 4; at. wt., 9.013)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

December 31, 1949

Formula	Substance		State	Description	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
					At 0°K	At 298.16°K (25°C)				
					kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
BeCl <sub>2</sub> ·2NH <sub>3</sub>		c			-206.4					
BeCl <sub>2</sub> ·4NH <sub>3</sub>		c			-265.2					
BeCl <sub>2</sub> ·6NH <sub>3</sub>		c			-303.1					
BeCl <sub>2</sub> ·12NH <sub>3</sub>		c			-414.9					
BeBr <sub>2</sub> ·4NH <sub>3</sub>		c			-239.4					
BeBr <sub>2</sub> ·6NH <sub>3</sub>		c			-278.8					
BeBr <sub>2</sub> ·10NH <sub>3</sub>		c			-353.6					
BeI <sub>2</sub> ·4NH <sub>3</sub>		c			-212.4					
BeI <sub>2</sub> ·6NH <sub>3</sub>		c			-252.0					
BeI <sub>2</sub> ·13NH <sub>3</sub>		c			-383.4					
Be <sub>2</sub> SiO <sub>4</sub>		c						15.4	22.84	
BeMoO <sub>4</sub>		c			-330.				25.	
Be(AlO <sub>2</sub> ) <sub>2</sub>		c								

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 86-1. Magnesium (at. no., 12; at. wt., 24.32)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance								
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
			At 298.16°K (25°C)					
			At 0°K kcal/mole	kcal/mole	kcal/mole			cal/deg mole
Mg		g		35.9	27.6	-20.23	35.504	4.9680
		c	0.000	0.000	0.000	0.0000	7.77	5.71
Mg +		g		213.663				
Mg ++		g		561.788				
	std. state, hyp. m = 1	aq		-110.41	-108.99	79.888	-28.2	
Mg +++		g		2410.91				
Mg ++++		g		4932.91				
Mg ++++		g		8191.4				
Mg ++++		g		12502.2				
Mg ++++		g		17699.6				
Mg ++++		g		23834.5				
Mg ++++		g		31398.				
Mg ++++		g		39872.				
Mg ++++		g		80490.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Table 66-2. Magnesium (at. no., 12; at. wt., 24.32)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

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Formula	Substance	Description	State	At 0°K					At 298.16°K (25°C)		
				$\Delta H_f^\circ$		$\Delta F_f^\circ$		$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	cal/deg mole
				kcal/mole	kcal/mole	kcal/mole	kcal/mole				
MgO			c		-143.84	-136.13		99.781	6.4		8.94
MgO <sub>2</sub>		finely divided	c		-142.95	-135.31		99.180	6.66		9.03
MgH			g				34.	-24.9	47.61		7.050
Mg(OH) <sub>2</sub>			c		-221.00	-199.27		146.062	15.09		18.41
MgF			g		-20.	-27.		19.8	54.85		7.82
MgF <sub>2</sub>			c		-263.5	-250.8		183.83	13.68		14.72
MgCl			g		2.						
MgCl <sub>2</sub>			c		-153.40	-141.57		103.769	21.4		17.04
		std. state, hyp. m = 1	aq		-190.46	-171.69		125.846	-1.9		
		in 10 H <sub>2</sub> O	aq		-185.65						
		12 H <sub>2</sub> O	aq		-186.06						
		15 H <sub>2</sub> O	aq		-187.25						
		20 H <sub>2</sub> O	aq		-188.05						



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table B6-3. Magnesium (at. no., 12; at. wt., 24.32)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	Description	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	At 298.16°K (25°C)									
	At 0°K kcal/mole			kcal/mole	kcal/mole		kcal/mole	cal/deg mole	cal/deg mole	
MgCl <sub>2</sub>	in	aq			-188.43					
		aq			-189.14					
		aq			-189.37					
		aq			-189.49					
		aq			-189.72					
		aq			-189.81					
		aq			-189.88					
		aq			-189.93					
		aq			-190.05					
		aq			-190.11					
		aq			-190.15					
		aq			-190.20					
		aq			-190.25					
		aq			-190.31					
		aq			-190.38					
		aq			-190.41					
		aq			-190.43					
		aq			-190.46					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 86-4. Magnesium (at. no., 12; at. wt., 24.32) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		State	Description	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				
$MgCl_2 \cdot H_2O$		c			-231.15	-206.11	151.076	32.8	27.48
$MgCl_2 \cdot 2H_2O$		c			-305.99	-267.32	195.942	43.0	38.02
$MgCl_2 \cdot 4H_2O$		c			-454.00	-390.49	286.224	63.1	57.70
$MgCl_2 \cdot 6H_2O$		c			-597.42	-505.65	370.635	87.5	75.46
$MgO \cdot MgCl_2$		c			-312.6				
$MgO \cdot MgCl_2 \cdot 6H_2O$		c			-742.1				
$MgO \cdot MgCl_2 \cdot 16H_2O$		c			-1440.8				
$Mg(ClO_4)_2$		c			-140.6				
		aq			-172.5				
$Mg(ClO_4)_2 \cdot 2H_2O$		c			-290.7				
$Mg(ClO_4)_2 \cdot 4H_2O$		c			-438.6				
$Mg(ClO_4)_2 \cdot 6H_2O$		c			-583.2				
$Mg(OH)Cl$		c			-191.3	-175.0	-128.27	19.8	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 86-5. Magnesium (at. no., 12; at. wt., 24.32)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	Description	At 298.16°K (25°C)					
Formula	$\Delta H_f^\circ$			$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
	At 0°K kcal/mole			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole
MgBr <sub>2</sub>	std. state, hyp m = 1	c							
	in	aq							
	25 H <sub>2</sub> O	aq							
	50 H <sub>2</sub> O	aq							
	75 H <sub>2</sub> O	aq							
	100 H <sub>2</sub> O	aq							
	250 H <sub>2</sub> O	aq							
	500 H <sub>2</sub> O	aq							
	1000 H <sub>2</sub> O	aq							
	2500 H <sub>2</sub> O	aq							
	5000 H <sub>2</sub> O	aq							
	10000 H <sub>2</sub> O	aq							
	50000 H <sub>2</sub> O	aq							
	∞ H <sub>2</sub> O	aq							
MgBr <sub>2</sub> ·6H <sub>2</sub> O		c							

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 86-6. Magnesium (at. no., 12; at. wt., 24.32)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Formula	Substance	Description	State	At 298.16°K (25°C)					$C_p^0$ cal/deg mole
				$\Delta H_f^0$ kcal/mole	$\Delta H_f^0$ kcal/mole	$\Delta F_f^0$ kcal/mole	$\log_{10} K_f$	$S^0$ cal/deg mole	
				At 0°K kcal/mole	kcal/mole	kcal/mole		cal/deg mole	
MgI <sub>2</sub>			c		-86.0				
		std. state, hyp. m = 1	aq		-137.15	-133.69	97.993	24.1	
MgS			c		-83.0				
MgSO <sub>3</sub>			c		-241.0				
MgSO <sub>3</sub> ·3H <sub>2</sub> O			c		-461.6				
MgSO <sub>3</sub> ·6H <sub>2</sub> O			c		-673.4				
MgSO <sub>4</sub>			c		-305.5	-280.5	152.83	21.9	23.01
		std. state, hyp. m = 1	aq		-327.31	-286.33	209.876	-24.0	
		in 200 H <sub>2</sub> O	aq		-326.23				
		400 H <sub>2</sub> O	aq		-326.34				
		500 H <sub>2</sub> O	aq		-326.37				
		1000 H <sub>2</sub> O	aq		-326.47				
		1500 H <sub>2</sub> O	aq		-326.54				
		2000 H <sub>2</sub> O	aq		-326.58				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 86-7. Magnesium (at. no., 12; at. wt., 24.32)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C)		cal/deg mole	
MgSO <sub>4</sub>	in	3000 H <sub>2</sub> O	aq					
		5000 H <sub>2</sub> O	aq					
		10000 H <sub>2</sub> O	aq					
		20000 H <sub>2</sub> O	aq					
		50000 H <sub>2</sub> O	aq					
		100000 H <sub>2</sub> O	aq					
		200000 H <sub>2</sub> O	aq					
		500000 H <sub>2</sub> O	aq					
		1000000 H <sub>2</sub> O	aq					
		∞ H <sub>2</sub> O	aq					
MgSO <sub>4</sub> · 2H <sub>2</sub> O			c					
MgSO <sub>4</sub> · 4H <sub>2</sub> O			c					
MgSO <sub>4</sub> · 6H <sub>2</sub> O			c					
MgSO <sub>4</sub> · 7H <sub>2</sub> O			c					
MgS <sub>2</sub> O <sub>3</sub>			aq					
MgS <sub>2</sub> O <sub>3</sub> · 3H <sub>2</sub> O			c					
MgS <sub>2</sub> O <sub>3</sub> · 6H <sub>2</sub> O			c					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 86-8. Magnesium (at. no., 12; at. wt., 24.32)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		State	Description	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K								
	kcal/mole			kcal/mole	kcal/mole	cal/deg mole	cal/deg mole		
MgTe		c			-50.				24.99
Mg <sub>3</sub> N <sub>2</sub>	$\alpha$	c			-110.24				
Mg(NO <sub>3</sub> ) <sub>2</sub>	std. state, hyp. m = 1	c			-188.72	-140.63	103.080	39.2	33.94
	in	aq			-209.15	-161.81	118.605	41.8	
	15 H <sub>2</sub> O	aq			-208.410				
	25 H <sub>2</sub> O	aq			-208.612				
	50 H <sub>2</sub> O	aq			-208.658				
	100 H <sub>2</sub> O	aq			-208.706				
	250 H <sub>2</sub> O	aq			-208.772				
	500 H <sub>2</sub> O	aq			-208.820				
	1000 H <sub>2</sub> O	aq			-208.865				
	2500 H <sub>2</sub> O	aq			-208.918				
	5000 H <sub>2</sub> O	aq			-208.958				
	10000 H <sub>2</sub> O	aq			-209.000				
	50000 H <sub>2</sub> O	aq			-209.074				
	100000 H <sub>2</sub> O	aq			-209.098				
	500000 H <sub>2</sub> O	aq			-209.124				
	$\infty$ H <sub>2</sub> O	aq			-209.15				

Table 86-9. Magnesium (at. no., 12; at. wt., 24.32)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description		At 0°K kcal/mole	kcal/mole			At 298.16°K (25°C) kcal/mole	cal/deg mole
$Mg(NO_3)_2 \cdot 6H_2O$		c		-624.36				
$Mg(NH_3)_2^{++}$		aq		-147.8				
$MgCl_2 \cdot NH_3$		c		-185.8				
$MgCl_2 \cdot 2NH_3$		c		-215.2				
		aq		-227.8				
$MgBr_2 \cdot NH_3$		c		-157.1				
$MgBr_2 \cdot 2NH_3$		c		-188.8				
$MgI_2 \cdot 2NH_3$		c		-156.0				
$Mg(NH_3)_2SO_4$		aq		-364.6				
$3MgSO_3 \cdot (NH_4)_2SO_3 \cdot 6H_2O$		c		-1380.2				
$3MgSO_3 \cdot (NH_4)_2SO_3 \cdot 18H_2O$		c		-2230.2				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 86-10. Magnesium (at. no., 12; at. wt., 24.32)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
$Mg_3(PO_4)_2$			c		-961.5				
$MgHPO_4$			aq		-422.2				
$Mg(NH_4)PO_4 \cdot 6H_2O$			c		-881.0				
$Mg_3(AsO_4)_2$			c		-731.3				
$MgHAsO_4$			aq		-349.2				
$Mg(H_2AsO_4)_2$			aq		-541.0				
$Mg(NH_4)AsO_4 \cdot 6H_2O$			c		-800.7				
$Mg_3Sb_2$			c		-68.1				
$Mg_3Bi_2$			c		-36.5				



Table 86-11. Magnesium (at. no., 12; at. wt., 24.32) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	Description		At 0°K kcal/mole	kcal/mole					
MgCO <sub>3</sub>		c		-266.		180.3	15.7	18.05	
Mg(C <sub>2</sub> H <sub>3</sub> O <sub>3</sub> ) <sub>2</sub>	magnesium glycolate	c		-415.9					
		aq		-420.3					
Mg(C <sub>2</sub> H <sub>3</sub> O <sub>3</sub> ) <sub>2</sub> ·2H <sub>2</sub> O		c		-558.7					
MgCl <sub>2</sub> ·6CH <sub>3</sub> OH		c		-509.					
MgCl <sub>2</sub> ·6C <sub>2</sub> H <sub>5</sub> OH		c		-610.					
MgCN <sub>2</sub>	magnesium cyanamide	c		-60.3					
Mg(CN) <sub>2</sub>		aq		-31.9					
Mg(NO <sub>3</sub> ) <sub>2</sub> ·6CH <sub>3</sub> OH		c		-560.					
Mg(NO <sub>3</sub> ) <sub>2</sub> ·6C <sub>2</sub> H <sub>5</sub> OH		c		-603.					
Mg <sub>2</sub> Si		c		-18.6					
MgSiO <sub>3</sub>		c		-337.9		247.16	16.2	19.56	
Mg <sub>2</sub> SiO <sub>4</sub>	forsterite	c		-488.2		337.03	22.7	28.21	
Mg <sub>2</sub> Sn		c		-17.0					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 86-12. Magnesium (at. no., 12; at. wt., 24.32)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)		
			kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
$Mg_2 \cdot Pb$ $2MgI_2 \cdot PbI_2$		c	-12.6			
		c	-205.0			
$MgTi$		c	-11.			
$MgZn_2$		c	-12.6			
$MgCd$		c, II	-9.2			
$MgHgBr_4$		aq	-207.6			
$MgHg_2Br_4$		aq	-244.9			
$Mg_2HgBr_6$		aq	376.7			
$Mg_4HgBr_{10}$		aq	-713.7			
$MgHg(CN)_4$		aq	21.8			
$MgHg_2(CN)_6$		aq	86.1			
$2Hg(CN)_2 \cdot MgCl_2$		aq	-52.7			
$2Hg(CN)_2 \cdot MgCl_2 \cdot 6H_2O$		c	-472.1			
$2Hg(CN)_2 \cdot MgBr_2$		aq	-31.6			
$2Hg(CN)_2 \cdot MgBr_2 \cdot 8H_2O$		c	-593.4			
$2Hg(CN)_2 \cdot MgI_2$		aq	-4.4			
$2Hg(CN)_2 \cdot MgI_2 \cdot 8H_2O$		c	-570.2			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Substance		HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)				
Formula	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$
			At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole		
MgCrO <sub>4</sub>		c				
MgCr <sub>2</sub> O <sub>4</sub>		aq				
		c				
MgTiO <sub>3</sub>		c				
Mg <sub>2</sub> Al <sub>3</sub>		c				
MgPr		c				
Mg <sub>3</sub> Pr		c				
MgLa		c				
Mg <sub>3</sub> La		c				
2La(NO <sub>3</sub> ) <sub>2</sub> ·3Mg(NO <sub>3</sub> ) <sub>2</sub> ·24H <sub>2</sub> O		c				
MgCe		c				
MgCe <sub>3</sub>		c				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 87-1. Calcium (at. no., 20; at. wt., 40.08) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$G_p^\circ$	
Formula	At 298.16°K (25°C)								
	kcal/mole			kcal/mole	kcal/mole	cal/deg mole	cal/deg mole		
Ca			g	45.94	46.04	37.98	-27.839	36.993	4.968
Ca <sup>+</sup>			c, II	0.000	0.000	0.000	0.0000	9.95	6.28
Ca <sup>++</sup>			g	186.878	188.459				
			g	460.56	463.64				
		std. state, hyp. m = 1	aq		-129.77	-132.18	96.886	-13.2	
Ca <sup>+++</sup>			g	1641.5	1646.0				
Ca <sup>++++</sup>			g	3191.	3197.				
Ca <sup>+++++</sup>			g	5137.	5145.				
CaO			c		-151.9	-144.4	105.84	9.5	10.23
CaO <sub>2</sub>			c		-157.5				
CaO <sub>2</sub> ·8H <sub>2</sub> O			c		-719.7				
CaH			g	59.	58.7				
CaH <sub>2</sub>			c		-45.1	-35.8	26.24	10.	
Ca(OH) <sub>2</sub>			c		-235.80	-214.33	157.101	18.2	20.2
		std. state, hyp. m = 1	aq		-239.68	-207.37	152.000	-18.2	
Ca(OH) <sub>2</sub> ·H <sub>2</sub> O <sub>2</sub>			c		-291.3				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 87-2. Calcium (at. no., 20; at. wt., 40.08) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950											
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
				At 0°K						At 298.16°K (25°C)	
				kcal/mole	kcal/mole					kcal/mole	cal/deg mole
CaF			g	-9.3	-9.3	-277.7	203.55	16.46	16.02		
CaF <sub>2</sub>			c	-290.3	-290.3	-264.34	193.758	-17.8			
		std. state, hyp. m = 1	aq	-287.09	-287.09						
CaCl			g	7	6.7	-179.3	131.42	27.2	17.36		
CaCl <sub>2</sub>			c	-190.0	-190.0	-194.68	142.844	13.1			
		std. state, hyp. m = 1	aq	-209.82	-209.82						
		in 25 H <sub>2</sub> O		-208.51	-208.51						
		50 H <sub>2</sub> O		-208.86	-208.86						
		100 H <sub>2</sub> O		-209.055	-209.055						
		200 H <sub>2</sub> O		-209.200	-209.200						
		400 H <sub>2</sub> O		-209.3	-209.3						
		570 H <sub>2</sub> O		-209.320	-209.320						
		800 H <sub>2</sub> O		-209.385	-209.385						
		1000 H <sub>2</sub> O		-209.41	-209.41						
		5000 H <sub>2</sub> O		-209.60	-209.60						
		∞ H <sub>2</sub> O		-209.82	-209.82						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 87-3. Calcium (at. no., 20; at. wt., 40.08) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		State	Description	At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	log <sub>10</sub> K <sub>f</sub>	S° cal/deg mole	C <sub>p</sub> <sup>o</sup>	
Formula									
CaCl <sub>2</sub> ·H <sub>2</sub> O		c			-265.1				
CaCl <sub>2</sub> ·2H <sub>2</sub> O		c			-335.5				
CaCl <sub>2</sub> ·4H <sub>2</sub> O		c			-480.2				
CaCl <sub>2</sub> ·6H <sub>2</sub> O		c			-623.15				
CaOCl <sub>2</sub>		c			-178.6				
		aq			-188.1				
CaOCl <sub>2</sub> ·H <sub>2</sub> O		c			-249.2				
Ca(OCl) <sub>2</sub>		aq			-180.0				
CaCl <sub>2</sub> ·2CaO		c			-505.				
CaCl <sub>2</sub> ·3CaO		c			-654.				
CaCl <sub>2</sub> ·3CaO·H <sub>2</sub> O		c			-910.6				
CaCl <sub>2</sub> ·3CaO·6H <sub>2</sub> O		c			-1833.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 87-4. Calcium (at. no., 20; at. wt., 40.08) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 298.16°K (25°C)								
	kcal/mole			kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole	
CaBr <sub>2</sub>			c						
		std. state, hyp. m = 1	aq		-156.8		114.93	31.	
		in 400 H <sub>2</sub> O	aq		-181.33		132.912	25.4	
		1000 H <sub>2</sub> O	aq		-161.3				
		2000 H <sub>2</sub> O	aq		-187.57				
		4000 H <sub>2</sub> O	aq		-187.0				
CaBr <sub>2</sub> ·6H <sub>2</sub> O		∞ H <sub>2</sub> O	aq		-187.17				
			aq		-187.24				
			aq		-187.30				
			aq		-187.57				
CaI <sub>2</sub>			c		-597.2				
			c		-1810.				
CaI <sub>2</sub> ·8H <sub>2</sub> O			c		-127.8		92.80	34.	
		std. state, hyp. m = 1	aq		-156.51		114.991	39.1	
		in 400 H <sub>2</sub> O	aq		-156.0				
		∞ H <sub>2</sub> O	aq		-156.51				
			c		-700.7				
			c		-1779.				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 87-5. Calcium (at. no., 20; at. wt., 40.08)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		State	Description	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula				At 298.16°K (25°C)						
	kcal/mole			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole		
CaS		c			-115.3	-114.1	83.63	13.5	11.33	
CaSO <sub>3</sub>		aq			-119.8					
CaSO <sub>3</sub> •2H <sub>2</sub> O		c						24.2	21.90	
CaSO <sub>4</sub>		c			-421.2	-374.1	274.21	44.	42.7	
	anhydrite	c			-342.42	-315.56	231.301	25.5	23.8	
	soluble α	c			-340.27	-313.52	229.806	25.9	23.9	
	soluble β	c			-339.21	-312.46	229.029	25.9	23.8	
	std. state, hyp. m = 1	aq			-346.67	-309.52	226.874	-9.1		
	in 2500 H <sub>2</sub> O	aq			-345.95					
	5000 H <sub>2</sub> O	aq			-346.040					
	7500 H <sub>2</sub> O	aq			-346.116					
	10000 H <sub>2</sub> O	aq			-346.174					
	20000 H <sub>2</sub> O	aq			-346.312					
	50000 H <sub>2</sub> O	aq			-346.446					
	100000 H <sub>2</sub> O	aq			-346.520					
	200000 H <sub>2</sub> O	aq			-346.574					
	500000 H <sub>2</sub> O	aq			-346.620					
	∞ H <sub>2</sub> O	aq			-346.67					
CaSO <sub>4</sub> •½H <sub>2</sub> O	α	c			-376.47	-343.02	251.429	31.2	28.6	
CaSO <sub>4</sub> •2H <sub>2</sub> O	β	c			-375.97	-342.78	251.253	32.1	29.6	
		c			-483.06	-429.19	314.590	46.36	44.5	



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 87-6. Calcium (at. no., 20; at. wt., 40.08) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				
$\text{CaS}_2\text{O}_3$	in 35 $\text{H}_2\text{O}$		aq		-282.86				
	100 $\text{H}_2\text{O}$		aq		-283.10				
	200 $\text{H}_2\text{O}$		aq		-283.25				
	400 $\text{H}_2\text{O}$		aq		-283.32				
	1000 $\text{H}_2\text{O}$		aq		-283.4				
$\text{CaS}_2\text{O}_3 \cdot 6\text{H}_2\text{O}$			c		-602.2		441.40		
CaSe			c		-74.7		53.87	16.	
CaTe			c					19.	
$\text{Ca}_3\text{N}_2$			c		75.8				
		calcium azide	c		-103.2		64.57	25.	22.5
$\text{CaM}_2\text{O}_7 \cdot 4\text{H}_2\text{O}$			c		-405.7				
$\text{Ca}(\text{NO}_2)_2$			c		-178.3				
			aq		-180.5				

National Bureau of Standards  
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Table 87-7. Calcium (at. no., 20; at. wt., 40.08)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		State	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description							
Ca(NO <sub>3</sub> ) <sub>2</sub>	std. state, hyp. m=1	c						
	in							
	10 H <sub>2</sub> O	aq		-224.00	-177.34	129,988	46.2	35.69
	25 H <sub>2</sub> O	aq		-228.51	-185.00	135,602	56.8	
	50 H <sub>2</sub> O	aq		-229.48				
	75 H <sub>2</sub> O	aq		-229.31				
	100 H <sub>2</sub> O	aq		-229.123				
	200 H <sub>2</sub> O	aq		-228.970				
	300 H <sub>2</sub> O	aq		-228.832				
	400 H <sub>2</sub> O	aq		-228.556				
	500 H <sub>2</sub> O	aq		-228.470				
	1000 H <sub>2</sub> O	aq		-228.426				
	2000 H <sub>2</sub> O	aq		-228.400				
	5000 H <sub>2</sub> O	aq		-228.345				
	10000 H <sub>2</sub> O	aq		-228.326				
∞ H <sub>2</sub> O	aq		-228.344					
5000 H <sub>2</sub> O	aq		-228.374					
10000 H <sub>2</sub> O	aq		-228.436					
50000 H <sub>2</sub> O	aq		-228.456					
∞ H <sub>2</sub> O	aq		-228.484					
∞ H <sub>2</sub> O	aq		-228.51					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 87-8. Calcium (at. no., 20; at. wt., 40.08)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$
Formula	Description		At 0°K			At 298.16°K (25°C)		
$\text{Ca}(\text{NO}_3)_2$	in 100 $\text{C}_2\text{H}_5\text{OH}$			-231.28				
$\text{Ca}(\text{NO}_3)_2 \cdot 2\text{H}_2\text{O}$	$\infty$ $\text{C}_2\text{H}_5\text{OH}$	c		-231.58	-293.51	215.139	64.3	
$\text{Ca}(\text{NO}_3)_2 \cdot 3\text{H}_2\text{O}$		c		-368.00	-349.0	255.81	74.	
$\text{Ca}(\text{NO}_3)_2 \cdot 4\text{H}_2\text{O}$		c, II		-437.18	-406.5	297.96	81.	
$\text{Ca}(\text{NH}_2)_2$		c		-509.37				
$\text{Ca}(\text{NO}_3)_2 \cdot \text{Ca}(\text{OH})_2$		c		-91.6				
$\text{Ca}(\text{NO}_3)_2 \cdot \text{Ca}(\text{OH})_2$		c		-461.7				
$\text{Ca}(\text{NO}_3)_2 \cdot \text{Ca}(\text{OH})_2 \cdot 2\text{H}_2\text{O}$		c		-641.1				
$\text{CaCl}_2 \cdot \text{NH}_3$		c		-217.3				
$\text{CaCl}_2 \cdot 2\text{NH}_3$		c		-242.8				
$\text{CaCl}_2 \cdot 4\text{NH}_3$		c		-285.0				
$\text{CaCl}_2 \cdot 8\text{NH}_3$		c		-368.5				
$\text{CaBr}_2 \cdot \text{NH}_3$		c		-189.6				
$\text{CaBr}_2 \cdot 2\text{NH}_3$		c		-217.1				
$\text{CaBr}_2 \cdot 6\text{NH}_3$		c		-307.3				
$\text{CaBr}_2 \cdot 8\text{NH}_3$		c		-349.0				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 87-9. Calcium (at. no., 20; at. wt., 40.08)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)		cal/deg mole	cal/deg mole
			kcal/mole	kcal/mole	kcal/mole		
$\text{CaI}_2 \cdot \text{NH}_3$		c		-157.7			
$\text{CaI}_2 \cdot 2\text{NH}_3$		c		-186.9			
$\text{CaI}_2 \cdot 6\text{NH}_3$		c		-285.3			
$\text{CaI}_2 \cdot 8\text{NH}_3$		c		-316.1			
$\text{Ca}_3\text{P}_2$		c		-120.5			
$\text{Ca}_3(\text{PO}_4)_2$		c, $\alpha$		-986.2	-929.7	57.6	55.35
		c, $\beta$		-988.9	-932.0	56.4	54.45
$\text{CaHPO}_4$		c		-435.2	-401.5	21.	
$\text{CaHPO}_4 \cdot 2\text{H}_2\text{O}$		c		-576.0	-514.6	40.	
$\text{Ca}(\text{H}_2\text{PO}_4)_2$		c		-744.4			
$\text{Ca}(\text{H}_2\text{PO}_4)_2 \cdot \text{H}_2\text{O}$		c		-618.0			
$\text{Ca}(\text{H}_2\text{PO}_4)_2 \cdot \text{H}_2\text{O} \cdot \text{NH}_3$		c		-840.6			
$\text{Ca}(\text{H}_2\text{PO}_4)_2 \cdot \text{H}_2\text{O} \cdot 2\text{NH}_3$		c		-860.7			
$\text{Ca}(\text{H}_2\text{PO}_4)_2 \cdot \text{H}_2\text{O} \cdot 4\text{NH}_3$		c		-898.0			
	precipitated						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 87-10. Calcium (at. no., 20; at. wt., 40.08)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance									
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\Delta P_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 298.16°K (25°C)						
			At 0°K kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole	
$\text{Ca}_3(\text{AsO}_4)_2$ $\text{CaHAsO}_4$ $\text{CaHAsO}_4 \cdot \text{H}_2\text{O}$ $\text{Ca}(\text{H}_2\text{AsO}_4)_2$		c		-796.					
		aq		-344.1					
		c		-410.		266.1	35.		
		aq		-560.8					
$\text{Ca}_3\text{Sb}_2$		c		-155.					
$\text{Ca}_3\text{Bi}_2$		c		-112.					
$\text{CaC}_2$		c		-15.0	-16.2	11.87	16.8	14.90	
$\text{CaCO}_3$	calcite	c		-286.45	-269.78	197.745	22.2	19.57	
	aragonite	c		-288.49	-269.53	197.562	21.2	19.42	
$\text{CaC}_2\text{O}_4 \cdot \text{H}_2\text{O}$	precipitated	c		-400.4	-361.9	265.27	37.28	36.40	
$\text{Ca}(\text{CH}_3\text{O}_2)_2$	calcium formate	c		-470.4	-418.2	306.53	47.		
$\text{Ca}(\text{HCO}_3)_2$ $\text{Ca}(\text{C}_2\text{H}_3\text{O}_2)_2$	in 400 $\text{H}_2\text{O}$	aq		-323.5					
	std. state, hyp. m = 1	aq		-324.6					
	calcium acetate	c		-460.13	-412.80	302.577	32.2		
$\text{Ca}(\text{C}_2\text{H}_3\text{O}_2)_2 \cdot \text{H}_2\text{O}$		aq		-355.0					
		c		-362.5					
		c		-425.1					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

Washington, D.C.

Table 87-11. Calcium (at. no., 20; at. wt., 40.08)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	Description	$\Delta H_f^\circ$					$S^\circ$	$C_p^\circ$
Formula	log <sub>10</sub> <i>f</i> <sup>o</sup>			$\Delta F_f^\circ$	At 298.16°K (25°C)		cal/deg mole	cal/deg mole		
					At 0°K	kcal/mole				
Ca(C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>2</sub>		c	calcium glycollate							
Ca(C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>2</sub> · 3H <sub>2</sub> O		aq								
Ca(C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>2</sub> · 5H <sub>2</sub> O		c								
Ca(C <sub>2</sub> H <sub>5</sub> O) <sub>2</sub>		c								
Ca(C <sub>2</sub> H <sub>5</sub> O) <sub>2</sub>		c	calcium ethylate							
Ca(C <sub>2</sub> H <sub>5</sub> O) <sub>2</sub> · 2C <sub>2</sub> H <sub>5</sub> OH		c								
3CaO · 4C <sub>2</sub> H <sub>5</sub> OH		c								
CaCl <sub>2</sub> · 3C <sub>2</sub> H <sub>5</sub> OH		c								
CaCl <sub>2</sub> · 4C <sub>2</sub> H <sub>5</sub> OH		c								
CaCN <sub>2</sub>		c								
Ca(CN) <sub>2</sub>		c	calcium cyanamide							
Ca(NO <sub>3</sub> ) <sub>2</sub> · 2CH <sub>3</sub> OH		c								
3CaO · Ca(CN) <sub>2</sub> · 15H <sub>2</sub> O		c								

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards		SERIES I						Washington, D.C.	
		Table 87-12. Calcium (at. no., 20; at. wt., 40.08) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950							
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole				
CaSi <sub>2</sub>			c		-36.				
Ca <sub>2</sub> Si			c		-50.				
Ca <sub>2</sub> Si <sub>2</sub>			c		-72.				
CaSiO <sub>3</sub>		pseudowollastonite	c, α		-377.4	-357.4	261.97	20.9	20.67
Ca <sub>2</sub> SiO <sub>4</sub>		wollastonite	c, β		-378.6	-358.2	262.55	19.6	20.38
Ca <sub>3</sub> SiO <sub>5</sub>			c, β		-538.0				
			c, γ		-539.0				
			c		-688.4				
CaSn			c		-38.				
CaSn <sub>3</sub>			c		-44.				
Ca <sub>2</sub> Sn			c		-75.				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 87-13. Calcium (at. no., 20; at. wt., 40.08) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	Description	State	At 298.16°K (25°C)					
				$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole	
CaPb			c		-25.				
CaPb <sub>3</sub>			c		-26.				
Ca <sub>2</sub> Pb			c		-47.				
CaI <sub>2</sub> · 2PbI <sub>2</sub>			c		-223.1				
CaI <sub>2</sub> · 2PbI <sub>2</sub> · 7H <sub>2</sub> O			c		-712.6				
CaTi			c		-35.				
CaZn			c		-17.4				
CaZn <sub>2</sub>			c		-22.				
CaZn <sub>5</sub>			c		-33.				
CaZn <sub>13</sub>			c		-57.				
Ca <sub>4</sub> Zn			c		-32.				
CaCd <sub>3</sub>			c		-30.				



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 87-14. Calcium (at. no., 20; at. wt., 40.08) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$
CaHgBr <sub>4</sub> CaHg <sub>2</sub> Br <sub>6</sub>	in 5000 H <sub>2</sub> O in 9000 H <sub>2</sub> O		aq			-226.9 -264.3			
CaHg(CN) <sub>4</sub> CaHg <sub>2</sub> (CN) <sub>6</sub>	in 700 H <sub>2</sub> O in 1000 H <sub>2</sub> O		aq			-4.2 60.2			
CaCl <sub>2</sub> ·2Hg(CN) <sub>2</sub> CaCl <sub>2</sub> ·2Hg(CN) <sub>2</sub> ·6H <sub>2</sub> O	in 1000 H <sub>2</sub> O		aq			-77.5 -501.0			
CaBr <sub>2</sub> ·2Hg(CN) <sub>2</sub> CaBr <sub>2</sub> ·2Hg(CN) <sub>2</sub> ·7H <sub>2</sub> O	in 1000 H <sub>2</sub> O		aq			-56.0 -553.1			
CaI <sub>2</sub> ·2Hg(CN) <sub>2</sub> CaI <sub>2</sub> ·2Hg(CN) <sub>2</sub> ·7H <sub>2</sub> O	in 1000 H <sub>2</sub> O		aq			-29.3 -529.0			
CaAg(CN) <sub>3</sub> CaAg <sub>2</sub> (CN) <sub>4</sub>	in 1200 H <sub>2</sub> O in 550 H <sub>2</sub> O		aq			-30.1 -1.6			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 87-15. Calcium (at. no., 20; at. wt., 40.08)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Formula	Substance	State	Description	At 0°K				At 298.16°K (25°C)			
				ΔH <sup>o</sup> kcal/mole		ΔF <sup>o</sup> kcal/mole		log <sub>10</sub> K <sup>o</sup>		S <sup>o</sup> cal/deg mole	
CaO·Fe <sub>2</sub> O <sub>3</sub>		c				-398.					
CaH <sub>2</sub> Fe(CN) <sub>6</sub>		aq				-2.9					
Ca <sub>2</sub> Fe(CN) <sub>6</sub>		aq				-132.8					
Ca <sub>2</sub> Fe(CN) <sub>6</sub> ·11H <sub>2</sub> O		c				-886.6					
CaCrO <sub>4</sub>		c				-329.6	-305.3	223.78		32.	
		aq				-336.0					
CaWO <sub>4</sub>		c				-392.5					
CaTiO <sub>3</sub>		c								22.4	23.34
CaO·B <sub>2</sub> O <sub>3</sub>		c				-483.3	-457.7	335.49		25.1	24.85
CaO·2B <sub>2</sub> O <sub>3</sub>		c				-798.8	-752.4	551.50		32.2	37.75
2CaO·B <sub>2</sub> O <sub>3</sub>		gls				-786.2					
2CaO·3B <sub>2</sub> O <sub>3</sub>		c				-651.6	-618.6	453.42		34.7	35.16
2CaO·3B <sub>2</sub> O <sub>3</sub>		c				-1280.					
2CaO·3B <sub>2</sub> O <sub>3</sub> ·13H <sub>2</sub> O		c				-2211.					
3CaO·B <sub>2</sub> O <sub>3</sub>	inyoite	c				-817.7	-777.1	569.60		43.9	44.90

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 87-16. Calcium (at. no., 20; at. wt., 40.08) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)									
March 31, 1950									
Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$		$\log_{10} K_f$		$C_p^\circ$
			At 0°K	At 298.16°K (25°C)	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
CaAl <sub>3</sub>		c							
CaO·Al <sub>2</sub> O <sub>3</sub>		gls							
2CaO·Al <sub>2</sub> O <sub>3</sub>		c							
		gls							
2CaO·Al <sub>2</sub> O <sub>3</sub> ·5H <sub>2</sub> O		c							
3CaO·Al <sub>2</sub> O <sub>3</sub>		c							
		gls							
3CaO·Al <sub>2</sub> O <sub>3</sub> ·6H <sub>2</sub> O		c							
4CaO·Al <sub>2</sub> O <sub>3</sub>		c							
12CaO·7Al <sub>2</sub> O <sub>3</sub>		c							
		gls							
3CaCl <sub>2</sub> ·4AlCl <sub>3</sub>		c							
CaO·Al <sub>2</sub> O <sub>3</sub> ·6SiO <sub>2</sub>		c							
3CaO·Al <sub>2</sub> O <sub>3</sub> ·2SiO <sub>2</sub>		c							
4CaO·Al <sub>2</sub> O <sub>3</sub> ·Fe <sub>2</sub> O <sub>3</sub>		c							
CaMg <sub>2</sub>		c							
CaCl <sub>2</sub> ·2MgCl <sub>2</sub> ·2H <sub>2</sub> O		c							
CaCO <sub>3</sub> ·MgCO <sub>3</sub>		c							

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 88-1. Strontium (at. no., 38; at. wt., 87.63) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K	At 298.16°K (25°C)				
	kcal/mole			kcal/mole	kcal/mole	cal/deg mole	cal/deg mole		
Sr			g		39.2	26.3	-19.28	39.325	4.9680
			c	0.000	0.000	0.000	0.0000	13.0	6.0
		anal gas	liq		-50.4				
Sr <sup>+</sup>			g		171.96				
	Sr <sup>++</sup>		g		427.75				
			aq		-130.38	-133.2	97.49	-9.4	
SrO			c		-141.1	-133.8	98.07	13.0	10.76
	SrO <sub>2</sub>		c		-153.6				
			c		-722.6				
			c		-154.0				
SrH			g		52.4	45.8	-33.57	49.43	7.179
	SrH <sub>2</sub>		c		-42.3				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 68-2. Strontium (at. no., 38; at. wt., 87.63) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	Description	State	At 298.16°K (25°C)						
			At 0°K kcal/mole	kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
Sr(OH) <sub>2</sub>	std. state, hyp. m=1 in 100 H <sub>2</sub> O	c							
		aq		-229.3	-208.2	152.61	-14.4		
		aq		-240.29					
Sr(OH) <sub>2</sub> ·H <sub>2</sub> O		c		-240.11					
		c		-302.3					
		c		-801.2					
SrF <sub>2</sub>		g		-5.					
		c		-290.3					
SrCl <sub>2</sub>		g		9.					
		c		-198.0	-186.7	136.85	28.	18.3	
Sr(OH) <sub>2</sub> ·8H <sub>2</sub> O	std. state, hyp. m=1 in 100 H <sub>2</sub> O	aq		-210.43	-195.7	143.45	16.9		
		aq		-209.70					
		aq		-209.83					
		aq		-209.9					
		aq		-209.97					
		aq		-210.06					
		aq		-210.10					
		aq		-210.22					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 88-3. Strontium (at. no., 38; at. wt., 87.63)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		Description	State	ΔH <sup>o</sup>		ΔF <sup>o</sup>	log <sub>10</sub> K <sub>f</sub>	S <sup>o</sup>	C <sub>p</sub> <sup>o</sup>
Formula	in			At 0°K	At 298.16°K (25°C)				
				kcal/mole	kcal/mole				
SrCl <sub>2</sub>	in 1000 H <sub>2</sub> O		aq		-210.27				
	5000 H <sub>2</sub> O		aq		-210.35				
	10000 H <sub>2</sub> O		aq		-210.37				
	50000 H <sub>2</sub> O		aq		-210.41				
	∞ H <sub>2</sub> O		aq		-210.43				
SrCl <sub>2</sub> ·H <sub>2</sub> O			c		-271.7				28.7
	SrCl <sub>2</sub> ·2H <sub>2</sub> O		c		-343.7				38.3
	SrCl <sub>2</sub> ·6H <sub>2</sub> O		c		-627.1				
SrOCl <sub>2</sub>			aq		-195.8				
	Sr(OCl) <sub>2</sub>		aq		-176.5				
SrCl <sub>2</sub> ·SrO·H <sub>2</sub> O			c		-437.4				
	SrCl <sub>2</sub> ·SrO·3H <sub>2</sub> O		c		-1006.8				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 88-4. Strontium (at. no., 38; at. wt., 87.63)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole				
SrBr <sub>2</sub>	std. state, hyp. m=1		c						
	in 100 H <sub>2</sub> O		aq		-171.1				19.0
	200 H <sub>2</sub> O		aq		-188.2		133.48	29.2	
	500 H <sub>2</sub> O		aq		-187.56				
	1000 H <sub>2</sub> O		aq		-187.65				
	1500 H <sub>2</sub> O		aq		-187.78				
	5000 H <sub>2</sub> O		aq		-187.86				
	10000 H <sub>2</sub> O		aq		-187.90				
	50000 H <sub>2</sub> O		aq		-188.01				
	100000 H <sub>2</sub> O		aq		-188.05				
	500000 H <sub>2</sub> O		aq		-188.12				
	∞ H <sub>2</sub> O		aq		-188.15				
			aq		-188.18				
SrBr <sub>2</sub> ·H <sub>2</sub> O SrBr <sub>2</sub> ·6H <sub>2</sub> O  SrBr <sub>2</sub> ·SrO·3H <sub>2</sub> O SrBr <sub>2</sub> ·SrO·9H <sub>2</sub> O			c						28.9
			c		-246.8				82.1
			c		-604.4				
			c		-552.8				
			c		-964.8				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 88-5. Strontium (at. no., 38; at. wt., 87.63) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950										
Substance		State	Description	$\Delta H_f^0$ kcal/mole	$\Delta H_f^0$ kcal/mole	$\Delta F_f^0$ kcal/mole	$\log_{10} K_f$	$S^0$	$C_p^0$	
Formula	State									
SrI <sub>2</sub>	c	std. state, hyp. m = 1 in 1000 H <sub>2</sub> O			-135.5	-157.7	115.59	42.9	19.5	
										aq
										aq
SrI <sub>2</sub> ·H <sub>2</sub> O	c				-212.2				28.5	
										c
										c
SrI <sub>2</sub> ·2H <sub>2</sub> O	c				-282.8				39.1	
										c
										c
SrI <sub>2</sub> ·6H <sub>2</sub> O	c				-571.2				84.9	
										c
										c
SrS	g				19.					
										c
										c
SrSO <sub>4</sub>	c	std. state, hyp. m = 1			-345.3	-318.9	233.75	29.1		
										aq
										aq
SrS <sub>2</sub> O <sub>6</sub>	aq				-410.8					
										c
										c
SrS <sub>2</sub> O <sub>6</sub> ·4H <sub>2</sub> O	c				-693.1					



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 88-6. Strontium (at. no., 38; at. wt., 87.63) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
$\text{SrI}_2 \cdot 2\text{SO}_2$			c		-299.5				
$\text{SrI}_2 \cdot 4\text{SO}_2$			c		-462.9				
SrSe			c		-78.7				
$\text{Sr}(\text{N}_3)_2$			c		48.9				
$\text{Sr}_3\text{N}_2$			c		-93.4				
$\text{SrN}_2\text{O}_2 \cdot 5\text{H}_2\text{O}$			c		-469.9				
$\text{Sr}(\text{NO}_2)_2$			c, II		-179.3				
			aq		-180.4				
$\text{Sr}(\text{NO}_3)_2$			c		-233.25				36.3
	std. state, hyp. m=1		aq		-229.02	-185.8	136.19	60.6	
	in 20 $\text{H}_2\text{O}$		aq		-231.950				
	50 $\text{H}_2\text{O}$		aq		-230.582				
	100 $\text{H}_2\text{O}$		aq		-229.736				
	200 $\text{H}_2\text{O}$		aq		-229.179				
	400 $\text{H}_2\text{O}$		aq		-228.895				
	500 $\text{H}_2\text{O}$		aq		-228.873				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 88-7. Strontium (at. no., 38; at. wt., 87.63)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		State	Description	log <sub>10</sub> K <sub>f</sub>	S <sup>o</sup>	C <sub>p</sub> <sup>o</sup>	
Formula	ΔF <sup>o</sup> kcal/mole						ΔF <sup>o</sup> kcal/mole
Sr(NO <sub>3</sub> ) <sub>2</sub>	in	aq					
	800 H <sub>2</sub> O	aq	-228.843				
	1000 H <sub>2</sub> O	aq	-228.838				
	1500 H <sub>2</sub> O	aq	-228.836				
	2000 H <sub>2</sub> O	aq	-228.840				
	3000 H <sub>2</sub> O	aq	-228.867				
	4000 H <sub>2</sub> O	aq	-228.894				
	5000 H <sub>2</sub> O	aq	-228.950				
	10000 H <sub>2</sub> O	aq	-228.968				
	500000 H <sub>2</sub> O	aq	-228.993				
Sr(NO <sub>3</sub> ) <sub>2</sub> ·4H <sub>2</sub> O	∞ H <sub>2</sub> O	ad	-229.02				
		c	-514.5				
		c	-83.2				
Sr(NH <sub>2</sub> ) <sub>2</sub>		c	-123.8				
		c					
SrCl <sub>2</sub> ·NH <sub>3</sub>		c	-220.5				
	SrCl <sub>2</sub> ·8NH <sub>3</sub>	c	-368.5				

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Table 88-8. Strontium (at. no., 38; at. wt., 87.63) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		State	$\Delta H^{\circ}$ kcal/mole	$\Delta F^{\circ}$ kcal/mole	$\log_{10} K_f$	$S^{\circ}$	$C_p^{\circ}$		
Formula	Description							At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole
$\text{SrBr}_2 \cdot \text{NH}_3$		c		-136.9					
$\text{SrBr}_2 \cdot 2\text{NH}_3$		c		-222.8					
$\text{SrBr}_2 \cdot 8\text{NH}_3$		c		-354.4					
$\text{SrI}_2 \cdot \text{NH}_3$		c		-164.8					
$\text{SrI}_2 \cdot 2\text{NH}_3$		c		-191.4					
$\text{SrI}_2 \cdot 6\text{NH}_3$		c		-285.9					
$\text{SrI}_2 \cdot 8\text{NH}_3$		c		-330.0					
$(\text{NH}_4)_2\text{SO}_4 \cdot \text{SrSO}_4$		c		-628.8					
$\text{Sr}_3(\text{PO}_4)_2$		c		-987.3					
$\text{SrHPO}_4$		c		-431.3					
$\text{Sr}(\text{H}_2\text{PO}_4)_2 \cdot \text{H}_2\text{O}$		c		-819.4					
$\text{Sr}_3(\text{AsO}_4)_2$		c		-800.7					
$\text{SrHAsO}_4$		aq		-344.8					
$\text{Sr}(\text{H}_2\text{AsO}_4)_2$		aq		-561.4					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 88-9. Strontium (at. no. 38):  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION,  
AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance									
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
			At 0°K	At 298.16°K (25°C)					
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole		
$\text{SrCO}_3$	strontianite	c, II		-291.2	-271.9	199.30	23.2	19.46	
$\text{SrC}_2\text{O}_4$	std. state, hyp. m = 1	aq		-327.4	-294.3	215.72	3		
$\text{SrC}_2\text{O}_4 \cdot \text{H}_2\text{O}$		c			-360.8	264.46			
$\text{SrC}_2\text{O}_4 \cdot 2\frac{1}{2}\text{H}_2\text{O}$		c		-504.2					
$\text{Sr}(\text{CHO}_2)_2$	strontium formate	c		-325.5					
		aq		-326.4					
$\text{Sr}(\text{CHO}_2)_2 \cdot 2\text{H}_2\text{O}$		c		-468.3					
$\text{Sr}(\text{HCO}_3)_2$	std. state, hyp. m = 1	aq		-460.7	-413.8	303.31	36.0		
$\text{Sr}(\text{C}_2\text{H}_3\text{O}_2)_2$	strontium acetate	c		-356.7					
		aq		-362.1					
$\text{Sr}(\text{C}_2\text{H}_3\text{O}_2)_2 \cdot \frac{1}{2}\text{H}_2\text{O}$		c		-391.2					
$\text{Sr}(\text{C}_2\text{H}_3\text{O}_3)_2$	strontium glycolate	c		-440.7					
		aq		-440.5					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 88-10. Strontium (at. no., 38; at. wt., 87.63) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		Description	State	Δ <i>F</i> °		log <sub>10</sub> <i>F</i> °	<i>S</i> °	<i>C</i> <sub><i>p</i></sub> °	
Formula				At 0°K kcal/mole	kcal/mole				
SrBr <sub>2</sub> ·½C <sub>2</sub> H <sub>5</sub> OH			c		-204.5				
Sr(CN) <sub>2</sub>			aq		-57.8				
Sr(CN) <sub>2</sub> ·4H <sub>2</sub> O			c		-335.2				
SrSi			c		-113.				
SrSi <sub>2</sub>			c		-150.				
SrSiO <sub>3</sub>			c		-371.2				
Sr <sub>2</sub> SiO <sub>4</sub>			gls		-367.4				
			c		-520.6				
SrI <sub>2</sub> ·2PbI <sub>2</sub>			c		-222.0				
SrI <sub>2</sub> ·2PbI <sub>2</sub> ·7H <sub>2</sub> O			c		-713.7				
SrHgBr <sub>4</sub>			aq		-227.5				
SrHg <sub>2</sub> Br <sub>6</sub>			aq		-265.0				
Sr <sub>2</sub> HgBr <sub>6</sub>			aq		-417.4				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 86-11. Strontium (at. no., 38; at. wt., 87.63) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				
$\text{SrHg}(\text{CN})_4$			aq		-4.3				
$\text{SrHg}_2(\text{CN})_6$			aq		60.2				
$\text{SrCl}_2 \cdot 2\text{Hg}(\text{CN})_2$			aq		-78.2				
$\text{SrCl}_2 \cdot 2\text{Hg}(\text{CN})_2 \cdot 6\text{H}_2\text{O}$			c		-504.6				
$\text{SrBr}_2 \cdot 2\text{Hg}(\text{CN})_2$			aq		-56.8				
$\text{SrBr}_2 \cdot 2\text{Hg}(\text{CN})_2 \cdot 6\text{H}_2\text{O}$			c		-486.3				
$\text{SrI}_2 \cdot 2\text{Hg}(\text{CN})_2$			aq		-30.1				
$\text{SrI}_2 \cdot 2\text{Hg}(\text{CN})_2 \cdot 7\text{H}_2\text{O}$			c		-531.1				
$\text{SrAg}(\text{CN})_3$			aq		-40.1				
$\text{SrAg}_2(\text{CN})_4$			aq		-1.1				
$\text{SrNi}(\text{CN})_4$			aq		-43.5				
$\text{Sr}_3[\text{FeCo}(\text{CN})_5]_2$			c		-257.5				
$\text{Sr}_3[\text{FeCo}(\text{CN})_5]_2 \cdot 4\text{H}_2\text{O}$			aq		-299.0				
$\text{SrWO}_4$			c		-538.7				
$3\text{SrCl}_2 \cdot 4\text{AlCl}_3$			c		-398.3				
					-1275.6				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 89-1. Barium (at. no., 56; at. wt., 137.36)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance									
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
			At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C)				
Ba		g c, II	0.000	41.96		-25.361	40.699	4.9680	
				0.000		0.0000	16.	6.30	
				163.582					
				395.705					
BaO	std. state, hyp. m = 1	aq		-128.67	-134.0	98.22	3.		
		g				55.70	7.38		
BaO <sub>2</sub>		c		-133.4	-126.3	92.58	16.8	11.34	
				-150.5					
				-223.5					
BaO <sub>2</sub> ·8H <sub>2</sub> O		c		-719.3					
Ba <sub>2</sub> O		c		-147.					
BaH		g		52.	4c.	-33.7	52.97	7.001	
		c		-40.9					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 89-2. Barium (at. no., 56; at. wt., 137.36) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance									
Formula	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 0°K kcal/mole	kcal/mole					
Ba(OH) <sub>2</sub>		c		-226.2					
	std. state, hyp. m = 1	aq		-238.58	-209.2	153.34	-2.		
Ba(OH) <sub>2</sub> •H <sub>2</sub> O		aq		-238.5					
Ba(OH) <sub>2</sub> •8H <sub>2</sub> O		c		-299.0					
	in 600 H <sub>2</sub> O	c		-799.5					
BaO <sub>2</sub> •H <sub>2</sub> O <sub>2</sub>		c		-209.6					
BaF		g		-9.					
BaF <sub>2</sub>		c		-286.9	-274.5	201.20	23.0	17.02	
		aq		-286.0					
BaCl		g		24.					
		c		-111.					
BaCl <sub>2</sub>		c		-205.56	-193.8	142.05	30.	18.0	
	std. state, hyp. m = 1	aq		-208.72	-196.7	144.18	29.		



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 89-3. Barium (at. no., 56; at. wt., 137.36) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				
BaCl <sub>2</sub>	in	31.1 H <sub>2</sub> O	aq		-208.072				
	50	H <sub>2</sub> O	aq		-208.085				
	75	H <sub>2</sub> O	aq		-208.097				
	100	H <sub>2</sub> O	aq		-208.110				
	150	H <sub>2</sub> O	aq		-208.135				
	200	H <sub>2</sub> O	aq		-208.162				
	300	H <sub>2</sub> O	aq		-208.2				
	400	H <sub>2</sub> O	aq		-208.231				
	500	H <sub>2</sub> O	aq		-208.257				
	700	H <sub>2</sub> O	aq		-208.306				
	1000	H <sub>2</sub> O	aq		-208.338				
	1200	H <sub>2</sub> O	aq		-208.373				
	1500	H <sub>2</sub> O	aq		-208.388				
	2000	H <sub>2</sub> O	aq		-208.419				
	3000	H <sub>2</sub> O	aq		-208.455				
	5000	H <sub>2</sub> O	aq		-208.496				
	7500	H <sub>2</sub> O	aq		-208.526				
	10000	H <sub>2</sub> O	aq		-208.544				
	20000	H <sub>2</sub> O	aq		-208.584				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

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Table 89-4. Barium (at. no., 56; at. wt., 137.36)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Formula	Substance	Description	State	At 298.16°K (25°C)				
				$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
BaCl <sub>2</sub>	in 30000 H <sub>2</sub> O	50000 H <sub>2</sub> O	aq		-208.604			
		10000 H <sub>2</sub> O	aq		-208.627			
		10000 H <sub>2</sub> O	aq		-208.650			
		20000 H <sub>2</sub> O	aq		-208.669			
		50000 H <sub>2</sub> O	aq		-208.686			
		∞ H <sub>2</sub> O	aq		-208.72			
BaCl <sub>2</sub> ·H <sub>2</sub> O			c		-278.4	-253.1	40.	28.2
BaCl <sub>2</sub> ·2H <sub>2</sub> O			c		-349.35	-309.7	48.5	37.10
BaOCl <sub>2</sub>			aq		-194.0			
Ba(OCl) <sub>2</sub>			aq		-176.1			
Ba(ClO <sub>2</sub> ) <sub>2</sub>			c		-158.2			
Ba(ClO <sub>3</sub> ) <sub>2</sub>			c		-181.7			
			aq		-175.6			
Ba(ClO <sub>3</sub> ) <sub>2</sub> ·H <sub>2</sub> O			c		-254.9			
Ba(ClO <sub>4</sub> ) <sub>2</sub>			c		-192.8			
			aq		-191.5			
Ba(ClO <sub>4</sub> ) <sub>2</sub> ·3H <sub>2</sub> O			c		-405.4			
								50.6

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 89-5. Barium (at. no., 56; at. wt., 137.36) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	Description		At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole					
$\text{BaCl}_2 \cdot \text{BaO} \cdot 3\text{H}_2\text{O}$	std. state, hyp. m = 1	c	-545.5						
$\text{BaCl}_2 \cdot \text{BaO} \cdot 5\text{H}_2\text{O}$		c	-727.5						
$\text{BaCl}_2 \cdot \text{BaO} \cdot 6\text{H}_2\text{O}$		c	-931.7						
$\text{BaBr}_2$		c	-180.4						
		aq	-186.47	-183.1		134.21	42.		
		aq	-186.047						
		aq	-186.074						
		aq	-186.093						
		aq	-186.123						
		aq	-186.153						
		aq	-186.186						
		aq	-186.211						
		aq	-186.243						
		aq	-186.281						
		aq	-186.326						
		aq	-186.364						
		aq	-186.399						
		aq	-186.417						
		aq	-186.47						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards Washington, D.C.

Table 89-6. Barium (at. no., 56; at. wt., 137.36) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	Description	State	$\Delta H^{\circ}$		$\Delta F^{\circ}$	$\log_{10} K_f$	$S^{\circ}$	$C_p^{\circ}$
				At 0°K kcal/mole	kcal/mole				
BaBr <sub>2</sub> ·H <sub>2</sub> O			c		-254.9				
BaBr <sub>2</sub> ·2H <sub>2</sub> O			-c		-326.3				
BaOBr <sub>2</sub>			aq		-181.6				
Ba(BrO <sub>3</sub> ) <sub>2</sub> ·H <sub>2</sub> O			c					68.8	52.90
BaBr <sub>2</sub> ·BaO·2H <sub>2</sub> O			c		-486.3				
BaBr <sub>2</sub> ·BaO·5H <sub>2</sub> O			c		-705.7				
BaI <sub>2</sub>			c		-144.0				
		std. state, hyp. m = 1	aq		-155.41	-158.7	116.32	55.	
		in 400 H <sub>2</sub> O	aq		-155.15				
		10/10 H <sub>2</sub> O	aq		-155.28				
BaI <sub>2</sub> ·H <sub>2</sub> O			c		-219.8				
BaI <sub>2</sub> ·2H <sub>2</sub> O			c		-290.9				
BaI <sub>2</sub> ·2½H <sub>2</sub> O			c		-326.0				
BaI <sub>2</sub> ·7H <sub>2</sub> O			c		-640.1				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 89-7. Barium (at. no., 56; at. wt., 137.36)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)		cal/deg mole	cal/deg mole
			kcal/mole	kcal/mole			
$Ba(IO_3)_2$		aq		-238.4			
$Ba(IO_3)_2 \cdot H_2O$		c		-319.6			
$BaI_2 \cdot BaO \cdot 2H_2O$		c		-455.9			
$BaI_2 \cdot BaO \cdot 9H_2O$		c		-956.0			
$BaS$		g		41.			
		c		-106.0			
		aq		-118.4			
$BaS_2O_3$		c		-282.6			
$BaS_2O_4$		c		-350.2	237.05	31.6	24.32
	std. state, hyp. m=1	aq		-345.57	228.18	7.	
$BaS_2O_3$		c					40.7
$BaS_2O_6$		aq		-409.3			
$BaS_2O_6 \cdot 2H_2O$		c		-552.5			
$BaS_2O_8$		aq		-454.0			
$BaS_2O_8 \cdot 4H_2O$		c		-738.7			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 89-8. Barium (at. no., 56; at. wt., 137.36) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	At 298.16°K (25°C)				
Formula				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	
BaSi <sub>2</sub> O <sub>6</sub>			aq		-419.				
BaSi <sub>2</sub> O <sub>6</sub> · 2H <sub>2</sub> O			c		-572.2				
Ba(HS) <sub>2</sub>			aq		-134.8				
Ba(HSO <sub>3</sub> ) <sub>2</sub>			aq		-430.7				
BaSO <sub>4</sub> · H <sub>2</sub> SO <sub>4</sub>			c		-549.1				
BaSO <sub>4</sub> · 2H <sub>2</sub> SO <sub>4</sub> · H <sub>2</sub> O			c		-821.0				
BaI <sub>2</sub> · 2SO <sub>2</sub>			c		-308.6				
BaI <sub>2</sub> · 4SO <sub>2</sub>			c		-470.3				
BaSe			c		-74.2				
BaSeO <sub>4</sub>			c		-280.0				
Ba(N <sub>3</sub> ) <sub>2</sub>			c		-8.0				
Ba <sub>3</sub> N <sub>2</sub>			aq		0.2				
			c		-86.9				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 89-9. Barium (at. no., 56; at. wt., 137.36)									
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)									
March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	kcal/mole				
$Ba(NO_2)_2$			c						
	in 80% $H_2O$		aq		-174.0				
			c		-178.6				
$Ba(NO_2)_2 \cdot H_2O$					-254.5				
			c		-237.06				
	std. state, hyp. m = 1		aq		-227.41		139.27	51.1	36.1
$Ba(NO_3)_2$	in 1% $H_2O$		aq		-226.27		136.92	73.	
	20% $H_2O$		aq		-226.57				
	25% $H_2O$		aq		-226.74				
	30% $H_2O$		aq		-226.86				
	40% $H_2O$		aq		-227.035				
	50% $H_2O$		aq		-227.104				
	60% $H_2O$		aq		-227.181				
	70% $H_2O$		aq		-227.229				
	80% $H_2O$		aq		-227.266				
	100% $H_2O$		aq		-227.317				
	120% $H_2O$		aq		-227.350				
	150% $H_2O$		aq		-227.381				
	180% $H_2O$		aq		-227.400				
	200% $H_2O$		aq		-227.415				

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Table 89-10. Barium (at. no., 56; at. wt., 137.36) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole			At 298.16°K (25°C) kcal/mole	cal/deg mole
Ba(NO <sub>3</sub> ) <sub>2</sub>	in	2500 H <sub>2</sub> O	aq		-227.449				
		3000 H <sub>2</sub> O	aq		-227.464				
		4000 H <sub>2</sub> O	aq		-227.476				
		5000 H <sub>2</sub> O	aq		-227.480				
		6000 H <sub>2</sub> O	aq		-227.484				
		7500 H <sub>2</sub> O	aq		-227.487				
		10000 H <sub>2</sub> O	aq		-227.488				
		15000 H <sub>2</sub> O	aq		-227.485				
		20000 H <sub>2</sub> O	aq		-227.482				
		30000 H <sub>2</sub> O	aq		-227.477				
		50000 H <sub>2</sub> O	aq		-227.468				
		100000 H <sub>2</sub> O	aq		-227.456				
		200000 H <sub>2</sub> O	aq		-227.445				
		500000 H <sub>2</sub> O	aq		-227.434				
		$\infty$ H <sub>2</sub> O	aq		-227.41				
BaNH			c		-53.8				
Ba(NH <sub>2</sub> ) <sub>2</sub>			c		-78.9				
Ba(NH <sub>2</sub> ) <sub>3</sub>			c		-124.1				



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 89-11. Barium (at. no., 56; at. wt., 137.36) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K					
				kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
$\text{BaCl}_2 \cdot 8\text{NH}_3$			c		-365.0				
$\text{BaBr}_2 \cdot \text{NH}_3$			c		-203.3				
$\text{BaBr}_2 \cdot 2\text{NH}_3$			c		-225.0				
$\text{BaBr}_2 \cdot 4\text{NH}_3$			c		-267.5				
$\text{BaBr}_2 \cdot 8\text{NH}_3$			c		-351.4				
$\text{BaI}_2 \cdot 2\text{NH}_3$			c		-193.1				
$\text{BaI}_2 \cdot 4\text{NH}_3$			c		-237.8				
$\text{BaI}_2 \cdot 6\text{NH}_3$			c		-282.0				
$\text{BaI}_2 \cdot 8\text{NH}_3$			c		-325.5				
$\text{BaI}_2 \cdot 9\text{NH}_3$			c		-346.5				
$\text{BaI}_2 \cdot 10\text{NH}_3$			c		-365.1				
$\text{Ba}_3(\text{PO}_4)_2$		colloidal	c		-998.0				
$\text{BaHPO}_4$			aq		-981.4				
$\text{Ba}(\text{H}_2\text{PO}_4)_2$			c		-465.8				
$\text{Ba}(\text{H}_2\text{PO}_4)_2 \cdot \text{H}_2\text{O}$		in 400 $\text{H}_2\text{O}$	aq		-423.6				
$\text{Ba}(\text{H}_2\text{PO}_4)_2$			c		-492.0				
$\text{Ba}(\text{H}_2\text{PO}_4)_2$			c		-749.6				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 89-12. Barium (at. no., 56; at. wt., 137.36)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	Description	ΔH <sup>o</sup>		log <sub>10</sub> K <sup>o</sup>	S <sup>o</sup>	C <sub>p</sub> <sup>o</sup>
Formula	ΔH <sup>o</sup>			At 298.16°K (25°C)				
				At 0°K				
Ba <sub>3</sub> (AsO <sub>4</sub> ) <sub>2</sub>		c						
BaHAsO <sub>4</sub> ·H <sub>2</sub> O		c						
Ba(H <sub>2</sub> AsO <sub>4</sub> ) <sub>2</sub> ·2H <sub>2</sub> O		c						
Ba <sub>3</sub> Sb <sub>2</sub>		c						
Ba <sub>3</sub> Bi <sub>2</sub>		c						
BaCO <sub>3</sub>	witherite	c. II						
	std. state, hyp. m=1	aq						
BaC <sub>2</sub> O <sub>4</sub> ·½H <sub>2</sub> O	barium oxalate	c						
BaC <sub>2</sub> O <sub>4</sub> ·2H <sub>2</sub> O		c						
BaC <sub>2</sub> O <sub>4</sub> ·3½H <sub>2</sub> O		c						
BaO·BaCO <sub>3</sub>		c						
Ba(CHO <sub>2</sub> ) <sub>2</sub>	barium formate	c						
	in 400 H <sub>2</sub> O	aq						
Ba(HCO <sub>3</sub> ) <sub>2</sub>	std. state, hyp. m=1	aq						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 89-13. Barium (at. no., 56; at. wt., 137.36)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Formula	Substance	State	Description	$\Delta H_f^\circ$		$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole			
$Ba(C_2H_3O_2)_2$	barium acetate in 400 $H_2O$	c			-355.1			
$Ba(C_2H_3O_2) \cdot 3H_2O$		aq			-361.5			
		c			-567.3			
$Ba(C_2H_3O_2)_2$	barium glycolate	c			-443.6			
		aq			-439.1			
$Ba(OC_2H_5)_2$	barium ethylate	c			-219.1			
$3BaO \cdot 4CH_3OH$		c			-713.1			
$3BaO \cdot 4C_2H_5OH$		c			-736.7			
$Ba(HSO_3)_2 \cdot C_2H_2O_2$	barium bisulfite-glyoxal	aq			-543.1			
$Ba(HSO_3)_2 \cdot C_2H_2O_2 \cdot 2\frac{1}{2}H_2O$		c			-722.1			
$Ba(OSO_3C_2H_5)_2$	barium ethyl sulfate	aq			-546.4			
$Ba(OSO_3C_2H_5)_2 \cdot 2H_2O$		c			-687.5			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 89-14. Barium (at. no., 56; at. wt., 137.36)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance										
Formula		Description		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
					At 0°K	At 298.16°K (25°C)				
					kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole
BaCN <sub>2</sub>			barium cyanamide	c		-63.8				
Ba(CN) <sub>2</sub>				c		-47.9				
				aq		-50.7				
Ba(CN) <sub>2</sub> ·H <sub>2</sub> O				c		-120.1				
Ba(CN) <sub>2</sub> ·2H <sub>2</sub> O				c		-191.1				
Ba(CNO) <sub>2</sub>			barium cyanate	c		-209.6				
				aq		-195.7				
BaSi <sub>3</sub>				c		-407.				
Ba <sub>2</sub> Si <sub>1/2</sub>				c		-367.				
BaSiO <sub>3</sub>				c		-359.5				
Ba <sub>2</sub> SiO <sub>4</sub>				c		-496.8				
BaSiF <sub>6</sub>				c		-691.6				
BaSn <sub>3</sub>				c		-44.				
Ba <sub>2</sub> Sn				c		-90.				

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Table 89-15. Barium (at. no., 56; at. wt., 137.36) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 0°K kcal/mole	kcal/mole					
Formula	Description								
BaPb									
BaPb <sub>2</sub>									
Ba <sub>2</sub> Pb									
BaI <sub>2</sub> · 2PbI <sub>2</sub>									
BaI <sub>2</sub> · 2PbI <sub>2</sub> · 7H <sub>2</sub> O									
BaHgBr <sub>4</sub>		aq							
BaHg <sub>2</sub> Br <sub>6</sub>		aq							
Ba <sub>2</sub> HgBr <sub>6</sub>		aq							
Ba <sub>4</sub> HgBr <sub>10</sub>		aq							
BaHg(CN) <sub>4</sub>		aq							
BaHg <sub>2</sub> (CN) <sub>6</sub>		aq							
BaCl <sub>2</sub> · Hg(CN) <sub>2</sub>		aq							
BaCl <sub>2</sub> · 2Hg(CN) <sub>2</sub> · 2H <sub>2</sub> O		c							
BaBr <sub>2</sub> · 2Hg(CN) <sub>2</sub>		aq							
BaBr <sub>2</sub> · 2Hg(CN) <sub>2</sub> · 2H <sub>2</sub> O		c							
BaI <sub>2</sub> · Hg(CN) <sub>2</sub>		aq							
PbI <sub>2</sub> · 2Hg(CN) <sub>2</sub> · (H <sub>2</sub> O)		c							

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 89-16. Barium (at. no., 56; at. wt., 137.36) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		State	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	Description								
BaAg(CN) <sub>3</sub>		aq		-24.3					
BaAg <sub>2</sub> (CN) <sub>4</sub>		aq		6.3					
BaPtCl <sub>6</sub>		c		-286.8					
BaPtCl <sub>2</sub> ·6H <sub>2</sub> O		aq		-296.1					
BaOsCl <sub>6</sub>		c		-707.2					
BaPdCl <sub>4</sub>		c		-270.0					
Ba <sub>3</sub> (RhCl <sub>6</sub> ) <sub>2</sub>		c		-229.9					
		c		-687.					
BaNi(CN) <sub>4</sub>		aq		-41.7					
Ba <sub>2</sub> Fe(CN) <sub>6</sub>		aq		-129.7					
Ba <sub>2</sub> Fe(CN) <sub>6</sub> ·6H <sub>2</sub> O		c		-551.0					
BaH <sub>2</sub> Fe(CN) <sub>6</sub>		aq		-2.5					
Ba <sub>3</sub> [FeCo(CN) <sub>5</sub> ] <sub>2</sub>		c		-273.7					
		aq		-292.0					
Ba <sub>3</sub> [FeCo(CN) <sub>5</sub> ] <sub>2</sub> ·11H <sub>2</sub> O		c		-1050.2					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} f_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)			
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
BaMnO <sub>4</sub>		c		-282.			
BaCrO <sub>4</sub>		c		-341.3			
BaMoO <sub>4</sub>		c		-373.8			
BaWO <sub>4</sub>		c		-407.7			
BaCl <sub>2</sub> ·2AlCl <sub>3</sub>		c		-543.3			
3BaCl <sub>2</sub> ·4AlCl <sub>3</sub>		c		-1090.5			
							33.6

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 90-1. Radium (at. no., 88; at. wt., 226.05) HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		State	Description	At 298.16°K (25°C)					C <sub>p</sub> <sup>o</sup> cal/deg mole
Formula	•			ΔH <sub>f</sub> <sup>o</sup> kcal/mole	ΔF <sub>f</sub> <sup>o</sup> kcal/mole	log <sub>10</sub> K <sub>f</sub>	S <sup>o</sup> cal/deg mole		
								At 0°K kcal/mole	
Ra		g		31.	27.	-16.3	42.15	4.968	
Ra <sup>+</sup>		g		0.000	0.000	0.0000	17.		
Ra <sup>++</sup>		g		153.1					
		g		368.5					
		aq		-126.	-134.5	98.53	13.		
RaO		c		-125.					
RaCl <sub>2</sub>		c					32.		
RaCl <sub>2</sub> · 2H <sub>2</sub> O		c		-351.	-311.7	228.47	50.		
RaSO <sub>4</sub>		c		-352.	-326.0	238.95	34.		
Ra(NO <sub>3</sub> ) <sub>2</sub>		c		-237.	-190.3	139.49	52.		



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 91-1. Lithium (at. no., 3; at. wt., 6.940)								
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)								
March 31, 1950								
Substance			$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)				
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole
Li		g	36.49	37.07	29.19	-21.39e	33.143	4.9e80
Li <sup>+</sup>		c	0.000	0.000	0.000	0.0000	6.8	5.6F
Li <sup>+</sup>		g	160.799	162.860				
Li <sup>+</sup>	std. state, hyp. m = 1	aq		-66.554	-71.22	-1.471	3.4	
Li <sup>++</sup>		g	1904.7	1908.2				
Li <sup>+++</sup>		g	4721.4	4732.7				
Li <sub>2</sub>		g	47.2	47.6	37.6	-27.56	41.06	8.52
Li <sub>2</sub>				-142.4				
Li <sub>2</sub>				-151.7				
Li <sub>2</sub>		aq		-159.4				
LiH		g	0.	40.7	26.2	-16.47	42.7	7.2
		c		-21.61	-16.2	12.255	5.7	6.7

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 91-2. Lithium (at. no., 3; at. wt., 6.940)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Formula	Substance	Description	State	$\Delta F^{\circ}$		$\Delta F^{\circ}$	$\log_{10} K_f$	$S^{\circ}$	$C_p^{\circ}$
				At 0°K	kcal/mole	At 298.16°K (25°C)			
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
LiOH	std. state, hyp. m = 1		c						
	in 11 H <sub>2</sub> O		aq				77.77	12.	
	25 H <sub>2</sub> O		aq		-116.45	-106.1	79.031	0.9	
	50 H <sub>2</sub> O		aq		-121.511	-107.82			
	100 H <sub>2</sub> O		aq		-120.51				
	100 H <sub>2</sub> O		aq		-120.930				
	100 H <sub>2</sub> O		aq		-121.090				
	200 H <sub>2</sub> O		aq		-121.198				
	400 H <sub>2</sub> O		aq		-121.27				
	800 H <sub>2</sub> O		aq		-121.38				
	1600 H <sub>2</sub> O		aq		-121.481				
	3200 H <sub>2</sub> O		aq		-121.47				
	6400 H <sub>2</sub> O		aq		-121.34				
	$\infty$ H <sub>2</sub> O		aq		-121.464				
LiOH•H <sub>2</sub> O Li <sub>2</sub> O <sub>2</sub> •H <sub>2</sub> O <sub>2</sub> •2H <sub>2</sub> O			aq		-121.511		120.80	22.	
			c		-188.77	-164.8			
			c		-414.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 91-3. Lithium (at. no., 3; at. wt., 6.940) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950										
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
				At 0°K	At 298.16°K (25°C)					
				kcal/mole	kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
LiF			c							
LiHF <sub>2</sub>	std. state, hyp. m = 1 in 400 H <sub>2</sub> O		aq		-146.3	-139.6	102.32	8.57	10.04	
			aq		-145.210	-136.30	99.906	1.1		
LiCl	std. state, hyp. m = 1 in 5 H <sub>2</sub> O		g	-53.		-56.	42.5	51.01	7.88	
			c						12.2	
			aq		-97.70	-106.577	-101.57	74.449	16.6	
			aq		-102.62					
			aq		-103.73					
			aq		-104.431					
			aq		-105.310					
			aq		-105.554					
			aq		-105.711					
			aq		-105.848					
			aq		-105.987					
			aq		-106.064					
	aq		-106.221							

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table I-4. Lithium (at. no., 3; at. wt., 6.940)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Formula	Substance		Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
					kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
LiCl	in 100 H <sub>2</sub> O		aq		-106.315				
	200 H <sub>2</sub> O		aq		-106.380				
	400 H <sub>2</sub> O		aq		-106.430				
	800 H <sub>2</sub> O		aq		-106.467				
	1600 H <sub>2</sub> O		aq		-106.537				
	3200 H <sub>2</sub> O		aq		-106.518				
	6400 H <sub>2</sub> O		aq		-106.536				
	$\infty$ H <sub>2</sub> O		aq		-106.577				
	in 25 CH <sub>3</sub> OH				-108.28				
	50 CH <sub>3</sub> OH				-108.56				
	100 CH <sub>3</sub> OH				-108.76				
	1000 CH <sub>3</sub> OH				-109.10				
	$\infty$ CH <sub>3</sub> OH				-109.25				
	in 20 C <sub>2</sub> H <sub>5</sub> OH				-108.96				
	25 C <sub>2</sub> H <sub>5</sub> OH				-109.05				
	50 C <sub>2</sub> H <sub>5</sub> OH				-109.24				
	100 C <sub>2</sub> H <sub>5</sub> OH				-109.38				
	1000 C <sub>2</sub> H <sub>5</sub> OH				-109.60				
	$\infty$ C <sub>2</sub> H <sub>5</sub> OH				-109.70				

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HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

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Table 9I-6. Lithium (at. no., 3; at. wt., 6.940) HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	kcal/mole				
LiBr	in 100 H <sub>2</sub> O		aq		-95.216				
	200 H <sub>2</sub> O		aq		-95.276				
	400 H <sub>2</sub> O		aq		-95.327				
	600 H <sub>2</sub> O		aq		-95.360				
	1600 H <sub>2</sub> O		aq		-95.384				
	3200 H <sub>2</sub> O		aq		-95.402				
	6400 H <sub>2</sub> O		aq		-95.414				
LiBr·H <sub>2</sub> O LiBr·2H <sub>2</sub> O LiBr·3H <sub>2</sub> O	H <sub>2</sub> O		aq		-95.45				22.6
			c		-158.34				
			c		-229.94				
			c		-301.9				
LiI			g		-16.		19.1	55.68	8.32
			c		-64.79				13.0
	std. state, hyp. m = 1		aq		-79.92		60.523	29.5	
	in 100 H <sub>2</sub> O		aq		-79.781				
	200 H <sub>2</sub> O		aq		-79.824				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 91-7. Lithium (at. no., 3; at. wt., 6.940)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°K)

March 31, 1950

Formula	Substance	Description	State	At 298.16°K (25°K)				
				$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
LiI	in	100 H <sub>2</sub> O	aq		-79.841			
		400 H <sub>2</sub> O	aq		-79.852			
		500 H <sub>2</sub> O	aq		-79.860			
		1000 H <sub>2</sub> O	aq		-79.878			
		$\infty$ H <sub>2</sub> O	aq		-79.92			
			c		-103.8			23.6
LiI·3H <sub>2</sub> O			c		-141.16			32.9
LiI·H <sub>2</sub> O			c		-213.03			43.2
LiI·2H <sub>2</sub> O			c		-285.02			
LiI·3H <sub>2</sub> O			c					
Li <sub>2</sub> SO <sub>4</sub>			c, II		-342.83			
	std. state, hyp. m=1		aq		-350.01	232.928	10.9	
	in	17.75 H <sub>2</sub> O	aq		-348.40			
		200 H <sub>2</sub> O	aq		-349.44			
		300 H <sub>2</sub> O	aq		-349.455			
		400 H <sub>2</sub> O	aq		-349.472			
		500 H <sub>2</sub> O	aq		-349.495			
		1000 H <sub>2</sub> O	aq		-349.590			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards Washington, D.C.

Table 91-8. Lithium (at. no., 3; at. wt., 6.940) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)										
March 31, 1950										
Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula				At 298.16°K (25°C)						
				At 0°K kcal/mole	kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
Li <sub>2</sub> SO <sub>4</sub>	in 2000 H <sub>2</sub> O		aq			-349.638				
	4000 H <sub>2</sub> O		aq			-349.790				
	10000 H <sub>2</sub> O		aq			-349.845				
	50000 H <sub>2</sub> O		aq			-349.930				
	100000 H <sub>2</sub> O		aq			-349.957				
Li <sub>2</sub> SO <sub>4</sub> ·H <sub>2</sub> O	∞ H <sub>2</sub> O		aq			-350.01				
			c			-414.20				
			c			-145.5				
			c			-225.9				
Li <sub>2</sub> Se			c			-91.1				
			aq			-101.6				
Li <sub>2</sub> Se·9H <sub>2</sub> O			c			-729.0				

Table 9I-8. Lithium (at. no., 3; at. wt., 6.940)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 91-9. Lithium (at. no., 3; at. wt., 6.940)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
$\text{Li}_3\text{N}$			c		-47.2				18.4
$\text{LiNO}_2$			c		-96.6				
$\text{LiNO}_3$			c		-115.279				
	std. state, hyp. m = 1		aq		-115.926				
	in 3 $\text{H}_2\text{O}$		aq		-115.1				
	10 $\text{H}_2\text{O}$		aq		-115.54				
	25 $\text{H}_2\text{O}$		aq		-115.623				
	50 $\text{H}_2\text{O}$		aq		-115.677				
	100 $\text{H}_2\text{O}$		aq		-115.721				
	200 $\text{H}_2\text{O}$		aq		-115.759				
	400 $\text{H}_2\text{O}$		aq		-115.795				
	800 $\text{H}_2\text{O}$		aq		-115.825				
	1600 $\text{H}_2\text{O}$		aq		-115.849				
	3200 $\text{H}_2\text{O}$		aq		-115.869				
	6400 $\text{H}_2\text{O}$		aq		-115.884				
	$\infty$ $\text{H}_2\text{O}$		aq		-115.926				
	in 400 $\text{C}_2\text{H}_5\text{OH}$		c		-120.0				
$\text{LiNO}_3 \cdot 3\text{H}_2\text{O}$					-328.6				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 9I-10. Lithium (at. no., 3; at. wt., 6.940)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance									
Formula	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 0°K						
			kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole	
LiNH <sub>2</sub>		c		-43.50					
Li(NH <sub>2</sub> ) <sub>4</sub>		liq		-78.9					
Li <sub>2</sub> NH		c		-52.9					
LiCl·NH <sub>3</sub>		c		-120.5					
LiCl·2NH <sub>3</sub>		c		-142.2					
LiCl·3NH <sub>3</sub>		c		-163.6					
LiCl·4NH <sub>3</sub>		c		-182.6					
LiCl·5NH <sub>3</sub>		c		-201.6					
LiBr·NH <sub>3</sub>		c		-107.6					
LiBr·2NH <sub>3</sub>		c		-130.3					
LiBr·3NH <sub>3</sub>		c		-152.4					
LiBr·4NH <sub>3</sub>		c		-173.5					
LiBr·5NH <sub>3</sub>		c		-192.6					
LiBr·6½NH <sub>3</sub>		c		-220.1					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 91-11. Lithium (at. no., 3; at. wt., 6.940) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	Description		At 0°K kcal/mole	kcal/mole					
LiI·NH <sub>3</sub>		c		-91.8					
LiI·2NH <sub>3</sub>		c		-116.1					
LiI·3NH <sub>3</sub>		c		-138.9					
LiI·4NH <sub>3</sub>		c		-161.3					
LiI·5NH <sub>3</sub>		c		-180.6					
LiI·5½NH <sub>3</sub>		c		-189.7					
LiI·7NH <sub>3</sub>		c		-216.9					
Li <sub>3</sub> Sb <sub>2</sub>		c		-43.5					
Li <sub>3</sub> Bi		c		-55.2					
Li <sub>2</sub> C <sub>2</sub>		c		-14.2					
Li <sub>2</sub> CO <sub>3</sub>		c		-290.54		198.390	21.60	23.28	
	std. state, hyp. m=1			-294.74	-270.66				
	in 1900 H <sub>2</sub> O	aq		-293.92	-266.66	195.459	-5.9		
	∞ H <sub>2</sub> O	aq		-294.74					
	std. state, hyp. m=1	aq		-231.73	-210.53	154.316	29.5		
	in 1000 H <sub>2</sub> O	aq		-231.4					
	∞ H <sub>2</sub> O	aq		-231.73					
LiHCO <sub>3</sub>									

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 91-12. Lithium (at. no., 3; at. wt., 6.940)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Formula	Substance	Description	State	At 298.16°K (25°C)				
				$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
LiOCH <sub>3</sub>		in 60 CH <sub>3</sub> OH		-112.1				
LiOC <sub>2</sub> H <sub>5</sub>		in 60 C <sub>2</sub> H <sub>5</sub> OH		-117.9				
LiCN		in 220 H <sub>2</sub> O	aq	-30.3				
LiCl · CH <sub>3</sub> NH <sub>2</sub>			c	-117.5				
LiCl · C <sub>2</sub> H <sub>5</sub> NH <sub>2</sub>			c	-123.0				
LiCl · (CH <sub>3</sub> ) <sub>2</sub> NH			c	-113.0				
LiCl · 2CH <sub>3</sub> NH <sub>2</sub>			c	-135.2				
LiCl · 3CH <sub>3</sub> NH <sub>2</sub>			c	-152.0				
LiCl · 2C <sub>2</sub> H <sub>5</sub> NH <sub>2</sub>			c	-145.6				
LiCl · 2(CH <sub>3</sub> ) <sub>2</sub> NH			c	-127.4				
LiCl · 4CH <sub>3</sub> NH <sub>2</sub>			c	-168.2				
LiCl · 3C <sub>2</sub> H <sub>5</sub> NH <sub>2</sub>			c	-167.7				
LiCl · 3(CH <sub>3</sub> ) <sub>2</sub> NH			c	-141.6				
LiBr · $\frac{1}{2}$ (CH <sub>3</sub> ) <sub>2</sub> NH			c	-92.1				
LiBr · CH <sub>3</sub> NH <sub>2</sub>			c	-104.9				
LiBr · (CH <sub>3</sub> ) <sub>2</sub> NH			c	-100.3				
LiBr · 2CH <sub>3</sub> NH <sub>2</sub>			c	-124.8				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 91-13. Lithium (at. no., 3; at. wt., 6.940)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description		At 0°K kcal/mole	kcal/mole				
			At 298.16°K (25°C)				cal/deg mole	cal/deg mole
$\text{LiBr} \cdot 3\text{CH}_3\text{NH}_2$		c		-144.4				
$\text{LiBr} \cdot 2(\text{CH}_3)_2\text{NH}$		c		-113.7				
$\text{LiBr} \cdot 4\text{CH}_3\text{NH}_2$		c		-161.2				
$\text{LiBr} \cdot 5\text{CH}_3\text{NH}_2$		c		-176.8				
$\text{LiBr} \cdot 3(\text{CH}_3)_2\text{NH}$		c		-127.9				
$\text{LiBr} \cdot 4(\text{CH}_3)_2\text{NH}$		c		-141.3				
$\text{LiBr} \cdot 5(\text{CH}_3)_2\text{NH}$		c		-154.1				
$\text{LiI} \cdot \frac{1}{2}(\text{CH}_3)_2\text{H}$		c		-74.7				
$\text{LiI} \cdot \text{CH}_3\text{NH}_2$		c		-87.8				
$\text{LiI} \cdot (\text{CH}_3)_2\text{NH}$		c		-83.8				
$\text{LiI} \cdot 2\text{CH}_3\text{NH}_2$		c		-107.6				
$\text{LiI} \cdot \frac{1}{2}(\text{CH}_3)_2\text{NH}$		c		-92.1				
$\text{LiI} \cdot 3\text{CH}_3\text{NH}_2$		c		-125.8				
$\text{LiI} \cdot \frac{3}{2}\text{CH}_3\text{NH}_2$		c		-133.7				
$\text{LiI} \cdot 2(\text{CH}_3)_2\text{H}$		c		-99.8				
$\text{LiI} \cdot 3(\text{CH}_3)_2\text{NH}$		c		-113.0				
$\text{LiI} \cdot 5(\text{CH}_3)_2\text{H}$		c		-138.6				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 91-14. Lithium (at. no., 3; at. wt., 6.940)									
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)									
March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta G_f^\circ$ kcal/mole	$\Delta P_f^\circ$ kcal/mole	At 298.16°K (25°C)		$C_p^\circ$
							$\log_{10} K_f$	$S^\circ$ cal/deg mole	
Li <sub>2</sub> SiO <sub>3</sub> Li <sub>2</sub> SiF <sub>6</sub>		in 15:0 H <sub>2</sub> O	gls						
			c		-376.7				
			aq		-688.9				
LiSn LiSn <sub>2</sub> Li <sub>2</sub> Sn Li <sub>4</sub> Sn Li <sub>5</sub> Sn <sub>2</sub> Li <sub>7</sub> Sn <sub>2</sub>					-691.2				
			c		-16.8				
			c		-17.1				
			c		-27.				
			c		-47.0				
			c		-66.				
LiPb Li <sub>3</sub> Pb Li <sub>4</sub> Pb Li <sub>5</sub> Pb <sub>2</sub> Li <sub>7</sub> Pb <sub>2</sub> 2LiI·PbI <sub>2</sub> 2LiI·PbI <sub>2</sub> ·4H <sub>2</sub> O			c		-86.				
			c		-14.6				
			c		-33.				
			c		-42.				
			c		-57.				
			c		-76.				
			c		-140.27				
			c		-488.65				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 91-15. Lithium (at. no., 3; at. wt., 6.940)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		Description	State	$\Delta H_f^\circ$				$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	$\Delta H_f^\circ$			$\Delta H_f^\circ$	$\Delta H_f^\circ$	At 298.16°K (25°C)				
						kcal/mole	cal/deg mole			
LiTl			c							
LiHg			c							
LiHg <sub>2</sub>			c							
LiHg <sub>3</sub>			c							
LiHg <sub>99</sub>			liq							
LiBr·HgBr <sub>2</sub>		in 4500 H <sub>2</sub> O	aq							
2LiBr·HgBr <sub>2</sub>		in 4700 H <sub>2</sub> O	aq							
4LiBr·HgBr <sub>2</sub>		in 5000 H <sub>2</sub> O	aq							
8LiBr·HgBr <sub>2</sub>		in 5900 H <sub>2</sub> O	aq							
LiCl·Hg(CN) <sub>2</sub>		in 550 H <sub>2</sub> O	aq							
2LiCl·Hg(CN) <sub>2</sub>		in 660 H <sub>2</sub> O	aq							
LiBr·Hg(CN) <sub>2</sub>		in 550 H <sub>2</sub> O	aq							
LiBr·Hg(CN) <sub>2</sub> ·3½H <sub>2</sub> O			c							
2LiBr·Hg(CN) <sub>2</sub>		in 660 H <sub>2</sub> O	aq							

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 91-16. Lithium (at. no., 3; at. wt., 6.940)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
$\text{LiI} \cdot \text{Hg}(\text{CN})_2$	in 50 H <sub>2</sub> O		aq		-16.5				
$\text{LiI} \cdot \text{Hg}(\text{CN})_2 \cdot 3\frac{1}{2}\text{H}_2\text{O}$			c		-265.6				
$2\text{LiI} \cdot \text{Hg}(\text{CN})_2$			aq		-97.3				
$\text{LiCN} \cdot \text{Hg}(\text{CN})_2$			aq		28.7				
$2\text{LiCN} \cdot \text{Hg}(\text{CN})_2$			aq		-7.1				
$\text{LiReO}_4$			c		-253.4				
$\text{LiReO}_4 \cdot \text{H}_2\text{O}$			c		-324.9				
$\text{LiReO}_4 \cdot 2\text{H}_2\text{O}$			c		-395.1				
$\text{LiCl} \cdot \text{ThCl}_4 \cdot 8\text{H}_2\text{O}$			c		-969.8				16.8
$2\text{LiCl} \cdot \text{ThCl}_4$			c		-485.8				
$\text{LiBH}_4$			c		-44.6				
$\text{LiAlH}_4$			c		-24.2				18.2



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 92-1. Sodium (at. no., 11; at. wt., 22.997) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950										
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 298.16°K (25°C)						
				At 0°K kcal/mole	kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	cal/deg mole
Na			g	26.05	25.98	18.67	-13.685	36.715	4.9680	
			c	0.000	0.000	0.000	0.0000	12.2	6.79	
Na <sup>+</sup>			g	144.534	146.015					
			aq		-57.279	-62.569	45.8769	14.4		
Na <sup>++</sup>			g	1235.14	1238.10					
Na <sup>+++</sup>			g	2887.45	2891.89					
Na <sup>++++</sup>			g	5167.8	5173.7					
Na <sup>+++++</sup>			g	8364.1	8371.5					
Na <sup>+++++</sup>			g	12339.0	12347.9					
Na <sup>+++++</sup>			g	17146.1	17156.5					
Na <sup>+++++</sup>			g	23238.	23250.					
Na <sup>+++++</sup>			g	30151.	30164.					
Na <sub>2</sub>			g	34.58	33.97	24.85	-18.215	55.02		
Na <sub>2</sub> O			c		-61.9					
Na <sub>2</sub> O			c		-99.4	-90.0	65.97	17.4	16.3	
Na <sub>2</sub> O <sub>2</sub>			c		-120.6					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 92-2. Sodium (at. no., 11; at. wt. 22.997)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance									
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 298.16°K (25°C)						
			At 0°K kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole	
NaH		g	30.4	29.88	24.78	-18.163	44.93	7.002	
		c		-13.7					
NaOH		c, II		-101.99		73.4335	11.9	19.2	
	std. state, hyp. m = 1	aq		-112.236					
	in 3 H <sub>2</sub> O	aq		-108.894	-100.184				
	4 H <sub>2</sub> O	aq		-110.219					
	5 H <sub>2</sub> O	aq		-111.015					
	6 H <sub>2</sub> O	aq		-111.520					
	7 H <sub>2</sub> O	aq		-111.836					
	8 H <sub>2</sub> O	aq		-112.011					
	10 H <sub>2</sub> O	aq		-112.148					
	12 H <sub>2</sub> O	aq		-112.200					
	15 H <sub>2</sub> O	aq		-112.228					
	18 H <sub>2</sub> O	aq		-112.237					
	20 H <sub>2</sub> O	aq		-112.235					
	25 H <sub>2</sub> O	aq		-112.221					
	30 H <sub>2</sub> O	aq		-112.203					
	40 H <sub>2</sub> O	aq		-112.175					
	50 H <sub>2</sub> O	aq		-112.154					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 92-3. Sodium (at. no., 11; at. wt., 22.987) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	At 298.16°K (25°C)				
NaOH	in	75 H <sub>2</sub> O	aq						
		100 H <sub>2</sub> O	aq		-112.123				
		200 H <sub>2</sub> O	aq		-112.108				
		300 H <sub>2</sub> O	aq		-112.1				
		500 H <sub>2</sub> O	aq		-112.105				
		1000 H <sub>2</sub> O	aq		-112.117				
		2000 H <sub>2</sub> O	aq		-112.139				
		5000 H <sub>2</sub> O	aq		-112.162				
		10000 H <sub>2</sub> O	aq		-112.186				
		50000 H <sub>2</sub> O	aq		-112.201				
NaOH·H <sub>2</sub> O		∞ H <sub>2</sub> O	aq		-112.220				
					-112.236				
NaHO <sub>2</sub> NaF			c		-175.17	-149.00	109.215	20.2	
			aq		-94.4				
			g	-71.	-72.				
		std. state, hyp. m = 1	c		-136.0	-129.3	94.78	14.0	11.0
		in 50 H <sub>2</sub> O	aq		-135.94	-128.67	94.313	12.1	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 92-4. Sodium (at. no., 11; at. wt., 22.987)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	Description	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	in			At 0°K					
				kcal/mole	kcal/mole				
NaF		aq		-135.74					
		aq		-135.78					
		aq		-135.81					
		aq		-135.82					
		aq		-135.84					
		aq		-135.88					
		aq		-135.94					
NaHF <sub>2</sub>		c		-216.6					
		aq		-211.2					
NaCl		g		-43.0		-43.50			
		c		-97.755		-98.232	67.2771	17.30	11.88
	std. state, hyp. m = 1	aq				-97.302	68.8560	27.6	
	in 8 H <sub>2</sub> O	aq				-97.780			
	9 H <sub>2</sub> O	aq				-97.774			
	10 H <sub>2</sub> O	aq				-97.768			
	15 H <sub>2</sub> O	aq				-97.662			
	20 H <sub>2</sub> O	aq				-97.572			
	25 H <sub>2</sub> O	aq				-97.506			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Substance		Table 92-5. Sodium (at. no., 11; at. wt., 22.997) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950				
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$C_p^\circ$
			At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole
NaCl	in 30 H <sub>2</sub> O	aq		-97.455		
	50 H <sub>2</sub> O	aq		-97.340		
	100 H <sub>2</sub> O	aq		-97.250		
	200 H <sub>2</sub> O	aq		-97.216		
	400 H <sub>2</sub> O	aq		-97.212		
	1000 H <sub>2</sub> O	aq		-97.228		
	2000 H <sub>2</sub> O	aq		-97.242		
	5000 H <sub>2</sub> O	aq		-97.260		
	10000 H <sub>2</sub> O	aq		-97.271		
	50000 H <sub>2</sub> O	aq		-97.288		
	100000 H <sub>2</sub> O	aq		-97.292		
	500000 H <sub>2</sub> O	aq		-97.297		
	$\infty$ H <sub>2</sub> O	aq		-97.302		
	in CH <sub>3</sub> OH			-100.1		
NaClO NaClO <sub>2</sub>		aq		-82.7		
		c		-72.65		
	in 100 H <sub>2</sub> O	aq		-73.80		

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 92-6. Sodium (at. no., 11; at. wt., 22.997)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
NaClO <sub>3</sub>	std. state, hyp. m = 1		C						
	in 400 H <sub>2</sub> O		aq		-85.73	-63.21	46.332	53.4	
	600 H <sub>2</sub> O		aq		-80.78				
	800 H <sub>2</sub> O		aq		-80.744				
	1000 H <sub>2</sub> O		aq		-80.739				
	2000 H <sub>2</sub> O		aq		-80.734				
	5000 H <sub>2</sub> O		aq		-80.733				
	10000 H <sub>2</sub> O		aq		-80.735				
	∞ H <sub>2</sub> O		aq		-80.744				
	std. state, hyp. m = 1		aq		-80.752				
	in 400 H <sub>2</sub> O		aq		-80.760				
	600 H <sub>2</sub> O		aq		-80.766				
	10000 H <sub>2</sub> O		aq		-80.770				
NaClO <sub>4</sub>	∞ H <sub>2</sub> O		aq		-80.78				24.1
	std. state, hyp. m = 1		C, II						
	in 400 H <sub>2</sub> O		aq		-92.18	-65.16	47.688	57.9	
	600 H <sub>2</sub> O		aq		-88.69				
	1000 H <sub>2</sub> O		aq		-88.76				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 92-7. Sodium (at. no., 11; at. wt., 22.997) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	Description	State	At 298.16°K (25°C)					
				$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole	
NaClO <sub>4</sub>	in 2000 H <sub>2</sub> O		aq	-88.68					
	5000 H <sub>2</sub> O		aq	-88.67					
	$\infty$ H <sub>2</sub> O		aq	-88.69					
	in CH <sub>3</sub> OH			-94.6					
	in C <sub>2</sub> H <sub>5</sub> OH			-92.6					
NaBr			g	-36.33					12.5
			c	-86.030					
	std. state, hyp. m=1		aq	-86.18	-87.163	63.8933	33.7		
	in 8 H <sub>2</sub> O		aq	-86.888					
	9 H <sub>2</sub> O		aq	-86.864					
	10 H <sub>2</sub> O		aq	-86.846					
	15 H <sub>2</sub> O		aq	-86.662					
	20 H <sub>2</sub> O		aq	-86.536					
	25 H <sub>2</sub> O		aq	-86.454					
	30 H <sub>2</sub> O		aq	-86.390					
	40 H <sub>2</sub> O		aq	-86.244					
	100 H <sub>2</sub> O		aq	-86.142					
	200 H <sub>2</sub> O		aq	-86.104					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards Washington, D.C.

Table 92-8. Sodium (at. no., 11; at. wt., 22.997) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		State	Description	At 0°K kcal/mole	$\Delta H^{\circ}$ kcal/mole	$\Delta F^{\circ}$ kcal/mole	$\log_{10} K_f$	$S^{\circ}$ cal/deg mole	$C_p^{\circ}$ cal/deg mole
Formula									
NaBr	in 400 H <sub>2</sub> O	aq			-86.098				
	1000 H <sub>2</sub> O	aq			-86.108				
	1000 H <sub>2</sub> O	aq			-86.121				
	1000 H <sub>2</sub> O	aq			-86.138				
	1000 H <sub>2</sub> O	aq			-86.149				
	1000 H <sub>2</sub> O	aq			-86.166				
	5000 H <sub>2</sub> O	aq			-86.170				
	10000 H <sub>2</sub> O	aq			-86.175				
	50000 H <sub>2</sub> O	aq			-86.18				
	$\infty$ H <sub>2</sub> O	aq			-90.2				
NaBr•2H <sub>2</sub> O NaBrO	in CH <sub>3</sub> OH				-88.7				
	in C <sub>2</sub> H <sub>5</sub> OH				-227.25				
		c			-79.1				
		aq							
NaI	std. state, hyp. m=1	g		-20.1	-20.94	-70.94	51.998	40.5	13.0
	in 8 H <sub>2</sub> O	c			-68.84				
	9 H <sub>2</sub> O	aq			-70.65				
	10 H <sub>2</sub> O	aq			-71.496				
		aq			-71.452				
		aq			-71.424				



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 92-9. Sodium (at. no., 11; at. wt., 22.997)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	At 298.16°K (25°C)				
NaI	in	15 H <sub>2</sub> O	aq		-71.186				
		20 H <sub>2</sub> O	aq		-71.037				
		25 H <sub>2</sub> O	aq		-70.940				
		30 H <sub>2</sub> O	aq		-70.816				
		50 H <sub>2</sub> O	aq		-70.708				
		100 H <sub>2</sub> O	aq		-70.598				
		200 H <sub>2</sub> O	aq		-70.564				
		400 H <sub>2</sub> O	aq		-70.560				
		1000 H <sub>2</sub> O	aq		-70.576				
		2000 H <sub>2</sub> O	aq		-70.590				
		5000 H <sub>2</sub> O	aq		-70.608				
		10000 H <sub>2</sub> O	aq		-70.619				
		50000 H <sub>2</sub> O	aq		-70.636				
		100000 H <sub>2</sub> O	aq		-70.640				
		500000 H <sub>2</sub> O	aq		-70.645				
		$\infty$ H <sub>2</sub> O	aq		-70.65				
NaI·2H <sub>2</sub> O		in CH <sub>3</sub> OH			-76.1				
		in C <sub>2</sub> H <sub>5</sub> OH			-74.5				
			c		-211.05				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 92-10. Sodium (at. no., 11; at. wt., 22.997) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	At 0°K kcal/mole			kcal/mole	kcal/mole	cal/deg mole	cal/deg mole		
NaIO <sub>3</sub>	in	100 H <sub>2</sub> O	aq		-112.568				
		50 H <sub>2</sub> O	aq		-112.501				
		100 H <sub>2</sub> O	aq		-112.461				
		300 H <sub>2</sub> O	aq		-112.415				
		400 H <sub>2</sub> O	aq		-112.387				
		500 H <sub>2</sub> O	aq		-112.367				
		700 H <sub>2</sub> O	aq		-112.342				
		1000 H <sub>2</sub> O	aq		-112.319				
		2000 H <sub>2</sub> O	aq		-112.291				
		5000 H <sub>2</sub> O	aq		-112.280				
		10000 H <sub>2</sub> O	aq		-112.281				
		20000 H <sub>2</sub> O	aq		-112.284				
Na <sub>2</sub> S	$\infty$	H <sub>2</sub> O	aq		-112.3				
			c		-89.2				
	in	20 H <sub>2</sub> O	aq		-105.42				
		25 H <sub>2</sub> O	aq		-105.22				
		30 H <sub>2</sub> O	aq		-105.08				
		50 H <sub>2</sub> O	aq		-104.79				
		100 H <sub>2</sub> O	aq		-104.55				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 92-11. Sodium (at. no., 11; at. wt., 22.997) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	kcal/mole				
Na <sub>2</sub> S	in 200 H <sub>2</sub> O		aq		-104.40				
	400 H <sub>2</sub> O		aq		-104.3				
	800 H <sub>2</sub> O		aq		-104.36				
Na <sub>2</sub> S·4½H <sub>2</sub> O			c		-416.9				
	Na <sub>2</sub> S·5H <sub>2</sub> O		c		-452.7				
	Na <sub>2</sub> S·9H <sub>2</sub> O		c		-736.7				
Na <sub>2</sub> S <sub>2</sub>			aq		-104.6				
	Na <sub>2</sub> S <sub>3</sub>		aq		-106.5				
	Na <sub>2</sub> S <sub>4</sub>		c		-98.4				
			aq		-108.5				
Na <sub>2</sub> SO <sub>3</sub>			c		-260.6				
			aq		-263.8				
	Na <sub>2</sub> SO <sub>3</sub> ·7H <sub>2</sub> O		c		-753.4				
					-239.5		175.55	34.9	28.7

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 9-12. Sodium (at. no., 11; at. wt., 22.997)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance			State	Description	At 298.16°K (25°C)				
Formula	$\Delta H_f^\circ$	$\Delta F_f^\circ$			$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
								At 0°K	kcal/mole
Na <sub>2</sub> SO <sub>4</sub>			c, II						
	std. state, hyp. m=1		aq						
	in		aq						
	100 H <sub>2</sub> O		aq						
	150 H <sub>2</sub> O		aq						
	200 H <sub>2</sub> O		aq						
	250 H <sub>2</sub> O		aq						
	300 H <sub>2</sub> O		aq						
	400 H <sub>2</sub> O		aq						
	500 H <sub>2</sub> O		aq						
	800 H <sub>2</sub> O		aq						
	1000 H <sub>2</sub> O		aq						
	1500 H <sub>2</sub> O		aq						
	2000 H <sub>2</sub> O		aq						
	3000 H <sub>2</sub> O		aq						
	5000 H <sub>2</sub> O		aq						
	7500 H <sub>2</sub> O		aq						
	10000 H <sub>2</sub> O		aq						
	20000 H <sub>2</sub> O		aq						
	50000 H <sub>2</sub> O		aq						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 92-13. Sodium (at. no., 11; at. wt., 22.997) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K					
				kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
$\text{Na}_2\text{SO}_4$	in 100000 $\text{H}_2\text{O}$		aq		-331.407				
	200000 $\text{H}_2\text{O}$		aq		-331.422				
	500000 $\text{H}_2\text{O}$		aq		-331.435				
	$\infty$ $\text{H}_2\text{O}$		aq		-331.46				
$\text{Na}_2\text{SO}_4 \cdot 10\text{H}_2\text{O}$			c		-1033.48	-870.93	638.380	141.7	140.4
$\text{Na}_2\text{S}_2\text{O}_3$			c		-267.0				34.9
			aq		-269.				
			aq		-270.7				
			c, I		-621.89				86.2
$\text{Na}_2\text{S}_2\text{O}_3 \cdot 5\text{H}_2\text{O}$		in 5 $\text{H}_2\text{O}$	c, II		-620.60				
$\text{Na}_2\text{S}_2\text{O}_5$			c		-349.1				
			aq		-345.				
			c		-399.9				
			aq		-394.6				
$\text{Na}_2\text{S}_2\text{O}_6 \cdot 2\text{H}_2\text{O}$			c		-542.5				
$\text{Na}_2\text{S}_3\text{O}_6$			aq		-409.				
$\text{Na}_2\text{S}_3\text{O}_6 \cdot 3\text{H}_2\text{O}$			c		-623.0				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 92-14. Sodium [at. no., 11; at. wt., 22.997]

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^{\circ}$		$\log_{10} K_f$	$S^{\circ}$	$C_p^{\circ}$
				At 0°K kcal/mole	kcal/mole			
					At 298.16°K (25°C)			
					kcal/mole		cal/deg mole	
$\text{Na}_2\text{S}_4\text{O}_6$			aq		-405.			
$\text{Na}_2\text{S}_4\text{O}_6 \cdot 2\text{H}_2\text{O}$			c		-550.0			
$\text{NaHS}$			c, II		-56.5			
	4 $\text{H}_2\text{O}$		aq		-61.60			
	5 $\text{H}_2\text{O}$		aq		-62.04			
	6 $\text{H}_2\text{O}$		aq		-62.25			
	8 $\text{H}_2\text{O}$		aq		-62.26			
	10 $\text{H}_2\text{O}$		aq		-62.20			
	12 $\text{H}_2\text{O}$		aq		-62.04			
	14 $\text{H}_2\text{O}$		aq		-61.94			
	20 $\text{H}_2\text{O}$		aq		-61.77			
	25 $\text{H}_2\text{O}$		aq		-61.66			
	50 $\text{H}_2\text{O}$		aq		-61.45			
	100 $\text{H}_2\text{O}$		aq		-61.36			
	200 $\text{H}_2\text{O}$		aq		-61.30			
	400 $\text{H}_2\text{O}$		aq		-61.26			
	800 $\text{H}_2\text{O}$		aq		-61.25			
$\text{NaHS} \cdot 2\text{H}_2\text{O}$			c		-199.27			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 92-15. Sodium (at. no., 11; at. wt., 22.987)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	At 298.16°K (25°C)				
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
NaHSO <sub>3</sub>			aq		-206.6				
NaHSO <sub>4</sub>			c		-269.2				
			aq		-270.6				
NaHSO <sub>4</sub> · H <sub>2</sub> O		in 200 H <sub>2</sub> O	c		-339.2				
NaI · 2SO <sub>2</sub>			c		-231.0				
NaI · 4SO <sub>2</sub>			c		-392.2				
Na <sub>2</sub> Se			c		-63.0				
Na <sub>2</sub> Se · 4½H <sub>2</sub> O			aq		-82.9				
Na <sub>2</sub> Se · 9H <sub>2</sub> O			c		-398.2				
Na <sub>2</sub> Se · 16H <sub>2</sub> O			c		-709.1				
					-1199.4				
Na <sub>2</sub> SeO <sub>3</sub>			aq		-236.6				
Na <sub>2</sub> SeO <sub>4</sub>			c		-258.				
			aq		-259.7				
NaHSe			c, II		-27.8				
			aq		-32.7				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards Washington, D.C.

## SERIES I

Table 92-16. Sodium (at. no., 11; at. wt., 22.997)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 298.16°K (25°C)								
	At 0°K kcal/mole			kcal/mole	kcal/mole		cal/deg mole	cal/deg mole	
NaHSO <sub>3</sub>		aq		-180.7					
NaHSO <sub>4</sub>		aq		-201.1					
Na <sub>2</sub> Te		c, II		-84.0					
Na <sub>2</sub> Te <sub>2</sub>		c		-101.5					
Na <sub>2</sub> TeO <sub>4</sub>		c		-313.					
NaN <sub>3</sub>	sodium azide	c			-85.9				19.1
NaNO <sub>2</sub>		c, II			-82.6				
		aq			-111.54				
NaNO <sub>3</sub>		c, II			-106.651	-87.45	64.100	27.8	22.24
	std. state, hyp. m = 1	aq			-108.47	-89.00	65.236	49.4	
	in 1 H <sub>2</sub> O	aq							
	7 H <sub>2</sub> O	aq			-108.31				
	10 H <sub>2</sub> O	aq			-108.10				
	15 H <sub>2</sub> O	aq			-107.81				
	25 H <sub>2</sub> O	aq			-107.45				
	50 H <sub>2</sub> O	aq			-107.02				



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 92-17. Sodium (at. no., 11; at. wt., 22.997) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole			At 298.16°K (25°C) kcal/mole	cal/deg mole
NaNO <sub>3</sub>	in	75 H <sub>2</sub> O	aq		-106.91				
		100 H <sub>2</sub> O	aq		-106.83				
		200 H <sub>2</sub> O	aq		-106.70				
		400 H <sub>2</sub> O	aq		-106.629				
		500 H <sub>2</sub> O	aq		-106.618				
		1000 H <sub>2</sub> O	aq		-106.604				
		2000 H <sub>2</sub> O	aq		-106.604				
		5000 H <sub>2</sub> O	aq		-106.614				
		10000 H <sub>2</sub> O	aq		-106.622				
		15000 H <sub>2</sub> O	aq		-106.626				
		25000 H <sub>2</sub> O	aq		-106.631				
		50000 H <sub>2</sub> O	aq		-106.636				
		100000 H <sub>2</sub> O	aq		-106.640				
		200000 H <sub>2</sub> O	aq		-106.643				
		500000 H <sub>2</sub> O	aq		-106.646				
NaNH <sub>2</sub>	$\infty$ H <sub>2</sub> O		aq		-106.651				
			c		-28.4				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

Table 92-18. Sodium [at. no., 11; at. wt., 22.98976928] <div>HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, <sup>10</sup> HEAT CAPACITY, AT 298.16°K (25°C)</div> <div>March 31, 1950</div>									
Formula	Substance	Description	State	$\Delta H_f^{\circ}$	$\Delta H_f^{\circ}$	$\Delta H_f^{\circ}$	$\log_{10} K_f$	$S^{\circ}$	$C_p^{\circ}$
				At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole		cal/deg mole	cal/deg mole
NaCl · 5NH <sub>3</sub>			c		-193.3				
NaBr · 5½NH <sub>3</sub>			c		-189.2				
NaBr · 5½NH <sub>3</sub>			c		-198.3				
NaI · 4½NH <sub>3</sub>			c		-161.1				
NaI · 6NH <sub>3</sub>			c		-189.2				
Na <sub>2</sub> SO <sub>4</sub> · (NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub> · H <sub>2</sub> O			c		-691.5				
NaPO <sub>3</sub>		in 600 H <sub>2</sub> O	c		-288.6				22.
Na <sub>3</sub> PO <sub>4</sub>		in 300 H <sub>2</sub> O	aq		-292.8				
		500 H <sub>2</sub> O	c		-460.				
		800 H <sub>2</sub> O	aq		-475.0				
		1000 H <sub>2</sub> O	aq		-474.8				
			aq		-474.4				
			aq		-473.9				
Na <sub>3</sub> PO <sub>4</sub> · 12H <sub>2</sub> O			c		-1309.0				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 92-19. Sodium (at. no., 11; at. wt., 22.997)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$
Formula	Description							
$\text{Na}_4\text{P}_2\text{O}_7$	in 1500 $\text{H}_2\text{O}$	c		-760.8				61.
$\text{Na}_4\text{P}_2\text{O}_7 \cdot 10\text{H}_2\text{O}$		aq		-772.9				
$\text{NaH}_2\text{PO}_3$		c		-1468.2				
		c		-289.4				
	in 600 $\text{H}_2\text{O}$	aq		-290.5				
$\text{NaH}_2\text{PO}_3 \cdot 2\frac{1}{2}\text{H}_2\text{O}$		c		-454.8				
$\text{NaH}_2\text{PO}_4$	in 300 $\text{H}_2\text{O}$	aq		-367.7				
$\text{NaH}_3\text{P}_2\text{O}_7$	in 120 $\text{H}_2\text{O}$	c		-602.7				
		aq		-603.9				
$\text{NaH}_3\text{P}_2\text{O}_7 \cdot \text{H}_2\text{O}$		c		-670.6				
$\text{Na}_2\text{HPO}_3$		c		-338.0				
	in 800 $\text{H}_2\text{O}$	aq		-347.5				
$\text{Na}_2\text{HPO}_3 \cdot 5\text{H}_2\text{O}$		c		-684.2				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 92-20. Sodium (at. no., 11; at. wt., 22.997)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
$\text{Na}_2\text{HPO}_4$	in 200 $\text{H}_2\text{O}$		c			-417.4			
			aq			-423.89			
		300 $\text{H}_2\text{O}$	aq			-423.72			
		500 $\text{H}_2\text{O}$	aq			-423.6			
$\text{Na}_2\text{HPO}_4 \cdot 12\text{H}_2\text{O}$	1000 $\text{H}_2\text{O}$		aq			-423.44			
			c			-560.2			86.6
			c			-913.3			133.3
			c			-1266.4			
$\text{Na}_2\text{H}_2\text{P}_2\text{O}_5$			c			-505.6			
			aq			-506.4			
$\text{Na}_2\text{H}_2\text{P}_2\text{O}_7$			c			-663.4			
		in 1500 $\text{H}_2\text{O}$	aq			-661.6			
$\text{Na}_2\text{H}_2\text{P}_2\text{O}_7 \cdot 6\text{H}_2\text{O}$			c			-1085.5			
			c			-711.4			
$\text{Na}_3\text{HP}_2\text{O}_7$			aq			-718.5			
		in 1500 $\text{H}_2\text{O}$	c			-788.2			
$\text{Na}_3\text{HP}_2\text{O}_7 \cdot \text{H}_2\text{O}$			c			-1135.7			
			c						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Washington, D.C.

Table 92-21. Sodium (at. no., 11; at. wt., 22.997) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)									
March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
$\text{NaNH}_4\text{HPO}_4$		in 500 $\text{H}_2\text{O}$	aq		-398.8				
$\text{NaNH}_4\text{HPO}_4 \cdot 4\text{H}_2\text{O}$			c		-682.7				
$\text{Na}_3\text{AsO}_4$			c		-365.				
$\text{Na}_3\text{AsO}_4 \cdot 12\text{H}_2\text{O}$		in 500 $\text{H}_2\text{O}$	aq		-381.5				
$\text{NaH}_2\text{AsO}_3$		in 400 $\text{H}_2\text{O}$	c		-1213.9				
$\text{NaH}_2\text{AsO}_4$		in 300 $\text{H}_2\text{O}$	aq		-227.7				
$\text{Na}_2\text{HASO}_4$		in 400 $\text{H}_2\text{O}$	aq		-273.4				
					-329.5				
$\text{Na}_3\text{Sb}$					-47.2				
$\text{Na}_3\text{SbO}_4$			c		-352.				
$\text{Na}_3\text{SnS}_3$			aq		-193.				
$\text{Na}_3\text{Bi}$					-45.6				
$\text{Na}_3\text{BiO}_4$			c		-288.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 92-22. Sodium (at. no., 11; at. wt., 22.997)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	Description	At 298.16°K (25°C)					
Formula	$\Delta H_f^\circ$			$\log_{10} K_f$	$S^\circ$	$C_p^\circ$			
	At 0°K kcal/mole						kcal/mole	kcal/mole	
$\text{Na}_2\text{C}_2$	sodium carbide	c			4.1				
$\text{Na}_2\text{CO}_3$	in 15 $\text{H}_2\text{O}$	c							
	18 $\text{H}_2\text{O}$	aq				-270.3		183.53	26.41
	20 $\text{H}_2\text{O}$	aq				-278.13			
	25 $\text{H}_2\text{O}$	aq				-278.00			
	30 $\text{H}_2\text{O}$	aq				-277.91			
	40 $\text{H}_2\text{O}$	aq				-277.72			
	50 $\text{H}_2\text{O}$	aq				-277.55			
	75 $\text{H}_2\text{O}$	aq				-277.30			
	100 $\text{H}_2\text{O}$	aq				-277.09			
	200 $\text{H}_2\text{O}$	aq				-276.77			
$\text{Na}_2\text{CO}_3 \cdot \text{H}_2\text{O}$ $\text{Na}_2\text{CO}_3 \cdot 7\text{H}_2\text{O}$ $\text{Na}_2\text{CO}_3 \cdot 10\text{H}_2\text{O}$	400 $\text{H}_2\text{O}$	aq				-276.57			
		aq				-276.17			
		aq				-275.9			
		c				-341.8			
		c				-765.1			128.
		c				-975.6			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 92-23. Sodium (at. no., 11; at. wt., 22.997)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Formula	Substance	State	Description	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} f_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				
$\text{Na}_2\text{C}_2\text{O}_4$	sodium oxalate in 600 $\text{H}_2\text{O}$	c aq			-314.3 -310.5				34.
$\text{NaHC}_2$		c			249.				
$\text{NaCHO}_2$	sodium formate in 400 $\text{H}_2\text{O}$	c aq			-155.03 -154.92				
$\text{NaCHO}_2 \cdot 2\text{H}_2\text{O}$		c			-296.6				
$\text{NaCHO}_2 \cdot 3\text{H}_2\text{O}$		c			-364.2				
$\text{NaHCO}_3$	sodium bicarbonate	c aq			-226.5 -222.5	-203.6	149.24	24.4	20.94
$\text{NaCH}_3\text{O}$	sodium methoxide in 60 $\text{CH}_3\text{OH}$				-105.1				
$\text{NaHC}_2\text{O}_4$	sodium acid oxalate in 400 $\text{H}_2\text{O}$	c aq			-257.8 -252.7				
$\text{NaHC}_2\text{O}_4 \cdot \text{H}_2\text{O}$		c			-330.2				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 97-24. Sodium (at. no., 11; at. wt., 22.997)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		State	Description	Formula	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
At 0°K					At 298.16°K (25°K)						
kcal/mole					kcal/mole	kcal/mole	cal/deg mole	cal/deg mole			
NaC <sub>2</sub> H <sub>3</sub> O <sub>2</sub>	sodium acetate	c				-169.8					
	in 3 H <sub>2</sub> O	aq				-171.20					
	4 H <sub>2</sub> O	aq				-171.51					
	5 H <sub>2</sub> O	aq				-172.09					
	6 H <sub>2</sub> O	aq				-172.57					
	7 H <sub>2</sub> O	aq				-172.86					
	8 H <sub>2</sub> O	aq				-173.01					
	10 H <sub>2</sub> O	aq				-173.24					
	12 H <sub>2</sub> O	aq				-173.36					
	15 H <sub>2</sub> O	aq				-173.48					
	20 H <sub>2</sub> O	aq				-173.58					
	25 H <sub>2</sub> O	aq				-173.63					
	50 H <sub>2</sub> O	aq				-173.752					
	100 H <sub>2</sub> O	aq				-173.827					
	200 H <sub>2</sub> O	aq				-173.890					
	400 H <sub>2</sub> O	aq				-173.941					
	800 H <sub>2</sub> O	aq				-173.995					
1600 H <sub>2</sub> O	aq				-174.022						
3200 H <sub>2</sub> O	aq				-174.049						
6400 H <sub>2</sub> O	aq				-174.072						
∞ H <sub>2</sub> O	aq				-174.122						



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 92-25. Sodium (at. no., 11; at. wt., 22.997) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		Description	State	ΔH <sup>o</sup> <sub>f</sub>		log <sub>10</sub> K <sup>o</sup>	S <sup>o</sup>	C <sup>o</sup> <sub>p</sub>	
Formula				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				cal/deg mole
NaC <sub>2</sub> H <sub>3</sub> O <sub>2</sub>	in 1100 C <sub>2</sub> H <sub>5</sub> OH		c		-171.1				
NaC <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ·3H <sub>2</sub> O					-383.50				
NaC <sub>2</sub> H <sub>3</sub> O <sub>3</sub>	sodium glycolate in 200 H <sub>2</sub> O		c		-213.9				
					-212.2				
NaC <sub>2</sub> H <sub>3</sub> O <sub>3</sub> ·½H <sub>2</sub> O					-249.3				
NaC <sub>2</sub> H <sub>5</sub> O	sodium ethoxide in 60 C <sub>2</sub> H <sub>5</sub> OH				-112.1				
NaC <sub>2</sub> H <sub>5</sub> O <sub>2</sub>	monosodium glycol		c		-148.2				
NaC <sub>2</sub> H <sub>5</sub> O <sub>2</sub> ·CH <sub>3</sub> OH	monosodium glycol methanol		c		-211.8				
NaC <sub>2</sub> H <sub>3</sub> O <sub>3</sub> ·HC <sub>2</sub> H <sub>3</sub> O <sub>3</sub>	sodium acid glycolate		c		-374.6				
NaC <sub>2</sub> H <sub>5</sub> O <sub>2</sub> ·C <sub>2</sub> H <sub>5</sub> OH	monosodium glycol ethanol		c		-219.6				
NaC <sub>2</sub> H <sub>5</sub> O <sub>2</sub> ·C <sub>2</sub> H <sub>6</sub> O <sub>2</sub>	monosodium glycol glycol		c		-263.8				
Na <sub>2</sub> C <sub>2</sub> H <sub>2</sub> O <sub>3</sub>	disodium glycolate in 400 H <sub>2</sub> O		c		-247.5				
			aq		-256.9				
Na <sub>2</sub> C <sub>2</sub> H <sub>2</sub> O <sub>3</sub> ·2H <sub>2</sub> O			c		-393.8				
Na <sub>2</sub> CO <sub>3</sub> ·NaHCO <sub>3</sub> ·2H <sub>2</sub> O	trona		c		-641.2				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 92-26. Sodium (at. no., 11; at. wt., 22.997)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance									
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$G_p^\circ$		
			At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole	cal/deg mole		
$\text{NaClO}_3$	sodium trichloroacetate in 400 $\text{H}_2\text{O}$	aq		-182.3					
$\text{NaCH}_2\text{ClCO}_2$	sodium chloroacetate in 400 $\text{H}_2\text{O}$	aq		-178.5					
$\text{NaCHCl}_2\text{CO}_2$	sodium dichloroacetate in 400 $\text{H}_2\text{O}$	aq		-180.7					
$\text{NaI} \cdot 3\text{CH}_3\text{OH}$		c		-249.3					
$\text{NaC}_2\text{H}_5\text{OSO}_3$	sodium ethylsulfate in 400 $\text{H}_2\text{O}$	aq		-265.9					
$2\text{NaHSO}_3 \cdot \text{C}_2\text{H}_2\text{O}_2 \cdot \text{H}_2\text{O}$	sodium bisulfite glyoxal	c		-580.4					
		aq		-572.0					
$\text{NaCN}$		c, III		-21.46					
	in 200 $\text{H}_2\text{O}$	aq		-21.2					
$\text{NaCN} \cdot \frac{1}{2}\text{H}_2\text{O}$		c		-56.19					
$\text{NaCN} \cdot 2\text{H}_2\text{O}$		c		-162.25					
$\text{NaCNO}$	sodium cyanate	c		-95.6					
	in 2000 $\text{H}_2\text{O}$	aq		-90.9					
$\text{NaCN}_2\text{H}$	sodium cyanamid	aq		-34.5					
$\text{NaCO}_2\text{NH}_2$	sodium carbamate	c		-179.4					

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Table 92-27. Sodium (at. no., 11; at. wt., 22.997) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta E_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	kcal/mole			At 298.16°K (25°C) kcal/mole	cal/deg mole
NaCNS	sodium thiocyanate in		c						
	3 H <sub>2</sub> O		aq		-41.73				
	4 H <sub>2</sub> O		aq		-41.13				
	5 H <sub>2</sub> O		aq		-41.25				
	6 H <sub>2</sub> O		aq		-41.30				
	8 H <sub>2</sub> O		aq		-41.23				
	10 H <sub>2</sub> O		aq		-41.21				
	12 H <sub>2</sub> O		aq		-41.11				
	15 H <sub>2</sub> O		aq		-41.01				
	20 H <sub>2</sub> O		aq		-40.88				
	25 H <sub>2</sub> O		aq		-40.70				
	30 H <sub>2</sub> O		aq		-40.59				
	40 H <sub>2</sub> O		aq		-40.30				
	100 H <sub>2</sub> O		aq		-40.16				
	200 H <sub>2</sub> O		aq		-40.08				
	400 H <sub>2</sub> O		aq		-40.05				
	800 H <sub>2</sub> O		aq		-40.04				
	1600 H <sub>2</sub> O		aq		-40.04				
	3200 H <sub>2</sub> O		aq		-40.05				
	6400 H <sub>2</sub> O		aq		-40.06				
	$\infty$ H <sub>2</sub> O		aq		-40.1				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 92-28. Sodium (at. no., 11; at. wt., 22.997) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	Description	State	$\Delta H^{\circ}$		$\log_{10} K^{\circ}$	$S^{\circ}$	$C_p^{\circ}$	
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				
NaCNS	in s	C <sub>2</sub> H <sub>5</sub> OH			-42.85				
	10	C <sub>2</sub> H <sub>5</sub> OH			-42.77				
	12	C <sub>2</sub> H <sub>5</sub> OH			-42.61				
	15	C <sub>2</sub> H <sub>5</sub> OH			-42.35				
	20	C <sub>2</sub> H <sub>5</sub> OH			-42.23				
	25	C <sub>2</sub> H <sub>5</sub> OH			-42.03				
	30	C <sub>2</sub> H <sub>5</sub> OH			-41.53				
	50	C <sub>2</sub> H <sub>5</sub> OH			-41.13				
	100	C <sub>2</sub> H <sub>5</sub> OH			-40.88				
	200	C <sub>2</sub> H <sub>5</sub> OH			-40.80				
			c		-363.	249.9	27.2	26.72	
Na <sub>2</sub> SiO <sub>3</sub>		g l s			-360.				
		c			-720.0				
		c			-1002.0				
		c			-677.		39.4	37.41	
Na <sub>2</sub> Si <sub>2</sub> O <sub>5</sub>		c			-671.2		46.8	43.79	
		aq			-614.1				
Na <sub>4</sub> SiO <sub>4</sub>		aq							
Na <sub>2</sub> SiF <sub>6</sub>	in 600 H <sub>2</sub> O								
	in 400 H <sub>2</sub> O								
NaHSiF <sub>6</sub>									

March 31, 1950

Substance						
Formula	Description	State	$\Delta H^\circ$ kcal/mole	$\Delta F^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole
			At 0°K	At 298.16°K (25°C)		
			kcal/mole	kcal/mole		cal/deg mole
NaSn		c		-12.		
NaSn <sub>2</sub>		c		-12.		
Na <sub>2</sub> Sn		c		-14.4		
Na <sub>4</sub> Sn		c		-13.5		
Na <sub>4</sub> Sn <sub>3</sub>		c		-38.		
Na <sub>2</sub> SnO <sub>3</sub>		c		-276.		
Na <sub>4</sub> SnO <sub>4</sub>	in 1200 H <sub>2</sub> O	aq		-455.5		
NaPb		c		-11.6		
Na <sub>2</sub> Pb		c		-15.6		
Na <sub>2</sub> Pb <sub>5</sub>		c		-22.		
Na <sub>4</sub> Pb		c		-20.		
Na <sub>5</sub> Pb <sub>2</sub>		c		-35.		
Na <sub>2</sub> PbO <sub>3</sub>		c		-205.		
2NaI·PbI <sub>2</sub>		c		-103.4		
2NaI·PbI <sub>2</sub> ·4H <sub>2</sub> O		c		-481.0		
2NaI·PbI <sub>2</sub> ·6H <sub>2</sub> O		c		-620.6		
2Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> ·PbS <sub>2</sub> O <sub>3</sub>		c		-698.7		
		aq		-696.		

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 92-30. Sodium (at. no., 11; at. wt., 22.997)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K kcal/mole			At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	At 298.16°K (25°C) cal/deg mole	At 298.16°K (25°C) cal/deg mole	At 298.16°K (25°C) cal/deg mole	
$\text{Na}_2\text{ZnO}_2$			c		-188.				
$\text{Na}_2\text{Zn(SO}_4)_2$			c		-567.0				
$\text{Na}_2\text{Zn(SO}_4)_2 \cdot 4\text{H}_2\text{O}$			c		-856.7				
$\text{NaCd}_2$			c		-8.4				
$\text{NaCd}_6$			c		-12.6				
$\text{Na}_2\text{CdO}_2$			c		-165.				
$\text{NaHg}$			c		-10.2				
$\text{NaHg}_2$			c		-18.3				
$\text{NaHg}_4$			c		-20.0				
$\text{NaHg}_{27.5}$			liq		-19.65				
$\text{NaHg}_{50}$			liq		-19.80				
$\text{NaHg}_{200}$			liq		-19.94				
$\text{NaHg}_{500}$			liq		-19.97				
$\text{Na}_3\text{Hg}$			c		-11.2				
$\text{Na}_3\text{Hg}_2$			c		-22.5				
$\text{Na}_5\text{Hg}_2$			c		-22.4				
$\text{Na}_7\text{Hg}_8$			c		-80.				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards Washington, D.C.

Table 92-31. Sodium (at. no., 11; at. wt., 22.997)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Formula	Substance	Description	State	At 298.16°K (25°C)				
				$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
$\text{Na}_2\text{HgBr}_4$		in 6000 $\text{H}_2\text{O}$	aq		-214.3			
$\text{Na}_2\text{HgS}_2$			c		-120.			
$\text{NaHg}(\text{CN})_3$		in 500 $\text{H}_2\text{O}$	aq		37.8			
$\text{Na}_2\text{Hg}(\text{CN})_4$		in 600 $\text{H}_2\text{O}$	aq		11.3			
$\text{NaCl} \cdot \text{Hg}(\text{CN})_2$			aq		-31.3			
$\text{NaCl} \cdot \text{Hg}(\text{CN})_2 \cdot 14\text{H}_2\text{O}$			c		-124.3			
$2\text{NaCl} \cdot \text{Hg}(\text{CN})_2$			aq		-128.6			
$\text{NaBr} \cdot \text{Hg}(\text{CN})_2$			aq		-20.5			
$\text{NaBr} \cdot \text{Hg}(\text{CN})_2 \cdot 2\text{H}_2\text{O}$			c		-168.9			
$2\text{NaBr} \cdot \text{Hg}(\text{CN})_2$			aq		-106.9			
$\text{NaI} \cdot \text{Hg}(\text{CN})_2$			aq		-7.2			
$\text{NaI} \cdot \text{Hg}(\text{CN})_2 \cdot 2\text{H}_2\text{O}$			c		-154.9			
$2\text{NaI} \cdot \text{Hg}(\text{CN})_2$			aq		-78.9			
$\text{Na}_2\text{O} \cdot \text{CuO}_2$			c		-168.			
$\text{Na}_2\text{CO}_3 \cdot \text{CuCO}_3$			c		-411.4			
$\text{Na}_2\text{CO}_3 \cdot \text{CuCO}_3 \cdot 3\text{H}_2\text{O}$			c		-625.6			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 92-32. Sodium (at. no., 11; at. wt., 22.997)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole		cal/deg mole	cal/deg mole
$\text{Na}_2\text{O} \cdot \text{Ag}_2\text{O}_2$			c					
$\text{Na}_3\text{Ag}(\text{S}_2\text{O}_3)_2$			aq					
$\text{NaAg}(\text{CN})_2$		in 600 $\text{H}_2\text{O}$	aq					
$\text{Na}_2\text{Ag}(\text{CN})_3$		in 100 $\text{H}_2\text{O}$	aq					
$\text{Na}_2\text{PtCl}_4$			aq					
$\text{Na}_2\text{PtCl}_6$			c					
$\text{Na}_2\text{PtCl}_6 \cdot 2\text{H}_2\text{O}$			aq					
$\text{Na}_2\text{PtCl}_6 \cdot 6\text{H}_2\text{O}$			c					
$\text{Na}_2\text{PtBr}_6$			c					
$\text{Na}_2\text{PtBr}_6 \cdot 6\text{H}_2\text{O}$			aq					
$\text{Na}_2\text{PtI}_6$			c					
$\text{Na}_2\text{IrCl}_6$			c					
$\text{Na}_3\text{IrCl}_6$			c					
$\text{Na}_2\text{OsCl}_6$			c					



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

Notional Bureau of Standards Washington, D.C.

SERIES I

Table 92-33. Sodium [at. no., 11; at. wt., 22.997]  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)		cal/deg mole	cal/deg mole
			kcal/mole	kcal/mole	kcal/mole		
NaReO <sub>4</sub>		c		-249.4			
Na <sub>3</sub> RhCl <sub>6</sub>		aq		-247.6			
Na <sub>3</sub> RhCl <sub>6</sub> ·12H <sub>2</sub> O		c		-371.7			
Na <sub>2</sub> Ni(CN) <sub>4</sub>		aq		-379.6			
		c		-1220.0			
Na <sub>2</sub> CoO <sub>3</sub>		aq		-27.9			
		c		-203.			
Na <sub>2</sub> Fe <sub>2</sub> O <sub>4</sub>		c		-251.			
Na <sub>3</sub> FeCO(CN) <sub>5</sub>		c		-117.8			
		aq		-124.0			
Na <sub>3</sub> FeCO(CN) <sub>5</sub> ·7H <sub>2</sub> O		c		-609.0			
Na <sub>2</sub> MnO <sub>4</sub>		c		-274.			
Na <sub>2</sub> SO <sub>4</sub> ·MnSO <sub>4</sub>		c		-565.5			
Na <sub>2</sub> SO <sub>4</sub> ·MnSO <sub>4</sub> ·2H <sub>2</sub> O		c		-732.7			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 92-34. Sodium (at. no., 11; at. wt., 22.997)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	$\Delta H_f^\circ$ kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$
Formula	Description		At 0°K	At 298.16°K (25°C)				
			kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
$\text{Na}_2\text{CrO}_4$		c						
	in 10 $\text{H}_2\text{O}$	aq		-317.6				
	600 $\text{H}_2\text{O}$	aq		-324.1				
$\text{Na}_2\text{CrO}_4 \cdot 4\text{H}_2\text{O}$		c		-320.6				
$\text{Na}_2\text{Cr}_2\text{O}_7$		aq		-601.3				
	in 600 $\text{H}_2\text{O}$	aq		-463.4				
$\text{Na}_2\text{MoO}_4$		c, II		-368.				
	in 800 $\text{H}_2\text{O}$	aq		-368.6				
$\text{Na}_2\text{WO}_4$		c		-395.				
	in 200 $\text{H}_2\text{O}$	aq		-380.9				
$\text{NaVO}_5$		aq		-252.2				
$\text{Na}_3\text{VO}_4$		c		-420.				
$\text{Na}_3\text{V}_3\text{O}_{19}$		aq		-849.9				
$\text{Na}_2\text{TiO}_3$		c, II					29.1	30.02
$\text{Na}_2\text{Ti}_2\text{O}_5$		c					41.5	41.68
$\text{Na}_2\text{Ti}_3\text{O}_7$		c					55.9	54.85

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Substance		Table 92-35. Sodium (at. no., 11; at. wt., 22.997) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950					
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
			At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
NaBO <sub>2</sub>	in 300 H <sub>2</sub> O	c		-253.			16.5
NaBO <sub>3</sub>		aq		-241.1			
NaBO <sub>3</sub> ·4H <sub>2</sub> O		aq		-220.0			
Na <sub>2</sub> B <sub>4</sub> O <sub>7</sub>		c		-504.8			48.
		c		-777.7			
		aq		-787.9			
Na <sub>2</sub> B <sub>4</sub> O <sub>7</sub> ·4H <sub>2</sub> O		c		-1072.9			
Na <sub>2</sub> B <sub>4</sub> O <sub>7</sub> ·5H <sub>2</sub> O		c		-1143.5			
Na <sub>2</sub> B <sub>4</sub> O <sub>7</sub> ·10H <sub>2</sub> O		c		-1497.2			147.
NaBH <sub>4</sub>	sodium borohydride	c, I		-43.82	-20.941	25.02	20.7
NaAlO <sub>2</sub>		c		-273.			
Na <sub>3</sub> AlF <sub>6</sub>		c, II		-759.6			56.2
Na <sub>3</sub> AlF <sub>6</sub> ·3H <sub>2</sub> O		c		-1021.0			
NaCl·AlCl <sub>3</sub>		c		-270.6			
3NaCl·AlCl <sub>3</sub>		c		-470.2			
3NaCl·2AlCl <sub>3</sub>		c		-642.4			
NaCl·AlCl <sub>3</sub> ·6NH <sub>3</sub>		c		-463.0			
Na <sub>2</sub> O·Al <sub>2</sub> O <sub>3</sub> ·3SiO <sub>2</sub>	natrolite	c		-1205.2			
Na <sub>2</sub> O·Al <sub>2</sub> O <sub>3</sub> ·4SiO <sub>2</sub>	dehydrated analcite	c		-1393.3			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Substance		HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)				
Formula	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\log_{10} K_f$	$C_p^\circ$
			At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole	cal/deg mole
$\text{Na}_2\text{UO}_4$		c		-501.		
$\text{Na}_2\text{U}_2\text{O}_7 \cdot 1\frac{1}{2}\text{H}_2\text{O}$		c		-880.		
$(\text{Na}_2\text{O})_2 \cdot 2\text{UO}_4$		aq		-596.		
$(\text{Na}_2\text{O}_2)_2 \cdot \text{UO}_4 \cdot 9\text{H}_2\text{O}$		c		-1225.		
$2\text{NaCl} \cdot \text{ThCl}_4$		c		-485.8		
$2\text{NaCl} \cdot \text{ThCl}_4 \cdot 10\text{H}_2\text{O}$		c		-1225.7		
$\text{Na}_2\text{SO}_4 \cdot \text{CaSO}_4$		c		-674.1		
$2\text{Na}_2\text{SO}_4 \cdot \text{CaSO}_4 \cdot 2\text{H}_2\text{O}$		c		-1146.1		
$\text{NaSrPO}_4 \cdot 9\text{H}_2\text{O}$		c		-1121.5		
$\text{NaSrAsO}_4 \cdot 9\text{H}_2\text{O}$		c		-1027.5		
$\text{NaBaPO}_4 \cdot 9\text{H}_2\text{O}$		c		-1121.2		
$\text{NaBaAsO}_4 \cdot 9\text{H}_2\text{O}$		c		-1026.3		
$\text{NaLiICl}$		c		-166.0		

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards									
Washington, D.C.									
Table 93-1. Potassium (at. no., 19; at. wt., 39.096)									
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)									
March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	At 298.16°K (25°C)				
Formula				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
K			g	21.73	21.51	14.52	-10.716	38.296	4.968
K <sup>+</sup>			c	0.000	0.000	0.000	0.0000	15.2	6.97
			g	121.81	123.07				
K <sup>++</sup>		std. state, hyp. m=1	aq		-60.04	-67.466	49.4516	24.5	
K <sup>+++</sup>			g	855.41	858.15				
K <sup>++++</sup>			g	1909.	1913.				
K <sup>*****</sup>			g	3313.	3318.				
K <sub>2</sub>			g	31.7	30.8	22.1	-16.20	59.69	
K <sub>2</sub> O			c		-86.4				
K <sub>2</sub> O <sub>2</sub>			c		-118.				
K <sub>2</sub> O <sub>3</sub>			c		-125.				
K <sub>2</sub> O <sub>4</sub>			c		-134.				
KH			g	30.4	30.0	25.1	-18.40	47.3	
			c		-13.6				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 93-2. Potassium (at. no., 19; at. wt., 39.096)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description							
KOH	std. state, hyp. m=1	c						
	in 3 H <sub>2</sub> O	aq		-101.78				
	4 H <sub>2</sub> O	aq		-115.00				
	5 H <sub>2</sub> O	aq		-111.77				
	6 H <sub>2</sub> O	aq		-112.72				
	8 H <sub>2</sub> O	aq		-113.31				
	10 H <sub>2</sub> O	aq		-113.700				
	12 H <sub>2</sub> O	aq		-114.143				
	15 H <sub>2</sub> O	aq		-114.363				
	20 H <sub>2</sub> O	aq		-114.493				
	25 H <sub>2</sub> O	aq		-114.594				
	50 H <sub>2</sub> O	aq		-114.677				
	100 H <sub>2</sub> O	aq		-114.702				
	200 H <sub>2</sub> O	aq		-114.767				
	500 H <sub>2</sub> O	aq		-114.796				
	1000 H <sub>2</sub> O	aq		-114.822				
	2000 H <sub>2</sub> O	aq		-114.864				
	5000 H <sub>2</sub> O	aq		-114.894				
	10000 H <sub>2</sub> O	aq		-114.923				
	50000 H <sub>2</sub> O	aq		-114.950				
	$\infty$ H <sub>2</sub> O	aq		-114.964				
		aq		-114.984				
		aq		-115.00				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 93-3. Potassium (at. no., 19; at. wt., 39.096) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950												
Substance				State	Description	Formula	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
At 0°K		At 298.16°K (25°C)										
kcal/mole	kcal/mole	kcal/mole	kcal/mole									
KOH·3H <sub>2</sub> O KOH·H <sub>2</sub> O KOH·2H <sub>2</sub> O	c		-161.7 -179.6 -251.2									
	c		-134.46 -137.35 -137.75 -138.34 -138.39 -138.44 -138.47 -138.48 -138.51			-127.42	93.397	15.91	11.73			
	aq											
KF	in 4 H <sub>2</sub> O											
	5 H <sub>2</sub> O											
	10 H <sub>2</sub> O											
	12 H <sub>2</sub> O											
	15 H <sub>2</sub> O											
	20 H <sub>2</sub> O											
	25 H <sub>2</sub> O											
	50 H <sub>2</sub> O											

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 93-4. Potassium (at. no., 19; at. wt., 39.096)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	Description	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	in			At 0°K					
				kcal/mole	kcal/mole				
KF	in 100 H <sub>2</sub> O	aq			-138.534				
	200 H <sub>2</sub> O	aq			-138.553				
	300 H <sub>2</sub> O	aq			-138.567				
	400 H <sub>2</sub> O	aq			-138.578				
	500 H <sub>2</sub> O	aq			-138.586				
	1000 H <sub>2</sub> O	aq			-138.609				
	5000 H <sub>2</sub> O	aq			-138.651				
	$\infty$ H <sub>2</sub> O	aq			-138.70				
KF·2H <sub>2</sub> O		c			-277.00	-242.7	177.90	36.	
		c			-418.0				
		c			-219.98	-203.73	149.331	24.92	18.37
	in 20 H <sub>2</sub> O	aq			-213.73				
	50 H <sub>2</sub> O	aq			-213.49				
KF·3HF	100 H <sub>2</sub> O	aq			-213.39				
	200 H <sub>2</sub> O	aq			-213.4				
		c			-296.7				
		c			-373.0				



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Table 93-5. Potassium (at. no., 19; at. wt., 39.098)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole			cal/deg mole	cal/deg mole
KCl	std. state, hyp. m=1 in 50 H <sub>2</sub> O 100 H <sub>2</sub> O 200 H <sub>2</sub> O 400 H <sub>2</sub> O 800 H <sub>2</sub> O 1600 H <sub>2</sub> O 3200 H <sub>2</sub> O 6400 H <sub>2</sub> O $\infty$ H <sub>2</sub> O		g	-51.2	-51.6	-56.2	41.19	57.24	8.65
					-104.175	-97.592	71.5336	19.70	12.31
					-100.06	-98.816	72.4302	37.7	
					-100.105				
					-100.015				
					-99.974				
					-99.973				
					-99.962				
					-99.994				
					-100.008				
KClO KClO <sub>3</sub>	std. state, hyp. m=1		aq		-100.026				
					-100.06				
					-85.4				
					-93.50	-69.29	50.769	34.17	23.96
					-83.54	-68.09	49.909	63.5	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 93-6. Potassium (at. no., 19; at. wt., 39.098)									
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)									
March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	At 298.16°K (25°C)				
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
KClO <sub>3</sub>	in	100 H <sub>2</sub> O	aq		-84.09				
		200 H <sub>2</sub> O	aq		-83.72				
		300 H <sub>2</sub> O	aq		-83.67				
		400 H <sub>2</sub> O	aq		-83.610				
		500 H <sub>2</sub> O	aq		-83.580				
		600 H <sub>2</sub> O	aq		-83.555				
		800 H <sub>2</sub> O	aq		-83.543				
		900 H <sub>2</sub> O	aq		-83.535				
		1000 H <sub>2</sub> O	aq		-83.531				
		2000 H <sub>2</sub> O	aq		-83.514				
		3000 H <sub>2</sub> O	aq		-83.511				
		4000 H <sub>2</sub> O	aq		-83.511				
		5000 H <sub>2</sub> O	aq		-83.512				
		10000 H <sub>2</sub> O	aq		-83.515				
		20000 H <sub>2</sub> O	aq		-83.520				
		50000 H <sub>2</sub> O	aq		-83.526				
		100000 H <sub>2</sub> O	aq		-83.530				
		200000 H <sub>2</sub> O	aq		-83.532				
		500000 H <sub>2</sub> O	aq		-83.535				
		∞ H <sub>2</sub> O	aq		-83.54				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 99-7. Potassium (at. no., 19; at. wt., 39.096)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Formula	Substance	Description	State	At 298.16°K (25°C)				
				$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
KC10 <sub>4</sub>	std. state, hyp. m = 1		c					
	in 500 H <sub>2</sub> O		aq					
	600 H <sub>2</sub> O		aq					
	700 H <sub>2</sub> O		aq					
	800 H <sub>2</sub> O		aq					
	900 H <sub>2</sub> O		aq					
	1000 H <sub>2</sub> O		aq					
	1500 H <sub>2</sub> O		aq					
	2000 H <sub>2</sub> O		aq					
	3000 H <sub>2</sub> O		aq					
	4000 H <sub>2</sub> O		aq					
	5000 H <sub>2</sub> O		aq					
	8000 H <sub>2</sub> O		aq					
	10000 H <sub>2</sub> O		aq					
	20000 H <sub>2</sub> O		aq					
	50000 H <sub>2</sub> O		aq					
	∞ H <sub>2</sub> O		aq					

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Table 93-8. Potassium (at. no., 19; at. wt., 39.096)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

Substance		State	Description	At 298.16°K (25°K)					
Formula	$\Delta H_f^\circ$			$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
	kcal/mole			kcal/mole		cal/deg mole	cal/deg mole		
KBr		g							
		c							
	std. state, hyp. m = 1	aq							
	in	aq							
	10 H <sub>2</sub> O	aq							
	12 H <sub>2</sub> O	aq							
	15 H <sub>2</sub> O	aq							
	20 H <sub>2</sub> O	aq							
	25 H <sub>2</sub> O	aq							
	30 H <sub>2</sub> O	aq							
	40 H <sub>2</sub> O	aq							
	50 H <sub>2</sub> O	aq							
	75 H <sub>2</sub> O	aq							
	100 H <sub>2</sub> O	aq							
	200 H <sub>2</sub> O	aq							
300 H <sub>2</sub> O	aq								
400 H <sub>2</sub> O	aq								
500 H <sub>2</sub> O	aq								
600 H <sub>2</sub> O	aq								
800 H <sub>2</sub> O	aq								
1000 H <sub>2</sub> O	aq								

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 93-9. Potassium (at. no., 19; at. wt., 39.098) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	Description	State	$\Delta H^{\circ}$	$\Delta F^{\circ}$	$\log_{10} K_f$	$S^{\circ}$	$C_p^{\circ}$	
				At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole	cal/deg mole	
KBr	in 2000 H <sub>2</sub> O		aq		-88.892				
	3000 H <sub>2</sub> O		aq		-88.898				
	4000 H <sub>2</sub> O		aq		-88.902				
	5000 H <sub>2</sub> O		aq		-88.905				
	8000 H <sub>2</sub> O		aq		-88.911				
	10000 H <sub>2</sub> O		aq		-88.915				
	20000 H <sub>2</sub> O		aq		-88.921				
	50000 H <sub>2</sub> O		aq		-88.928				
	100000 H <sub>2</sub> O		aq		-88.931				
	200000 H <sub>2</sub> O		aq		-88.934				
	$\infty$ H <sub>2</sub> O		aq		-88.94				
	$\infty$ CH <sub>3</sub> OH		aq		-94.4				
KBr <sub>3</sub> KBr <sub>5</sub>			aq		-92.0				
			aq		-96.5				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 93-10. Potassium (at. no., 19; at. wt., 39.096) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole				
KBrO			aq						
KBrO <sub>3</sub>			c						
	std. state, hyp. m = 1		aq						
	in 500 H <sub>2</sub> O		aq	-62.0	-58.2		42.66	35.65	25.07
	1000 H <sub>2</sub> O		aq	-79.4	-56.6		41.49	63.4	
	10000 H <sub>2</sub> O		aq	-69.6					
			aq	-69.64					
			aq	-69.59					
			aq	-69.58					
	$\infty$ H <sub>2</sub> O		aq	-69.6					
KI			g						
	std. state, hyp. m = 1		c						
	in 6 H <sub>2</sub> O		aq	-78.31	-77.03		56.462	61.64	8.78
	8 H <sub>2</sub> O		aq	-73.41	-79.816		58.5040	24.94	13.16
	10 H <sub>2</sub> O		aq	-74.94				50.6	
	15 H <sub>2</sub> O		aq	-74.71					
	20 H <sub>2</sub> O		aq	-74.55					
	25 H <sub>2</sub> O		aq	-74.23					
	50 H <sub>2</sub> O		aq	-74.04					
	100 H <sub>2</sub> O		aq	-73.910					
			aq	-73.625					
			aq	-73.471					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 93-11. Potassium (at. no., 19; at. wt., 39.096) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	Description		At 0°K kcal/mole	kcal/mole					
KI	in 200 H <sub>2</sub> O	aq		-73.397					
	400 H <sub>2</sub> O	aq		-73.365					
	800 H <sub>2</sub> O	aq		-73.355					
	1600 H <sub>2</sub> O	aq		-73.357					
	3200 H <sub>2</sub> O	aq		-73.365					
	6400 H <sub>2</sub> O	aq		-73.373					
	$\infty$ H <sub>2</sub> O	aq		-73.41					
	CH <sub>3</sub> OH, methanol			-78.14					
	C <sub>2</sub> H <sub>5</sub> OH, ethanol			-77.6					
	200 CH <sub>3</sub> COCH <sub>3</sub> , acetone			-82.5					
KI <sub>3</sub>	200 CH <sub>3</sub> CN, acetonitrile			-75.6					
		c		-76.6					
	in 100 H <sub>2</sub> O	aq		-72.6					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards Washington, D.C.

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Table 93-12. Potassium (at. no., 19; at. wt., 39.096)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K kcal/mole			kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole	
KIO <sub>3</sub>		std. state, hyp. m=1	c						
		in 200 H <sub>2</sub> O	aq		-121.5	-101.7	74.54	36.20	25.42
		300 H <sub>2</sub> O	aq		-115.0	-99.9	73.22	52.2	
		400 H <sub>2</sub> O	aq		-115.16				
		500 H <sub>2</sub> O	aq		-115.12				
		1000 H <sub>2</sub> O	aq		-115.09				
		5000 H <sub>2</sub> O	aq		-115.07				
		10000 H <sub>2</sub> O	aq		-115.02				
		$\infty$ H <sub>2</sub> O	aq		-114.98				
		in 1200 H <sub>2</sub> O	aq		-114.98				
			aq		-115.0				
KIO <sub>3</sub> ·HIO <sub>3</sub>			aq		-97.6				
			c		-180.7				
			aq		-170.1				
			aq		-234.2				
			aq		-302.2				
KH <sub>4</sub> IO <sub>6</sub>									
K <sub>2</sub> H <sub>3</sub> IO <sub>6</sub>									
KI·KCl		fresh melt	c		-181.95				



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 93-13. Potassium (at. no., 19; at. wt., 39.096) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance		State	ΔH <sup>o</sup> <sub>f</sub>	ΔH <sup>o</sup> <sub>f</sub>	ΔF <sup>o</sup> <sub>f</sub>	log <sub>10</sub> K <sub>f</sub>	S <sup>o</sup>	C <sub>p</sub> <sup>o</sup>
	Description	At 0°K kcal/mole		kcal/mole	kcal/mole	At 298.16°K (25°C) cal/deg mole	cal/deg mole	cal/deg mole	
K <sub>2</sub> S	in	7 H <sub>2</sub> O	c		-100.				
		8 H <sub>2</sub> O	aq		-107.04				
		10 H <sub>2</sub> O	aq		-108.05				
		12 H <sub>2</sub> O	aq		-108.95				
		15 H <sub>2</sub> O	aq		-109.49				
		20 H <sub>2</sub> O	aq		-109.90				
		25 H <sub>2</sub> O	aq		-110.17				
		50 H <sub>2</sub> O	aq		-110.27				
		100 H <sub>2</sub> O	aq		-110.17				
		200 H <sub>2</sub> O	aq		-110.02				
		400 H <sub>2</sub> O	aq		-109.94				
			aq		-109.9				
K <sub>2</sub> S·2H <sub>2</sub> O K <sub>2</sub> S·5H <sub>2</sub> O K <sub>2</sub> S <sub>4</sub>		c			-243.0				
		c			-456.7				
		c			-113.0				
K <sub>2</sub> S <sub>4</sub> ·4H <sub>2</sub> O K <sub>2</sub> S <sub>4</sub> ·2H <sub>2</sub> O		aq			-114.6				
		c			-150.5				
		c			-258.7				

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Table 93-14. Potassium (at. no., 19; at. wt., 39.096)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K kcal/mole			kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole	
K <sub>2</sub> SO <sub>3</sub>			c		-266.9				
			aq		-269.1				
K <sub>2</sub> SO <sub>4</sub>			c, II		-342.66				
		in 600 H <sub>2</sub> O	aq		-336.731				
		800 H <sub>2</sub> O	aq		-336.738				
		1000 H <sub>2</sub> O	aq		-336.748				
		2000 H <sub>2</sub> O	aq		-336.757				
		3000 H <sub>2</sub> O	aq		-336.776				
		5000 H <sub>2</sub> O	aq		-336.808				
		8000 H <sub>2</sub> O	aq		-336.830				
		10000 H <sub>2</sub> O	aq		-336.852				
		20000 H <sub>2</sub> O	aq		-336.876				
		50000 H <sub>2</sub> O	aq		-336.910				
		100000 H <sub>2</sub> O	aq		-336.930				
	500000 H <sub>2</sub> O	aq		-336.958					
	∞ H <sub>2</sub> O	aq		-336.98					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 93-15. Potassium (at. no., 19; at. wt., 39.096)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)			
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
$K_2S_2O_3$		aq		-274.			
$K_2S_2O_5$		c		-362.6			
		aq		-351.9			
$K_2S_2O_5 \cdot 1\frac{1}{2}H_2O$		c		-397.0			
		c		-413.6			
$K_2S_2O_6$		aq		-401.0			
	in 400 $H_2O$	aq		-400.8			
	600 $H_2O$						
		c		-458.3			
$K_2S_2O_8$		aq		-445.3			
		c		-422.			
$K_2S_4O_6$		aq		-410.			
$K_2S_5O_6$		c					
		aq		-401.			
$K_2S_5O_6 \cdot 1\frac{1}{2}H_2O$		c		-515.4			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 99-1b. Potassium (at. no., 19; at. wt., 39.096)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K		At 298.16°K (25°C)		
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
KHS	in 10 H <sub>2</sub> O	c		-83.2			
	20 H <sub>2</sub> O	aq		-64.35			
	30 H <sub>2</sub> O	aq		-64.21			
	40 H <sub>2</sub> O	aq		-64.16			
	50 H <sub>2</sub> O	aq		-64.05			
	100 H <sub>2</sub> O	aq		-64.00			
	200 H <sub>2</sub> O	aq		-63.99			
		aq		-64.1			
KHS·4H <sub>2</sub> O		c		-80.1			
KHSO <sub>3</sub>	in 385 H <sub>2</sub> O	aq		-209.7			
KHSO <sub>4</sub>	in 20 H <sub>2</sub> O	c		-276.8			
	25 H <sub>2</sub> O	aq		-273.38			
	50 H <sub>2</sub> O	aq		-273.36			
	100 H <sub>2</sub> O	aq		-273.37			
	200 H <sub>2</sub> O	aq		-273.44			
	400 H <sub>2</sub> O	aq		-273.59			
	800 H <sub>2</sub> O	aq		-273.90			
		aq		-274.3			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 93-17. Potassium (at. no., 19; at. wt., 39.096) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950							
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} f$
				At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole
KI·4SO <sub>2</sub>			c		-400.8		
K <sub>2</sub> Se			c		-79.3		
K <sub>2</sub> Se·9H <sub>2</sub> O			aq		-88.5		
K <sub>2</sub> Se·14H <sub>2</sub> O			c		-722.0		
K <sub>2</sub> Se·19H <sub>2</sub> O			c		-1066.		
			c		-1417.		
K <sub>2</sub> SeO <sub>4</sub>		in 440 H <sub>2</sub> O	aq		-265.3		
KHSe			c		-35.9		
			aq		-35.4		
KHSeO <sub>4</sub>		in 220 H <sub>2</sub> O	aq		-202.8		
K <sub>2</sub> TeO <sub>3</sub>			aq		-261.6		
K <sub>2</sub> TeO <sub>4</sub>			aq		-290.3		

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Substance		HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)				
Formula	Description	State	$\Delta H_f^\circ$		$\log_{10} K_f$	$C_p^\circ$
			At 0°K	At 298.16°K (25°C)		
			kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
KNO <sub>2</sub>	in 400 H <sub>2</sub> O	c				
		aq				
KNO <sub>3</sub>	std. state, hyp. m = 1	c				
	in 400 H <sub>2</sub> O	aq				
	500 H <sub>2</sub> O	aq				
	600 H <sub>2</sub> O	aq				
	800 H <sub>2</sub> O	aq				
	1000 H <sub>2</sub> O	aq				
	2500 H <sub>2</sub> O	aq				
	5000 H <sub>2</sub> O	aq				
	10000 H <sub>2</sub> O	aq				
	20000 H <sub>2</sub> O	aq				
	50000 H <sub>2</sub> O	aq				
	100000 H <sub>2</sub> O	aq				
	∞ H <sub>2</sub> O	aq				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 93-19. Potassium (at. no., 19; at. wt., 39.096)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K	At 298.16°K (25°C)				
				kcal/mole	kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
KNH <sub>2</sub>			c		-28.3				
KBr·4NH <sub>3</sub>			c		-166.5				
KI·NH <sub>3</sub>			c		-153.1				
KI·6NH <sub>3</sub>			c		-188.9				
KH <sub>2</sub> PO <sub>4</sub>			c		-374.9				
	in 35 H <sub>2</sub> O		aq		-370.70				
	50 H <sub>2</sub> O		aq		-370.63				
	100 H <sub>2</sub> O		aq		-370.51				
	200 H <sub>2</sub> O		aq		-370.40				
	300 H <sub>2</sub> O		aq		-370.34				
	755 H <sub>2</sub> O		aq		-370.4				
KH <sub>2</sub> AsO <sub>4</sub>			c		-271.5	-237.0	173.72	37.08	30.29
			aq		-276.2				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Substance		HEAT OF FORMATION AT 0°K: HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)				
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$C_p^\circ$
			At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole	cal/deg mole
$K_2CO_3$	in 50 $H_2O$	c		-273.93		
	100 $H_2O$	aq		-281.02		
	200 $H_2O$	aq		-281.03		
	400 $H_2O$	aq		-281.08		
	800 $H_2O$	aq		-281.20		
	1000 $H_2O$	aq		-281.37		
$K_2CO_3 \cdot \frac{1}{2}H_2O$ $K_2CO_3 \cdot 1\frac{1}{2}H_2O$		aq		-281.56		
		c		-210.43		
		c		-283.40		
		c		-320.8		
$K_2C_2O_4$	potassium oxalate	aq		-316.88		
	in 100 $H_2O$	aq		-316.80		
	200 $H_2O$	aq		-316.77		
	300 $H_2O$	aq		-316.76		
	400 $H_2O$	aq		-316.76		
	500 $H_2O$	aq		-316.783		
	1000 $H_2O$	aq		-316.829		
	2500 $H_2O$	aq		-316.845		
	5000 $H_2O$	aq		-316.897		



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 93-21. Potassium (at. no., 19; at. wt., 39.096) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	Description		At 0°K kcal/mole	kcal/mole					
$K_2C_2O_4$	in 10000 $H_2O$	aq		-316.951					
	50000 $H_2O$	aq		-317.032					
	100000 $H_2O$	aq		-317.051					
	$\infty H_2O$	aq		-317.1					
$K_2C_2O_4 \cdot H_2O$		c		-392.17					
$KCHO_2$	potassium formate	c		-158.0					
		aq		-157.7					
$KHCO_3$	in 1500 $H_2O$	c		-229.3					
		aq		-224.5					
$KOCH_3$	potassium methylate in $CH_3OH$			-108.0					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 93-22. Potassium (at. no., 19; at. wt., 39.096) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
$KC_2H_3O_2$	potassium acetate in 1% $H_2O$	1% $H_2O$	c		-173.2				
					-175.6				
					-175.9				
					-176.08				
					-176.18				
					-176.40				
					-176.54				
					-176.64				
					-176.71				
					-176.76				
					-176.88				
					-215.2				
$KCN$	potassium cyanide	c	c		-116.0				
					-2833.				
					-26.90				
					-24.1				
$KCNO$	potassium isocyanate	c	c		-98.5				
					-93.5				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards SERIES I Washington, D.C.

Table 93-23. Potassium (at. no., 19; at. wt., 39.096) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950										
Substance				State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula	Description	At 0°K kcal/mole	kcal/mole		At 298.16°K (25°C) kcal/mole	cal/deg mole				
KCO <sub>2</sub> CONH <sub>2</sub>	potassium oxamate	in	aq		-214.5					
		2 H <sub>2</sub> O	c		-48.62					
KCNS		3 H <sub>2</sub> O	aq		-45.46					
		4 H <sub>2</sub> O	aq		-45.22					
		5 H <sub>2</sub> O	aq		-45.000					
		6 H <sub>2</sub> O	aq		-44.760					
		8 H <sub>2</sub> O	aq		-44.600					
		10 H <sub>2</sub> O	aq		-44.320					
		15 H <sub>2</sub> O	aq		-44.110					
		20 H <sub>2</sub> O	aq		-43.780					
		25 H <sub>2</sub> O	aq		-43.570					
		50 H <sub>2</sub> O	aq		-43.425					
		100 H <sub>2</sub> O	aq		-43.105					
		200 H <sub>2</sub> O	aq		-42.925					
		400 H <sub>2</sub> O	aq		-42.830					
		∞ H <sub>2</sub> O	aq		-42.78					
	KCN· $\frac{1}{2}$ SO <sub>2</sub> KCN·SO <sub>2</sub>		aq		-42.8					
			c		-69.7					
		c		-130.2						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 93-24. Potassium (at. no., 19; at. wt., 39.096)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°C)			
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
$K_2SiF_6$		c		-671.			
$K_2SnCl_4 \cdot H_2O$		c		-365.3			
$K_2SnCl_6$		c		-362.9			
		aq		-360.0			
$KCl \cdot PbCl_2 \cdot \frac{3}{2}H_2O$		c		-214.1			
$KCl \cdot 2PbCl_2$		c		-278.1			
$2KI \cdot PbI_2$		c		-199.0			
$2KI \cdot PbI_2 \cdot 2H_2O$		c		-340.4			
$4KI \cdot 3PbI_2$		c		-437.1			
$4KI \cdot 3PbI_2 \cdot 6H_2O$		c		-859.4			
$K_2SO_4 \cdot PbSO_4$		c		-567.4			
$K_2SO_4 \cdot ZnSO_4$		c		-580.31			
$K_2SO_4 \cdot ZnSO_4 \cdot 2H_2O$		c		-724.6			
$K_2SO_4 \cdot ZnSO_4 \cdot 6H_2O$		c		-1011.8			
$2KCN \cdot Zn(CN)_2$		c		-44.2			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 93-25. Potassium (at. no., 19; at. wt., 39.098)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		State	Description	At 298.16°K (25°C)					
Formula	$\Delta H_f^\circ$			$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
	At 0°K kcal/mole							kcal/mole	kcal/mole
KHg		c		-11.6					
KHg <sub>2</sub>		c		-17.7					
KHg <sub>12</sub>		c		-32.83					
KHg <sub>50</sub>		liq		-26.06					
KHg <sub>100</sub>		liq		-26.08					
KHg <sub>200</sub>		liq		-26.30					
KHg <sub>1000</sub>		liq		-26.39					
KHgCl <sub>3</sub>		c		-160.9					
KHgCl <sub>3</sub> ·H <sub>2</sub> O		aq		-152.2					
K <sub>2</sub> HgCl <sub>4</sub>		c		-231.1					
		c		-267.6					
K <sub>2</sub> HgCl <sub>4</sub> ·H <sub>2</sub> O		aq		-253.3					
		c		-337.6					
4KCl·3HgCl <sub>2</sub>		c		-590.3					
4KCl·3HgCl <sub>2</sub> ·3H <sub>2</sub> O		c		-801.2					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 93-26. Potassium (at. no., 19; at. wt., 39.096) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K					
				kcal/mole	kcal/mole				
KHBr <sub>3</sub>			c		-137.3				
KHgBr <sub>3</sub> ·H <sub>2</sub> O			aq		-127.3				
			c		-209.1				
KHaI <sub>3</sub>			c		-105.7				
KHgI <sub>3</sub> ·H <sub>2</sub> O			c		-174.2				
K <sub>2</sub> HgI <sub>4</sub>			c		-184.9				
			aq		-175.7				
KHg(CN) <sub>3</sub>			aq		34.9				
K <sub>2</sub> Hg(CN) <sub>4</sub>			c		-7.7				
			aq		5.5				
KCN·2Hg(CN) <sub>2</sub>			aq		100.8				
KCl·Hg(CN) <sub>2</sub>			c		-42.3				
KCl·Hg(CN) <sub>2</sub> ·H <sub>2</sub> O			c		-112.2				
KBr·Hg(CN) <sub>2</sub>			c		-34.5				
			aq		-23.3				
KBr·Hg(CN) <sub>2</sub> · $\frac{1}{2}$ H <sub>2</sub> O			c		-138.0				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 93-27. Potassium (at. no., 19; at. wt., 39.096) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	Description	State	At 0°K			At 298.16°K (25°C)		
				kcal/mole			kcal/mole		
				$\Delta H^{\circ}$	$\Delta F^{\circ}$	$\log_{10} K$	$S^{\circ}$	$C_p^{\circ}$	
$KI \cdot Hg(CN)_2 \cdot 4H_2O$			c	-21.7					
			aq	-10.1					
			c	-39.2					
$KCl \cdot CuCl_2$			c	-153.0					
$2KCl \cdot CuCl$			c	-242.3					
$2KCl \cdot CuCl_2$			c	-261.6					
$KCl \cdot CuCl_2 \cdot 2H_2O$			c	-400.7					
$K_2^{2+} \cdot SO_4^{2-}$			c	-547.4					
			c	-550.8					
			c	-548.7					
			c	-694.8					
			c	-980.7					
$K_2^{2+} \cdot SO_4^{2-} \cdot 4H_2O$			c, V	-418.6					
$K_2^{2+} \cdot SO_4^{2-} \cdot 6H_2O$			c, IV	-418.1					
$K_2^{2+} \cdot SO_4^{2-}$			c, II	-416.					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 93-28. Potassium (at. no., 19; at. wt., 39.096)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°K)

March 31, 1950

Substance		$\Delta H_f^\circ$	$\Delta G_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State	At 0°K	At 298.16°K (25°K)		
			kcal/mole	kcal/mole	kcal/mole	cal/deg mole
KCl·AgCl		c		-134.54		
KBr·AgBr		c		-117.11		
3KBr·AgBr· $\frac{1}{2}$ H <sub>2</sub> O		c		-337.53		
KI·AgI		c		-91.42		
KI·AgI· $\frac{1}{2}$ H <sub>2</sub> O		c		-109.95		
2KI·AgI· $\frac{1}{2}$ H <sub>2</sub> O		c		-205.69		
3KI·AgI		c		-248.94		
3KI·AgI· $\frac{1}{2}$ H <sub>2</sub> O		c		-282.8		
3KI·2AgI·H <sub>2</sub> O		c		-332.37		
KAg(CN) <sub>2</sub>		c		-3.9		
K <sub>2</sub> Ag(CN) <sub>3</sub>		aq		4.2		
KAuCl <sub>4</sub>		aq		-21.4		
KAu(CN) <sub>2</sub>		c				
KAu(CN) <sub>2</sub>		aq		-2.1		



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards

Washington, D.C.

Table 93-29. Potassium (at. no., 19; at. wt., 39.096)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	Description	At 298.16°K (25°C)				
Formula	ΔHf°			log10ff	S°	Cp°		
	At 0°K						At 0°K	
				kcal/mole	kcal/mole		cal/deg mole	
K2PtCl4		c		-254.2				
K2PtCl6		aq		-242.8				
	in 9400 H2O	c		-301.0	194.31	79.8	49.1	
K2PtBr4		aq		-287.5				
	in 1000 H2O	c		-220.7				
K2PtBr6		aq		-210.7				
	in 2000 H2O	c		-248.6				
		aq		-236.9				
K2IrCl6		c		-286.				
		aq		-274.				
K3IrCl6		c		-374.				
		aq		-369.				
K2OsCl6		c		-280.				
KReO4		c		-264.02				
	in 10000 H2O	aq		-250.3				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards SERIES I Table 93-30. Potassium (at. no., 19; at. wt., 39.096) Washington, D.C.

Substance		HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)				
Formula	Description	State	$\Delta H_f^{\circ}$	$\Delta F_f^{\circ}$	$\log_{10} K_f$	$C_p^{\circ}$
			At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	At 298.16°K (25°C) cal/deg mole	At 298.16°K (25°C) cal/deg mole
$K_2PdCl_4$		c		-261.6		
		aq		-248.4		
$K_2PdCl_6$		c		-283.8		
		aq		-272.		
$K_2PtBr_4$		c		-217.2		
		aq		-205.6		
$K_3PtCl_6$		c		-343.		
$K_2Ni(CN)_4$		aq		-33.3		
$K_2CO_3 \cdot CaCO_3$		c		-450.5		
$K_2CO_3 \cdot CaCO_3 \cdot 4H_2O$		c		-735.2		
$KFe(SO_4)_2$		aq		-496.3		
$K_2Fe(SO_4)_2$		aq		-572.9		
$K_2Fe(SO_4)_2 \cdot 4H_2O$		c		-836.		
$K_2Fe(SO_4)_2 \cdot 6H_2O$		c		-972.1		

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

Washington, D.C.

National Bureau of Standards

Table 93-31. Potassium (at. no., 19; at. wt., 39.096) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		State		$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	State		At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole		cal/deg mole	
$K_3Fe(CN)_6$		c			-41.4				80.3
$K_4Fe(CN)_6$		aq			-28.2				
		c			-125.1				
		aq			-113.5				
$K_4Fe(CN)_6 \cdot 3H_2O$		c			-334.3				
$KH_2Fe(CN)_6$		aq			92.3				
$KH_3Fe(CN)_6$		aq			67.1				
$K_2HFe(CN)_6$		aq			31.6				
$K_2H_2Fe(CN)_6$		aq			7.2				
$K_3HFe(CN)_6$		aq			-53.1				
$K_3FeCO(CN)_5$		c			-137.4				
		aq			-133.0				
$K_3FeCO(CN)_5 \cdot 3\frac{1}{2}H_2O$		c			-382.5				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 93-32. Potassium (at. no., 19; at. wt., 39.096)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		State	Description	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	kcal/mole			kcal/mole	kcal/mole	At 298.16°K (25°C)	cal/deg mole	cal/deg mole
KMnO <sub>4</sub>	in 140 H <sub>2</sub> O	c		-194.4	-170.6	125.05	41.04	28.5
	175 H <sub>2</sub> O	aq		-184.44				
	220 H <sub>2</sub> O	aq		-184.32				
	660 H <sub>2</sub> O	aq		-184.29				
	1000 H <sub>2</sub> O	aq		-184.04				
	4000 H <sub>2</sub> O	aq		-183.98				
K <sub>2</sub> Mn(SO <sub>4</sub> ) <sub>2</sub>		aq		-183.9				
		c		-597.7				
		c		-743.7				
K <sub>2</sub> Mn(SO <sub>4</sub> ) <sub>2</sub> ·2H <sub>2</sub> O		c		-884.0				
		c		-330.49				
		c		-328.2				
K <sub>2</sub> CrO <sub>4</sub>	in 17.2 H <sub>2</sub> O	aq		-327.88				
	21.5 H <sub>2</sub> O	aq		-327.63				
	27 H <sub>2</sub> O	aq		-327.32				
	36 H <sub>2</sub> O	aq		-327.00				
	54 H <sub>2</sub> O	aq		-326.00				
	1080 H <sub>2</sub> O	aq		-325.79				
	1700 H <sub>2</sub> O	aq		-325.7				
	2000 H <sub>2</sub> O	aq		-325.7				
	∞ H <sub>2</sub> O	aq		-326.0				
		aq						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 93-33. Potassium (at. no., 19; at. wt., 39.096)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	in			At 0°K					
				kcal/mole	kcal/mole				
$K_2Cr_2O_7$			c		-485.90				
	135 H <sub>2</sub> O		aq		-470.41				
	150 H <sub>2</sub> O		aq		-470.36				
	200 H <sub>2</sub> O		aq		-470.22				
	300 H <sub>2</sub> O		aq		-470.05				
	400 H <sub>2</sub> O		aq		-469.95				
	800 H <sub>2</sub> O		aq		-469.50				
	1000 H <sub>2</sub> O		aq		-469.22				
	1500 H <sub>2</sub> O		aq		-468.87				
	2000 H <sub>2</sub> O		aq		-468.7				
$K_2Cr_2O_7 \cdot CrO_3$			c		-598.				
$KCl \cdot CrO_3$			c		-245.5				
$KCr(SO_4)_2$			c		-562.				
$KCr(SO_4)_2 \cdot H_2O$			aq		-553.8				
$KCr(SO_4)_2 \cdot 2H_2O$			c		-642.				
$KCr(SO_4)_2 \cdot 6H_2O$			c		-720.				
$KCr(SO_4)_2 \cdot 12H_2O$			c		-959.3				
			c		-1383.1				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

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Table 93-34. Potassium (at. no., 19; at. wt., 39.096)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description						
$\text{KNH}_4\text{CrO}_4$  $\text{K}_2\text{MnO}_4$  $\text{KVO}_3$ $\text{KVO}_4$  $\text{KVO}_5$  $3\text{KF} \cdot \text{AlF}_3$  $3\text{KF} \cdot \text{AlF}_3 \cdot 3\frac{1}{2}\text{H}_2\text{O}$ $\text{KCl} \cdot \text{AlCl}_3$ $3\text{KCl} \cdot \text{AlCl}_3$ $3\text{KCl} \cdot 2\text{AlCl}_3$	  						

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 93-35. Potassium (at. no., 19; at. wt., 39.096)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	Description	At 298.16°K (25°C)				
Formula	$\Delta H_f^\circ$			$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
								kcal/mole
KAl(SO <sub>4</sub> ) <sub>2</sub>		c		-589.24	-534.29	391.627	48.9	46.12
KAl(SO <sub>4</sub> ) <sub>2</sub> ·H <sub>2</sub> O		aq		-616.9				
KAl(SO <sub>4</sub> ) <sub>2</sub> ·2H <sub>2</sub> O		c		-667.5				
KAl(SO <sub>4</sub> ) <sub>2</sub> ·3H <sub>2</sub> O		c		-742.				
KAl(SO <sub>4</sub> ) <sub>2</sub> ·12H <sub>2</sub> O		c		-814.				
KAl(SO <sub>4</sub> ) <sub>2</sub> ·12H <sub>2</sub> O		c		-1447.74	-1227.8	699.96	164.3	155.6
KCl·AlCl <sub>3</sub> ·6NH <sub>3</sub>		c		-472.4				
K <sub>2</sub> O·Al <sub>2</sub> O <sub>3</sub> ·4SiO <sub>2</sub>	leucite	c		-1406.4				
K <sub>2</sub> O·Al <sub>2</sub> O <sub>3</sub> ·6SiO <sub>2</sub>	microcline	gls		-1395.0				
	adularia	c		-1816.				
		c		-1842.				
		gls		-1780.				
2KCl·UO <sub>2</sub> Cl <sub>2</sub> ·2H <sub>2</sub> O		c		-665.				
KCl·ThCl <sub>4</sub> ·9H <sub>2</sub> O		c		-1055.0				
2KCl·ThCl <sub>4</sub>		c		-502.5				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards Washington, D.C.

Table 93-36. Potassium (at. no., 19; at. wt., 39.096) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)										
March 31, 1950										
Substance					State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	Description	$\Delta H_f^\circ$								
			At 0°K	At 298.16°K (25°C)						
			kcal/mole	kcal/mole						
KCl·MgCl <sub>2</sub>					c					
KCl·MgCl <sub>2</sub> ·6H <sub>2</sub> O	carallite			-260.75	c					
2KCl·MgCl <sub>2</sub>	fresh melt			-702.0	c					
4KCl·MgCl <sub>2</sub>	fresh melt			-364.36	c					
				-574.76	c					
K <sub>2</sub> SO <sub>4</sub> ·MgSO <sub>4</sub>					c					
				-651.0	c					
K <sub>2</sub> SO <sub>4</sub> ·MgSO <sub>4</sub> ·2H <sub>2</sub> O	fresh melt			-655.2	c					
K <sub>2</sub> SO <sub>4</sub> ·MgSO <sub>4</sub> ·H <sub>2</sub> O				-798.6	c					
K <sub>2</sub> SO <sub>4</sub> ·MgSO <sub>4</sub> ·5H <sub>2</sub> O				-946.4	c					
K <sub>2</sub> SO <sub>4</sub> ·MgSO <sub>4</sub> ·H <sub>2</sub> O				-1012.4	c					
K <sub>2</sub> SO <sub>4</sub> ·MgSO <sub>4</sub> ·H <sub>2</sub> O				-1082.8	c					
K <sub>2</sub> SO <sub>4</sub> ·2MgSO <sub>4</sub>	langbeinite			-985.8	c					
KCl·MgSO <sub>4</sub>	kainite			-424.5	c					
2KCl·CaCl <sub>2</sub>	fresh melt				c					
				-400.9	c					
K <sub>2</sub> SO <sub>4</sub> ·CaSO <sub>4</sub> ·H <sub>2</sub> O				-757.96	c					
K <sub>2</sub> SO <sub>4</sub> ·5CaSO <sub>4</sub> ·H <sub>2</sub> O				-2133.6	c					
KCaFeO(CN) <sub>5</sub> ·5H <sub>2</sub> O				933.	c					



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards Washington, D.C.

Table 93-37. Potassium (at. no., 19; at. wt., 39.096) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		State	At 298.16°K (25°C)				log <sub>10</sub> K <sub>f</sub>	S°	C <sub>p</sub> °
Formula	Description		ΔH <sub>f</sub> ° kcal/mole	ΔH <sub>f</sub> ° kcal/mole	ΔF <sub>f</sub> ° kcal/mole	ΔF <sub>f</sub> ° kcal/mole			
K <sub>2</sub> SO <sub>4</sub> ·SrSO <sub>4</sub>		c							
KI·LiCl	fresh melt	c							
KNa		g	33.6						
KNa <sub>2</sub>		liq							
K <sub>2</sub> Na		liq							
K <sub>3</sub> Na		liq							
KCl·NaCl	fresh melt	c							
KI·NaCl	fresh melt	c							
KNa <sub>2</sub> PO <sub>4</sub>	in 330 H <sub>2</sub> O	aq							
3KCNS·NaCNS		c							

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 94-1. Rubidium (at. no., 37; at. wt., 85.46) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula				At 0°K kcal/mole	kcal/mole				
Rb			g		20.51	13.35	-9.785	40.628	4.9680
Rb <sup>+</sup>		std. state, hyp. m = 1	c, I	0.000	0.000	0.000	0.0000	16.6	7.27
Rb <sup>++</sup>			aq		-58.9	-67.45	47.974	29.7	
Rb <sup>+++</sup>			g		118.297				
Rb <sub>2</sub>			g		753.95				
			g		1670.4				
			g		29.6				
Rb <sub>2</sub> O			c		-78.9				
Rb <sub>2</sub> O <sub>2</sub>			c		-101.7				
Rb <sub>2</sub> O <sub>3</sub>			c		-116.7				
Rb <sub>2</sub> O <sub>4</sub>			c		-126.2				
RbH			g		33.				
RbOH			q, II		-98.9				
		std. state, hyp. m = 1	aq		-113.9	-105.05	77.000	27.2	
		in 5.2 H <sub>2</sub> O	aq		-110.5				
		200 H <sub>2</sub> O	aq		-113.7				
RbOH·H <sub>2</sub> O			c		-177.8				
RbOH·2H <sub>2</sub> O			c		-250.8				

March 31, 1950

Substance		State	$\Delta H^\circ$ kcal/mole	$\Delta F^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole
Formula	Description						
RbF	std. state, hyp. m=1	c					12.2
	in 100 H <sub>2</sub> O	aq		-131.28			
	200 H <sub>2</sub> O	aq		-137.6	-133.53	97.876	27.4
	500 H <sub>2</sub> O	aq		-137.45			
	1000 H <sub>2</sub> O	aq		-137.476			
	2000 H <sub>2</sub> O	aq		-137.512			
	5000 H <sub>2</sub> O	aq		-137.534			
	10000 H <sub>2</sub> O	aq		-137.551			
	20000 H <sub>2</sub> O	aq		-137.567			
	50000 H <sub>2</sub> O	aq		-137.576			
	100000 H <sub>2</sub> O	aq		-137.583			
		aq		-137.588			
		aq		-137.591			
		aq		-137.6			
RbH <sub>2</sub>		c		-156.12			
		c		-240.33			
		c		-217.3			
		aq		-212.5			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 94-3. Rubidium (at. no., 37; at. wt., 85.48) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	Description	State	At 298.16°K (25°C)					
				$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
				At 0°K kcal/mole	kcal/mole		cal/deg mole	cal/deg mole	
RbCl	std. state, hyp. m = 1 in 100 H <sub>2</sub> O 200 H <sub>2</sub> O 500 H <sub>2</sub> O 1000 H <sub>2</sub> O 2000 H <sub>2</sub> O 5000 H <sub>2</sub> O 10000 H <sub>2</sub> O 20000 H <sub>2</sub> O 50000 H <sub>2</sub> O $\infty$ H <sub>2</sub> O in CH <sub>3</sub> OH	c aq aq aq aq aq aq aq aq aq aq aq			-102.91				12.3
					-98.9	72.419	42.9		
					-98.90				
					-98.840				
					-98.836				
					-98.840				
					-98.849				
					-98.863				
					-98.872				
					-98.880				
					-98.887				
					-98.9				
					-98.6				
RbClO <sub>3</sub>	std. state, hyp. m = 1		c		-69.8	51.16	36.3		24.66
					-82.4	49.894	68.7		
RbClO <sub>4</sub>	std. state, hyp. m = 1		c, II		-103.87	53.647	38.4		
					-90.3	51.324	73.2		

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards  
 SERIES I  
 Table 94-4. Rubidium (at. no., 37; at. wt., 85.48)  
 HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
 March 31, 1950  
 Washington, D.C.

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole	kcal/mole		cal/deg mole	cal/deg mole
RbBr	std. state, hyp. m = 1		c						
	in 200 H <sub>2</sub> O		aq		-93.03	-90.38	66.247	25.88	12.68
	500 H <sub>2</sub> O		aq		-87.8	-92.02	67.449	49.0	
	1000 H <sub>2</sub> O		aq		-87.766				
	2000 H <sub>2</sub> O		aq		-87.745				
	5000 H <sub>2</sub> O		aq		-87.748				
	10000 H <sub>2</sub> O		aq		-87.754				
	20000 H <sub>2</sub> O		aq		-87.766				
	50000 H <sub>2</sub> O		aq		-87.774				
	$\infty$ H <sub>2</sub> O		aq		-87.781				
RbI	std. state, hyp. m = 1		c						
			aq		-87.787	-77.8	57.03	28.21	12.50
			aq		-87.8	-79.80	58.492	55.8	
Rb <sub>2</sub> S			c						
	in 50% H <sub>2</sub> O		aq		-83.2				
					-107.8				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards Washington, D.C.

Table 94-5. Rubidium (at. no., 37; at. wt., 85.46)									
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)									
March 31, 1950									
Substance		Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
Formula				At 0°K	At 298.16°K (25°C)			cal/deg mole	
				kcal/mole	kcal/mole	kcal/mole	cal/deg mole		
Rb <sub>2</sub> SO <sub>4</sub>		std. state, hyp. m = 1	c						
		in 400 H <sub>2</sub> O	aq		-340.50				
		500 H <sub>2</sub> O	aq		-334.7	228.866	33.8		
		800 H <sub>2</sub> O	aq		-334.508				
		1000 H <sub>2</sub> O	aq		-334.489				
		2000 H <sub>2</sub> O	aq		-334.478				
		3000 H <sub>2</sub> O	aq		-334.479				
		5000 H <sub>2</sub> O	aq		-334.494				
		10000 H <sub>2</sub> O	aq		-334.507				
		20000 H <sub>2</sub> O	aq		-334.529				
		50000 H <sub>2</sub> O	aq		-334.565				
		100000 H <sub>2</sub> O	aq		-334.598				
		∞ H <sub>2</sub> O	aq		-334.637				
			aq		-334.652				
			aq		-334.677				
RbHS		c		-62.4					
RbHSO <sub>4</sub>		aq		-63.1					
RbI·4SO <sub>2</sub>		c		-273.7					
		aq		-270.4					
		c		-402.6					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards Washington, D.C.

Table 94-6. Rubidium (at. no., 37; at. wt., 85.48)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance		State	Description	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	State			At 0°K					
				kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
RbH <sub>2</sub> Se		c			-35.5				
		aq			-34.3				
RbNO <sub>3</sub>		c, II			-117.04				
	std. state, hyp. m = 1	aq			-108.3		68.798	64.7	
	in 60 H <sub>2</sub> O	aq			-109.00				
	100 H <sub>2</sub> O	aq			-108.76				
	200 H <sub>2</sub> O	aq			-108.54				
	400 H <sub>2</sub> O	aq			-108.40				
	500 H <sub>2</sub> O	aq			-108.36				
	800 H <sub>2</sub> O	aq			-108.316				
	1000 H <sub>2</sub> O	aq			-108.305				
	2000 H <sub>2</sub> O	aq			-108.280				
	5000 H <sub>2</sub> O	aq			-108.277				
	10000 H <sub>2</sub> O	aq			-108.278				
	50000 H <sub>2</sub> O	aq			-108.287				
	∞ H <sub>2</sub> O	aq			-108.3				
	RbOH		c			-25.7			
LiClO <sub>4</sub> ·5RbCl <sub>3</sub>		c			-148.0				
60I <sub>2</sub> ·RbI <sub>2</sub>		c			-190.7				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards Washington, D.C.

Table 94-7. Rubidium (at. no., 37; at. wt., 85.46) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance		State	$\Delta H_f^\circ$		$\Delta H_f^\circ$		$\Delta F_f^\circ$		$C_p^\circ$
Formula	Description		At 0°K kcal/mole	kcal/mole	At 298.16°K (25°C) kcal/mole	cal/deg mole	log <sub>10</sub> K <sub>f</sub>	S° cal/deg mole	
Rb <sub>2</sub> CO <sub>3</sub>	in 200 H <sub>2</sub> O 2000 H <sub>2</sub> O	c aq aq							
Rb <sub>2</sub> CO <sub>3</sub> ·H <sub>2</sub> O		c							
Rb <sub>2</sub> CO <sub>3</sub> · $\frac{1}{2}$ H <sub>2</sub> O		c							
Rb <sub>2</sub> CO <sub>3</sub> · $\frac{3}{2}$ H <sub>2</sub> O		c							
RbHCO <sub>3</sub>		c							
3Rb <sub>2</sub> CO <sub>3</sub> ·2RbHCO <sub>3</sub> ·4 $\frac{1}{2}$ H <sub>2</sub> O		c							
RbCNS	rubidium thiocyanate	c							
RbCNS· $\frac{1}{2}$ SO <sub>2</sub>		aq							
Rb <sub>2</sub> SiF <sub>6</sub>		c							
2RbI·PbI <sub>2</sub>		c							
2RbI·PbI <sub>2</sub> ·4H <sub>2</sub> O		c							



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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## SERIES I

Washington, D.C.

Table 94-6. Rubidium (at. no., 37; at. wt., 85.48)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		$\Delta H_f^\circ$ At 0°K kcal/mole	$\Delta H_f^\circ$ kcal/mole	$\Delta F_f^\circ$ kcal/mole	$\log_{10} K_f$ At 298.16°K (25°C)	$S^\circ$ cal/deg mole	$C_p^\circ$
Formula	Description						
$Rb_2CuCl_4$	c		-261.0				
$Rb_2CuCl_4 \cdot 2H_2O$							
$Rb_2IrCl_6$	c		-290.9				
$RbReO_4$	c		-256.9				
	aq		-249.2				
$RbAl(SO_4)_2$	c		-569.				
$RbAl(SO_4)_2 \cdot H_2O$	c		-647.				
$RbAl(SO_4)_2 \cdot 2H_2O$	c		-722.				
$RbAl(SO_4)_2 \cdot 3H_2O$	c		-797.				
$RbAl(SO_4)_2 \cdot 11H_2O$	c		-1448.0				
$2RbCl \cdot ThCl_4$	c		-512.1				
$2RbCl \cdot ThCl_4 \cdot 9H_2O$	c		-1154.6				
$4RbCl \cdot ThCl_4$	c		-725.9				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 95-1. Cesium (at. no., 55; at. wt., 132.91)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	Description	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
Formula	At 0°K			At 298.16°K (25°C)				
				kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole
$f^-$		g		18.83	12.24	-8.972	41.944	4.9680
$Cs^+$		c		0.000	0.000	0.0000	19.8	7.42
		g		110.081				
$Cs^{++}$	std. state, hyp. m = 1	aq		-59.2	-67.41	49.411	31.6	
$=$		g		*52.5				
		g		27.0				
$Cs_2O$		c		-75.9				
$Cs_2O_2$		c		-96.2				
$Cs_2O_3$		c		-111.2				
$Cs_2O_4$		c		-124.2				
$CSH$		g		29.0	24.3	-17.81	51.25	7.133
$CSOH$		c, II		-97.2				
	std. state, hyp. m = 1	aq		-114.2	-105.00	76.964	29.3	
	in 2.2 $H_2O$	aq		-111.5				
	20% $H_2O$	aq		-114.0				
	$H_2O$	aq		-114.2				
$CSOH \cdot H_2O$		c		-186.9				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 95-2. Cesium (at. no., 55; at. wt., 132.91) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Description	State	Substance				At 298.16°K (25°C)		
			$\Delta H_f^\circ$ kcal/mole	$\Delta G_f^\circ$ kcal/mole	$\Delta F^\circ$ kcal/mole	$\log_{10} K_f$	$S^\circ$ cal/deg mole	$C_p^\circ$ cal/deg mole	
CsF	std. state, hyp. m=1	c							
	in 500 H <sub>2</sub> O	aq		-126.9	-133.49	97.846	29.5		
	1000 H <sub>2</sub> O	aq		-135.9					
	5000 H <sub>2</sub> O	aq		-135.85					
	10000 H <sub>2</sub> O	aq		-135.86					
	100000 H <sub>2</sub> O	aq		-135.87					
	$\infty$ H <sub>2</sub> O	aq		-135.86					
		aq		-135.89					
		aq		-135.9					
		c		-176.8					
CsF· $\frac{3}{2}$ H <sub>2</sub> O		c		-237.2					
CsF·1½H <sub>2</sub> O		c							
CsHF <sub>2</sub>		c		-216.1					
		aq		-212.8					

## SERIES I

Washington, D.C.

Table 95-3. Cesium (at. no., 55; at. wt., 132.91)  
HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Substance								
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$		
			At 298.16°K (25°C)					
			At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole	
CsCl	c, II							
	std. state, hyp. m = 1	aq	-103.5					
	in 50 H <sub>2</sub> O	aq	-99.2	-88.76	65.060	45.0		
	75 H <sub>2</sub> O	aq	-99.55					
	100 H <sub>2</sub> O	aq	-99.46					
	200 H <sub>2</sub> O	aq	-99.346					
	500 H <sub>2</sub> O	aq	-99.205					
	1000 H <sub>2</sub> O	aq	-99.152					
	2000 H <sub>2</sub> O	aq	-99.150					
	5000 H <sub>2</sub> O	aq	-99.155					
	10000 H <sub>2</sub> O	aq	-99.165					
	50000 H <sub>2</sub> O	aq	-99.174					
	$\infty$ H <sub>2</sub> O	aq	-99.188					
		aq	-99.2					
CsClO <sub>4</sub>	c		-103.86	-73.28	53.713	41.89		
	std. state, hyp. m = 1	aq	-90.6	-69.98	51.294	75.3		

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 95-4. Cesium (at. no., 55; at. wt., 132.91) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Substance									
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$		
			At 0°K kcal/mole	kcal/mole	kcal/mole	cal/deg mole	cal/deg mole		
CsBr	std. state, hyp. m = 1	c							
	in 500 H <sub>2</sub> O	aq		-94.3	67.142	29.	12.4		
	1000 H <sub>2</sub> O	aq		-88.1	67.420	51.1			
	2000 H <sub>2</sub> O	aq		-88.065					
	5000 H <sub>2</sub> O	aq		-88.060					
	10000 H <sub>2</sub> O	aq		-88.062					
	50000 H <sub>2</sub> O	aq		-88.069					
	$\infty$ H <sub>2</sub> O	aq		-88.076					
CsI	std. state, hyp. m = 1	c							
	in 500 H <sub>2</sub> O	aq		-88.088					
	1000 H <sub>2</sub> O	aq		-88.1					
	2000 H <sub>2</sub> O	aq		-80.5	58.419	31.	12.4		
	5000 H <sub>2</sub> O	aq		-72.6	58.463	57.9			
	10000 H <sub>2</sub> O	aq		-72.57					
	50000 H <sub>2</sub> O	aq		-72.56					
	$\infty$ H <sub>2</sub> O	aq		-72.57					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

National Bureau of Standards Washington, D.C.

Table 95-5. Cesium (at. no., 55; at. wt., 132.91) HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C) March 31, 1950									
Formula	Substance	State	$\Delta H_f^\circ$		$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 0°K kcal/mole	kcal/mole					
Cs <sub>2</sub> S	c	aq		-61.1					
				-108.4					
Cs <sub>2</sub> SO <sub>4</sub>	c, II	aq		-339.38					
		aq		-335.3	-312.16	228.809	67.7		
		aq		-335.261					
		aq		-335.191					
		aq		-335.176					
		aq		-335.145					
		aq		-335.138					
		aq		-335.157					
		aq		-335.184					
		aq		-335.211					
		aq		-335.239					
		aq		-335.268					
CsHS	c	aq		-335.3					
		aq		-62.9					
CsHSO <sub>4</sub>	c	aq		-63.2					
		aq		-274.0					
	in 300 H <sub>2</sub> O	aq		-270.7					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 95-6. Cesium (at. no., 55; at. wt., 132.91)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
March 31, 1950

Formula	Substance	Description	State	$\Delta H_f^\circ$	$\Delta G_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$
				At 0°K kcal/mole	At 298.16°K (25°C) kcal/mole		cal/deg mole	cal/deg mole
CsI·4SO <sub>2</sub>			c					
CsHSe			c					
			aq					
CsNO <sub>3</sub>			c, II					
	std. state, hyp. m = 1		aq					
	in 60 H <sub>2</sub> O		aq					
	100 H <sub>2</sub> O		aq					
	200 H <sub>2</sub> O		aq					
	400 H <sub>2</sub> O		aq					
	500 H <sub>2</sub> O		aq					
	1000 H <sub>2</sub> O		aq					
	2000 H <sub>2</sub> O		aq					
	5000 H <sub>2</sub> O		aq					
	10000 H <sub>2</sub> O		aq					
	20000 H <sub>2</sub> O		aq					
	100000 H <sub>2</sub> O		aq					
	∞ H <sub>2</sub> O		aq					
CsNH <sub>2</sub>			c					

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

Washington, D.C.

Table 95-7. Cesium (at. no., 55; at. wt., 132.91)

HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)

March 31, 1950

Substance		State	At 298.16°K (25°C)				C <sub>p</sub> <sup>o</sup>
Formula	Description		ΔH <sub>f</sub> <sup>o</sup> kcal/mole	ΔF <sub>f</sub> <sup>o</sup> kcal/mole	log <sub>10</sub> K <sub>f</sub>	S <sup>o</sup> cal/deg mole	
Cs <sub>2</sub> CO <sub>3</sub>	in 100 H <sub>2</sub> O 2(100 H <sub>2</sub> O)	c	-267.4				
		aq	-280.0				
		aq	-278.6				
Cs <sub>2</sub> CO <sub>3</sub> ·3 $\frac{1}{2}$ H <sub>2</sub> O		c	-522.3				
CsHCO <sub>3</sub>		c	-228.4				
		aq	-224.4				
5Cs <sub>2</sub> CO <sub>3</sub> ·2CsHCO <sub>3</sub> ·10H <sub>2</sub> O		c	-2538.3				
5Cs <sub>2</sub> CO <sub>3</sub> ·2CsHCO <sub>3</sub> ·17 $\frac{1}{2}$ H <sub>2</sub> O		c	-3069.1				
Cs <sub>2</sub> SiF <sub>6</sub>		c	-669.5				
Cs <sub>2</sub> CuCl <sub>4</sub>		c	-263.5				
Cs <sub>2</sub> CuCl <sub>4</sub> ·2H <sub>2</sub> O		c	-404.3				
CsReO <sub>4</sub>		c	-257.2				
		aq	-249.5				



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards  
 SERIES I  
 Washington, D.C.

Table 95-8. Cesium (at. no., 55; at. wt., 132.91)  
 HEAT OF FORMATION AT 0°K; HEAT, FREE ENERGY, AND EQUILIBRIUM CONSTANT OF FORMATION, ENTROPY, AND HEAT CAPACITY, AT 298.16°K (25°C)  
 March 31, 1950

Substance									
Formula	Description	State	$\Delta H_f^\circ$	$\Delta F_f^\circ$	$\Delta F_f^\circ$	$\log_{10} K_f$	$S^\circ$	$C_p^\circ$	
			At 0°K kcal/mole			At 298.16°K (25°C)			
				kcal/mole	kcal/mole		cal/deg mole	cal/deg mole	
CsAl(SO <sub>4</sub> ) <sub>2</sub>		c			-570.				
CsAl(SO <sub>4</sub> ) <sub>2</sub> ·H <sub>2</sub> O		c			-649.				
CsAl(SO <sub>4</sub> ) <sub>2</sub> ·2H <sub>2</sub> O		c			-725.				
CsAl(SO <sub>4</sub> ) <sub>2</sub> ·3H <sub>2</sub> O		c			-798.				
CsAl(SO <sub>4</sub> ) <sub>2</sub> ·12H <sub>2</sub> O		c			-1449.5	-1218.5	893.14	164.	
								148.1	
2CsCl·ThCl <sub>4</sub>		c			-5022.				
2CsCl·ThCl <sub>4</sub> ·8H <sub>2</sub> O		c			-1084.5				
4CsCl·ThCl <sub>4</sub>		c			-718.2				
CsLiICl	fresh melt	c			-178.3				
CsNaICl	fresh melt	c			-178.4				
CsKICl	fresh melt	c			-184.0				



### III. TABLES OF SELECTED VALUES OF PROPERTIES OF SERIES II:

Heat, Temperature, and Entropy of  
Transition, Fusion, and Vaporization  
(including values of pressure)

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 1-1. Oxygen (at. no., 8; at. wt., 16.0000)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
O <sub>2</sub>	Transition	c, III		c, II		23.66	-249.50	0.022	0.93	0.3
	Transition	c, II		c, I		43.77	-229.39	0.178	4.07	-0.14
	Fusion	c, I		liq	1.1	54.40	-218.76	0.106	1.95	1.74
	Vaporization	liq		g	1.1	54.40	-218.76			
	Vaporization	liq		g	760	90.19	-182.97	1.630	18.07	-6.00
O <sub>3</sub>	Vaporization	liq		g	760	162.65	-110.51	2.59	15.92	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 2-1. Hydrogen (at. no., 1; at. wt., 1.0080)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
H <sub>2</sub>	Fusion Vaporization Vaporization		c	liq	54.0	13.96	-259.20	0.028	2.0	1.9
			liq	g	54.0	13.96	-259.20			
			liq	g	760	20.39	-252.77	0.216	10.6	
H <sub>2</sub> O	Fusion Vaporization Vaporization Vaporization Vaporization		c	liq	760	273.16	0.00	1.4363	5.2581	8.911
			liq	g	4.58	273.16	0.00	10.767	39.416	-10.184
			liq	g	23.75	298.16	25.00	10.514	35.263	-9.971
			liq	g	760	373.16	100.00	9.7171	26.0400	-10.021
			liq, std.	g, std.	760	298.16	25.00	10.520	28.390	
			c	liq		271.2	-2.0	2.52	9.29	
H <sub>2</sub> O <sub>2</sub>	Vaporization		liq	g	2.1	298.16	25.00	13.01	43.64	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 2a-1. Hydrogen isotope of mass 1 (at. no., 1; at. wt., 1.0078)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$^1H_2$		Fusion Vaporization Vaporization	c	liq	54.0	13.96	-259.20	0.028	2.0	1.9
			liq	g	54.0	13.96	-259.20			
			liq	g	760	20.39	-252.77	0.216	10.6	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

SERIES II

Washington, D.C.

Table 2b-1. Hydrogen isotope of mass 2 (at. no., 1; at. wt., 2.0142)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$^2\text{H}_2$	Fusion		c	liq	128	18.72	-254.44	0.047	2.5	2.2
	Vaporization		liq	g	128	18.72	-254.44			
	Vaporization		liq	g	760	23.57	-249.59			
$^3\text{H}_2$	Fusion		c	liq	93	16.60	-256.56	0.038	2.3	2.0
	Vaporization		liq	g	93	16.60	-256.56	0.265	16.	
	Vaporization		liq	g	760	22.13	-251.03			
$^2\text{H}_2\text{O}$	Fusion		c	liq		276.98	3.82	1.501	5.419	9.48
	Vaporization		liq	g	20.78	298.16	25.00	10.193	34.185	
	Vaporization		liq, std.	g, std.	760	298.16	25.00	10.350	29.22	
$^3\text{H}_2\text{O}$	Vaporization		liq	g	22.0	298.16	25.00	10.652	35.726	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 3-1. Helium (at. no., 2; at. wt., 4.003)

HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION

June 30, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
He		Fusion	c	liq. I	78200	3.5	-269.7	0.005	1.5	-19
		Transition	liq. II	liq. I	38.3	2.186	-270.974	0.00	0.00	
		Vaporization	liq. I	g	760	4.216	-268.944	0.020	4.7	



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}\text{K}$	$^{\circ}\text{C}$			
Ne		Fusion Vaporization	c	liq	324	24.57	-248.59	0.080	3.26	2.42
			liq	g	324	24.57	-248.59	0.431	17.54	-3.42

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 5-1. Argon (at. no., 18; at. wt., 39.944) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION June 30, 1947; March 31, 1950										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
A		Fusion	c	liq	516.5	83.85	-189.31	0.281	3.35	2.11
		Vaporization	liq	g	516.5	83.85	-189.31			
		Vaporization	liq	g	760	87.29	-185.87	1.558	17.85	-5.08

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 6-1. Krypton [at. no., 36; at. wt., 83.7]  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Kr		Fusion	c	liq	549	115.95	-157.21	0.391	3.37	2.01
		Vaporization	liq	g	549	115.95	-157.21			
		Vaporization	liq	g	760	119.93	-153.23	2.158	17.99	-5.67

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 7-1. Xenon (at. no., 54, at. wt., 131.3)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Xe		Fusion	c	liq	611	161.3	-111.9	0.549	3.40	2.13
		Vaporization	liq	g	611	161.3	-111.9			
		Vaporization	liq	g	760	165.1	-108.1	3.021	18.29	-5.71

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 8-1. Radon (at. no., 86; at. wt., 222)

HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Rn		Fusion Vaporization	c liq	liq g	760	202. 211.	-71. -62.	0.693 3.92	3.43 18.6	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 9-1. Fluorine (at. no., 9; at. wt., 19.00) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION June 30, 1947; March 31, 1950										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$F_2$		Fusion	c	liq		55.20	-217.96	0.372	6.74	1.86
		Vaporization	liq	g	760	85.24	-187.92	1.51	17.7	-4.27
$F_2O$		Vaporization	liq	g	760	128.3	-144.9	2.65	20.7	
HF		Fusion	c	liq		190.09	-65.7	1.094	5.756	2.55
		Vaporization	liq	g	760	293.1	13.9	1.6	6.1	-10.9

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Substance		Type of Process		State		Pressure	Temperature		$\Delta H$	$\Delta S$	$\Delta C_p$
Formula	Description	Process		Initial	Final	mm Hg	$^{\circ}K$	$^{\circ}C$	kcal/mole	cal/deg mole	cal/deg mole
Cl <sub>2</sub>		Fusion		c	liq		172.16	-101.00	1.531	8.89	2.75
		Vaporization		liq	g		172.16	-101.00			
		Vaporization		liq	g	760	239.10	-34.06	4.878	20.40	-8.76
ClO <sub>2</sub>		Vaporization		liq	g	760	284.	11.	7.13	25.1	
Cl <sub>2</sub> O		Vaporization		liq	g	760	275.1	1.9	6.11	22.2	
Cl <sub>2</sub> O <sub>7</sub>		Vaporization		liq	g	85.31	298.16	25.00	7.96	26.7	
		Vaporization		liq	g	760	354.7	81.5	7.88	22.2	
HCl		Transition		c, II	c, I		98.38	-174.78	0.284	2.89	1.15
		Fusion		c, I	liq		158.94	-114.22	0.476	2.99	2.10
		Vaporization		liq	g	760	188.11	-85.05	3.86	20.5	-7.14
HCl·2H <sub>2</sub> O		Fusion		c	liq		258.	-18.	2.5	9.7	
ClF		Vaporization		liq	g	760	172.9	-100.3	5.34	30.88	
ClF <sub>3</sub>		Vaporization		liq	g	760	284.6	11.4	5.74	20.2	
ClO <sub>2</sub> F		Vaporization		liq	g	760	267.4	-5.8	5.60	20.9	

SERIES II

National Bureau of Standards

Washington, D.C.

Table 11-1. Bromine (at. no., 35; at. wt., 79.91)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Br <sub>2</sub>		Fusion Vaporization	c	liq	214	265.9	-7.3	2.52	9.48	2.0
			liq	g		298.16	25.00	7.34	24.6	
HBr		Fusion Vaporization Vaporization	c	liq	760	186.28	-86.82	0.575	3.09	1.64
			liq	g		186.28	-86.82			
			liq	g		206.43	-66.73	4.210	20.39	-1.37
HBr·H <sub>2</sub> O		Fusion	c	liq		552.	-11.	3.0	11.	
BrF <sub>5</sub>		Fusion Vaporization Vaporization	c	liq	2.3	210.7	-62.5	1.74	8.25	
			liq	g		210.7	-62.5	7.43	35.2	
			liq	g		315.5	40.3	7.20	23.0	



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 12-1. Iodine (at. no., 53; at. wt., 126.92)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
I <sub>2</sub>		Fusion Sublimation	c	liq	0.31	386.8	113.6	3.74	9.67	
			c	g		238.16	25.00	14.88	49.91	
HI		Fusion Vaporization	c	liq	760	222.36	-50.80	0.686	3.94	1.17
			liq	g		222.36	-50.80			
IF <sub>5</sub>		Vaporization	liq	g	760	237.80	-35.36	4.724	19.86	-7.14
IF <sub>7</sub>		Fusion Vaporization	c	liq	10.1	282.	9.	3.83	13.5	
			liq	g		282.	9.	10.12	35.7	
ICl		Sublimation	c	g	760	276.6	3.4	7.37	7.6	
ICl		Fusion Vaporization	c	liq	33.2	300.5	27.3	1.83	17.4	
			liq	g		300.5	27.3	9.93	33.06	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 14-1. Sulfur [at. no., 16; at. wt., 32.066]

HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1947; March 31, 1950

Formula	Substance	Type of Process	State		Pressure		Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final	mm Hg		$^{\circ}K$	$^{\circ}C$			
S		Transition	c, II, rhombic	c, I, monoclinic	0.0047		368.6	95.4	0.09	0.25	0.24
		Sublimation	c, II, rhombic	g, S <sub>8</sub>	0.0047		368.6	95.4	3.01	8.17	
		Sublimation	c, I, monoclinic	g, S <sub>8</sub>	0.0047		368.6	95.4	2.92	7.93	
		Fusion	c, I, monoclinic	liq, $\lambda$			392.	119.	0.293	0.75	1.6
		Vaporization	equilibrium liq, $\lambda$ , $\mu$	g	760		717.76	444.60	2.5	3.5	
SO <sub>2</sub>		Fusion	c	liq	1.256		197.68	-75.48	1.769	8.95	4.50
		Vaporization	liq	g	1.256		197.68	-75.48			
		Vaporization	liq	g	760		263.14	-10.02	5.955	22.63	-11.84
SO <sub>3</sub>		Fusion	c, III, ice-like	liq	160		290.0	16.8	0.47	1.6	
		Sublimation	c, III, ice-like	g	160		290.0	16.8	11.76	40.69	
		Fusion	c, II, wool-like	liq	400		305.7	32.5	2.47	8.08	
		Sublimation	c, II, wool-like	g	400		305.7	32.5	12.96	42.40	
		Fusion	c, I, wool-like	liq	1760		335.4	62.2	6.09	18.16	
		Sublimation	c, I, wool-like	g	1760		335.4	62.2			
		Sublimation	c, I, wool-like	g	760		324.8	51.6	15.91	48.98	
		Vaporization	liq	g	760		316.5	43.3	9.99	31.55	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 14-2. Sulfur (at. no., 16; at. wt., 32.066) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION September 30, 1947; March 31, 1950										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$S_2O_7$		Fusion	c	liq		273.	0.			
$H_2S$		Transition	c, III	c, II		103.54	-169.62	0.366	3.53	1.20
		Transition	c, II	c, I		126.24	-146.92	0.108	0.86	-0.65
		Fusion	c, I	liq		173.9	-85.53	0.568	3.03	1.59
		Vaporization	liq	g		173.9	-85.53			
		Vaporization	liq	g		760	-60.34	4.463	20.97	-8.34
$H_2S_2$		Fusion	c	liq		183.4	-89.8	1.8	9.7	12.9
		Vaporization	liq	g		748	70.3	8.4	24.4	
$H_2S_3$		Fusion	c	liq		220.	-53.			
		Vaporization	liq	g		760	69.			
$H_2SO_4$		Fusion	c	liq		283.51	10.35	2.36	8.32	6.218
$H_2SO_4 \cdot H_2O$		Fusion	c	liq		281.69	8.53	4.63	16.44	25.04
$H_2SO_4 \cdot 2H_2O$		Fusion	c	liq		234.3	-38.9			
$H_2SO_4 \cdot 4H_2O$		Fusion	c	liq		248.	-25.			
$H_2SO_5$		Fusion	c	liq		318.	45.			
$H_2S_2O_7$		Fusion	c	liq		308.1	34.9	3.19	10.4	30.1

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 14-3. Sulfur (at. no., 16; at. wt., 32.066)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1947; March 31, 1950

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
			Initial	Final		$^{\circ}K$	$^{\circ}C$			
$H_2S_2O_8$		Fusion	c	liq		438.	65.			
$SF_4$		Fusion	c	liq	1.3	149.	-124.			
		Vaporization	liq	g	1.3	149.	-124.			
		Vaporization	liq	g	760	233.	-40.	5.20	22.2	
		Vaporization	liq	g						
$SF_6$		Transition	c, II	c, I		94.26	-178.90	0.384	4.07	-1.98
		Sublimation	c, I	g	760	209.5	-63.7	5.46	26.04	
		Fusion	c, I	liq	1700	222.5	-50.7	1.20	5.40	-1.7
		Vaporization	liq	g	1700	222.5	-50.7	4.08	18.34	
		Vaporization	liq	g						
$S_2F_{10}$		Fusion	c	liq		167.7	-105.5			
		Vaporization	liq	g	760	174.	-99.			
		Vaporization	liq	g						
$SO_2F_2$		Fusion	c	liq	760	181.	-92.		20.33	
		Vaporization	liq	g	760	302.3	29.1	6.15		
		Vaporization	liq	g						
$SO_2F_2$		Fusion	c	liq	0.5	143.7	-129.5			
		Vaporization	liq	g	760	229.31	-43.85	5.08	22.14	
		Vaporization	liq	g						
$SO_2F_2$		Fusion	c	liq	65	153.	-120.			
		Vaporization	liq	g	760	221.	-52.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 14-4. Sulfur (at. no., 16; at. wt., 32.066)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1947; March 31, 1950

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		$^{\circ}K$	$^{\circ}C$			
$HSO_3F$		Vaporization	liq	g	760	435.8	162.6			
$S_2Cl_2$		Fusion	c	liq		193.	-80.			
		Vaporization	liq	g	760	411.2	138.0	8.61	20.92	
$SOCl_2$		Fusion	c	liq		168.7	-104.5			
		Vaporization	liq	g	760	348.9	75.7	7.41	21.24	
$SO_2Cl_2$		Fusion	c	liq		227.	-46.			
		Vaporization	liq	g	760	342.4	69.2	7.50	21.90	
$S_2O_5Cl_2$		Fusion	c	liq		234.	-39.			
		Vaporization	liq	g	730.5	423.9	150.7	13.2	31.1	
$SOFCI$		Fusion	c	liq		135.7	-137.5			
		Vaporization	liq	g	760	285.4	12.2	5.67	19.9	
$SO_2FCI$		Fusion	c	liq		148.5	-124.7			
		Vaporization	liq	g	760	280.4	7.2	5.69	20.3	
$S_2Br_2$		Fusion	c	liq		227.	-46.			
		Vaporization	liq	g	0.18	327.	54.			
$SOBr_2$		Fusion	c	liq		221.	-52.			
		Vaporization	liq	g	760	411.	138.	10.2	24.8	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 14a-1. Sulfur (at. no., 16) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION September 30, 1947; March 31, 1950										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$^{32}_{16}S$	Transition		c, III	c, II		107.82	-165.34	0.402	3.73	0.78
	Transition		c, II	c, I		132.85	-140.31	0.124	0.93	-0.75
	Fusion		c, I	liq	163.0	187.14	-86.02	0.566	3.02	3.50
$^{34}_{16}S$	Transition		c, III	c, II		105.8	-167.4	0.386	3.64	1.00
	Transition		c, II	c, I		129.6	-143.6	0.116	0.90	-0.70
	Fusion		c, I	liq	168	187.40	-85.76	0.567	3.02	2.56

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Se		Transition	vitreous	c, I, gray	0.0043 0.0043 760 760 760 760	398.	125.	1.05	2.64	2.2
		Transition	c, III, red	c, I, gray		423.	150.	0.18	0.43	0.2
		Fusion	c, I, gray	liq		490.6	217.4	1.25	2.56	2.7
		Vaporization	liq	g, Se <sub>6</sub>		490.6	217.4	4.31	8.79	
		Vaporization	liq	g, Se <sub>6</sub>		1009.	736.	3.43	3.40	
		Vaporization	liq	g, Se <sub>4.37</sub>		958.0	684.8	5.10	5.32	
SeO <sub>2</sub>		Vaporization	liq	g, Se <sub>2</sub>	760	1027.	754.	12.80	12.46	
		Sublimation	c	g		595.	322.	21.1	35.4	
H <sub>2</sub> Se		Transition	c, III	c, II	205.4 205.4 760	82.3	-190.9	0.309	3.75	1.5
		Transition	c, II	c, I		172.54	-100.62	0.267	1.55	-1.9
		Fusion	c, I	liq		207.43	-65.73	0.601	2.90	1.95
		Vaporization	liq	g		207.43	-65.73	5.34	25.72	
		Vaporization	liq	g		231.9	-41.3	4.62	19.93	
H <sub>2</sub> SeO <sub>4</sub>		Fusion	c	liq	760	333.	60.	3.45	10.4	
H <sub>2</sub> SeO <sub>4</sub> ·H <sub>2</sub> O		Fusion	c	liq		299.	26.	4.75	15.9	
H <sub>2</sub> SeO <sub>4</sub> ·4H <sub>2</sub> O		Fusion	c	liq		221.5	-51.7			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 15-- Selenium (at. no., 34; at. wt., 78.96)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1947; March 31, 1950

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		°K	°C			
SeF <sub>4</sub>			Fusion Vaporization	c liq	liq g	760	260.0 366.	-13.2 93.			
SeF <sub>6</sub>			Sublimation Fusion Vaporization	c c liq	g liq g	760 1500 1500	236.6 336.6 236.6	-46.6 -34.6 -34.6	6.27 1.70 4.38	27.68 7.14 18.34	
SeOF <sub>2</sub>			Fusion Vaporization	c liq	liq g	760	277.8 397.	4.6 124.			
SeCl <sub>4</sub>			Sublimation Fusion Vaporization	c c liq	g liq g	760	469. 578. 578.	196. 305. 305.			
SeOCl <sub>2</sub>			Fusion Vaporization	c liq	liq g	760	284.0 450.8	10.8 177.6	1.01 10.2	3.56 22.6	
Se(OH) <sub>3</sub> ClO <sub>4</sub>			Fusion	c	liq		306.	33.			
SeOBr <sub>2</sub>			Fusion Vaporization	c liq	liq g	740	314.8 490.	41.6 217.			



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 15a-1. Selenium (at. no., 34) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION September 30, 1947; March 31, 1950										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$24_2\text{Se}$		Transition	c, III	c, II		90.5	-182.7	0.378	4.19	1.1
		Transition	c, II	c, I		176.02	-97.14	0.283	1.61	-0.5
		Fusion	c, I	liq	193.4	206.24	-66.92	0.596	2.89	1.9
		Vaporization	liq	g	193.4	206.24	-66.92			
$144_4\text{Se}$		Transition	c, III	c, II		87.0	-186.2	0.349	4.01	1.3
		Transition	c, II	c, I		174.35	-98.81	0.275	1.58	-1.2
		Fusion	c, I	liq	199.4	206.80	-66.36	0.598	2.89	1.9
		Vaporization	liq	g	199.4	206.80	-66.36			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 16-1. Tellurium (at. no., 52; at. wt., 127.61)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1947; March 31, 1950

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		$^{\circ}K$	$^{\circ}C$			
Te		Transition	c, II	c, I	0.18	621.	348.	0.13	0.21	2.0
		Fusion	c, I	liq		723.	450.	4.28	5.92	
		Vaporization	liq	g, Te <sub>2</sub>		723.	450.	13.3	18.4	
		Vaporization	liq	g, Te <sub>2</sub>		1360.	1087.	11.9	8.75	
H <sub>2</sub> Te		Fusion	c	liq	760	222.	-51.	1.0	4.5	
		Vaporization	liq	g		270.9	-2.3	5.55	20.49	
TeF <sub>6</sub>		Transition	c, II	c, I	60	199.7	-73.5	0.5	2.5	
		Sublimation	c, I	g		234.6	-38.6	6.47	27.6	
		Fusion	c, I	liq		800	-37.7	2.1	9.0	
		Vaporization	liq	g		800	-37.7	4.3	18.4	
TeCl <sub>4</sub>		Fusion	c	liq	760	497.3	224.1	4.51	9.07	21.9
		Vaporization	liq	g		661.	388.	18.4	27.9	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 18-1. Nitrogen (at. no., 7; at. wt., 14.008)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$t_K$	$t_C$			
N <sub>2</sub>	Transition	c, II		c, I		35.62	-237.54	0.055	1.54	
	Fusion	c, I		liq	94	63.15	-210.01	0.172	2.72	
	Vaporization	liq		g	94	63.15	-210.01			
	Vaporization	liq		g	760	77.34	-195.82	1.333	17.24	
NO	Fusion	c		liq	164.4	109.51	-163.65	0.550	5.02	6.0
	Vaporization	liq		g	164.4	109.51	-163.65			
	Vaporization	liq		g	760	121.39	-151.77	3.293	27.13	11.8
N <sub>2</sub> O	Fusion	c		liq	658.9	182.30	-90.86	1.563	8.574	4.67
	Vaporization	liq		g	658.9	182.30	-90.86			
	Vaporization	liq		g	760	184.68	-88.48	3.956	21.42	
N <sub>2</sub> O <sub>3</sub>	Fusion	c		liq		162.	-111.			
	Vaporization	liq		g	760	275.	2.	9.4	34.	
N <sub>2</sub> O <sub>4</sub>	Fusion	c		liq	139.78	261.96	-11.20	3.502	13.37	6.12
	Vaporization	liq		g	139.78	261.96	-11.20			
	Vaporization	liq		g	760	294.31	21.15	9.110	30.95	
N <sub>2</sub> O <sub>5</sub>	Sublimation	c		g	760	305.6	32.4	13.6	44.5	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 18-2. Nitrogen (at. no., 7; at. wt., 14.008)

HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
NH <sub>3</sub>		Fusion	c	liq	45.57	195.40	-77.76	1.351	6.914	
		Vaporization	liq	g	45.57	195.40	-77.76			
		Vaporization	liq	g	760	239.73	-33.43	5.551	23.28	
NH <sub>3</sub> · $\frac{1}{2}$ H <sub>2</sub> O		Fusion	c	liq		194.3	-78.9			
NH <sub>3</sub> ·H <sub>2</sub> O		Fusion	c	liq		194.1	-79.1			
N <sub>2</sub> H <sub>4</sub>		Fusion	c	liq		274.7	1.5			
		Vaporization	liq	g	764	386.7	113.5	10.	25.9	
N <sub>2</sub> H <sub>4</sub> ·H <sub>2</sub> O		Fusion	c	liq		233.	-40.			
		Vaporization	liq	g	739.5	391.7	118.5			
HN <sub>3</sub>		Fusion	c	liq		193.	-80.			
		Vaporization	liq	g	760	309.	36.	7.1	23.	
NH <sub>4</sub> N <sub>3</sub>		Sublimation	c	g	760	407.	134.	15.1	37.1	
		Fusion	c	liq		433.	160.			
N <sub>2</sub> H <sub>4</sub> ·HN <sub>3</sub>		Fusion	c	liq		338.	65.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

SERIES II

Washington, D.C.

Table 18-3. Nitrogen (at. no., 7; at. wt., 14.006)										
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION										
December 31, 1947; March 31, 1950										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
HNO <sub>3</sub>		Fusion Vaporization	c	liq	48	231.56	-41.60	2.503	10.81	10.55
			liq	g		293.	20.	9.43	32.2	
HNO <sub>3</sub> ·H <sub>2</sub> O		Fusion	c	liq		235.53	-37.63	4.184	17.76	21.51
HNO <sub>3</sub> ·3H <sub>2</sub> O		Fusion	c	liq		254.69	-18.47	6.954	27.3	34.2
NH <sub>2</sub> OH		Fusion Vaporization	c	liq	22	306.3	33.1			
			liq	g		331.	58.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 18-4. Nitrogen (at. no., 7; at. wt., 14.008)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
NH <sub>4</sub> NO <sub>3</sub>	Transition	c, V	c, IV	c, IV	760	255.	-18.	0.13	0.51	
	Transition	c, IV	c, III	c, III	760	305.3	32.1	0.38	1.23	
	Transition	c, III	c, II	c, II	6.32×10 <sup>5</sup>	336.5	63.3	0.20	0.59	
	Transition	c, IV	c, III	c, III	6.32×10 <sup>5</sup>	336.5	63.3	0.32	0.95	
	Transition	c, IV	c, II	c, II	6.32×10 <sup>5</sup>	336.5	63.3	0.32	1.54	
	Transition	c, III	c, II	c, II	760	357.4	84.2	0.32	0.90	
	Transition	c, II	c, I	c, I	760	398.4	125.2	1.01	2.54	
	Transition	c, IV	c, II	c, II	6.73×10 <sup>6</sup>	442.4	169.2	0.96	2.16	
	Transition	c, IV	c, VI	c, VI	6.73×10 <sup>6</sup>	442.4	169.2	0.98	2.21	
	Transition	c, II	c, VI	c, VI	6.73×10 <sup>6</sup>	442.4	169.2	0.02	0.05	
	Fusion	c, I	liq	liq	760	442.8	169.6	1.3	2.94	
	Transition	c, VI	c, I	c, I	6.63×10 <sup>6</sup>	459.9	186.7	0.99	2.15	
	Transition	c, II	c, VII	c, VII	6.63×10 <sup>6</sup>	459.9	186.7	0.03	0.06	
	Transition	c, II	c, I	c, I	6.63×10 <sup>6</sup>	459.9	186.7	1.02	2.21	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 18-5. Nitrogen (at. no., 7; at. wt., 14.008)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$T_K$	$T_C$			
$N_2H_4 \cdot HNO_2$	metastable stable		Fusion	c	liq		316.	43.			
$N_2H_4 \cdot HNO_3$			Fusion	c, II	liq		335.3	62.1			
			Fusion	c, I	liq		343.9	70.7			
$N_2H_4 \cdot 2HNO_3$			Fusion	c	liq		377.	104.			
$NH_4NO_3 \cdot 2HNO_3$			Fusion	c	liq		302.7	29.5			
$NF_3$			Transition	c, II	c, I		54.	-219.			
			Fusion	c, I	liq		64.7	-208.5			
			Vaporization	liq	g	760	144.1	-129.1	2.93	20.3	
$NOF$			Fusion	c	liq		140.7	-132.5			
			Vaporization	liq	g	760	213.3	-59.9	4.61	21.6	
$NO_2F$			Fusion	c	liq		107.2	-166.0			
			Vaporization	liq	g	760	200.8	-72.4	4.31	21.5	
$NO_3F$			Fusion	c	liq		92.	-181.			
			Vaporization	liq	g	103	193.	-80.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 18-6. Nitrogen (at. no., 7; at. wt., 14.008)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$NH_2F$		Sublimation	c	g	760	196.	-77.			
$NHF_2$		Fusion	c	liq		148.	-125.			
		Vaporization	liq	g	760	208.	-65.			
$NH_4F \cdot HF$		Fusion	c	liq		397.8	126.6			
$N_2H_4 \cdot 2HF$		Fusion	c	liq		378.	105.			
$NCI_3$		Vaporization	liq	g	760	344.	71.			
$NOCl$		Fusion	c	liq		211.7	-61.5			
		Vaporization	liq	g	38.6	211.7	-61.5			
		Vaporization	liq	g	760	267.4	-5.8	6.0	22.4	
$NO_2Cl$		Fusion	c	liq		128.	-145.			
		Vaporization	liq	g	760	257.9	-15.3	5.0	19.4	
$NH_2Cl$		Fusion	c	liq		207.	-66.			
$NH_4Cl$		Transition	c, II	c, I	760	457.6	184.4	1.0	2.2	
		Fusion	c, I	liq	$2.62 \times 10^4$	793.	520.			
		Vaporization	liq	g	$2.62 \times 10^4$	793.	520.			
$NH_4Cl \cdot 3NH_3$		Fusion	c	liq		283.9	10.7			



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 18-7. Nitrogen (at. no., 7; at. wt., 14.008)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$N_2H_4 \cdot HCl$		Fusion	c	liq		365.8	92.6	3.64	9.95	
$N_2H_4 \cdot 2HCl$		Fusion	c	liq		471.	198.			
$NH_2OH \cdot HCl$		Fusion	c	liq		430.	157.			
$NH_4ClO_4$		Transition	c, II	c, I		513.	240.			
$N_3Br$		Fusion	c	liq		228.	-45.			
$NOBr$		Fusion	c	liq		217.7	-55.5			
$NOBr_3$		Fusion	c	liq		233.	-40.			
$NH_4Br$		Transition Fusion	c, II c, I	c, I liq	735	411.0 815.	137.8 542.	0.77	1.87	
$NH_4Br \cdot 3NH_3$		Fusion	c	liq		286.9	13.7			
$N_2H_4 \cdot HBr$		Fusion	c	liq		353.	80.			
$N_2H_4 \cdot 2HBr$		Fusion	c	liq		468.	195.			
$NH_4I$		Transition Fusion	c, II c, I	c, I liq	760	260. 824.	-13. 551.	0.70	2.7	
$NH_4I \cdot 3NH_3$		Fusion	c	liq		265.2	-8.0			
$NH_4I \cdot 4NH_3$		Fusion	c	liq		268.1	-5.1			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$N_2H_4 \cdot HI$		Fusion	c	liq		400.	127.			
$N_2H_4 \cdot 2HI$		Fusion	c	liq		493.	220.			
$NH_4Cl_2I$		Fusion	c	liq		435.	162.			
$NH_4Cl_4I$		Fusion	c	liq		398.	125.			
$NH_4IBr_2$		Fusion	c	liq		471.	198.			
$N_2S_5$		Fusion	c	liq		283.	10.			
$N_4S_4$		Fusion	c	liq		453.	180.			
$N_2O_3 \cdot 2SO_3$		Fusion Vaporization	c liq	liq g	760	490. 630.	217. 357.			
$NO_2SO_2OH$		Fusion	c	liq		346.7	73.5			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 18-9. Nitrogen (at. no., 7; at. wt., 14.008)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$\text{NH}_4\text{HSO}_4$		Transition	c, I	c, II	$1.368 \times 10^6$	399.4	126.2	3.42	8.20	
		Transition	c, II	c, III		399.4	126.2			
		Fusion	c, I	liq	$4.165 \times 10^6$	417.	144.			
		Transition	c, II	c, III		450.1	176.9			
$\text{N}_2\text{H}_4 \cdot \text{H}_2\text{SO}_4$		Transition	c, III	c, IV	$4.165 \times 10^6$	450.1	176.9			
		Fusion	c	liq		527.	254.			
		Fusion	c	liq		786.	513.			
		Fusion	c	liq		443.	170.			
$(\text{NH}_4)_2\text{SO}_4$		Fusion	c	liq		441.	168.			
		Fusion	c	liq		390.	117.			
		Fusion	c	liq		413.	140.			
		Fusion	c	liq						
$(\text{NH}_2\text{OH}) \cdot \text{H}_2\text{SO}_4$										
$(\text{H}_2\text{NSO}_2)_2\text{NH}$										
$(\text{N}_2\text{H}_4)_2 \cdot \text{H}_2\text{SO}_4$										
$\text{NOSO}_3\text{F}$										

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
P		Transition	c, IV, white	c, III, white	760	196.	-77.			
		Transition	c, IV, white	c, III, white	4.41x10 <sup>6</sup>	270.8	-2.4	1.35	4.09	
		Transition	c, III, white	c, II, red		298.16	25.00	-4.4		
		Transition	c, III, white	c, I, black		298.16	25.00	-10.3		
		Fusion	c, III, white	liq	760	317.4	44.2	0.15	0.47	
		Vaporization	liq	g, P <sub>4</sub>	760	553.	280.	2.97	5.37	
		Fusion	c, II, red	liq	32760	863.	590.	4.85	5.62	
		Vaporization	liq	g, P <sub>4</sub>	32760	863.	590.	2.50	2.90	
P <sub>4</sub> O <sub>6</sub>		Fusion	c	liq		297.0	23.8			
		Vaporization	liq	g	760	448.5	175.3	9.0	20.1	
P <sub>4</sub> O <sub>10</sub>		Sublimation	c, II	g	760	632.	359.	17.6	27.8	
		Fusion	c, II	liq	3700	693.	420.	5.	7.2	
		Fusion	c, I	liq	570	845.	572.	11.5	13.6	
		Vaporization	liq	g	570	845.	572.	16.8	19.9	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
PH <sub>3</sub>		Transition	c, IV	c, III		30.31	-242.85	0.0197	0.650	
		Transition	c, III	c, II		49.44	-223.72	0.186	3.76	
		Transition	c, II	c, I		88.12	-185.04	0.116	1.32	
		Fusion	c, I	liq	27.33	139.38	-133.78	0.270	1.94	
		Vaporization	liq	g	27.33	139.38	-133.78			
		Vaporization	liq	g	760	185.42	-87.74	3.490	18.82	
P <sub>2</sub> H <sub>4</sub>		Vaporization	liq	g	735	330.	57.			
H <sub>3</sub> PO <sub>2</sub>		Fusion	c	liq		299.7	26.5	2.32	7.74	
H <sub>3</sub> PO <sub>3</sub>		Fusion	c	liq		343.3	70.1	3.07	8.94	
H <sub>3</sub> PO <sub>4</sub>		Fusion	c	liq		315.51	42.35	2.52	7.99	
H <sub>3</sub> PO <sub>4</sub> · $\frac{1}{3}$ H <sub>2</sub> O		Fusion	c	liq		302.51	29.35	3.64	12.03	
H <sub>4</sub> P <sub>2</sub> O <sub>6</sub>		Fusion	c	liq		328.	55.			
H <sub>4</sub> P <sub>2</sub> O <sub>7</sub> · 2H <sub>2</sub> O		Fusion	c	liq		335.	62.	0.84	2.5	
H <sub>4</sub> P <sub>2</sub> O <sub>7</sub>		Fusion	c	liq		334.	61.	2.20	6.6	
H <sub>4</sub> P <sub>2</sub> O <sub>7</sub> · $\frac{1}{3}$ H <sub>2</sub> O		Fusion	c	liq		299.	26.	3.14	10.5	
H <sub>6</sub> P <sub>4</sub> O <sub>13</sub>		Fusion	c	liq		307.	34.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 19-3. Phosphorus (at. no., 15; at. wt., 30.98)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
PF <sub>3</sub>		Fusion	c	liq	760	121.7	-151.5	3.43	19.9	
		Vaporization	liq	g		172.0	-101.2			
PF <sub>5</sub>		Fusion	c	liq	427	179.4	-93.8	2.8	15.6	
		Sublimation	c	g	427	179.4	-93.8	6.9	38.5	
		Vaporization	liq	g	760	188.7	-84.5	4.1	21.7	
POF <sub>3</sub>		Sublimation	c	g	760	233.7	-39.5	9.0	38.5	
		Fusion	c	liq	778	234.1	-39.1	3.7	15.8	
		Vaporization	liq	g	760	233.5	-39.7	5.3	22.7	
		Vaporization	liq	g	760	383.	110.			
PCl <sub>3</sub>		Fusion	c	liq	760	181.	-92.	7.28	20.9	
		Vaporization	liq	g		349.	76.			
PCl <sub>5</sub>		Sublimation	c	g, mixt. equil.	760	432.	159.	16.1	37.3	
		Fusion	c	liq		433.	160.			
P <sub>2</sub> Cl <sub>4</sub>		Fusion	c	liq		245.	-28.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 19-4. Phosphorus (at. no., 15; at. wt., 30.98)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$POCl_3$		Fusion Vaporization	c liq	liq g	760	274.33 378.5	1.17 105.3	3.11 8.21	11.33 21.69	
$PH_4Cl$		Transition Fusion	c, II c, I	c, I liq	34790	232. 301.7	-41. 28.5			
$PF_2Cl$		Fusion Vaporization	c liq	liq g	760	108.4 226.0	-164.8 -47.3	4.21	18.6	
$PFCl_2$		Fusion Vaporization	c liq	liq g	760	129.2 286.9	-144.0 13.7	5.93	20.7	
$POF_2Cl$		Fusion Vaporization	c liq	liq g	760	176.8 276.1	-96.4 2.9	5.93	21.5	
$POFCl_2$		Fusion Vaporization	c liq	liq g	760	193.1 326.0	-80.1 52.8	6.98	21.4	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 19-5. Phosphorus (at. no., 15; at. wt. 30.98)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$PBr_3$		Fusion Vaporization	c liq	liq g	760	232.7 446.4	-40.5 173.2	9.28	20.79	
$PBr_5$		Sublimation	c	g	210	357.0	83.8	13.0	36.4	
$P O B r _3$		Fusion Vaporization	c liq	liq g	760	328. 464.9	55. 131.7	9.08	19.53	
$P F _2 B r$		Fusion Vaporization	c liq	liq g	760	139.4 257.1	-133.8 -16.1	5.45	21.2	
$P F B r _2$		Fusion Vaporization	c liq	liq g	760	158.2 351.5	-115.0 78.3	7.34	20.88	
$P F _3 B r _2$		Fusion	c	liq		253.	-20.			
$P O F _2 B r$		Fusion Vaporization	c liq	liq g	760	188.4 305.1	-84.8 31.9	6.73	22.06	
$P O F B r _2$		Fusion Vaporization	c liq	liq g	760	156.0 383.0	-117.2 109.8	7.38	19.27	



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 19-6. Phosphorus (at. no., 15; at. wt., 30.98)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$POCl_2Br$		Fusion	c	liq		286.	13.			
$POClBr_2$		Fusion	c	liq		303.	30.			
$POClBr$		Vaporization	liq	g	765	352.	79.			
$PI_3$		Fusion	c	liq		334.7	61.5			
$P_2I_4$		Fusion	c	liq		397.7	124.5			
$P_2S_3$		Fusion	c	liq		569.	236.			
$P_4S_3$		Fusion	c	liq		445.	172.			
		Vaporization	liq	g	760	680.	407.			
$P_4S_7$		Fusion	c	liq		580.	307.			
$P_4S_{10}$		Fusion	c	liq		561.	288.			
$P_4O_6S_4$		Fusion	c	liq		375.	102.			
		Vaporization	liq	g	760	568.	295.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 19-7. Phosphorus (at. no., 15; at. wt., 30.98)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
PSF <sub>3</sub>		Fusion Vaporization	c liq	liq g	760	124. 221.0	-149. -52.2	4.68	21.2	
PSCl <sub>3</sub>		Fusion Fusion Vaporization	c, II c, I liq	liq liq g	760	232.4 237.0 398.	-40.8 -36.2 125.			
PSF <sub>2</sub> Cl		Fusion Vaporization	c liq	liq g	760	118. 279.5	-155. 6.3	5.82	20.82	
PSFCl <sub>2</sub>		Fusion Vaporization	c liq	liq g	760	177.2 337.9	-96.0 64.7	7.1	21.0	
PSBr <sub>3</sub>		Fusion	c	liq		311.	38.			
PSF <sub>2</sub> Br		Fusion Vaporization	c liq	liq g	760	136.3 308.7	-136.9 35.5	6.77	21.93	
PSFBr <sub>2</sub>		Fusion Vaporization	c liq	liq g	760	198.0 398.5	-75.2 125.3	8.35	20.95	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 19-8. Phosphorous (at. no., 15; at. wt., 30.98)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$PSCl_2Br$		Fusion	c	liq		243.	-30.			
		Vaporization	liq	g	60	353.	80.			
$PSClBr_2$		Fusion	c	liq		267.	-6.			
		Vaporization	liq	g	60	368.	95.			
$NH_2OH \cdot H_3PO_2$		Fusion	c	liq		365.	92.			
$NH_4H_2PO_3$		Fusion	c	liq		393.	120.			
$N_2H_4 \cdot H_3PO_3$		Fusion	c	liq		309.	36.			
$NH_4NH_2HPO_3$		Fusion	c	liq		578.	305.			
$N_2H_4 \cdot H_3PO_4$		Fusion	c	liq		355.	82.			
$N_2H_4 \cdot H_4P_2O_6$		Fusion	c	liq		425.	152.			
$N_2H_4 \cdot (H_3PO_3)_2$		Fusion	c	liq		365.	82.			
$(NH_4)_2H_2P_2O_6$		Fusion	c	liq		443.	170.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 19-9. Phosphorus (at. no., 15; at. wt., 30.98)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
$N_3P_3F_8H_{12} \cdot 2H_2O$			Fusion	c	liq		305.7	32.5			
$NH_4HPO_3F$			Fusion	c	liq		498.7	225.5			
$NH_4PO_2F_2$			Fusion	c	liq		486.	213.			
$(PNCI_2)_3$			Fusion	c	liq	9	388.1	114.9	5.0	12.9	
			Vaporization	liq	g	9	388.1	114.9	13.2	34.0	
$(PNCI_2)_4$			Vaporization	liq	g	760	526.	253.	13.	25.	
			Fusion	c	liq		396.7	123.5			
$(PNCI_2)_5$			Vaporization	liq	g	760	601.7	328.5	15.	25.	
			Fusion	c	liq		314.	41.			
$(PNCI_2)_6$			Vaporization	liq	g	13	501.	224.			
			Fusion	c	liq		364.	91.			
$N_4P_4Cl_2F_6$			Vaporization	liq	g	13	595.	262.			
			Fusion	c	liq		261.0	-12.2			
$PNBr_2$			Vaporization	liq	g	760	379.0	105.8	8.75	23.09	
			Fusion	c	liq		463.	190.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 20-1. Arsenic (at. no., 33; at. wt., 74.91)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
As		Sublimation	c, metallic	g, As <sub>4</sub>	760	883.	610.			
		Fusion	c, metallic	liq	27200	1090.	817.	7.75	8.78	
		Vaporization	liq	g	27200	1090.	817.			
As <sub>4</sub> O <sub>6</sub>		Sublimation	c, octahedral	g	28	547.	274.	26.1	47.7	
		Fusion	c, octahedral	liq	28	547.	274.	11.9	21.8	
		Sublimation	c, monoclinic	g	67	586.	313.	23.0	39.2	
		Fusion	c, monoclinic	liq	67	586.	313.	8.8	15.0	
		Vaporization	liq	g	760	733.	460.	14.2	19.4	
AsH <sub>3</sub>		Fusion	c	liq	24.6	156.9	-116.3	0.56	3.6	
		Vaporization	liq	g	24.6	156.9	-116.3			
		Vaporization	liq	g	760	210.7	-62.5	4.18	19.8	
H <sub>3</sub> AsO <sub>4</sub> · $\frac{3}{2}$ H <sub>2</sub> O		Fusion	c	liq		309.30	36.14			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards SERIES II φ Washington, D.C.

Table 20-2. Arsenic (at. no., 33; at. wt., 74.91)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
AsF <sub>3</sub>		Fusion Vaporization	c liq	liq g		267.21 292.50	-5.95 19.34	2.486 8.566	9.304 29.285	
AsF <sub>5</sub>		Fusion Vaporization	c liq	liq g	149 149	192.9 192.9	-80.3 -80.3	2.71	14.0	
		Vaporization	liq	g	760	220.6	-52.6	4.96	22.5	
AsCl <sub>3</sub>		Fusion Vaporization	c liq	liq g	760	257. 403.	-16. 130.	2.42 7.5	9.42 18.6	
AsBr <sub>3</sub>		Fusion Vaporization	c liq	liq g	760	304.4 494.	31.2 221.	2.81 10.	9.23 20.	
AsI <sub>2</sub>		Fusion	c	liq		403.	130.			
AsI <sub>3</sub>		Fusion Vaporization	c liq	liq g	1.1 1.1	417. 417.	144. 144.	20.	48.	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Substance		Type of Process		State		Pressure	Temperature		$\Delta H$	$\Delta S$	$\Delta C_p$
Formula	Description	Process		Initial	Final	mm Hg	°K	°C	kcal/mole	cal/deg mole	cal/deg mole
$As_2S_3$		Transition Fusion Vaporization		c, II, red	c, I, black	760	540.	267.			
				c, I, black	liq		580.	307.			
				liq	g		838.	565.			
$As_2S_3$		Transition Fusion		c, II, yellow	c, I, red		443.	170.			
				c, I, red	liq		573.	300.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 21-1. Antimony (at. no., 51; at. wt., 121.76)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
Sb			Transition	c, III	c, II	760	367.8	94.6	4.8	5.3	
			Transition	c, II	c, I		690.	413.			
			Fusion	c, I	liq		903.7	630.5			
			Vaporization	liq	g, equil		1713.	1440.			
Sb <sub>2</sub> O <sub>3</sub>			Transition	c, II	c, I	0.525	830.	557.	3.24	3.91	
			Fusion	c, I	liq		928.	655.			
			Vaporization	liq	g		928.	655.			
			Vaporization	liq	g		1729.	1456.			
SbH <sub>3</sub>			Fusion	c	liq	760	185.	-88.			
			Vaporization	liq	g		256.	-17.			
SbF <sub>3</sub>			Fusion	c	liq	760	563.	290.			
			Vaporization	liq	g		649.	376.			
SbF <sub>5</sub>			Fusion	c	liq	760	279.	6.			
			Vaporization	liq	g		422.	149.			
2SbF <sub>3</sub> ·SbF <sub>5</sub>			Vaporization	liq	g	760	663.	390.			



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 21-2. Antimony [at. no., 51; at. wt., 121.76]  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		°K	°C			
SbCl <sub>3</sub>			Fusion Vaporization	c	liq	760	246.4	73.2	3.03	8.74	
				liq	g		494.	221.	10.80	21.86	
SbCl <sub>5</sub>			Fusion Vaporization	c	liq	21	276.2	3.0	2.4	8.7	
				liq	g		349.7	76.5	11.5	32.9	
SbCl <sub>3</sub> F <sub>2</sub>			Fusion	c	liq		328.	55.			
SbBr <sub>3</sub>			Fusion Vaporization	c	liq	749	369.8	96.6	3.51	9.49	
				liq	g		561.	288.			
SbI <sub>3</sub>			Fusion	c	liq		443.7	170.5			
Sb <sub>4</sub> S <sub>6</sub>			Fusion	c	liq		820.	547.	11.2	13.7	
				c	liq						
Sb <sub>2</sub> Se <sub>3</sub>			Fusion	c	liq		890.	617.			
Sb <sub>2</sub> Te <sub>3</sub>			Fusion	c	liq		898.	625.			
				c	liq						
3(NH <sub>4</sub> Cl)·2SbCl <sub>3</sub>			Fusion	c	liq		564.	291.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 22-1. Bismuth (at. no., 83; at. wt., 209.00)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1947; March 31, 1950

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
Bi			Fusion	c	liq		544.2	271.0			
			Vaporization	liq	g	760	1693.	1420.	2.63	4.83	
Bi <sub>2</sub> O <sub>3</sub>			Fusion	c	liq		1090.	817.	6.8	6.2	
BiF <sub>3</sub>			Fusion	c	liq		1000.	727.			
BiF <sub>5</sub>			Sublimation	c	g	760	823.	550.			
BiCl <sub>3</sub>			Fusion	c	liq		505.	232.	2.6	5.1	
			Vaporization	liq	g	760	714.	441.	17.35	24.30	
BiCl <sub>4</sub>			Fusion	c	liq		498.	225.			
BiBr <sub>3</sub>			Transition	c, II	c, I		428.	155.			
			Fusion	c, I	liq		491.	218.			
			Vaporization	liq	g	760	734.	461.	18.0	24.5	
BiS			Fusion	c	liq		958.	685.			
BiSe			Transition	c, II	c, I		695.	422.			
Bi <sub>2</sub> Se <sub>3</sub>			Fusion	c	liq		979.	706.			
Bi <sub>2</sub> Te <sub>3</sub>			Fusion	c	liq		853.	580.			
Bi <sub>2</sub> Te <sub>3</sub> ·Bi <sub>2</sub> S <sub>3</sub>			Fusion	c	liq		888.	615.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 23-1. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1948; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
C		Sublimation Sublimation	c, graphite c, graphite	g, std. state equilibrium gas	760 760	298.16 4620.	25.00 4347.	171.698	36.4002	
CO	carbon monoxide	Transition Fusion Vaporization Vaporization	c, II c, I liq liq	c, I liq g g		61.53 68.10 68.10 81.66	-211.63 -205.06 -205.06 -191.50	0.151 0.200 1.444	2.45 2.94 17.68	-3.2 1.9
CO <sub>2</sub>	carbon dioxide	Sublimation Fusion	c c	g liq	760	194.68 217.0	-78.48 -56.2	6.031 1.99	30.98 9.2	
CH <sub>4</sub>	methane	Fusion Vaporization Vaporization	c liq liq	liq g g	87.7 87.7 760	90.68 90.68 111.67	-182.48 -182.48 -161.49	0.225 1.955	2.48 17.51	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 23-2. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1948; March 31, 1950

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		°K	°C			
CH <sub>2</sub> O	formaldehyde		Fusion Vaporization	c liq	liq g	760	154.9 253.9	-118.3 -19.3	5.85	23.0	
CH <sub>2</sub> O <sub>2</sub>	formic acid		Fusion Vaporization Vaporization	c liq liq	liq g g	18 18 760	281.46 281.46 373.7	8.30 8.30 100.5	3.03 5.32	10.8 14.24	8.8
CH <sub>4</sub> O	methanol		Transition Fusion Vaporization Vaporization Vaporization	c, II c, I liq liq, std. liq	c, I liq g g, std. g		157.4 175.26 298.16 298.16 337.9	-115.8 -97.90 25.00 25.00 64.7	0.154 0.757 8.94 8.94 8.43	0.98 4.32 29.98 26.48 24.95	4.2
CH <sub>4</sub> O <sub>2</sub>	methyl hydrogen peroxide		Vaporization	liq	g	34	298.	25.	7.9	26.5	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 23-3. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1948; March 31, 1950

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		°K	°C			
CF <sub>4</sub>	tetrafluoromethane	Transition	c, II	c, I		76.23	-196.93	0.95	4.6	
		Fusion	c, I	liq		89.47	-183.69	0.167	1.87	
		Vaporization	liq	g	760	145.14	-128.02	3.01	20.7	
COF <sub>2</sub>	carbonyl fluoride	Fusion	c	liq	106	159.2	-114.0			
		Vaporization	liq	g	760	189.9	-83.3	3.86	20.3	
CH <sub>3</sub> F	fluoromethane	Fusion	c	liq						
		Vaporization	liq	g	760	195.1	-78.1	4.23	21.7	
CH <sub>2</sub> F <sub>2</sub>	difluoromethane	Fusion	c	liq						
		Vaporization	liq	g	760	221.	-52.			
CHF <sub>3</sub>	trifluoromethane	Fusion	c	liq						
		Vaporization	liq	g	760	113. 189.0	-160. -84.2	4.4	23.	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 23-4. Carbon (at. no., 6; at. wt., 12.010) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION March 31, 1948; March 31, 1950										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$CCl_4$	tetrachloromethane	Transition	c, II c, I liq	c, I liq g		225.5	-47.7	1.09	4.8	1.1
		Fusion				250.3	-22.9	0.60	2.4	
		Vaporization				349.9	76.7	7.17	20.5	
$COCl_2$	carbonyl chloride	Fusion	c liq	liq g	760	140.37	-127.79	1.371	9.43	
		Vaporization				280.72	7.56	5.832	20.78	
$CH_3Cl$	chloromethane	Fusion	c liq liq	liq g g	65.66 65.66 760	175.44	-97.72	1.537	8.76	1.5
		Vaporization				175.44	-97.72			
		Vaporization				248.94	-24.22	5.15	20.7	
$CH_2Cl_2$	dichloromethane	Fusion	c liq	liq g	760	176.	-97.	1.1	6.	
		Vaporization				313.	40.	6.69	21.4	
$CHCl_3$	trichloromethane	Fusion	c liq	liq g	760	209.7	-63.5	2.2	10.5	
		Vaporization				334.4	61.2	7.02	20.99	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 23-5. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1948; March 31, 1950

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
$CF_3Cl$		trifluorochloromethane	Fusion Vaporization	c liq	liq g	760	91.6 192.0	-181.6 -81.2	3.92	20.4	
$CF_2Cl_2$		difluorodichloromethane	Fusion Vaporization	c liq	liq g	760	118. 242.7	-155. -30.5	0.99 4.85	8.4 20.0	
$CFCl_3$		fluorotrichloromethane	Fusion Vaporization	c liq	liq g	760	162.68 296.8	-110.68 23.6	1.648 5.96	10.13 20.1	-10.4
$COFCl$		carbonyl fluorochloride	Fusion Vaporization	c liq	liq g	760	135. 231.	-138. -42.	5.3	23.	
$CH_2FCl$		fluorochloromethane	Fusion Vaporization	c liq	liq g	760	264.	-9.	5.6	21.	
$CHF_2Cl$		difluorochloromethane	Fusion Vaporization	c liq	liq g	760	113. 232.4	-160. -40.8	5.1	22.	
$CHFCI_2$		fluorodichloromethane	Fusion Vaporization	c liq	liq g	760	138. 282.1	-135. 8.9	6.4	23.	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 23-6. Carbon (at. no., 6; at. wt. 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1948; March 31, 1950

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		°K	°C			
CBr <sub>4</sub>	tetrabromomethane	Transition	c, II	c, I		320.1	46.9	1.5	4.7	
		Fusion	c, I	liq		363.3	90.1	0.98	2.70	
		Vaporization	liq	g	760	460.	187.	10.4	22.6	
COBr <sub>2</sub>	carbonyl bromide	Fusion	c	liq						
		Vaporization	liq	g	760	333.	60.	7.2	21.6	
CH <sub>3</sub> Br	bromomethane	Transition	c, II	c, I		173.79	-99.37	0.113	0.650	
		Fusion	c, I	liq		179.48	-93.68	1.429	7.96	
		Vaporization	liq	g	760	276.72	3.56	5.715	20.65	
CH <sub>2</sub> Br <sub>2</sub>	dibromomethane	Fusion	c	liq		220.5	-52.7	1.	4.	
		Vaporization	liq	g	760	370.4	97.2	8.6	23.	
CHBr <sub>3</sub>	tribromomethane	Fusion	c	liq		281.21	8.05	2.65	9.42	
		Vaporization	liq	g	22	323.	50.	10.4	32.2	
		Vaporization	liq	g	760	422.7	149.5			



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

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Table 23-7. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1948; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
CF <sub>3</sub> Br	trifluorobromomethane	Fusion Vaporization	c liq	liq g	740	213.	-60.			
CF <sub>2</sub> Br <sub>2</sub>	difluorodibromomethane	Fusion Vaporization	c liq	liq g	760	24.5	297.7			
CFBr <sub>3</sub>	fluorotribromomethane	Fusion Vaporization	c liq	liq g	760	380.	107.			
CHF <sub>2</sub> Br	difluorobromomethane	Fusion Vaporization	c liq	liq g	760	258.7	-14.5			
CHBrF <sub>2</sub>	fluorodibromomethane	Fusion Vaporization	c liq	liq g	757	338.1	64.9			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

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Table 23-8. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1948; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$CI_4$	tetraiodomethane	Fusion	c	liq		444.	171.			
$CH_3I$	iodomethane	Fusion	c	liq		206.7	-66.5			
		Vaporization	liq	g	760	315.6	42.4	6.7	21.2	
$CH_2I_2$	diiodomethane	Fusion	c II	liq		278.76	5.60	3.02	10.8	
		Fusion	c, I	liq		279.26	6.10	2.88	10.3	
		Vaporization	liq	g	330	425.	152.			
$CHI_3$	triiodomethane	Fusion	c	liq		398.	125.	3.9	9.8	
$CHF_2I$	difluoroiodomethane	Fusion	c	liq		151.2	-122.0			
		Vaporization	liq	g	760	294.7	21.5	6.29	21.3	
$CHF_2I_2$	fluorodiiodomethane	Fusion	c	liq		238.7	-34.5			
		Vaporization	liq	g	100	329.2	56.0	7.8	24.	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

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Table 23-9. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1948; March 31, 1950

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		°K	°C			
CS <sub>2</sub>	carbon disulfide	Fusion Vaporization	c liq	liq g	760	161.1 319.41	-112.1 46.25	1.05 6.40	6.52 20.0	
COS	carbon oxysulfide	Fusion Vaporization	c liq	liq g	760	134.34 222.92	-138.82 -50.24	1.130 4.423	8.41 19.84	5.6
CH <sub>4</sub> S	methanethiol	Transition Fusion Vaporization	c, II c, I liq	c, I liq g	760	137.6 150.16 279.12	-135.6 -123.00 5.96	0.052 1.411 5.872	0.38 9.40 21.04	4.8
H <sub>2</sub> CS <sub>3</sub>	trithiocarbonic acid	Fusion	c	liq		242.7	-30.5			
CS <sub>2</sub>	carbon diselenide	Fusion Vaporization	c liq	liq g	760	227.7 397.	-45.5 124.		22.	
COSe	carbon oxyselenide	Fusion Vaporization	c liq	liq g	760	151.1 251.6	-122.1 -21.6	5.4	21.	
CSSe	carbon sulfoselenide	Fusion Vaporization	c liq	liq g	760	188. 357.5	-85. 84.3	7.56	21.2	
CSTe	carbon sulfotelluride	Fusion	c	liq		219.	-54.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 23-10. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1948; March 31, 1950

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		°K	°C			
$C(NO_2)_4$	tetranitromethane		Fusion Vaporization	c liq	liq g	760	286. 398.9	13. 125.7	9.2	23.	
HCN	hydrogen cyanide		Transition Fusion Vaporization Vaporization	c, II c, I liq liq	c, I liq g g	140.4 140.4 760	170.41 259.92 259.92 298.86	-102.75 -13.24 -13.24 25.70	0.004 2.009 6.027	0.02 7.73 20.17	1.7
$CH_5N$	methyl amine		Fusion Vaporization	c liq	liq g	760	179.70 266.84	-93.46 -6.32	1.466 6.17	8.16 23.1	
$CH_2N_2$	cyanamide		Fusion	c	liq		316.0	42.8	2.1	6.6	
HNCO	hydrogen isocyanate		Fusion Vaporization	c liq	liq g	202	193. 267.3	-80. -5.9	7.2	27.	
$CH_3ON$	formamide		Fusion	c	liq		275.71	2.55	1.6	5.8	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$0_K$	$0_C$			
$\text{CH}_3\text{O}_2\text{N}$	nitromethane	Fusion	c	liq		244.78	-28.38	2.319	9.47	
		Vaporization	liq	g	36.7	298.16	25.00	9.147	30.68	
		Vaporization	liq	g	760	374.0	100.8			
$\text{CH}_3\text{O}_2\text{N}$	methyl nitrite	Fusion	c	liq						
		Vaporization	liq	g	760	255.	-18.	5.0	19.7	
$\text{CH}_3\text{O}_3\text{N}$	methyl nitrate	Fusion	c	liq						
		Vaporization	liq	c	760	339.7	66.5	7.8	23.0	
$\text{CH}_4\text{ON}_2$	urea	Fusion	c	liq		405.8	132.6	3.60	8.9	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 23-12. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1948; March 31, 1950

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta T_p$ cal/deg mole
				Initial	Final		°K	°C			
CFN	cyanogen fluoride		Sublimation	c	g	760	201.	-72.	5.66	28.2	
COF <sub>3</sub> N	N,N-di fluoro fluoroformamide		Fusion Vaporization	c liq	liq g	760	120.7 190.9	-152.5 -82.3	4.46	21.8	
COF <sub>3</sub> N	trifluoronitrosomethane		Fusion Vaporization	c liq	liq g	760	77.2 189.1	-196.0 -84.1	4.20	22.2	
CCl <sub>3</sub> N	cyanogen chloride		Fusion Vaporization	c liq	liq g	760	286.3 286.1	-6.9 12.9	2.72 6.29	10.2 22.0	
CO <sub>2</sub> Cl <sub>3</sub> N	trichloronitromethane		Fusion Vaporization	c liq	liq g	760	385.56	112.40	8.71	22.6	
CH <sub>6</sub> Cl <sub>3</sub> N	methylamine hydrochloride		Transition Transition	c, III c, II	c, II c, I		220.4 264.5	-52.8 -8.7	0.425 0.674	1.93 2.55	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 23-13. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1948; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
CBrl	cyanogen bromide	Sublimation	c	g	119.5	298.	25.	11.3	37.9	-
		Fusion	c	liq		324.5	51.3			
		Vaporization	liq	g	760	334.5	61.3			
CIN	cyanogen iodide	Sublimation	c	g	760	413.0	139.8	14.2	34.4	
		Fusion	c	liq		419.	146.			
HCNS	hydrogen thiocyanate	Fusion	c	liq		173.	-110.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 23-14. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1948; March 31, 1950

Formula	Substance	Type, of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		°K	°C			
NH <sub>4</sub> SCN	ammonium thiocyanate	Transition	c, II	c, I	760	360.9	87.7	0.789	2.13	
COF <sub>2</sub> NP	phosphorus difluoroisocyanate	Fusion Vaporization	c liq	liq g	760	165. 285.5	-108. 12.3			
CF <sub>2</sub> SNP	phosphorus difluoroisothiocyanate	Vaporization	liq	g	760	363.5	90.3	8.0	22.	
CCl <sub>2</sub> SNP	phosphorus dichloroethiocyanate	Fusion Vaporization	c liq	liq g	760	197. 421.	-76. 148.	12.6	29.9	
CH <sub>3</sub> As	methyl arsine	Fusion Vaporization	c liq	liq g	760	130. 275.0	-143. 1.8	5.34	19.4	
CH <sub>3</sub> F <sub>2</sub> As	methyl difluoroarsine	Fusion Vaporization	c liq	liq g	760	243.5 349.7	-29.7 76.5	8.5	24.	
CH <sub>3</sub> Cl <sub>2</sub> As	methyl dichloroarsine	Fusion Vaporization	c liq	liq g	760	230.7 405.7	-42.5 132.5	8.9	22.	



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 23-15. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948 (Corrected)

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$0^\circ K$	$^\circ C$			
$C_2H_2$		ethyne (acetylene)	Sublimation	c	g	760	189.2	-84.0	5.1	27.	
			Fusion	c	liq	900	191.7	-81.5	0.9	5.	
			Vaporization	liq	g	900	191.7	-81.5	4.2	22.	
$C_2H_4$		ethene (ethylene)	Fusion	c	liq	0.9	103.97	-169.19	0.8008	7.702	
			Vaporization	liq	g	0.9	103.97	-169.19			
			Vaporization	liq	g	760	169.45	-103.71	3.237	19.10	
$C_2H_6$		ethane	Fusion	c	liq	0.006	89.89	-183.27	0.6834	7.603	2.2
			Vaporization	liq	g	0.006	89.89	-183.27			
			Vaporization	liq	g	760	184.53	-88.63	3.517	19.06	-11.5
$C_2H_2O$		ketene	Fusion	c	liq		122.	-151.			
			Vaporization	liq	g	760	217.	-56.			
$C_2H_2O_2$		glyoxal	Fusion	c	liq		288.	15.			
			Vaporization	liq	g	776	324.	51.	9.	28.	
$C_2H_2O_4$		oxalic acid	Sublimation	c	g	0.35	373.	100.	21.7	58.2	
$C_2H_4O$		acetaldehyde	Fusion	c	liq		155.	-118.	0.77	5.0	
			Vaporization	liq	g	760	293.3	20.1	6.5	22.2	-9.3
$C_2H_4O \cdot H_2O$		acetaldehyde monohydrate	Fusion	c	liq		319.2	46.0	1.0	3.1	8.1
$C_2H_4O$		ethylene oxide	Fusion	c	liq	760	160.71	-112.45	1.236	7.69	3.45
			Vaporization	liq	g	760	283.72	10.56	6.101	21.50	-9.7

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

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Table 23-16. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2H_4O_2$	acetic acid	Fusion	c, I	liq	760	289.77	16.61	2.80	9.66	9.4
		Transition	c, I	c, II	455000	328.9	55.7	0.11	0.3	
		Fusion	c, I	liq	455000	328.9	55.7	2.78	8.5	
		Fusion	c, II	liq	455000	328.9	55.7	2.89	8.8	
		Vaporization	liq	g, equilibrium	760	391.4	118.2	5.83	14.9	
$C_2H_4O_2$	methyl formate	Fusion	c	liq		174.	-99.	1.6	9.2	
		Vaporization	liq	g	760	304.7	31.5	6.75	22.2	
$C_2H_4O_2$	hydroxyacetaldehyde	Fusion	c	liq		370.	97.			
$C_2H_4O_3$	hydroxyacetic acid (glycollic acid)	Fusion	c, II	liq		336.	63.	2.0	6.	
		Fusion	c, I	liq		351.	78.	2.1	6.	
$C_2H_6O$	ethanol	Fusion	c	liq		158.6	-114.6	1.200	7.57	5.70
		Vaporization	liq	g	58.6	298.16	25.00	10.12	33.94	
		Vaporization	liq, std.	g, std.	760	298.16	25.00	10.12	28.99	
		Vaporization	liq	g	760	351.7	78.5	9.22	26.22	
$C_2H_6O$	dimethyl ether	Fusion	c	liq		131.66	-141.50	1.180	8.96	6.8 -10.6
		Vaporization	liq	g	760	248.34	-24.82	5.141	20.70	
$C_2H_6O_2$	1,2-ethanediol (ethylene glycol)	Fusion	c	liq.		260.0	-13.2			
		Vaporization	liq	g	760	470.4	197.2	13.6	28.9	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 23-17. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2F_4$		tetrafluoroethene	Fusion Vaporization	c liq	liq g	760	130.7 197.	-142.5 -76.	4.3	22.	
$C_2F_6$		hexafluoroethane	Transition Fusion Vaporization	c, II c, I liq	c, I liq g	760	103.98 173.10 194.9	-169.18 -100.06 -78.3	0.893 0.642 3.857	8.59 3.71 19.79	1.2 1.78
$C_2OF_4$		trifluoroacetyl fluoride	Vaporization	liq	g	760	214.	-59.			
$C_2H_3F$		fluoroethene	Fusion Vaporization	c liq	liq g	760	112.7 201.0	-160.5 -72.2	4.1	20.	
$C_2H_5F$		fluoroethane	Vaporization	liq	g	760	241.	-32.			
$C_2H_4F_2$		1,1-difluoroethane	Fusion Vaporization	c liq	liq g	760	156. 246.7	-117. -26.5	5.7	23.	
$C_2HF_3$		trifluoroethene	Vaporization	liq	g	760	221.	-52.			
$C_2H_3F_3$		1,1,1-trifluoroethane	Transition Fusion Vaporization	c, II c, I liq	c, I liq g	709.1	156.35 161.82 224.40	-116.81 -111.34 -48.76	0.071 1.48 4.583	0.45 9.1 20.42	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

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Table 23-18. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$C_2H_3OF$	acetyl fluoride	Vaporization	liq	g	760	293.6	20.4	6.2	21.	
$C_2H_3O_2F$	fluoroacetic acid	Fusion	c	liq		308.4	35.2			
$C_2H_5OF$	2-fluoroethanol	Fusion	c	liq		246.7	-26.5			
		Vaporization	liq	g	760	376.6	103.4	10.5	27.9	
$C_2H_2O_2F_2$	difluoroacetic acid	Fusion	c	liq		272.1	-1.1			
		Vaporization	liq	g	760	407.4	134.2			
$C_2H_4OF_2$	2,2-difluoroethanol	Fusion	c	liq		245.0	-28.2			
		Vaporization	liq	g	750	368.7	95.5			
$C_2H_2O_2F_3$	trifluoroacetic acid	Fusion	c	liq		257.9	-15.3			
		Vaporization	liq	g	760	345.6	72.4			
$C_2H_3OF_3$	2,2,2-trifluoroethanol	Fusion	c	liq		229.7	-43.5			
		Vaporization	liq	g	760	347.3	74.1			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

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Table 23-19. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2Cl_2$	dichloroethyne		Vaporization	liq	g	740	302.	29.			
$C_2Cl_4$	tetrachloroethene		Fusion	c	liq		250.8	-22.4	2.5	10.	
			Vaporization	liq	g	760	394.2	121.0	8.3	21.	
$C_2Cl_6$	hexachloroethane		Transition	c, II	c, I	9.5	344.4	71.2	1.8	5.	
			Sublimation	c, I	g	760	457.6	184.4	12.2	26.7	
$C_2O_2Cl_4$	trichloromethyl chloroformate		Vaporization	liq	g	760	400.	127.	11.	28.	
$C_2H_3Cl$	chloroethene		Fusion	c	liq		113.5	-159.7			
			Vaporization	liq	g	760	259.3	-13.9	5.5	21.	
$C_2H_5Cl$	chloroethane		Fusion	c	liq		134.83	-138.33	1.064	7.89	1.52
			Vaporization	liq	g	760	285.43	12.27	5.9	21.	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

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Table 23-20. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2H_2Cl_2$	1,1-dichloroethene	Fusion Vaporization	c liq	liq g	760	150.7 304.9	-122.5 31.7	6.3	21.	
$C_2H_2Cl_2$	cis-1,2-dichloroethene	Fusion Vaporization	c liq	liq g	760	192.7 333.0	-80.5 59.8	7.15	21.5	
$C_2H_2Cl_2$	trans-1,2-dichloroethene	Fusion Vaporization	c liq	liq g	760	223. 321.7	-50. 48.5	7.0	21.8	
$C_2H_4Cl_2$	1,1-dichloroethane	Fusion Vaporization	c liq	liq g	760	176.5 330.6	-96.7 57.4	1.3 7.3	7. 22.	
$C_2H_4Cl_2$	1,2-dichloroethane	Fusion Vaporization	c liq	liq g	760	237.5 356.7	-35.7 83.5	2.11 7.7	8.9 22.	5.9
$C_2HCl_3$	trichloroethene	Fusion Vaporization	c liq	liq g	760	186.8 360.3	-86.4 87.1	7.6	21.	
$C_2H_3Cl_3$	1,1,1-trichloroethane	Transition Fusion Vaporization	c, II c, I liq liq	c, I liq g g		224.20 240.4 286.53 347.2	-48.96 -32.8 13.37 74.0	1.786 0.8 7.96	7.97 3.3 27.8	

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Table 23-21. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2H_3Cl_3$	1,1,2-trichloroethane	Fusion Vaporization	c liq	liq c	760	236.5 387.1	-36.7 113.9	8.8	22.7	
$C_2H_2Cl_4$	1,1,1,2-tetrachloroethane	Fusion Vaporization	c liq	liq g	760	204.5 403.7	-68.7 130.5	8.8	21.8	
$C_2H_2Cl_4$	1,1,2,2-tetrachloroethane	Fusion Vaporization	c liq	liq g	760	230.7 418.	-42.5 145.	9.2	22.	
$C_2HCl_5$	pentachloroethane	Fusion Vaporization	c liq	liq g	760	244.2 432.5	-29.0 159.3	2.7 9.7	11. 22.4	
$C_2H_3OCl$	acetyl chloride	Fusion Vaporization	c liq	liq g	760	161. 323.6	-112. 50.4	6.93	21.2	
$C_2H_3OCl$	chloroacetaldehyde	Vaporization	liq	g	760	358.	85.			
$C_2H_3O_2Cl$	chloroacetic acid	Fusion Fusion Fusion Vaporization	c, III c, II c, I liq	liq liq liq g		324. 329. 334. 403.	51. 56. 61. 130.	3.8 4.5 4.6 13.	12. 14. 14. 32.	6.

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-22. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2H_5OCl$		2-chloroethanol	Fusion Vaporization	c liq	liq g		205.7. 401.8	-67.5 128.6		24.64	
$C_2H_5OCl$		chloromethyl methyl ether	Vaporization	liq	g	760	332.7	59.5			
$C_2H_2OCl_2$		chloroacetyl chloride	Vaporization	liq	g	760	378.	105.			
$C_2H_2OCl_2$		dichloroacetaldehyde	Vaporization	liq	g	760	362.	89.			
$C_2H_2O_2Cl_2$		dichloroacetic acid	Fusion Vaporization	c liq	liq g		284.0 405.	10.8 132.	1.83 14.	6.44 35.	-3.
$C_2H_4O_2Cl_2$		dichloroacetaldehyde monohydrate	Fusion	c	liq		330.	57.			
$C_2HOCl_3$		trichloroacetaldehyde (chloral)	Fusion Vaporization	c liq	liq g		216. 370.9	-57. 97.7		21.8	
$C_2HO_2Cl_3$		trichloroacetic acid	Fusion Vaporization	c liq	liq g	760	332.3 468.	59.1 195.	1.4 13.4	4.2 28.6	-16.7
$C_2H_3OCl_3$		2,2,2-trichloroethanol	Fusion Vaporization	c liq	liq g		291.0 424.	17.8 151.			
$C_2H_3O_2Cl_3$		trichloroacetaldehyde monohydrate	Fusion	c	liq	740	324.9	51.7	5.50	17.	44.



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-23. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2F_3Cl$	trifluorochloroethene	Fusion Vaporization	c liq	liq g	760	115.7 245.3	-157.5 -27.9	5.23	21.3	
$C_2F_5Cl$	pentafluorochloroethane	Vaporization	liq	g	760	235.	-38.			
$C_2F_2Cl_2$	1,1-difluoro-2,2-dichloroethene	Fusion Vaporization	c liq	liq g	760	158. 292.	-115. 19.			
$C_2F_2Cl_2$	cis-1,2-difluoro-1,2-dichloroethene	Fusion Vaporization	c liq	liq g	760	142.7 294.3	-130.5 21.1			
$C_2F_2Cl_2$	trans-1,2-difluoro-1,2-dichloroethene	Fusion Vaporization	c liq	liq g	760	162.9 295.2	-110.3 22.0			
$C_2F_4Cl_2$	1,1,1,2-tetrafluoro-2,2-dichloroethane	Vaporization	liq	g	760	271.	-2.			
$C_2F_4Cl_2$	1,1,2,2-tetrafluoro-1,2-dichloroethane	Fusion Vaporization	c liq	liq g	1 760	179. 277.0	-94. 3.8	6.09	22.0	
$C_2FCl_3$	fluorotrichloroethene	Fusion Vaporization	c liq	liq g	740.8	191. 343.2	-82. 70.0			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-24. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2F_3Cl_3$		1,1,2-trifluoro-1,2,2-trichloroethane	Fusion Vaporization	c liq	liq g	760	238. 320.4	-35. 47.2			
$C_2F_2Cl_3$		1,1,1-trifluoro-2,2,2-trichloroethane	Fusion Vaporization	c liq	liq g	760	287.4 319.1	14.2 45.9	6.8	21.2	-13.1
$C_2F_2Cl_4$		1,2-difluoro-1,1,2,2-tetrachloroethane	Fusion Vaporization Vaporization	c liq liq	liq g g	63 63 760	301.3 301.3 365.1	28.1 28.1 91.9	8.59	23.5	
$C_2F_2Cl_4$		1,1-difluoro-1,2,2,2-tetrachloroethane	Fusion Vaporization	c liq	liq g	760	313.8 364.7	40.6 91.5			
$C_2FCl_5$		fluoropentachloroethane	Fusion Vaporization	c liq	liq g	760	374.5 410.7	101.3 137.5			
$C_2OF_3Cl$		trifluoroacetyl chloride	Fusion Vaporization	c liq	liq g	760	127. 245.	-146. -28.	4.7	19.	
$C_2OF_2Cl_2$		difluorochloroacetyl chloride	Vaporization	liq	g	760	307.	34.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-25. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$C_2H_3F_2Cl$	1,1-difluoro-1-chloroethane	Fusion Vaporization	c liq	liq g	760	142.4 263.95	-130.8 -9.21	0.642 5.35	4.51 20.27	1.3
$C_2H_3F_2Cl$	1,1-difluoro-2-chloroethane	Vaporization	liq	g	760	309.	36.			
$C_2H_2F_3Cl$	1,1,2-trifluoro-2-chloroethane	Vaporization	liq	g	760	290.	17.			
$C_2HFC1_2$	2-fluoro-1,1-dichloroethene	Vaporization	liq	g	760	310.7	37.5			
$C_2H_2F_2Cl_2$	1,1-difluoro-2,2-dichloroethane	Vaporization	liq	g	760	333.	60.			
$C_2H_2FCl_3$	1-fluoro-1,2,2-trichloroethane	Vaporization	liq	g	760	376.	103.			
$C_2HFC1_4$	2-fluoro-1,1,1,2-tetrachloroethane	Vaporization	liq	g	760	389.7	116.5			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-26. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2H_2OFCl$	fluoroacetyl chloride	Vaporization	liq	g	760	343.	70.	8.8	26.	
$C_2H_2OFCl$	chloroacetyl fluoride	Vaporization	liq	g	760	348.	75.	9.1	26.	
$C_2HOF_2Cl$	difluoroacetyl chloride	Vaporization	liq	g	760	298.	25.			
$C_2H_2O_2F_2Cl$	difluorochloroacetic acid	Fusion	c	liq		296.1	22.9			
		Vaporization	liq	g	760	394.7	121.5			
$C_2H_2OFCl_2$	dichloroacetyl fluoride	Vaporization	liq	g	760	344.	71.			
$C_2Br_4$	tetrabromoethene	Sublimation	c	g	7.5	298.2	25.0	7.3	24.	
		Fusion	c	liq		329.7	56.5			
$C_2H_3Br$	bromoethene	Fusion	c	liq		135.	-138.	5.5	19.0	
		Vaporization	liq	g	760	288.9	15.7			
$C_2H_3Br$	bromoethane	Fusion	c	liq		155.0	-118.2	1.4	9.	-7.
		Vaporization	liq	g	760	311.5	38.3	7.3	23.	
$C_2H_2Br_2$	1,1-dibromoethene	Vaporization	liq	g	760	349.	76.			
$C_2H_2Br_2$	cis-1,2-dibromoethene	Fusion	c	liq		220.	-53.			
		Vaporization	liq	g	760	385.7	112.5			
$C_2H_2Br_2$	trans-1,2-dibromoethene	Fusion	c	liq		266.7	-6.5			
		Vaporization	liq	g	760	381.	108.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-27. Carbon (at. no., 6; at. wt., 12.010)

HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION

June 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2H_4Br_2$	1,1-dibromoethane	Vaporization	liq	g	760	383.7	110.5			
$C_2H_4Br_2$	1,2-dibromoethane	Transition	c, II	c, I		249.65	-23.51	0.464	1.86	-11.1
		Fusion	c, I	liq		283.1	9.9	2.616	9.24	7.9
		Vaporization	liq	g	760	404.9	131.7	8.7	21.	
$C_2H_3Br_3$	1,1,2-tribromoethane	Fusion	c	liq		244.0	-29.2			
		Vaporization	liq	g	760	461.	188.	10.	22.	
$C_2H_2Br_4$	1,1,1,2-tetrabromoethane	Vaporization	liq	g	18	385.7	112.5			
$C_2H_2Br_4$	1,1,2,2-tetrabromoethane	Fusion	c	liq		273.3	0.1	2.4	8.8	
		Vaporization	liq	g	12	387.	114.	13.	34.	
$C_2HBr_5$	pentabromoethane	Fusion	c	liq		330.	57.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-28. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		O <sub>K</sub>	O <sub>C</sub>			
$C_2H_3OBr$	acetyl bromide	Fusion Vaporization	c liq	liq g	760	177. 349.1	-96. 75.9	7.25	20.8	
$C_2H_3OBr$	bromoacetaldehyde	Vaporization	liq	g	750	378.	105.			
$C_2H_3O_2Br$	bromoacetic acid	Fusion Vaporization	c liq	liq g	760	322.7 481.	49.5 208.	13.	27.	
$C_2H_5OBr$	2-bromoethanol	Vaporization	liq	g	20	330.	57.			
$C_2H_5OBr$	bromomethyl methyl ether	Vaporization	liq	g	760	360.	87.			
$C_2H_2OBr_2$	bromoacetyl bromide	Vaporization	liq	g	760	422.	149.			
$C_2H_2O_2Br_2$	dibromoacetic acid	Fusion Vaporization	c liq	liq g	760	321. 503.	48. 230.			
$C_2H_4OBr_2$	bis(bromomethyl) ether	Vaporization	liq	g	760	416.	143.	12.	29.	
$C_2H_4OBr_3$	tribromoacetaldehyde (bromal)	Vaporization	liq	g	760	447.	174.	8.8	20.	
$C_2H_2O_2Br_3$	tribromoacetic acid	Fusion	c	liq		404.	131.			
$C_2H_3OBr_3$	2,2,2-tribromoethanol	Fusion	c	liq		353.	80.			
$C_2H_3O_2Br_3$	tribromoacetaldehyde monohydrate	Fusion	c	liq		326.7	53.5	5.05	15.	39.

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-29. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2F_3Br$	trifluorobromoethene	Vaporization	liq	g	760	270.7	-2.5			
$C_2F_5Br$	pentafluorobromoethane	Vaporization	liq	g	740	251.	-22.			
$C_2F_4Br_2$	1,1,2,2-tetrafluoro-1,2-dibromoethane	Fusion	c	liq		161.7	-111.5			
		Vaporization	liq	g	760	320.5	47.3	7.2	22.5	
$C_2FBr_3$	fluorotribromoethene	Vaporization	liq	g	760	420.4	147.2			
$C_2F_2Br_4$	1,1-difluoro-1,2,2,2-tetrabromoethane	Fusion	c	liq		372.	99.			
		Vaporization	liq	g	760	458.	185.			
$C_2OF_3Br$	trifluoroacetyl bromide	Fusion	c	liq		137.	-136.			
		Vaporization	liq	g	760	268.	-5.	4.9	18.	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-30. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2H_2FBr$		1-fluoro-1-bromoethene	Vaporization	liq	g	760	285.7	12.5			
$C_2H_4FBr$		1-fluoro-2-bromoethane	Vaporization	liq	g	760	345.	72.			
$C_2HF_2Br$		1,1-difluoro-2-bromoethene	Vaporization	liq	g	760	279.4	6.2			
$C_2H_3F_2Br$		1,1-difluoro-2-bromoethane	Vaporization	liq	g	760	330.	57.			
$C_2H_2F_3Br$		1,1,2-trifluoro-1-bromoethane	Vaporization	liq	g	760	298.	25.			
$C_2H_2F_3Br$		1,1,1-trifluoro-2-bromoethane	Vaporization	liq	g	760	299.7	26.5			
$C_2H_3FBr_2$		2-fluoro-1,1-dibromoethane	Vaporization	liq	g	760	390.7	117.5			
$C_2H_3FBr_2$		1-fluoro-1,2-dibromoethane	Fusion	c	liq		219.	-54.			
$C_2H_3FBr_2$		1-fluoro-1,2-dibromoethane	Vaporization	liq	g	760	395.7	122.5			
$C_2H_2F_2Br_2$		1,1-difluoro-2,2-dibromoethane	Vaporization	liq	g	760	380.7	107.5			
$C_2H_2F_2Br_2$		1,1-difluoro-1,2-dibromoethane	Vaporization	liq	g	760	365.	92.			
$C_2HF_3Br_2$		1,1,2-trifluoro-1,2-dibromoethane	Vaporization	liq	g	760	349.7	76.5			



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Table 23-31. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2H_2FBr_3$		1-fluoro-1,2,2-tribromoethane	Vaporization	liq	g	760	446.	173.			
$C_2H_2FBr_3$		1-fluoro-1,1,2-tribromoethane	Vaporization	liq	g	760	436.	163.			
$C_2HF_2Br_3$		1,1-difluoro-1,2,2-tribromoethane	Vaporization	liq	g	755	416.7	143.5			
$C_2HFBri_4$		1-fluoro-1,1,2,2-tetrabromoethane	Vaporization	liq	g	760	484.	211.			
$C_2H_2O_2FBr$		fluorobromoacetic acid	Fusion	c	liq		322.	49.			
			Vaporization	liq	g	760	456.	183.			
$C_2HOFBr_2$		fluorobromoacetyl bromide	Vaporization	liq	g	760	385.7	112.5			

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Table 23-32. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
	$C_2Cl_2Br_2$	1,2-dichloro-1,2-dibromoethene	Fusion Vaporization	c liq	liq g	760	278. 445.	5. 172.			
	$C_2Cl_4Br_2$	1,1,1,2-tetrachloro-1,2-dibromoethane	Sublimation	c	g	6.0	373.	100.	15.	40.	
	$C_2OCl_3Br$	trichloroacetyl bromide	Vaporization	liq	g	760	416.	143.	9.5	23.	
	$C_2OCl_2Br_2$	chlorodibromoacetyl chloride	Vaporization	liq	g	760	440.	167.			
	$C_2H_2ClBr$	cis-1-chloro-2-bromoethene	Fusion Vaporization	c liq	liq g	760	186. 357.8	-87. 84.6			
	$C_2H_2ClBr$	trans-1-chloro-2-bromoethene	Fusion Vaporization	c liq	liq g	760	232. 348.5	-41. 75.3			
	$C_2H_4ClBr$	1-chloro-1-bromoethane	Fusion Vaporization	c liq	liq g	760	289.8. 355.9	16.6 82.7	7.9	22.2	
	$C_2H_4ClBr$	1-chloro-2-bromoethane	Fusion Vaporization	c liq	liq g	760	256.6 379.9	-16.6 106.7	2.3 8.3	9.0 21.8	4.3
	$C_2HCl_2Br$	1,1-dichloro-2-bromoethene	Fusion Vaporization	c liq	liq g	760	214.7 380.	-58.5 107.			
	$C_2HCl_2Br$	cis-1,2-dichloro-1-bromoethene	Fusion Vaporization	c liq	liq g	760	189.7 387.	-83.5 114.			
	$C_2HCl_2Br$	trans-1,2-dichloro-1-bromoethene	Fusion	c	liq		185.	-88.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 23-33. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$C_2H_3Cl_2Br$	1,1-dichloro-2-bromoethane	Vaporization	liq	g	760	411.	138.			
$C_2H_2Cl_3Br$	1,2,2-trichloro-1-bromoethane	Fusion	c	liq		317.8	44.6			
		Vaporization	liq	g	760	445.	172.			
$C_2H_2Cl_4Br_2$	1,1-dichloro-1,2-dibromoethane	Fusion	c	liq		206.4	-66.8			
		Vaporization	liq	g	18	343.	70.			
$C_2H_2Cl_2Br_2$	1,2-dichloro-1,2-dibromoethane	Fusion	c	liq		246.	-27.			
		Vaporization	liq	g	24	363.2	90.0	10.8	29.7	
		Vaporization	liq	g	760	468.	195.			
$C_2HCl_3Br_2$	1,1,2-trichloro-1,2-dibromoethane	Fusion	c	liq		264.3	-8.9			
$C_2HCl_3Br_2$	1,1,1-trichloro-2,2-dibromoethane	Fusion	c	liq		268.7	-4.5			
$C_2H_2ClBr_3$	2-chloro-1,1,2-tribromoethane	Fusion	c	liq		252.6	-20.6			
		Vaporization	liq	g	14	370.	97.			
$C_2HCl_2Br_3$	1,2-dichloro-1,1,2-tribromoethane	Fusion	c	liq		279.2	6.0			
		Vaporization	liq	g	16	385.	112.			
$C_2HCl_2Br_3$	1,1-dichloro-1,2,2-tribromoethane	Fusion	c	liq		290.0	16.8			
		Vaporization	liq	g	16	379.	106.			
$C_2H_2OClBr$	bromoacetyl chloride	Vaporization	liq	g	760	400.	127.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 23-34. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2F_3ClBr_2$		1,1,2-trifluoro-2-chloro-1,2-dibromoethane	Vaporization	liq	g	760	366.1	92.9			
$C_2F_2Cl_2Br_2$		1,1-difluoro-2,2-dichloro-1,2-dibromoethane	Fusion Vaporization	c liq	liq g	760	319. 390.3	46. 117.1			
$C_2F_2Cl_2Br_2$		1,2-difluoro-1,2-dichloro-1,2-dibromoethane	Fusion Vaporization	c liq	liq g	760	306.1 413.0	32.9 139.8			
$C_2FCl_3Br_2$		1-fluoro-1,2,2-trichloro-1,2-dibromoethane	Fusion	c	liq		395.7	122.5			
$C_2OF_2ClBr$		fluorochlorobromoacetyl fluoride	Vaporization	liq	g	760	324.	51.			
$C_2HFC_2Br_2$		2-fluoro-1,1-dichloro-1,2-dibromoethane	Vaporization	liq	g	760	436.	163.			
$C_2HOFClBr$		fluorobromoacetyl chloride	Vaporization	liq	g	765	371.	98.			
$C_2H_2FC_2Br$		fluorochlorobromoacetic acid	Fusion Vaporization	c liq	liq g	760	268. 454.	-5. 181.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

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Table 23-35. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$C_2I_4$	tetraiodoethene	Fusion	c	liq		465.	192.			
$C_2H_5I$	iodoethane	Fusion	c	liq		162.	-111.			
		Vaporization	liq	g	760	345.5	72.3	7.15	20.7	
$C_2H_2I_2$	1,1-diiodoethene	Fusion	c	liq		329.	56.			
$C_2H_2I_2$	cis-1,2-diiodoethene	Fusion	c	liq		259.4	-13.8			
		Vaporization	liq	g	760	461.	188.			
$C_2H_4I_2$	trans-1,2-diiodoethene	Sublimation	c	g	0.04	293.	20.	10.	34.	
		Fusion	c	liq		345.	72.			
		Vaporization	liq	g	760	463.	190.			
$C_2H_4I_2$	1,2-diiodoethane	Sublimation	c	g	1	318.3	45.1	15.	47.	
		Fusion	c	liq		354.	81.			
$C_2H_3OI$	acetyl iodide	Vaporization	liq	g	760	381.	108.			
$C_2H_3O_2I$	iodoacetic acid	Fusion	c	liq		356.	83.			
$C_2H_5OI$	2-iodoethanol	Vaporization	liq	g	25	359.	86.			
$C_2H_5OI$	iodomethyl methyl ether	Vaporization	liq	g	760	397.	124.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

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Table 23-36. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		°K	°C			
$C_2H_3F_2I$	1,1-difluoro-2-iodoethane	Vaporization	liq	g	760	354.7	81.5			
$C_2H_2F_3I$	1,1,1-trifluoro-2-iodoethane	Vaporization	liq	g	730	328.	55.			
$C_2H_2O_2FI$	fluoroiodoacetic acid	Fusion	c	liq		365.7	92.5			
$C_2OCl_3I$	trichloroacetyl iodide	Vaporization	liq	g	760	453.	180.			
$C_2H_2ClI$	<i>cis</i> -1-chloro-2-iodoethene	Fusion Vaporization	c liq	liq g	760	232.2 386.	-41.0 113.			
$C_2H_2ClI$	<i>trans</i> -1-chloro-2-iodoethene	Fusion Vaporization	c liq	liq g	760	236.8 390.	-36.4 117.			
$C_2H_2OClI$	iodoacetyl chloride	Vaporization	liq	g	15	323.	50.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

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Table 23-37. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2H_6S$		ethanethiol	Fusion Vaporization	c liq	liq g	760	126. 307.6	-147. 34.4	6.9	22.	
$C_2H_6S$		dimethyl sulfide	Fusion Vaporization	c liq	liq g	365.5 760	174.86 291.06 309.1	-88.30 17.90 35.9	1.908 6.688	10.91 22.98	6.95
$C_2H_6OS$		dimethyl sulfoxide	Fusion Vaporization	c liq	liq g	0.6	281. 298.2	8. 25.0	12.64	42.39	
$C_2H_6O_2S$		dimethyl sulfone	Fusion Vaporization	c liq	liq g	760	382. 507.	109. 234.			
$C_2H_6O_3S$		dimethyl sulfite	Vaporization	liq	g	760	395.	122.			
$C_2H_6O_4S$		dimethyl sulfate	Fusion Vaporization	c liq	liq g	760	241. 461.7	-32. 188.5			
$C_2H_6O_4S$		ethylsulfuric acid	Fusion	c	liq		256.	-17.			
$C_2H_6O_4S \cdot H_2O$		ethylsulfuric acid monohydrate	Fusion	c	liq		278.6	5.4			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 23-38. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$C_2H_5O_3FS$	ethyl fluorosulfonate	Vaporization	liq	g	755	385.	112.	9.2	24.	
$C_2H_5O_3ClS$	ethyl chlorosulfonate	Vaporization	liq	g	760	425.	152.			
$C_2H_4Br_2S$	bis(bromomethyl) sulfide	Vaporization	liq	g	608	468.	195.	12.	26.	
$C_2H_6Se$	dimethyl selenide	Vaporization	liq	g	55	327.	54.			
$C_2H_6O_4Se$	dimethyl selenate	Vaporization	liq	g	760	341.	68.			
$C_2H_6Te$	dimethyl telluride	Vaporization	liq	g	770	367.	94.			
$C_2H_6TeF_2$	dimethyl tellurium difluoride	Fusion	c	liq		357.	84.			
$C_2H_6TeCl_2$	dimethyl tellurium dichloride	Fusion	c	liq		365.	92.			
$C_2H_6TeBr_2$	dimethyl tellurium dibromide	Fusion	c	liq		365.	92.			
$C_2H_6TeI_2$	dimethyl tellurium diiodide	Fusion	c	liq		403.	130.			



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-39. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance		Type of Process	State		Pressure	Temperature		$\Delta H$	$\Delta S$	$\Delta G_p$
Formula	Description		Initial	Final	mm Hg	$^{\circ}K$	$^{\circ}C$	kcal/mole	cal/deg mole	cal/deg mole
$C_2N_2$	Cyanogen	Fusion Vaporization Vaporization	c liq liq	liq g g	553.6 553.6 760	245.32 245.32 252.01	-27.84 -27.84 -21.15	1.938 5.576	7.900 22.126	5.29
$C_2H_3N$	acetonitrile	Fusion Vaporization	c liq	liq g	760	228.3 354.7	-44.9 81.5	2.13 7.83	9.3 22.1	
$C_2H_7N$	ethylamine	Fusion Vaporization	c liq	liq g	760	192.2 289.7	-81.0 16.5	6.7	23.	-14.4
$C_2H_7N$	dimethylamine	Fusion Vaporization	c liq	liq g	760	180.97 280.0	-92.19 6.8	1.420 6.33	7.85 22.6	9.81 -17.1
$C_2H_8N_2$	1,2-ethanediamine (ethylenediamine)	Fusion Vaporization	c liq	liq g	760	281. 390.6	8. 117.4	4.6 9.3	16.4 23.8	
$C_2H_8N_2$	1,1-dimethylhydrazine	Vaporization	liq	g	760	335.7	62.5			
$C_2H_8N_2$	1,2-dimethylhydrazine	Vaporization	liq	g	760	354.	81.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 23-40. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		°K	°C			
$C_2H_5ON$		acetamide	Fusion	c, II	liq	760	343.	71.	3.4	10.	
			Fusion	c, I	liq	760	355.1	81.9	3.75	10.56	15.1
			Transition	c, I	c, III	396000	400.2	127.0	0.14	0.35	0.7
			Fusion	c, I	liq	396000	400.2	127.0	3.60	9.00	
			Fusion	c, III	liq	396000	400.2	127.0	3.46	8.65	
$C_2H_5O_2N$		nitroethane	Fusion	c	liq		183.	-90.			
			Vaporization	liq	g	17	293.	20.	9.1	31.	
$C_2H_5O_2N$		ethyl nitrite	Vaporization	liq	g	760	290.1	16.9	6.64	22.9	
$C_2H_5O_3N$		ethyl nitrate	Fusion	c	liq		171.	-102.			
			Vaporization	liq	g	67.5	298.2	25.0	9.20	30.9	
			Vaporization	liq	g	760	361.9	88.7			
$C_2H_7ON$		2-aminoethanol	Fusion	c	liq		284.68	11.52			
			Vaporization	liq	g	760	444.3	171.1			
$C_2H_4O_3N_2$		2-nitroacetamide	Fusion	c	liq		380.	107.			
$C_2H_4O_6N_2$		glycol dinitrate	Fusion	c	liq		250.9	-22.3	4.5	18.	
			Vaporization	liq	g	19	378.	105.			
$C_2H_6O_2N_2$		ethylnitramine	Fusion	c	liq		279.	6.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

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Washington, D.C.

Table 23-41. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
	$C_2F_3N$	trifluoroacetonitrile	Vaporization	liq	g	760	212.	-61.			
	$C_2F_6N_2$	hexafluoroazomethane	Fusion	c	liq		140.	-133.			
			Vaporization	liq	g	760	241.6	-31.6	5.5	23.	
	$C_2H_2FN$	fluoroacetonitrile	Vaporization	liq	g	760	351.	78.	9.1	26.	
	$C_2H_5F_2N$	2,2-difluoroethylamine	Vaporization	liq	g	760	341.	68.			
	$C_2HF_6N$	bis(trifluoromethyl) amine	Vaporization	liq	g	760	267.1	-6.1	5.6	21.0	
	$C_2H_4OFN$	fluoroacetamide	Fusion	c	liq		381.	108.			
	$C_2H_4O_2FN$	2-fluoroethyl nitrite	Vaporization	liq	g	760	332.	59.	9.	27.	
	$C_2H_3OF_2N$	difluoroacetamide	Fusion	c	liq		325.0	51.8			
			Vaporization	liq	g	35	381.8	108.6			
	$C_2H_2OF_3N$	trifluoroacetamide	Fusion	c	liq		348.0	74.8			
			Vaporization	liq	g	760	435.7	162.5			
	$C_2H_4O_2F_2N_2$	2,2-difluoroethylnitramine	Fusion	c	liq		295.7	22.5			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards Washington, D.C.

Table 23-42. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		°K	°C			
chloroacetonitrile	$C_2H_2ClN$		Vaporization	liq	g	760	397.	124.			
<i>p</i> -chlorodimethylamine	$C_2H_6ClN$		Vaporization	liq	g	760	319.	46.			
chloroacetamide	$C_2H_4OCN$		Fusion	c	liq		390.	117.			
<i>p</i> -chloroacetamide	$C_2H_4OCN$		Fusion	c	liq		383.	110.			
difluorochloroacetamide	$C_2H_2OF_2ClN$		Fusion	c	liq		351.7	78.5			
			Vaporization	liq	g	760	407.	134.			
bromoacetamide	$C_2H_4OBrN$		Fusion	c	liq		381.	108.			
<i>p</i> -bromoacetamide	$C_2H_4OBrN$		Fusion	c	liq		379.	106.			
tribromoacetamide	$C_2H_2OBr_3N$		Fusion	c	liq		392.	119.			
fluorochlorobromoacetamide	$C_2H_2OFClBrN$		Fusion	c	liq		404.7	131.5			
methyl thiocyanate	$C_2H_3SN$		Fusion	c	liq		218.7	-54.5			
			Vaporization	liq	g	760	406.1	132.9	9.2	23.	
methyl isothiocyanate	$C_2H_3SN$		Fusion	c	liq		308.7	35.5			
			Vaporization	liq	g	760	392.2	119.0	8.9	23.	
dimethyltellurium dinitrate	$C_2H_6OTeV_2$		Fusion	c	liq		415.	142.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

SERIES II

Washington, D.C.

Table 23-43. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2H_7P$	dimethylphosphine	Vaporization	liq	g	760	294.8	21.6	6.0	20.4	
$C_2H_6O_3FP$	dimethylfluorophosphonate	Vaporization	liq	g	4.4	298.2	25.0	10.6	35.5	
$C_2O_2FN_2P$	phosphorous fluorodiiisocyanate	Fusion	c	liq		218.	-55.			
		Vaporization	liq	g	760	371.9	98.7	9.2	24.7	
$C_2O_2ClN_2P$	phosphorous chlorodiiisocyanate	Fusion	c	liq		223.	-50.			
		Vaporization	liq	g	760	407.8	134.6	11.2	27.5	
$C_2H_7As$	dimethylarsine	Vaporization	liq	g	747	308.8	35.6			
$C_2H_7As$	ethylarsine	Vaporization	liq	g	760	309.	36.			
$C_2H_5OAs$	ethylarsenic oxide	Vaporization	liq	g	168	489.	216.	17.	35.	
$C_2H_5F_2As$	ethyldifluoroarsine	Fusion	c	liq		234.5	-38.7			
		Vaporization	liq	g	760	367.5	94.3	8.3	22.5	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 23-44. Carbon (at. no., 6; at. wt., 12.010)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$C_2H_6ClAs$	dimethylchloroarsine	Vaporization	liq	g	760	380.	107.			
$C_2H_5Cl_2As$	ethylchloroarsine	Fusion	c	liq		208.	-65.			
		Vaporization	liq	g	760	428.5	155.3	9.2	21.5	
$C_2H_2Cl_3As$	2-chlorovinylidichloroarsine	Vaporization	liq	g	0.57	298.2	25.0	12.7	42.6	
$C_2H_2OClAs$	2-chlorovinylarsenic oxide	Fusion	c	liq		416.	143.			
$C_2H_4O_3ClAs$	2-chlorovinylarsonic acid	Fusion	c	liq		403.	130.			
$C_2H_2F_2Cl$	2-chlorovinylidifluoroarsine	Vaporization	liq	g	14.5	316.7	43.5			
$C_2H_6BrAs$	dimethylbromoarsine	Vaporization	liq	g	760	403.	130.			
$C_2H_5Br_2As$	ethylbromoarsine	Vaporization	liq	g	0.56	298.2	25.0	11.9	39.9	
$C_2H_4O_3BrAs$	2-bromovinylarsonic acid	Fusion	c	liq		416.	143.			
$C_2H_2ClBr_2As$	2-chlorovinylbromoarsine	Vaporization	liq	g	15	274.7	1.5			
$C_2H_6IAs$	dimethyliodoarsine	Vaporization	liq	g	760	428.	155.			
$C_2H_2ClI_2As$	2-chlorovinyl-diiodoarsine	Fusion	c	liq		311.	38.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 24-1. Silicon (at. no., 14; at. wt., 28.06)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Si		Fusion	c	liq		1683.	1410.	11.	6.5	
SiO <sub>2</sub>		Transition	quartz, c, III	quartz, c, II		91.	-182.			
		Transition	quartz, c, II	quartz, c, I		846.	573.	0.15	0.18	
		Transition	quartz, c, I	tridymite, c, I		1140.	867.	0.12	0.11	
		Fusion	quartz, c, I	liq		1883.	1610.	2.04	1.08	
		Transition	tridymite, c, IV	tridymite, c, III		390.	117.	0.07	0.18	
		Transition	tridymite, c, III	tridymite, c, II		436.	163.	0.04	0.09	
		Transition	tridymite, c, II	tridymite, c, I		598.	225.	0.05	0.08	
		Transition	tridymite, c, I	cristobalite, I		1743.	1470.	0.05	0.03	
		Fusion	tridymite, c, I	liq		1953.	1680.			
		Transition	cristobalite, II	cristobalite, I		515.	242.	0.31	0.60	
		Fusion	cristobalite, I	liq		2001.	1728.	1.84	0.92	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 24-2. Silicon (at. no., 14; at. wt., 28.06)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$SiH_4$	silane	Transition Fusion Vaporization	c, II c, I liq	c, I liq g		63.5 88.5 161.8	-209.7 -184.7 -111.4	0.147 0.159 2.9	2.32 1.80 18.	1.3
$Si_2H_6$	disilane	Fusion Vaporization	c liq	liq q	760	141. 258.9	-132. -14.3	5.2	20.1	
$Si_3H_8$	trisilane	Fusion Vaporization	c liq	liq q	760	156. 326.2	-117. 53.0	6.8	20.8	
$Si_4H_{10}$	tetrasilane	Fusion Vaporization	c liq	liq q	760	188.9 381.6	-84.3 108.4	8.2	21.5	
$Si_2H_6O$	disiloxane	Fusion Vaporization	c liq	liq g	760	129. 257.7	-144. -15.5	5.3	20.6	



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Table 24-3. Silicon (at. no., 14; at. wt., 28.06)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deq mole	$\Delta C_p$ cal/deq mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$SiF_4$	tetrafluorosilane	Sublimation	c	g	760	177.7	-95.5	6.15	34.6	
		Fusion	c	liq	1320	182.9	-90.3	1.69	9.2	
		Vaporization	liq	g	1320	182.9	-90.3	4.46	24.4	
$Si_2F_6$	hexafluorodisilane	Sublimation	c	g	760	254.1	-19.1	10.1	39.7	
		Fusion	c	liq	780	254.5	-18.7	3.5	13.8	
		Vaporization	liq	g	780	254.5	-18.7	6.6	25.9	
$Si_2OF_6$	hexafluorodisiloxane	Fusion	c	liq		225.4	-47.8			
		Vaporization	liq	g	760	249.9	-23.3	5.1	20.4	
		Vaporization	liq	g	760	175.	-98.	4.2	24.	
$SiH_3F$	difluorosilane	Fusion	c	liq		151.	-122.	1.	7.	
		Vaporization	liq	g	760	197.	-76.	3.9	20.	
$SiHF_3$	trifluorosilane	Fusion	c	liq		141.9	-131.3			
		Vaporization	liq	g	760	178.8	-94.4	3.8	21.	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 24-4. Silicon (at. no., 14; at. wt., 28.06)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		°K	°C			
$\text{SiCl}_4$	tetrachlorosilane		Fusion Vaporization	c liq	liq g	760	205. 330.2	-68. 57.0	1.84 7.0	9.0 21.2	
$\text{Si}_2\text{Cl}_6$	hexachlorodisilane		Fusion Vaporization	c liq	liq g	760	272. 418.	-1. 145.	10.	24.	
$\text{Si}_3\text{Cl}_8$	octachlorotrisilane		Vaporization	liq	g	760	485.	212.	15.	31.	
$\text{Si}_2\text{OCl}_6$	hexachlorodisiloxane		Fusion Vaporization	c liq	liq g	760	240. 409.5	-33. 136.3	9.3	22.7	
$\text{SiH}_3\text{Cl}$	chlorosilane		Fusion Vaporization	c liq	liq g	760	155. 242.8	-118. -30.4	5.1	21.0	
$\text{SiH}_2\text{Cl}_2$	dichlorosilane		Fusion Vaporization	c liq	liq g	760	151. 281.5	-122. 8.3	6.0	21.3	
$\text{SiHCl}_3$	trichlorosilane		Fusion Vaporization	c liq	liq g	760	146.7 304.9	-126.5 31.7	6.35	20.83	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 24-5. Silicon (at. no., 14; at. wt., 28.06)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$SiF_3Cl$	trifluorochlorosilane	Fusion Vaporization	c liq	liq g	760	133. 203.2	-140. -70.0	4.4	21.7	
$SiF_2Cl_2$	difluorodichlorosilane	Fusion Vaporization	c liq	liq g	760	133.5 241.3	-139.7 -31.9	5.0	20.7	
$SiFCl_3$	fluorotrichlorosilane	Fusion Vaporization	c liq	liq g	760	152. 285.5	-121. 12.3	6.0	21.0	
$Si_2O_2F_4Cl_2$	1,1,1,3-tetrafluoro-3,3-dichlorodisiloxane	Fusion Vaporization	c liq	liq g	760	213. 290.0	-60. 16.8	6.8	23.4	
$Si_2O_2F_3Cl_3$	1,1,1,1-trifluoro-3,3,3-trichlorodisiloxane	Fusion Vaporization	c liq	liq g	760	173. 316.2	-100. 43.0	7.3	23.1	
$SiHFC1_2$	fluorodichlorosilane	Fusion Vaporization	c liq	liq g	760	124. 254.5	-149. -18.7	4.8	18.9	

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Table 24-6. Silicon (at. no., 14; at. wt., 28.06)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
				Initial	Final		°K	°C			
$\text{SiBr}_4$		tetrabromosilane	Fusion Vaporization	c liq	liq g	760	278.4 426.0	5.2 152.8	9.1	21.4	
$\text{Si}_2\text{Br}_6$		hexabromodisilane	Fusion Vaporization	c liq	liq g	760	368. 538.	95. 265.			
$\text{Si}_3\text{Br}_8$		octabromotrisilane	Fusion	c	liq		406.	133.			
$\text{SiH}_3\text{Br}$		bromosilane	Fusion Vaporization	c liq	liq g	760	179. 275.1	-94. 1.9	5.8	21.1	
$\text{SiH}_2\text{Br}_2$		di-bromosilane	Fusion Vaporization	c liq	liq g	760	203. 286.8	-70. 13.6	7.5	26.2	
$\text{SiHBr}_3$		tribromosilane	Fusion Vaporization	c liq	liq g	760	200. 385.4	-73. 112.2	8.1	21.0	
$\text{Si}_2\text{H}_5\text{Br}$		bromodisilane	Fusion Vaporization	c liq	liq g	10	173. 243.	-100. -30.	7.	29.	

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Table 24-7. Silicon (at. no., 14; at. wt., 28.06)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Formula	Substance Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
			Initial	Final		°K	°C			
SiF <sub>3</sub> Br	trifluorobromosilane	Fusion Vaporization	c liq	liq g	760	202.7 231.6	-70.5 -41.6			
SiF <sub>2</sub> Br <sub>2</sub>	difluorodibromosilane	Fusion Vaporization	c liq	liq g	760	206.3 287.1	-66.9 13.9	4.2	18.	
SiFBr <sub>3</sub>	fluorotribromosilane	Fusion Vaporization	c liq	liq g	760	190.7 357.1	-82.5 83.9	5.9	20.6	
SiCl <sub>3</sub> Br	trichlorobromosilane	Vaporization	liq	g	760	353.	80.	7.8	21.8	
SiCl <sub>2</sub> Br <sub>2</sub>	dichlorodibromosilane	Vaporization	liq	g	760	377.	104.			
SiClBr <sub>3</sub>	chlorotribromosilane	Fusion Vaporization	c liq	liq g	760	234. 400.	-39. 127.			
SiFCl <sub>2</sub> Br	fluorodichlorobromosilane	Fusion Vaporization	c liq	liq g	760	160.9 308.5	-112.3 35.3	6.0	19.4	
SiFClBr <sub>2</sub>	fluorochlorodibromosilane	Fusion Vaporization	c liq	liq g	760	173.9 332.6	-99.3 59.4	7.1	21.4	

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Table 24-8. Silicon (at. no., 14; at. wt., 28.06)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		°K	°C			
$\text{SiI}_4$		tetraiodosilane	Fusion Vaporization	c liq	liq g	760	393. 563.	120. 290.			
$\text{Si}_2\text{I}_6$		hexaiododisilane	Fusion	c	liq		523.	250.			
$\text{SiH}_3\text{I}$		iodosilane	Fusion Vaporization	c liq	liq g	760	216. 318.8	-57. 45.6	6.9	21.6	
$\text{SiH}_2\text{I}_2$		diiodosilane	Fusion Vaporization	c liq	liq g	760	272. 423.	-1. 150.	8.8	20.8	
$\text{SiHI}_3$		triiodosilane	Fusion Vaporization	c liq	liq g	45	281. 400.	8. 127.	15.	38.	
$\text{SiCl}_3\text{I}$		trichloriodosilane	Vaporization	liq	g	760	386.	113.			
$\text{SiCl}_2\text{I}_2$		dichlorodiiodosilane	Vaporization	liq	g	760	445.	172.			
$\text{SiClI}_3$		chlorotriiodosilane	Fusion Vaporization	c liq	liq g	760	275. 508.	2. 235.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$SiBr_3I$	tribromiodosilane	Fusion Vaporization	c liq	liq g	760	287. 465.	14. 192.			
$SiBr_2I_2$	dibromodiodosilane	Fusion Vaporization	c liq	liq g	760	311. 503.	38. 230.			
$SiBrI_3$	bromotriiodosilane	Fusion Vaporization	c liq	liq g	760	326. 528.	53. 255.			
$SiS_2$		Fusion	c	liq		1363.	1090.			
$SiSCl_2$		Fusion	c	liq		348.	75.			
$SiSBr_2$		Fusion	c	liq		366.	93.			
$Si_3H_9N$	trisilylamine	Fusion Vaporization	c liq	liq g	100	167.6 271.4	-105.6 -1.8	6.8	25.1	
$3Bi_2O_3 \cdot SiO_2$		Fusion	c	liq		1095.	822.			
$3Bi_2O_3 \cdot 2SiO_2$		Fusion	c	liq		1064.	791.			
$Bi_2O_3 \cdot 6SiO_2$		Fusion	c	liq		1265.	992.			

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Table 24-10. Silicon (at. no., 14; at. wt., 28.06)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		°K	°C			
$\text{SiH}_6\text{C}$		methylsilane	Fusion Vaporization	c liq	liq g	760	116.8 216.3	-156.4 -56.9	4.2	19.	
$\text{SiH}_8\text{C}_2$		ethylsilane	Vaporization	liq	g	760	259.	-14.	5.4	20.9	
$\text{SiH}_8\text{C}_2$		dimethylsilane	Fusion Vaporization	c liq	liq g	760	123.4 253.0	-149.8 -20.2	4.9	19.4	
$\text{SiH}_{10}\text{C}_3$		n-propylsilane	Vaporization	liq	g	760	296.	23.	6.4	22.	
$\text{SiH}_{10}\text{C}_3$		trimethylsilane	Vaporization	liq	g	760	283.	10.			
$\text{SiH}_{12}\text{C}_4$		diethylsilane	Vaporization	liq	g	760	327.	54.	7.4	23.	
$\text{SiH}_{12}\text{C}_4$		tetramethylsilane	Fusion Fusion Vaporization	c, I c, II liq	liq liq g	760	171.04 174.12 299.80	-102.12 -99.04 26.64	1.427 1.648 5.79	8.34 9.46 19.31	5.6
$\text{SiH}_{14}\text{C}_5$		trimethylsilylsilane	Vaporization	liq	g	760	335.	62.	7.2	21.5	
$\text{SiH}_{16}\text{C}_7$		methyltriethylsilane	Vaporization	liq	g	760	400.	127.	8.6	21.5	
$\text{SiH}_{20}\text{C}_8$		tetraethylsilane	Vaporization	liq	g	760	426.	153.	9.5	22.3	
$\text{Si}_2\text{H}_8\text{C}_6$		hexamethyldisilane	Fusion Vaporization	c liq	liq g	100	287.4 325.2	14.2 52.0	10.7	32.9	



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Table 24-11. Silicon (at. no., 14; at. wt., 28.06)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		O <sub>K</sub>	O <sub>C</sub>			
SiH <sub>10</sub> O <sub>3</sub>	trimethylsilanol	Vaporization	liq	g	752	371.8	98.6			
SiH <sub>14</sub> O <sub>5</sub>	ethoxytrimethylsilane	Vaporization	liq	g	760	349.	76.	8.	23.	
SiH <sub>16</sub> O <sub>2</sub> C <sub>6</sub>	diethoxydimethylsilane	Vaporization	liq	g	760	387.	114.	9.	23.	
SiH <sub>18</sub> O <sub>3</sub> C <sub>7</sub>	triethoxymethylsilane	Vaporization	liq	g	760	417.	144.	10.	24.	
SiH <sub>20</sub> O <sub>4</sub> C <sub>8</sub>	tetraethoxysilane	Fusion Vaporization	c liq	liq g	760	190.7 441.7	-82.5 168.5	2.1 10.19	11. 23.07	
Si <sub>2</sub> H <sub>18</sub> O <sub>6</sub>	hexamethyl disiloxane	Vaporization	liq	g	760	373.	100.	8.	21.	
Si <sub>2</sub> H <sub>22</sub> O <sub>3</sub> C <sub>8</sub>	1,3-diethoxytetramethyldisiloxane	Vaporization	liq	g	760	444.	161.	11.	25.	
Si <sub>2</sub> H <sub>30</sub> O <sub>6</sub> C <sub>12</sub>	hexaethoxydisilane	Vaporization	liq	g	29	410.	137.	12.	29.	
Si <sub>3</sub> H <sub>24</sub> O <sub>2</sub> C <sub>8</sub>	octamethyl trisiloxane	Vaporization	liq	g	760	425.	152.	10.	23.	
Si <sub>3</sub> H <sub>28</sub> O <sub>4</sub> C <sub>10</sub>	1,5-diethoxyhexamethyl trisiloxane	Vaporization	liq	g	760	470.	197.	12.	26.	
Si <sub>4</sub> H <sub>24</sub> O <sub>4</sub> C <sub>8</sub>	octamethyl cyclotetrasiloxane	Vaporization	liq	g	760	449.	176.	11.	24.	
Si <sub>4</sub> H <sub>30</sub> O <sub>3</sub> C <sub>10</sub>	decamethyl tetrasiloxane	Vaporization	liq	g	760	467.	194.	12.	26.	

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Table 24-12. Silicon (at. no., 14; at. wt., 28.06)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
$SiH_3F_3C$		methyltrifluorosilane	Fusion Vaporization	c liq	liq g	760	200.4 243.0	-72.8 -30.2	5.7	23.5	
$SiH_6F_2C_2$		dimethyldifluorosilane	Fusion Vaporization	c liq	liq g	760	185.7 275.9	-87.5 2.7	6.12	22.18	
$SiH_5F_3C_2$		ethyltrifluorosilane	Fusion Vaporization	c liq	liq g	760	159.9 269.0	-113.3 -4.2	6.1	22.7	
$SiH_9FC_3$		trimethylfluorosilane	Fusion Vaporization	c liq	liq g	760	198.9 289.6	-74.3 16.4	6.4	22.1	
$SiH_9F_2C_4$		diethyldifluorosilane	Fusion Vaporization	c liq	liq g	760	194.5 334.1	-78.7 60.9	7.2	21.6	
$SiH_5ClC$		methylchlorosilane	Fusion Vaporization	c liq	liq g	250	139. 254.6	-134. -18.6	5.8	22.8	
$SiH_4Cl_2C$		methyldichlorosilane	Fusion Vaporization	c liq	liq g	100	180. 266.8	-93. -6.4	6.8	25.5	
$SiH_3Cl_3C$		methyltrichlorosilane	Fusion Vaporization	c liq	liq g	760	195.4 338.9	-77.8 65.7	7.4	21.8	

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Table 24-13. Silicon (at. no., 14; at. wt., 28.06) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION September 30, 1948										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}\text{K}$	$^{\circ}\text{C}$			
$\text{SiH}_6\text{Cl}_2\text{C}_2$	dimethyldichlorosilane	Fusion Vaporization	c liq	liq g	760	197. 343.2	-76. 70.0	7.5	21.9	
$\text{SiH}_5\text{Cl}_3\text{C}_2$	ethyltrichlorosilane	Fusion Vaporization	c liq	liq g	760	168. 371.1	-105. 97.9	7.7	20.7	
$\text{SiH}_3\text{ClC}_3$	trimethylchlorosilane	Fusion Vaporization	c liq	liq g	760	215.5 331.	-57.7 58.	7.	21.	
$\text{SiH}_{10}\text{Cl}_2\text{C}_4$	diethyldichlorosilane	Vaporization	liq	g	760	403.	130.	10.	25.	
$\text{SiH}_{15}\text{ClC}_6$	triethylchlorosilane	Vaporization	liq	g	760	419.	146.	9.	22.	
$\text{SiH}_5\text{OCl}_3\text{C}_2$	ethoxytrichlorosilane	Vaporization	liq	g	760	375.	102.	8.	21.	
$\text{SiH}_3\text{OCl}_2\text{C}_3$	ethoxymethyldichlorosilane	Vaporization	liq	g	760	373.	100.	8.	21.	
$\text{Si}_2\text{H}_{12}\text{OCl}_2\text{C}_4$	tetramethyl-1,3-dichlorodisiloxane	Fusion Vaporization	c liq	liq g	760	236. 411.	-37. 138.	9.	22.	

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Table 24-14. Silicon (at. no., 14; at. wt., 28.06)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		$^{\circ}K$	$^{\circ}C$			
$SiH_3FCl_2C$	methylfluorodichlorosilane	Fusion Vaporization	c liq	liq g	760	174.5 302.7	-98.7 29.5	6.7	22.1	
$SiH_6FCl_2C_2$	dimethylfluorochlorosilane	Fusion Vaporization	c liq	g	760	188. 309.6	-85. 36.4	6.6	21.3	
$SiH_5F_2ClC_2$	ethylfluorochlorosilane	Vaporization	liq	g	760	300.3	27.1	6.7	22.3	
$SiH_5FCl_2C_2$	ethylfluorodichlorosilane	Vaporization	liq	g	760	335.3	62.1	7.4	22.1	
$SiO_4N_4C_4$	silicon tetracyanate	Fusion Vaporization	c liq	liq g	760	307.9 520.	34.5 247.	16.5	31.7	
$SiO_4N_4C_4$	silicon tetraisocyanate	Fusion Vaporization	c liq	liq g	760	299.2 459.	26.0 185.	12.9	28.1	
$Si_2H_9NC$	2-methyldisilazane	Vaporization	liq	g	760	305.5	32.3	6.9	22.6	
$Si_2H_{11}NC_2$	2-ethyldisilazane	Fusion Vaporization	c liq	liq g	68	146. 273.	-127. 0.	8.0	29.3	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$SiH_3ONC_4$	trimethylsilicon isocyanate	Fusion Vaporization	c liq	liq g	760	224.2 364.1	-49.0 90.9	8.3	22.8	
$SiH_6O_2N_2C_4$	dimethylsilicon diisocyanate	Fusion Vaporization	c liq	liq g	760	242.0 412.3	-31.2 139.1	9.9	24.0	
$SiH_3O_2N_3C_4$	methylsilicon triisocyanate	Fusion Vaporization	c liq	liq g	760	275.9 444.0	2.7 170.8	11.2	25.2	
$SiH_5O_4N_3C_5$	ethoxysilicon triisocyanate	Vaporization	liq	g	760	452.7	179.5	11.5	25.4	
$SiH_7O_4N_2C_6$	diethoxysilicon diisocyanate	Vaporization	liq	g	760	448.5	175.3	11.0	24.5	
$SiH_{15}O_4NC_7$	triethoxysilicon isocyanate	Vaporization	liq	g	760	446.0	172.8	10.9	24.4	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 24-16. Silicon (at. no., 14; at. wt., 28.06)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		°K	°C			
$\text{SiOF}_3\text{NC}$		trifluorosilicon isocyanate	Vaporization	liq	g	760	267.	-6.			
$\text{SiO}_2\text{F}_2\text{N}_2\text{C}_2$		difluorosilicon diisocyanate	Fusion Vaporization	c liq	liq g	760	198. 341.8	-75. 68.6		24.9	
$\text{SiO}_3\text{FN}_3\text{C}_3$		fluorosilicon triisocyanate	Fusion Vaporization	c liq	liq g	760	244.0 407.	-29.2 134.		26.	
$\text{Si}_3\text{N}_4\text{C}_4$		silicon tetraisoisocyanate	Fusion Vaporization	c liq	liq g	760	417. 586.	144. 313.	15.	26.	
$\text{SiH}_9\text{SiNC}_4$		trimethylsilicon isothiocyanate	Fusion Vaporization	c liq	liq g	760	240.4 416.2	-32.8 143.0	9.3	22.3	
$\text{SiH}_6\text{S}_2\text{N}_2\text{C}_4$		dimethylsilicon diisothiocyanate	Fusion Vaporization	c liq	liq g	760	291.2 490.	18.0 217.	12.2	24.9	
$\text{SiH}_3\text{S}_3\text{N}_3\text{C}_4$		methylsilicon triisothiocyanate	Fusion Vaporization	c liq	liq g	760	345.6 540.	72.4 267.	13.7	25.4	
$\text{SiH}_9\text{O}_3\text{SiNC}_4$		trimethoxysilicon isothiocyanate	Vaporization	liq	g	760	443.7	170.5	13.4	30.2	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 25-1. Germanium (at. no., 32; at. wt., 72.60)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
				Initial	Final		$0^\circ K$	$0^\circ C$			
Ge			Fusion	c	liq		1233.	960.			
GeO			Sublimation	c	g	760	983.	710.	8.3	6.7	
GeO <sub>2</sub>			Transition	c, II, insoluble	c, I, soluble		1306.	1033.			
			Fusion	c, II, insoluble	liq		1359.	1086.			
			Fusion	c, I, soluble	liq		1389.	1116.			
GeH <sub>4</sub>			Transition	c, III	c, II		73.2	-200.0	0.050	0.68	
			Transition	c, II	c, I		76.5	-196.7	0.086	1.12	
			Fusion	c, I	liq		107.26	-165.90	0.200	1.86	2.0
			Vaporization	liq	g	760	184.80	-88.36	3.361	18.19	
Ge <sub>2</sub> H <sub>6</sub>			Fusion	c	liq		164.	-109.			
			Vaporization	liq	g	760	304.0	30.8	6.0	19.8	
Ge <sub>3</sub> H <sub>8</sub>			Fusion	c	liq		168.	-105.			
			Vaporization	liq	g	760	384.3	111.1	7.7	20.0	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 25-2. Germanium (at. no., 32; at. wt., 72.60)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta f$ kcal/mole	$\Delta s$ cal/deg mole	$\Delta c_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
GeF <sub>4</sub>		Sublimation	c	g	760	236.4	-36.8	7.8	33.0	
		Fusion	c	liq	3032	258.2	-15.0			
GeCl <sub>4</sub>		Fusion	c	liq		223.7	-49.5			
		Vaporization	liq	g	760	356.3	83.1	7.9	22.2	
GeHCl <sub>3</sub>		Fusion	c	liq		202.	-71.			
		Vaporization	liq	g	760	348.5	75.3	8.2	23.5	
GeF <sub>3</sub> Cl		Fusion	c	liq		207.	-66.			
		Vaporization	liq	g	760	253.	-20.			
GeF <sub>2</sub> Cl <sub>2</sub>		Fusion	c	liq		221.	-52.			
		Vaporization	liq	g	760	269.	-4.	5.9	21.9	
GeFCl <sub>3</sub>		Fusion	c	liq		223.	-50.			
		Vaporization	liq	g	760	310.4	37.2	6.6	21.3	
GeBr <sub>4</sub>		Fusion	c	liq		299.3	26.1			
		Vaporization	liq	g	760	460.3	187.1	9.9	21.5	
GeI <sub>4</sub>		Fusion	c	liq		417.	144.			



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 25-3. Germanium (at. no., 32; at. wt., 72.60)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
GeS		Fusion	c, I, black	liq		898.	625.			
GeSe		Fusion	c	liq		940.	667.			
GeSe <sub>2</sub>		Fusion	c	liq		980.	707.			
GeH <sub>12</sub> Cl <sub>4</sub>	tetramethylgermane	Fusion	c	liq		185.	-88.			
		Vaporization	liq	g	760	316.5	43.3	6.7	21.2	
GeH <sub>20</sub> Cl <sub>8</sub>	tetraethylgermane	Fusion	c	liq		183.	-90.			
		Vaporization	liq	g	760	436.7	163.5			
GeH <sub>6</sub> OCl <sub>2</sub>	dimethylgermanium oxide	Fusion	c	liq		406.6	133.4			
		Vaporization	liq	g	760	484.	211.			
GeH <sub>5</sub> F <sub>3</sub> Cl <sub>2</sub>	ethylgermanium trifluoride	Fusion	c	liq		257.	-16.			
		Vaporization	liq	g	750	385.	112.			
GeH <sub>3</sub> Cl <sub>3</sub> Cl <sub>3</sub>	methylgermanium trichloride	Vaporization	liq	g	760	384.	111.			
GeH <sub>6</sub> Cl <sub>2</sub> Cl <sub>2</sub>	dimethylgermanium dichloride	Fusion	c	liq		251.	-22.			
		Vaporization	liq	g	760	397.	124.			
GeH <sub>5</sub> Cl <sub>3</sub> Cl <sub>2</sub>	ethylgermanium trichloride	Vaporization	liq	g	760	417.	144.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 25-4. Germanium (at. no., 32; at. wt., 72.60)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		$^{\circ}K$	$^{\circ}C$			
$GeH_5Br_3C_2$	ethylgermanium tribromide	Vaporization	liq	g	760	473.	200.			
$GeH_5I_3C_2$	ethylgermanium triiodide	Fusion	c	liq		271.	-2.			
		Vaporization	liq	g	760	555.	282.			
$GeH_6SC_2$	dimethylgermanium sulfide	Fusion	c	liq		328.7	55.5			
		Vaporization	liq	g	760	575.	302.			
$GeO_4N_4C_4$	germanium tetraisocyanate	Fusion	c	liq		265.	-8.			
		Vaporization	liq	g	760	477.	204.	12.6	26.4	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 26-1. Tin (at. no., 50; at. wt., 118.70)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Sn		Transition	c, III, gray	c, II, white	760	291.	18.	0.6	2.1	
		Transition	c, II, white	c, I		476.0	202.8	0.002	0.004	
		Fusion	c, I	liq		505.1	231.9	1.69	3.35	
		Vaporization	liq	g		2600.	2337.			
SnO <sub>2</sub>		Transition	c, II	c, I		683.	410.	0.45	0.66	
SnH <sub>4</sub>	stannane	Fusion	c	liq	760	123.	-150.			
		Vaporization	liq	g		221.4	-51.8	4.4	19.9	
		Vaporization	liq	g	760	978.	705.			
SnF <sub>4</sub>										
SnCl <sub>2</sub>		Fusion	c	liq	760	520.	247.	3.0	5.8	
		Vaporization	liq	g		896.	623.	21.	23.	
SnCl <sub>4</sub>		Fusion	c	liq	760	239.9	-33.3	2.19	9.13	
		Vaporization	liq	g		386.	113.	8.3	21.5	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 26-2. Tin (at. no., 50; at. wt., 118.70)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$SnBr_2$		Fusion Vaporization	c liq	liq g	760	505. 911.	232. 638.	1.7 22.	3.4 24.	
$SnBr_4$		Transition Fusion Vaporization	c, II c, I liq	c, I liq g	760	267. 303. 478.	-6. 30. 205.	3.0 10.	9.9 21.	
$SnCl_3Br$		Fusion	c	liq		242.	-31.			
$SnCl_2Br_2$		Fusion	c	liq		253.	-20.			
$SnClBr_3$		Fusion	c	liq		274.	1.			
$SnI_2$		Fusion Vaporization	c liq	liq g	760	593. 991.	320. 718.	24.	24.	
$SnI_4$		Fusion Vaporization	c liq	liq g	760	417.7 617.	144.5 344.	4.48 13.6	10.73 22.	5.7
$SnS$		Fusion	c	liq		1153.	880.			
$SnSe$		Fusion	c	liq		1133.	860.			
$SnTe$		Fusion	c	liq		1073.	800.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

SERIES II

Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$SnH_6C$	methyl stannane	Vaporization	liq	g	760	273.	0.	5.7	20.9	
$SnH_8C_2$	dimethyl stannane	Vaporization	liq	g	760	308.	35.	6.8	22.1	
$SnH_{10}C_3$	trimethyl stannane	Vaporization	liq	g	760	332.	59.	7.2	21.6	
$SnH_{12}C_4$	tetramethyl stannane	Vaporization	liq	g	760	351.2	78.0	7.4	21.1	
$SnH_{14}C_5$	ethyltrimethyl stannane	Vaporization	liq	g	760	381.4	108.2	8.4	22.0	
$SnH_{16}C_6$	n-propyltrimethyl stannane	Vaporization	liq	g	760	404.0	130.8	8.8	21.8	
$SnH_{20}C_8$	tetraethyl stannane	Fusion Vaporization	c liq	liq g	760	161. 448.	-112. 175.			
$SnH_3Cl_3C$	methyltin trichloride	Fusion	c	liq		300.	27.			
$SnH_6Cl_2C_2$	dimethyltin dichloride	Fusion	c	liq		380.	107.			
$SnH_9ClC_3$	trimethyltin chloride	Fusion	c	liq		316.	43.			
$SnH_{15}ClC_6$	triethyltin chloride	Fusion Vaporization	c liq	liq g		289. 367.	16. 94.			
					13					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 26-4. Tin (at. no., 50; at. wt., 118.70)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$SnH_3Br_3C$	methyltin tribromide	Fusion	c	liq		326.	53.			
$SnH_6Br_2C_2$	dimethyltin dibromide	Fusion	c	liq		349.	76.			
$SnH_9BrC_3$	trimethyltin bromide	Fusion	c	liq		300.	27.			
$SnH_{15}BrC_6$	triethyltin bromide	Fusion	c	liq		260.	-13.			
		Vaporization	liq	g	13	370.	97.			
$SnH_3I_3C$	methyltin triiodide	Fusion	c	liq		357.	84.			
$SnH_6I_2C_2$	dimethyltin diiodide	Fusion	c	liq		315.	42.			
		Fusion	c	liq		276.6	3.4			
$SnH_9IC_3$	trimethyltin iodide	Vaporization	liq	g	13	343.	70.			
		Fusion	c	liq		239.	-34.			
$SnH_{15}IC_6$	triethyltin iodide	Vaporization	liq	g	13	390.	117.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 27-1. Lead (at. no., 82; at. wt., 207.21)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		°K	°C			
Pb			Fusion Vaporization	c liq	liq g	760	600.6 2023.	327.4 1750.	1.22 43.0	2.03 21.3	
PbO			Transition Fusion Vaporization	c, II, red c, I, yellow liq	c, I, yellow liq g	760	762. 1159. 1745.	489. 886. 1472.	2.8 51.	2.4 29.	
PbF <sub>2</sub>			Fusion Vaporization	c, I liq	liq g	760	1095. 1563.	822. 1290.	1.8 38.3	1.6 24.5	
PbF <sub>4</sub>			Fusion	c	liq		873.	600.			
PbCl <sub>2</sub>			Fusion Vaporization	c liq	liq g	760	771. 1227.	498. 954.	5.7 29.6	7.4 24.1	
2PbO·PbCl <sub>2</sub>			Fusion	c	liq		966.	693.			
4PbO·PbCl <sub>2</sub>			Fusion	c	liq		984.	711.			
PbFCI			Fusion	c	liq		874.	601.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 27-2. Lead (at. no., 82; at. wt., 207.21)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		$^{\circ}K$	$^{\circ}C$			
PbBr <sub>2</sub>		Fusion Vaporization	c liq	liq g	760	643.2 1187.	370.0 914.	5. 27.7	7.8 23.3	
2PbO·PbBr <sub>2</sub>		Fusion	c	liq		985.	712.			
PbBrF		Fusion	c	liq		834.	561.			
PbI <sub>2</sub>		Fusion Vaporization	c liq	liq g	760	695. 1145.	412. 872.	5.2 24.8	7.6 21.7	
PbS		Sublimation Fusion	c c	g liq	10	1238. 1387.	965. 1114.	55. 4.2	44. 3.0	
PbSO <sub>4</sub>		Transition Fusion	c, II c, I	c, I liq		1139. 1360.	866. 1087.	4.06 9.6	3.56 7.1	
PbO·PbSO <sub>4</sub>		Fusion	c	liq		1243.	970.			
2PbO·PbSO <sub>4</sub>		Fusion	c	liq		1223.	950.			
PbSe		Fusion	c	liq		1338.	1065.			
PbTe		Fusion	c	liq		1178.	905.			



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards Washington, D.C.

Table 27-3. Lead (at. no., 82; at. wt., 207.21)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Pb(Pb_3)_2$		Fusion	c	liq		1073.	800.			
$Pb_3(Pb_4)_2$		Fusion	c	liq		1287.	1014.			
$PbO \cdot Pb_3(Pb_4)_2$		Fusion	c	liq		1253.	980.			
$5PbO \cdot Pb_3(Pb_4)_2$		Fusion	c	liq		1133.	860.			
$PbF_2 \cdot 3Pb_3(Pb_4)_2$		Fusion	c	liq		1373.	1100.			
$PbCl_2 \cdot 3Pb_3(Pb_4)_2$		Fusion	c	liq		1429.	1156.			
$Pb_3(AsO_4)_2$		Fusion	c	liq		1315.	1042.			
$5PbO \cdot Pb_3(AsO_4)_2$		Fusion	c	liq		1135.	862.			
$PbF_2 \cdot 3Pb_3(AsO_4)_2$		Fusion	c	liq		1315.	1042.			
$PbCl_2 \cdot 3Pb_3(AsO_4)_2$		Fusion	c	liq		1413.	1140.			
$PbO \cdot Sb_2O_3$		Fusion	c	liq		831.	558.			
$2PbO \cdot Bi_2O_3$		Fusion	c	liq		898.	625.			
$2PbO \cdot 3Bi_2O_3$		Fusion	c	liq		959.	686.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 27-4. Lead (at. no., 82; at. wt., 207.21)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$PbH_{12}C_4$	tetramethyllead	Fusion Vaporization	c liq	liq g	760	245.7 379.3	-27.5 106.1	8.0	21.1	
$PbH_{14}C_5$	ethyltrimethyllead	Vaporization	liq	g	760	407.8	134.6	9.1	22.3	
$PbH_{16}C_6$	diethyldimethyllead	Vaporization	liq	g	760	427.7	154.5	10.2	23.8	
$PbH_{18}C_7$	triethylmethyllead	Vaporization	liq	g	760	443.6	170.4	11.2	25.2	
$PbH_{20}C_8$	tetraethyllead	Fusion Vaporization	c liq	liq g	609.7	137. 450.0	-136. 176.8	11.87	26.38	
$PbO \cdot SiO_2$		Fusion	c	liq		1037.	764.			
$2PbO \cdot SiO_2$		Fusion	c	liq		1016.	743.			
$4PbO \cdot SiO_2$		Transition Transition	c, III c, II	c, II c, I		413. 993.	140. 720.			
$3PbS \cdot 2SiS_2$		Fusion	c	liq		1089.	816.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Ga		Transition	c, I	c, II	8.86×10 <sup>6</sup>	275.6	2.4	0.51	1.85	0.38
		Fusion	c, I	liq		275.6	2.4	1.23	4.46	
		Fusion	c, II	liq		275.6	2.4	0.72	2.61	
		Fusion	c, I	liq		309.940	29.780	1.336	4.31	
		Vaporization	liq	g		1210.	937.	63.8	52.7	
Ga <sub>2</sub> O <sub>3</sub> GaF <sub>3</sub>		Fusion	c	liq	760	2013.	1740.			
		Vaporization	liq	g		1223.	950.			
		Fusion	c	liq		443.7	170.5			
		Sublimation	c	g		350.7	77.5	17.4	49.6	
		Fusion	c	liq		350.7	77.5	5.1	14.5	
GaCl <sub>2</sub> Ga <sub>2</sub> Cl <sub>6</sub>		Vaporization	liq	g, equilibrium	760	473.	200.	11.0	23.3	
		Fusion	c	liq		397.7	124.5			
		Vaporization	liq	g, equilibrium		551.	278.	18.6	33.8	
		Fusion	c	liq		283.	10.			
		Vaporization	liq	g, equilibrium		619.	346.	27.	44.	
Ga <sub>2</sub> Br <sub>6</sub> Ga <sub>2</sub> I <sub>6</sub>		Fusion	c	liq	760	397.7	124.5			
		Vaporization	liq	g, equilibrium		551.	278.	18.6	33.8	
		Fusion	c	liq		283.	10.			
		Vaporization	liq	g, equilibrium		619.	346.	27.	44.	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 28-2. Gallium (at. no., 31; at. wt., 69.72)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1948 (Corrected)

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
GaS		Fusion	c	liq		1218.	965.			
Ga <sub>2</sub> S <sub>3</sub>		Fusion	c	liq		1523.	1250.			
GaSe		Fusion	c	liq		1213.	960.			
GaTe		Fusion	c	liq		1097.	824.			
Ga <sub>2</sub> Te <sub>3</sub>		Fusion	c	liq		1063.	790.			
GaCl <sub>3</sub> ·NH <sub>3</sub>		Fusion	c	liq		397.	124.			
GaBr <sub>3</sub> ·NH <sub>3</sub>		Fusion	c	liq		397.	124.			
GaI <sub>3</sub> ·NH <sub>3</sub>		Fusion	c	liq		413.	140.			
Ga(CH <sub>3</sub> ) <sub>3</sub>		Fusion	c	liq		257.5	-15.7			
		Vaporization	liq	g	760	328.8	55.6	7.8	23.7	
Ga(C <sub>2</sub> H <sub>5</sub> ) <sub>3</sub>		Fusion	c	liq		189.9	-82.3			
		Vaporization	liq	g	760	415.8	142.6	9.9	23.8	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$(CH_3)_3Ga \cdot NH_3$		Fusion	c	liq		304.	31.			
$(CH_3)_3Ga \cdot N(CH_3)_3$		Fusion	c	liq	89.74	369.4	96.2	2.3	6.2	
		Vaporization	liq	g	89.74	369.4	96.2	10.3	27.9	
$(CH_3)_3Ga \cdot N(C_2H_5)_3$		Sublimation	c	g	40.7	369.	96.	14.2	38.5	
		Fusion	c	liq	40.7	369.	96.	0.9	2.4	
		Vaporization	liq	g	760	440.	167.	13.3	30.2	
$Ga(CH_3)_2Cl \cdot NH_3$		Fusion	c	liq		327.	54.			
$Ga(CH_3)_2Cl \cdot 2NH_3$		Fusion	c	liq		385.	112.			

Table 28-3. Gallium (at. no., 31; at. wt., 69.72)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1948

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 29-1. Indium (at. no., 49; at. wt., 114.76)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
In		Fusion	c	liq		430.	157.	0.78	1.81	
		Vaporization	liq	g	0.007	1200.	927.	55.7	46.4	
InF <sub>3</sub>		Fusion	c	liq		1443.	1170.			
InCl		Transition	c, II, yellow	c, I, red		393.	120.			
		Fusion	c, I, red	liq		498.	225.			
		Vaporization	liq	g	760	882.	609.	21.5	24.4	
InCl <sub>2</sub>		Fusion	c	liq		508.	235.			
		Vaporization	liq	g	760	758.	485.	46.0	60.7	
InCl <sub>3</sub>		Sublimation	c	g	100	715.	442.	38.	53.	
		Fusion	c	liq	10000	859.	586.			
InBr		Fusion	c	liq		493.	220.			
		Vaporization	liq	g	760	929.	656.	22.0	23.7	
InBr <sub>2</sub>		Fusion	c	liq		513.	240.			
		Vaporization	liq	g	760	911.	638.	19.7	21.6	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
InBr <sub>3</sub>		Sublimation	c	g	760	645.	372.	25.1	38.9	
		Fusion	c	liq	4900	709.	436.			
InI		Fusion	c	liq		624.	351.			
InI <sub>3</sub>		Vaporization	liq	g	760	985.	712.	21.7	22.0	
		Fusion	c	liq		483.	210.			
InS		Fusion	c	liq		965.	692.			
In <sub>2</sub> S		Fusion	c	liq		926.	653.			
In <sub>2</sub> S <sub>3</sub>		Fusion	c	liq		1323.	1050.			
InSe		Fusion	c	liq		933.	660.			
In <sub>2</sub> Se <sub>3</sub>		Fusion	c	liq		1163.	890.			
InTe		Fusion	c	liq		969.	696.			
In <sub>2</sub> Te <sub>3</sub>		Fusion	c	liq		940.	667.			
In(CH <sub>3</sub> ) <sub>3</sub>		Sublimation	c	g	153	361.6	88.4	13.8	30.2	
		Fusion	c	liq	153	361.6	88.4	3.8	8.3	
		Vaporization	liq	g	760	409.0	135.8	10.0	24.4	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Tl		Transition	c, I	c, III	$2.87 \times 10^7$	426.	153.	-0.19	-0.45	
		Transition	c, II	c, III	$2.87 \times 10^7$	426.	153.	0.01	0.02	
		Transition	c, II	c, I	$2.87 \times 10^7$	426.	153.	0.20	0.47	
		Transition	c, II	c, I		508.3	235.1	0.082	0.16	0.18
		Fusion	c, I	liq		576.8	303.6	1.03	1.79	0.43
		Vaporization	liq	g	28	1350.	1077.	40.1	29.7	
Tl <sub>2</sub> O		Vaporization	liq	g	760	1730.	1457.			
		Fusion	c	liq		573.	300.			
		Vaporization	liq	g	760	773.	500.			
Tl <sub>2</sub> O <sub>3</sub>		Fusion	c	liq		990.	717.			
		Fusion	c	liq		600.	327.			
TlF		Vaporization	liq	g	760	928.	655.			
TlF <sub>3</sub>		Fusion	c	liq		823.	550.			



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 30-2. Thallium (at. no., 81; at. wt., 204.39) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION December 31, 1948											
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole	
Formula	Description		Initial	Final		°K	°C				
TlCl		Sublimation	c	g	1.12	702.	429.	29.	41.		
		Fusion	c	liq	1.12	702.	429.	4.0	5.6		
		Vaporization	liq	g	760	1079.	806.	24.8	23.0		
		Fusion	c	liq		308.	35.				
TlCl <sub>3</sub> ·4H <sub>2</sub> O	TlClO <sub>4</sub>	Transition	c, II	c, I		539.	266.				
		Fusion	c, I	liq		774.	501.				
TlBr		Sublimation	c	g	2.11	732.	459.	31.4	42.9		
		Fusion	c	liq	2.11	732.	459.	3.6	4.9		
		Vaporization	liq	g	760	1089.	816.	24.6	22.7		
TlI		Transition	c, II, yellow	c, I, red		438.	165.				
		Sublimation	c, I, red	g	1.00	713.	440.	30.0	42.1		
		Fusion	c, I, red	liq	1.00	713.	440.	2.7	3.8		
		Vaporization	liq	g	760	1098.	825.	24.9	22.6		

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 30-3. Thallium (at. no., 81; at. wt., 204.39)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1948 (Corrected)

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Tl_2S$			Fusion	c	liq		721.	448.	3.0	4.2	
$Tl_2S_3$			Fusion	c	liq		533.	260.			
$Tl_2S_5$			Fusion	c	liq		583.	310.			
$Tl_2SO_4$			Fusion	c	liq		905.	632.	5.5	6.1	
$Tl_2Se$			Fusion	c	liq		671.	398.			
$Tl_2Se_3$			Transition	c, II	c, I		438.	165.			
$Tl_2Te_3$			Fusion	c	liq		701.	428.			
$TlNO_2$			Fusion	c	liq		455.	182.			
$TlNO_3$			Transition	c, III	c, II		334.2	61.0	0.24	0.72	
			Transition	c, II	c, I		416.7	143.5	0.76	1.82	
			Fusion	c, I	liq		479.7	206.5	2.29	4.77	
			Vaporization	liq	g	760	706.	433.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 30-4. Thallium (at. no., 81; at. wt., 204.39)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1948

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
			Initial	Final		$^{\circ}K$	$^{\circ}C$			
TlPO <sub>3</sub>		Fusion	c	liq		678.	405.			
TlH <sub>2</sub> PO <sub>2</sub>		Fusion	c	liq		383.	110.			
TlH <sub>2</sub> PO <sub>3</sub>		Fusion	c	liq		343.	70.			
TlH <sub>2</sub> PO <sub>4</sub>		Fusion	c	liq		437.	164.			
TlAsS <sub>2</sub>		Fusion	c	liq		573.	300.			
Tl <sub>3</sub> Bi <sub>5</sub>		Fusion	c	liq		487.6	214.4			
3TlCl·8lCl <sub>3</sub>		Fusion	c	liq		686.	413.			
Tl <sub>2</sub> CO <sub>3</sub>		Transition	c, II	c, I		501.	228.			
Tl(C <sub>2</sub> H <sub>5</sub> ) <sub>3</sub>		Fusion	c, I	liq		546.	273.	4.4	8.1	
HCOOTl	thallous formate	Fusion	c	liq		210.	-63.			
CH <sub>3</sub> COOTl	thallous acetate	Fusion	c	liq		377.	104.			
CF <sub>3</sub> COOTl	thallous trifluoroacetate	Fusion	c	liq		403.	130.			
		Vaporization	liq	g	1.5	393.7	120.5			
						433.	160.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 30-5. Thallium (at. no., 81; at. wt., 204.39)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Tl_2CO_3 \cdot TlNO_3$		Fusion	c	liq		487.	214.			
$Tl_2CO_3 \cdot 2TlNO_3$		Fusion	c	liq		491.	218.			
$TlCl \cdot SnCl_2$		Fusion	c	liq		517.	244.			
$3TlCl \cdot SnCl_2$		Fusion	c	liq		583.	310.			
$TlCl \cdot 2PbCl_2$		Fusion	c	liq		708.	435.			
$3TlCl \cdot PbCl_2$		Fusion	c	liq		680.	407.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 31-1. Zinc (at. no., 30; at. wt., 65.38)

HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$0^\circ K$	$0^\circ C$			
Zn		Fusion	c	liq		692.7	419.5	1.595	2.303	
		Vaporization	liq	g	760	1180.	907.	27.43	23.24	
ZnO		Fusion	c	liq		2248.	1975.			
ZnF <sub>2</sub>		Fusion	c	liq		1145.	872.			
		Vaporization	liq	g	760	1775.	1502.	43.1	24.3	
ZnCl <sub>2</sub>		Fusion	c	liq		548.	275.	5.5	10.	
		Vaporization	liq	g	760	1029.	756.	30.9	30.0	
ZnCl <sub>2</sub> ·1½H <sub>2</sub> O		Fusion	c	liq		299.	26.			
ZnCl <sub>2</sub> ·2H <sub>2</sub> O		Fusion	c	liq		292.	19.			
ZnCl <sub>2</sub> ·2½H <sub>2</sub> O		Fusion	c	liq		286.	13.			
ZnCl <sub>2</sub> ·3H <sub>2</sub> O		Fusion	c	liq		280.	7.			
ZnBr <sub>2</sub>		Sublimation	c	g	0.00316	542.	269.	30.1	55.4	
		Fusion	c	liq		667.	394.			
		Vaporization	liq	g	760	970.	697.			
ZnBr <sub>2</sub> ·2H <sub>2</sub> O		Fusion	c	liq		310.	37.			
ZnBr <sub>2</sub> ·3H <sub>2</sub> O		Fusion	c	liq		268.	-5.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 31-2. Zinc (at. no., 30; at. wt., 65.38)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1948 (Corrected)

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
ZnI <sub>2</sub>		Sublimation	c	g	0.00316	518	245.	27.6	53.3	
		Fusion	c	liq		719.	446.			
ZnI <sub>2</sub> · 2H <sub>2</sub> O		Fusion	c	liq		300.	27.			
ZnS		Sublimation	c, II	g	0.01	1127.	854.	64.3	57.1	
		Transition	c, II	c, I		1293.	1020.			
		Sublimation	c, I	g	760	1455.	1182.			
ZnTe		Fusion	c	liq		1511.	1238.			
Zn(NO <sub>3</sub> ) <sub>2</sub> · H <sub>2</sub> O		Fusion	c	liq		343.9	70.7			
Zn(NO <sub>3</sub> ) <sub>2</sub> · 2H <sub>2</sub> O		Fusion	c	liq		338.6	55.4			
Zn(NO <sub>3</sub> ) <sub>2</sub> · 4H <sub>2</sub> O		Fusion	c	liq		317.9	44.7			
Zn(NO <sub>3</sub> ) <sub>2</sub> · 6H <sub>2</sub> O		Fusion	c	liq		309.3	36.1			
2ZnCl <sub>2</sub> · NH <sub>4</sub> Cl		Fusion	c	liq		522.	249.			
ZnAs <sub>2</sub>		Fusion	c	liq		1044.	771.			
ZnSb		Fusion	c	liq		817.	544.			
Zn <sub>3</sub> Sb <sub>2</sub>		Fusion	c	liq		839.	566.	12.1	14.5	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 31-3. Zinc (at. no., 30; at. wt., 65.38)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Zn(CH_3)_2$	dimethylzinc	Vaporization	liq	g	760	321.	48.	6.7	20.9	
$Zn(C_2H_5)_2$	diethylzinc	Vaporization	liq	g	760	397.	124.	8.4	21.2	
$ZnBr_2 \cdot 2CH_3OH$		Fusion	c	liq		312.	39.			
$ZnSiO_3$		Fusion	c	liq		1702.	1429.			
$Zn_2SiO_4$		Fusion	c	liq		1785.	1512.			
$2TiCl \cdot ZnCl_2$		Fusion	c	liq		625.	352.			
$TiCl \cdot 2ZnCl_2$		Fusion	c	liq		499.	226.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 32-1. Cadmium (at. no., 48; at. wt., 112.411)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Cd		Sublimation	c	g	0.11	594.1	320.9	26.28	44.23	
		Fusion	c	liq	0.11	594.1	320.9	1.46	2.46	
		Vaporization	liq	g	760	1040.	767.	23.86	22.94	
CdO		Sublimation	c	g	760	1832.	1559.	53.18	29.4	
CdF <sub>2</sub>		Fusion	c	liq		1383.	1110.	5.4	3.9	
		Vaporization	liq	g	760	2020.	1747.	56.0	27.7	
CdCl <sub>2</sub>		Sublimation	c	g	0.966	841.	568.	41.2	49.0	
		Fusion	c	liq	0.966	841.	568.	5.3	6.3	
		Vaporization	liq	g	760	1253.	980.	29.4	23.5	
CdBr <sub>2</sub>		Sublimation	c	g	0.0032	638.	365.	38.2	59.9	
		Fusion	c	liq		841.	568.	5.0	6.0	
		Vaporization	liq	g	760		1136.	863.		
CdI <sub>2</sub>		Sublimation	c	g	0.48	660.	387.	32.	48.	
		Fusion	c	liq	0.48	660.	387.	8.	12.	



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 32-2. Cadmium (at. no., 48; at. wt., 112.41)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
CdS		Sublimation	c	g	0.0126	958.	685.	51.4	53.7	
CdSO <sub>4</sub>		Fusion	c	liq		1273.	1000.			
CdTe		Fusion	c	liq		1315.	1042.			
Cd(NO <sub>3</sub> ) <sub>2</sub> ·4H <sub>2</sub> O		Fusion	c	liq		332.7	59.5	7.8	23.	
2CdCl <sub>2</sub> ·NH <sub>4</sub> Cl		Fusion	c	liq		639.	366.			
CdSb		Fusion	c	liq		729.	456.	7.66	10.51	
Cd <sub>3</sub> Sb <sub>2</sub>		Fusion	c	liq		694.	421.			
Cd(HCOO) <sub>2</sub> ·3NH <sub>3</sub>	cadmium formate triamine	Fusion	c	liq		343.	70.			
CdSiO <sub>3</sub>		Fusion	c	liq		1428.	1155.			
CdCl <sub>2</sub> ·TiCl		Fusion	c	liq		703.	430.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 33-1. Mercury (at. no., 80; at. wt., 200.61)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1948

Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
		Initial	Final		°K	°C			
Hg	Sublimation	c	g	$2.5 \times 10^{-6}$	234.29	-38.87	15.20	64.9	
	Fusion	c	liq	$2.5 \times 10^{-6}$	234.29	-38.87	0.557	2.37	
	Vaporization	liq	g	0.00209	298.16	25.00	14.54	48.8	
	Vaporization	liq	g, equilibrium	760	629.73	356.57	13.89	22.06	
HgF <sub>2</sub>	Sublimation	c	g	178	575.	302.	16.	28.	
	Fusion	c	liq		918.	645.			
HgCl <sub>2</sub>	Sublimation	c	g	418	550.	277.	18.50	33.6	
	Fusion	c	liq	418	550.	277.	4.15	7.5	
	Vaporization	liq	g	760	577.	304.	14.08	24.4	
HgBr <sub>2</sub>	Sublimation	c	g	116	514.	241.	18.82	36.6	
	Fusion	c	liq	116	514.	241.	3.96	7.7	
	Vaporization	liq	g	760	592.	319.	14.08	23.8	
HgI <sub>2</sub>	Transition	c, II, red	c, I, yellow	0.195	402.	129.	0.601	1.50	
	Sublimation	c, I, yellow	g	8.8	530.	257.	19.86	37.5	
	Fusion	c, I, yellow	liq	8.8	530.	257.	4.53	8.6	
	Vaporization	liq	g	760	627.	354.	14.26	22.7	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 33-2. Mercury (at. no., 80; at. wt., 200.61)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1948

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
HgS		Transition	c, II, red	c, I, black	80	659.	386.	1.0	1.5	
HgI <sub>2</sub> ·2N <sub>2</sub> H <sub>4</sub> ·H <sub>2</sub> O		Fusion	c	liq		346.	73.			
Hg(CH <sub>3</sub> ) <sub>2</sub>	dimethylmercury	Vaporization	liq	g	760	365.8	92.6	8.1	22.1	
Hg(C <sub>2</sub> H <sub>5</sub> ) <sub>2</sub>	diethylmercury	Vaporization	liq	g	760	432.	159.	10.1	23.4	
Hg <sub>2</sub> Tl		Fusion	c	liq		288.	15.			
HgCl <sub>2</sub> ·TlCl		Fusion	c	liq		497.	224.			
HgCl <sub>2</sub> ·2TlNO <sub>3</sub>		Fusion	c	liq		468.	195.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 34-1. Copper (at. no., 29; at. wt., 63.54)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Cu		Fusion Vaporization	c liq	liq g	760	1356.2 2855.	1083.0 2582.	3.11 72.8	2.29 25.4	0.5
Cu <sub>2</sub> O		Fusion	c	liq		1502.	1229.	13.4	8.9	
CuCl		Fusion	c	liq		703.	430.	2.4	3.4	
CuBr		Fusion	c	liq		761.	488.			
CuI		Fusion	c	liq		861.	588.			
CuI <sub>2</sub>		Transition	c, II	c, I		863.	590.			
Cu <sub>2</sub> S		Transition Fusion	c, II c, I	c, I liq		376. 1400.	103. 1127.	1.34 5.5	3.6 3.9	0.2
CuSO <sub>4</sub> ·5H <sub>2</sub> O		Transition	c, II	c, I		327.	54.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 34-2. Copper (at. no., 29; at. wt., 63.54) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION March 31, 1949										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Cu_2Se$		Transition Fusion	c, II c, I	c, I liq		383. 1386.	110. 1113.	1.13	2.95	0.5
$Cu_2Te$		Fusion	c	liq		1128.	855.			
$Cu_4Te_3$		Transition Fusion	c, II c, I	c, I liq		638. 896.	365. 623.			
$Cu(NO_3)_{1/2} \cdot 6H_2O$		Fusion	c	liq		297.6	24.4			
$Cu_3P$		Fusion	c	liq		1303.	1030.			
$Cu_3As$		Fusion	c	liq		1103.	830.			
$Cu_5As_2$		Fusion	c	liq		984.	711.			
$Cu_3Sb$		Fusion	c	liq		960.	687.			
$Cu_5Sb_2$		Fusion	c	liq		956.	683.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards Washington, D.C.

Table 34-3. Copper (at. no., 29; at. wt., 63.54)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1949

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
CuCN			Fusion	c	liq		746.	473.			
$Cu(C_2H_3O_2)_2 \cdot 2NH_3$		cupric acetate diammine	Fusion	c	liq		427.	154.			
$Cu(C_2H_3O_2Cl)_2 \cdot 2NH_3$		cupric chloroacetate diammine	Fusion	c	liq		453.	180.			
$Cu(C_2O_2Cl)_3 \cdot 5NH_3$		cupric trichloroacetate pentammine	Fusion	c	liq		371.	98.			
$CuBr_2 \cdot C_2H_4(NH_2)_2 \cdot 2HBr$		ethylenediammonium cupric bromide	Fusion	c	liq		515.	242.			
$(C_2H_5)_3PO \cdot CuCl_2$		triethylphosphine oxide cupric chloride	Fusion	c	liq		506.	233.			
$Cu_5Si$			Fusion	c	liq		1126.	853.			
$4CuS \cdot SiS_2$			Fusion	c	liq		1468.	1195.			
$Cu_2Ga$			Fusion	c	liq		1108.	835.			
$Cu_2Cd_3$			Fusion	c	liq		895.	562.	11.05	13.23	-3.50

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 35-1. Silver (at. no., 47; at. wt., 107.880)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1949; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Ag		Fusion Vaporization	c liq	liq g	760	1234.0 2456.	960.8 2193.	2.70 60.72	2.19 24.62	0.7
AgF		Fusion	c	liq		708.	435.			
AgCl		Fusion Vaporization	c liq	liq g	760	728. 1830.	455. 1557.	3.16 43.7	4.34 23.9	-2.31
AgClO <sub>3</sub>		Fusion	c	liq		504.	231.			
AgBr		Transition Fusion Vaporization	c, II c, I liq	c, I liq g		532. 703. 1806.	259. 430. 1533.	2.18 37.0	3.10 20.5	-3.6
AgI		Transition Fusion Vaporization	c, II c, I liq	c, I liq g		420. 831. 1777.	147. 558. 1504.	1.47 2.25 34.4	3.50 2.71 19.4	
Ag <sub>2</sub> H <sub>3</sub> IO <sub>6</sub>		Transition	c, II	c, I		227.2	-46.0	0.38	1.7	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 35-2. Silver (at. no., 47; at. wt., 107.880)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Ag_2S$		Transition Fusion	c, II c, I	c, I liq		452. 1115.	179. 842.	1.05 3.36	2.32 3.01	1.2
$Ag_2SO_4$		Transition Fusion	c, II c, I	c, I liq		685. 933.	412. 660.	1.9 4.	2.8 4.	
$Ag_2SO_4 \cdot H_2SO_4$		Transition	c, II	c, I		339.	66.			
$Ag_2Se$		Transition	c, II	c, I		406.	133.	1.7	4.2	0.2
$Ag_2Te$		Transition	c, II	c, I		685.	412.			
$Ag_2Te$		Transition Fusion	c, II c, I	c, I liq		411. 1232.	138. 959.			
$Ag_3Te_2$		Transition Transition	c, III c, II	c, II c, I		559. 698.	286. 425.			



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 35-3. Silver (at. no., 47; at. wt., 107.880)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
AgNO <sub>3</sub>		Transition	c, II	c, I		433.	160.			
		Fusion	c, I	liq		483.	210.	0.66	1.52	-0.06
AgPO <sub>3</sub>		Fusion	c	liq		482.	209.	2.76	5.71	4.5
Ag <sub>3</sub> PO <sub>4</sub>		Fusion	c	liq		1122.	849.			
Ag <sub>4</sub> P <sub>2</sub> O <sub>7</sub>		Fusion	c	liq		858.	585.			
Ag <sub>2</sub> S·As <sub>2</sub> S <sub>3</sub>		Fusion	c	liq		690.	417.			
3Ag <sub>2</sub> S·As <sub>2</sub> S <sub>3</sub>		Fusion	c	liq		763.	490.			
Ag <sub>2</sub> S·Sb <sub>2</sub> S <sub>3</sub>		Fusion	c	liq		782.	509.			
2Ag <sub>2</sub> S·Sb <sub>2</sub> S <sub>3</sub>		Fusion	c	liq		755.	482.			
3Ag <sub>2</sub> S·Sb <sub>2</sub> S <sub>3</sub>		Fusion	c	liq		756.	483.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards Washington, D.C.

Table 35-4. Silver (at. no., 47; at. wt., 107.880)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
AgCN		Fusion	c	liq		623.	350.	2.8	4.5	
Ag <sub>2</sub> HgI <sub>4</sub>		Transition	c, II	c, I		323.9	50.7	1.04	3.21	
AgNO <sub>3</sub> ·HgI <sub>2</sub>		Transition	c, II	c, I		325.	52.			
2AgNO <sub>3</sub> ·HgI <sub>2</sub>		Fusion	c	liq		380.	107.			
Ag <sub>3</sub> Cu <sub>2</sub>		Fusion	c	liq		1052.	779.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 36-1. Gold (at. no., 79; at. wt., 197.2)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Au		Fusion Vaporization	c liq	liq g	760	1336.46 2933.	1063.00 2660.	3.03 74.21	2.27 25.30	-0.53 -2.0
AuTe <sub>2</sub>		Fusion	c	liq		737.	464.			
AuSb <sub>2</sub>		Transition Transition	c, III c, II	c, II c, I		628. 678.	355. 405.	1.12	1.78	2.46
Au <sub>2</sub> Bi		Fusion	c	liq		646.	373.			
AuSn		Fusion	c	liq		691.	418.	6.12	8.86	-0.2
AuGa		Fusion	c	liq		741.	468.			
AuGa <sub>2</sub>		Fusion	c	liq		765.	492.			
AuZn		Fusion	c	liq		1033.	760.	5.9	5.7	-2.6
AuCd		Fusion	c	liq		900.	627.	4.28	4.76	-1.8

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

SERIES II

Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Pt		Fusion Vaporization	c liq	liq g		2042.5 4283.	1769.3 4010.	5.2	2.5	
PtTe <sub>2</sub>		Fusion	c	liq		1523.	1250.			
PtSb		Transition	c, II	c, I		933.	660.			
PtSb <sub>2</sub>		Fusion	c	liq		1483.	1210.			
PtCl <sub>2</sub> ·CO		Fusion	c	liq		468.	195.			
PtCl <sub>2</sub> ·2CO		Fusion	c	liq		415.	142.			
2PtCl <sub>2</sub> ·3CO		Fusion	c	liq		403.	130.			
PtBr <sub>2</sub> ·CO		Fusion	c	liq		451.	178.			
PtI <sub>2</sub> ·CO		Fusion	c	liq		418.	145.			
PtSi		Fusion	c	liq		1501.	1228.			
Pt <sub>2</sub> Si		Transition Fusion	c, II c	c, I liq		973. 1372.	700. 1099.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

SERIES II

Washington, D.C.

Table 38-1. Iridium (at. no., 77; at. wt., 193.1)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Ir		Fusion Vaporization	c liq	liq g	760	2716. 4623.	2443. 4350.			
IrF <sub>6</sub>		Vaporization	liq	g	760	326.	53.	8.6	26.4	
IrCl <sub>5</sub> ·4(C <sub>2</sub> H <sub>5</sub> ) <sub>2</sub> S		Fusion	c	liq		480.	207.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 39-1. Osmium (at. no., 76; at. wt., 190.2)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1949

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
Os			Fusion Vaporization	c liq	liq g	760	2973. 4673.	2700. 4400.			
OsO <sub>4</sub>			Fusion Vaporization	c, l liq	liq g	760	313.3 403.	40.1 130.	3.41 9.5	10.9 23.6	
OsF <sub>6</sub>			Vaporization	liq	g	760	476.	203.			
OsF <sub>8</sub>			Fusion Vaporization	c liq	liq g	760	307.6 320.5	34.4 47.3	6.8	21.2	
Os(CO) <sub>3</sub> Cl <sub>2</sub>			Fusion	c	liq		525.	252.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 40-1. Rhenium (at. no., 75; at. wt., 186.31)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Re		Fusion	c	liq		3420.	3147.			
ReO <sub>3</sub>		Fusion	c	liq		433.	160.			
ReO <sub>4</sub>		Fusion	c	liq		423.	150.			
Re <sub>2</sub> O <sub>7</sub>		Fusion	c	liq		569.	296.	15.3	26.9	
		Vaporization	liq	g	760	635.5	362.3	18.1	28.5	
ReF <sub>4</sub>		Fusion	c	liq		397.7	124.5			
ReF <sub>6</sub>		Fusion	c	liq		292.0	18.8	5.	1.7	
		Vaporization	liq	g	760	321.	48.	6.9	21.5	
ReOF <sub>4</sub>		Sublimation	c	g	331	312.9	39.7	9.5	30.4	
		Fusion	c	liq	331	312.9	39.7			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 40-2. Rhenium (at. no., 75; at. wt., 186.31)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta f$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
ReO <sub>3</sub> Cl		Fusion	c	liq		688.	415.			
		Vaporization	liq	g	760	404.	131.			
ReO <sub>2</sub> Cl <sub>3</sub>		Fusion	c	liq		296.6	23.4			
ReOCl <sub>4</sub>		Fusion	c	liq		302.5	29.3			
TlReO <sub>4</sub>		Fusion	c	liq		600.	527.			
AgReO <sub>4</sub>		Fusion	c	liq		703.	430.			



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$O_K$	$O_C$			
Pd		Fusion Vaporization	c liq	liq g	760	1823. 3833.	1550. 3560.	4.	2.	
PdCl <sub>2</sub>		Fusion	c	liq		951.	678.	9.7	10.2	
PdS		Fusion	c	liq		1243.	970.			
PdSb		Fusion	c	liq		1078.	805.			
Pd <sub>3</sub> Sb		Transition Fusion	c, II c	c, I liq		1223. 1493.	950. 1220.	2.2	1.8	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 42-1. Rhodium (at. no., 45; at. wt., 102.91)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Rh		Fusion Vaporization	c liq	liq g		2233. 4233.	1960. 3960.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 43-1. Ruthenium (at. no., 44; at. wt., 101.7)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Ru		Transition	c, IV	c, III	760	1308.	1035.	0.034	0.026	
		Transition	c, III	c, II		1473.	1200.			
		Transition	c, II	c, I		1773.	1500.			
		Fusion	c, I	liq		2773.	2500.			
		Vaporization	liq	g		4383.	4110.			
RuO <sub>4</sub>		Fusion	c, II	liq		298.	25.			
		Fusion	c, I	liq		300.	27.			
RuF <sub>5</sub>		Fusion	c	liq	190	374.	101.			
		Vaporization	liq	g		500.	227.			
Ru(CO) <sub>5</sub>		Fusion	c	liq		251.	-22.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 45-1. Nickel (at. no., 28; at. wt., 58.69)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Ni		Transition Fusion Vaporization	c, II c, I liq	c, I liq g		626. 1728. 3073.	353. 1455. 2800.	0.092 4.2 91.0	0.15 2.4 29.6	-0.71 0.0
NiO		Sublimation Fusion	c c	g liq	$1.14 \times 10^{-5}$	1500. 2223.	1227. 1950.	111.4	74.2	
NiCl <sub>2</sub>		Sublimation Fusion	c c	g liq	760	1260. 1274.	987. 1001.	48.36	38.38	
NiBr <sub>2</sub>		Fusion	c	liq		1236.	963.			
NiI <sub>2</sub>		Fusion	c	liq		1070.	797.			
NiS		Fusion	c	liq		1070.	797.			
Ni <sub>2</sub> S		Fusion	c	liq		948.	645.	3.0	3.3	
Ni <sub>3</sub> S <sub>2</sub>		Fusion	c	liq		1063.	790.	5.8	5.5	
NiSO <sub>4</sub> ·6H <sub>2</sub> O		Transition	c, II	c, I		326.5	53.3			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 45-2. Nickel [at. no., 28; at. wt., 58.69]  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1949 (Corrected)

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Ni(NO_3)_2 \cdot 6H_2O$		Fusion	c	liq		329.9	56.7			
$Ni(NO_3)_2 \cdot 6NH_3$		Transition	c, II	c, I		243.	-30.	1.7	7.	
$Ni_2P$		Fusion	c	liq		1383.	1110.			
$Ni_5P_2$		Fusion	c	liq		1454.	1181.			
$NiSb$		Fusion	c	liq		1413.	1140.			
$Ni_5Sb_2$		Fusion	c	liq		1423.	1150.			
$Ni(CO)_4$		Fusion	c	liq		248.	-25.			
		Vaporization	liq	g	760	315.6	42.4	7.0	22.2	
$Ni(SCN)_4 \cdot C_2H_4(NH_3)_2$	nickel ethylenediammonium thiocyanate	Fusion	c	liq		368.	95.			
$NiSi$		Fusion	c	liq		1263.	990.			
$Ni_2Si$		Fusion	c	liq		1573.	1300.			
$NiZn_3$		Fusion	c	liq		1153.	880.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 46-1. Cobalt (at. no., 27; at. wt., 58.94)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
June 30, 1949; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Co		Transition	c, III	c, II		723.	450.	0.005	0.007	
		Fusion	c, II	c, I		1398.	1125.	0.07	0.05	
		Vaporization	c, I	g		1766.	1493.	3.7	2.1	
			liq		760	3373.	3100.			
CoO		Fusion	c	liq		2078.	1805.			
CoCl <sub>2</sub>		Fusion	c	liq		997.	724.	7.4	7.4	
		Vaporization	liq	g	760	1323.	1050.	27.2	20.6	
CoBr <sub>2</sub>		Fusion	c	liq		960.	687.			
CoI <sub>2</sub>		Fusion	c	liq		793.	520.			
CoS		Fusion	c	liq		1373.	1100.			
Co(NO <sub>3</sub> ) <sub>2</sub> ·3H <sub>2</sub> O		Fusion	c	liq		364.	91.	8.8	26.7	
Co(NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O		Fusion	c	liq		330.	57.			
CoI <sub>3</sub> ·6NH <sub>3</sub>		Transition	c, II	c, I		277.4	4.2	0.42	1.51	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 46-2. Cobalt (at. no., 27; at. wt., 58.94) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION June 30, 1949										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Co_2P$		Fusion	c	liq			1653.	1380.		
$CoSb$		Fusion	c	liq			1463.	1190.		
$Co(SCN)_4 \cdot C_2H_4(NH_3)_2$	cobalt ethylenediamonium thiocyanate	Fusion	c	liq			355.	82.		
$CoSi$		Fusion	c	liq			1663.	1390.		
$CoSi_2$		Fusion	c	liq			1548.	1275.		
$CoSi_3$		Fusion	c	liq			1578.	1305.		
$Co_2Si$		Fusion	c	liq			1598.	1325.		
$Co_2Sn$		Fusion	c	liq			1438.	1165.		

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 47-1. Iron (at. no., 26; at. wt., 55.85)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
Fe			Transition	c, $\alpha$	c, $\beta$	$2.1 \times 10^{-7}$	1033.	760.	0.0	0.0	0.0
			Transition	c, $\beta$	c, $\gamma$		1180.	907.	0.217	0.184	1.40
			Sublimation	c, $\gamma$	g		1473.	1200.	92.5	62.7	
			Transition	c, $\gamma$	c, $\delta$		1673.	1400.	0.15	0.09	1.
			Fusion	c, $\delta$	liq		1808.	1535.	3.6	2.0	-2.8
			Vaporization	liq	g		3073.	2800.			
Fe <sub>0.95</sub> O		wüstite	Fusion	c	liq		1641.	1368.	7.5	4.6	
Fe <sub>2</sub> O <sub>3</sub>			Transition	c, II	c, I		1303.	1030.			
Fe <sub>3</sub> O <sub>4</sub>			Fusion	c	liq		1867.	1594.	33.	18.	



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 47-2. Iron (at. no., 26; at. wt., 55.85)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$FeF_2$		Fusion	c	liq		1373.	1100.			
$FeCl_2$		Fusion	c	liq		950.	677.	10.28	10.82	3.63
		Vaporization	liq	g	760	1299.	1026.	30.21	23.26	
$FeCl_3$		Sublimation	c	$\frac{1}{2}Fe_2Cl_6(g)$	582	577.	304.	16.5	28.6	
		Fusion	c	$\frac{1}{2}Fe_2Cl_6(liq)$	582	577.	304.	10.3	17.8	-4.
		Vaporization	liq	$\frac{1}{2}Fe_2Cl_6(g)$	582	577.	304.	6.24	10.81	
		Vaporization	liq	$\frac{1}{2}Fe_2Cl_6(g)$	760	592.	319.	6.02	10.17	
$FeCl_3 \cdot 2H_2O$		Fusion	c	liq		346.7	73.5			
$FeCl_3 \cdot 2\frac{1}{2}H_2O$		Fusion	c	liq		329.	56.			
$FeCl_3 \cdot 3\frac{1}{2}H_2O$		Fusion	c	liq		305.7	32.5			
$FeCl_3 \cdot 6H_2O$		Fusion	c	liq		310.	37.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards Washington, D.C.

Table 47-3. Iron (at. no., 26; at. wt., 55.85)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$\text{FeCl}_3 \cdot \text{HCl} \cdot 2\text{H}_2\text{O}$		Fusion	c	liq		318.9	45.7			
$\text{FeCl}_3 \cdot \text{HCl} \cdot 4\text{H}_2\text{O}$		Fusion	c	liq		270.	-3.			
$\text{FeCl}_3 \cdot \text{HCl} \cdot 6\text{H}_2\text{O}$		Fusion	c	liq		267.	-6.			
$\text{FeBr}_2$		Fusion	c	liq		957.	684.			
$\text{FeI}_2$		Fusion	c	liq		860.	587.			
$\text{FeI}_2 \cdot 4\text{H}_2\text{O}$		Fusion	c	liq		388.	95.			
$\text{FeI}_2 \cdot 9\text{H}_2\text{O}$		Fusion	c	liq		274.	1.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 47-4. Iron (at. no., 26; at. wt., 55.85)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
FeS		Transition	c, II	c, I		411.	138.	1.05	2.55	-4.89
FeSO <sub>4</sub> ·7H <sub>2</sub> O		Fusion	c	liq		337.	64.			
Fe(NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O		Fusion	c	liq		333.7	60.5			
Fe(NO <sub>3</sub> ) <sub>3</sub> ·9H <sub>2</sub> O		Fusion	c	liq		323.3	50.1			
FeCl <sub>3</sub> ·NH <sub>4</sub> Cl		Fusion	c	liq		570.	297.			
NH <sub>4</sub> Fe(SO <sub>4</sub> ) <sub>2</sub> ·12H <sub>2</sub> O		Fusion	c	liq		313.	40.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 47-5. Iron (at. no., 26; at. wt., 55.85)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949 (Corrected)

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Fe_2P$		Fusion	c	liq		1633.	1360.			
$Fe_2O_3 \cdot P_2O_5$		Fusion	c	liq		1510.	1237.			
FeAs		Fusion	c	liq		1303.	1030.			
$Fe_2As$		Fusion	c	liq		1192.	919.			
$Fe_3Sb_2$		Fusion	c	liq		1287.	1014.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 47-6. Iron (at. no., 26; at. wt., 55.85)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$\text{Fe}(\text{CO})_5$		Fusion Vaporization	c liq	liq g	760	252. 378.	-21. 105.	3.25 8.9	12.90 23.5	
$\text{Fe}(\text{CO})_4\text{H}_2$		Fusion Vaporization	c liq	liq g	4.2	203.0 243.	-70.2 -30.	6.1	25.	
$\text{Fe}(\text{CO})_2(\text{NO})_2$		Sublimation Fusion Vaporization	c c liq	g liq g	18.7 18.7 760	291.7 291.7 380.	18.5 18.5 107.	12. 2.5 8.6	41. 8.6 22.6	
$\text{FeCl}_3 \cdot \text{CH}_3\text{NH}_3\text{Cl}$	methylammonium ferric chloride	Fusion	c	liq		484.	211.			
$\text{FeCl}_3 \cdot (\text{CH}_3)_2\text{NH}_2\text{Cl}$	dimethylammonium ferric chloride	Fusion	c	liq		373.	100.			
$\text{FeCl}_3 \cdot \text{C}_2\text{H}_5\text{NH}_3\text{Cl}$	ethylammonium ferric chloride	Fusion	c	liq		358.	85.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 47-7. Iron (at. no., 26; at. wt., 55.85)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949 (Corrected)

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		$^{\circ}K$	$^{\circ}C$			
FeSi		Fusion	c	liq		1703.	1430.			
FeO·SiO <sub>2</sub>	FeSiO <sub>3</sub>	Fusion	c	liq		1413.	1140.			
2FeO·SiO <sub>2</sub>	Fe <sub>2</sub> SiO <sub>4</sub>	Fusion	c	liq		1513.	1240.			
4FeO·3SiO <sub>2</sub>	Fe <sub>4</sub> Si <sub>3</sub> O <sub>10</sub>	Fusion	c	liq		1443.	1170.			
Fe <sub>2</sub> O <sub>3</sub> ·PbO		Fusion	c	liq		1803.	1530.			
FeGa <sub>2</sub>		Fusion	c	liq		1063.	790.			
FeCl <sub>3</sub> ·2TiCl		Fusion	c	liq		563.	290.			
FeTi(SO <sub>4</sub> ) <sub>2</sub> ·12H <sub>2</sub> O		Fusion	c	liq		310.	37.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 47-8. Iron (at. no., 26; at. wt., 55.85)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Fe_2O_3 \cdot ZnO$		Fusion	c	liq		1863.	1590.			
$Fe_2O_3 \cdot CdO$		Fusion	c	liq		1813.	1540.			
$Fe_2O_3 \cdot CuO$		Fusion	c	liq		1833.	1560.			
$2FeCl_3 \cdot Cu_2Cl_2$		Fusion	c	liq		593.	320.			
$Fe_2O_3 \cdot NiO$		Fusion	c	liq		1933.	1660.			
$2FeS \cdot Ni_2S$		Fusion	c	liq		1159.	886.			
$Fe_2O_3 \cdot CoO$		Fusion	c	liq		1843.	1570.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 48-1. Manganese (at. no., 25; at. wt., 54.93)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm. Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Mn		Transition	c, IV	c, III	760	1000.	727.	0.54	0.54	
		Transition	c, III	c, II		1374.	1101.	0.54	0.39	
		Transition	c, II	c, I		1410.	1137.	0.43	0.30	
		Fusion	c, I	liq		1517.	1244.	3.50	2.31	
		Vaporization	liq	g		2360.	2087.	53.7	22.8	
MnO		Fusion	c	liq		2053.	1780.			
Mn <sub>2</sub> O <sub>3</sub>		Transition	c, II	c, I		873.	600.			
Mn <sub>2</sub> O <sub>4</sub>		Transition	c, II	c, I		1445.	1172.	4.5	3.1	
		Fusion	c, I	liq		1833.	1560.			



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 48-2. Manganese (at. no., 25; at. wt., 54.93)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION

September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
MnF <sub>2</sub>		Fusion	c	liq		1129.	856.			
MnCl <sub>2</sub>		Fusion	c	liq		923.	650.	9.0	9.7	
		Vaporization	liq	g	760	1463.	1190.	28.8	19.7	
MnBr <sub>2</sub>		Fusion	c	liq		971.	698.			
MnI <sub>2</sub>		Fusion	c	liq		911.	638.			
MnS		Fusion	c	liq		1888.	1615.			
MnSO <sub>4</sub>		Fusion	c	liq		973.	700.			
Mn(NO <sub>3</sub> ) <sub>2</sub> ·3H <sub>2</sub> O		Fusion	c	liq		308.7	35.5	6.5	21.0	
Mn(NO <sub>3</sub> ) <sub>2</sub> ·4H <sub>2</sub> O		Fusion	c	liq		310.3	37.1			
Mn(NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O		Fusion	c	liq		299.0	25.8	9.61	32.14	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 48-3. Manganese (at. no., 25; at. wt., 54.93)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Mn_5P_2$		Fusion	c	liq		1638.	1365.			
$Mn_2Sb$		Fusion	c	liq		1221.	948.			
$Mn_3C$		Transition	c, II	c, I		1310.	1037.	3.1	2.8	
$Mn(SCN)_2 \cdot C_2H_4(NH_3)_2(SCN)_2$	manganese ethylenediammonium thiocyanate	Fusion	c	liq		459.	186.			
$MnSiO_3$		Transition	c, III	c, II		1393.	1120.			
		Transition	c, II	c, I		1481.	1208.			
		Fusion	c, I	liq		1545.	1272.	8.	5.	
$Mn_2SiO_4$		Fusion	c	liq		1563.	1290.			
$MnAu$		Fusion	c	liq		1503.	1230.			
$MnPd$		Fusion	c	liq		1788.	1515.			
$MnFe_2O_4$		Fusion	c	liq		1843.	1570.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 49-1. Chromium (at. no., 24; at. wt., 52.01)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Cr		Fusion	c	liq		2173.	1900.			
Cr <sub>2</sub> O <sub>3</sub>		Transition Fusion	c, II c, I	c, I liq		306.0 2538.	32.8 2265.	3.5	1.6	
CrCl <sub>2</sub>		Sublimation Fusion	c c	g liq	0.61 0.61	1088. 1088.	815. 815.	60.1 7.7	55.2 7.1	
CrO <sub>2</sub> Cl <sub>2</sub>		Vaporization	liq	g	760	380.	117.	8.4	21.5	
CrBr <sub>2</sub>		Fusion	c	liq		1115.	842.			
CrI <sub>2</sub>		Fusion	c	liq		1068.	795.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 49-2. Chromium (at. no., 24; at. wt., 52.01)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		°K	°C			
$\text{NH}_4\text{Cr}(\text{SO}_4)_2 \cdot 12\text{H}_2\text{O}$	chromium	Fusion	c	liq		367.	94.			
CrSb		Fusion	c	liq		1383.	1110.			
$\text{Cr}(\text{CO})_6$		Sublimation	c	g	13	350.	77.	17.2	49.1	
CrSi		Fusion	c	liq		1903.	1630.			
$\text{CrSi}_2$		Fusion	c	liq		1773.	1500.			
$\text{PbCrO}_4$		Transition Transition	c, III c, II	c, II c, I		980. 1056.	707. 783.			
$\text{PbCrO}_4 \cdot \text{PbO}$		Fusion	c	liq		1193.	920.			
$\text{TiCr}(\text{SO}_4)_2 \cdot 12\text{H}_2\text{O}$		Fusion	c	liq		365.	92.			
$\text{Cr}_3\text{Pd}_2$		Fusion	c	liq		1673.	1400.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 50-1. Molybdenum (at. no., 42; at. wt., 95.95)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Mo		Fusion	c	liq		2883.	2610.			
MoO <sub>3</sub>		Sublimation	c	g	0.3	973.	700.	65.	67.	
		Fusion	c	liq	10	1088.	795.			
		Vaporization	liq	g	760	1428.	1155.	33.	31.	
MoF <sub>6</sub>		Sublimation	c	g	406	290.7	17.5	8.3	28.6	
		Fusion	c	liq	406	290.7	17.5	2.2	7.6	
		Vaporization	liq	g	760	308.	35.	6.0	19.5	
MoOF <sub>4</sub>		Fusion	c	liq		370.	97.			
MoCl <sub>5</sub>		Fusion	c	liq		467.	194.			
		Vaporization	liq	g	760	541.	268.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 50-2. Molybdenum (at. no., 42; at. wt., 95.95)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		°K	°C			
$8\frac{1}{2} \text{MoO}_4)_3$		Fusion	c	liq		916.	643.			
MoC		Fusion	c	liq		2950.	2677.			
Mo <sub>2</sub> C		Fusion	c	liq		2945.	2672.			
Mo(CO) <sub>6</sub>		Sublimation	c	g	48	375.	102.	16.3	43.5	
PbMoO <sub>4</sub>		Fusion	c	liq		1338.	1065.			
PbMoO <sub>4</sub> ·PbO		Fusion	c	liq		1224.	951.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 51-1. Tungsten (at. no., 74; at. wt., 183.92)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
W		Fusion	c	liq		3653.	3380.			
WO <sub>3</sub>		Sublimation Fusion	c c	g liq	0.0058	1375. 1743.	1102. 1470.	112.	61.	
WF <sub>6</sub>		Transition Fusion	c, II c, I	c, I liq		265.0 275.7	-8.2 2.5	1.6 0.5	6.0 1.8	
WOF <sub>4</sub>		Fusion Vaporization	c liq	liq g		383. 461.	110. 183.			
WCl <sub>5</sub>		Fusion Vaporization	c liq	liq g		521. 549.	248. 276.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 51-2. Tungsten (at. no., 74; at. wt., 183.92)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
WCl <sub>6</sub>	Transition	c, III		c, II		442.	169.			
	Transition	c, II		c, I	38	500.1	226.9	3.4	6.8	
	Sublimation	c, II		g	38	500.1	226.9	21.0	42.0	
	Sublimation	c, I		g	233	557.2	284.0	17.4	31.2	
	Fusion	c, I		liq	233	557.2	284.0	2.3	4.1	
WCl <sub>4</sub>	Vaporization	liq		g	760	609.7	336.5	14.9	24.4	
	Sublimation	c		g	355	483.	210.	18.1	37.5	
	Fusion	c		liq	355	483.	210.	1.4	2.9	
	Vaporization	liq		g	760	505	232	16.6	32.9	
	Fusion	c		liq		549.	276.			
WBr <sub>5</sub>	Vaporization	liq		g	760	606.	333.			



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 51-3. Tungsten (at. no., 74; at. wt., 183.92)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
WO <sub>3</sub>			Fusion	c	liq		550.	277.			
			Vaporization	liq	g	760	600.	327.			
WCl <sub>6</sub> ·3WBr <sub>6</sub>			Fusion	c	liq		505.	232.			
Bi <sub>2</sub> (WO <sub>4</sub> ) <sub>3</sub>			Fusion	c	liq		1105.	832.			
WC			Fusion	c	liq		3140.	2867.			
W <sub>2</sub> C			Fusion	c	liq		3130.	2857.			
W(CO) <sub>6</sub>			Sublimation	c	g	70	400.	127.	17.7	44.2	
PbWO <sub>4</sub>			Transition	c, II	c, I		1150.	877.			
			Fusion	c, I	liq		1403.	1130.			
PbWO <sub>4</sub> ·PbO			Fusion	c	liq		1178.	905.			
W <sub>2</sub> Re <sub>3</sub>			Fusion	c	liq		3260.	2987.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 52-1. Vanadium (at. no., 23; at. wt., 50.95)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
V		Fusion	c	liq		2003.	1730.			
V <sub>2</sub> O <sub>3</sub>		Fusion	c	liq		2250.	1977.			
V <sub>2</sub> O <sub>4</sub>		Transition Fusion	c, II c, I	c, I liq		345. 1815.	72. 1542.	2.05 27.21	5.94 15.0	6.9 10.1
V <sub>2</sub> O <sub>5</sub>		Fusion	c	liq		943.	670.	15.56	16.50	2.9
VF <sub>5</sub>		Sublimation	c	g	758	384.4	111.2			
VC1 <sub>4</sub>		Fusion Vaporization	c liq	liq g	760	247.5 425.	-25.7 152.	9.1	21.4	
VOC1 <sub>3</sub>		Fusion Vaporization	c liq	liq g	760	194. 400.	-79. 127.	8.3	20.8	
VN		Fusion	c	liq		2300.	2027.			
NH <sub>4</sub> V(SO <sub>4</sub> ) <sub>2</sub> ·12H <sub>2</sub> O		Fusion	c	liq		318.	45.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

SERIES II

Washington, D.C.

Table 52-2. Vanadium (at. no., 23; at. wt., 50.95)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
VC		Fusion	c	liq		3043.	2770.			
VS <sub>i</sub> <sub>2</sub>		Fusion	c	liq		1923.	1650.			
V <sub>2</sub> O <sub>5</sub> ·2PbO		Fusion	c	liq		995.	722.			
V <sub>2</sub> O <sub>5</sub> ·3PbO		Fusion	c	liq		1225.	952.			
V <sub>2</sub> O <sub>5</sub> ·8PbO		Fusion	c	liq		1067.	794.			
3Pb <sub>3</sub> (VO <sub>4</sub> ) <sub>2</sub> ·PbF <sub>2</sub>		Fusion	c	liq		1189.	916.			
3Pb <sub>3</sub> (VO <sub>4</sub> ) <sub>2</sub> ·PbCl <sub>2</sub>		Transition Fusion	c, II c, I	c, I liq		988. 1263.	715. 990.			
TiV(SO <sub>4</sub> ) <sub>2</sub> ·12H <sub>2</sub> O		Fusion	c	liq		321.	48.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 53-1. Columbium (at. no., 41; at. wt., 92.91)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
				Initial	Final		°K	°C			
Cb			Sublimation	c	g	0.00017	2500.	2227.	181.	72.	
			Fusion	c	liq		2760.	2487.			
Cb <sub>2</sub> O <sub>3</sub>			Fusion	c	liq		2045.	1772.			
Cb <sub>2</sub> O <sub>5</sub>			Fusion	c	liq		1733.	1460.			
CbF <sub>5</sub>			Fusion	c	liq		348.7	75.5			
			Vaporization	liq	g	760	498.	225.	11.1	22.3	
CbCl <sub>5</sub>			Sublimation	c	g	330	483.	210.	20.4	42.2	
			Fusion	c	liq	330	483.	210.	8.5	17.6	
			Vaporization	liq	g	760	519.	246.	11.8	22.7	
CbN			Fusion	c	liq		2300.	2027.			
CbC			Fusion	c	liq		3730.	3457.			
Cb <sub>2</sub> Fe <sub>3</sub>			Fusion	c	liq		1933.	1660.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 54-1. Tantalum (at. no., 73; at. wt., 180.88) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION September 30, 1949										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$\alpha_K$	$\alpha_C$			
Ta		Sublimation	c	g	8.6x10 <sup>-6</sup>	2500.	2227.	180.	72.	
Ta <sub>2</sub> O <sub>5</sub>		Fusion	c	liq		3250.	2977.			
TaF <sub>5</sub>		Fusion	c	liq		2163.	1890.			
		Fusion	c	liq		370.0	96.8			
		Vaporization	liq	g	760	402.	229.	6.6	16.4	
TaCl <sub>5</sub>		Sublimation	c	g	415	484.	211.	22.7	46.9	
		Fusion	c	liq	415	484.	211.	11.1	22.9	
		Vaporization	liq	g	760	508.	235.	11.5	22.6	
TaBr <sub>5</sub>		Sublimation	c	g	130	540.	267.	25.5	47.2	
		Fusion	c	liq	130	540.	267.	10.7	19.8	
		Vaporization	liq	g	760	618.	345.	14.6	23.6	
TaI <sub>5</sub>		Fusion	c	liq		638.	365.			
TaN		Fusion	c	liq		3340.	3067.			
TaC		Fusion	c	liq		4100.	3827.			
TaFe		Fusion	c	liq		1973.	1700.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 55-1. Titanium (at. no., 22; at. wt., 47.90)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
Ti			Transition	c, II	c, I	0.036 0.036	1155.	882.	106.5	51.0	
			Sublimation	c, I	g		2085.	1812.			
			Fusion	c, I	liq		2085.	1812.			
TiO			Transition	c, II	c, I		1264.	991.	0.82	0.65	
TiO <sub>2</sub>			Transition	c, III, $\beta$ anatase	c, II, $\alpha$ anatase		915.	642.	0.3	0.3	
			Fusion	c, I, rutile	liq		2108.	1835.			
Ti <sub>2</sub> O <sub>3</sub>			Transition	c, II	c, I		473.	200.	0.22	0.46	
Ti <sub>3</sub> O <sub>5</sub>			Transition	c, II	c, I		450.	177.	0.22	0.49	
TiF <sub>4</sub>			Vaporization	liq	g	760	557.	284.			
TiCl <sub>4</sub>			Fusion	c	liq	760	250.	-23.	2.24	9.0	
			Vaporization	liq	g		409.0	135.8	8.4	20.5	
TiBr <sub>4</sub>			Transition	c, II	c, I	760	258.	-15.	2.1	6.7	
			Fusion	c, I	liq		311.7	38.5			
			Vaporization	liq	g		503.	230.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 55-2. Titanium (at. no., 22; at. wt., 47.90)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
TiI <sub>4</sub>		Fusion Vaporization	c liq	liq g	760	423. 650.3	150. 377.1	13.6	20.9	
TiCl <sub>4</sub> ·SCl <sub>4</sub>		Fusion	c	liq		336.	63.			
TiN		Fusion	c	liq		3200.	2927.			
TiCl <sub>4</sub> ·POCl <sub>3</sub>		Fusion Vaporization	c liq	liq g	760	383. 413.	110. 140.			
TiC		Fusion	c	liq		2400.	2127.			
TiFe <sub>2</sub>		Fusion	c	liq		1803.	1530.			
FeTiO <sub>3</sub>		Fusion	c	liq		1640.	1367.	21.7	13.2	
Fe <sub>2</sub> TiO <sub>4</sub>		Fusion	c	liq		1645.	1372.			
MnTiO <sub>3</sub>		Fusion	c	liq		1430.	1157.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 56-1. Zirconium (at. no., 40; at. wt., 91.22)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949; March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Zr		Transition Sublimation Fusion	c, II c, I c, I	c, I g liq		1135. 2125.	862. 1852.	0.7	0.6	
ZrO		Transition	c, II	c, I		973.	700.			
ZrO <sub>2</sub>		Transition Transition Fusion	c, III c, II c, I	c, II c, I liq		1273. 2173. 2950.	1000. 1900. 2677.		7.0	
ZrCl <sub>4</sub>		Sublimation Fusion	c c	g liq	760	604. 710.	331. 437.	25.3	41.9	
ZrBr <sub>4</sub>		Sublimation Fusion	c c	g liq	760	630. 723.	357. 450.	25.8	41.0	
ZrI <sub>4</sub>		Sublimation Fusion	c c	g liq	760	704. 772.	431. 499.	29.0	41.2	



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 56-2. Zirconium (at. no., 40; at. wt., 91.22)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
ZrN			Fusion	c	liq		3225.	2952.			
Zr (NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O			Fusion	c	liq		309.6	36.4			
2ZrCl <sub>4</sub> ·PCl <sub>5</sub>			Fusion	c	liq		437.7	164.5			
			Vaporization	liq	g	760	689.	416.			
2ZrCl <sub>4</sub> ·POCl <sub>3</sub>			Vaporization	liq	g	760	636.	363.			
ZrC			Fusion	c	liq		3765.	3492.			
ZrSiO <sub>4</sub>			Fusion	c	liq		2693.	2420.			
ZrCu <sub>3</sub>			Fusion	c	liq		1388.	1115.			
ZrAu <sub>3</sub>			Fusion	c	liq		1828.	1555.			
Zr <sub>2</sub> Fe <sub>3</sub>			Fusion	c	liq		1908.	1635.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 57-1. Hafnium (at. no., 72; at. wt., 178.6)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Hf		Transition Fusion	c, II c, I	c, I liq		1773. 2500.	1500. 2227.			
HfO <sub>2</sub>		Fusion	c	liq		3050.	2777.			
HfCl <sub>4</sub>		Sublimation Fusion	c c	g liq	760	590. 705.	317. 432.	24.	40.	
HfBr <sub>4</sub>		Sublimation Fusion	c c	g liq	760	595. 693.	322. 420.	24.	40.	
HfC		Fusion	c	liq		4110.	3837.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 58-1. Boron (at. no., 5; at. wt., 10.82)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
B			Fusion	c	liq		2313.	2040.			
B <sub>2</sub> O <sub>3</sub>			Fusion Vaporization	c liq	liq g	2	723. 1523.	450. 1250.	5.27 77.	7.29 50.	
B <sub>2</sub> H <sub>6</sub>		diborane	Fusion Vaporization	c liq	liq g	760	107.7 180.63	-165.5 -92.53	1.06 3.45	9.8 19.1	
B <sub>4</sub> H <sub>10</sub>		tetraborane	Fusion Vaporization	c liq	liq g	760	153.0 289.	-120.0 16.	6.47	22.4	
B <sub>5</sub> H <sub>9</sub>		pentaborane	Fusion Vaporization	c liq	liq g	760	226.3 331.	-46.9 58.	7.7	23.3	
B <sub>5</sub> H <sub>11</sub>			Fusion Vaporization	c liq	liq g	760	149.9 338.	-123.3 65.	7.61	22.5	
B <sub>6</sub> H <sub>10</sub>		hexaborane	Fusion	c	liq		208.1	-65.1			
B <sub>10</sub> H <sub>14</sub>		decaborane	Sublimation Fusion Vaporization	c c liq	g liq g	20.6 20.6 20.6	372.7 372.7 372.7	99.5 99.5 99.5	19.4 7.8 11.6	52.0 20.9 31.1	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

Washington, D.C.

Table 58-2. Boron (at. no., 5; at. wt., 10.82)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
H <sub>2</sub> O <sub>2</sub>		Fusion	c, I	liq		509.	236.			
BF <sub>3</sub>		Sublimation	c	g	54	144.5	-128.7	5.7	39.	
		Fusion	c	liq	54	144.5	-128.7	1.0	7.	
		Vaporization	liq	g	54	144.5	-128.7	4.7	33.	
		Vaporization	liq	g	760	174.	-99.	4.3	25.	
BF <sub>3</sub> ·H <sub>2</sub> O		Fusion	c	liq		278.7	5.5			
BF <sub>3</sub> ·2H <sub>2</sub> O		Fusion	c	liq		278.	5.			
		Vaporization	liq	g	1.2	332.	59.			
BCl <sub>3</sub>		Fusion	c	liq		166.	-107.			
		Vaporization	liq	g	760	285.6	12.4	5.7	20.0	
B <sub>2</sub> H <sub>6</sub> Cl		Fusion	c	liq		131.2	-142.0			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 58-3. Boron (at. no., 5; at. wt., 10.82)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
BBr <sub>3</sub>			Fusion Vaporization	c liq	liq g		225.7 364.4	-47.5 91.2	7.30	20.03	
B <sub>2</sub> H <sub>5</sub> Br			Fusion Vaporization	c liq	liq g	760	169. 289.	-104. 16.	6.2	21.4	
BI <sub>3</sub>			Fusion Vaporization	c liq	liq g	760	316. 483.	43. 210.			
B <sub>2</sub> S <sub>3</sub>			Fusion	c	liq		583.	310.			
B <sub>2</sub> S <sub>5</sub>			Fusion	c	liq		663.	390.			
B(HSO <sub>4</sub> ) <sub>3</sub>			Fusion	c	liq		488.	215.			
SCl <sub>4</sub> ·BCl <sub>3</sub>			Fusion	c	liq		250.	-23.			
B <sub>3</sub> N <sub>3</sub> H <sub>6</sub>		hexahydro-s-triazatriborine, borazole	Fusion Vaporization	c liq	liq g	760	215. 323.	-58. 50.	7.7	23.8	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 58-4. Boron (at. no., 5; at. wt., 10.82)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949 (Corrected)

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		$^{\circ}K$	$^{\circ}C$			
$B(CH_3)_3$	trimethylborane	Fusion Vaporization	c liq	liq g	180	120. 223.	-153. -50.	5.7	25.6	
$(CH_3)_2B_2H_4$	unsym. dimethyldiborane	Fusion Vaporization	c liq	liq g	760	123.0 271.	-150.2 -2.	5.5	20.3	
$(CH_3)_3B_2H_3$	1,1,2-trimethyldiborane	Fusion Vaporization	c liq	liq g	760	150.3 318.	-122.9 45.	7.0	22.0	
$(CH_3)_4B_2H_2$	sym. tetramethyldiborane	Fusion Vaporization	c liq	liq g	760	200.7 345.	-72.5 72.	7.3	21.1	
$(C_2H_5)_2B_2H_4$	unsym. diethyldiborane	Vaporization	liq	g	760	340.	67.	8.1	23.8	
$(C_2H_5)_4B_2H_2$	sym. tetraethyldiborane	Fusion	c	liq		216.9	-56.3			
$(CH_3O)_2BH$	dimethoxyborane	Fusion Vaporization	c liq	liq g	760	142.6 298.	-130.6 25.	6.6	22.1	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 58-5. Boron (at. no.. 5; at. wt., 10.82)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		°K	°C			
$B_3O_3(CH_3)_3$	trimethylboroxole		Fusion Vaporization	c liq	liq g	7.9	235. 298.	-38. 25.	8.7	29.2	
$B(CH_3)_2F$	methyl difluoroborine		Fusion Vaporization	c liq	liq g	760	142.8 210.	-130.4 -63.	5.0	23.8	
$B(CH_3)_2F$	dimethyl fluoroborine		Fusion Vaporization	c liq	liq g	760	125.9 231.	-147.3 -42.	5.1	22.1	
$B(CH_3)_2Cl$	methyl dichloroborine		Fusion Vaporization	c liq	liq g	760	146.1 284.3	-127.1 11.1			
$B(CH_3)_2Cl$	dimethyl chloroborine		Fusion Vaporization	c liq	liq g	760	278.1 188.6	4.9 -84.6	5.7	20.5	
$B(C_2H_5)_2Cl$	diethyl chloroborine		Fusion Vaporization	c liq	liq g	760	351.7 162.6	78.5 -110.6	7.9	22.4	
$B(CH_3)_2Br$	methyl dibromoborine		Fusion Vaporization	c liq	liq g	760	333.2 144.3	60.0 -128.9	7.3	22.0	
$B(CH_3)_2Br$	dimethyl bromoborine		Fusion Vaporization	c liq	liq g	760	304.8 192.2	31.6 -81.0	6.3	20.6	
$B(C_2H_5)_2Br$	diethyl bromoborine		Fusion Vaporization	c liq	liq g	760	374.2 165.7	101.0 -107.5	8.4	22.5	
$B(CH_3)_2I$	dimethyl iodoborine		Fusion Vaporization	c liq	liq g	760	342.0	68.8	6.6	19.3	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

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Table 58-6. Boron (at. no., 5; at. wt., 10.82)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$BH_2N(CH_3)_2$	dimethylaminoborine	Fusion	c	liq		346.7	73.5			
$BH_3 \cdot NH(CH_3)_2$	dimethylamine borine	Fusion	c	liq		284.	11.			
$B(CH_3)_2NH(CH_3)$	methylaminodimethylborine	Vaporization	liq	g	760	313.	40.	6.1	19.5	
$B(CH_3)_2NH_2$	aminodimethylborine	Fusion	c	liq		282.	9.			
$B(CH_3)_3 \cdot NH_2CH_3$	methylamine trimethylborine	Fusion	c	liq		299.6	26.4			
$BH[N(CH_3)_2]_2$	bis(dimethylamino)borine	Fusion	c	liq		228.	-45.			
$B(CH_3)_3 \cdot NH(CH_3)_2$	dimethylamine trimethylborine	Fusion	c	liq		307.	34.			
$B(CH_3)_2N(C_2H_5)_2$	diethylaminodimethylborine	Vaporization	liq	g	760	371.	98.	8.8	23.7	
$B(C_2H_5)_2N(CH_3)_2$	dimethylaminodiethylborine	Vaporization	liq	g	760	396.	123.	9.0	22.8	
$B[N(CH_3)_2]_3$	tris(dimethylamino)borine	Fusion	c	liq		233.	-40.			
		Vaporization	liq	g	760	420.	147.	10.5	25.0	
$[B(CH_3)_2NH(CH_3)]_2$	bis(methylaminodimethylborine)	Vaporization	liq	g	760	312.	39.	6.2	19.9	
$[B(N(CH_3)_2)_3]_2$		Vaporization	liq	g	760	420.	147.			



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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SERIES II

Washington, D.C.

Table 58-7. Boron (at. no., 5; at. wt., 10.82)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$(CH_3)_3B_3N_3H_5$	1-methylborazole (N-methylborazole)	Vaporization	liq	g	760	357.	84.	8.6	24.1	
$(CH_3)_2B_3N_3H_5$	2-methylborazole (B-methylborazole)	Fusion Vaporization	c liq	liq g	760	214. 360.	-59. 87.	9.0	25.0	
$(CH_3)_2B_3N_3H_4$	1,3-dimethylborazole	Vaporization	liq	g	760	381.	108.	8.3	21.8	
$(CH_3)_2B_3N_3H_4$	2,4-dimethylborazole	Fusion Vaporization	c liq	liq g	760	225. 382.	-48. 109.	9.0	23.6	
$(CH_3)_2B_3N_3H_4$	1,2-dimethylborazole	Vaporization	liq	g	760	388.	115.	8.3	21.4	
$(CH_3)_3B_3N_3H_3$	1,3,5-trimethylborazole	Fusion Vaporization	c liq	liq g	760	284. 406.	-9. 133.	9.3	22.9	
$(CH_3)_3B_3N_3H_3$	2,4,6-trimethylborazole	Fusion Vaporization	c liq	liq g	760	305.0 402.	31.8 129.	9.3	23.1	
$(CH_3)_3B_3N_3H_3$	1,2,4-trimethylborazole	Vaporization	liq	g	760	445.	142.	8.6	20.7	
$(CH_3)_4B_3N_3H_2$	1,2,4,6-tetramethylborazole	Vaporization	liq	g	760	431.	158.	10.4	24.1	
$(CH_3)_6B_3N_3$	hexamethylborazole	Fusion Vaporization	c liq	liq g	760	370.3 496.	97.1 223.	11.	22.	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

Washington, D.C.

Table 58-8. Boron (at. no., 5; at. wt., 10.82)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
$(CH_3)_3B_3O_3 \cdot NH_2(CH_3)$	trimethylboroxole methylamine		Fusion	c	liq		455.	182.			
$BCl_2N(CH_3)_2$	dimethylaminodichloroborine		Fusion Vaporization	c liq	liq g	760	227.1 385.	-46.1 112.	8.9	23.1	
$BCl[N(CH_3)_2]_2$	bis(dimethylamino)chloroborine		Fusion Vaporization	c liq	liq g	760	219. 418.	-54. 145.	9.8	23.4	
$[BCl_2N(CH_3)_2]_2$	bis(dimethylaminodichloroborine)		Fusion	c	liq		415.	142.			
$BBr_2N(CH_3)_2$	dimethylaminodibromoborine		Fusion	c	liq		251.7	-21.5			
$[BBr_2N(CH_3)_2]_2$	bis(dimethylaminodibromoborine)		Fusion	c	liq		425.	152.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

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Table 58-9. Boron (at. no., 5; at. wt., 10.82)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$PbO \cdot B_2O_3$		Fusion	c	liq		773.	500.			
$PbO \cdot 2B_2O_3$		Fusion	c	liq		843.	570.			
$PbO \cdot 3B_2O_3$		Fusion	c	liq		833.	560.			
$2PbO \cdot 5B_2O_3$		Fusion	c	liq		793.	520.			
$3ZnO \cdot 2B_2O_3$		Fusion	c	liq		1253.	980.			
$CdO \cdot B_2O_3$		Fusion	c	liq		1123.	850.			
$2CdO \cdot 3B_2O_3$		Fusion	c	liq		1128.	855.			
$Ni_2B$		Fusion	c	liq		1503.	1230.			
$Ni_3B_2$		Transition Fusion	c, II c, I	c, I liq		1323. 1433.	1050. 1160.			
$Fe_2B$		Fusion	c	liq		1693.	1420.			
$MnO \cdot B_2O_3$		Fusion	c	liq		1113.	840.			
$MnO \cdot 2B_2O_3$		Fusion	c	liq		1213.	940.			
$MnO \cdot 3B_2O_3$		Fusion	c	liq		1193.	920.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 58-10. Boron (at. no., 5; at. wt., 10.82)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
September 30, 1949

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
ZrB			Fusion	c	liq		3245.	2972.			
Zr(BH <sub>4</sub> ) <sub>4</sub>		zirconium borohydride	Sublimation	c	g	19.0	301.9	28.7	13.	43.	
			Fusion	c	liq	19.0	301.9	28.7	4.	13.	
			Vaporization	liq	g	19.0	301.9	28.7	9.3	30.8	
HfB			Fusion	c	liq		3315.	3042.			
Hf(BH <sub>4</sub> ) <sub>4</sub>		hafnium borohydride	Sublimation	c	g	20.0	302.	29.	13.	43.	
			Fusion	c	liq	20.0	302.	29.	3.	10.	
			Vaporization	liq	g	20.0	302.	29.	9.7	32.1	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 59-1. Aluminum (at. no., 13; at. wt., 26.97)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		$^{\circ}K$	$^{\circ}C$			
Al		Fusion	c	liq		933.	660.	2.6	2.8	
		Vaporization	liq	g	760	2600.	2327.	67.9	26.	
Al <sub>2</sub> O <sub>3</sub>		Fusion	c	liq		2313.	2040.	26.	11.	
AlF <sub>3</sub>		Sublimation	c	g	760	1530.	1257.	77.	50.	
Al <sub>2</sub> Cl <sub>6</sub>		Sublimation	c	g	760	453.3	180.1	26.7	58.9	
		Sublimation	c	g	1625	465.	192.	26.5	57.0	
		Fusion	c	liq	1625	465.	192.	16.9	36.3	
Al <sub>2</sub> Br <sub>6</sub>		Fusion	c	liq		370.7	97.5	5.4	14.6	
		Vaporization	liq	g	760	530.	257.	11.	21.	
Al <sub>2</sub> I <sub>6</sub>		Fusion	c	liq		464.	191.	3.8	8.2	
		Vaporization	liq	g	760	659.	386.	15.4	23.4	
AlBr <sub>3</sub> ·H <sub>2</sub> S		Fusion	c	liq		377.	104.			
Al <sub>3</sub> Se <sub>4</sub>		Fusion	c	liq		1223.	950.			
Al <sub>2</sub> Te <sub>3</sub>		Fusion	c	liq		1173.	900.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 59-2. Aluminum (at. no., 13; at. wt., 26.97)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$0^\circ\text{C}$	$0^\circ\text{C}$			
$\text{AlCl}_3 \cdot \text{NH}_3$		Fusion Vaporization	c liq	liq g	760	398. 696.9	125. 423.7	17.6	25.2	
$\text{AlBr}_3 \cdot \text{NH}_3$		Fusion	-c	liq		397.	124.			
$\text{AlI}_3 \cdot \text{NH}_3$		Fusion	c	liq		399.	126.			
$\text{NH}_4\text{Al}(\text{SO}_4)_2 \cdot 12\text{H}_2\text{O}$		Transition Fusion	c, II c, I	c, I liq		71.0 368.	-202.2 95.	0.194	2.73	
$\text{AlCl}_3 \cdot \text{PH}_3$		Fusion	c	liq		355.	82.			
$\text{AlCl}_3 \cdot \text{POCl}_3$		Fusion	c	liq		438.	165.			
$\text{AlBr}_3 \cdot \text{PH}_3$		Fusion	c	liq		390.	117.			
$\text{AlI}_3 \cdot \text{PH}_3$		Fusion	c	liq		422.	149.			
$\text{AlSb}$		Fusion	c	liq		1353.	1080.			
$\text{AlBr}_3 \cdot \text{SbBr}_3$		Fusion	c	liq		357.	84.			
$\text{AlBr}_3 \cdot \text{BiBr}_3$		Fusion	c	liq		424.	151.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 59-3. Aluminum (at. no., 13; at. wt., 26.97)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$\text{Al}(\text{CH}_3)_3$	trimethylaluminum	Fusion Vaporization	c liq	liq g	43	288.6 323.	15.4 50.	9.7	30.0	
$\text{Al}(\text{C}_2\text{H}_5)_3$	triethylaluminum	Vaporization	liq	g	5.9	353.	80.	13.	37.	
$(\text{CH}_3)_2\text{O} \cdot \text{Al}(\text{CH}_3)_3$	trimethylaluminum dimethyl ether	Fusion Vaporization	c liq	liq g	25	243.3 333.	-29.9 60.	10.6	31.8	
$[(\text{CH}_3)_2\text{AlO}(\text{CH}_3)]_3$	tris(dimethylaluminum methoxide)	Fusion Vaporization	c liq	liq g	30	308. 364.	35. 91.	13.1	36.0	
$(\text{CH}_3\text{AlCl}_2)_2$	bis(methylaluminum dichloride)	Fusion	c	liq		344.7	71.5			
$[(\text{CH}_3)_2\text{AlCl}]_2$	bis(dimethylaluminum chloride)	Fusion Fusion Vaporization	c, II c, I liq	liq liq g	760	228. 252.2 397.	-45. -21.0 124.	9.1	22.9	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 59-4. Aluminum (at. no., 13; at. wt., 26.97)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		°K	°C			
$(CH_3)_2O \cdot Al(CH_3)_2Cl$	dimethylaluminum chloride dimethyl ether		Vaporization	liq	g	40	390.	117.	10.6	27.2	
$(C_2H_5)_2O \cdot AlCl_3$	aluminum chloride diethyl ether		Fusion	c	liq		310.	37.			
$[(CH_3)_2AlBr]_2$	bis(dimethylaluminum bromide)		Fusion	c	liq		253.	-20.			
			Vaporization	liq	g	40	341.	68.	10.3	30.2	
$(C_2H_5)_2O \cdot AlBr_3$	aluminum bromide diethyl ether		Fusion	c	liq		320.	47.			
$(CH_3)_2S \cdot Al(CH_3)_3$	trimethylaluminum dimethyl sulfide		Fusion	c	liq		253.	-20.			
			Vaporization	liq	g	32	340.	67.	11.3	33.2	
$[(CH_3)_2AlSCH_3]_2$	bis(methylmercaptodimethylaluminum)		Fusion	c	liq		376.	103.			
			Vaporization	liq	g	26	404.	131.	13.6	33.7	



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 59-5. Aluminum (at. no., 13; at. wt., 26.97)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		O <sub>K</sub>	O <sub>C</sub>			
$\text{AlH}_3 \cdot \text{N}(\text{CH}_3)_3$	aluminum hydride trimethylamine	Fusion	c	liq		338.	65.			
$\text{Al}(\text{CH}_3)_3 \cdot \text{NH}(\text{CH}_3)_2$	trimethylaluminum dimethylamine	Fusion	c	liq		324.	51.			
		Vaporization	liq	g	10	346.	73.	12.0	34.7	
$\text{Al}(\text{CH}_3)_3 \cdot \text{N}(\text{CH}_3)_3$	trimethylaluminum trimethylamine	Sublimation	c	g	79	378.	105.	12.	32.	
		Vaporization	liq	g	79	378.	105.	11.	29.	
$\text{AlH}_3 \cdot 2\text{N}(\text{CH}_3)_3$	aluminum hydride di(trimethylamine)	Fusion	c	liq		363.	90.			
$[(\text{CH}_3)_2\text{AlN}(\text{CH}_3)_2]_2$	bis(dimethylamino dimethylaluminum)	Sublimation	c	g	16	358.	85.	12.9	36.0	
		Fusion	c	liq		428.	155.			
$\text{AlCl}_3 \cdot \text{C}_2\text{H}_5\text{NH}_2$	aluminum chloride ethylamine	Fusion	c	liq		396.	63.			
$\text{Al}(\text{CH}_3)_2\text{Cl} \cdot (\text{CH}_3)_3\text{N}$	dimethylaluminum chloride trimethylamine	Sublimation	c	g	7.1	370.	97.	13.3	35.9	
		Fusion	c	liq		397.	124.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

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Washington, D.C.

Table 59-6. Aluminum (at. no., 13; at. wt., 26.97) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION December 31, 1949											
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole	
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$				
$(CH_3)_2PH \cdot Al(CH_3)_3$	trimethylaluminum dimethylphosphine	Vaporization	liq	g	50	363.	90.	11.3	31.1		
$(CH_3)_3P \cdot Al(CH_3)_3$	trimethylaluminum trimethylphosphine	Sublimation	c	g	7.9	335.7	62.5	14.5	43.2		
		Vaporization	liq	g	7.9	335.7	62.5	11.7	34.8		
$[(CH_3)_2AlP(CH_3)_2]_3$	tris (dimethylphosphinodimethylaluminum)	Sublimation	c	g	20	460.	187.	14.	30.		
$Al_2O_3 \cdot SiO_2$		Fusion	c	liq		2085.	1812.				
$Al_2Cl_6 \cdot SnCl_2$		Fusion	c	liq		482.5	209.3				
$Al_2Cl_6 \cdot 2SnCl_2$		Fusion	c	liq		431.7	158.5				
$Al_2Br_6 \cdot SnBr_2$		Fusion	c	liq		478.	205.				
$Al_2Br_6 \cdot 2SnBr_2$		Fusion	c	liq		456.	183.				
$Al_2Br_6 \cdot PbBr_2$		Fusion	c	liq		547.	274.				

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 59-7. Aluminum (at. no., 13; at. wt., 26.97)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
AlGa		Fusion	c	liq		647.	374.			
AlGa <sub>2</sub>		Fusion	c	liq		554.	281.			
Al <sub>2</sub> Ga		Fusion	c	liq		740.	467.			
Al <sub>2</sub> Cl <sub>6</sub> ·2TlCl		Fusion	c	liq		570.	297.			
Al <sub>2</sub> Br <sub>6</sub> ·TlBr		Fusion	c	liq		385.	112.			
Al <sub>2</sub> Br <sub>6</sub> ·2TlBr		Fusion	c	liq		483.	210.			
AlTl(SO <sub>4</sub> ) <sub>2</sub> ·12H <sub>2</sub> O		Fusion	c	liq		364.	91.			
Al <sub>2</sub> O <sub>3</sub> ·ZnO		Fusion	c	liq		2218.	1945.			
Al <sub>2</sub> Br <sub>6</sub> ·ZnBr <sub>2</sub>		Fusion	c	liq		384.7	111.5			
Al <sub>2</sub> Br <sub>6</sub> ·CdBr <sub>2</sub>		Fusion	c	liq		497.	224.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 59-8. Aluminum (at. no., 13; at. wt., 26.97)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$Al_2Br_6 \cdot HgBr_2$		Fusion	c	liq		377.	104.			
$Al_2Br_6 \cdot Hg_2Br_2$		Fusion	c	liq		534.	261.			
$AlCu$		Fusion	c	liq		898.	625.			
$AlCu_3$		Fusion	c	liq		1323.	1050.			
$Al_2Cu$		Fusion	c	liq		863.	590.			
$Al_2Cl_6 \cdot Cu_2Cl_2$		Fusion	c	liq		506.	233.			
$AlAg_2$		Fusion	c	liq		994.	721.			
$AlAg_3$		Fusion	c	liq		1044.	771.			
$Al_2Br_6 \cdot 2AgBr$		Fusion	c	liq		488.8	215.6			
$AlAu_2$		Fusion	c	liq		896.	623.			
$Al_2Au$		Fusion	c	liq		1330.	1057.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 59-9. Aluminum (at. no., 13; at. wt., 26.97)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		°K	°C			
NiAl			Fusion	c	liq		1945.	1642.			
Al <sub>2</sub> O <sub>3</sub> ·NiO			Fusion	c	liq		2286.	2015.			
CoAl			Fusion	c	liq		1900.	1627.			
Al <sub>2</sub> O <sub>3</sub> ·CoO			Fusion	c	liq		2228.	1955.			
FeAl <sub>3</sub>			Fusion	c	liq		1423.	1150.			
Al <sub>3</sub> Ti			Fusion	c	liq		1628.	1355.			
Al <sub>4</sub> Ti			Fusion	c	liq		1598.	1325.			
Al <sub>2</sub> O <sub>3</sub> ·TiO <sub>2</sub>			Fusion	c	liq		2128.	1855.			
Al <sub>2</sub> O <sub>3</sub> ·2TiO <sub>2</sub>			Fusion	c	liq		2168.	1895.			
Al (BH <sub>4</sub> ) <sub>3</sub>		aluminum borohydride	Fusion	c	liq		208.7	-64.5			
			Vaporization	liq	g	760	348.	45.	7.2	22.6	
Al (BH <sub>4</sub> ) <sub>3</sub> ·N(CH <sub>3</sub> ) <sub>3</sub>		aluminum borohydride trimethylamine	Fusion	c	liq		352.	79.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 60-1. Scandium (at. no., 21; at. wt., 45.10)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Sc	Fusion Vaporization		c liq	liq g	760	1673. 4173.	1400. 3900.			
ScCl <sub>3</sub>	Sublimation Fusion		c c	g liq	684 684	1233. 1233.	960. 960.	61.1	49.6	
ScBr <sub>3</sub>	Sublimation Fusion		c c	g liq	760	1200. 1233.	927. 960.	62.4	52.0	
ScI <sub>3</sub>	Sublimation Fusion		c c	g liq	760	1183. 1218.	910. 945.	60.5	51.1	
ScN	Fusion		c	liq		2920.	2647.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

SERIES II

Washington, D.C.

Table 61-1. Yttrium (at. no., 39; at. wt., 88.92)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Y		Fusion	c	liq		1773.	1500.			
		Vaporization	liq	g	760	4373.	4100.			
YCl <sub>3</sub>		Fusion	c, y	liq		973.	700.			
YBr <sub>3</sub>		Fusion	c	liq		1177.	904.			
YI <sub>3</sub>		Fusion	c	liq		1773.	1000.			
Y <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>		Fusion	c	liq		1620.	1347.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 62-1. Lutetium (at. no., 71; at. wt., 174.99)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$\text{LuCl}_3$		Fusion	c, γ	liq		1165.	892.			
$\text{LuI}_3$		Fusion	c	liq		1318.	1045.			



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

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Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
YbCl <sub>3</sub>		Fusion	c, γ	liq		1127.	854.			
YbBr <sub>3</sub>		Fusion	c	liq		1213.	940.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 64-1. Thulium (at. no., 69; at. wt., 169.4)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$T_m$		Vaporization	liq	g	760	3773.	3500.			
$T_mCl_3$		Fusion	c, γ	liq		1118.	845.			
$T_mI_3$		Fusion	c	liq		1288.	1015.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 65-1. Erbium (at. no., 68; at. wt., 167.2)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
ErCl <sub>3</sub>		Fusion	c, γ	liq		1047.	774.			
ErBr <sub>3</sub>		Fusion	c	liq		1223.	950.			
ErI <sub>3</sub>		Fusion	c	liq		1293.	1020.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 66-1. Holmium (at. no., 67; at. wt., 164.94)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}\text{K}$	$^{\circ}\text{C}$			
$\text{HoCl}_3$		Fusion	c, y	liq		991.	718.			
$\text{HoBr}_3$		Fusion	c	liq		1187.	914.			
$\text{HoI}_3$		Fusion	c	liq		1283.	1010.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 67-1. Dysprosium (at. no., 66; at. wt., 162.46)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$DyCl_3$		Fusion	c, $\beta$	liq		928.	655.			
$DyBr_3$		Fusion	c	liq		1154.	881.			
$DyI_3$		Fusion	c	liq		1228.	955.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta G$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
TbCl <sub>3</sub>		Fusion	c, β	liq		861.	588.			

Table 68-1. Terbium (at. no., 65; at. wt., 159.2)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
GdCl <sub>3</sub>		Fusion	c, $\alpha$	liq		882.	609.			
GdBr <sub>3</sub>		Fusion	c	liq		1038.	765.			
GdI <sub>3</sub>		Fusion	c	liq		1199.	926.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 70-1. Europium (at. no., 63; at. wt., 152.0)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
EuCl <sub>3</sub>		Fusion	c, $\alpha$	liq		896.	623.			



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Washington, D.C.

Table 71-1. Samarium (at. no., 62; at. wt., 150.43)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$0^\circ K$	$0^\circ C$			
Sm		Fusion	c	liq		1623.	1350.			
SmCl <sub>2</sub>		Fusion	c	liq		1013.	740.			
SmCl <sub>3</sub>		Fusion	c, $\alpha$	liq		955.	682.			
SmBr <sub>3</sub>		Fusion	c	liq		937.	664.			
SmI <sub>3</sub>		Fusion	c	liq		1093.	820.			
Sm <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>		Fusion	c	liq		1347.	1074.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 73-1. Neodymium (at. no., 60; at. wt., 144.27)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Nd		Transition	c, IV	c, III		109.	-164.			
		Transition	c, III	c, II		781.	508.			
		Transition	c, II	c, I		987.	714.			
		Fusion	c, I	liq		1113.	840.			
NdF <sub>3</sub>		Fusion	c	liq		1683.	1410.			
NdCl <sub>3</sub>		Fusion	c, α	liq		1033.	760.			
NdCl <sub>3</sub> ·6H <sub>2</sub> O		Fusion	c	liq		399.	126.			
NdBr <sub>3</sub>		Fusion	c	liq		957.	684.			
NdI <sub>3</sub>		Fusion	c	liq		1048.	775.			
Nd <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>		Fusion	c	liq		1449.	1176.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards Washington, D.C.

Table 74-1. Praseodymium (at. no., 59; at. wt., 140.92)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Pr		Fusion	c, I	liq		1205.	932.	2.7	2.2	
PrF <sub>3</sub>		Fusion	c	liq		1643.	1370.			
PrCl <sub>3</sub>		Fusion	c, $\alpha$	liq		1096.	823.			
PrCl <sub>3</sub> ·7H <sub>2</sub> O		Fusion	c	liq		379.	106.			
PrBr <sub>3</sub>		Fusion	c	liq		966.	693.			
PrI <sub>3</sub>		Fusion	c	liq		1006.	733.			
PrSn <sub>3</sub>		Fusion	c	liq		1423.	1150.			
Pr <sub>2</sub> Sn		Fusion	c	liq		1683.	1410.			
Pr <sub>2</sub> Sn <sub>3</sub>		Fusion	c	liq		1433.	1160.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 74-2. Praseodymium (at. no., 59; at. wt., 140.92)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
PrPb		Fusion	c	liq		1458.	1185.			
PrPb <sub>3</sub>		Fusion	c	liq		1385.	1112.			
Pr <sub>2</sub> Pb		Fusion	c	liq		1585.	1312.			
PrGa <sub>2</sub>		Fusion	c	liq		1743.	1470.			
PrTi		Fusion	c	liq		1423.	1150.			
PrTi <sub>3</sub>		Fusion	c	liq		1333.	1060.			
PrCu <sub>2</sub>		Fusion	c	liq		1114.	841.			
PrCu <sub>6</sub>		Fusion	c	liq		1235.	962.			
PrAg <sub>4</sub>		Fusion	c	liq		1201.	928.			
PrAg <sub>3</sub>		Fusion	c	liq		1228.	955.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 74-3. Praseodymium (at. no., 59; at. wt., 140.92)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
PrAu		Fusion	c	liq		1523.	1350.			
PrAu <sub>2</sub>		Fusion	c	liq		1483.	1210.			
PrAu <sub>3</sub>		Fusion	c	liq		1473.	1200.			
PrAu <sub>4</sub>		Fusion	c	liq		1473.	1200.			
PrNi		Fusion	c	liq		1003.	730.			
PrNi <sub>5</sub>		Fusion	c	liq		1638.	1365.			
Pr <sub>3</sub> Ni		Fusion	c	liq		803.	530.			
Pr <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>		Fusion	c	liq		1303.	1030.			
PrAl <sub>2</sub>		Fusion	c	liq		1713.	1440.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Ce		Transition	c, IV	c, III		140.	-133.			
		Transition	c, III	c, II		666.	393.			
		Transition	c, II	c, I		713.	440.			
		Fusion	c, I	liq		1048.	775.	2.1	2.0	
Ce <sub>2</sub> O <sub>3</sub>		Fusion	c	liq		1960.	1687.			
CeF <sub>3</sub>		Fusion	c	liq		1733.	1460.			
CeCl <sub>3</sub>		Fusion	c, α	liq		1095.	822.	8.	7.	
CeBr <sub>3</sub>		Fusion	c	liq		1005.	732.			
CeI <sub>3</sub>		Fusion	c	liq		1025.	752.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 75-2. Cerium (at. no., 58; at. wt., 140.13)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
CeBi		Fusion	c	liq		1798.	1525.			
Ce <sub>4</sub> Bi <sub>3</sub>		Fusion	c	liq		1903.	1630.			
CeSn <sub>2</sub>		Fusion	c	liq		1408.	1135.			
CeSn <sub>3</sub>		Fusion	c	liq		1436.	1163.			
Ce <sub>2</sub> Sn		Fusion	c	liq		1673.	1400.			
Ce <sub>2</sub> Sn <sub>3</sub>		Fusion	c	liq		1438.	1165.			
CePb <sub>3</sub>		Fusion	c	liq		1453.	1180.			
Ce <sub>2</sub> Pb		Fusion	c	liq		1663.	1390.			
CeTi		Fusion	c	liq		1508.	1235.			
CeTi <sub>3</sub>		Fusion	c	liq		1353.	1080.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 75-3. Cerium (at. no., 58; at. wt., 140.13)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949 (Corrected)

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
CeZng		Fusion	c	liq		1245.	.972.			
CeCu <sub>2</sub>		Fusion	c	liq		1093.	820.			
CeCu <sub>6</sub>		Fusion	c	liq		1213.	940.			
CeAg		Fusion	c	liq		1128.	855.			
CeAg <sub>3</sub>		Fusion	c	liq		1263.	990.			
CeAu		Fusion	c	liq		1623.	1350.			
CeAu <sub>2</sub>		Fusion	c	liq		1403.	1130.			
CeAu <sub>3</sub>	•	Fusion	c	liq		1423.	1150.			



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 75-4. Cerium (at. no., 58; at. wt., 140.13)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949; March 31, 1950

Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
			Initial	Final		$^{\circ}K$	$^{\circ}C$			
CeNi		Fusion	c	liq		913.	640.			
CeNi <sub>5</sub>		Fusion	c	liq		1593.	1320.			
Ce <sub>3</sub> Ni		Fusion	c	liq		758.	485.			
Ce <sub>3</sub> Co		Fusion	c	liq		753.	480.			
Ce <sub>2</sub> O <sub>3</sub> ·Cr <sub>2</sub> O <sub>3</sub>		Fusion	c	liq		2713.	2440.			
Ce <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>		Fusion	c	liq		1146.	873.			
Ce <sub>2</sub> (WO <sub>4</sub> ) <sub>3</sub>		Fusion	c	liq		1362.	1089.			
CeAl <sub>2</sub>		Fusion	c	liq		1733.	1460.			
CeAl <sub>4</sub>		Transition	c, II	c, I		1278.	1005.			
Ce <sub>3</sub> Al		Fusion	c	liq		888.	615.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 76-1. Lanthanum (at. no., 57, at. wt., 138.92)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
La		Transition	c, IV	c, III		110.	-163.			
		Transition	c, III	c, II		821.	548.			
		Transition	c, II	c, I		982.	709.			
		Fusion	c, I	liq		1153.	880.			
La <sub>2</sub> O <sub>3</sub>		Fusion	c	liq		2578.	2305.			
LaCl <sub>3</sub>		Fusion	c, α	liq		1143.	870.			
LaBr <sub>3</sub>		Fusion	c	liq		1056.	783.			
LaI <sub>3</sub>		Fusion	c	liq		1034.	761.			
La(NO <sub>3</sub> ) <sub>3</sub> ·6H <sub>2</sub> O		Transition	c, II	c, I		316.	43.			
		Fusion	c, I	liq		339.7	66.5			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 76-2. Lanthanum [at. no., 57; at. wt., 138.92]  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
LaSn <sub>3</sub>		Fusion	c	liq		1408.	1135.			
La <sub>2</sub> Sn		Fusion	c	liq		1694.	1421.			
La <sub>2</sub> Sn <sub>3</sub>		Fusion	c	liq		1465.	1192.			
LaPb		Fusion	c	liq		1518.	1245.			
LaPb <sub>3</sub>		Fusion	c	liq		1363.	1090.			
La <sub>2</sub> Pb		Fusion	c	liq		1590.	1317.			
LaTl		Fusion	c	liq		1455.	1182.			
LaTl <sub>3</sub>		Fusion	c	liq		1369.	1096.			
LaCu <sub>2</sub>		Fusion	c	liq		1107.	834.			
LaCu <sub>6</sub>		Fusion	c	liq		1198.	925.			
LaZn <sub>9</sub>		Fusion	c	liq		1235.	962.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 76-3. Lanthanum (at. no., 57; at. wt., 138.92)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
December 31, 1949

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
LaAg		Fusion	c	liq		1159.	886.			
LaAg <sub>3</sub>		Fusion	c	liq		1228.	955.			
LaAu		Fusion	c	liq		1632.	1359.			
LaAu <sub>2</sub>		Fusion	c	liq		1486.	1213.			
LaAu <sub>3</sub>		Fusion	c	liq		1476.	1203.			
LaNi		Fusion	c	liq		958.	685.			
LaNi <sub>5</sub>		Fusion	c	liq		1598.	1325.			
La <sub>3</sub> Ni		Fusion	c	liq		788.	515.			
La <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>		Fusion	c	liq		1454.	1181.			
LaAl <sub>2</sub>		Fusion	c	liq		1695.	1422.			
LaAl <sub>4</sub>		Transition	c, II	c, I		1089.	816.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 79-1. Plutonium (at. no., 94; at. wt., 239.)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
PuF <sub>3</sub>		Sublimation Vaporization	c	g	0.00036	1329.	1056.	96.9	72.9	
			liq	g	0.079	1562.	1289.	80.9	51.8	
PuCl <sub>3</sub>		Sublimation Vaporization	c	g	0.0017	1033.	760.	65.	63.	
			liq	g	0.0017	1033.	760.	59.7	57.8	
PuBr <sub>3</sub>		Sublimation Vaporization	c	g	0.0019	954.	681.	70.6	74.0	
			liq	g	0.0019	954.	681.	57.2	60.0	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 80-1. Neptunium (at. no., 93; at. wt., 237.) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION March 31, 1950										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}\text{K}$	$^{\circ}\text{C}$			
$\text{NpF}_6$		Fusion	c	liq		326.	53.			
$\text{NpCl}_3$		Fusion	c	liq		1075.	802.			
$\text{NpCl}_4$		Fusion	c	liq		811.	538.			
$\text{NpBr}_4$		Fusion	c	liq		740.	467.			
$\text{NpI}_3$		Fusion	c	liq		1040.	767.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 81-1. Uranium (at. no., 92; at. wt., 238.07) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION March 31, 1950										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
U		Transition Transition Fusion	c, III c, II c, I	c, II c, I liq		938. 1049. 1405.	665. 776. 1132.	0.665 1.165	0.709 1.110	
UF <sub>4</sub>		Sublimation Vaporization Vaporization	c liq liq	g g g	3.8 3.8 760	1309. 1309. 1690.	1036. 1036. 1417.	68.9 63.2 57.5	52.6 48.3 34.1	
UF <sub>6</sub>		Sublimation Sublimation Fusion Vaporization	c c c liq	g g liq g	760 1133 1133 1133	329. 337.2 337.2 337.2	56. 64.0 64.0 64.0	11.8 11.8 4.59 7.2	35.9 35.0 13.61 21.4	8.81

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 81-2. Uranium (at. no., 92; at. wt., 238.07)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
UCl <sub>3</sub>		Sublimation Vaporization	c	g	0.027	1108.	835.	63.	57.	
			liq	g	0.027	1108.	835.	54.	49.	
UCl <sub>4</sub>		Sublimation Vaporization	c	g	15.5	863.	590.	46.3	53.6	
			liq	g	15.5	863.	590.	36.0	41.7	
UBr <sub>3</sub>		Sublimation Fusion Vaporization	c	g	0.0065	1025.	752.	68.6	66.9	
			c	liq	0.0065	1025.	752.	11.0	10.7	
			liq	g	0.0065	1025.	752.	57.6	56.2	
UBr <sub>4</sub>		Sublimation Vaporization Vaporization	c	g	4.78	792.	519.	41.9	52.9	
			liq	g	4.78	792.	519.	43.8	43.8	
			liq	g	760	1039.	766.	31.0	29.8	



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 81-3. Uranium (at. no., 92; at. wt., 238.07)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$UCl_2$		Fusion	c	liq		2700.	2427.			
$U_2Cl_3$		Fusion	c	liq		2700.	2427.			
$(UO_2)Cl_4 \cdot (CH_3NH_2)_2$	ethylene diammonium uranyl chloride	Fusion	c	liq		492.	219.			
$(UO_2)(NO_3)_4 \cdot (CH_3NH_2)_2$	ethylene diammonium uranyl nitrate	Fusion	c	liq		488.	215.			
$U(BH_4)_4$	uranium borohydride	Vaporization	liq	g	0.866	318.	45.	19.	60.	
$U(BH_4)_3(BH_3CH_3)$		Vaporization	liq	g	5.8	318.	45.	14.	45.	
$U(BH_3CH_3)_4$		Vaporization	liq	g	0.30	318.	45.	13.	41.	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Substance		Type of Process	State		Pressure	Temperature		$\Delta H$	$\Delta S$	$\Delta G_p$
Formula	Description		Initial	Final	mm Hg	°K	°C	kcal/mole	cal/deg mole	cal/deg mole
$\text{PaCl}_5$		Fusion	c	liq		574.	301.			

Table 82-1. Protactinium (at. no., 59; at. wt., 140.92)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 83-1. Thorium (at. no., 90; at. wt., 232.12)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Th		Transition Fusion	c, I c	c, II liq		498. 2173.	225. 1840.			
ThO <sub>2</sub>		Fusion	c	liq		3323.	3050.			
ThCl <sub>4</sub>		Sublimation Vaporization Vaporization	c liq liq	g g g	81.3 81.3 760	1043. 1043. 1194.	770. 770. 921.	58.0 37.9 35.0	55.6 36.3 29.3	
ThBr <sub>4</sub>		Sublimation Vaporization Vaporization	c liq liq	g g g	42.7 42.7 760	952. 952. 1130.	679. 679. 857.	45.0 33.8 30.4	47.2 35.5 26.9	
ThI <sub>4</sub>		Fusion Vaporization	c liq	liq g	760	839. 1110.	566. 837.	28.8	25.9	
Th(BH <sub>4</sub> ) <sub>4</sub>		Sublimation	c	g	0.1	412.	139.	23.	56.	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 85-1. Beryllium (at. no., 4; at. wt., 9.013)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Be		Sublimation	c	g	0.034	1556.	1283.	75.2	48.3	
		Fusion	c	liq		1556.	1283.	2.3	1.5	
BeO		Sublimation	c	g	0.00076	2327.	2054.	147.4	63.3	
		Sublimation	c	g		2823.	2550.	145.4	51.5	
		Fusion	c	liq		2823.	2550.			
BeCl <sub>2</sub>		Sublimation	c	g, BeCl <sub>2</sub>	2	678.	405.	29.2	43.1	
		Vaporization	liq	g, BeCl <sub>2</sub>		678.	405.	26.2	38.6	
		Vaporization	liq	g, BeCl <sub>2</sub>		820.	547.	25.	30.	
		Sublimation	c	g, Be <sub>2</sub> Cl <sub>4</sub>		678.	405.	15.4	22.7	
BeBr <sub>2</sub>		Sublimation	c	g, BeBr <sub>2</sub>	405	761.	488.	27.4	36.0	
		Fusion	c	liq		761.	488.			
		Sublimation	c	g, Be <sub>2</sub> Br <sub>4</sub>		761.	488.	13.2	17.4	
BeI <sub>2</sub>		Sublimation	c	g, BeI <sub>2</sub>	543	753.	480.	24.5	32.5	
		Fusion	c	liq		753.	480.			
		Sublimation	c	g, Be <sub>2</sub> I <sub>4</sub>		753.	480.	27.6	36.6	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 85-2. Beryllium (at. no., 4; at. wt., 9.013)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$\text{BeO} \cdot 3\text{Be}(\text{C}_2\text{H}_3\text{O}_2)_2$		Sublimation	c, II	g	0.096	421.	148.	31.7	75.3	
		Sublimation	c, I	g	0.096	421.	148.	27.1	64.4	
		Transition	c, II	c, I	0.096	421.	148.	4.6	10.9	
		Fusion	c, I	liq	112	559.9	286.7	6.5	11.6	
$\text{BeCl}_2 \cdot 2(\text{C}_2\text{H}_5)_2\text{O}$		Fusion	c	liq		317.	44.			
$\text{BeBr}_2 \cdot 2(\text{C}_2\text{H}_5)_2\text{O}$		Fusion	c	liq		322.	49.			
$\text{Be}_2\text{H}_8$	beryllium borohydride	Sublimation	c	g	29.6	315.0	41.8	14.8	47.0	
$\text{Be}_6\text{H}_8 \cdot \text{N}(\text{CH}_3)_3$		Vaporization	liq	g	3.7	373.8	100.6	13.4	35.8	
$3\text{BeO} \cdot 2\text{ZrO}_2$		Fusion	c	liq		308.	35.			
$\text{BeO} \cdot \text{Al}_2\text{O}_3$		Fusion	c	liq		2808.	2535.			
		Fusion	c	liq		2143.	1870.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 86-1. Magnesium (at. no., 12; at. wt., 24.32)

HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Mg		Fusion	c	liq	3	923.	650.	2.2	2.4	
		Vaporization	liq	g	760	1393.	1120.	31.5	22.6	
MgO		Fusion	c	liq		3173.	2800.	18.5	5.8	
MgF <sub>2</sub>		Fusion	c	liq		1536.	1263.	13.9	9.0	
		Vaporization	liq	g	760	2500.	2227.	65.	26.	
MgCl <sub>2</sub>		Fusion	c	liq		987.	714.	10.3	10.4	
		Vaporization	liq	g	760	1691.	1418.	32.7	19.3	
MgCl <sub>2</sub> ·6H <sub>2</sub> O		Fusion	c	liq		390.	117.	8.2	21.0	
MgBr <sub>2</sub>		Fusion	c	liq		984.	711.	8.3	8.4	
MgBr <sub>2</sub> ·6H <sub>2</sub> O		Fusion	c	liq		445.6	172.4			
MgSO <sub>4</sub>		Fusion	c	liq		1400.	1127.	3.5	2.5	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 86-2. Magnesium (at. no., 12; at. wt., 24.32)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Mg_3N_2$		Transition	c, II	c, I		1061.	788.	0.26	0.24	
		Transition	c, III	c, II		823.	550.	0.22	0.27	
$Mg(NO_3)_2 \cdot 2H_2O$		Fusion	c	liq		403.	130.			
$Mg(NO_3)_2 \cdot 6H_2O$		Fusion	c	liq		363.1	89.9	9.8	27.0	
$Mg_2P_2O_7$		Transition	c, II	c, I		341.	68.			
		Fusion	c, I	liq		1653.	1380.			
$Mg_3(PO_4)_2$		Fusion	c	liq		1457.	1184.	11.	9.	
$Mg_3As_2$		Fusion	c	liq		1073.	800.			
$Mg_3Sb_2$		Transition	c, II	c, I		1203.	930.			
$3Mg(NO_3)_2 \cdot 28H_2O$		Fusion	c	liq		344.	71.			
$Mg(CH_3COO)_2 \cdot 4H_2O$		Fusion	c	liq		341.	68.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 88-3. Magnesium (at. no., 12; at. wt., 24.32)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		O <sub>K</sub>	O <sub>C</sub>			
$MgBr_2 \cdot 6HCOOH$		Fusion	c	liq		361.	88.			
$MgBr_2 \cdot 6CH_3OH$		Fusion	c	liq		463.	190.			
$MgBr_2 \cdot 6CH_3COOH$		Fusion	c	liq		385.	112.			
$MgBr_2 \cdot 6C_2H_5OH$		Fusion	c	liq		381.7	108.5			
$MgI_2 \cdot 6CH_3OH$		Fusion	c	liq		483.	210.			
$MgI_2 \cdot 6CH_3COOH$		Fusion	c	liq		415.	142.			
$MgI_2 \cdot 6C_2H_5OH$		Fusion	c	liq		419.7	146.5			
$MgBr_2 \cdot 3CH_3CN$		Fusion	c	liq		405.	132.			
$MgBr_2 \cdot 6CH_3CONH_2$		Fusion	c	liq		442.	169.			
$MgI_2 \cdot 6CH_3CONH_2$		Fusion	c	liq		450.	177.			



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 86-4. Magnesium (at. no., 12; at. wt., 24.32)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Mg_2Si$			Fusion	c	liq		1375.	1102.			
$MgSiO_3$			Fusion	c	liq		1798.	1525.	14.7	8.2	
$Mg_2SiO_4$			Fusion	c	liq		2158.	1885.			
$Mg_2Ge$			Fusion	c	liq		1388.	1175.			
$Mg_5Ti_2$			Fusion	c	liq		682.	409.			
$MgZn_2$			Fusion	c	liq		863.	590.	8.3	9.6	
$MgCu_2$			Fusion	c	liq		1092.	819.			
$Mg_2Cu$			Fusion	c	liq		841.	568.			
$MgCuSb$			Fusion	c	liq		1203.	830.			
$AgMg$			Fusion	c	liq		1093.	820.			
$Ni_2Mg$			Fusion	c	liq		1418.	1145.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 86-5. Magnesium (at. no., 12; at. wt., 24.32)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$MgO \cdot Fe_2O_3$		Fusion	c	liq		2033.	1760.			
$MgO \cdot 2TiO_2$		Fusion	c	liq		1953.	1680.			
$2MgO \cdot TiO_2$		Fusion	c	liq		2108.	1835.			
$MgO \cdot ZrO_2$		Fusion	c	liq		2293.	2120.			
$MgO \cdot B_2O_3$		Fusion	c	liq		1464.	1191.			
$2MgO \cdot B_2O_3$		Fusion	c	liq		1653.	1380.			
$3MgO \cdot B_2O_3$		Fusion	c	liq		1673.	1400.			
$MgCl_2 \cdot 5MgO \cdot 7B_2O_3$	boracite	Transition	c, II	c, I		538.	265.			
$Mg_2Al_3$		Fusion	c	liq		723.	450.			
$Mg_3Al_2$		Fusion	c	liq		728.	455.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 86-6. Magnesium (at. no., 12; at. wt., 24.32)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$Mg_4Al_3$		Fusion	c	liq		738.	465.			
$MgO \cdot Al_2O_3$		Fusion	c	liq		2383.	2110.			
$MgPr$		Fusion	c	liq		1159.	886.			
$Mg_3Pr$		Fusion	c	liq		1025.	752.			
$MgCe$		Fusion	c	liq		1003.	730.			
$MgCe_4$		Fusion	c	liq		893.	620.			
$Mg_3Ce$		Fusion	c	liq		1063.	790.			
$MgLa$		Fusion	c	liq		1016.	743.			
$Mg_3La$		Fusion	c	liq		1071.	798.			
$MgO \cdot La_2O_3$		Fusion	c	liq		2303.	2030.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 87-1. Calcium (at. no., 20; at. wt. 40.08)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Formula	Substance	Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
			Initial	Final		°K	°C			
Ca		Transition	c, II	c, I	0.35	723.	450.	0.2	0.3	-0.5
		Sublimation	c, I	g		1123.	850.			
		Fusion	c, I	liq		1123.	850.			
CaO		Fusion	c	liq		2873.	2600.	12.	4.2	
CaF <sub>2</sub>		Transition	c, II	c, I		1424.	1151.	1.14	0.80	
		Fusion	c, I	liq		1691.	1418.			
		Vaporization	liq	g		2145.	1872.			
CaCl <sub>2</sub>		Sublimation	c	g	0.018	934.6	661.4	54.	58.	
		Fusion	c	liq		1055.	782.			
CaCl <sub>2</sub> ·6H <sub>2</sub> O		Fusion	c	liq		303.4	30.2			
4CaCl <sub>2</sub> ·CaO		Fusion	c	liq		1112.	839.			
CaBr <sub>2</sub> ·6H <sub>2</sub> O		Fusion	c	liq		307.4	34.2			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 87-2. Calcium (at. no., 20; at. wt., 40.08)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta G$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
CaSO <sub>4</sub>		Transition Fusion	c, II c, I	c, I liq		1466. 1573.	1193. 1400.	6.7	4.3	
Ca <sub>3</sub> N <sub>2</sub>		Fusion	c	liq		1468.	1195.			
Ca(NO <sub>3</sub> ) <sub>2</sub>		Fusion	c	liq		834.	561.	5.1	6.1	
Ca(NO <sub>3</sub> ) <sub>2</sub> ·3H <sub>2</sub> O		Fusion	c	liq		324.3	51.1			
Ca(NO <sub>3</sub> ) <sub>2</sub> ·4H <sub>2</sub> O		Fusion	c, II	liq		312.9	39.7			
		Fusion	c, I	liq		315.8	42.6			
CaO·P <sub>2</sub> O <sub>5</sub>		Fusion	c	liq		1248.	975.			
2CaO·P <sub>2</sub> O <sub>5</sub>		Fusion	c	liq		1573.	1300.			
3CaO·P <sub>2</sub> O <sub>5</sub>		Transition Fusion	c, II c, I	c, I liq		1373. 2003.	1100. 1730.	3.7	2.7	
4CaO·P <sub>2</sub> O <sub>5</sub>		Fusion*	c	liq		1973.	1700.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}\text{K}$	$^{\circ}\text{C}$			
$\text{Ca}_3(\text{AsO}_4)_2$		Fusion	c	liq		1723.	1450.			
$\text{Ca}_3\text{Bi}_2$		Fusion	c	liq		1201.	928.			
$\text{CaC}_2$		Transition	c, II	c, I		720.	447.	1.33	1.85	-1.20
$\text{CaC}_2 \cdot \text{CaO}$		Fusion	c	liq		2243.	1970.			
$\text{CaCl}_2 \cdot 3\text{CH}_3\text{OH}$		Fusion	c	liq		450.	177.			
$\text{CaCl}_2 \cdot 3\text{C}_2\text{H}_5\text{OH}$		Fusion	c	liq		370.	97.			
$\text{CaSi}_2$		Fusion	c	liq		1493.	1220.			
$\text{CaSiO}_3$		Transition Fusion	c, II c, I	c, I liq		1463. 1803.	1190. 1530.			
$\text{Ca}_2\text{SiO}_4$		Transition Transition Fusion	c, III c, II c, I	c, II c, I liq		948. 1673. 2393.	675. 1400. 2120.	0.35 0.77	0.37 0.46	0.58

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

SERIES II

Washington, D.C.

Table 87-4. Calcium (at. no., 20; at. wt., 40.08)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
CaSn <sub>3</sub>		Fusion	c	liq		897.	624.			
CaPb <sub>3</sub>		Fusion	c	liq		923.	650.			
Ca <sub>2</sub> Pb		Fusion	c	liq		1423.	1150.			
CaTi		Fusion	c	liq		1243.	970.			
CaCl <sub>2</sub> · TiCl		Fusion	c	liq		956.	683.			
CaZn <sub>10</sub>		Fusion	c	liq		1003.	730.			
Ca <sub>2</sub> Zn <sub>3</sub>		Fusion	c	liq		963.	690.			
CaCd <sub>3</sub>		Fusion	c	liq		892.	619.			
CaCu <sub>4</sub>		Fusion	c	liq		1208.	935.			
CaAg		Fusion	c	liq		938.	665.			
CaAg <sub>2</sub>		Fusion	c	liq		868.	595.			
CaAg <sub>3</sub>		Fusion	c	liq		998.	725.			
CaAg <sub>4</sub>		Fusion	c	liq		956.	683.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 87-5. Calcium (at. no., 20; at. wt., 40.08)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$CaAu_2$		Fusion	c	liq		1137.	864.			
$CaAu_4$		Fusion	c	liq		1153.	880.			
$CaO \cdot 9FeO$		Fusion	c	liq		1583.	1310.			
$2CaO \cdot 3Fe_2O_3$		Fusion	c	liq		1673.	1400.			
$CaO \cdot FeO \cdot SiO_2$		Fusion	c	liq		1481.	1208.			
$CaCrO_4$		Fusion	c	liq		2433.	2160.			
$CaTiO_3$		Transition	c, II	c, I		1530.	1257.	0.55	0.36	-0.23
$CaO \cdot TiO_2 \cdot SiO_2$		Fusion	c	liq		1658.	1385.			
$CaO \cdot ZrO_2$		Fusion	c	liq		2618.	2345.			
$CaO \cdot B_2O_3$		Fusion	c	liq		1435.	1162.	17.67	12.31	17.06
$CaO \cdot 2B_2O_3$		Fusion	c	liq		1260.	987.	27.06	21.48	31.90
$2CaO \cdot B_2O_3$		Transition Fusion	c, II c, I	c, I liq		804. 1585.	531. 1312.	1.1 24.09	1.4 15.20	2.87 12.11



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

SERIES II

Washington, D.C.

Table 87-6. Calcium (at. no., 20; at. wt., 40.08)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$3CaO \cdot B_2O_3$		Fusion	c	liq		1760.	1487.			
$5CaO \cdot B_2O_3 \cdot SiO_2$		Fusion	c	liq		1690.	1417.	35.49	20.16	19.64
$CaAl_2$		Fusion	c	liq		1353.	1080.			
$CaAl_3$		Fusion	c	liq		968.	695.			
$CaO \cdot Al_2O_3$		Fusion	c	liq		1873.	1600.			
$12CaO \cdot 7Al_2O_3$		Fusion	c	liq		1723.	1450.			
$CaBr_2 \cdot AlBr_3$		Fusion	c	liq		579.	306.			
$CaO \cdot Al_2O_3 \cdot 2SiO_2$		Fusion	c	liq		1823.	1550.	29.4	16.1	
$2CaO \cdot Al_2O_3 \cdot SiO_2$		Fusion	c	liq		1857.	1584.			
$CaMg_2$		Fusion	c	liq		1003.	730.			
$CaMg(SiO_3)_{1/2}$	diopside	Fusion	c	liq		1665.	1392.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 88-1. Strontium (at. no., 38; at. wt., 87.63)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}\text{K}$	$^{\circ}\text{C}$			
Sr		Fusion	c	liq		1043.	770.	2.2	2.1	
		Vaporization	liq	g	760	1657.	1384.	33.8	20.4	
SrO		Fusion	c	liq		2688.	2415.			
SrF <sub>2</sub>		Fusion	c	liq		1673.	1400.	4.3	2.6	
SrCl <sub>2</sub>		Fusion	c	liq		1148.	875.	4.1	3.6	
4SrCl <sub>2</sub> ·SrO		Fusion	c	liq		1268.	995.			
SrCl <sub>2</sub> ·SrF <sub>2</sub>		Fusion	c	liq		1235.	962.			
SrBr <sub>2</sub>		Fusion	c	liq		916.	643.	4.8	5.2	
SrI <sub>2</sub>		Fusion	c	liq		788.	515.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

Table 88-2. Strontium (at. no., 38; at. wt., 87.63)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		°K	°C			
$\text{SrSO}_4$			Transition Fusion	c, II c, I	c, I liq		1425. 1878.	1152. 1605.			
$\text{Sr}(\text{NO}_3)_2$			Fusion	c	liq		918.	645.			
$\text{Sr}(\text{AsO}_4)_2$			Fusion	c	liq		1903.	1630.			
$\text{Sr}_3(\text{PO}_4)_2$			Fusion	c	liq		2040.	1767.	18.5	9.1	
$\text{SrCO}_3$			Transition Fusion	c, II c, I	c, I liq		1198. 1770.	925. 1497.			
$\text{SrO} \cdot \text{SiO}_2$			Fusion	c	liq		1848.	1575.			
$2\text{SrO} \cdot \text{SiO}_2$			Fusion	c	liq		1861.	1588.			
$\text{SrPb}_3$			Fusion	c	liq		949.	676.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards Washington, D.C.

Table 88-3. Strontium (at. no., 38; at. wt., 87.63)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
SrAg		Fusion	c	liq		953.	680.			
SrAg <sub>11</sub>		Fusion	c	liq		1054.	781.			
Sr <sub>3</sub> Ag <sub>2</sub>		Fusion	c	liq		939.	666.			
Sr <sub>3</sub> Ag <sub>5</sub>		Fusion	c	liq		1030.	757.			
SrO·B <sub>2</sub> O <sub>3</sub>		Fusion	c	liq		1373.	1100.			
SrO·2B <sub>2</sub> O <sub>3</sub>		Fusion	c	liq		1213.	940.			
2SrO·B <sub>2</sub> O <sub>3</sub>		Fusion	c	liq		1403.	1130.			
SrO·Al <sub>2</sub> O <sub>3</sub>		Fusion	c	liq		2283.	2010.			
SrMg <sub>2</sub>		Fusion	c	liq		953.	680.			
SrMg <sub>9</sub>		Fusion	c	liq		876.	603.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 89-1. Barium (at. no., 56; at. wt., 137.36)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Ba	Transition	c, II		c, I		648.	375.			
	Sublimation	c, I		g	0.0063	977.	704.	41.1	42.1	
	Fusion	c, I		liq		977.	704.			
	Vaporization	liq		g	760	1911.	1638.	35.7	18.7	
BaO	Sublimation	c		g	0.00076	1650.	1377.	89.	54.	
	Fusion	c		liq		2190.	1917.			
Ba(OH) <sub>2</sub>	Fusion	c		liq		681.	408.	3.4	5.0	
BaF <sub>2</sub>	Fusion	c		liq		1593.	1320.	3.0	1.9	
	Vaporization	liq		g	42	2072.	1799.	83.	40.	
BaCl <sub>2</sub>	Transition	c, II		c, I		1198.	925.			
	Fusion	c, I		liq		1235.	962.	5.4	4.4	
	Vaporization	liq		g	6.3	1462.	1189.	57.	39.	
Ba(ClO <sub>4</sub> ) <sub>2</sub>	Transition	c, II		c, I		557.	284.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 89-2. Barium (at. no., 56; at. wt., 137.36)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}\text{K}$	$^{\circ}\text{C}$			
BaBr <sub>2</sub>		Fusion	c	liq		1123.	850.	6.	5.	
BaI <sub>2</sub>		Fusion	c	liq		1013.	740.			
BaSO <sub>4</sub>		Transition	c, II	c, I		1422.	1149.			
		Fusion	c, I	liq		1623.	1350.	9.7	6.0	
Ba(NO <sub>3</sub> ) <sub>2</sub>		Fusion	c	liq		868.7	595.5	6.	7.	
Ba <sub>3</sub> (PO <sub>4</sub> ) <sub>2</sub>		Fusion	c	liq		2000.	1727.	18.6	9.3	
Ba <sub>3</sub> (AsO <sub>4</sub> ) <sub>2</sub>		Fusion	c	liq		1878.	1605.			
BaCO <sub>3</sub>		Transition	c, II	c, I		1083.	810.	3.55	3.28	7.93
BaSiO <sub>3</sub>		Fusion	c	liq		1878.	1605.			
BaO·2SiO <sub>2</sub>		Fusion	c	liq		1698.	1425.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 89-3. Barium (at. no., 56; at. wt., 137.36)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
BaAg <sub>4</sub>		Fusion	c	liq		1002.	729.			
Ba <sub>2</sub> Ag <sub>3</sub>		Fusion	c	liq		1121.	848.			
BaO·Fe <sub>2</sub> O <sub>3</sub>		Fusion	c	liq		1663.	1390.			
Ba(MoO <sub>4</sub> )		Fusion	c	liq		1753.	1480.			
2BaO·V <sub>2</sub> O <sub>5</sub>		Fusion	c	liq		1136.	863.			
BaTiO <sub>3</sub>		Transition	c, II	c, I		385.8	112.6			
BaO·ZrO <sub>2</sub>		Fusion	c	liq		2973.	2700.			
BaO·B <sub>2</sub> O <sub>3</sub>		Fusion	c	liq		1323.	1050.			
BaO·2B <sub>2</sub> O <sub>3</sub>		Fusion	c	liq		1083.	810.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table B9-4. Barium (at. no., 56; at. wt., 137.36)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$BaO \cdot 4B_2O_3$		Fusion	c	liq		1023.	750.			
$2BaO \cdot B_2O_3$		Fusion	c	liq		1263.	990.			
$3BaO \cdot B_2O_3$		Fusion	c	liq		1593.	1320.			
$BaO \cdot 3B_2O_3$		Fusion	c	liq		1013.	740.			
$BaO \cdot Al_2O_3$		Fusion	c	liq		2273.	2000.			
$BaMg_2$		Fusion	c	liq		880.	607.			
$BaMg_3$		Fusion	c	liq		980.	707.			
$BaBeF_4$		Fusion	c	liq		1268.	995.			
$BaBe_2F_6$		Fusion	c	liq		1173.	900.			



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

Table 90-1. Radium (at. no., 88; at. wt., 226.05) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION March 31, 1950										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Ra		Fusion	c	liq			973. 700.			
RaCl <sub>3</sub>		Fusion	c	liq			1173. 900.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 91-1. Lithium (at. no., 3; at. wt., 6.940)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Li		Transition Fusion Vaporization	c, II c, I liq	c, I liq equil g		77. 353. 1599.	-196. 180. 1326.	1.	3.	
LiOH		Fusion	c	liq		735.	462.			
LiF		Fusion Vaporization	c liq	liq g	760	1118. 1954.	845. 1681.	2.4 51.0	2.1 26.1	
LiCl		Sublimation Fusion Vaporization	c c liq	g liq g	0.014 760	883. 883. 1655.	610. 610. 1382.	46.2 3.2 36.0	52.3 3.6 21.8	
LiClO <sub>3</sub>		Transition Transition Fusion	c, III c, II c, I	c, II c, I liq		314.7 372. 400.8	41.5 99. 127.6			
LiClO <sub>3</sub> ·3H <sub>2</sub> O		Fusion	c	liq		281.	8.			
LiClO <sub>4</sub>		Fusion	c	liq		509.	236.			
LiBr		Fusion Vaporization	c liq	liq g	760	823. 1583.	550. 1310.	2.9 35.4	3.5 22.4	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 91-2. Lithium (at. no., 3; at. wt., 6.940) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION March 31, 1950										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
LiI		Fusion Vaporization	c liq	liq g		722. 1444.	449. 1171.	40.8	28.2	
Li <sub>2</sub> SO <sub>4</sub>		Transition Fusion	c, II c, I	c, I liq		848. 1132.	575. 859.	6.8 3.0	8.0 2.6	
LiNO <sub>3</sub>		Fusion	c	liq		527.	254.	6.1	11.6	0.45
LiNO <sub>3</sub> ·3H <sub>2</sub> O		Fusion	c	liq		303.1	29.9	8.7	29.	
Li <sub>3</sub> PO <sub>4</sub>		Fusion	c	liq		1130.	857.			
Li <sub>3</sub> Bi		Fusion	c	liq		1418.	1145.			
Li <sub>2</sub> CO <sub>3</sub>		Fusion	c	liq		1008.	735.			
LiC <sub>2</sub> H <sub>3</sub> O <sub>2</sub>	lithium acetate	Fusion	c	liq		553.	280.			
LiC <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ·2H <sub>2</sub> O	lithium acetate dihydrate	Fusion	c	liq		331.	58.			
Li <sub>2</sub> SiO <sub>3</sub>		Fusion	c	liq		1461.	1188.	7.2	4.9	
Li <sub>4</sub> SiO <sub>4</sub>		Fusion	c	liq		1523.	1250.	7.4	4.8	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

SERIES II

Washington, D.C.

Table 91-3. Lithium (at. no., 3; at. wt., 6.940)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Li_2GeO_3$		Fusion	c	liq		1513.	1240.			
$Li_4GeO_4$		Fusion	c	liq		1568.	1295.			
$LiReO_4$		Fusion	c	liq		699.	426.			
$LiCl \cdot CoCl_2$		Fusion	c	liq		836.	565.			
$Li_2MnO_4$		Fusion	c	liq		978.	705.	4.2	4.3	
$Li_2WO_4$		Fusion	c	liq		1015.	742.	7.	7.	
$Li_2W_2O_7$		Transition Fusion	c, II c, I	c, I liq		960. 1018.	687. 745.			
$LiVO_3$		Fusion	c	liq		894.	621.			
$2LiSiO_4 \cdot 3ZrSiO_4$		Fusion	c	liq		1425.	1152.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 91-4. Lithium [at. no., 3; at. wt., 6.940]  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$LiBO_2$		Fusion	c	liq		1118.	845.	6.	5.	
$Li_2O \cdot B_2O_3$		Fusion	c	liq		1033.	760.			
$Li_2O \cdot 2B_2O_3$		Fusion	c	liq		1198.	925.			
$Li_2O \cdot 3B_2O_3$		Fusion	c	liq		1023.	750.			
$Li_2O \cdot 4B_2O_3$		Fusion	c	liq		1003.	730.			
$Li_2O \cdot 5B_2O_3$		Fusion	c	liq		953.	680.			
$3LiF \cdot AlF_3$		Fusion	c	liq		1063.	790.			
$LiCl \cdot AlCl_3$		Fusion	c	liq		417.	144.			
$LiBr \cdot AlBr_3$		Fusion	c	liq		470.	197.			
$LiBr \cdot 7AlBr_3$		Fusion	c	liq		388.	115.			
$Li_2BeF_4$		Fusion	c	liq		748.	475.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

SERIES II

Washington, D.C.

Table 92-1. Sodium (at. no., 11; at. wt., 22.997)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Na		Fusion Vaporization	c liq	liq g, equil	760	371. 1162.	98. 889.	0.63	1.7	0.51
Na <sub>2</sub> O <sub>2</sub>		Fusion	c	liq		733.	460.			
NaOH		Transition Fusion	c, II c, I	c, I liq		572.8 593.	299.6 320.	1. 1.7	2. 2.9	1.2
NaOH·H <sub>2</sub> O		Fusion	c	liq		337.4	64.2			
NaOH·3 1/2H <sub>2</sub> O		Fusion	c	liq		288.7	15.5			
NaF		Sublimation Fusion Vaporization	c c liq	g liq g	0.5 0.5 760	1268. 1268. 1977.	995. 995. 1704.	63. 7.8 50.	49. 6.2 25.	0.65
NaCl		Sublimation Fusion Vaporization	c c liq	g liq g	0.5 0.5 760	1081. 1081. 1738.	808. 808. 1465.	51.5 6.8 40.8	47.6 6.3 23.5	0.8
NaClO <sub>3</sub>		Fusion	c	liq		528.	255.	5.4	10.2	4.9
NaClO <sub>4</sub>		Transition	c, II	c, I		581.	308.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 92-2. Sodium (at. no., 11; at. wt., 22.997)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
NaBr		Sublimation	c	g	0.40	1023.	750.	49.3	48.2	
		Fusion	c	liq		1023.	750.	6.1	5.9	
		Vaporization	liq	g		1665.	1392.	38.7	23.2	
NaI		Fusion	c	liq	760	935.	662.	5.2	5.6	
		Vaporization	liq	g		1577.	1304.	38.2	24.2	
Na <sub>2</sub> S		Fusion	c	liq		1223.	950.	1.6	1.3	
Na <sub>2</sub> S <sub>2</sub>		Fusion	c	liq		718.	445.			
Na <sub>2</sub> S <sub>3</sub>		Fusion	c	liq		593.	320.			
Na <sub>2</sub> S <sub>4</sub>		Fusion	c	liq		548.	275.			
Na <sub>2</sub> S <sub>5</sub>		Fusion	c	liq		493.	220.			
Na <sub>4</sub> S <sub>3</sub>		Fusion	c	liq		1045.	772.			
Na <sub>4</sub> S <sub>5</sub>		Fusion	c	liq		618.	345.			
Na <sub>4</sub> S <sub>7</sub>		Fusion	c	liq		568.	295.			
Na <sub>4</sub> S <sub>9</sub>		Fusion	c	liq		463.	210.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 92-3. Sodium (at. no., 11; at. wt., 22.997)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance	Formula	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Na_2SO_4$			Transition Fusion	c, II	c, I		513.	240.	5.8	5.0	
				c, I	liq		1163.	890.			
				c, I	liq		321.7	48.5	5.6	17.	20.4
$NaSH$			Transition	c, II	c, I		358.	85.	0.7	2.	
$NaHSO_4$			Fusion	c	liq		455.	182.			
$NaSeH$			Transition	c, II	c, I		359.	86.	0.7	2.	
$Na_2Te$			Transition	c, II	c, I		725.	452.			
$NaNO_2$			Transition	c, II	c, I		435.	162.			
$NaNO_3$			Transition Fusion	c, II c, I	c, I liq		548. 583.	275. 310.	3.8	6.5	0.3
$NaNH_2$			Fusion	c	liq		481.	208.			
$NaPO_3$			Fusion	c	liq		898.	625.			
$Na_4P_2O_7$			Fusion	c	liq		1243.	970.	14.	11.	



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 92-4. Sodium (at. no., 11; at. wt., 22.997)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$NaH_2PO_3$		Fusion	c	liq		345.	42.			
$NaBi$		Fusion	c	liq		748.	445.			
$Na_3Bi$		Fusion	c	liq		1046.	775.			
$Na_2CO_3$		Transition	c, IV	c, III		629.	356.			
		Transition	c, III	c, II		759.	486.			
		Transition	c, II	c, I		891.	618.			
		Fusion	c, I	liq		1127.	854.	8.	7.	
$NaCHO_2$	sodium formate	Fusion	c	liq		573.	300.			
$NaC_2H_3O_2$	sodium acetate	Transition	c, II	c, I		471.	198.			
$NaCN$		Transition	c, III	c, II		172.1	-101.1	0.15	0.87	
		Transition	c, II	c, I		288.5	15.3	0.70	2.43	
		Fusion	c, I	liq		835.	562.	4.	5.	
		Vaporization	liq	g	760	1770.	1497.	37.	21.	
$NaCNS$		Fusion	c	liq		596.	323.	4.4	7.4	
$Na_2SiO_3$		Fusion	c	liq		1362.	1089.	12.5	9.2	-1.07
$Na_5Si_2O_5$		Fusion	c	liq		1147.	874.	8.5	7.4	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$\text{Na}_2\text{GeO}_3$		Fusion	c	liq		1353.	1080.			
$\text{Na}_2\text{Ge}_2\text{O}_5$		Fusion	c	liq		1072.	799.			
$\text{Na}_2\text{Ge}_4\text{O}_9$		Fusion	c	liq		1325.	1052.			
$\text{NaSn}$		Transition Fusion	c, II c, I	c, I liq		756. 849.	483. 576.			
$\text{Na}_2\text{Sn}$		Fusion	c	liq		750.	477.			
$\text{NaPb}$		Fusion	c	liq		641.	368.			
$\text{Na}_2\text{Pb}$		Fusion	c	liq		678.	405.			
$\text{Na}_2\text{Pb}_5$		Fusion	c	liq		592.	319.			
$\text{Na}_4\text{Pb}$		Fusion	c	liq		653.	380.			
$\text{Na}_5\text{Pb}_2$		Fusion	c	liq		673.	400.			
$\text{NaCd}_2$		Fusion	c	liq		658.	385.			
$\text{NaCd}_5$		Fusion	c	liq		633.	360.			
$\text{NaReO}_4$		Fusion	c	liq		687.	414.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 92-6. Sodium (at. no., 11; at. wt., 22.997)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta f$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Na_2O \cdot Fe_2O_3$		Fusion	c	liq		1618.	1345.			
$Na_2CrO_4$		Transition Fusion	c, II c, I	c, I liq		686. 1065.	413. 792.			
$Na_2CrO_4 \cdot 10H_2O$		Fusion	c	liq		293.1	19.9	15.	51.	
$Na_2Cr_2O_7$		Fusion	c	liq		629.9	356.7			
$Na_2MoO_4$		Transition Fusion	c, II c, I	c, I liq		713. 960.	440. 687.	14.6 3.6	20.5 3.8	
$Na_2WO_4$		Transition Transition Fusion	c, III c, II c, I	c, II c, I liq		860.8 862.0 968.7	587.6 588.8 695.5	7.4 1.0 5.7	8.6 1.2 5.9	
$Na_2TiO_3$		Transition Fusion	c, II c, I	c, I liq		560. 1303.	287. 1030.	0.4 16.8	0.7 12.9	-1.31 -1.2
$Na_2Ti_2O_5$		Fusion	c	liq		1258.	985.			
$Na_2Ti_3O_4$		Fusion	c	liq		1401.	1128.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$NaBO_2$		Fusion	c	liq		1239.	966.			
$NaO \cdot B_2O_3$		Fusion	c	liq		1238.	965.			
$Na_2O \cdot B_2O_3 \cdot 8H_2O$		Fusion	c	liq		326.7	53.5			
$Na_2O \cdot 2B_2O_3$		Fusion	c, I	liq		1015.	742.			
$Na_2O \cdot 4B_2O_3$		Fusion	c	liq		1089.	816.			
$2Na_2O \cdot B_2O_3$		Fusion	c	liq		898.	625.			
$Na_3AlF_6$		Transition Fusion	c, II c, I	c, I liq		838. 1273.	565. 1000.	1.5 20.8	1.8 16.3	5.5 12.8
$NaAlSiO_4$		Transition Fusion	c, II c, I	c, I liq		1521. 1796.	1248. 1526.			
$Na_2SO_4 \cdot MgSO_4$		Transition Transition Fusion	c, III c, II c, I	c, II c, I liq		563. 783. 943.	290. 510. 670.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 93-1. Potassium (at. no., 19; at. wt., 39.076)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
K		Fusion Vaporization	c liq	liq g, equil.	760	336.6 1030.	63.4 757.	0.57	1.69	
KO <sub>2</sub>		Transition	c, II	c, I		197.7	-75.5			
K <sub>2</sub> O <sub>2</sub>		Fusion	c	liq		763.	490.			
K <sub>2</sub> O <sub>3</sub>		Fusion	c	liq		703.	430.			
KOH		Transition Fusion Vaporization	c, II c, I liq	c, I liq g	760	522. 673. 1600.	249. 400. 1327.	1.52 1.8 30.8	2.91 2.6 19.3	
KOH·H <sub>2</sub> O		Fusion	c	liq		416.	143.			
KF		Fusion Vaporization	c liq	liq g	760	1129. 1775.	856. 1502.	6.8 41.3	6.0 23.3	1.5
KF·4H <sub>2</sub> O		Fusion	c	liq		290.7	17.5			
KHF <sub>2</sub>		Transition Fusion	c, II c, I	c, I liq		469.2 511.9	196.0 238.7	2.659 1.58	5.670 3.09	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 93-2. Potassium (at. no., 19; at. wt., 39.096)

HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
KF·2HF		Fusion	c	liq		344.9	71.7			
KF·3HF		Fusion	c	liq		339.0	65.8			
KF·4HF		Fusion	c	liq		345.2	72.0			
2KF·5HF		Fusion	c	liq		337.5	64.3			
KCl		Sublimation	c	g	0.40	1045.	772.	49.5	47.4	
		Fusion	c	liq	0.40	1045.	772.	6.1	5.8	0.61
		Vaporization	liq	g	760	1680.	1407.	38.8	23.1	
KClO <sub>4</sub>		Transition	c, II	c, I		572.7	299.5	3.29	5.74	
KBr		Sublimation	c	g	0.3	1008.	735.	48.9	48.5	
		Fusion	c	liq	0.3	1008.	735.	7.	7.	
		Vaporization	liq	g	760	1656.	1383.	37.1	22.4	
KI		Sublimation	c	g	0.36	958.	685.	47.2	49.3	
		Fusion	c	liq		958.	685.			
		Vaporization	liq	g	760	1597.	1324.	34.7	21.7	
KI <sub>3</sub>		Fusion	c	liq		352.	79.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

National Bureau of Standards

## SERIES II

Washington, D.C.

Table 93-3. Potassium (at. no., 19; at. wt., 39.096) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION March 31, 1950										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
KIO <sub>3</sub>		Fusion	c	liq		833.	560.			
KIBr <sub>2</sub>		Fusion	c	liq		331.	58.			
K <sub>2</sub> S		Transition	c, II	c, I		419.6	146.4	0.085	0.20	
K <sub>2</sub> S <sub>5</sub>		Fusion	c	liq		479.	206.			
K <sub>2</sub> SO <sub>4</sub>		Transition	c, II	c, I		856.	563.	1.94	2.27	-4.79
		Fusion	c, I	liq		1342.	1069.	8.76	6.53	-6.49
K <sub>2</sub> SO <sub>4</sub> ·3K <sub>2</sub> S		Fusion	c	liq		1033.	760.			
KSH		Transition	c, II	c, I		453.	180.	0.55	1.2	
KHSO <sub>4</sub>		Transition	c, III	c, II		437.4	164.2	0.49	1.12	
		Transition	c, II	c, I		453.7	180.5	0.095	0.21	
		Fusion	c	liq		491.8	218.6			
K <sub>2</sub> SO <sub>4</sub> ·3H <sub>2</sub> SO <sub>4</sub>		Fusion	c	liq		364.7	91.5			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 93-4. Potassium (at. no., 19; at. wt., 39.096)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
KSeH		Transition	c, II	c, I		445.	175.	0.45	1.0	
KN <sub>3</sub>		Fusion	c	liq		628.	354.			
KNO <sub>2</sub>		Fusion	c	liq		713.	440.			
KNO <sub>3</sub>		Transition	c, IV	c, III	2.84x10 <sup>6</sup>	294.5	21.3	-0.512	-1.74	
		Transition	c, IV	c, II	2.84x10 <sup>6</sup>	294.5	21.3	-0.381	-1.29	
		Transition	c, III	c, II	2.84x10 <sup>6</sup>	294.5	21.3	0.131	0.445	
		Transition	c, III	c, II	61500	401.18	128.02	-0.502	-1.25	
		Transition	c, III	c, I	61500	401.18	128.02	0.558	1.39	
		Transition	c, II	c, I	61500	401.18	128.02	1.060	2.64	
		Transition	c, II	c, I	760	400.9	127.7	1.3	3.2	2.9
		Fusion	c, I	liq		610.	337.	2.8	4.6	0.7
		Fusion	c	liq		603.	330.			
		Fusion	c	liq		295.	22.			
KNO <sub>3</sub> ·2HNO <sub>3</sub>		Fusion								



# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 93-5. Potassium (at. no., 19; at. wt., 39.096)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$KPO_3$		Transition Fusion	c, II c, I	c, I liq		723. 1083.	450. 810.			
$K_3PO_4$		Fusion	c	liq		1613.	1340.	2.1	1.9	
$K_4P_2O_7$		Transition Fusion	c, II c, I	c, I liq		551. 1363.	278. 1090.	8.9	5.5	
$KH_2PO_4$		Transition	c, II	c, I		121.97	-151.19	14.	10.	
$K_2HPO_4 \cdot 6H_2O$		Fusion	c	liq		286.01	14.85	0.085	0.70	
$KPO_3 \cdot KF$		Transition Fusion	c, II c, I	c, I liq		813. 1153.	540. 880.			
$KH_2AsO_4$		Transition	c, II	c, I		95.57	-177.59	0.084	0.88	
KSb		Fusion	c	liq		876.	605.			
$K_3Sb$		Fusion	c	liq		1085.	812.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 93-6. Potassium (at. no., 19; at. wt., 39.096)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		OK	OC			
KBi <sub>2</sub>			Fusion	c	liq		825.	552.			
K <sub>2</sub> Bi			Fusion	c	liq		827.	554.			
K <sub>3</sub> Bi			Fusion	c	liq		944.	671.			
K <sub>3</sub> Bi <sub>2</sub>			Fusion	c	liq		715.	442.			
K <sub>2</sub> CO <sub>3</sub>			Transition	c, IV	c, III		523.	250.			
			Transition	c, III	c, II		701.	428.			
			Transition	c, II	c, I		895.	622.			
			Fusion	c, I	liq		1169.	896.	7.8	6.8	
KCHO <sub>2</sub>	potassium formate		Fusion	c	liq		440.7	167.5			
KC <sub>2</sub> H <sub>3</sub> O <sub>2</sub>	potassium acetate		Fusion	c	liq		568.	295.			
KCHO <sub>2</sub> ·HCO <sub>2</sub> H			Fusion	c	liq		555.0	381.8			
KC <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ·HC <sub>2</sub> H <sub>3</sub> O <sub>2</sub>			Fusion	c	liq		421.	148.			
KCN			Transition	c, II	c, I		168.3	-104.9	0.30	1.8	
			Fusion	c, I	liq		883.	610.	3.5	4.0	

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 93-7. Potassium (at. no., 19; at. wt., 39.096) HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION March 31, 1950										
Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
KCNs		Transition Fusion	c, II c, I	c, I liq		413.8 450.	140.6 177.	0.030 2.5	0.075 5.6	
K <sub>2</sub> SiO <sub>3</sub>		Fusion	c	liq		1249.	976.			
K <sub>2</sub> O•2SiO <sub>2</sub>		Fusion	c	liq		1318.	1045.			
K <sub>2</sub> O•4SiO <sub>2</sub>		Transition Fusion	c, II c, I	c, I liq		865. 1038.	592. 765.	0.78 11.7	0.90 11.3	
K <sub>2</sub> O•GeO <sub>2</sub>		Fusion	c	liq		1115.	842.			
K <sub>2</sub> O•4GeO <sub>2</sub>		Fusion	c	liq		1311.	1038.			
2Kf•3eF <sub>4</sub>		Fusion	c	liq		1003.	730.			
KPtI <sub>3</sub>		Fusion	c	liq		622.	349.			
K <sub>2</sub> SO <sub>4</sub> •2PbSO <sub>4</sub>		Transition Fusion	c, II c, I	c, I liq		817. 1221.	544. 948.			
KTi		Fusion	c	liq		608.	335.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 93-8. Potassium (at. no., 19; at. wt., 39.096)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
KCd		Fusion	c	liq		760.	487.			
KReO <sub>4</sub>		Sublimation Fusion	c c	g liq	3.2x10 <sup>-3</sup>	824. 824.	551. 551.	72.8 20.4	88.3 23.9	
K <sub>2</sub> SO <sub>4</sub> · 2CoSO <sub>4</sub>		Fusion	c	liq		1009.	736.			
KFe(SO <sub>4</sub> ) <sub>2</sub> · 12H <sub>2</sub> O		Fusion	c	liq		301.	28.			
KCl · MnCl <sub>2</sub>		Fusion	c	liq		768.	495.			
K <sub>2</sub> SO <sub>4</sub> · 2MnSO <sub>4</sub>		Fusion	c	liq		1118.	845.			
K <sub>2</sub> CrO <sub>4</sub>		Transition Fusion	c, II c, I	c, I liq		938. 1253.	665. 980.	2.45 6.9	2.61 5.5	
K <sub>2</sub> Cr <sub>2</sub> O <sub>7</sub>		Transition Fusion	c, II c, I	c, I liq		514.6 671.	241.6 398.	8.5	13.	23.9

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$K_2MoO_4$		Transition Transition Transition Fusion	c, IV c, III c, II c, I	c, III c, II c, I liq		596. 731. 753. 1199.	323. 458. 480. 926.			
$K_2Mo_2O_7$		Fusion	c	liq		757.	484.		3.	
$K_2O \cdot 3MoO_3$		Fusion	c	liq		844.	571.			
$K_2WO_4$		Transition Fusion	c, II c, I	c, I liq		623. 1203.	350. 930.	4.4	3.6	
$K_2W_2O_7$		Fusion	c	liq		828.	555.			
$KVO_3$		Fusion	c	liq		793.	520.			
$KBO_2$		Fusion	c	liq		1220.	947.	5.7	4.7	
$K_2O \cdot 8_2O_3$		Fusion	c	liq		1243.	970.			
$K_2O \cdot 2B_2O_3$		Fusion	c	liq		1088.	815.			
$K_2O \cdot 4B_2O_3$		Fusion	c	liq		1130.	857.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
KBr <sub>4</sub>		Transition Fusion	c, II c, I	c, I liq		548. 803.	275. 530.			
KPO <sub>3</sub> ·KBO <sub>2</sub>		Fusion	c	liq		1145.	872.			
3KF·AlF <sub>3</sub>		Fusion	c	liq		1308.	1035.			
KBr·AlBr <sub>3</sub>		Fusion	c	liq		454.7	191.5			
KBr·2AlBr <sub>3</sub>		Fusion	c	liq		369.0	95.8			
KAl(SO <sub>4</sub> ) <sub>2</sub> ·12H <sub>2</sub> O		Transition Fusion	c, II c, I	c, I liq		57.9 364.	-215.3 91.	0.047 6.7	0.8 18.4	
KAlSi <sub>2</sub> O <sub>6</sub>	lincite	Transition Fusion	c, II c, I	c, I liq		893. 1958.	620. 1685.			
K <sub>2</sub> SO <sub>4</sub> ·BeSO <sub>4</sub>		Fusion	c	liq		1188.	915.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 93-11. Potassium (at. no., 19; at. wt., 39.096)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
KF·MgF <sub>2</sub>		Fusion	c	liq		1363.	1090.			
KCl·MgCl <sub>2</sub>		Fusion	c	liq		760.	487.	12.69	16.7	
2KCl·MgCl <sub>2</sub>		Fusion	c	liq		706.	433.			
K <sub>2</sub> SO <sub>4</sub> ·2MgSO <sub>4</sub>		Fusion	c	liq		1203.	930.			
KCl·CaCl <sub>2</sub>		Fusion	c	liq		1027.	754.			
K <sub>2</sub> SO <sub>4</sub> ·2CaSO <sub>4</sub>		Transition	c, II	c, I		1211.	938.			
		Fusion	c	liq		1277.	1004.			
KCaPO <sub>4</sub>		Transition	c, II	c, I		978.	705.			
K <sub>2</sub> O·CaO·SiO <sub>2</sub>		Fusion	c	liq		1903.	1630.			
K <sub>2</sub> O·3CaO·6SiO <sub>3</sub>		Fusion	c	liq		1293.	1020.			
2K <sub>2</sub> O·CaO·3SiO <sub>2</sub>		Fusion	c	liq		1232.	959.			
2K <sub>2</sub> O·CaO·6SiO <sub>2</sub>		Fusion	c	liq		1278.	1005.			
4K <sub>2</sub> O·CaO·10SiO <sub>2</sub>		Fusion	c	liq		1219.	946.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards Washington, D.C.

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$KBr \cdot 2SrBr_2$		Fusion	c	liq		847.	574.			
$2KBr \cdot SrBr_2$		Fusion	c	liq		832.	559.			
$K_2SO_4 \cdot 2SrSO_4$		Transition	c, II	c, I		1048.	775.			
$2KBr \cdot BaBr_2$		Fusion	c	liq		907.	634.			
$K_2SO_4 \cdot Li_2SO_4$		Fusion	c	liq		989.	716.			
$K_2MoO_4 \cdot Li_2MoO_4$		Transition	c, II	c, I		685.	412.			
		Fusion	c, I	liq		844.	571.			
$K_2WO_4 \cdot Li_2WO_4$		Fusion	c	liq		905.	632.			



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 94-1. Rubidium (at. no., 37; at. wt., 85.48)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
Rb	Transition	c, II	c, I			243.	-30.			
	Sublimation	c, I	q		.01	312.0	38.8	20.47	65.61	
	Fusion	c, I	liq			312.0	36.8	0.52	1.67	
	Vaporization	liq	g		760	952.	679.	18.11	19.02	
RbOH	Transition	c, II	c, I			518.	245.	1.70	3.28	
	Fusion	c, I	liq			574.	301.	1.62	2.82	
RbF	Fusion	c	liq			1048.	775.	4.13	3.94	
	Vaporization	liq	q		760	1681.	1408.	39.51	23.50	
RbF·HF	Fusion	c	liq			478.	205.			
	Fusion	c	liq			324.8	51.6			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 94-2. Rubidium (at. no., 37; at. wt., 85.48)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
RbCl		Fusion Vaporization	c liq	liq g	760	990. 1654.	717. 1381.	4.40 36.92	4.44 22.32	
RbBr		Fusion Vaporization	c liq	liq g	760	953. 1625.	680. 1352.	3.7 37.12	3.9 22.84	
RbI		Fusion Vaporization	c liq	liq g	760	913. 1577.	640. 1304.	2.99 35.96	3.27 22.80	
Rb <sub>2</sub> S <sub>2</sub>		Fusion	c	liq		693.	420.			
Rb <sub>2</sub> S <sub>3</sub>		Fusion	c	liq		486.	213.			
Rb <sub>2</sub> S <sub>5</sub>		Fusion	c	liq		504.	231.			
Rb <sub>2</sub> S <sub>6</sub>		Fusion	c	liq		474.	201.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

National Bureau of Standards

Washington, D.C.

Table 94-3. Rubidium (at. no., 37; at. wt., 85.48)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Formula	Substance	Description	Type of Process	State		Pressure mm Hg	Temperature		$\Delta F$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
				Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Rb_2SO_4$			Transition Fusion	c, II c, I	c, I liq		923. 1347.	650. 1074.			
$RbSH$			Transition	c, II	c, I		403.	130.	0.4	1.0	
$RbSeH$			Transition	c, II	c, I		420.	147.			
$RbN_3$			Fusion	c	liq		594.	321.			
$RbNO_3$			Transition Transition Transition	c, IV c, III c, II	c, III c, II c, I		438. 498. 564.	165. 225. 291.			
			Fusion	c, I	liq		589.	316.	1.34	2.28	
$RbNO_3 \cdot HNO_3$			Fusion	c	liq		335.	62.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 94-4. Rubidium (at. no., 37; at. wt., 85.48)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
$Rb_2CO_3$		Fusion	c	liq		1108.	835.			
$RbCHO_2$	rubidium formate	Fusion	c	liq		443.	170.			
$RbC_2H_3O_2$	rubidium acetate	Fusion	c	liq		519.	246.			
$RbCNS$		Fusion	c	liq		458.	195.			
$RbBF_4$		Fusion	c	liq		663.	590.			
$3RbF \cdot AlF_3$		Fusion	c	liq		1258.	985.			
$RbNO_3 \cdot LiNO_3$		Fusion	c	liq		464.	191.			
$RbOH \cdot 2NaOH$		Fusion	c	liq		551.	278.			
$2RbOH \cdot KOH$		Fusion	c	liq		672.	399.			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES II

National Bureau of Standards

Washington, D.C.

Table 95-1. Cesium (at. no., 55; at. wt., 132.91)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		°K	°C			
Cs	Sublimation		c	g	$1.2 \times 10^{-6}$	301.9	28.7	18.82	62.34	
	Fusion		c	liq	$1.2 \times 10^{-6}$	301.9	28.7	0.50	1.6	
	Vaporization		liq	g	760	963.	690.	16.32	16.95	
Cs <sub>2</sub> O	Fusion		c	liq		763.	490.			
Cs <sub>2</sub> O	Fusion		c	liq		276.	3.			
CsOH	Transition		c, II	c, I		496.	223.	1.76	3.55	
	Fusion		c, I	liq		545.5	272.3	1.61	2.93	
CsF	Fusion		c	liq		955.	682.	2.45	2.56	
CsF•HF	Fusion		c	liq		449.	176.			
CsF•2HF	Fusion		c	liq		323.4	50.2			
CsF•3HF	Fusion		c	liq		305.8	32.6			
CsF•6HF	Fusion		c	liq		230.9	-42.3			

# SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta G_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
CsCl		Transition	c, II	c, I	760	748.	445.	1.8	2.5	
		Fusion	c, I	liq		948.	645.	3.60	3.92	
		Vaporization	liq	g		1573.	1300.	35.69	22.69	
CsBr		Fusion	c	liq		909.	636.	1.7	1.9	
		Vaporization	liq	g		1573.	1300.	35.99	22.88	
CsBrO <sub>3</sub>		Fusion	c	liq		693.	420.			
CsI		Fusion	c	liq		894.	621.			
		Vaporization	liq	g		1573.	1300.	35.93	22.84	
Cs <sub>2</sub> S <sub>2</sub>		Fusion	c	liq		733.	460.			
Cs <sub>2</sub> S <sub>3</sub>		Fusion	c	liq		490.	217.			
Cs <sub>2</sub> S <sub>5</sub>		Fusion	c	liq		483.	210.			
Cs <sub>2</sub> S <sub>6</sub>		Fusion	c	liq		459.	186.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES II

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Table 95-3. Cesium (at. no., 55; at. wt., 132.91)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$Cs_2SO_4$		Transition Fusion	c, II c, I	c, I liq		933. 1292.	660. 1019.			
$CsNH_3$		Fusion	c	liq		599.	326.			
$CsNH_2$		Fusion	c	liq		535.	262.			
$CsNO_3$		Transition Fusion	c, II c, I	c, I liq		429. 690.	156. 417.	3.25	4.71	
$CsNO_3 \cdot 4H_2O$		Fusion	c	liq		315.9	42.7			
$CsNO_3 \cdot HNO_3$		Fusion	c	liq		373.	100.			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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National Bureau of Standards

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Table 95-4. Cesium (at. no., 55; at. wt., 132.91)  
HEAT, TEMPERATURE, AND ENTROPY OF TRANSITION, FUSION, AND VAPORIZATION  
March 31, 1950

Substance		Type of Process	State		Pressure mm Hg	Temperature		$\Delta H$ kcal/mole	$\Delta S$ cal/deg mole	$\Delta C_p$ cal/deg mole
Formula	Description		Initial	Final		$^{\circ}K$	$^{\circ}C$			
$CsHCO_2$	cesium formate	Fusion	c	liq		538.	265.			
$CsHCO_2 \cdot H_2O$		Transition	c, II	c, I		314.	41.			
		Fusion	c, I	liq		318.	45.			
$CsC_2H_3O_2$	cesium acetate	Fusion	c	liq		467.	194.			
$CsCl \cdot 2CuCl$		Fusion	c	liq		547.	274.			
$Cs_2TiF_6$		Fusion	c	liq		963.	690.			
$CsBF_4$		Fusion	c	liq		823.	550.			
$3CsF \cdot AlF_3$		Fusion	c	liq		1096.	823.			



#### IV. SPECIFIC REFERENCES FOR THE TABLES OF PROPERTIES OF SERIES I

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

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Table 1  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
O	ozone	g	1,2,3		4	4
O <sup>n+</sup>		g	5,6,7,8,9,10			
O <sup>-</sup>		g	11,12,13,14			
O <sub>2</sub>		g			15,16,17	15,16,17,18
		aq	19,20,21			
O <sub>2</sub> <sup>+</sup>		g	22,23,24			
O <sub>3</sub>		g	25,26,27		28,29,30	4
		aq	31,32			
O <sub>4</sub>		g	33,34			
REFERENCES						
1. Knauss and Ballard-1			19. Winkler-1,7,9			
2. Herzberg-9			20. Cassuto-1			
3. Bichowsky and Rossini-1			21. Washburn and Strachan-1			
4. Calculated			22. Tanaka and Takamine-1			
5. Moore-1			23. Mulliken and Stevens-1			
6. Edlen-5			24. Hagstrum and Tate-1			
7. Edlen-7			25. Kailan and Jahn-1			
8. Kruger and Shoupp-1			26. Mulder and van der Meulen-2			
9. Cady-4			27. Günther, Wassmuth, and Schryver-1			
10. Robinson-5			28. Kelley-24			
11. Senftleben-1			29. Shand and Spurr-1			
12. Born and Gerlach-1			30. Kassel-3			
13. Hanson-2			31. Fischer and Tropsch-1			
14. Vier and Mayer-1			32. Rothmund-1			
15. Wagman, Kilpatrick, Taylor, Pitzer, and Rossini-1			33. Lewis-4			
16. Johnston and Walker-2			34. Kondrat'ev-1			
17. Woolley-1						
18. Trautz and Ader-1						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 2  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
H	water	g	1,2,3,4	25,26,27	5,6	5,6
H <sup>+</sup>		g	7			
		aq				
H <sub>2</sub>		g			6,8,9,10, 11,12	6,8,9,10, 12
OH		g	13,14,15,16,17,18, 19,20,21		5,11,17	5,17
OH <sup>-</sup>		g	22			
		aq	1,23,24		5,31,32,11, 33	23,28 5,31,32,34
H <sub>2</sub> O		g	8,29,30			
		liq	1,8,29			
H <sub>2</sub> O <sub>2</sub>		g	37			
		liq	38,39			
		aq	40,41,42		35,36	30,35

REFERENCES

- |  |   |
|--|---|
| <ol style="list-style-type: none"> <li>1. Bichowsky and Rossini-1</li> <li>2. Beutler-2</li> <li>3. Beutler and Junger-2</li> <li>4. Herzberg-9</li> <li>5. Calculated</li> <li>6. Giauque-2</li> <li>7. Moore-1</li> <li>8. Wagman, Kilpatrick, Taylor, Pitzer, and Rossini-1</li> <li>9. Woolley, Scott, and Brickwedde-1</li> <li>10. Davis and Johnston-1</li> <li>11. Zeise-1</li> <li>12. Johnston et al-1</li> <li>13. Bonhoffer and Haber-1</li> <li>14. Bonhoffer and Reichardt-1,2</li> <li>15. Bates-6</li> <li>16. Tanaka and Koana-1</li> <li>17. Johnston and Dawson-1</li> <li>18. Senftleben and Riechemeier-1</li> <li>19. Lewis and von Elbe-3</li> <li>20. Dwyer and Oldenberg-1</li> <li>21. Avramenko and Kondrat'ev-1</li> <li>22. Lederle-2</li> <li>23. Pitzer-3</li> <li>24. Rossini-5</li> </ol> | <ol style="list-style-type: none"> <li>25. Harned and Hamer-1</li> <li>26. Harned and Copsone-1</li> <li>27. Harned and Owen-1</li> <li>28. Rossini-6</li> <li>29. Rossini-16</li> <li>30. Osborne, Stimson, and Ginnings-2</li> <li>31. Gordon-2</li> <li>32. Wilson, Jr.-1</li> <li>33. Giauque and Ashley-1</li> <li>34. Trautz and Ader-1</li> <li>35. Giauque and Stout-1</li> <li>36. Kelley-24</li> <li>37. Maass and Hiebert-1</li> <li>38. Matheson and Maass-1</li> <li>39. de Forcrand-33</li> <li>40. Thomsen-16</li> <li>41. Berthelot-28,61</li> <li>42. Roth, Grau, and Meichsner-1</li> </ol> |
|--|---|

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

SERIES I

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Table 2a  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$^1\text{H}$		g	1		2	2
$^1\text{H}^+$		g	3			
$^1\text{H}^-$		g	4			
$^1\text{H}_2$		g			2,5	2
$^1\text{H}_2^+$		g	6,7			
$^1\text{H}_2\text{O}$		g	8,9,10		2,5	2
		liq	9,10		2	

REFERENCES

- |   |  |
|---|--|
| <p>1. Herzberg-9<br/>2. Calculated<br/>3. Moore-1<br/>4. Hylleraas-5<br/>5. Urey and Rittenberg-1<br/>6. Beutler and Junger-2<br/>7. Hagstrum and Tate-1<br/>8. Rossini, Knowlton, and Johnston-1</p> | <p>9. Wagman, Kilpatrick, Taylor, Pitzer, and Rossini-1<br/>10. Rossini-16</p> |
|---|--|

Table 2b

$^2\text{H}$		g	1,2,3		4	4
$^2\text{H}^+$		g	5,6			
$^2\text{H}^-$		g	7			
$^2\text{H}_2$		g			8,9,10	8,9
$^1\text{H}^2\text{H}$		g	3,4,8		9,11	8,9
$^2\text{H}_2\text{O}$		g	11		11,12,13	4
		liq	11		11,13,14	13,15
$^1\text{H}^2\text{HO}$		g	11		4	4
		liq	16	4,16		4

REFERENCES

- |   |  |
|---|--|
| <p>1. Beutler-1<br/>2. Beutler-2<br/>3. Herzberg-9<br/>4. Calculated<br/>5. Birge-12<br/>6. Moore-1<br/>7. Hylleraas-5<br/>8. Johnston and Long-2<br/>9. Woolley, Scott, and Brickwedde-1</p> | <p>10. Clusius and Bartholomé-3<br/>11. Rossini, Knowlton, and Johnston-1<br/>12. Libby-1<br/>13. Long and Kemp-1<br/>14. Kelley-24<br/>15. Bartholomé and Clusius-1<br/>16. Wahl and Urey-1</p> |
|---|--|

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 3  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
He		g	2, 3		1	1
He <sup>n+</sup>		g				

REFERENCES

1. Calculated
2. Moore-1
3. Bacher and Goudsmit-1

Table 4

Ne		g	2, 3		1	1
Ne <sup>n+</sup>		g				

REFERENCES

1. Calculated
2. Moore-1
3. Paul and Polster-1

Table 5

A		g	2 3		1	1
A·5H <sub>2</sub> O		c				
A <sup>n+</sup>		g				

REFERENCES

1. Calculated
2. deForcrand-69
3. Moore-1

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 6  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Kr		g			1	1
Kr·5H <sub>2</sub> O		c	2			
Kr <sup>n+</sup>		g	3,4,5,6			

REFERENCES

- |  |                            |
|--|----------------------------|
| <p>1. Calculated</p> <p>2. de Forcrand-69</p> <p>3. Boyce-1</p> <p>4. de Bruin, Humphreys, and Meggers-1</p> <p>5. Humphreys-3</p> | <p>6. Tate and Smith-2</p> |
|--|----------------------------|

Table 7

Xe		g			1,2	1
Xe·6H <sub>2</sub> O		c	3			
Xe <sup>n+</sup>		g	4,5,6,7			

REFERENCES

- |  |   |
|--|---|
| <p>1. Calculated</p> <p>2. Clusius and Riccoboni-1</p> <p>3. de Forcrand-71</p> <p>4. Tate and Smith-2</p> <p>5. Humphreys and Meggers-1</p> | <p>6. Humphreys-1</p> <p>7. Humphreys-2</p> |
|--|---|

Table 8

Rn		g			1	1
Rn <sup>n+</sup>		g	2,3			

REFERENCES

- |  |  |
|--|--|
| <p>1. Calculated</p> <p>2. Rasmussen-1</p> <p>3. Rasmussen-4</p> |  |
|--|--|

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 9  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
F		g	1, 2, 6		3, 4, 5	3
$F^{n+}$		g	4			
$F^-$		g	3, 7, 8			
		aq	3, 9, 14		10	16, 17, 18
$F_2$		g			3	3
$F_2O$		g	11, 15, 19		12, 13	
HF		g	19, 20, 21		3, 5, 25	3, 5
		aq	9, 17, 20, 22, 23, 24		10	16, 17, 18, 19
$HF_2^-$		aq	3			
$(HF)_6$		g	26			

REFERENCES-

- |  |   |
|--|---|
| 1. von Wartenberg and Taylor-1           | 15. von Wartenberg-8                    |
| 2. Schmitz and Schumacher-2,3,4,5        | 16. Thorvaldson and Bailey-1            |
| 3. Calculated                            | 17. Roth, Pahlke, Bertram, and Börger-1 |
| 4. Moore-1                               | 18. Pranschke and Schwiete-1            |
| 5. Murphy and Vance-1                    | 19. Ruff and Menzel-1                   |
| 6. Desai- 5                              | 20. von Wartenberg and Fitzner-1        |
| 7. Dukel'skii and Ionov-1                | 21. von Wartenberg and Schütza-2        |
| 8. Wu-2                                  | 22. Petersen-7                          |
| 9. Roth-2                                | 23. Guntz-1                             |
| 10. Latimer, Pitzer, and Smith-1         | 24. Roth-25                             |
| 11. von Wartenberg and Klinkott-1        | 25. Herzberg-9                          |
| 12. Kelley-24                            | 26. Long, Hildebrand, and Morrell-1     |
| 13. Potter-2                             |   |
| 14. Elmore, Hatfield, Mason, and Jones-1 |   |

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 10  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Cl		g	1, 2, 3		4, 5	4
Cl <sup>n+</sup>		g	5			
Cl <sup>-</sup>		g	6, 7, 8, 9			
		aq	4			10, 34
Cl <sub>2</sub>		g			4, 11, 12	4, 12, 13
		aq	14, 15, 16, 17			
	in CCl <sub>4</sub>		18			
Cl <sub>2</sub> ·8H <sub>2</sub> O		c	19, 20, 54			
Cl <sub>3</sub>		g	21			
ClO		g	22			
ClO <sup>-</sup>		aq	4		4, 23	
ClO <sub>2</sub>		g	24, 25, 26		27	
		aq	24			
ClO <sub>2</sub> <sup>-</sup>		aq	4		4, 28	
ClO <sub>3</sub>		g	29			
ClO <sub>3</sub> <sup>-</sup>		aq	4		4	
ClO <sub>4</sub> <sup>-</sup>		aq	4		4	
Cl <sub>2</sub> O		g	25, 30, 31, 32	33	27	
		aq	30, 31			
Cl <sub>2</sub> O <sub>7</sub>		g	34			
HCl		g	35, 36, 37, 38		4, 12, 39, 40, 41	4, 41
		aq	30, 37, 42, 43, 44, 45, 46, 47, 48, 49, 50, 51, 52			10, 45, 51, 52
	in CCl <sub>4</sub>		18			
	in C <sub>2</sub> H <sub>5</sub> OH		53			
HClO		aq	30, 31			
HClO <sub>2</sub>		aq	45			
HClO <sub>3</sub>		aq	30			
HClO <sub>4</sub>		liq	55			
		aq	55			
HClO <sub>4</sub> ·H <sub>2</sub> O		c	55, 56			
HClO <sub>4</sub> ·2H <sub>2</sub> O		liq	55			
ClF		g	57, 58, 59		60	60
ClF <sub>3</sub>		g	59, 61, 62		62	62



## SERIES I

National Bureau of Standards

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Table 10 (Continued)  
SPECIFIC REFERENCES

## REFERENCES

- |                                  |                                      |
|----------------------------------|--------------------------------------|
| 1. Kuhn-1                        | 43. Berthelot and Louguinine-1       |
| 2. Wohl-2,3                      | 44. Becker and Roth-4                |
| 3. Herzberg-9                    | 45. Pitzer-3                         |
| 4. Calculated                    | 46. Roth and Bertram-2               |
| 5. Moore-1                       | 47. Roth, Börger, and Siemonsen-1    |
| 6. Mitchell and Mayer-1          | 48. Sturtevant-1,4                   |
| 7. McCallum and Mayer-1          | 49. Slansky-1                        |
| 8. Beutler and Levi-1            | 50. Rossini-4                        |
| 9. Dukel'skii and Ionov-1        | 51. Thorvaldson, Brown, and Peaker-2 |
| 10. Rossini-6                    | 52. Sturtevant-7                     |
| 11. Giauque and Powell-1         | 53. Berthelot-15B                    |
| 12. Giauque and Overstreet-1     | 54. Harris-1                         |
| 13. Trautz and Ader-2            | 55. Berthelot-79,B0                  |
| 14. Thomsen-16                   | 56. Arlman-1                         |
| 15. Baker-2                      | 57. Ruff and Laass-1                 |
| 16. Berthelot-24                 | 58. Wahrhaftig-1                     |
| 17. von Wartenberg and Werth-3   | 59. Schmitz and Schumacher-4         |
| 18. Noyes and Tuley-1            | 60. Potter-2                         |
| 19. de Forcrand-39               | 61. Schmitz and Schumacher-3         |
| 20. Roozeboom-2,5                | 62. Schäfer and Wicke-1              |
| 21. Bodenstein and Plaut-1       |                                      |
| 22. Finkelnburg and Schumacher-1 |                                      |
| 23. Latimer, Pitzer, and Smith-1 |                                      |
| 24. Booth and Bowen-1            |                                      |
| 25. Wallace and Goodeve-1        |                                      |
| 26. Mayer-4                      |                                      |
| 27. Kelley-24                    |                                      |
| 28. Smith, Pitzer, and Latimer-1 |                                      |
| 29. Goodeve and Marsh-3          |                                      |
| 30. Thomsen-16                   |                                      |
| 31. Neumann and Muller-2,3       |                                      |
| 32. Gunther and Wekua-1          |                                      |
| 33. Yost and Felt-1              |                                      |
| 34. Goodeve and Marsh-1          |                                      |
| 35. Rossini-9                    |                                      |
| 36. von Wartenberg and Hanisch-1 |                                      |
| 37. Roth and Richter-2           |                                      |
| 38. Bichowsky and Rossini-1      |                                      |
| 39. Giauque and Wiebe-1          |                                      |
| 40. Linnett-2                    |                                      |
| 41. Gordon and Barnes-1          |                                      |
| 42. Vrevskii and Savaritskii-1   |                                      |

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Table 11  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Br		g	1,2,3,4		5	5
$\text{Br}^{n+}$		g	6,7,8,9,10			
$\text{Br}^-$		g	11,12,13,14,15			
		aq	5	16	5	17
$\text{Br}_2$		g	18,19,20		5,21,22	5,22
		liq			5,23,38	
		aq	24,25,26,27			
	in $\text{CCl}_4$		26,28			
	in $\text{CHCl}_3$		26			
	in $\text{CS}_2$		26			
$\text{Br}_2 \cdot 10\text{H}_2\text{O}$		c	29,30			
$\text{Br}_3^-$		aq	31,32,33,34,35			
$\text{Br}_5^-$		aq	32			
$\text{BrO}_3^-$		aq	5		5	5
HBr		g	33,36,37,39		5,22	5,22
		aq	33,39,40,41,11		5	17
$\text{HBrO}_3$		aq	39		5	
$\text{BrCl}$		g	28,43		5,43	5
	in $\text{CCl}_4$		28			

REFERENCES

1. Brown-4	15. Beutler and Levi-1
2. Darbyshire-3	16. Jones and Baekstrom-1
3. Kuhn-1	17. Rossini-6
4. DeVries and Rodebush-1	18. Smits and Cannegieter-1
5. Calculated	19. Berthelot and Ogier-6
6. Kiess and DeBruin-1	20. Lewis and Randall-7
7. Bloch, Bloch, and Lacroute-1	21. Brown-1
8. Rao and Krishnamurty-1	22. Gordon and Barnes-2
9. Rao and Krishnamurty-2	23. Kelley-24
10. Bacher and Goudsmit-1	24. Berthelot-99
11. Bichowsky and Rossini-1	25. Thomsen-2
12. Dukel'skii and Ionov-1	26. Pickering-9
13. Doty and Mayer-1	27. Bray and Connolly-1
14. Glockler and Calvin-1	28. Blair and Yost-1

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 11 (Continued)  
SPECIFIC REFERENCES

REFERENCES

29. Roozeboom-3,5
30. Tammann and Krige-1
31. Berthelot-104
32. Linhart-1
33. Roth, Börger, and Siemonsen-1
34. von Wartenberg and Klinkott-1
35. Hieber and Woerner-3
36. Roth and Börger-2
37. Roth-17
38. Latimer and Hoenschel-1
39. Thomsen-16
40. Roth and Bertram-2
41. Harned and Hamer-2
42. Giauque and Wiebe-2,3
43. Beeson and Yost-1

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Table 12  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
I		g	1, 2, 3, 4, 5, 7		6, 8, 9	6
$I^{n+}$		g	10, 11, 12			
$I^-$		g	13, 14, 15, 16, 17			
		aq	6	20, 21	6	6, 18
$I_2$		g	7		6, 7	6
		c			7, 19	19
		aq	22, 23			
	in $C_6H_6$		24, 25			
	in $C_2H_5OH$		25			
	in $(C_2H_5)_2O$		24, 25			
	in $CCl_4$		24			
	in $CHCl_3$		24, 25			
	in $CS_2$		24, 25			
$I_3^-$		aq	24, 26, 27, 28	29		
$IO^-$		aq	6			
$IO_3^-$		aq	6		6	6
$I_2O_5$		c	30, 31, 32, 33			
$3I_2O_5 \cdot H_2O$		c	30			
HI		g	31, 34, 35, 36, 37, 38		6, 37, 41	6
		aq	4, 31, 39, 40, 42	20, 21		18
HIO		aq	43, 44			
$HIO_3$		c	30, 31, 32, 33			
		aq	31, 45			
$HIO_6^{----}$		aq	31			
$H_2IO_6^{---}$		aq	31			
$H_3IO_6^{--}$		aq	31			
$H_4IO_6^-$		aq	31			
$H_5IO_6$		c	31			
		aq	31			
$I_2O_5 \cdot HIO_3$		c	33			
ICl		g	46, 47, 48, 49		6, 47	6
	in $CCl_4$		50			
$ICl_3$		c	31, 49, 51	51		
IBr		g	52, 53, 54, 55, 56, 57		6, 46	6
	in $CCl_4$		50	58		

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| 29. Lewis and Randall-7           |                                 |
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
S	in CS <sub>2</sub>	g	24		23, 3	23, 3
		liq	12, 10, 13, 14, 15, 8, 1, 11			
		c	6, 7, 8, 9, 10, 11, 4, 12		4	4
			5, 16, 17			
S <sup>n+</sup>		g	23			
S <sup>--</sup>		g	25			
		aq	3	26		
S <sub>2</sub>		g	20, 31, 21, 64, 65, 66		1, 2	2, 3
S <sub>2</sub> <sup>--</sup>		aq	27, 3			
S <sub>3</sub> <sup>--</sup>		aq	27, 3			
S <sub>4</sub> <sup>--</sup>		aq	27, 3			
S <sub>6</sub>		g	20, 21			
S <sub>8</sub>		g	11, 18, 19, 21, 22			
SO		g	28, 29		28	
SO <sub>2</sub>		g	30, 31		32, 36	32, 36
		aq	33, 34, 35, 37, 31			
SO <sub>2</sub> ·7H <sub>2</sub> O		c	38, 39			
SO <sub>3</sub>		g	44, 42, 43, 41, 40, 47		44	44
		c	48, 42, 43, 49, 50			
		liq	45, 46, 40			
SO <sub>3</sub> <sup>--</sup>		aq	45, 51, 52, 3, 34	20		
SO <sub>4</sub> <sup>--</sup>		aq	45, 81, 41, 82, 83, 84, 85, 86, 87, 92	90, 91	89	88
S <sub>2</sub> O <sub>3</sub> <sup>--</sup>		aq	3, 45, 55	3		
S <sub>2</sub> O <sub>4</sub> <sup>--</sup>		aq	53, 3	54		
S <sub>2</sub> O <sub>5</sub> <sup>--</sup>		aq	31, 56, 77			
S <sub>2</sub> O <sub>6</sub> <sup>--</sup>		aq	45, 3			
S <sub>2</sub> O <sub>7</sub>		c	58			
S <sub>2</sub> O <sub>8</sub> <sup>--</sup>		aq	59, 60			
S <sub>3</sub> O <sub>6</sub> <sup>--</sup>		aq	55, 3			
S <sub>4</sub> O <sub>6</sub> <sup>--</sup>		aq	45, 55, 3, 57	54		
S <sub>5</sub> O <sub>6</sub> <sup>--</sup>		aq	55, 3			
HS <sup>-</sup>		aq	67	26		
H <sub>2</sub> S		g	67, 31		1, 69, 70, 71	1
		aq	67, 45	54		
H <sub>2</sub> S·6H <sub>2</sub> O		c	68, 72, 73, 74, 75			
H <sub>2</sub> S <sub>2</sub>		liq	3			
H <sub>2</sub> S <sub>5</sub>		liq	76			
HSO <sub>3</sub> <sup>-</sup>		aq	45, 51, 78, 52, 35	35		
HSO <sub>4</sub> <sup>-</sup>		aq	79	80		
H <sub>2</sub> SO <sub>3</sub>		aq	33, 34, 35, 37, 31	35		

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$H_2SO_4$	in $C_2H_5OC_2H_5$	liq	82,83	90,91	89	95
		aq	45,81,41,82,83,84, 85,86,87,92			88
			93			
$H_2SO_4 \cdot H_2O$		liq	94			
$H_2S_2O_4$		aq	3			
$H_2S_2O_6$		aq	45,3			
$H_2S_2O_7$		liq	45,96,97			
$H_2S_2O_8$		aq	98			
$H_2S_4O_6$		aq	3			
$SF_6$		g	99		100,101,3	
$SCl_4$		liq	102,103			
$S_2Cl_2$		liq	45,61,103,14			103
$S_2Cl_4$		liq	103,14			
$SOCl_2$		liq	62,103			62
$SO_2Cl_2$		liq	45,62			62
$S_2O_5Cl_2$		liq	62			62
$HSO_3Cl$		liq	103			
$S_2Br_2$		liq	63			

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14. Trautz-6	28. Herzberg-9

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| 46. Giran-3                            | 88. Craig and Vinal-1             |
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$	
Formula	Description	State					
Se		g	4, 2, 23, 24, 25		1	1	
		c			7	7	
		gls	5, 1			7	
		amorp	5, 1			7	
Se <sup>n+</sup>		g	3	15	1	1	
Se <sup>--</sup>		aq	13, 1				
Se <sub>2</sub>		g	1, 2, 4, 5, 6, 8, 7, 23, 24, 25				
SeO		g	9				
SeO <sub>2</sub>		c	10, 11				
		aq	10				
SeO <sub>3</sub> <sup>--</sup>		aq	10, 1				15
SeO <sub>4</sub> <sup>--</sup>		aq	10, 1				15
HSe <sup>-</sup>		aq	13, 1	15	19		
H <sub>2</sub> Se		g	13	20			
		aq	20, 13				
HSeO <sub>3</sub> <sup>-</sup>		aq	10, 1				15
HSeO <sub>4</sub> <sup>-</sup>		aq	10, 1		15		
H <sub>2</sub> SeO <sub>3</sub>		c	11, 12	15			
		aq	10, 11				
H <sub>2</sub> SeO <sub>4</sub>		c	14				
		aq	10, 13, 1				
H <sub>2</sub> SeO <sub>4</sub> ·H <sub>2</sub> O		c	14			17, 1	17, 1
SeF <sub>6</sub>		g	16				
SeCl <sub>2</sub>		g	21				
SeCl <sub>4</sub>		c	10				
Se <sub>2</sub> Cl <sub>2</sub>		liq	10, 18				
Se(OH) <sub>3</sub> ClO <sub>4</sub>		c	22				
SeO <sub>2</sub> ·SO <sub>3</sub>		c	14				

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Table 16  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Te		g	17,1		1	1
		c			2,16	2,16
		amorp	28			
Te <sup>n+</sup>		g	4,5,6,7,8,9,10,11,12			
Te <sub>2</sub>		g	13,29,30,31,1		3,1	3,1
TeO		g	14,15			
TeO <sub>2</sub>		c	18,21	18		18,19
TeO <sub>3</sub> <sup>--</sup>		aq	21,1			
TeO <sub>4</sub> <sup>--</sup>		aq	23,1,21			
H <sub>2</sub> Te		g	22		1	
H <sub>2</sub> TeO <sub>3</sub>		c	21	20		
		aq	21			
H <sub>2</sub> TeO <sub>4</sub>		aq	21,22			
H <sub>2</sub> TeO <sub>4</sub> ·2H <sub>2</sub> O		c	23		26	
TeF <sub>6</sub>		g	25		27,1	
TeCl <sub>4</sub>		c	21			
TeBr <sub>4</sub>		c	24			
(TeO <sub>2</sub> ) <sub>2</sub> SO <sub>3</sub>		c	23			

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Table 17  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Po		c				
Po <sup>++</sup>		aq		1,2		
PoO <sub>2</sub>		c		1,2		
REFERENCES						
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Table 18  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
N		g	1, 2, 3, 4, 5, 6		7, 8	7
N <sup>n+</sup>		g	9, 10, 11, 12			
N <sub>2</sub>		g			13, 14, 15	13, 14, 16
N <sub>2</sub> <sup>n+</sup>		g	4, 5, 17, 18, 19, 20			
N <sub>3</sub> <sup>-</sup>		aq	7			
NO		g	1, 4, 5, 13, 21		13, 22, 23, 24	13, 22, 23, 24
NO <sub>2</sub>		g	25, 26, 27, 28		27	28
NO <sub>2</sub> <sup>-</sup>		aq	7			
NO <sub>3</sub>		g	29			
NO <sub>3</sub> <sup>-</sup>		aq	7		7, 30	
N <sub>2</sub> O		g	31, 32, 33, 34, 35		36, 37, 38	36
N <sub>2</sub> O <sub>2</sub> <sup>==</sup>		aq	7, 39, 40		7, 40	
N <sub>2</sub> O <sub>3</sub>		g	41, 42			
N <sub>2</sub> O <sub>4</sub>		g	26, 27, 28, 43, 44		27	27, 28
N <sub>2</sub> O <sub>5</sub>		g	45			
		c	45, 46, 47			
NH		g	5, 48, 49			
NH <sub>3</sub>		g	1, 50		51, 52, 53, 54, 55	51, 55, 56, 57
		aq	1, 58	59		
NH <sub>4</sub> <sup>+</sup>		g	60, 71			
		aq	32, 50, 61, 62	63, 64		
N <sub>2</sub> H <sub>4</sub>		liq	65, 66			
		aq	66, 67, 68, 69, 70			
N <sub>2</sub> H <sub>4</sub> ·H <sub>2</sub> O		liq	65, 70			
N <sub>2</sub> H <sub>4</sub> H <sup>+</sup>		aq	7			
N <sub>2</sub> H <sub>4</sub> H <sub>2</sub> <sup>++</sup>		aq	7			
HN <sub>3</sub>		g	72		73	73
		aq	72, 74			
NH <sub>4</sub> N <sub>3</sub>		c	75			
		aq	74, 75			
HNO <sub>2</sub>		aq	1, 76, 32			
HNO <sub>3</sub>		g			77	
		liq	77		77	77
		aq	77, 32, 78, 79, 80	59	7, 81, 30	
HNO <sub>3</sub> ·nH <sub>2</sub> O		liq	77		77	77
NH <sub>2</sub> OH		c	83			
		aq	32, 82			
NH <sub>2</sub> OH·H <sup>+</sup>		aq	7			
NH <sub>4</sub> OH		aq	7, 1, 58			
HN <sub>2</sub> O <sub>2</sub> <sup>-</sup>		aq	40	40		

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$H_2N_2O_2$	in $CCl_4$	aq	40	107	64	86
$NH_4NO_2$		c	32,84			
		aq	85			
$NH_4NO_3$		c	50			
		aq	1,32,50			
$NH_4NO_3 \cdot 1/2NH_3$		liq	87			
$NH_2OH \cdot HNO_3$		c	88			
		aq	32,88			
$N_2H_4 \cdot nHNO_3$		aq	74			
$NF_3$		g	89			
$NH_4F$		c	90			
		aq	7,90			
$NCI_3$			91,92			99
$NOCl$		g	1,93,94		93,95	
$NH_4Cl$		c	1,96		99	
		aq	1,97,32,98			
$NH_4Cl \cdot nNH_3$		liq	100			
$N_2H_4 \cdot nHCl$		c	74,103			
		aq	74,101,102			
$NH_2OH \cdot HCl$		c	82			
		aq	82,104			
$NH_4ClO_4$		c	105			
		aq	7,105			
$NOBr$		g	106,107			
$NH_4Br$		c	32,108			
		aq	7			
$NH_4Br \cdot nNH_3$		liq	109,110			
$NH_4I$		c	32,111			
		aq	7			
$NH_4I \cdot nNH_3$		liq	109			
$N_4S_4$		c	112			
$NH_4HS$		c	113			
		aq	32,114			
$NH_4S_4$		c	115			
		aq	115			
$(NH_4)_2S$		aq	32,114			
$(NH_4)_2S_5$		c	115			
		aq	115			
$NH_4HSO_3$		c	116,117			
		aq	7,116,117			

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Table 1B (Continued)  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$\text{NH}_4\text{HSO}_4$		c	32			119
		aq	32, 11B			
$\text{NH}_2\text{OH} \cdot \text{H}_2\text{SO}_4$		c	32			
		aq	104			
$\text{N}_2\text{H}_4 \cdot \text{H}_2\text{SO}_4$		c	75, 66, 67			
		aq	74, 75, 122			
$(\text{NH}_4)_2\text{SO}_3$		c	116, 117			
		aq	116, 117			
$(\text{NH}_4)_2\text{SO}_3 \cdot \text{H}_2\text{O}$		c	117, 121			
$(\text{NH}_4)_2\text{SO}_4$		c	32, 62		119	99, 119, 125
		aq	32, 62, 123			
$(\text{NH}_2\text{OH})_2 \cdot \text{H}_2\text{SO}_4$		c	82			
		aq	32, B2			
$(\text{NH}_4)_2\text{S}_2\text{O}_5$		aq	7			
$(\text{NH}_4)_2\text{S}_2\text{O}_8$		c	126			
		aq	7			
$(\text{N}_2\text{H}_4)_2\text{H}_2\text{SO}_4$		aq	74			
$\text{N}_4\text{Se}_4$		c	127			
$\text{NH}_4\text{HSe}$		aq	124			
$(\text{NH}_4)_2\text{Se}$		aq	7			
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3. Herzberg and Sponer-1			25. Berthelot-58			
4. Herzberg-9			26. Thomsen-14			
5. Gaydon-1			27. Giauque and Kemp-1			
6. van der Ziel-1			28. Kelley-27			
7. Calculated			29. Schumacker and Sprenger-2			
B. Giauque and Clayton-1			30. Brown, Smith, and Latimer-3			
9. Moore-1			31. Carlton-Sutton, Ambler, and Williams-1			
10. Bacher and Goudsmit-1			32. Thomsen-16			
11. Edlen-5			33. Sutton-3			
12. Cady-4			34. Awbery and Griffiths-2			
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14. Johnston and Davis-1			36. Kassel-2			
15. Giauque and Clayton-1			37. Badger and Woo-2			
16. Trautz and Ader-1			38. Blue and Giauque-1			
17. Hagstrum and Tate-1			39. Berthelot-106			
18. Worley and Jenkins-1			40. Latimer and Zimmerman-1			
19. Mulliken-2			41. Abel and Proisl-1			
20. Stille-1			42. Rideal-4			
21. Hanson-2			43. Wourtsel-1			
22. Johnston and Chapman-1			44. Bodenstein-5			

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Table 18 (Continued)  
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| 48. Villars-1                             | 90. Guntz-1                    |
| 49. Bates-6                               | 91. Noyes and Tuley-1          |
| 50. Becker and Roth-2                     | 92. Deville and Hautefeuille-1 |
| 51. Thompson-7                            | 93. Beeson and Yost-2          |
| 52. Stephenson and McMahon-1              | 94. Trautz and Schlueter-1     |
| 53. Zeise-1                               | 95. Jahn-2                     |
| 54. Overstreet and Giauque-1              | 96. Stephenson-3               |
| 55. Wilson, Jr.-1                         | 97. Nacken-3                   |
| 56. Osborne, Stimson, Sligh, and Cragoe-1 | 98. Streeck-1                  |
| 57. Haupt and Teller-1                    | 99. Ziegler and Messer-1       |
| 58. Calvet-3                              | 100. Troost-1                  |
| 59. Lewis and Randall-7                   | 101. Hieber and Woerner-2      |
| 60. Grimm-2                               | 102. Roth-25                   |
| 61. Pitzer-3                              | 103. Gilbert and Cobb-1        |
| 62. Roth and Zeumer-1                     | 104. Ellingson-1               |
| 63. Stephenson-1                          | 105. Berthelot-79              |
| 64. Latimer-1                             | 106. Trautz and Dalal-1        |
| 65. Hughes, Corruccini, and Gilbert-1     | 107. Blair, Brass, and Yost-1  |
| 66. Roth-17                               | 108. Andre-1                   |
| 67. Berthelot-110                         | 109. Troost-3                  |
| 68. Jirsa-4                               | 110. Roozeboom-7               |
| 69. Bushnell, Hughes, and Gilbert-1       | 111. Varet-2                   |
| 70. Bach-1                                | 112. Berthelot and Vieille-3   |
| 71. Sherman-1                             | 113. Randall and White-2       |
| 72. Gunther, Meyer, and Müller-Skjold-1   | 114. Berthelot-13              |
| 73. Eyster and Gillette-1                 | 115. Sabatier-1                |
| 74. Bach-1                                | 116. de Forcrand-16            |
| 75. Berthelot and Matignon-3              | 117. de Forcrand-5             |
| 76. Klemenc and Hayek-2                   | 118. Berthelot-4               |
| 77. Forsythe and Giauque-1,2              | 119. Shomate-3                 |
| 78. Berthelot-26                          | 120. Hartog-1                  |
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| 80. Rossini-4                             | 122. Gilbert and Bushnell-1    |
| 81. Latimer and Ahlberg-2                 | 123. Bouzat-1                  |
| 82. Berthelot-41                          | 124. Fabre-1                   |
| 83. Berthelot and Matignon-4              | 125. Nitta and Suenaga-1       |
| 84. Berthelot-9                           | 126. Berthelot-109             |
| 85. Berthelot-26                          | 127. Berthelot and Vieille-5   |
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SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
P		g	3, 5, 7		6	6
		c	1, 2, 4		9, 6, 12	8
pn+		g	10			
P <sub>2</sub>		g	7, 11		3, 6	3, 6
P <sub>4</sub>		g	12, 13, 14, 15, 16		7	7, 6
		liq	8, 17, 18, 19, 20			8
PO		g	3			
PO <sub>3</sub> <sup>-</sup>		aq	21, 4, 6	22, 23		
PO <sub>4</sub> <sup>---</sup>		aq	22			
P <sub>2</sub> O <sub>7</sub> <sup>----</sup>		aq	4, 6, 21			
P <sub>4</sub> O <sub>10</sub>		c	24, 4			
		amorp	4			
PH <sub>3</sub>		g	7, 25, 26, 27		7, 28, 29, 30 31	
PH <sub>3</sub> ·6H <sub>2</sub> O		c	32, 52			
P <sub>2</sub> H		c	25			
HPO <sub>3</sub>		c	33, 4			
		aq	33, 4			
HPO <sub>3</sub> <sup>--</sup>		aq	21, 6			
HPO <sub>4</sub> <sup>--</sup>		aq	34, 22	34, 35, 36		
H <sub>2</sub> PO <sub>3</sub> <sup>-</sup>		aq	21, 6			
H <sub>2</sub> PO <sub>4</sub> <sup>-</sup>		aq	35, 22	35	37	
H <sub>3</sub> PO <sub>2</sub>		c	21			
		liq	21			
		aq	21			
H <sub>3</sub> PO <sub>3</sub>		c	21			
		liq	21			
		aq	21			
H <sub>3</sub> PO <sub>4</sub>		c	21			
		aq	21, 4, 33, 38, 39, 22			
H <sub>3</sub> PO <sub>4</sub> ·1/2H <sub>2</sub> O		c	39			
		liq	39			
HP <sub>2</sub> O <sub>7</sub> <sup>---</sup>		aq	4, 6			
H <sub>2</sub> P <sub>2</sub> O <sub>7</sub> <sup>--</sup>		aq	21, 6, 4			
H <sub>3</sub> P <sub>2</sub> O <sub>7</sub> <sup>-</sup>		aq	21, 6, 4			
H <sub>4</sub> P <sub>2</sub> O <sub>5</sub>		aq	66			
H <sub>4</sub> P <sub>2</sub> O <sub>7</sub>		c	4			
		aq	4			
H <sub>4</sub> P <sub>2</sub> O <sub>7</sub> ·1 1/2H <sub>2</sub> O		c	65			
		liq	65			
PF <sub>3</sub>		g			7, 31, 40	

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
PCl <sub>3</sub>		liq	21,45			
		g	41		7,31,40	
PCl <sub>5</sub>		c	21			
		g	7		7	
POCl <sub>3</sub>		liq	21,42			
		g	43		7	
PH <sub>4</sub> Cl		g	1,3,4			
PBr <sub>3</sub>		g	46		7	
		liq	42			
	in CS <sub>2</sub>		2			
PBr <sub>5</sub>		c	47			
POBr <sub>3</sub>		c	48,47			
PH <sub>4</sub> Br		c	25,49			
PI <sub>3</sub>		c	42,47			
P <sub>2</sub> I <sub>4</sub>		c	47			
PH <sub>4</sub> I		c	25,50			51
PSCl <sub>3</sub>		g			7	
PN		g	54,55		53	53
1/n(PN) <sub>n</sub>		c	56			
P <sub>3</sub> N <sub>5</sub>		c	58			57
NH <sub>4</sub> H <sub>2</sub> PO <sub>4</sub>		c	62,61,59		63	63
		aq	59,60,61,62			
(NH <sub>4</sub> ) <sub>2</sub> HPO <sub>4</sub>		c	59,62,64			64
		aq	62,59			
(NH <sub>4</sub> ) <sub>3</sub> PO <sub>4</sub>		c	59			
		aq	59			
(NH <sub>4</sub> ) <sub>3</sub> PO <sub>4</sub> ·3H <sub>2</sub> O		c	59			
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6. Calculated			16. Centnerszwer-7			
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8. Young and Hildebrand-1			18. Desains-2			
9. Kelley-24			19. Petersen-5			
10. Moore-1			20. Tammann-3			

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| 23. Latimer-1                               | 65. Giran-2                      |
| 24. Roth, Meichsner, and Richter-1          | 66. Amat-1                       |
| 25. Ogier-1                                 |                                  |
| 26. Lemoult-10                              |                                  |
| 27. Ipat'v and Frost-1                      |                                  |
| 28. Stephenson and Giauque-1                |                                  |
| 29. Clusius and Frank-3                     |                                  |
| 30. Frank and Clusius-2                     |                                  |
| 31. Yost and Anderson-1                     |                                  |
| 32. de Forcrand and Taboury-1               |                                  |
| 33. Frandsen-1                              |                                  |
| 34. Bates and Acree-1                       |                                  |
| 35. Nims-1                                  |                                  |
| 36. Harned and Owen-1                       |                                  |
| 37. Stephenson-1                            |                                  |
| 38. Mixter-9                                |                                  |
| 39. Joly-1                                  |                                  |
| 40. Yost-1                                  |                                  |
| 41. Kelley-19                               |                                  |
| 42. Berthelot and Louguinine-2              |                                  |
| 43. Arai-4                                  |                                  |
| 44. Briner-1,3,4                            |                                  |
| 45. Fischer and Jüßermann-2                 |                                  |
| 46. van Driel and Gerding-2                 |                                  |
| 47. Ogier-4                                 |                                  |
| 48. Berger-1,2                              |                                  |
| 49. Johnson-2                               |                                  |
| 50. Smith and Calvert-1                     |                                  |
| 51. Crenshaw and Ritter-1                   |                                  |
| 52. Caillaudet and Bordet-1                 |                                  |
| 53. McCallum and Leifer-1                   |                                  |
| 54. Ghosh and Datta-1                       |                                  |
| 55. Curry, Herzberg, and Herzberg-2         |                                  |
| 56. Wettruff-1                              |                                  |
| 57. Sato-14                                 |                                  |
| 58. Stock and Wrede-1                       |                                  |
| 59. de Passillé and Séon-1                  |                                  |
| 60. Berthelot and Louguinine-3              |                                  |
| 61. Perreu-13                               |                                  |
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Table 20  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
As		g	10, 2, 17		2	2
		c	19, 21		16	16
		amorp	19, 20			
As <sup>n+</sup>		g	11, 12, 13, 14, 15			
As <sub>2</sub>		g	1, 2		3	4, 2
As <sub>4</sub>		g	1, 6, 7, 8, 9, 5		5	
AsO		g	18, 4, 2			
AsO <sup>+</sup>		aq		22		
AsO <sub>2</sub> <sup>-</sup>		aq		22		
AsO <sub>4</sub> <sup>---</sup>		aq	23, 2, 24	23		
As <sub>2</sub> O <sub>5</sub>		c	24		3, 16	
		aq	24			
As <sub>2</sub> O <sub>5</sub> ·4H <sub>2</sub> O		c	23			
3As <sub>2</sub> O <sub>5</sub> ·5H <sub>2</sub> O		c	23			
As <sub>4</sub> O <sub>6</sub>		c	32, 24, 26	26	3, 16	
		aq	24, 25			
As <sub>2</sub> O <sub>3</sub> ·As <sub>2</sub> O <sub>5</sub>		c	27, 28			
AsH <sub>3</sub>		g	29			
AsH <sub>3</sub> ·6H <sub>2</sub> O		c	30			
HAsO <sub>2</sub>		aq	24, 25	26		
HAsO <sub>3</sub> <sup>--</sup>		aq	24			
HAsO <sub>4</sub> <sup>--</sup>		aq	23, 24	22		
H <sub>2</sub> AsO <sub>3</sub> <sup>-</sup>		aq	24			
H <sub>2</sub> AsO <sub>4</sub> <sup>-</sup>		aq	23, 24	31		
H <sub>3</sub> AsO <sub>3</sub>		aq	24, 25			
H <sub>3</sub> AsO <sub>4</sub>		c	24			
		aq	24	22		
AsF <sub>3</sub>		g	33		33, 39	
		liq	34		33	33
AsCl <sub>2</sub>		g	35		38, 39	
AsCl <sub>3</sub>		g	24			
		liq	5, 36, 37	5, 36, 37		
AsBr <sub>3</sub>		c	40			
AsI <sub>3</sub>		c	40, 41			
As <sub>2</sub> S <sub>2</sub>		g	43, 44			
		c	27, 42			
As <sub>2</sub> S <sub>3</sub>		c	27, 42, 44			
As <sub>2</sub> O <sub>3</sub> ·SO <sub>3</sub>		c	27, 44			
AsN		g	4, 45			4, 2
NH <sub>4</sub> H <sub>2</sub> AsO <sub>4</sub>		c	23		46	
		aq	23			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$(\text{NH}_4)_2\text{HAsO}_4$		c	23			
		aq	23			
		c	23			
		aq	23			
		c	23			
		aq	23			
$(\text{NH}_4)_3\text{AsO}_4 \cdot 3\text{H}_2\text{O}$		c	23			

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5. Kelley-19	29. Ogier-1,2
6. Gibson-1	30. de Forcrand-50
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9. Horiba-2	33. Russell, Rundle, and Yost-1
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13. Rao-8	37. Baxter, Bezenberger, and Wilson-1
14. Kruger and Shoupp-2	38. Yost and Anderson-1
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16. Anderson-1	40. Berthelot-50
17. Almy and Kinzer-1	41. Mosnier-1
18. Jenkins and Strait-1	42. Britzke, Kapustinskiĭ, and Chentzova-2
19. Petersen-6	43. Britzke and Kapustinskiĭ-4
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21. Lashchenko-7	45. Gaydon-1
22. Latimer-1	46. Stephenson and Adams-1
23. de Passillé-1	
24. Thomsen-16	

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Table 21  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Sb		g	1,3		1,2	1
		c	5		4	4
Sb <sup>n+</sup>		g	6,7			
Sb <sub>2</sub>		g	8		1,2	1,2
Sb <sub>4</sub>		g	9,10			
SbO		g	3,12			
SbO <sup>+</sup>		aq		13		
Sb <sub>2</sub> O <sub>3</sub>		aq	15			
Sb <sub>2</sub> O <sub>4</sub>		c	14,19		4	4
Sb <sub>2</sub> O <sub>5</sub>		c	14		4	4
		aq	15			
Sb <sub>4</sub> O <sub>6</sub>		c	16,17,14	11	4	4
Sb <sub>6</sub> O <sub>13</sub>		c	19			
H <sub>3</sub> SbO <sub>4</sub>		aq	15			
SbF		g	3,12			
SbF <sub>3</sub>		c	16,18			
		aq	16,17,18			
H <sub>3</sub> SbF <sub>6</sub>		aq	16,18			
SbCl		g	3,12			
SbCl <sub>3</sub>		g	1		2,1	2,1
		c	14		1	
SbCl <sub>5</sub>		g	20			
		liq	14			
SbOCl		c	16,30			
Sb <sub>4</sub> O <sub>5</sub> Cl <sub>2</sub>		c	14,16,30			
SbBr <sub>3</sub>		c	5,16,18,22			
	in CS <sub>2</sub>		5			
SbI <sub>3</sub>		c	24			
		aq	25			
SbS <sub>3</sub> ----		aq	27,1			
Sb <sub>2</sub> S <sub>3</sub>		c	27,28,29			
		amorp	26			
Sb <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c	23			
SbN		g	3,12			
SbF <sub>3</sub> ·nNH <sub>3</sub>		c	21			

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Table 21 (Continued)  
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9. Niwa and Yoshiyama-1
10. Kelley-18
11. Schuhmann-3
12. Latimer-1
13. Herzberg-9
14. Mixter-9
15. Thomsen-16
16. Guntz-1
17. Guntz-18
18. Guntz-6
19. Simon and Thaler-1
20. Braune and Tiedje-1
21. Biltz and Rahlfs-1
22. Berthelot and Petit-1
23. Beck-1
24. Guntz-8
25. Mosnier-1
26. Berthelot-100
27. Fricke and Dönges-1
28. Britzke and Kapustinskii-3
29. Britzke and Kapustinskii-1
30. Guntz-7

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Table 22  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Bi		g	1,2,7,8		6	6
		c			5	5
Bi <sup>n+</sup>		g	10,11,12,13,14,15			
Bi <sub>2</sub>		g	1,2,3,4,6,7,8		4,6,9	4,6,9
BiO		g	3			
		c				
BiO <sup>+</sup>		aq		21		
Bi <sub>2</sub> O <sub>3</sub>		c	16,1B		5	5
BiH		g	3			
Bi(OH) <sub>3</sub>		c	16			
BiF		g	3			
BiCl		g	3		20	
BiCl <sub>3</sub>		g	B,6		22,6	22,6
		c	16		B,6	
		aq	19			
BiOCl		c	16,17	17		
BiBr		g	3		20	
BiI		g	3		20	
Bi <sub>2</sub> S <sub>3</sub>		c	23,24,25,26	23,24,25,26		
Bi <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c	27			

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5. Anderson-4	19. Tanatar-B
6. Calculated	20. Stevenson-2
7. Brewer et al-1	21. Latimer-1
B. Kelley-19	22. Kelley-24
9. Herzberg-9	23. Kelley-22
10. Bacher and Goudsmit-1	24. Britzke and Kapustinskiĭ-1
11. Crawford and McLay-1	25. Britzke and Kapustinskiĭ-3
12. Schoepfle-1	26. Jellinek and Zakowski-1
13. Mack and Fromer-1	27. Beck-1
14. Thorsen-3	



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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
C		g	4,2,7,8		1,5	1,5
		c	3,6,4		3,39,40,41	3,39,40,41
$C^{n+}$		g	1			
CO		g	9,10,11,12,13,14,4		3,5,42	3,5
$CO^+$		g	15,16,17,18,19			
$CO_2$		g	12,6,20,21		3,5,43	3,5
		aq	25,26,27,28,66	25		
$CO_2 \cdot 6H_2O$		c	22,23			
$CO_2^+$		g	24			
$CO_3^{--}$		aq	69,63	69		
CH		g	15			
$CH_3$		g	29,31,32,33,34,35			
$CH_4$		g	9,13,14,30		30,44,45,46	30
$CH_4 \cdot 6H_2O$		c	22			
$HCOO^-$	formate ion	aq	60,62,26,27,5	60		
$HCO_3^-$	bicarbonate ion	aq	63,5,64,65,66,67,68			
$CH_2O$	formaldehyde	g	70,71		72	72
		aq	73			
	in $CH_3OH$		73			
$CH_2O_2$	formic acid	g	26,36,49,51,52		47,49	
		liq	49,47,51,52,53,54,55,56		48,49	47,50
		aq	57,58,26,59	61		
$H_2CO_3$	carbonic acid	aq	66,25,26,27,28			
$CH_3OH$	methanol	g	74,75,76,77,4		89,5,90,91	
		liq	77,78,4,5		48,79,80	48,79,80
		aq	82,83,84	81		
$CF_4$	tetrafluoromethane	g	85,92,93		86,87	
$CH_3F$	fluoromethane	g			94	94
$CH_2F_2$	difluoromethane	g				94
$CHF_3$	trifluoromethane	g				94
$CCl_4$	tetrachloromethane	g	26,97,96		98,99,88,101,102	103
		liq	104,105,106,107,4			98
$COCl_2$	carbonyl chloride	g	26,111,112,113,114,115,116		108,110	110
$CH_3Cl$	chloromethane	g	26,117		109,102	103
$CH_3Cl \cdot 9H_2O$		c	118			
$CH_2Cl_2$	dichloromethane	g	123,4,5		102	103,121
		liq	119,5		122,123,4,5	120

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
CHCl <sub>3</sub>	trichloromethane	g	26,124,97		102	103
		liq	95,123,5,4		123,5,4	125
CF <sub>3</sub> Cl	trifluorochloromethane	g				126
CF <sub>2</sub> Cl <sub>2</sub>	difluorodichloromethane	g				126
CFCl <sub>3</sub>	fluorotrichloromethane	g			127	126,128
		liq				127,128
CH <sub>2</sub> FCl	fluorochloromethane	g				128,121
CHF <sub>2</sub> Cl	difluorochloromethane	g				129,128
CHFC1 <sub>2</sub>	fluorodichloromethane	g				128,129
CBr <sub>4</sub>	tetrabromomethane	g	5		102	102
		c				130
COBr <sub>2</sub>	carbonyl bromide	g	8			
		liq	133			
CH <sub>3</sub> Br	bromomethane	g	26,134		102,131	102
CH <sub>2</sub> Br <sub>2</sub>	dibromomethane	g	5		94,102	94,102
CHBr <sub>3</sub>	tribromomethane	g	5		102	102
		liq	95		5	
CF <sub>3</sub> Br	trifluorobromomethane	g				135,5
CF <sub>2</sub> Br <sub>2</sub>	difluorodibromomethane	g				135,5
CFBr <sub>3</sub>	fluorotribromomethane	g				135,5
CH <sub>2</sub> FBr	fluorobromomethane	g				135,5
CHF <sub>2</sub> Br	difluorobromomethane	g				135,5
CHFBBr <sub>2</sub>	fluorodibromomethane	g				121
CCl <sub>3</sub> Br	trichlorobromomethane	g				121
CCl <sub>2</sub> Br <sub>2</sub>	dichlorodibromomethane	g				121
CClBr <sub>3</sub>	chlorotribromomethane	g				121
CH <sub>2</sub> ClBr	chlorobromomethane	g				135,5
CHCl <sub>2</sub> Br	dichlorobromomethane	g				121
CHClBr <sub>2</sub>	chlorodibromomethane	g				121
CHFC1Br	fluorochlorobromomethane	g				135,5
CH <sub>3</sub> I	iodomethane	g	26		94	94
		liq	136,5		5	
CH <sub>2</sub> I <sub>2</sub>	diiodomethane	g				121
		liq	136			
CHI <sub>3</sub>	triiodomethane	c	136			
CS		g			15,5	15,5
1/n(CS) <sub>n</sub>		c	137			
CS <sub>2</sub>	carbon disulfide	g	26,140,141	140,141	138,141	141
		liq	97,139,5,4		138	138

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Formula	Description	State				
COS	carbon oxysulfide	g	140,141,26,111,4	140,141	141,142	141
CH <sub>4</sub> S	methanethiol	g	26		5,143,144	5,143,144
CSCl <sub>2</sub>	carbon sulfochloride	g			110	110
CSe <sub>2</sub>	carbon diselenide	liq	145			
CN		g			15,5,146	15,5
CN <sup>-</sup>	cyanide ion	aq	172,26,173,149	168,170		
CN <sub>4</sub>	cyanogen azide	c	147			
1/n(CN <sub>4</sub> ) <sub>n</sub>		c	147			
CNO <sup>-</sup>	cyanate ion	aq	161,164,5,169	165,166,167, 168,169		
C(NO <sub>2</sub> ) <sub>4</sub>	tetranitromethane	liq	148			
HCN	hydrogen cyanide	g	26,150,151		151,152,153	5,151,152, 153
		liq	149,5,4,152		152,87	152
	hydrocyanic acid	aq	149,171	168,170		
CH <sub>5</sub> N	methyl amine	g	26,154		159,160	158,160
		aq	155,156,157			
CH <sub>6</sub> N <sup>+</sup>	methyl ammonium ion	aq	156			
CH <sub>2</sub> N <sub>2</sub>	cyanamide	c	161,162			
		aq	161,162			
NH <sub>4</sub> CN	ammonium cyanide	c	149			
		aq	149			
CH <sub>5</sub> N <sub>3</sub>	guanidine	c	174			
		aq	174			
CH <sub>6</sub> N <sub>3</sub> <sup>+</sup>	guanidine ion	aq	174,5			
HCNO	cyanic acid	aq	149,161	168		
CH <sub>2</sub> O <sub>2</sub> N <sup>-</sup>	nitromethane ion	aq	177,182			
CH <sub>3</sub> ON	formamide	liq	176			
		aq	175			
CH <sub>3</sub> O <sub>2</sub> N	nitromethane	liq	26,177,123,178, 180,181		178,179	178
		aq	177			
CH <sub>5</sub> O <sub>2</sub> N	ammonium formate	c	183			
		aq	63,183,5			
NH <sub>4</sub> HCO <sub>3</sub>	ammonium bicarbonate	c	184			
		aq	5,184,185	5		
CH <sub>4</sub> ON <sub>2</sub>	urea	c	186,187,4		186,188	186,188
		aq	26,174,189,163, 190,191,192,194	193		
NH <sub>4</sub> CNO	ammonium cyanate	c	163			
		aq	163			
CH <sub>6</sub> O <sub>2</sub> N <sub>2</sub>	ammonium carbamate	c	196,197,198	198		
		aq	195,196			

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Formula	Description	State				
$(\text{NH}_4)_2\text{CO}_3$	ammonium carbonate	aq	63,5,26,185	63,5		
$\text{CHO}_6\text{N}_3$	trinitromethane	liq	180			
$\text{CH}_5\text{O}_4\text{N}_3$	urea nitrate	c	26,174			
$\text{CH}_2\text{O}_2\text{N}_4$	nitroguanidine	c	100			
$\text{CH}_6\text{O}_3\text{N}_4$	guanidine nitrate	c	174			
		aq	174			
$\text{CNCI}$	cyanogen chloride	g	149		199,5	199,5
$\text{CH}_6\text{NCI}$	methylamine hydrochloride	c			156	156
		aq	156,201,202			
$\text{CH}_6\text{ON}_3\text{Cl}$	semicarbazide hydrochloride	c				203
$\text{CN}_8\text{r}$	cyanogen bromide	g			199,5	199,5
$\text{CNI}$	cyanogen iodide	g	5		199,5	199,5
		c	149		5	
		aq	149	204		
$\text{CNS}^-$	thiocyanate ion	aq	205,5			
$\text{HCNS}$	thiocyanic acid	aq	205			
$\text{CH}_4\text{N}_2\text{S}$	thiourea	c	206,174			
		aq	174			
$\text{NH}_4\text{CNS}$	ammonium thiocyanate	c	174			
		aq	205			
$\text{CH}_5\text{O}_3\text{N}_3\text{S}$	thiourea nitrate	c	174			
$\text{CP}$		g			15,5	15,5
$\text{C}_2$		g	2,207,208,209,210, 211,212,213,214			
$\text{C}_2\text{O}_4^{--}$	oxalate ion	aq	5	5		
$\text{C}_2\text{H}_2$	ethyne (acetylene)	g	218,4	218	218,221	218,221
		aq	215,216			
$\text{C}_2\text{H}_2 \cdot 6\text{H}_2\text{O}$		c	217			
$\text{C}_2\text{H}_4$	ethene (ethylene)	g	218,225,4,226	218	218,222, 219,223	218,222, 223
$\text{C}_2\text{H}_4 \cdot 6\text{H}_2\text{O}$		c	217			
$\text{C}_2\text{H}_6$	ethane	g	218,227,4	218	218,224, 220,223	218,224, 223
$\text{HC}_2\text{O}_4^-$	bioxalate ion	aq	5			
$\text{C}_2\text{H}_2\text{O}$	ketene	g	229			
$\text{C}_2\text{H}_2\text{O}_2$	glyoxal	c	230			
		aq	230			
$\text{C}_2\text{H}_2\text{O}_4$	oxalic acid	c	231,4,232,233		234	235,236
		aq	26,57,238			
	in $\text{CH}_3\text{OH}$		237			
	in $\text{C}_2\text{H}_5\text{OH}$		237			
	in $\text{C}_3\text{H}_7\text{OH}$		237			

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Formula	Description	State				
$C_2H_2O_4 \cdot 2H_2O$	oxalic acid dihydrate	c	233, 237, 26, 238, 239, 240, 241			
$C_2H_3O_2^-$	acetate ion	aq	242, 243			
$C_2H_3O_3^-$	glycollate ion	aq	244, 5			
$C_2H_4O$	acetaldehyde	g	245, 26, 4		246	
		aq	247			
$C_2H_4O$	ethylene oxide	g	248, 26, 249, 253		250, 251	252
$C_2H_4O_2$	acetic acid	liq	254, 255, 256		234	257
		aq	26, 258, 259, 260, 261, 262, 263, 264, 265			
	in $H_2SO_4$		266			
	in $CH_3COCH_3$		266			
	in $C_2H_5OC_2H_5$		266			
	in $C_5H_{12}$		266			
	in $C_6H_6$		266			
$C_2H_4O_2$	methyl formate	g	26, 267			
		liq	268, 5			269
		aq	267			
$C_2H_4O_3$	hydroxyacetic acid (glycollic acid)	c	270			
		aq	244			
$C_2H_4O_4$	dihydroxyacetic acid (glyoxylic acid)	c	271			
		aq	272			
$C_2H_4O_4$	formic acid dimer	g	26, 36, 49, 51, 52		49, 48, 47	
$C_2H_6O$	ethanol	g	4, 75, 77		281, 5	
		liq	5, 77		279, 280	279, 280
		aq	265, 83, 273, 274, 275, 276			
	in $H_2SO_4$		266			
	in $C_6H_6$		277			
$C_2H_6O$	dimethyl ether	g	26, 278		282	283
		aq	278			
$C_2H_6O_2$	1,2-ethanediol (ethylene glycol)	liq	284, 285, 253, 288		280	
		aq	286			
$C_2H_6O_2$	ethyl hydrogen peroxide	liq	287			
$C_2F_6$	hexafluoroethane	g	290			
$C_2H_3F_3$	1,1,1-trifluoroethane	g			291	
$C_2H_3O_2F$	fluoroacetic acid	c	292			
$C_2H_5OF$	2-fluoroethanol	liq	292, 293			
$C_2H_2O_2F_2$	difluoroacetic acid	liq	292			
$C_2H_4OF_2$	2,2-difluoroethanol	liq	292, 293			
$C_2Cl_4$	tetrachloroethene	liq	26			294
$C_2Cl_6$	hexachloroethane	c	97, 5			236

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$C_2O_2Cl_3^-$	trichloroacetate ion	aq	5, 26, 295			
$C_2OCl_4$	trichloroacetyl chloride	liq	296			
$C_2H_3Cl$	chloroethene	g	5, 26			
$C_2H_5Cl$	chloroethane	g	26, 97		250, 297	236
$C_2H_2Cl_2$	cis-1,2-dichloroethene	liq				269
$C_2H_2Cl_2$	trans-1,2-dichloroethene	liq				269
$C_2H_4Cl_2$	1,1-dichloroethane	liq	26, 5, 267			
$C_2H_4Cl_2$	1,2-dichloroethane	liq	95, 299, 26		298	298
$C_2HCl_3$	trichloroethene	liq				294
$C_2H_3Cl_3$	1,1,1-trichloroethane	liq			300	300
$C_2H_2Cl_4$	1,1,2,2-tetrachloroethane	liq				294
$C_2HCl_5$	pentachloroethane	liq				294
$C_2H_2O_2Cl^-$	chloroacetate ion	aq	265, 295			
$C_2H_3OCl$	acetylchloride	liq	259, 296, 301			236
$C_2H_3OCl$	chloroacetaldehyde	liq	296			
$C_2H_3O_2Cl$	chloroacetic acid	c	254, 302, 303, 305			303
		aq	295, 303			
$C_2H_5OCl$	2-chloroethanol	liq	304, 306			
$C_2H_7OCl$	dimethyl ether hydrochloride	g	307, 308, 309	307, 308, 309		
$C_2HO_2Cl_2$	dichloroacetate ion	aq	26, 5			
$C_2H_2OCl_2$	chloroacetyl chloride	liq	296			
$C_2H_2O_2Cl_2$	dichloroacetic acid	liq	254			303, 295
		aq	303, 295			
$C_2HOCl_3$	trichloroacetaldehyde (chloral)	liq	310, 311			310
		aq	310, 311			
$C_2HO_2Cl_3$	trichloroacetic acid	c	254, 305			303
		aq	295, 303, 266			
	in $C_2H_5OH$		266			
	in $CH_3COOCH_3$		266			
	in $C_2H_5OC_2H_5$		266			
	in $C_5H_{12}$		266			
$C_2H_3O_2Cl_3$	trichloroacetaldehyde monohydrate	c	310, 311, 312, 313			310
		aq	5			
	in $CHCl_3$		312			
$C_2F_3Cl_3$	1,1,2-trifluoro-1,2,2-trichloroethane	liq				5
$C_2H_5Br$	bromoethane	g	26, 314			5
		liq	5			236
$C_2H_4Br_2$	1,2-dibromoethane	liq	95, 299, 304, 316		315	315

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$C_2H_3OBr$	acetyl bromide	liq	259			
$C_2H_3O_2Br_3$	tribromoacetaldehyde monohydrate	c	316			
		aq	316			
$C_2H_4ClBr$	1-chloro-2-bromoethane	liq				317
$C_2I_4$	tetraiodoethene	c	136			
$C_2H_5I$	iodoethane	liq	26,136			236
$C_2H_4I_2$	1,2-diiodoethane	c	136,318,319			
$C_2H_3OI$	acetyl iodide	liq	259			
$C_2H_4BrI$	1-bromo-2-iodoethane	c	314			
$C_2H_6S$	ethanethiol	liq	26,289			
$C_2H_6S$	dimethyl sulfide	g	26		320,321	
		liq	321		321	321
$C_2H_6OS$	dimethyl sulfoxide	liq	322			
$C_2H_6O_2S$	dimethyl sulfone	c	322			
$C_2H_6O_4S$	ethylsulfuric acid	aq	83			
$C_2N_2$	cyanogen	g	200,323,26,324,325		326	327,199,72
$C_2H_3N$	acetonitrile	g	26,161		72,328	72,328
		liq	5		5	
$C_2H_3N$	methyl isocyanide	g	5		328	328
		liq	329,330			
$C_2H_7N$	ethylamine	g	26,162,331			332
		aq	331,333,334			
$C_2H_7N$	dimethylamine	g	26,154,338		335	332
		aq	336			
$C_2H_8N^+$	ethylammonium ion	aq	5			
$C_2H_8N^+$	dimethylammonium ion	aq	5			
$C_2H_8N_2$	1,2-ethanediamine (ethylenediamine)	liq	339,340			
		aq	340			
$C_2H_8N_2 \cdot H_2O$	1,2-ethanediamine monohydrate	c	340			
$C_2H_4N_4$	dicyandiamide	c	161,341			236
		aq	161			
$C_2H_2O_3N^-$	oxamate ion	aq	174,5			
$C_2H_3ON$	methyl isocyanate	liq	342			
$C_2H_3ON$	glycollonitrile	liq	343			
$C_2H_3O_3N$	oxamic acid	c	174,344			
		aq	174			
$C_2H_4O_2N^-$	aminoacetate ion	aq	345,5,295			
$C_2H_4O_2N^-$	nitroethane ion	aq	271,182			
$C_2H_5ON$	acetamide	c	176,175,346			348
		aq	175,176,346,191			
	in $C_2H_5OH$		191			

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Formula	Description	State				
$C_2H_5ON$	acetaldehyde oxime	c	349			
$C_2H_5O_2N$	aminoacetic acid (glycine)	c	350,351,352		188	188
		aq	295,353,345,354, 355,356			
$C_2H_5O_2N$	nitroethane	liq	271,26,5			271
		aq	5,182			
$C_2H_5O_2N$	ethyl nitrite	g	26			
$C_2H_5O_3N$	ethyl nitrate	liq	26,5			
$C_2H_7O_2N$	ammoniumacetate	c	348			
		aq	26,358,359,360,175			
$C_2H_7O_3N$	methylammonium bicarbonate	aq	361			
$C_2H_7O_3N$	ammonium glycollate	c	244			
		aq	244			
$C_2H_5O_4N$	ammonium bioxalate	c				235
		aq	348			
$C_2H_6O_2N^+$	glycinium ion	aq	354			
$C_2H_7O_4N$	ammonium glyoxylate	aq	174			
$C_2H_4O_2N_2$	glyoxime	c	362			
$C_2H_4O_2N_2$	oxamide	c	344			
$C_2H_4O_2N_2$	formylurea	c	174			
		aq	174			
$C_2H_4O_6N_2$	glycol dinitrate	liq	363			
$C_2H_6ON_2$	dimethylnitrosamine	liq	181			
$C_2H_6O_2N_2$	ethylnitramine	liq	364,365			
$C_2H_8O_4N_2$	ammonium oxalate	c	348,5			235,366
		aq	348,233			
$C_2H_8O_4N_2 \cdot H_2O$	ammonium oxalate monohydrate	c	348,233			
$C_2H_7O_4N_5$	guanylurea nitrate	c	341			
$C_2H_5F_2N$	2,2-difluoroethylamine	liq	292,293			
$C_2H_4OFN$	fluoroacetamide	c	292,293			
$C_2H_3OF_2N$	difluoroacetamide	c	292,293			
$C_2H_4O_2F_2N_2$	2,2-difluoroethyl- nitramine	liq	292,293			
$C_2H_8ClN$	ethylamine hydro- chloride	aq	331,26			
$C_2H_8ClN$	dimethylamine hydrochloride	aq	26,201,202			
$C_2H_{10}Cl_2N_2$	1,2-ethanediamine dihydrochloride	c	339,340			
		aq	367,368,369,370, 371,340,339			
$C_2H_4OCIN$	chloroacetamide	c	357,254,364			
		aq	175			



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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$C_2H_6O_2ClN$	glycine hydrochloride	aq	5			
$C_2H_6O_2ClN$	ammonium chloroacetate	aq	357,5			
$C_2H_3OCl_2N$	dichloroacetamide	aq	175			
$C_2H_2OCl_3N$	trichloroacetamide	c	364,365,254			
		aq	175			
$C_2H_4O_2Cl_3N$	ammonium trichloroacetate	aq	357,5			
$C_2H_3SN$	methyl thiocyanate	liq	26,5,372			
$C_2H_3SN$	methyl isothiocyanate	c	26,5,372			
$C_2H_7O_3SN$	2-aminoethylsulfonic acid (taurine)	c	97		373	373
		aq	355			

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7. Herzberg-2	26. Thomsen-16
8. Gaydon-1	27. Berthelot-7
9. Roth and Banse-1	28. Adolph and Henderson-1
10. Awbery and Griffiths-2	29. Kistiakowsky and van Artsdalen-1
11. Fenning and Cotton-1	30. Rossini, Pitzer, Taylor, Ebert, Kilpatrick, Seckett, Williams, and Werner-1
12. Rossini-16	31. Andersen and Kistiakowsky-1
13. Rossini-4	32. Stevenson-3
14. Rossini-23	33. van Artsdalen-1
15. Herzberg-9	34. Andersen, Kistiakowsky, and van Artsdalen-1
16. Savard-1	35. Ruedy-1
17. Savard and de Hemptinne-2	36. Favre and Silbermann-1
18. Savard, de Hemptinne, and Capron-1	37. Berthelot and Matignon-5,12
19. Hagstrum and Tate-1	38. Jahn-3

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| 46. Storch-1                                  | BB. Yost-1                                 |
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| 48. Parks, Kelley, and Huffman-1              | 90. Halford and Pecherer-1                 |
| 49. Halford-3                                 | 91. Kassel-9                               |
| 50. Radulescu and Jula-1                      | 92. von Wartenberg-21                      |
| 51. Ramsperger and Porter-2                   | 93. Ruff and Bretschneider-2               |
| 52. Coolidge-1                                | 94. Edgell and Glockler-1                  |
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| 55. Ogier-5                                   | 97. Berthelot-110                          |
| 56. Favre and Silbermann-10                   | 98. Hicks, Hooley, Stephenson-1            |
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| 58. Faucon-1                                  | 100. Stull-1                               |
| 59. Glagoleva-1                               | 101. Yost and Blair-1                      |
| 60. Harned and Embree-1                       | 102. Stevenson and Beach-1                 |
| 61. Kaye and Parks-1                          | 103. Vold-1                                |
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| 66. Roughton-1                                | 108. Giauque and Jones-1                   |
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| 76. Roth and Banse-2                          | 118. Villard and de Forcrand-1             |
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| 80. Kelley-4                                  | 122. Dzung-1                               |

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| 127. Osborne, Garner, Doescher, and Yost-1     | 169. Walker-2                       |
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| 133. Schumacher and Lenher-1                   | 175. Calvet-1,3                     |
| 134. Berthelot-118                             | 176. Stohmann and Schmidt-4         |
| 135. Glockler and Edgell-2                     | 177. Berthelot and Matignon-9       |
| 136. Berthelot-123                             | 178. Jones and Giauque-1            |
| 137. Dewar-2                                   | 179. Pitzer and Gwinn-1             |
| 138. Brown and Manov-1                         | 180. Holcomb and Dorsey-1           |
| 139. Berthelot-73                              | 181. Swietoslawski-6                |
| 140. Terres and Wesemann-1                     | 182. Turnbull and Maron-1           |
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| 149. Berthelot-25                              | 191. Campbell and Campbell-2        |
| 150. Berthelot-76                              | 192. Gucker and Pickard-1           |
| 151. Gordon-5                                  | 193. Parks and Huffman-4            |
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| 154. Muller-4                                  | 196. Raabe-1                        |
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| 164. Berthelot-111                             | 206. Becker and Roth-6              |

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Si		g	1,2,3,4,5		1,5	1,5
		c			1,8,9	1,8,9
		amorp	10			
Si <sup>n+</sup>		g	6,7			
SiO		g	13,14		1,12	1,12
SiO <sub>2</sub>		c	10,15,16,18,19,21, 22,54		22,23	23,24
		gls	20,25,26		24	24
		aq	26			
SiH <sub>4</sub>		g	10,27,28		24	1,24
H <sub>2</sub> SiO <sub>3</sub>		c	29,26			
H <sub>4</sub> SiO <sub>4</sub>		c	29,26			
SiF <sub>4</sub>		g	30,31,32		24,33	1,24
SiF <sub>6</sub> <sup>--</sup>		aq	1,46			
H <sub>2</sub> SiF <sub>6</sub>		aq	18,25,26,31,32,46, 53			
SiCl <sub>4</sub>		g	40,41,42,43,44		24,38,39	37
		liq	11,34,35,36		24	24
SiBr <sub>4</sub>		liq	35,45			
SiI <sub>4</sub>		c	35			
SiS <sub>2</sub>		c	51			
Si <sub>3</sub> N <sub>4</sub>		c	17,47,52			
(NH <sub>4</sub> ) <sub>2</sub> SiF <sub>6</sub>		c	46			
		aq	46			
SiC		c	19,30,48,49,50		50	50
Si(C <sub>2</sub> H <sub>5</sub> O) <sub>4</sub>	tetraethoxy silane	liq	27			

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| 37. Herman-1                       |  |
| 38. Yost-1                         |  |
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Table 25  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Ge		g	1		2, 3	2, 3
		c	6, 1		1, 4, 6	1, 4, 6
Ge <sup>n+</sup>		g	5, 7			
GeO		g	9, 10, 11, 12		3, 8	3, B
GeO <sub>2</sub>		amorp	13, 14			
		aq	15, 16			
GeH <sub>4</sub>		g			3, 17	3, 17
H <sub>2</sub> GeO <sub>3</sub>		aq	15, 16			
GeCl		g	3, 12, 18			3, 12, 1B
GeCl <sub>4</sub>		liq	19			
GeBr		g	3, 12, 1B			3, 12, 18
GeS		g	3, 12, 20			3, 12, 20
GeSe		g				3, 21
GeTe		g				3, 21
Ge <sub>3</sub> N <sub>4</sub>		c	14, 22			
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6. Cristescu and Simon-1			20. Shapiro, Gibbs, and Laubengayer-1			
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8. Barrow-3			22. Morey and Johnson-1			
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10. Jevons, Bashford, and Briscoe-2						
11. Shaw-1						
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13. Becker and Roth-3						
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Table 26  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Sn		g	1,2,3,4			
		c	6,7,8		5	5
$\text{Sn}^{n+}$		g	9,10			
SnO		g			1,9	1,9
		c	11,13,14	12,13	5,15	5,15
$\text{SnO}_2$		c	14,17,18,19,20,21,22		5,15	5,15
$\text{Sn}(\text{OH})^+$		aq		11		
$\text{HSnO}_2^-$		aq		11		
$\text{Sn}(\text{OH})_2$		c	23	11		
$\text{Sn}(\text{OH})_4$		c	23			
$\text{SnF}_6^{--}$		aq	23,1			
$\text{H}_2\text{SnF}_6$		aq	23			
$\text{SnCl}_2$		c	23,29,30			
		aq	23,25,26,27			
$\text{SnCl}_2 \cdot 2\text{H}_2\text{O}$		c	23,28			
$\text{SnCl}_4$		liq	23,30		5	5
		aq	23,31			
$\text{SnOCl}_2$		aq	32			
$\text{SnBr}_2$		c	30			
		aq	30,29			
$\text{SnBr}_4$		c	30			
		aq	30,33,34			
$\text{SnBr}_4 \cdot 8\text{H}_2\text{O}$		liq	33			
$\text{SnI}_2$		c	30			
		aq	35,29			
SnS		c	36,37,38,39,40,41	40,41		
$\text{Sn}(\text{SO}_4)_2$		c	42			
$\text{SnX}_2 \cdot n\text{NH}_3$		c	29			
$\text{SnCl}_4 \cdot 1 \frac{1}{2} \text{PH}_3$		c	43			
$\text{SnH}_{20}\text{C}_8$	tetraethyltin	liq	44			
$\text{SnH}_{28}\text{C}_{12}$	tetra-n-propyltin	liq	44			
$\text{SnH}_{36}\text{C}_{16}$	tetra-n-butyltin	liq	44			
$\text{SnH}_{44}\text{C}_{20}$	tetra-n-amyltin	liq	44			

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| 14. Mixter-7                    |  |
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| 18. Eastman and Robinson-1      |  |
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| 21. Meyer and Scheffer-2        |  |
| 22. Kelley-18                   |  |
| 23. Thomsen-16                  |  |
| 24. Latimer-1                   |  |
| 25. Noyes and Chow-1            |  |
| 26. Kapustinskiĭ-8              |  |
| 27. Kapustinskiĭ-2              |  |
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| 29. Biltz and Fischer-2         |  |
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| 31. Berthelot-24                |  |
| 32. Berthelot-155               |  |
| 33. Pickering-13                |  |
| 34. Herschkowitsch-2            |  |
| 35. Mosnier-1                   |  |
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| 37. Britzke and Kapustinskiĭ-3  |  |
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| 42. Beck-1                      |  |

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Table 27  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Pb		g	1,2,3	1,2,3		3,4
		c			3	3
Pb <sup>n+</sup>		g	5,6			
Pb <sup>++</sup>		aq	7,8	9,10		
Pb <sub>2</sub>		g	11			
PbO		g			12	
		c	13,16,17,21	13,14,15,18,19	12,17	12
PbO <sub>2</sub>		c	21	17,20,22		12
Pb <sub>2</sub> O		c	23			
Pb <sub>3</sub> O <sub>4</sub>		c	25,26,27	17,24	12	12
HPbO <sub>2</sub> <sup>-</sup>		aq		10,28		
Pb(OH) <sub>2</sub>		c	10,29		30	
PbF <sub>2</sub>		c	31,32	10,33		
		aq	34			
PbF <sub>4</sub>		c	32			
PbCl <sub>2</sub>		c	37,38	15,35,36		37
		aq	7,8,34		34	
PbCl <sub>2</sub> ·nPbO		c	39			
PbBr <sub>2</sub>		c	42	15,36,38,41,43	12	12
		aq	34		34	
PbBr <sub>2</sub> ·nPbO		c	39			
PbI <sub>2</sub>		c	15,34,36,44	15,36,44		
		aq	34	15,45		
PbI <sub>3</sub> <sup>-</sup>		aq		45		
PbI <sub>4</sub> <sup>--</sup>		aq		45		
PbI <sub>2</sub> ·HI·5H <sub>2</sub> O		c	46			
PbS		c	47,48,50,51	49,50	12	12
PbSO <sub>4</sub>		c	20,34		9,12	12
PbS <sub>2</sub> O <sub>3</sub>		c	52			
PbS <sub>3</sub> O <sub>6</sub>		c	52			
		aq	52			
PbSO <sub>4</sub> ·nPbO		c	53,54,55			
PbSe		c	56			
PbSeO <sub>4</sub>		c	57			
PbTe		c	56			
Pb(N <sub>3</sub> ) <sub>2</sub>		c	58			
Pb(NO <sub>3</sub> ) <sub>2</sub>		c	8			
		aq	34,59,60,61,67		34	
Pb(OH)NO <sub>3</sub>		c	67			
PbCl <sub>2</sub> ·nNH <sub>3</sub>		c	62			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$2\text{PbCl}_2 \cdot \text{NH}_4\text{Cl}$		c	63			
$\text{PbBr}_2 \cdot n\text{NH}_3$		c	62			
$\text{PbI}_2 \cdot n\text{NH}_3$		c	62			
$3\text{PbI}_2 \cdot 4\text{NH}_4\text{I}$		c	64			
$3\text{PbI}_2 \cdot 4\text{NH}_4\text{I} \cdot 6\text{H}_2\text{O}$		c	64			
$\text{PbSO}_4 \cdot (\text{NH}_4)_2\text{SO}_4$		c	65			
$\text{Pb}_3(\text{PO}_4)_2$		c	61	10		12
$\text{PbHPO}_3$		c	66			
$3\text{PbI}_2 \cdot \text{PI}_3$		c	64			
$3\text{PbI}_2 \cdot \text{PI}_3 \cdot 12\text{H}_2\text{O}$		c	64			
$3\text{PbI}_2 \cdot \text{AsI}_3$		c	64			
$3\text{PbI}_2 \cdot \text{AsI}_3 \cdot 12\text{H}_2\text{O}$		c	64			
$3\text{PbI}_2 \cdot \text{SbI}_3$		c	64			
$3\text{PbI}_2 \cdot \text{SbI}_3 \cdot 12\text{H}_2\text{O}$		c	64			
$\text{PbCO}_3$		c	7, 67, 69	68	70	12, 70
$\text{PbC}_2\text{O}_4$	lead oxalate	c	71			
$\text{PbO} \cdot \text{PbCO}_3$		c	69, 72		12	12
$2\text{PbO} \cdot \text{PbCO}_3$		c	72		34	
$\text{Pb}(\text{C}_2\text{H}_5)_4$	tetraethyllead	liq	73			
$\text{Pb}(\text{CHO}_2)_2$	lead formate	c	74			
		aq	74			
$\text{Pb}(\text{C}_2\text{H}_3\text{O}_2)_2$	lead acetate	c	74			
		aq	8, 51			
$\text{Pb}(\text{C}_2\text{H}_3\text{O}_2)_2 \cdot 3\text{H}_2\text{O}$		c	74, 75			
$\text{Pb}(\text{C}_2\text{H}_3\text{O}_3)_2$	lead glycollate	c	76			
		aq	76			
$\text{Pb}(\text{CN})_2 \cdot 2\text{PbO} \cdot \text{H}_2\text{O}$		c	77			
$\text{Pb}(\text{CNS})_2$		c	77			
$\text{PbSiO}_3$		c	78		34	
		amorp	78			
$\text{Pb}_2\text{SiO}_4$		c	78		34	
		amorp	78			
$\text{PbI}_2 \cdot \text{SnI}_2$		c	64			
$\text{PbI}_2 \cdot \text{SnI}_2 \cdot 8\text{H}_2\text{O}$		c	64			

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Table 27 (Continued)  
SPECIFIC REFERENCES

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5. Bacher and Goudsmit-1	47. Korshunov-2
6. Earls and Sawyer-1	48. Britzke and Kapustinskiĭ-1
7. Zeumer and Roth-4	49. Kapustinskiĭ and Makolkin-3
8. Thomsen-16	50. Watanabe-7
9. Kelley-22	51. Berthelot-13
10. Latimer-1	52. Fogh-1
11. Gaydon-1	53. Schenk and Rassback-1,2
12. Kelley-24	54. Schenk and Borkenstein-1
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18. Garrett, Vellenga, and Fontana-1	60. Plake-1
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21. Tscheltzow-1	63. Bronsted-9
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25. Krustinsons-1	67. Roth-6
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27. Biltz-15	69. Marshall and Bruzs-1
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36. Gerke-1	78. Nacken-4
37. Wachter-2	
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Table 28  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Ga		g		1,2	3	2
		c				5
Ga <sup>n+</sup>		g	6			
Ga <sup>+++</sup>		aq	8,9	3,7		
GaO		g	10,11			
GaO <sub>3</sub> <sup>---</sup>		aq		12		
Ga <sub>2</sub> O		c	13			
Ga <sub>2</sub> O <sub>3</sub>		c	13,14			15
Ga(OH) <sup>++</sup>		aq		12		
HGaO <sub>3</sub> <sup>--</sup>		aq		12		
Ga(OH) <sub>2</sub> <sup>+</sup>		aq		12		
H <sub>2</sub> GaO <sub>3</sub> <sup>-</sup>		aq		12		
Ga(OH) <sub>3</sub>		c		12		
GaCl		g	11,16			
GaCl <sub>3</sub>		c	17			
		aq	7,9	3	3	
GaBr		g	11			
GaBr <sub>3</sub>		c	9			
		aq	17			
GaI		g	11			
GaI <sub>3</sub>		c	9			
		aq	11			
Ga <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c				15
GaN		c	18			
GaX <sub>3</sub> ·nNH <sub>3</sub>		c	9			
Ga <sub>2</sub> (C <sub>2</sub> O <sub>4</sub> ) <sub>3</sub>	gallium oxalate	c	19			
Ga <sub>2</sub> (C <sub>2</sub> O <sub>4</sub> ) <sub>3</sub> ·nH <sub>2</sub> O		c	19			
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Table 29  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
In		g	3	1	2	3
		c			5,6	4
In <sup>n+</sup>		g	2			
In <sup>+++</sup>		aq	7,8	8	3	3
InO		g	9			
In <sub>2</sub> O <sub>3</sub>		c	10,11			12
InH		g	9	3	3,13,14	3
In(OH) <sup>++</sup>		aq		7		
In(OH) <sub>3</sub>		c	3,15	15	3,15	
InCl		g	9	3	3,13,14	3
		c	16			
InCl <sub>2</sub>		c	16			
InCl <sub>3</sub>		c	16			
		aq	8,17			
InBr		g	9	3	3,13,14	3
InBr <sub>3</sub>		c	17			
		aq	3			
InI		g	9	3	3,13,14	3
InI <sub>3</sub>		c	17			
		aq	3			
In <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c	18			12
InN		c	19			
InX <sub>3</sub> ·nNH <sub>3</sub>		c	17			
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9. Gaydon-1						
10. Ditte-6						
11. Becker and Roth-3						
12. Nilson and Pettersson-1						
13. Herzberg-9						
14. Stevenson-2						



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Table 30  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Tl		g	1	1,2,3		4
		c			5	6
Tl <sup>n+</sup>		g	7,8			
Tl <sup>+</sup>		aq	4,9	9,10,11		
Tl <sup>+++</sup>		aq	4	10		
Tl <sub>2</sub> O		c	12,13,14	10		
TlH		g	15		16	
TlOH		c	12	10		
		aq	12			
Tl(OH) <sub>3</sub>		c	17			
TlF		g	15			
		aq	18,19			
HTlF <sub>2</sub>		aq	19			
TlCl		g	15		16,21	4
		c	12,14,24,25	21,22,23	5,23	
		aq	9,14,24	9,10		
TlCl <sub>3</sub>		c	26			
		aq	4			
TlCl <sub>3</sub> ·4H <sub>2</sub> O		c	26			
TlBr		g	15		16,20	20
		c	14	25,27,28		
		aq	4	25,27,28		
TlBr <sub>3</sub>		c	14			
TlBr <sub>3</sub> ·4H <sub>2</sub> O		c	26			
TlBrO <sub>3</sub>		c	25			
		aq	4	4		
TlClBr <sub>2</sub>		aq	26			
TlClBr <sub>2</sub> ·4H <sub>2</sub> O		c	26			
TlI		c	14,15	27,29,30	16,20	20
		aq	4	4		
TlIO <sub>3</sub>		c		26		
		aq	4	4		
Tl <sub>2</sub> S		c	10,14			
Tl <sub>2</sub> SO <sub>4</sub>		c	12,14,31			
		aq	4	4		
TlHSO <sub>4</sub>		aq	12,14			
Tl <sub>2</sub> Se		c	32			
Tl <sub>2</sub> Te		c	32			
TlN <sub>3</sub>	thallous azide	c	33			
TlNO <sub>3</sub>		c	11,14		4	
		aq	11,14		4	

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
TlCl·3NH <sub>3</sub>		c	34			
TlCl <sub>3</sub> ·3NH <sub>3</sub>		c	34, 35			
TlBr·3NH <sub>3</sub>		c	34			
TlI·3NH <sub>3</sub>		c	34			
Tl(C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> )	thallous acetate	aq	18			
TlOC <sub>2</sub> H <sub>5</sub>	thallous ethoxide	liq	18			
	in C <sub>2</sub> H <sub>5</sub> OH		18			
TlONC	thallous fulminate	c	33			

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19. Petersen-1	
20. Stevenson-2	

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Zn	in Hg	g	1, 2		3	3
		c			4	4
			5	5		
Zn <sup>n+</sup>		g	6			
Zn <sup>++</sup>		aq	3	3		
ZnO		c	7, 9, 10, 11, 12, 13	8, 13	4	4
ZnO <sub>2</sub> ·2H <sub>2</sub> O		c	14			
Zn <sub>3</sub> O <sub>5</sub> ·nH <sub>2</sub> O		c	14			
ZnH		g	16		3, 15	3, 15
Zn(OH) <sub>2</sub>		c	11, 17, 18, 19			7, 34
ZnF <sub>2</sub>		aq	20, 17, 21, 19, 22			
ZnCl		g	16		3, 15	3, 15
ZnCl <sub>2</sub>		c	24, 26, 18, 39, 40, 29, 28, 27	26, 28, 27		34
		aq	18, 23, 25, 5, 30, 31, 32, 33, 35, 36, 37, 38, 7, 44	25, 5		
	in C <sub>2</sub> H <sub>5</sub> OH		39			
ZnCl <sub>2</sub> ·3ZnO·5H <sub>2</sub> O		c	41			
ZnCl <sub>2</sub> ·4ZnO·11H <sub>2</sub> O		c	41			
ZnCl <sub>2</sub> ·5ZnO·8H <sub>2</sub> O		c	41			
ZnCl <sub>2</sub> ·8ZnO·10H <sub>2</sub> O		c	41			
ZnBr		g			42, 3	42, 3
ZnBr <sub>2</sub>		c	18, 28	28		
		aq	43	43		
ZnBr <sub>2</sub> ·2H <sub>2</sub> O		c	45	45		
ZnBr <sub>2</sub> ·4ZnO·13H <sub>2</sub> O		c	41			
ZnI		g	16		15, 3, 42	15, 3
ZnI <sub>2</sub>		c	28, 46	28, 46		
		aq	47, 48, 3, 49	3		
ZnI <sub>2</sub> ·5ZnO·11H <sub>2</sub> O		c	50			
ZnS		g	16			
		c	18, 51, 57, 54, 58, 59, 60, 55, 56	57	61, 4	34
ZnSO <sub>4</sub>		c	64, 18, 66, 67, 71	64		68
		aq	18, 14, 62, 63, 4, 65, 69, 70, 71	4		
ZnSO <sub>4</sub> ·nH <sub>2</sub> O		c	18, 66, 72, 63, 73, 74, 75	64		34
ZnS <sub>2</sub> O <sub>4</sub>		aq	3			
ZnS <sub>2</sub> O <sub>6</sub>		aq	3			
ZnS <sub>2</sub> O <sub>6</sub> ·6H <sub>2</sub> O		c	18			
Zn(HSO <sub>3</sub> ) <sub>2</sub>		aq	3			
ZnSe		c	77			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
ZnTe		g	16			
		c	77, 78	78		
Zn(N <sub>3</sub> ) <sub>2</sub>	zinc azide	c	79, 80			
Zn <sub>3</sub> N <sub>2</sub>		c	81			82
Zn(NO <sub>3</sub> ) <sub>2</sub>		c	83			
		aq	18, 83			
Zn(NO <sub>3</sub> ) <sub>2</sub> · nH <sub>2</sub> O		c	18, 83			
Zn(NH <sub>2</sub> ) <sub>2</sub>		c	84			
ZnCl <sub>2</sub> · nNH <sub>3</sub>		c	86, 87, 88			
ZnCl <sub>2</sub> · 2N <sub>2</sub> H <sub>4</sub>		c	85			
ZnCl <sub>2</sub> · 5NH <sub>3</sub> · 2H <sub>2</sub> O		c	41			
2ZnCl <sub>2</sub> · 4NH <sub>3</sub> · 1/2H <sub>2</sub> O		c	41			
3ZnCl <sub>2</sub> · 6NH <sub>4</sub> Cl · H <sub>2</sub> O		c	41			
2ZnCl <sub>2</sub> · 8NH <sub>4</sub> Cl · ZnO		c	41			
3ZnCl <sub>2</sub> · 10NH <sub>4</sub> Cl · ZnO		c	41			
6ZnCl <sub>2</sub> · 12NH <sub>3</sub> · ZnO		c	41			
ZnBr <sub>2</sub> · nNH <sub>3</sub>		c	50, 86, 87			
ZnBr <sub>2</sub> · 2N <sub>2</sub> H <sub>4</sub>		c	85			
ZnI <sub>2</sub> · nNH <sub>3</sub>		c	50, 86			
ZnI <sub>2</sub> · 2N <sub>2</sub> H <sub>4</sub>		c	85			
ZnSO <sub>4</sub> · nNH <sub>3</sub>		c	76, 89, 90			
Zn(NH <sub>4</sub> ) <sub>2</sub> (SO <sub>4</sub> ) <sub>2</sub> · nH <sub>2</sub> O		c	91, 92			
ZnSb		c	93			34
ZnCO <sub>3</sub>		c	94, 95, 97		96	96
Zn(CH <sub>3</sub> ) <sub>2</sub>	dimethylzinc	liq	99			
Zn(C <sub>2</sub> H <sub>5</sub> ) <sub>2</sub>	diethylzinc	liq	98, 99			
ZnC <sub>2</sub> O <sub>4</sub> · 2H <sub>2</sub> O	zinc oxalate dihydrate	c	100			
Zn(CHO <sub>2</sub> ) <sub>2</sub>	zinc formate	c	101, 102			
		aq	101, 102			
Zn(CHO <sub>2</sub> ) <sub>2</sub> · 2H <sub>2</sub> O	zinc formate dihydrate	c	101, 102			
Zn(C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>2</sub>	zinc acetate	c	101, 102			
		aq	18, 103			
Zn(C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>2</sub> · nH <sub>2</sub> O		c	101, 102			
Zn(C <sub>2</sub> H <sub>3</sub> O <sub>3</sub> ) <sub>2</sub>	zinc glycollate	c	104			
		aq	104			
Zn(C <sub>2</sub> H <sub>3</sub> O <sub>3</sub> ) <sub>2</sub> · 2H <sub>2</sub> O	zinc glycollate dihydrate	c	104			
Zn(CN) <sub>2</sub>		c	105, 106			
Zn(CN) <sub>2</sub> · ZnO		c	105, 106			
Zn(CN) <sub>4</sub> <sup>2-</sup>		aq	106			
ZnCl <sub>2</sub> · C <sub>2</sub> H <sub>4</sub> (NH <sub>2</sub> ) <sub>2</sub>	zinc chloride ethylenediamine	c	95			

Table 31 (Continued)  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$\text{ZnCl}_2 \cdot 3\text{C}_2\text{H}_4(\text{NH}_2)_2$	zinc chloride tris-ethylenediamine	c	95			
$\text{ZnBr}_2 \cdot \text{C}_2\text{H}_4(\text{NH}_2)_2$	zinc bromide ethylenediamine	c	95			
$\text{ZnBr}_2 \cdot 3\text{C}_2\text{H}_4(\text{NH}_2)_2$	zinc bromide tris-ethylenediamine	c	95			
$\text{ZnI}_2 \cdot \text{C}_2\text{H}_4(\text{NH}_2)_2$	zinc iodide ethylenediamine	c	95			
$\text{ZnI}_2 \cdot 3\text{C}_2\text{H}_4(\text{NH}_2)_2$	zinc iodide tris-ethylenediamine	c	95			
$\text{Zn}(\text{CHO}_2)_2 \cdot n\text{NH}_3$		c	89, 90			
$\text{ZnSiO}_3$		c	20			
$\text{Zn}_2\text{SiO}_4$		c	20			
		gls	20			
$2\text{ZnI}_2 \cdot \text{PbI}_2$		c	48			

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Table 32  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Cd	in Hg	g	1	2, 3		4
		c		6, 7, 8	5	4, 9
				8		
Cd <sup>n+</sup>		g	10			
Cd <sup>++</sup>		aq	4	11		
Cd <sub>2</sub>		g	12, 13			
CdO		c	14, 15		5	5
CdH		g	12, 13		4, 12	
(CdH) <sup>+</sup>		g	13			
Cd(OH) <sub>2</sub>		c	16, 17	18		
CdF <sub>2</sub>		c	11, 19, 20, 21, 23	11, 19, 21, 23		
		aq	24			
CdCl		g	13		4, 25, 26	
CdCl <sub>2</sub>		c	27, 28, 30, 31	8, 18, 27, 28		
		aq	27, 32	8, 18, 27, 28		
CdCl <sub>2</sub> •nH <sub>2</sub> O		c	30, 33	34		
CdCl <sub>2</sub> •2HCl•7H <sub>2</sub> O		c	35			
CdCl <sub>2</sub> •CdO•H <sub>2</sub> O		c	36			
CdBr		g	13	26		
CdBr <sub>2</sub>		c	39, 40, 41	37, 38, 42		
		aq	32			
CdBr <sub>2</sub> •4H <sub>2</sub> O		c	17, 38	38		
CdBr <sub>2</sub> •CdO•H <sub>2</sub> O		c	36			
CdI		g	13, 43	4, 12, 13		
CdI <sub>2</sub>		c	17, 28, 44	28, 37		
		aq	32			
CdI <sub>2</sub> •CdO•H <sub>2</sub> O		c	36			
CdS		c	45, 47, 48, 49	46		
CdSO <sub>4</sub>		c	17, 50	51		
		aq	17, 52	8, 11		
CdSO <sub>4</sub> •nH <sub>2</sub> O		c	17, 50	51, 54		
CdSO <sub>4</sub> •2 1/2H <sub>2</sub> SO <sub>4</sub>		c	53			
CdS <sub>2</sub> O <sub>6</sub>		aq	4, 17			
CdSe		g	55			
CdTe		c	56	56		
Cd(N <sub>3</sub> ) <sub>2</sub>	cadmium azide	c	57			
Cd <sub>3</sub> N <sub>2</sub>		c	58			
Cd(NO <sub>3</sub> ) <sub>2</sub>		c	59			
		aq	17, 59	4		
Cd(NO <sub>3</sub> ) <sub>2</sub> •nH <sub>2</sub> O		c	59			

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Table 32 (Continued)  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$\text{Cd}(\text{NH}_2)_2$		c	60			
$\text{CdCl}_2 \cdot n\text{NH}_3$		c	36, 61			
$\text{CdCl}_2 \cdot \text{NH}_4\text{Cl} \cdot 1/2\text{H}_2\text{O}$		c	36			
$\text{CdCl}_2 \cdot 4\text{NH}_4\text{Cl}$		c	36			
$\text{CdBr}_2 \cdot n\text{NH}_3$		c	36, 61			
$\text{CdBr}_2 \cdot \text{NH}_4\text{Br} \cdot 1/2\text{H}_2\text{O}$		c	36			
$\text{CdI}_2 \cdot \text{NH}_4\text{I} \cdot 1/2\text{H}_2\text{O}$		c	36			
$\text{CdI}_2 \cdot n\text{NH}_3$		c	36, 61, 62			
$\text{CdSO}_4 \cdot n\text{NH}_3$		c	62			
$\text{CdSb}$		c	63, 64			
$\text{Cd}_3\text{Sb}_2$		c	63	63		
$\text{CdCO}_3$		c	17, 65	4		
		amorp	17			
$\text{Cd}(\text{CH}_3)_2$		liq	66			
$\text{Cd}(\text{C}_2\text{H}_5)_2$		liq	66			
$\text{Cd}(\text{CN})_2$		c	29			
		aq	4, 17, 29			
$\text{Cd}(\text{ONC})_2$	cadmium fulminate	c	57			
$2\text{Cd}(\text{CN})_2 \cdot \text{CdO} \cdot 5\text{H}_2\text{O}$		c	29			
$2\text{CdI}_2 \cdot \text{PbI}_2$		c	22			

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Table 33  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Hg		g	3, 12		3	3
		liq			9	9
Hg <sup>n+</sup>		g	2			
Hg <sup>++</sup>		aq		4, 5		
Hg <sub>2</sub>		g	6			
Hg <sub>2</sub> <sup>++</sup>		aq		4, 7		
HgO		c	10, 11, 12, 13, 14, 34	5, 8, 15		9
Hg <sub>2</sub> O		c	34, 16			
HgH		g	6, 17		1, 6	1, 6
HHgO <sub>2</sub> <sup>-</sup>		aq		5		
Hg(OH) <sub>2</sub>		aq		5		
HgF		g	6			
HgCl		g	6, 17, 18		1, 6, 18	1, 6, 18
HgCl <sub>2</sub>		c	34, 20			
		aq	34, 21, 22, 23, 24, 25, 26			
HgCl <sub>3</sub> <sup>-</sup>		aq	1, 19			
Hg <sub>2</sub> Cl <sub>2</sub>		c	16, 27, 32, 28, 29, 30	27, 28, 29, 30, 33, 69		9
HgCl <sub>2</sub> ·nHgO		c	21, 35			
Hg <sub>8</sub> r		g	6, 17, 18		1, 6, 18	1, 6, 18
Hg <sub>8</sub> r <sub>2</sub>		c	16, 36			
		aq	36			
HgBr <sub>4</sub> <sup>--</sup>		aq	34, 19			
Hg <sub>2</sub> Br <sub>2</sub>		c	37, 38, 39, 16	37, 38		
Hg <sub>8</sub> r <sub>2</sub> ·nHgO		c	21, 35			
HgI		g	6, 17, 18		1, 6, 18	1, 6, 18
HgI <sub>2</sub>		c	16			
HgI <sub>4</sub> <sup>--</sup>		aq	34, 19			
Hg <sub>2</sub> I <sub>2</sub>		c	41, 42, 16	41, 42		40
HgS		g	6			
		c	34, 16, 43, 44, 45	44, 45		
HgSO <sub>4</sub>		c	16, 46			
Hg <sub>2</sub> SO <sub>4</sub>		c	47	47, 48, 50, 69	9	
HgSO <sub>4</sub> ·2HgO		c	16			
HgSe		g	6			
		c	49			
Hg <sub>2</sub> (N <sub>3</sub> ) <sub>2</sub>	mercurous azide	c	51, 52			
Hg(NO <sub>3</sub> ) <sub>2</sub>		aq	34, 16			
Hg(NO <sub>3</sub> ) <sub>2</sub> ·1/2H <sub>2</sub> O		c	16			
Hg <sub>2</sub> (NO <sub>3</sub> ) <sub>2</sub>		aq	34			

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Table 33 (Continued)  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$\text{Hg}_2(\text{NO}_3)_2 \cdot 2\text{H}_2\text{O}$		c	16			
$\text{Hg}(\text{NO}_3)_2 \cdot 2\text{HgO} \cdot \text{H}_2\text{O}$		c	16			
$(\text{Hg}_2\text{N})_2\text{O}$	millon's oxide	c	53			
$\text{Hg}_2\text{NOH}$	millon's hydroxide	c	53, 54			
$\text{Hg}_2\text{NOH} \cdot 1 \frac{1}{2}\text{H}_2\text{O}$		c	53, 54			
$\text{Hg}_2\text{NOH} \cdot 2\text{H}_2\text{O}$	millon's base	c	53, 54			
$\text{Hg}_2\text{NCl} \cdot n\text{H}_2\text{O}$		c	53, 54			
$\text{Hg}_2\text{NCl} \cdot 1/2\text{HgCl}_2$		c	53			
$\text{HgCl}_2 \cdot n\text{NH}_3$		c	55, 56			
$\text{Hg}_2\text{NCl} \cdot n\text{NH}_3$		c	53, 54			
$\text{Hg}_2\text{NCl} \cdot n\text{NH}_4\text{Cl}$		c	53, 54			
$\text{Hg}_2\text{NBr}$	millon's bromide	c	53			
$\text{HgBr}_2 \cdot n\text{NH}_3$		c	55, 56			
$\text{Hg}_2\text{NBr} \cdot n\text{NH}_4\text{Br}$		c	53, 54			
$\text{Hg}_2\text{NBr} \cdot n\text{HgBr}_2$		c	54			
$\text{HgI}_2 \cdot n\text{NH}_3$		c	56			
$\text{Hg}(\text{CH}_3)_2$	dimethylmercury	liq	57, 58			
$\text{Hg}(\text{C}_2\text{H}_5)_2$	diethylmercury	liq	57			
$\text{HgC}_2\text{O}_4$	mercuric oxalate	c	59			
$\text{Hg}(\text{C}_2\text{H}_3\text{O}_2)_2$	mercuric acetate	c	59			
		aq	59			
$\text{Hg}_2(\text{C}_2\text{H}_3\text{O}_2)_2$	mercurous acetate	c	16			
$\text{HgCl}_2 \cdot \text{CH}_3\text{OH}$		c	60			
$\text{Hg}(\text{CN})_2$		c	34, 61, 62			
		aq	34, 16			
$\text{Hg}(\text{CN})_4^{--}$		aq	53, 61			
$\text{Hg}(\text{ONC})_2$	mercuric fulminate	c	51, 63			
$\text{Hg}(\text{CN})_2 \cdot \text{HgO}$		c	64			
$3\text{Hg}(\text{CN})_2 \cdot \text{HgO}$		c	64			
$\text{HgX}_2 \cdot \text{C}_2\text{H}_4(\text{NH}_2)_2$		c	65			
$\text{Hg}(\text{CN})_2 \cdot \text{NH}_4\text{Cl} \cdot 3/4\text{H}_2\text{O}$		c	66			
$\text{Hg}(\text{CN})_2 \cdot \text{NH}_4\text{Br} \cdot 1/2\text{H}_2\text{O}$		c	66			
$\text{Hg}(\text{CN})_2 \cdot \text{NH}_4\text{I} \cdot 1/4\text{H}_2\text{O}$		c	66			
$\text{Hg}(\text{CNS})_2$		c	64			
$\text{Hg}_5\text{I}_2$		c	67			
$\text{Hg}(\text{CN})_2 \cdot \text{ZnCl}_2 \cdot 7\text{H}_2\text{O}$		c	66			
$2\text{Hg}(\text{CN})_2 \cdot \text{ZnBr}_2 \cdot 8\text{H}_2\text{O}$		c	66			
$\text{Hg}(\text{CN})_2 \cdot \text{CdCl}_2 \cdot 2\text{H}_2\text{O}$		c	66, 68			
$\text{Hg}(\text{CN})_2 \cdot \text{CdBr}_2 \cdot 3\text{H}_2\text{O}$		c	66, 68			
$2\text{Hg}(\text{CN})_2 \cdot \text{CdI}_2 \cdot 8\text{H}_2\text{O}$		c	66, 68			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Cu		g	3,4		6	6
		c			1	1
	in Hg			2		
Cu <sup>n+</sup>		g	5			
Cu <sup>+</sup>		aq	8	12, 9, 22		
Cu <sup>++</sup>		aq	6	9, 10, 11, 12		
CuO		g	13			
		c	15, 16, 17, 18, 19, 20	21, 19	14	14, 15
CuO <sub>2</sub> <sup>--</sup>		aq		27		
Cu <sub>2</sub> O		c	15, 23, 26	21, 19, 23, 25, 26	14	14, 15
CuH		g	13		28, 6	28, 6
HCuO <sub>2</sub>		aq		27		
Cu(OH) <sub>2</sub>		c	17, 18, 29, 30, 31			
CuF		g	13			
CuF <sub>2</sub>		c	32, 33			
CuF <sub>2</sub> ·2H <sub>2</sub> O		c	34	34		
CuCl		g	13, 35		28, 36, 6	28, 36, 6
		c	37, 38, 18, 39, 22, 40, 41	37, 38, 9		
		aq	22			
CuCl <sub>2</sub>		c	18, 42, 44, 45, 46			
		aq	39, 17, 31, 18, 43, 47, 48, 49, 50, 51			
	in C <sub>2</sub> H <sub>5</sub> OH		42			
CuCl <sub>2</sub> ·2H <sub>2</sub> O		c	52, 53, 18			
CuCl <sub>2</sub> <sup>-</sup>		aq	22	22		
Cu(ClO <sub>3</sub> ) <sub>2</sub>		aq	18, 6			
Cu(ClO <sub>4</sub> ) <sub>2</sub>		aq	12, 6			
CuCl <sub>2</sub> ·nCuO		c	48, 77			
CuCl <sub>2</sub> ·3CuO·4H <sub>2</sub> O		c	77			
CuBr		g	13		28, 36, 6	28, 36, 6
		c	55, 18, 56, 57	56		
CuBr <sub>2</sub>		c	59, 18, 60			
		aq	58, 18			
CuBr <sub>2</sub> ·4H <sub>2</sub> O		c	59			
CuBr <sub>2</sub> <sup>-</sup>		aq	59	61		
CuBr <sub>2</sub> ·3Cu(OH) <sub>2</sub>		c	62			
CuI		g	13		28, 36, 6	28, 36, 6
		c	61, 56, 18, 65, 41	64, 56	14	
CuI <sub>2</sub>		c	6			63
		aq	6			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
CuS	cuprous azide	c	18, 67, 70, 69	67	68	68
Cu <sub>2</sub> S		c	18, 69, 71, 72, 73, 75, 74	72, 75	68	68
CuSO <sub>4</sub>		c	79, 67, 80, 81, 18, 82	19, 79, 67		63
		aq	10, 11, 18, 76, 31, 77, 78	10, 11		
CuSO <sub>4</sub> ·nH <sub>2</sub> O		c	18, 80, 81, 82, 19, 83, 79, 84, 85	19, 67, 79		63
CuS <sub>2</sub> O <sub>6</sub>		aq	6			
CuS <sub>2</sub> O <sub>6</sub> ·5H <sub>2</sub> O		c	18			
Cu <sub>2</sub> SO <sub>4</sub>		c	86			
		aq	12			
CuSO <sub>4</sub> ·3CuO·4H <sub>2</sub> O		c	62			
CuSe		c	87			
Cu <sub>2</sub> Se		c	87, 88			92
CuSeO <sub>4</sub>		aq	89			
CuSeO <sub>4</sub> ·5H <sub>2</sub> O		c	89			
CuN <sub>3</sub>		c	90			
Cu <sub>3</sub> N		c	91			92
Cu(NO <sub>3</sub> ) <sub>2</sub>		c	94			
		aq	18, 93, 30			
Cu(NO <sub>3</sub> ) <sub>2</sub> ·nH <sub>2</sub> O		c	62, 18			
Cu(NO <sub>3</sub> ) <sub>2</sub> ·3CuO·3H <sub>2</sub> O		c	62			
4CuO·3N <sub>2</sub> O <sub>5</sub>		c	95			
Cu(NH <sub>3</sub> ) <sup>+</sup>		aq		96		
Cu(NH <sub>3</sub> ) <sup>++</sup>		aq		96		
Cu(NH <sub>3</sub> ) <sub>2</sub> <sup>+</sup>		aq		96		
Cu(NH <sub>3</sub> ) <sub>2</sub> <sup>++</sup>		aq	31	96		
Cu(NH <sub>3</sub> ) <sub>3</sub> <sup>++</sup>		aq		96		
Cu(NH <sub>3</sub> ) <sub>4</sub> <sup>++</sup>		aq	31	96		
Cu(NH <sub>3</sub> ) <sub>5</sub> <sup>++</sup>		aq		96		
Cu(NO <sub>3</sub> ) <sub>2</sub> ·nNH <sub>3</sub>		c	97, 98			
CuCl·nNH <sub>3</sub>		c	102			
CuCl <sub>2</sub> ·nNH <sub>3</sub>		c	99, 31, 100			
CuCl <sub>2</sub> ·2NH <sub>3</sub> ·1/4H <sub>2</sub> O		c	31			
CuCl <sub>2</sub> ·2NH <sub>4</sub> Cl		c	31, 101			
CuCl <sub>2</sub> ·2NH <sub>4</sub> Cl·2H <sub>2</sub> O		c	31, 101			
CuCl <sub>2</sub> ·4NH <sub>3</sub> ·2H <sub>2</sub> O		c	31			
CuCl <sub>2</sub> ·5NH <sub>3</sub> ·nH <sub>2</sub> O		c	31			
CuBr·nNH <sub>3</sub>		c	102			
CuBr <sub>2</sub> ·nNH <sub>3</sub>		c	100			
CuI·nNH <sub>3</sub>		c	102			

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Substance			$\Delta H_f^\circ$	$\Delta P_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$\text{CuI}_2 \cdot n\text{NH}_3$		c	100, 6		-	
$\text{CuSO}_4 \cdot n\text{NH}_3$		c	31, 103, 104			
$\text{CuSO}_4 \cdot (\text{NH}_4)_2\text{SO}_4$		c	31			
$\text{CuSO}_4 \cdot (\text{NH}_4)_2\text{SO}_4 \cdot n\text{H}_2\text{O}$		c	31, 105			
$\text{CuSO}_4 \cdot 4\text{NH}_3 \cdot 1 \frac{1}{2}\text{H}_2\text{O}$		c	31			
$\text{CuP}_2$		c	106			
$\text{Cu}_3\text{P}$		c	107			
$\text{CuX} \cdot n\text{PH}_3$		c	108			
$\text{Cu}_2\text{Sb}$		c				92
$\text{Cu}_3\text{Sb}$		c	109			92
$\text{CuCO}_3$		c	112	110, 111	6	
$\text{Cu}(\text{CHO}_2)_2$	cupric formate	aq	113, 114			
$\text{Cu}(\text{CHO}_2)_2 \cdot 4\text{H}_2\text{O}$		c	113, 114			
$\text{Cu}(\text{C}_2\text{H}_3\text{O}_2)_2$	cupric acetate	c	113, 93			
		aq	113, 93, 18			
$\text{Cu}(\text{C}_2\text{H}_3\text{O}_2)_2 \cdot \text{H}_2\text{O}$		c	18, 93, 113			
$\text{Cu}(\text{C}_2\text{H}_3\text{O}_3)_2$	cupric glycollate	c	115			
		aq	115			
$2\text{CuCl} \cdot \text{CO}$		aq	116			
$2\text{CuCl} \cdot \text{CO} \cdot 2\text{H}_2\text{O}$		c	116			
$2\text{CuCl} \cdot \text{C}_2\text{H}_2$		c	117			
$\text{CuCl}_2 \cdot 2\text{CH}_3\text{OH}$		c	118			
$\text{Cu}(\text{C}_2\text{H}_5\text{SO}_4)_2$	cupric ethylsulfate	aq	18			
$\text{CuCN}$		aq	119			
$\text{CuONC}$	cuprous fulminate	c	90			
$\text{Cu}(\text{CHO}_2)_2 \cdot n\text{NH}_3$		c	120			
$\text{Cu}(\text{C}_2\text{H}_3\text{O}_2)_2 \cdot n\text{NH}_3$		c	120			
$\text{Cu}(\text{C}_2\text{H}_3\text{O}_3)_2 \cdot n\text{NH}_3$		c	120			
$\text{CuCl}_2 \cdot n\text{C}_2\text{H}_4(\text{NH}_2)_2$		c	121			
$\text{CuBr}_2 \cdot n\text{C}_2\text{H}_4(\text{NH}_2)_2$		c	121			
$\text{Cu}_3\text{Si}$		c				122
$\text{CuSiO}_3 \cdot \text{H}_2\text{O}$		c				92
$\text{Cu}_3\text{Sn}$		c	51, 66			
$\text{Cu}_2\text{Zn}_3$		c	51, 66			
$\text{Cu}_2\text{Cd}_3$		c	109, 51			

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| 3. Kelley-19                         | 45. Agostini-1                                |
| 4. Washburn et al-1                  | 46. Sabatier-2                                |
| 5. Baur and Brunner-1                | 47. Krustinsons-3                             |
| 6. Calculated                        | 48. Korveze-1                                 |
| 7. Bacher and Goudsmit-1             | 49. Sieverts and Gotta-1                      |
| 8. Fenwick-1                         | 50. Jellinek and Koop-1                       |
| 9. Nielsen and Brown-2               | 51. Biltz, Wagner, Pieper, and Holverscheid-1 |
| 10. Müller and Reuther-3             | 52. van Deventer and van de Stadt-1           |
| 11. Wetmore and Gordon-1             | 53. Reicher and van Deventer-2                |
| 12. Heinerth-1                       | 54. Berthelot-77                              |
| 13. Gaydon-1                         | 55. Klein-2                                   |
| 14. Kelley-24                        | 56. Ishikawa, Yamoaki, and Murooka-1          |
| 15. Wöhler and Jochum-1              | 57. Kapustinskiĭ and Panteev-1                |
| 16. von Wartenberg and Werth-2       | 58. Rolla-1                                   |
| 17. Fricke, Gwinner, and Feichtner-1 | 59. Sabatier-7                                |
| 18. Thomsen-16                       | 60. McMorris and Yost-1                       |
| 19. Randall, Nielsen, and West-1     | 61. Rodlander and Storbeck-1                  |
| 20. Wöhler and Salz-2                | 62. Sabatier-9                                |
| 21. Maier-5                          | 63. Kelley-18                                 |
| 22. Noyes and Chow-1                 | 64. Seidell-1                                 |
| 23. Makolkin-4                       | 65. Joannis-6                                 |
| 24. Maier-1                          | 66. Herschkowitsch-1                          |
| 25. Allmand-1,2                      | 67. Kelley-22                                 |
| 26. Ishikawa and Kimura-1            | 68. Anderson-6                                |
| 27. McDowell and Johnston-1          | 69. von Wartenberg-19                         |
| 28. Herzberg-9                       | 70. Berthelot-13                              |
| 29. Sabatier-9                       | 71. Vanyukov and Kiseleva-1                   |
| 30. de Forcrand-64                   | 72. Kapustinskiĭ and Makolkin-2               |
| 31. Bouzat-1                         | 73. Britzke and Kapustinskiĭ-4                |
| 32. von Wartenberg-9                 | 74. Britzke and Kapustinskiĭ-1                |
| 33. Domange-2                        | 75. Kapustinskiĭ and Makolkin-1               |
| 34. Jahn-Held and Jellinek-1         | 76. Quinten-2                                 |
| 35. Brewer et al-1                   | 77. Lange, Monheim and Robinson-1             |
| 36. Stevenson-2                      | 78. Birnthalder and Lange-1                   |
| 37. Kapustinskiĭ-5                   | 79. Ishikawa and Murooka-1                    |
| 38. Watanabe-2,4                     | 80. Pickering-8                               |
| 39. von Wartenberg and Werth-3       | 81. Schottky-1                                |
| 40. Jellinek and Uloth-1             | 82. Donnan and Hope-1                         |
| 41. Berthelot-157                    | 83. Sano-21                                   |
| 42. Partington and Soper-1           | 84. Scholz-1                                  |



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| 89. Metzner-1                           |  |
| 90. Wöhler and Martin-1,2               |  |
| 91. Juza and Hahn-1                     |  |
| 92. Kelley-25                           |  |
| 93. Berthelot-5                         |  |
| 94. Guntz and Martin-1                  |  |
| 95. Neumann and Sonntag-2               |  |
| 96. Bjerrum-3                           |  |
| 97. Ephraim and Bolle-2                 |  |
| 98. Portillo-2                          |  |
| 99. Biltz, Klatte, and Rahlfs-1         |  |
| 100. Biltz, Brohan, and Wein-1          |  |
| 101. Bouzat and Chauvenet-1             |  |
| 102. Biltz and Stollenwerk-2            |  |
| 103. Ephraim-9,13                       |  |
| 104. Hart and Partington-1              |  |
| 105. Caven and Ferguson-1               |  |
| 106. Haraldsen-1                        |  |
| 107. Weibke and Schrag-1                |  |
| 108. Höltje and Schlegel-1              |  |
| 109. Biltz and Haase-1                  |  |
| 110. Kelley-20                          |  |
| 111. Latimer-1                          |  |
| 112. Berthelot-11                       |  |
| 113. Berthelot-B                        |  |
| 114. Berthelot-9                        |  |
| 115. de Forcrand-32                     |  |
| 116. Hammerl-1                          |  |
| 117. Gilliland, Bliss and Kip-1         |  |
| 118. Lloyd, Brown, Bonnell, and Jones-1 |  |
| 119. Varet-3                            |  |
| 120. Spacu and Voichescu-2              |  |
| 121. Hieber and Feder-1                 |  |
| 122. Schimpff-1                         |  |

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Formula	Description	State				
Ag		g	1		2	2
		c			2	2
$\text{Ag}^{n+}$		g	3, 4, 5			
$\text{Ag}^+$		aq	7, 10	8, 9		6
$\text{Ag}_2\text{O}$		c	7, 13, 14	11, 12		7
$\text{Ag}_2\text{O}_2$		c	15			
AgH		g	12, 17		12	12
AgF		c	18, 19, 34		6	
		aq	6, 20, 21, 22, 23			
$\text{AgF} \cdot n\text{H}_2\text{O}$		c	19, 23, 34	23		
$\text{AgF}_2$		c	24			
$\text{Ag}_2\text{F}$		c	25			
$\text{AgHF}_2$		aq	26			
AgCl		g	1		6, 16, 27, 28, 29	6, 16, 27, 28, 29
		c	31	30, 32, 33	35	35
$\text{Ag}_2\text{Cl}$		c	112			
$\text{AgClO}_2$		c	37	8, 37		38
$\text{AgClO}_3$		c	40			
		aq	6			
$\text{AgClO}_4$		c	41			
		aq	42			
AgBr		c	31, 47	35, 43, 44, 45, 46	38	38
AgI		c	31, 49, 50, 51, 54	48, 52, 53		38
$\text{AgIO}_3$		c			55, 56, 57	57
$\text{Ag}_2\text{H}_3\text{IO}_6$		c	58			
$3\text{AgI} \cdot \text{HI} \cdot 7\text{H}_2\text{O}$		c			59	6, 59
$\text{Ag}_2\text{S}$		c	60, 61, 63, 64, 65, 67, 68	61, 62	66	
$\text{Ag}_2\text{SO}_4$		c	70	8	69	69
		aq	51, 72			
$\text{Ag}_2\text{S}_2\text{O}_6$		aq	6			
$\text{Ag}_2\text{S}_2\text{O}_6 \cdot 2\text{H}_2\text{O}$		c	51			
$\text{Ag}(\text{S}_2\text{O}_3)_2^{---}$		aq	35, 39			
$\text{Ag}_2\text{SO}_4 \cdot 2\text{HCl}$		c	73			
$\text{Ag}_2\text{Se}$		c	74			
$\text{Ag}_2\text{SeO}_4$		c	8, 75			
$\text{AgN}_3$	silver azide	c	76			
$\text{AgNO}_2$		c	77, 78		79	79
		aq	6	6, 80		
$\text{AgNO}_3$		c	81, 82, 83, 84, 85		85	38
		aq	6	8, 46		

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$\text{Ag}_2\text{N}_2\text{O}_2$		c	86			
$\text{Ag}(\text{NH}_3)_2^+$		aq	85			
$\text{AgNO}_3 \cdot n\text{NH}_3$		c	87, 88, 89, 90, 91			
		aq	88, 89			
$\text{AgCl} \cdot n\text{NH}_3$		c	92, 93			
$\text{AgClO}_4 \cdot n\text{NH}_3$		c	89			
		aq	89			
$\text{AgBr} \cdot n\text{NH}_3$		c	92, 93			
$\text{AgI} \cdot n\text{NH}_3$		c	92			
$\text{AgP}_2$		c	94			
$\text{AgP}_3$		c	94			
$\text{Ag}_2\text{C}_2$	silver acetylide	c	88			
$\text{Ag}_2\text{CO}_3$		c	51, 95, 96, 98	97	99	99
$\text{Ag}_2\text{C}_2\text{O}_4$	silver oxalate	c	100			
$\text{AgC}_2\text{H}_3\text{O}_2$	silver acetate	c	101			
		aq	102, 103			
$\text{Ag}_2\text{C}_2 \cdot \text{AgCl}$		c	88			
$(\text{Ag}_2\text{C}_2)_2 \cdot \text{AgCl}$		c	88			
$\text{Ag}_2\text{C}_2 \cdot \text{AgI}$		c	88			
$\text{Ag}_2\text{C}_2 \cdot 2\text{AgI}$		c	88			
$\text{Ag}_2\text{C}_2 \cdot \text{Ag}_2\text{SO}_4$		c	88			
$(\text{Ag}_2\text{C}_2)_2 \cdot \text{Ag}_2\text{SO}_4$		c	88			
$\text{AgCN}$		c	51	8, 104		
$\text{Ag}(\text{CN})_2^-$		aq	51	104		
$\text{Ag}_2\text{CN}_2$	silver cyanamide	c	90			
$\text{AgCNO}$	silver cyanate	c	8, 90			
$\text{AgONC}$	silver fulminate	c	106			
$\text{Ag}_2\text{C}_2 \cdot \text{AgNO}_3$		c	88			
$\text{AgCl} \cdot \text{CH}_3\text{NH}_2$		c	107			
$\text{AgBr} \cdot \text{CH}_3\text{NH}_2$		c	107			
$\text{AgI} \cdot n\text{CH}_3\text{NH}_2$		c	107			
$\text{AgSCN}$		c	6			
		aq	8, 108			
$\text{Ag}_2\text{Zn}_3$		c	109			
$\text{Ag}_3\text{Hg}_4$		c	110			
$\text{Ag}_2\text{HgI}_4$		c	105			

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| 3. Shenstone-7                       | 45. Woitinek-1                             |
| 4. Gilbert-1                         | 46. Harned and Owen-1                      |
| 5. Rasmussen-6                       | 47. Webb-1,2                               |
| 6. Calculated                        | 48. Owen-4                                 |
| 7. Pitzer and Smith-1                | 49. Klein-2                                |
| 8. Seidell-1                         | 50. Braune and Koref-1                     |
| 9. Owen and Brinkley Jr.-1           | 51. Thomsen-16                             |
| 10. Latimer, Schutz, and Hicks Jr.-1 | 52. Gerke-1                                |
| 11. Makolkin-4                       | 53. Taylor-2                               |
| 12. Fried-1                          | 54. Gerth-2                                |
| 13. Favre and Silbermann-2           | 55. Pearce and Wirth-1                     |
| 14. Mixer-12                         | 56. Li and Lo-1                            |
| 15. Jirsa-3                          | 57. Greensfelder and Latimer-2             |
| 16. Herzberg-9                       | 58. Berthelot-69                           |
| 17. Farkas-2                         | 59. Stephenson and Adams-2                 |
| 18. Guntz-1                          | 60. Zeumer and Roth-4                      |
| 19. Guntz and Guntz-1                | 61. Kapustinskiĭ and Makolkin-1            |
| 20. Schwiete and Pranschke-1         | 62. Kimura-1,5                             |
| 21. Pranschke and Schwiete-1         | 63. Sano-23                                |
| 22. Petersen-4                       | 64. Watanabe-1                             |
| 23. Jahn-Held and Jellinek-2         | 65. Voevodskiĭ and Gol'bert-1              |
| 24. von Wartenberg-9                 | 66. Kelley-22                              |
| 25. Guntz-10                         | 67. Keyes and Felsing-1                    |
| 26. Guntz-4                          | 68. Kapustinskiĭ and Korshunov-6           |
| 27. Jenkins and Rochester-1          | 69. Latimer, Hicks Jr., and Schutz-1       |
| 28. Brice-1                          | 70. Ishikawa-5                             |
| 29. Stevenson-2                      | 71. Ishikawa and Hagisawa-1                |
| 30. Harned and Ehlers-1              | 72. Berthelot-152                          |
| 31. Bertram and Roth-1               | 73. Ephraim-8                              |
| 32. Salstrom-5                       | 74. Fabre-2                                |
| 33. Grube and Rau-1                  | 75. Metzner-1,2                            |
| 34. Guntz and Guntz-2                | 76. Wöhler and Martin-1                    |
| 35. Eastman and Milner-1             | 77. Randall, Manov, and Brown-1            |
| 36. Guntz-11                         | 78. Centnerszwer and Chesinski-1           |
| 37. Smith, Pitzer, and Latimer-1     | 79. Brown, Smith, and Latimer-3            |
| 38. Kelley-24                        | 80. Abegg and Pick-1                       |
| 39. Sichowsky and Rossini-1          | 81. Roth-25                                |
| 40. Foote and Saxton-1               | 82. Lange and Martin-3                     |
| 41. Bruni and Levi-2                 | 83. Jackson, Smith, Gatty, and Wolfenden-1 |
| 42. Jirsa-4                          | 84. Centnerszwer and Blumenthal-5          |

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90. Lemoult-1
91. Joannis and Croizier-1
92. Biltz and Stollenwerk-1
93. Isambert-5
94. Haraldsen and Biltz-1
95. Centnerszwer and Krustinson-1
96. Watanabe-12
97. Walker, Bray, and Johnston-1
98. Berthelot-11
99. Anderson-7
100. Berthelot-9
101. van Laar-9
102. Goldschmidt and Maarseveen-1
103. Berthelot-8
104. Randall and Halford-1
105. Ketalaar-1
106. Wöhler and Martin-2
107. Jarry-1
108. Kirschner-1
109. Samson and Himmelstjerna-1
110. Berthelot-125

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Formula	Description	State				
Au		g	1, 2, 3, 4, 5, 6		1	1
		c			7	6
Au <sup>+</sup>		g	8			
AuO <sub>3</sub> <sup>---</sup>		aq		9		
Au <sub>2</sub> O <sub>3</sub>		c	10	11, 12		
HAuO <sub>3</sub> <sup>--</sup>		aq		9		
H <sub>2</sub> AuO <sub>3</sub> <sup>-</sup>		aq		9		
Au(OH) <sub>3</sub>		c	13	9		
		aq		9		
AuCl		c	14, 15, 16			
AuCl <sub>3</sub>		c	14			
		aq	14			
AuCl <sub>3</sub> ·2H <sub>2</sub> O		c	14			
AuCl <sub>4</sub> <sup>-</sup>		aq		11, 17		
HAuCl <sub>4</sub>		aq	13, 14	13, 14		
HAuCl <sub>4</sub> ·nH <sub>2</sub> O		c	13			
AuBr		c	13, 14, 18			
AuBr <sub>2</sub> <sup>-</sup>		aq		11		
AuBr <sub>3</sub>		c	14, 18			
		aq	13			
AuBr <sub>4</sub> <sup>-</sup>		aq	13	11		
HAuBr <sub>4</sub>		aq	13			
HAuBr <sub>4</sub> ·5H <sub>2</sub> O		c	13			
AuI		c	13, 14, 18			
AuCl·nNH <sub>3</sub>		c	16, 19			
AuBr·nNH <sub>3</sub>		c	19			
AuI·nNH <sub>3</sub>		c	19			
Au <sub>2</sub> P <sub>3</sub>		c	20			
AuI·PH <sub>3</sub>		c	21			
AuSb <sub>2</sub>		c	22			
Au(CN) <sub>2</sub> <sup>-</sup>		aq	19	11		
AuSn		c	23			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 36 (Continued)  
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5. Ruff and Konschak-1
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7. Kelley-24
8. Bacher and Goudsmit-1
9. Johnston and Leland-1
10. Mixter-12
11. Latimer-1
12. Gerke and Rourke-1
13. Thomsen-16
14. Fischer and Biltz-1
15. Maignon-12
16. Ephraim-6
17. Bjerrum and Kirschner-1
18. Meyer-2
19. Biltz-10,11
20. Haraldsen and Biltz-1
21. Holtje and Schlegel-1
22. Weibke and Schrag-1
23. Biltz, Rohlf, and Vogel-1

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Table 37  
SPECIFIC REFERENCES

Substance			$\Delta H_f^0$	$\Delta F_f^0$	$S^0$	$C_p^0$
Formula	Description	State				
Pt		g	4	6	2,3	2,3
		c			1	1
Pt <sup>+</sup>		g	2			
Pt(OH) <sub>2</sub>		c	5			
PtCl		c	7			
PtCl <sub>2</sub>		c	7			
PtCl <sub>3</sub>		c	7			
PtCl <sub>4</sub>		c	7,8			
		aq	7,8			
PtCl <sub>4</sub> ·5H <sub>2</sub> O		c	8			
PtCl <sub>4</sub> <sup>--</sup>		aq	5,9	6		
PtCl <sub>6</sub> <sup>--</sup>		aq	3	10	3	
HPtCl <sub>5</sub> ·2H <sub>2</sub> O		c	8			
H <sub>2</sub> PtCl <sub>6</sub>		aq	5,8,11			
H <sub>2</sub> PtCl <sub>6</sub> ·6H <sub>2</sub> O		c	8			
		c	8			
PtBr <sub>4</sub>		c	8			
		aq	8			
PtBr <sub>4</sub> <sup>--</sup>		aq	5			
PtBr <sub>6</sub> <sup>--</sup>		aq	5,8			
H <sub>2</sub> PtBr <sub>6</sub>		aq	8			
H <sub>2</sub> PtBr <sub>6</sub> ·9H <sub>2</sub> O		c	8			
PtI <sub>4</sub>		c	8			
PtI <sub>6</sub> <sup>--</sup>		aq	8			
PtS		c	12,13			
PtS <sub>2</sub>		c	12,13			
(NH <sub>4</sub> ) <sub>2</sub> PtCl <sub>4</sub>		c	5			
		aq	5			
Pt(OH) <sub>2</sub> ·4NH <sub>3</sub>		c	5			
PtCl <sub>2</sub> ·2NH <sub>3</sub>		c	14,15			
PtCl <sub>2</sub> ·4NH <sub>3</sub>		c	15			
		aq	3			
PtCl <sub>2</sub> ·4NH <sub>3</sub> ·H <sub>2</sub> O		c	5			
PtCl <sub>2</sub> ·5NH <sub>3</sub>		c	15			
PtSO <sub>4</sub> ·4NH <sub>3</sub>		c	5			
PtSb <sub>2</sub>		c	16			
PtSn		c	17			
Ag <sub>2</sub> PtCl <sub>6</sub>		c	8			
Ag <sub>2</sub> PtBr <sub>6</sub>		c	8			



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Table 37 (Continued)  
SPECIFIC REFERENCES

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2. Bacher and Goudsmit-1
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4. Jones, Langmuir, and Mackey-1
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6. Latimer-1
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8. Pigeon-1
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10. Grünberg, Lawrentiew, and Ptizyn-1
11. Gire-1,2
12. Biltz and Juza-1
13. Kelley-22
14. Isambert-1
15. Ephraim and Millman-1
16. Poppema and Jaeger-1
17. Jaeger and Bottema-1

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Table 38  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Ir		g		3	1,2,3,4	3
		c			5	6
Ir <sup>+</sup>		g	4			
IrO <sub>2</sub>		c	7,8,9			7,10
IrF <sub>6</sub>		liq	11			
IrCl		c	12			
IrCl <sub>2</sub>		c	12			
IrCl <sub>3</sub>		c	12,13			
IrCl <sub>6</sub> <sup>--</sup>		aq	1			
IrCl <sub>6</sub> <sup>---</sup>		aq	1			
IrS <sub>2</sub>		c	14			
Ir <sub>2</sub> S <sub>3</sub>		c	14			

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9. Biltz-2
10. Jaeger-1
11. Ruff and Fischer-1
12. Wohler and Streicher-1
13. Remy-1
14. Biltz, Laar, Ehrlich, and Meisel-1

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Table 39  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Os		g	2		2	1
		c			3	3
Os <sup>+</sup>		g	4			
OsO <sub>4</sub>		g	6		5	
		c	6,7	6,7		
		aq		5		
HOsO <sub>5</sub>		aq		8		
H <sub>2</sub> OsO <sub>5</sub>		aq		8		
OsS <sub>2</sub>		c	9			
OsP <sub>2</sub>		c	10			
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Table 40  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Re		g	2		1	
		c				
ReO <sub>3</sub>		c			3	4
ReO <sub>4</sub> <sup>-</sup>		aq	5			
Re <sub>2</sub> O <sub>7</sub>		c	5			
Re <sub>2</sub> O <sub>8</sub>		c	6			
HReO <sub>4</sub>		aq	7			
ReF <sub>6</sub>		g	8			
ReS <sub>2</sub>		c	9			
ReAs <sub>2</sub>		c	10			
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1. Kelley-24 2. Brewer et al-1 3. Calculated 4. Kelley-18 5. Roth and Becker-1,4 6. Ogawa-4 7. Roth and Becker-4 8. Ruff and Kwasnik-2 9. Juza and Biltz-2 10. Wiechmann, Hamburg, and Biltz-1						

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Table 41  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Pd		g	3		3	15
		c			1	2
Pd <sup>n+</sup>		g	4			
PdO		c	5, 6, 7			5
Pd <sub>2</sub> H		c	8, 9, 10			
Pd(OH) <sub>2</sub>		c	11, 12			
Pd(OH) <sub>4</sub>		c	11			
PdCl <sub>2</sub>		c	11, 12, 13			
PdCl <sub>4</sub> <sup>--</sup>		aq		14		
PdCl <sub>6</sub> <sup>--</sup>		aq		14		
H <sub>2</sub> PdCl <sub>4</sub>		aq	11			
H <sub>2</sub> PdCl <sub>6</sub>		aq	15			
PdBr <sub>2</sub>		c	12			
PdBr <sub>4</sub> <sup>--</sup>		aq	15			
PdI <sub>2</sub> ·H <sub>2</sub> O		c	11, 12			
PdCl <sub>2</sub> ·nNH <sub>3</sub>		c	16			
PdSb		c				2
PdSb <sub>2</sub>		c				2
PdSb <sub>3</sub>		c				2
Pd(CN) <sub>2</sub>		c	12			
PdCu		c				2
PdCu <sub>3</sub>		c				2
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1. Kelley-24			15. Calculated			
2. Jaeger and Poppema-1			16. Isambert-6			
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12. Joannis-5						
13. Puche-1						
14. Templeton, Watt, and Gaines-1						

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Table 42  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Rh		g	1		2	3
		c			2	4
Rh <sup>n+</sup>		g	5			
RhO		c	6,7			6
Rh <sub>2</sub> O		c	6,7			6
Rh <sub>2</sub> O <sub>3</sub>		c	6,7			6
RhCl		c	7			
RhCl <sub>2</sub>		c	7			
RhCl <sub>3</sub>		c	7			
RhCl <sub>6</sub> <sup>----</sup>		aq	3			
REFERENCES						
1. Brewer et al-1 2. Kelley-24 3. Calculated 4. Jaeger-2 5. Bacher and Goudsmit-1 6. Wohler and Jocheim-1 7. Wohler and Muller-2						

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Table 43  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Ru		g	1		2, 4	3
		c			2	
RuO <sub>2</sub>		c	5			
RuCl <sub>3</sub>		c	5			
RuS <sub>2</sub>		c	6			

REFERENCES

1. Brewer et al-1
2. Kelley-24
3. Kelley-18
4. Calculated
5. Remy-1
6. Kelley-22

Table 44

Tc		c			1	
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REFERENCES

1. Calculated

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Table 45  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Ni		g	1, 2		1	1
		c			3, 4, 5, 6, 7	3, 4, 5, 6, 7
Ni <sup>0+</sup>		g	8			
Ni <sup>++</sup>		aq	7		7	
NiO		g	9		9	
		c	10, 11, 12, 13, 14, 15, 16, 17, 18, 19, 20		21	21
NiH		g	22			
		c	23			
NiH <sub>2</sub>		c	23			
Ni(OH) <sub>2</sub>		c	26, 27	24, 25		
Ni(OH) <sub>3</sub>		c	26			
NiF <sub>2</sub>		c	28, 29			
		aq	30, 31, 32, 33			
NiF <sub>2</sub> ·4H <sub>2</sub> O		c		32		
NiCl		g	22			
NiCl <sub>2</sub>		c	26, 34, 35		24	24
		aq	26, 37			
NiCl <sub>2</sub> ·nH <sub>2</sub> O		c	26, 32, 38, 39			
NiBr <sub>2</sub>		c	40, 41			
		aq	42			
NiBr <sub>2</sub> ·3H <sub>2</sub> O		c	7, 40, 41			
NiI <sub>2</sub>		c	42			
		aq	7			
Ni(IO <sub>3</sub> ) <sub>2</sub>		c	43			
		aq	43			
Ni(IO <sub>3</sub> ) <sub>2</sub> ·2H <sub>2</sub> O		c	43			
Ni(IO <sub>3</sub> ) <sub>2</sub> ·4H <sub>2</sub> O		c	43			
NiS		c	26, 44, 45			
Ni <sub>3</sub> S <sub>2</sub>		c	46			
NiSO <sub>4</sub>		c	47, 48		47, 48	49
		aq	26, 37		50	
NiSO <sub>4</sub> ·6H <sub>2</sub> O		c	48, 51, 52		48	49
NiSO <sub>4</sub> ·7H <sub>2</sub> O		c	26, 33	33, 53		
Ni <sub>2</sub> S <sub>2</sub> O <sub>6</sub>		aq	7			
Ni <sub>2</sub> S <sub>2</sub> O <sub>6</sub> ·6H <sub>2</sub> O		c	28			
NiSe		c	40			
NiTe		c	54			
Ni(N <sub>3</sub> ) <sub>2</sub> ·H <sub>2</sub> O	nickel azide hydrate	c	55			
Ni(NO <sub>3</sub> ) <sub>2</sub>		c	56			
		aq	7			



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Table 45 (Continued)  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$G_p^\circ$
Formula	Description	State				
$\text{Ni}(\text{NO}_3)_2 \cdot 6\text{H}_2\text{O}$		c	7,28			
$\text{Ni}(\text{NH}_3)_4^{++}$		aq		45		
$\text{Ni}(\text{NH}_3)_6^{++}$		aq		45		
$\text{Ni}(\text{NO}_3)_2 \cdot 6\text{NH}_3$		c	57			
$\text{NiCl}_2 \cdot n\text{NH}_3$		c	58,59,60,61,62			
$\text{NiBr}_2 \cdot n\text{NH}_3$		c	58,60,61,62			
$\text{NiI}_2 \cdot n\text{NH}_3$		c	58,60,61,62			
$\text{NiSO}_4 \cdot n\text{NH}_3$		c	62			
$\text{Ni}_2\text{P}$		c	63			
$\text{Ni}_3\text{P}$		c	63			
$\text{Ni}_5\text{P}_2$		c	63			
$\text{NiSb}$		c	64			
$\text{Ni}_5\text{Sb}_2$		c	64			
$\text{Ni}_3\text{C}$		c	65			
$\text{NiCO}_3$		c		66,67		
$\text{Ni}(\text{CO})_4$		g			36	
		liq			36	
$\text{NiBr}_2 \cdot 6\text{CH}_3\text{OH}$		c			68	
$\text{Ni}(\text{CN})_2$		c	69			
$\text{Ni}(\text{CN})_4^{--}$		aq	69			
$\text{NiSi}$		c	70			67
$\text{Ni}_2\text{Si}$		c	70			67
$\text{NiSn}$		c	71			
$\text{Ni}_3\text{Sn}$		c				67
$\text{Ni}_3\text{Sn}_2$		c	71			
$2\text{NiI}_2 \cdot \text{PbI}_2$		c	72			
$2\text{NiI}_2 \cdot \text{PbI}_2 \cdot 3\text{H}_2\text{O}$		c	72			

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| 7. Calculated                    | 17. Roth and Muller-6            |
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| 9. Johnston and Marshall-1       | 19. Ruff and Gersten-4           |
| 10. Kapustinskiĭ and Silberman-1 | 20. Mixer-11                     |

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 Table 45 (Continued)  
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| 22. Gaydon-1                      | 64. Oelsen-1                           |
| 23. Ray and Sahai-1               | 65. Roth-3                             |
| 24. Näsänen-1                     | 66. Dobychn-1                          |
| 25. Britton-2                     | 67. Kelley-20                          |
| 26. Thomsen-16                    | 68. Lloyd, Brown, Bonnell, and Jones-1 |
| 27. Giordoni and Mattias-1        | 69. Varet-4,5                          |
| 28. Domange                       | 70. Oelsen and Samson-Himmelstjerna-1  |
| 29. Jellinek and Rudak-1          | 71. Oelsen-1                           |
| 30. Mulert-1                      | 72. Mosnier-1                          |
| 31. Petersen-1                    |  |
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| 35. Sano-3                        |  |
| 36. Kelley-24                     |  |
| 37. Plake-1                       |  |
| 38. Derby and Yngve-1             |  |
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| 54. Fabre-3                       |  |
| 55. Wöhler and Martin-1           |  |
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| 57. Long and Toettcher-1          |  |
| 58. Biltz and Hüttig-2            |  |
| 59. Hart and Partington-1         |  |
| 60. Ephraim-1,12                  |  |
| 61. Biltz and Felkenheuer-3       |  |
| 62. Ephraim-2,8                   |  |

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Table 46  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Co		g	1		1, 2	1
		c			4	3
Co <sup>n+</sup>		g	1			
Co <sup>++</sup>		aq	5		6	
Co <sup>+++</sup>		aq		5, 7		
CoO		c	8, 12, 13	9, 10, 11, 13		
Co <sub>3</sub> O <sub>4</sub>		c	12, 14, 15			
CoH		c	16			
CoH <sub>2</sub>		c	16			
Co(OH) <sub>2</sub>		c	17			
Co(OH) <sub>3</sub>		c	17			
CoF <sub>2</sub>		c	18, 19, 20			
		aq	21			
CoF <sub>2</sub> ·4H <sub>2</sub> O		c	22			
CoF <sub>3</sub>		c	20, 23, 24			
CoCl <sub>2</sub>		c	17, 25, 26, 28		24	24
		aq	17, 29			
CoCl <sub>2</sub> ·nH <sub>2</sub> O		c	17, 25, 26, 28			
CoBr <sub>2</sub>		c	30, 31, 32			
		aq	5			
CoBr <sub>2</sub> ·6H <sub>2</sub> O		c	30			
CoI <sub>2</sub>		c	32, 33, 34			
		aq	35			
Co(IO <sub>3</sub> ) <sub>2</sub>		c	36			
		aq	36			
Co(IO <sub>3</sub> ) <sub>2</sub> ·nH <sub>2</sub> O		c	36			
CoS		c	6, 17, 37, 38			3
Co <sub>2</sub> S <sub>3</sub>		c	5			
CoSO <sub>4</sub>		c	39, 40		39, 40	
		aq	17			
CoSO <sub>4</sub> ·nH <sub>2</sub> O		c	17, 28, 41			40
CoSe		c	31			
CoTe		c	31			
Co(NO <sub>3</sub> ) <sub>2</sub>		c	5, 42			
		aq	17			
Co(NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O		c	17			
Co(NH <sub>3</sub> ) <sub>6</sub> <sup>++</sup>		aq		43		
Co(NH <sub>3</sub> ) <sub>6</sub> <sup>+++</sup>		aq		43		
[Co(NH <sub>3</sub> ) <sub>5</sub> H <sub>2</sub> O] <sup>+++</sup>		aq	5	43		
[Co(NH <sub>3</sub> ) <sub>5</sub> NO <sub>3</sub> ] <sup>++</sup>		aq	44			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$[\text{Co}(\text{NH}_3)_5\text{NO}_3](\text{NO}_3)_2$		c	44			
		aq	44			
$[\text{Co}(\text{NH}_3)_5\text{H}_2\text{O}](\text{NO}_3)_3$		c	44			
		aq	44			
$[\text{Co}(\text{NH}_3)_5\text{H}_2\text{O}]\text{F}_3$		c	45			
$\text{CoCl}_2 \cdot n\text{NH}_3$		c	46, 47, 48			
$\text{CoCl}_2 \cdot 2\text{N}_2\text{H}_4$		c	49, 50			
$[\text{Co}(\text{NH}_3)_4\text{Cl}_2]^+$		aq	51			
$[\text{Co}(\text{NH}_3)_4\text{Cl}_2]\text{Cl}$		c	51			
		aq	51			
$[\text{Co}(\text{NH}_3)_5\text{Cl}]^{++}$		aq	52	52		
$[\text{Co}(\text{NH}_3)_5\text{Cl}]\text{Cl}_2$		c	44			
		aq	44			
$\text{CoCl}_3 \cdot 6\text{NH}_3$		c	53			
$[\text{Co}(\text{NH}_3)_5\text{H}_2\text{O}]\text{Cl}_3$		c	44			
		aq	44			
$\text{CoBr}_2 \cdot n\text{NH}_3$		c	46, 47, 48			
$\text{CoBr}_2 \cdot 2\text{N}_2\text{H}_4$		c	49			
$[\text{Co}(\text{NH}_3)_5\text{Br}]^{++}$		aq	44			
$[\text{Co}(\text{NH}_3)_5\text{Br}]\text{Br}_2$		c	44			
		aq	44			
$[\text{Co}(\text{NH}_3)_5\text{H}_2\text{O}]\text{Br}_3$		c	44			
		aq	44			
$\text{CoI}_2 \cdot n\text{NH}_3$		c	46, 48			53
$\text{CoI}_3 \cdot 6\text{NH}_3$		c				53
$\text{CoSO}_4 \cdot n\text{NH}_3$		c	47			
$[\text{Co}(\text{NH}_3)_5\text{SO}_4]^+$		aq		52		
$\text{CoP}$		c	54			
$\text{CoP}_3$		c	54			
$\text{Co}_2\text{P}$		c	55			
$\text{CoAs}_2 \cdot \text{CoS}_2$		c				3
$\text{CoSb}$		c	56			3
$\text{CoSb}_2$		c	56			
$\text{Co}_3\text{C}$		c	57		57	
$\text{CoCO}_3$		c	58			
$\text{CoCl}_2 \cdot n\text{C}_2\text{H}_5\text{OH}$		c	50, 59			
$\text{CoCl}_2 \cdot 3\text{C}_2\text{H}_4(\text{OH})_2$		c	50			
$\text{CoBr}_2 \cdot 2\text{CH}_3\text{OH}$		c	50			
$\text{CoBr}_2 \cdot n\text{C}_2\text{H}_5\text{OH}$		c	50, 59			
$\text{CoBr}_2 \cdot n\text{C}_2\text{H}_4(\text{OH})_2$		c	50			
$\text{CoCl}_2 \cdot \text{C}_2\text{H}_4(\text{NH}_2)_2$		c	49, 50			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$[\text{Co}[\text{C}_2\text{H}_4(\text{NH}_2)_2]_2\text{Cl}_2]^+$	cobalt ethylsulfate	aq	51			
$[\text{Co}[\text{C}_2\text{H}_4(\text{NH}_2)_2]_2\text{Cl}_2]\text{Cl}$		c	51			
		aq	51			
$[\text{Co}[\text{C}_2\text{H}_4(\text{NH}_2)_2]_2\text{Cl}_2]$ $\text{Cl} \cdot n\text{NH}_3$		c	60			
$\text{CoCl}_2 \cdot 3\text{C}_2\text{H}_4(\text{NH}_2)_2$		c	49, 50			
$\text{CoBr}_2 \cdot n\text{C}_2\text{H}_4(\text{NH}_2)_2$		c	49, 50			
$\text{CoI}_2 \cdot 3\text{C}_2\text{H}_4(\text{NH}_2)_2$		c	49, 50			
$\text{Co}(\text{C}_2\text{H}_5\text{SO}_4)_2$		aq	17			
$\text{CoSi}$		c	61			
$\text{CoSi}_2$		c	61			
$\text{CoSi}_3$		c	61			
$\text{Co}_2\text{Si}$		c	61			
$\text{Co}_2\text{Sn}$		c	62, 63			
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B. Roth and Havekoss-1			28. Bell-1			
9. Kapustinskiĭ and Hofmann-1			29. Biltz-9			
10. Watanabe-10			30. Crut-1			
11. Emmett and Schultz-2			31. Fabre-1			
12. Mixter-11			32. Jellinek and Uloth-2			
13. Shibata and Mori-1			33. Mosnier-1			
14. Biltz-15			34. Devoto and Guzzi-1			
15. Watanabe-5			35. Pigeon-1			
16. Ray and Sahai-1			36. Meusser-1			
17. Thomsen-16			37. Vanyukov and Kiseleva-1			
18. Domange-1			38. Jellinek and Zakowski-1			
19. Jellinek and Rudat-1			39. Marchal-B			
20. Jellinek and Koop-1						

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47. Ephraim-9
48. Biltz and Huttig-2
49. Hieber and Mühlbauer-1
50. Hieber and Woerner-2
51. Ovenston-1
52. Adell-1
53. Ziegler-1
54. Biltz and Heimbrecht-1
55. Weibke and Schrag-1
56. Oelson-1
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Formula	Description	State				
Fe		g	3	1	1, 9	1
		c			4, 41	9
Fe <sup>n+</sup>		g	5, 6, 28			
Fe <sup>++</sup>		aq	1	7		
Fe <sup>+++</sup>		aq	1, 8, 12, 13	8, 11, 10, 15, 14		
Fe <sub>0.95</sub> O	"FeO" wustite	c	16, 17, 20, 21	18, 19, 21		
Fe <sub>2</sub> O <sub>3</sub>	hematite	c	16, 23, 24, 25	22	9	9
Fe <sub>3</sub> O <sub>4</sub>	magnetite	c	23, 20, 9, 7, 27, 2, 16, 25	26, 7, 27		
Fe(OH) <sup>++</sup>		aq	29	10, 30		
Fe(OH) <sub>2</sub>		c	34, 31	7		
Fe(OH) <sub>2</sub> <sup>+</sup>		aq		30		
Fe(OH) <sub>3</sub>		c	32, 34			
FeF <sub>2</sub>		aq	33			
FeF <sub>3</sub>		aq	33			
FeCl <sup>++</sup>		aq	29	35		
FeCl <sub>2</sub>		c	34, 38, 39, 40	40	36, 37	36, 37
		aq	1, 42			
FeCl <sub>2</sub> ·nH <sub>2</sub> O		c	42, 43, 44			
FeCl <sub>3</sub>		c	44, 45, 46			
		aq	34, 16			
FeCl <sub>3</sub> ·6H <sub>2</sub> O		c	43, 44			
FeBr <sup>++</sup>		aq	29	29		
FeBr <sub>2</sub>		c	47			
		aq	48			
FeBr <sub>3</sub>		aq	47			
FeI <sub>2</sub>		c	49, 47			
		aq	48, 50			
FeS		c	50, 52, 51, 53, 54, 55, 56, 57, 58, 59, 60, 7		61	61
FeS <sub>2</sub>		c	7, 62, 63, 64		65	65
FeSO <sub>4</sub>		c	66			
		aq	34, 32			
FeSO <sub>4</sub> ·nH <sub>2</sub> O		c	66, 67			
Fe <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		aq	34, 32, 68			
Fe(HSO <sub>4</sub> ) <sub>3</sub>		aq	32			
FeSe		c	69			
		amorp	69			
FeTe		c	70			
Fe <sub>2</sub> N		c	80, 71		80	72

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$\text{Fe}_4\text{N}$		c	80, 73, 74		80	80, 72
$\text{Fe}(\text{NO}_3)_3$		aq	34			
$\text{Fe}(\text{NO}_3)_3 \cdot 9\text{H}_2\text{O}$		c	75			
$\text{Fe}(\text{NO})\text{Cl}_2$		aq	76, 77			
$\text{FeCl}_2 \cdot n\text{NH}_3$		c	78, 79, 81, 82			
$\text{FeCl}_3 \cdot 6\text{NH}_3$		c	83			
$\text{FeBr}_2 \cdot n\text{NH}_3$		c	79, 78, 84, 83			
$\text{FeI}_2 \cdot n\text{NH}_3$		c	79, 78, 84			
$\text{Fe}(\text{NO})\text{SO}_4$		aq	76, 77			
$\text{FeSO}_4 \cdot n\text{NH}_3$		c	85			
$\text{FeP}$		c	86, 87			
$\text{FeP}_2$		c	90			
$\text{Fe}_2\text{P}$		c	89, 90			
$\text{Fe}_3\text{P}$		c	90			
$\text{FePO}_4$		c	89			
$\text{FePO}_4 \cdot n\text{H}_2\text{O}$		c	91			
$\text{FeSb}$		c	92			
$\text{FeSb}_2$		c	92			
$\text{Fe}_3\text{C}$	cementite	c	23, 50, 8, 94, 95, 96, 97		93	9
$\text{FeCO}_3$	siderite	c	23, 99, 100, 101		98, 100	98
$\text{Fe}(\text{CO})_5$		liq	23, 102, 103			
$\text{Fe}_2(\text{C}_2\text{O}_4)_3$	ferric oxalate	aq	104			
$\text{Fe}(\text{HC}_2\text{O}_4)_3$	ferric bioxalate	aq	104			
$\text{Fe}(\text{C}_2\text{H}_3\text{O}_2)_3$	ferric acetate	aq	34, 32			
$\text{Fe}(\text{CO})_4\text{Br}_2$		c	47			
		aq	48			
$\text{Fe}(\text{CO})_4\text{I}_2$		c	47			
$\text{Fe}(\text{CN})_6^{---}$		aq	106			
$\text{Fe}_4[\text{Fe}(\text{CN})_6]_3$		c	105			
$\text{FeCO}(\text{CN})_5^{---}$		aq	106			
$\text{Fe}_2\text{CO}(\text{CN})_5$		c	109			
$\text{HFe}(\text{CN})_6^{---}$		aq	106			
$\text{H}_2\text{Fe}(\text{CN})_6^{--}$		aq	106			
$\text{H}_3\text{Fe}(\text{CN})_6$		aq	1			
$\text{H}_3\text{Fe}(\text{CN})_6^-$		aq	106			
$\text{H}_4\text{Fe}(\text{CN})_6$		c	107			
		aq	105			
$(\text{NH}_4)_4\text{Fe}(\text{CN})_6$		aq	108			
$(\text{NH}_4)_4\text{Fe}(\text{CN})_6 \cdot 6\text{H}_2\text{O}$		c	108			
$\text{HFeCO}(\text{CN})_5^{--}$		aq	106			



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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$H_2FeCO(CN)_5^-$		aq	106			
$H_3FeCO(CN)_5$		aq	109			
$H_3FeCO(CN)_5 \cdot H_2O$		c	109			
FeSi		c	110			111
Fe <sub>3</sub> Si		c	112,113			
FeSiO <sub>3</sub>		c	115,116,114			
Fe <sub>2</sub> SiO <sub>4</sub>		c	22	88		88
2FeI <sub>2</sub> ·PbI <sub>2</sub>		c	49			
Zn <sub>2</sub> Fe(CN) <sub>6</sub>		c	108			

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| 55. Parravano and de Cesaris-3        | 97. Troost and Hautefeuille-3                               |
| 56. Mannheimer-1                      | 98. Anderson-9  |
| 57. Mixter-15                         | 99. Berthelot-7   |
| 58. Jellinek and Zakowski-1           | 100. Kelley-20  |
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| 60. Berthelot-13                      | 102. Hieber, Behrens, and Teller-1                          |
| 61. Anderson-5                        | 103. Mittasch-1   |
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| 64. Lukes, Prutton, and Turnball-1    | 106. Muller-2   |
| 65. Anderson-14                       | 107. Chretien and Guinchant-1                               |
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| 67. Biltz-6                           | 109. Muller-3   |
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| 75. Berthelot-8                       |   |
| 76. Gay-1                             |   |
| 77. Manchot-2                         |   |
| 78. Biltz-10                          |   |
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Formula	Description	State				
Mn		g	2, 3		1	1
		c	2, 5		2	2
Mn <sup>n+</sup>		g	4, 5			
Mn <sup>++</sup>		aq	1	1		
Mn <sup>+++</sup>		aq	1			
MnO		g	7			
		c	8, 9, 10, 12, 13, 1, 14, 15, 16		19	19
MnO <sub>2</sub>		c	20, 22, 23, 16, 10, 15, 24	17, 21	17, 21, 25	25
MnO <sub>4</sub> <sup>-</sup>		aq	1	1		
Mn <sub>2</sub> O <sub>3</sub>		c	10, 16, 24, 26, 27, 28			32
Mn <sub>3</sub> O <sub>4</sub>		c	9, 16, 22, 29, 15		9	
Mn(OH) <sub>2</sub>		amorp	20, 33	34, 35		
Mn(OH) <sub>3</sub>		amorp	36			
MnF <sub>2</sub>		c	38		37	37, 55
		aq	36			
MnF <sub>3</sub>		aq	1			
MnCl <sub>2</sub>		c	20, 39, 40		41	41
		aq	20, 39, 42, 43			
MnCl <sub>2</sub> ·nH <sub>2</sub> O		c	20, 44, 45, 46, 47, 48, 49, 50			
H <sub>2</sub> MnCl <sub>6</sub>		aq	51			
MnBr <sub>2</sub>		c	52			
		aq	1			
MnBr <sub>2</sub> ·nH <sub>2</sub> O		c	53			
MnBr <sub>3</sub>		aq	54			
MnI <sub>2</sub>		c	52, 56			
		aq	1			
MnI <sub>2</sub> ·nH <sub>2</sub> O		c	57			
MnS		c	39, 20, 33, 58, 59, 60, 70, 71, 72		72	72
MnSO <sub>4</sub>		c	8, 20, 74		73	73
		aq	22, 8, 78	1, 75, 77, 76		
MnSO <sub>4</sub> ·nH <sub>2</sub> O		c	74, 20, 79, 80, 81, 82			32
MnS <sub>2</sub> O <sub>6</sub>		aq	1			
MnS <sub>2</sub> O <sub>6</sub> ·nH <sub>2</sub> O		c	20		83	83
Mn <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c	84			
		aq	1			
MnSe		c	54		85	85
MnTe		c			85	85
Mn(N <sub>3</sub> ) <sub>2</sub>	manganese azide	c	86			
Mn <sub>5</sub> N <sub>2</sub>		c	87, 88, 89			89

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Formula	Description	State				
$Mn_8N_2$		c	89			90
$Mn(NO_3)_2$		c	91, 92			
		aq	20, 92, 93			
$Mn(NO_3)_2 \cdot nH_2O$		c	94			
$Mn(NO_3)_2 \cdot 6H_2O$		liq	94			95
$MnCl_2 \cdot nNH_3$		c	96, 97			
$MnCl_2 \cdot 2NH_4Cl \cdot 2H_2O$		c	48			
$MnBr_2 \cdot nNH_3$		c	96, 98, 99			
$MnI_2 \cdot nNH_3$		c	96, 98, 99			
$MnSO_4 \cdot nNH_3$		c	100			
$MnSO_4 \cdot (NH_4)_2SO_4 \cdot nH_2O$		c	101, 102			
$Mn_2(PO_4)_3$		c	103			
$Mn_3C$		c	15, 10, 11, 29, 9		25	25
$MnCO_3$		c	20, 104, 105, 10, 42, 106, 11, 15		19, 107	19
		aq	1, 61			
$MnC_2O_4$	manganese oxalate	c	62			
$MnC_2O_4 \cdot nH_2O$		c	62, 63			
$Mn(CHO_2)_2$	manganese formate	c	63			
		aq	63			
$Mn(CHO_2)_2 \cdot 2H_2O$		c	63			
$Mn(C_2H_3O_2)_2$	manganese acetate	c	63, 64			
		aq	63, 64			
$Mn(C_2H_3O_2)_2 \cdot 4H_2O$		c	63, 64			
$MnSiO_3$		c	15, 65, 66, 67		68	68
		gls	65			
$MnI_2 \cdot PbI_2$		c	69			
$MnI_2 \cdot PbI_2 \cdot 3H_2O$		c	69			
$ZnMnO_4$		aq				30
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7. Gaydon-1			17. Hutchinson-1			
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9. Roth and Muller-6			19. Kelley-24			
10. Ulich and Siemonsen-1			20. Thomsen-16			

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| 23. Mixter-11                       | 65. Mulert-1                                 |
| 24. Kapustinskiĭ and Bayushkina-1   | 66. Wologdine-1                              |
| 25. Kelley and Moore-1              | 67. Wologdine and Pankiewitsch-1             |
| 26. Drucker and Hüttner-1           | 68. Kelley-14                                |
| 27. Meyer and Rotger-1              | 69. Mosnier-1                                |
| 28. Biltz-15                        | 70. Vanyukov and Kiseleva-1                  |
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| 31. Millar-3                        | 73. Moore and Kelley-1                       |
| 32. Kelley-18                       | 74. Perreu-10                                |
| 33. Berthelot-13                    | 75. Seidell-1                                |
| 34. Näsänen-4                       | 76. Harned and Owen-1                        |
| 35. Fox, Swinehart, and Garrett-1   | 77. Robinson and Jones-1                     |
| 36. Petersen-4                      | 78. Plake-1                                  |
| 37. Stout and Adams-1               | 79. Perreu-1                                 |
| 38. Jellinek and Koop-1             | 80. Perreu-6                                 |
| 39. Könnecker and Biltz-1           | 81. Biltz-6                                  |
| 40. Sano-14                         | 82. Biltz-19                                 |
| 41. Kelley and Moore-2              | 83. Kelley and Moore-4                       |
| 42. Kapustinskiĭ-7                  | 84. Beck-1                                   |
| 43. Walkley-1                       | 85. Kelley-9                                 |
| 44. Sano-16                         | 86. Wöhler and Martin-1                      |
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| 46. Voskresenskaya and Ponomareva-2 | 88. Neumann, Kröger, and Haebler-2           |
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| 48. Foote and Saxton-1              | 90. Sato-16                                  |
| 49. Sabatier-2                      | 91. Guntz and Martin-1                       |
| 50. Lescoeur-3                      | 92. Ewing, Glick, and Rasmussen-1            |
| 51. Berthelot-68                    | 93. Morgan and Owen-1                        |
| 52. Devoto and Guzzi-1              | 94. Shomate and Young-1                      |
| 53. Lescoeur-7                      | 95. Kelley-27                                |
| 54. Fabre-1                         | 96. Biltz and Hüttig-2                       |
| 55. Kelley-25                       | 97. Hart and Partington-1                    |
| 56. Mosnier-1                       | 98. Ephraim-1                                |
| 57. Lescoeur-1                      | 99. Ephraim-12                               |
| 58. Jellinek and Podjaski-1         | 100. Ephraim-9                               |
| 59. Kelley-22                       | 101. Graham-2                                |
| 60. Kapustinskiĭ and Korshunov-1    | 102. Ephraim and Wagner-1                    |
| 61. Kelley-20                       | 103. Berthelot-102                           |
| 62. Smith and Topley-1              | 104. Bichowsky and Rossini-1                 |

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Table 49  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Cr		g	1, 2, 4		3	3
		c	2, 3		2	2
$\text{Cr}^{n+}$		g	5, 6			
$\text{Cr}^{++}$		aq	3			
$[\text{Cr}(6\text{H}_2\text{O})]^{+++}$		aq	3			
CrO		g	7			
CrO <sub>3</sub>		c	8, 9, 10, 11			
		aq	21, 22, 23			
CrO <sub>4</sub> <sup>--</sup>		aq	3		3	
Cr <sub>2</sub> O <sub>3</sub>		c	8, 10, 11, 12, 13, 14, 15		2	2
Cr <sub>2</sub> O <sub>3</sub> ·nH <sub>2</sub> O		c	16			
Cr <sub>2</sub> O <sub>7</sub> <sup>--</sup>		aq	3	3, 17		
Cr <sub>2</sub> O <sub>9</sub> <sup>--</sup>		aq	20			
Cr <sub>7</sub> H <sub>2</sub>		c	18			
$[\text{Cr}(5\text{H}_2\text{O})(\text{OH})]^{++}$		aq	35			
HCrO <sub>4</sub> <sup>-</sup>		aq	3	17		
$[\text{Cr}(4\text{H}_2\text{O})(\text{OH})_2]^+$		aq	35			
H <sub>2</sub> CrO <sub>4</sub>		aq	3	17		
Cr(OH) <sub>3</sub>		c	19			
$[\text{Cr}(4\text{H}_2\text{O})(\text{OH})_2](\text{OH})$		c	19			
$[\text{Cr}(5\text{H}_2\text{O})(\text{OH})](\text{OH})_2$		c	19			
HCr <sub>2</sub> O <sub>9</sub> <sup>--</sup>		aq	20			
CrF <sub>2</sub>		c	24			
CrF <sub>3</sub>		c	25, 26			
$[\text{Cr}(6\text{H}_2\text{O})]\text{F}_3$		aq	27			
H <sub>3</sub> $[\text{Cr}(6\text{H}_2\text{O})]\text{F}_6$		aq	27			
$[\text{Cr}(5\text{H}_2\text{O})\text{Cl}]^{++}$		aq	3			
CrCl <sub>2</sub>		c	3, 28, 19		2, 30	2, 30
		aq	19			
CrCl <sub>2</sub> ·nH <sub>2</sub> O		c	19, 29			
$[\text{Cr}(4\text{H}_2\text{O})\text{Cl}_2]^+$		aq	3			
CrCl <sub>3</sub>		c	19, 28, 25		2, 30, 31	2, 30
$[\text{Cr}(4\text{H}_2\text{O})\text{Cl}_2]\text{Cl}$		c	36			
		aq	32, 33, 34			
$[\text{Cr}(4\text{H}_2\text{O})\text{Cl}_2]\text{Cl} \cdot \frac{n}{2}\text{H}_2\text{O}$		c	19, 36			
$[\text{Cr}(5\text{H}_2\text{O})\text{Cl}]\text{Cl}_2$		aq	3			
$[\text{Cr}(6\text{H}_2\text{O})]\text{Cl}_3$		c	19			
		aq	19, 35			
CrCl <sub>4</sub>		g	28			
CrO <sub>2</sub> Cl <sub>2</sub>		liq	37, 38, 39			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$[\text{Cr}(4\text{H}_2\text{O})(\text{OH})_2]\text{Cl}$		aq	35			
$[\text{Cr}(5\text{H}_2\text{O})(\text{OH})]\text{Cl}_2$		aq	35			
$[\text{Cr}(4\text{H}_2\text{O})\text{Br}_2]\text{Br}$		aq	19			
$[\text{Cr}(4\text{H}_2\text{O})\text{Br}_2]\text{Br} \cdot 2\text{H}_2\text{O}$		c	19			
$[\text{Cr}(6\text{H}_2\text{O})]\text{Br}_3$		c	19			
		aq	19			
$\text{CrI}_2$		c	40			
		aq	3			
$[\text{Cr}_2(6\text{H}_2\text{O})(\text{SO}_4)_3]$		aq	41, 43			
$[\text{Cr}_2(6\text{H}_2\text{O})(\text{SO}_4)_3] \cdot 2\text{H}_2\text{O}$		c	43			
$[\text{Cr}_2(8\text{H}_2\text{O})(\text{SO}_4)_2] (\text{SO}_4)$		aq	41, 43			
$[\text{Cr}_2(10\text{H}_2\text{O})(\text{SO}_4)_2]$		aq	41, 43			
$[\text{Cr}_2(12\text{H}_2\text{O})(\text{SO}_4)_3]$		aq	41, 32, 43			
$[\text{Cr}_2(12\text{H}_2\text{O})](\text{SO}_4)_3 \cdot n\text{H}_2\text{O}$		c	43, 42			
$\text{CrN}$		c	44, 45			
$\text{Cr}_2\text{N}$		c	45			
$(\text{NH}_4)_2\text{CrO}_4$		c	23			
		aq	21, 95, 3			
$(\text{NH}_4)_2\text{Cr}_2\text{O}_7$		c	47, 46			
		aq	47, 46, 21			
$\text{CrSb}$		c				48
$\text{CrSb}_2$		c				48
$\text{Cr}_3\text{C}_2$		c	15, 49		15	15
$\text{Cr}_4\text{C}$		c	15		15	15
$\text{Cr}_7\text{C}_3$		c	15		15	15
$\text{PbCrO}_4$		c	50, 51			
$2\text{CrI}_2 \cdot \text{PbI}_2$		c	40			
$2\text{CrI}_2 \cdot \text{PbI}_2 \cdot 3\text{H}_2\text{O}$		c	40			
$\text{Ag}_2\text{CrO}_4$		c	54	52, 53	54	54
$\text{FeCr}_2\text{O}_4$		c	55		56	56



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| 3. Calculated                                       | 45. Sano-4                         |
| 4. Maier-2  | 46. Berthelot-95                   |
| 5. Bacher and Goudsmit-1                            | 47. Moles and Gonzales-1           |
| 6. Phillips-1                                       | 48. Schimpff-1                     |
| 7. Gaydon-1   | 49. Heusler-1                      |
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| 10. Mixter-16                                       | 52. Cann and Mueller-1             |
| 11. Roth and Becker-2                               | 53. Murgulescu-1                   |
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| 23. Sabatier-5                                      |                                    |
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| 25. von Wartenberg-12                               |                                    |
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| 32. Thomsen-16                                      |                                    |
| 33. Berthelot-96                                    |                                    |
| 34. Neumann, Kröger, and Kunz-1                     |                                    |
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| 40. Mosnier-1                                       |                                    |
| 41. Colson-2  |                                    |
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Table 50  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Mo		g	1		2	
		c			3	3
Mo <sup>+</sup>		g	4			
MoO <sub>2</sub>		c	5, 6, 7			
MoO <sub>3</sub>		c	5, 8, 9, 10		11	
		aq	12			
MoO <sub>4</sub>		aq	14			
MoO <sub>4</sub> <sup>--</sup>		aq	13, 14			
MoO <sub>5</sub>		aq	14			
H <sub>2</sub> MoO <sub>4</sub>		c	13			
		aq	13			
H <sub>2</sub> MoO <sub>4</sub> ·H <sub>2</sub> O		c	13			
MoCl <sub>2</sub>		c	2			
MoCl <sub>3</sub>		c	2			
MoCl <sub>4</sub>		c	2			
MoCl <sub>5</sub>		c	2			
MoCl <sub>6</sub>		c	2			
MoBr <sub>2</sub>		c	2			
MoBr <sub>3</sub>		c	2			
MoBr <sub>4</sub>		c	2			
MoBr <sub>5</sub>		c	2			
MoI <sub>2</sub>		c	2			
MoI <sub>3</sub>		c	2			
MoI <sub>4</sub>		c	2			
MoI <sub>5</sub>		c	2			
MoS <sub>2</sub>		c		15, 17	16	15
MoS <sub>3</sub>		c	17			
Mo <sub>2</sub> C		c	18	18		
PbMoO <sub>4</sub>		c	19			
CuMoO <sub>4</sub>		c	19			
FeMoO <sub>4</sub>		c	19			
Fe <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>		c	19			

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Table 50 (Continued)  
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13. Pechard-1
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17. Kelley-22
18. Kelley-23
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Table 51  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
W		g	2		1, 3	3
		c			4	4
W <sup>+</sup>		g	1			
WO <sub>2</sub>		c	5, 6, 7			
WO <sub>3</sub>		c	7, 8, 9, 10, 11, 12, 13		13	
WO <sub>4</sub> <sup>--</sup>		aq	15			
W <sub>2</sub> O <sub>5</sub>		c	5, 16			
H <sub>2</sub> WO <sub>4</sub>		c	15			
WCl <sub>2</sub>		c	2			
WCl <sub>4</sub>		c	2			
WCl <sub>5</sub>		c	2			
WCl <sub>6</sub>		c	2, 18			
WBr <sub>2</sub>		c	2			
WBr <sub>4</sub>		c	2			
WBr <sub>5</sub>		c	2			
WBr <sub>6</sub>		c	2			
WI <sub>2</sub>		c	2			
WI <sub>4</sub>		c	2			
WI <sub>5</sub>		c	2			
WS <sub>2</sub>		c	17			
WC		c	19			
CuWO <sub>4</sub>		c	20			
CuWO <sub>4</sub> ·2H <sub>2</sub> O		c	20			
FeWO <sub>4</sub>		c	20			
FeWO <sub>4</sub> ·3H <sub>2</sub> O		c	20			
Fe <sub>2</sub> (WO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O		c	20			
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Table 52  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
V		g	1	1,2	2	2
		c			2	2
$Vn^+$		g	3			
VO		g	4			
$VO_4^-$		aq	5			
$VO_5^-$		aq	2			
$V_2O_2$		c	6,7			
$V_2O_3$		c	8,11	10	9	2
$V_2O_4$		c	7,8,12	13,14,15	9	2
$V_2O_5$		c	7,8,11		9	2
$V_3O_9^{---}$		aq	2			
$HV_6O_{17}^{---}$		aq		16		
$H_2V_6O_{17}^{--}$		aq		16		
$VCl_2$		c	11		17	17
$VCl_3$		c	11	18	17	
$VCl_4$		liq	11			
$VOCl_3$		c	11			
$VOSO_4$		c	14,19			
VN		c	21		20	20
$NH_4VO_3$		c	22		23	23
VC		c			20	20

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Table 53  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Cb  Cb <sub>2</sub> O <sub>4</sub> Cb <sub>2</sub> O <sub>5</sub>		g	3		1, 2	1, 2
		c			1	
		c	4, 5, 6			
		c	7			8
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Table 54  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Ta		g	1,2,3		1,4	1,4
		c			5	5
Ta <sub>2</sub> O <sub>5</sub>		c	6,7,8		10	10
TaN		c	7			9
TaC		c			10	10

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8. Moose and Parr-1
9. Sato-12
10. Kelley-12

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Table 55  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Ti		g	2,27	1,3	1,3	1,3
		c			4	4
Ti <sup>n+</sup>		g	3,5			
TiO		g	6			
		c			7	7
TiO <sub>2</sub>	rutile, III	c	8,9,10,11,12		13	13
	anatase, III	c			13	13
		amorp	14			
Ti <sub>2</sub> O <sub>3</sub>		c	2	2	7	7
Ti <sub>3</sub> O <sub>5</sub>		c	2	2	7	7
TiF <sub>2</sub>		c	2			
TiF <sub>3</sub>		c	2			
TiF <sub>4</sub>		c	2			
H <sub>2</sub> TiF <sub>6</sub>		aq	16			
TiCl		g	6			
TiCl <sub>2</sub>		c	2			
TiCl <sub>3</sub>		c	2			
TiCl <sub>4</sub>		g			1,2	1,2
		liq	15		16,17,19	16,18,20
		aq	15			
TiBr <sub>2</sub>		c	2			
TiBr <sub>3</sub>		c	2			
TiBr <sub>4</sub>		c	2			
TiI <sub>2</sub>		c	2			
TiI <sub>3</sub>		c	2			
TiI <sub>4</sub>		c	2			
TiCl <sub>4</sub> ·nH <sub>2</sub> S		c	21			
TiBr <sub>4</sub> ·nH <sub>2</sub> S		c	21			
TiN		c	9	7	7	7
TiCl <sub>4</sub> ·nPH <sub>3</sub>		c	22			
TiBr <sub>4</sub> ·nPH <sub>3</sub>		c	22,2			
TiC		c	23,24	24,4	4	4
FeTiO <sub>3</sub>		c	26,25	26,25	26,25	26,25



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Table 55 (Continued)  
SPECIFIC REFERENCES

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4. Kelley-2
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6. Gaydon-1
7. Shomate-6
8. Roth and Becker-3
9. Neumann, Kröger, and Kunz-2
10. Roth and Wolf-1
11. Mixter-8
12. Sieverts and Gotta-3
13. Shomate-8
14. Roth and Richter-1
15. Thomsen-16
16. Latimer-2
17. Yost and Blair-1
18. Herman-1
19. Kelley-19
20. Kelley-20
21. Biltz and Keunecke-1
22. Hóltje-1
23. Brantly and Beckman-1
24. Naylor-3
25. Shomate, Naylor and Boericke-1
26. Shomate-4
27. Blocher and Campbell-2

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Table 56  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Zr		g	3	3	1,2	1,2
		c			1	
Zr <sup>n+</sup>		g	2			
ZrO <sup>++</sup>		aq	1,4			
ZrO <sub>2</sub>		c	6,7,8,9	6,7,8,9	5	
ZrO <sub>2</sub> ·5H <sub>2</sub> O		amorp	1,4			
ZrO(OH) <sub>2</sub>		c	4			
Zr(OH) <sub>4</sub>		c	4			
Zr(OH) <sub>4</sub> ·nH <sub>2</sub> O		c	4			
ZrF <sub>2</sub>		c	3			
ZrF <sub>3</sub>		c	3			
ZrF <sub>4</sub>		c	3			
ZrCl <sub>2</sub>		c	3			
ZrCl <sub>3</sub>		c	3			
ZrCl <sub>4</sub>		c	3	3	1	
ZrOCl <sub>2</sub>		aq	10			
ZrOCl <sub>2</sub> ·nH <sub>2</sub> O		c	10,14			
ZrBr <sub>2</sub>		c	3			
ZrBr <sub>3</sub>		c	3			
ZrBr <sub>4</sub>		c	3			
ZrOBr <sub>2</sub>		aq	1			
ZrOBr <sub>2</sub> ·nH <sub>2</sub> O		c	1,12			
ZrI <sub>2</sub>		c	3			
ZrI <sub>3</sub>		c	3			
ZrI <sub>4</sub>		c	3			
ZrOSO <sub>4</sub>		aq	11,12,7,15			
Zr(SO <sub>4</sub> ) <sub>2</sub>		c	11,12			
Zr(SO <sub>4</sub> ) <sub>2</sub> ·nH <sub>2</sub> O		c	11,12			
ZrN		c	6	1,6	1	
ZrO(NO <sub>3</sub> ) <sub>2</sub>		aq	1			
ZrO(NO <sub>3</sub> ) <sub>2</sub> ·nH <sub>2</sub> O		c	13			
ZrC		c	16			
ZrSiO <sub>4</sub>		c			17	17

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Table 56 (Continued)  
SPECIFIC REFERENCES

REFERENCES

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3. Brewer et al-1
4. Simon and Fischer-1
5. Kelley-2
6. Neumann, Kröger, and Kunz-2
7. Roth, Börger and Siemonsen-2
8. Roth and Becker-3
9. Weiss and Neumann-1
10. Chauvenet-5
11. Beck-1
12. Chauvenet-6
13. Chauvenet and Nicolle-1
14. Chauvenet-3
15. Chauvenet-7
16. Prescott-1
17. Kelley-14

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Table 57  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Hf		g	3		1	2
		c			2	
HfO <sub>2</sub>		c				
REFERENCES						
1. Calculated						
2. Kelley-24						
3. Roth and Becker-3						

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Table 58  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
B		g	2	2	1, 2	2
		c			3	3
		amorp	2		3	3
B <sup>n+</sup>		g	1			
B <sub>2</sub>		g	5, 6			
BO		g	7	2	2	2
BO <sub>2</sub> <sup>-</sup>		aq	8			
B <sub>2</sub> O <sub>3</sub>		c	8		3, 9	
		gls	8		8	
BH		g	8	2	2	2
B <sub>2</sub> H <sub>6</sub>	diborane	g	11		12	12
B <sub>5</sub> H <sub>9</sub>	pentaborane	g	11		3	3
		liq	3		3	
B <sub>10</sub> H <sub>14</sub>	decaborane	c	2			
HBO <sub>2</sub>		c	14, 15	14		
		aq	8			
H <sub>2</sub> BO <sub>3</sub> <sup>-</sup>		aq	8	8		
H <sub>3</sub> BO <sub>3</sub>		c	14, 16		3	3
		aq	2, 9, 19	17, 18		
H <sub>2</sub> B <sub>4</sub> O <sub>7</sub>		c	15			
BF		g	7			
BF <sub>3</sub>		g	20		21	
		aq	22			
BF <sub>4</sub> <sup>-</sup>		aq	2	23		
HBFe <sub>4</sub>		aq	22			
BCl		g	2, 5, 7			2
BCl <sub>3</sub>		g	24	21		2
		liq	25, 26	21		
BBr		g	2			
BBr <sub>3</sub>		g	2, 27, 24	21		2
		liq	24, 27	24		
B <sub>2</sub> S <sub>3</sub>		c	28, 29			
BN		g	5			
		c		30		30
B <sub>3</sub> N <sub>3</sub> H <sub>6</sub>	hexahydro-s-triaza-triborane, borazole	g			31	31
NH <sub>4</sub> BO <sub>2</sub>		aq	32			
NH <sub>4</sub> BO <sub>3</sub>		aq	32			
NH <sub>4</sub> BO <sub>3</sub> ·H <sub>2</sub> O		c	32			
(NH <sub>4</sub> ) <sub>2</sub> HBO <sub>3</sub>		aq	10			
B <sub>4</sub> C		c			4	4
B(CH <sub>3</sub> ) <sub>3</sub>		liq	13			

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Table 58 (Continued)  
SPECIFIC REFERENCES

## REFERENCES

1. Moore-1
2. Calculated
3. Johnston et al-1
4. Kelley-24
5. Gaydon-1
6. Douglas and Herzberg-1
7. Herzberg-9
8. Owen-1
9. Southard-2
10. Berthelot-4
11. Prosen, Johnson, and Yenchius-1
12. Webb, Neu, and Pitzer-1
13. Long and Norrish-1
14. von Stackelberg, Quatram, and Dressel-1
15. Gilbert and Levi-1
16. Roth, Börger, and Bertram-1
17. Seidell-1
18. Blasdale and Slansky-1
19. Torgeson and Shomate-1
20. Hammerl-2
21. Spencer-2
22. Thomsen-16
23. Latimer-1
24. Kelley-19
25. Berthelot-50,155
26. Troost and Hautefeuille-4,5
27. Berthelot-50
28. Sabatier-3,6
29. Stock and Popenberg-1
30. Kelley-23
31. Crawford and Edsall-1
32. Tanatar-10

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Table 59  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Al		g		2,3,4	1,5	1,5
		c			6	6
Al <sup>n+</sup>		g	1			
Al <sup>+++</sup>		aq	7	5,7		
AlO		g	8,9			
AlO <sub>2</sub> <sup>-</sup>		aq	10			
Al <sub>2</sub> O <sub>3</sub>	corundum	c	11,13		12	12
Al <sub>2</sub> O <sub>3</sub> ·H <sub>2</sub> O		c	5		14	14
Al <sub>2</sub> O <sub>3</sub> ·3H <sub>2</sub> O	hydrargillite	c	10,15		14	14
AlH		g	9,16			
Al(OH) <sub>3</sub>		amorp	17,18,19			
AlF		g	9			
AlF <sub>3</sub>		c	5,19		3	
		aq	20,21			
AlF <sub>3</sub> ·nH <sub>2</sub> O		c	19	22,23	5	
H <sub>3</sub> AlF <sub>6</sub>		aq	24			
AlCl		g	9,16			
AlCl <sub>3</sub>		c	25,26		3	3
		aq	27,28,29,30,31			
AlCl <sub>3</sub> ·6H <sub>2</sub> O		c	33	23,32		
Al <sub>2</sub> Cl <sub>6</sub>		g	4,34			
AlBr		g	9,16			
AlBr <sub>3</sub>		c	34,35		3	34
		aq	17			
AlI		g	9,16			
AlI <sub>3</sub>		c	34,35		5	
		aq	17			
Al <sub>2</sub> S <sub>3</sub>		c	36		5	
Al <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c	25		37	37
		aq	18			
Al <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·nH <sub>2</sub> O		c	13		37	37
AlCl <sub>3</sub> ·nSO <sub>2</sub>		c	19			
AlCl <sub>3</sub> ·nH <sub>2</sub> S		c	19,38			
AlBr <sub>3</sub> ·H <sub>2</sub> S		c	19,38			
AlI <sub>3</sub> ·nH <sub>2</sub> S		c	19,38			
AlN		c	39		41	40
Al(NO <sub>3</sub> ) <sub>3</sub>		aq	29	23,32		
Al(NO <sub>3</sub> ) <sub>3</sub> ·nH <sub>2</sub> O		c	25		5,42	42
AlF <sub>3</sub> ·2NH <sub>4</sub> F·3/2H <sub>2</sub> O		c	19			
AlCl <sub>3</sub> ·nNH <sub>3</sub>		c	19,35			19
AlCl <sub>3</sub> ·NH <sub>4</sub> Cl		c	5,19			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 59 (Continued)  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$\text{AlCl}_3 \cdot \text{NH}_4\text{Cl} \cdot 6\text{NH}_3$		c	5, 19			
$\text{AlBr}_3 \cdot n\text{NH}_3$		c	35			
$\text{AlI}_3 \cdot n\text{NH}_3$		c	35			
$(\text{NH}_4)\text{Al}(\text{SO}_4)_2$		c	13		13	13
		aq	13, 43			
$(\text{NH}_4)\text{Al}(\text{SO}_4)_2 \cdot 12\text{H}_2\text{O}$		c	13		13	13
$(\text{NH}_4)_2\text{O} \cdot 3\text{Al}_2\text{O}_3 \cdot 4\text{SO}_3 \cdot 6\text{H}_2\text{O}$	ammonium alunite	c	13			
$(\text{NH}_4)_2\text{O} \cdot 3\text{Al}_2\text{O}_3 \cdot 5\text{SO}_3 \cdot 9\text{H}_2\text{O}$	ammonium basic alum	c	13			
$\text{AlCl}_3 \cdot \text{PH}_3$		c	44			
$\text{AlBr}_3 \cdot \text{PH}_3$		c	44			
$\text{AlI}_3 \cdot \text{PH}_3$		c	44			
$\text{Al}_4\text{C}_3$		c	45		46	46
$\text{Al}(\text{CH}_3)_3$		liq	47			
$\text{Al}_2\text{SiO}_5$	andalusite	c	48		49	49
	disthene	c	48		49	49
	sillimanite	c	48		49	49
$2\text{AlI}_3 \cdot 3\text{PbI}_2$		c	50			
$2\text{AlI}_3 \cdot 3\text{PbI}_2 \cdot 10\text{H}_2\text{O}$		c	50			
$\text{AlCu}$		c	51, 52, 53, 54			
$\text{AlCu}_2$		c	51, 52			
$\text{Al}_2\text{Cu}$		c	51, 53			
$\text{AlCl}_3 \cdot \text{AgCl}$		c	19			
$\text{AlNi}$		c	51			
$\text{AlNi}_3$		c	51			
$\text{Al}_2\text{Ni}$		c	51			
$\text{Al}_3\text{Ni}$		c	51			
$\text{AlCo}$		c	18, 29, 51, 54, 55			
$\text{Al}_4\text{Co}$		c	51, 54			
$\text{Al}_5\text{Co}_2$		c	18, 29, 51, 54, 55			
$\text{AlFe}$		c	51			
$\text{Al}_2\text{Fe}$		c	51			
$\text{Al}_3\text{Fe}$		c	51, 29, 56, 57			



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Table 59 (Continued)  
SPECIFIC REFERENCES

REFERENCES

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|--|---|
| 1. Moore-1   | 42. Shomate and Kelley-1                      |
| 2. Baur and Brunner-1                                      | 43. Perreu-13                                 |
| 3. Brewer et al-1  | 44. Höltje-1                                  |
| 4. Kelley-19   | 45. Meichsner and Roth-1                      |
| 5. Calculated  | 46. Kelley-23                                 |
| 6. Giauque and Meads-1                                     | 47. Long and Norrish-1                        |
| 7. Latimer and Greensfelder-1                              | 48. Neumann-2                                 |
| 8. Roy-1   | 49. Kelley-24                                 |
| 9. Gaydon-1  | 50. Mosnier-1                                 |
| 10. Roth, Wirths, and Berendt-1                            | 51. Oelsen and Middel-1                       |
| 11. Snyder and Seltz-1                                     | 52. Oelsen-1                                  |
| 12. Johnston et al-1                                       | 53. Rolla-2                                   |
| 13. Kelley, Shomate, Young, Naylor, Salo, and<br>Huffman-1 | 54. Weibke and Kubaschewski-1                 |
| 14. Shomate and Cook-1                                     | 55. Biltz, Wagner, Pieper, and Holverscheid-1 |
| 15. Roth and Richter-1                                     | 56. Biltz and Hasse-1                         |
| 16. Herzberg-9   | 57. Roth, Umbach, and Chale-1                 |
| 17. Berthelot-50   |   |
| 18. Thomsen-16   |   |
| 19. Baud-1   |   |
| 20. Bichowsky and Rossini-1                                |   |
| 21. Fricke and Wullhorst-1                                 |   |
| 22. Ehret and Frere-1                                      |   |
| 23. Latimer-1  |   |
| 24. Mulert-1   |   |
| 25. Young-2  |   |
| 26. Roth and Büchner-1                                     |   |
| 27. Canneri and Rossi-1                                    |   |
| 28. Roth and Wolf-3  |   |
| 29. Young-1  |   |
| 30. Biltz and Hohorst-1                                    |   |
| 31. Richard, Rowe, and Burgess-1                           |   |
| 32. Seidell-1  |   |
| 33. Sabatier-2   |   |
| 34. Fischer-2  |   |
| 35. Klemm and Tanke-1                                      |   |
| 36. Korshunov-4  |   |
| 37. Shomate-2  |   |
| 38. Biltz and Keunecke-1                                   |   |
| 39. Neumann, Kröger, and Haebler-2                         |   |
| 40. Sato-6   |   |
| 41. Kelley-22  |   |

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Table 60  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Sc		g	1		2,3	2,3
		c				
Sc <sup>n+</sup>		g	3			
Sc <sup>+++</sup>		aq	7	5		
Sc <sub>2</sub> O <sub>3</sub>		c				6
ScCl <sub>3</sub>		c	8			
		aq	7	2		
ScBr <sub>3</sub>		c	8			
		aq	2	2		
Sc <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c				6
ScF <sub>3</sub> ·3NH <sub>4</sub> F		c				4
Sc <sub>2</sub> (C <sub>2</sub> O <sub>4</sub> ) <sub>3</sub>	scandium oxalate	c				9
Sc(CHO <sub>2</sub> ) <sub>3</sub>	scandium formate	c				9
REFERENCES						
1. Brewer et al-1 2. Calculated 3. Sitterly-1 4. Kelley-25 5. Noddack and Brukl-1 6. Nilson and Pettersson-2 7. Bommer and Hohmann-2 8. Bommer and Hohmann-3 9. Turska-1						

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Table 61  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Y		g	4		2,3	2,3
		c				
$Y^{n+}$		g	3			
$Y^{+++}$		aq	8	5		
$Y_2O_3$		c				6
$Y(OH)_3$		c	7	9		
$YCl_3$		c	10			
		aq	8	2		
$YI_3$		c	12			
		aq	2	2		
$Y_2(SO_4)_3$		c				6
		aq	7,2	2		
$Y_2(SO_4)_3 \cdot 8H_2O$		c	7	1,11,13		
$Y(NO_3)_2$		c				1
$Y_2(MoO_4)_3$		c				1
REFERENCES						
1. Kelley-25 2. Calculated 3. Sitterly-1 4. Brewer et al-1 5. Noddack and Brühl-1 6. Nilson and Pettersson-2 7. Thomsen-16 8. Bommer and Hohmann-2 9. Endres-1 10. Bommer and Hohmann-3 11. Seidell-1 12. Hohmann and Bommer-1 13. Jackson and Reinacker-1						

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Table 62  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Lu		g	2		2,3	2,3
		c				
Lu <sup>n+</sup>		g	3			
Lu <sup>+++</sup>		aq	4	5		
LuCl <sub>3</sub>		c	6			
		aq	4	2		
LuI <sub>3</sub>		c	1			
		aq	2	2		
Lu <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		aq	2	2		
Lu <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O		c		7		
REFERENCES						
1. Hohmann and Bommer-1						
2. Calculated						
3. Sitterly-1						
4. Bommer and Hohmann-2						
5. Noddack and Brukl-1						
6. Bommer and Hohmann-2						
7. Jackson and Reinäcker-1						

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Table 63  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Yb		g	2	2,3	2,3	
		c				
Yb <sup>n+</sup>		g	3			
Yb <sup>++</sup>		aq		1,4,5		
Yb <sup>+++</sup>		aq	2	5		
YbCl <sub>3</sub>		c	6			
		aq	2	5		
Yb <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		aq	2	2		
Yb <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O		c		7		
REFERENCES						
1. Laitinen and Taebel-1 2. Calculated 3. Sitterly-1 4. Laitinen-1 5. Noddack and Brukl-1 6. Bommer and Hohmann-3 7. Jackson and Reinäcker-1						

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Table 64  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Tm		g			2,3	
		c				
Tm <sup>+</sup>		g				
Tm <sup>+++</sup>		aq	4	5		
TmCl <sub>3</sub>		c	6			
		aq	4	2		
TmI <sub>3</sub>		c	1			
		aq	2	2		
REFERENCES						
1. Hohmann and Bommer-1 2. Calculated 3. Sitterly-1 4. Bommer and Hohmann-2 5. Noddack and Brukl-1 6. Bommer and Hohmann-3						

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Table 65  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Er	erbium acetate	g	4	5		3
		c				
Er <sup>+</sup>		g				
Er <sup>+++</sup>		aq				
Er <sub>2</sub> O <sub>3</sub>		c				
Er(OH) <sub>3</sub>		c				
ErCl <sub>3</sub>		c				
		aq				
ErI <sub>3</sub>		c				
		aq				
Er <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c				
		aq				
Er <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O		c				
Er(C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>3</sub>		aq				
Er(C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>3</sub> ·4H <sub>2</sub> O		c				
REFERENCES						
1. Thomsen-16						
2. Calculated						
3. Nilson and Pettersson-2						
4. Bommer and Hohmann-2						
5. Noddack and Brukl-1						
6. Bommer and Hohmann-3						
7. Jackson and Reinäcker-1						
8. Hohmann and Bommer-1						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 66  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta \bar{R}f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Ho		g				
		c				
Ho <sup>+</sup>		g				
Ho <sup>+++</sup>		aq	4	5		
HoCl <sub>3</sub>		c	6			
		aq	4	2		
HoI <sub>3</sub>		c	1			
		aq	2	2		
Ho <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		aq	2	2		
Ho <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O		c		3		
REFERENCES						
1. Hohmann and Bommer-1 2. Calculated 3. Jackson and Reinäcker-1 4. Bommer and Hohmann-2 5. Noddack and Brukl-1 6. Bommer and Hohmann-3						



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Table 67  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Dy		g	2			
		c				
Dy <sup>n+</sup>		g	3			
Dy <sup>+++</sup>		aq	4	5		
Dy(OH) <sub>3</sub>		c		7		
DyCl <sub>3</sub>		c	6			
		aq	4	2		
DyI <sub>3</sub>		c	8			
		aq	2	2		
Dy <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		aq	2	2		
Dy <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O		c		1		
REFERENCES						
1. Jackson and Reinacker-1 2. Calculated 3. Sitterly-1 4. Bommer and Hohmann-2 5. Noddack and Brukl-1 6. Bommer and Hohmann-3 7. Endres-1 8. Hohmann and Bommer-1						

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Table 68  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Tb		g	2			
		c				
Tb <sup>n+</sup>		g	3			
Tb <sup>+++</sup>		aq	2	5		
TbCl <sub>3</sub>		c	1			
		aq	2	2		
Tb <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		aq	2	2		
Tb <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O		c		4		
REFERENCES						
1. Bommer and Hohmann-3 2. Calculated 3. Sitterly-1 4. Jackson and Reinäcker-1 5. Noddack and Brukl-1						

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Table 69  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Gd		g	2		2, 3	2, 3
		c			2	
Gd <sup>h+</sup>		g	3			
Gd <sup>+++</sup>		aq	4	5	2	
Gd(OH) <sub>3</sub>		c		7		
GdCl <sub>3</sub>		c	6			
		aq	4	2		
GdI <sub>3</sub>		c	8			
		aq	2	2		
Gd <sub>3</sub> (SO <sub>4</sub> ) <sub>2</sub>		aq	2	2		
Gd <sub>3</sub> (SO <sub>4</sub> ) <sub>2</sub> ·8H <sub>2</sub> O		c		10, 11	1, 12	9, 12, 13
REFERENCES						
1. Kelley-24 2. Calculated 3. Sitterly-1 4. Bommer and Hohmann-2 5. Noddack and Brukl-1 6. Bommer and Hohmann-3 7. Endres-1 8. Hohmann and Bommer-1 9. Ahlberg and Clark-1 10. Jackson and Reinäcker-1 11. Seidell-1 12. Coulter and Latimer-1 13. Giauque and Clark-1						

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Table 70  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Eu		g	2		2,3	2,3
		c				
Eu <sup>n+</sup>		g	3			
Eu <sup>++</sup>		aq		5,7,8		
Eu <sup>+++</sup>		aq	2	5		
EuCl <sub>3</sub>		c	6			
		aq	2	2		
Eu <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		aq	2	2		
Eu <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O		c		4		1
REFERENCES						
1. Kelley-25 2. Calculated 3. Sitterly-1 4. Jackson and Reinäcker-1 5. Noddack and Brukl-1 6. Bommer and Hohmann-3 7. McCoy-1 8. Laitinen and Taebel-1						

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Table 71  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Sm		g	2		2, 3	2, 3
		c				
Sm <sup>n+</sup>		g	3			
Sm <sup>+++</sup>		aq	2	5		
Sm(OH) <sub>3</sub>		c		7		
SmCl <sub>3</sub>		c	6			
		aq	2	2		
SmI <sub>3</sub>		c	8			
		aq	2	2		
Sm <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		aq	2	2		
Sm <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·8H <sub>2</sub> O		c		9		
SmCl <sub>3</sub> ·nNH <sub>3</sub>		c	1, 4			
REFERENCES						
1. Matignon and Trannoy-2						
2. Calculated						
3. Sitterly-1						
4. Matignon-8						
5. Noddack and Brukl-1						
6. Bommer and Hohmann-3						
7. Endres-1						
8. Hohmann and Bommer-1						
9. Jackson and Reinäcker-1						

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Table 72  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Pm		g				
		c				
Pm <sup>+</sup>		g				
Pm <sup>+++</sup>		aq	1			
PmCl <sub>3</sub>		c	1			
PmCl <sub>3</sub>		aq	1			
REFERENCES						
1. Calculated						

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Table 73  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$					
Formula	Description	State									
Nd		g	2	5	7	1					
		c									
		g	3								
		aq	4								
		c	13, 14								
		c									
		c	6, 13								
		aq	4								
		c	13								
		c	8, 15, 9								
		aq	2								
		c	15								
		c	15, 9								
		aq	15, 9, 2								
		c	15, 9								
		c	15, 9								
		c	12								
		REFERENCES									
		1. Kelley-25 2. Calculated 3. Sitterly-1 4. Bommer and Hohmann-2 5. Noddack and Brukl-1 6. Bommer and Hohmann-3 7. Endres-1 8. Hohmann and Bommer-1 9. Matignon-22 10. Jackson and Reinäcker-1 11. Seidell-1 12. Matignon and Trannoy-2 13. Matignon-5 14. Muthmann and Weiss-1					15. Matignon-11				

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Table 74  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Pr		g	2			12
		c				
Pr <sup>0+</sup>		g	3			
Pr <sup>++</sup>		aq	4,1,9	5		
PrO <sub>2</sub>		c	15			
Pr <sub>2</sub> O <sub>3</sub>		c	13,14			
Pr <sub>6</sub> O <sub>11</sub>		c	15			
Pr(OH) <sub>3</sub>		c		7		
PrCl <sub>3</sub>		c	1,6			
		aq	4,9	2		
PrCl <sub>3</sub> ·nH <sub>2</sub> O		c	13			
PrI <sub>3</sub>		c	8			
		aq	2	2		11,10
Pr <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		aq	2	2		
Pr <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·BH <sub>2</sub> O		c				
Pr(NO <sub>3</sub> ) <sub>2</sub>		aq	15			
PrAl <sub>4</sub>		c	9			
REFERENCES						
1. Sieverts and Gotta-2 2. Calculated 3. Sitterly-1 4. Bommer and Hohmann-2 5. Noddack and Brukl-1 6. Bommer and Hohmann-3 7. Endres-1 8. Hohmann and Bommer-1 9. Canneri and Rossi-3 10. Jackson and Reinäcker-1 11. Seidell-1 12. Kelley-25 13. Matignon-6 14. Muthmann and Weiss-1			15. Prandtl and Hüttner-1			



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Table 75  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Ce		g	1		2	3
		c				
Ce <sup>n+</sup>		g	4	5, 6, 7, 9	8	
Ce <sup>+++</sup>		aq				
Ce <sup>++++</sup>		aq	10, 11			15
CeO <sub>2</sub>		c	12, 13, 14			
CeO <sub>3</sub> ·2H <sub>2</sub> O		c	16			
Ce <sub>3</sub> H <sub>8</sub>		c	17			
Ce(OH) <sup>+++</sup>		aq		9		
Ce(OH) <sub>2</sub> <sup>++</sup>		aq		9		
CeCl <sub>3</sub>		c	17, 18, 19, 20		22	
		aq	19			
CeI <sub>3</sub>		c	19		22	
		aq	19			
CeS <sub>2</sub>		c	23			
Ce <sub>2</sub> S <sub>3</sub>		c	23			
Ce(SO <sub>4</sub> ) <sub>2</sub>		c	16, 24			15
Ce <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c				
		aq	25	26		
Ce <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·nH <sub>2</sub> O		c	25	26		
CeN		c	20			27
CeCl <sub>3</sub> ·nNH <sub>3</sub>		c	28			
CeHg <sub>4</sub>		c	29			
Ce <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>		c	30			
CeAl <sub>4</sub>		c	18, 21, 31			
Ce <sub>3</sub> Al		c	18, 21, 31			

REFERENCES

1. Brewer et al-1	11. Evans-1
2. Kelley-24	12. Moose and Parr-1
3. Kelley-25	13. Hirsch-1
4. Sitterly-1	14. Muthmann and Weiss-1
5. Noddack and Brukl-1	15. Nilson and Pettersson-2
6. Müller and Schmidt-1	16. Pissarjewsky-1
7. Walters and de Vries-1	17. Sieverts and Gotta-2
8. Kapustinskiĭ-12	18. Biltz and Pieper-1
9. Sherrill, King, and Spooner-1	19. Bommer and Hohmann-2
10. Fontana-1	20. Neumann, Kröger, and Kunz-1

## SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

## SERIES I

National Bureau of Standards

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Table 75 (Continued)  
SPECIFIC REFERENCES

## REFERENCES

- |                              |  |
|------------------------------|--|
| 21. Muthmann and Beck-1      |  |
| 22. Calculated               |  |
| 23. Biltz-1                  |  |
| 24. Beck-1                   |  |
| 25. Thomsen-16               |  |
| 26. Seidell-1                |  |
| 27. Kellenberger and Kraft-1 |  |
| 28. Barre-2                  |  |
| 29. Biltz and Meyer-2        |  |
| 30. Cane-1                   |  |
| 31. Biltz and Hohorst-1      |  |

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Table 76  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$		
Formula	Description	State						
La	lanthanum formate lanthanum cyanamide	g	2	5	1	4		
		c			3			
La <sup>n+</sup>		g	1					
La <sup>+++</sup>		aq	1		6			
La <sub>2</sub> O <sub>3</sub>		c	7, 8, 9, 10, 11, 12, 13				14	
La <sub>3</sub> H <sub>8</sub>		c	15				16	
La(OH) <sub>3</sub>		c			17			
LaCl <sub>3</sub>		c	15, 18, 19, 20					
		aq	12, 18				21	
LaI <sub>3</sub>		c	18					
		aq	18				21	
LaS <sub>2</sub>		c	13					
La <sub>2</sub> S <sub>3</sub>		c	13					
La <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub>		c						14
		aq	22		21			
La <sub>2</sub> (SO <sub>4</sub> ) <sub>3</sub> ·9H <sub>2</sub> O		c			21			14
LaN		c	19, 23					16
La(NO <sub>3</sub> ) <sub>3</sub> ·6H <sub>2</sub> O		c						4
La(CHO <sub>2</sub> ) <sub>3</sub>		c						24
La <sub>2</sub> (CN <sub>2</sub> ) <sub>3</sub>		c	25					
La <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>	c				26			
LaAl <sub>2</sub>	c	20, 27						
LaAl <sub>4</sub>	c	20, 27						
REFERENCES								
1. Sitterly-1			15. Sieverts and Gotta-2					
2. Brewer et al-1			16. Kellenberger and Kraft-1					
3. Kelley-24			17. Sadolin-1					
4. Kelley-25			18. Bommer and Hohmann-2					
5. Noddack and Brukl-1			19. Neumann, Kröger, and Kunz-1					
6. Kapuētinskiĭ-12			20. Canneri and Rossi-1					
7. Roth, Wolf, and Fritz-1			21. Calculated					
8. Moose and Parr-1			22. Thomsen-16					
9. Muthmann and Weiss-1			23. Neumann, Kröger, and Hoebler-1					
10. Kremers and Stevens-1			24. Turska-1					
11. Beck-1			25. Hartmann, Eckelmann, and Beerman-1					
12. Matignon-9			26. Cane-1					
13. Biltz-4			27. Muthmann and Beck-1					
14. Nilson and Pettersson-2								

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Table 79  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Pu		c				
Pu <sup>+++</sup>		aq	1,3			
Pu <sup>++++</sup>		aq	1,2,3,4			
PuO <sub>2</sub>		c	1,5,6			
PuO <sub>2</sub> <sup>+</sup>		aq	5			
PuO <sub>2</sub> <sup>++</sup>		aq	3,5			
PuH <sub>2</sub>		c	7			
PuF <sub>3</sub>		aq	1,2			
PuCl <sub>3</sub>		c	2			
PuOCl		c	2			
PuBr <sub>3</sub>		c	1,2			
		aq	1			
PuI <sub>3</sub>		c	5			
REFERENCES						
1. Calculated 2. Seaborg, Manning, and Katz-1 3. Evans-1 4. Connick et al-1 5. Brewer et al-1 6. Abraham and Davidson-1 7. Johns-1						

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Table 80  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Np		c				
Np <sup>+++</sup>		aq	1, 2, 3			
Np <sup>++++</sup>		aq	1, 2, 3			
NpO <sub>2</sub> <sup>+</sup>		aq	1, 2, 3			
NpO <sub>2</sub> <sup>++</sup>		aq	1, 2, 3			
NpF <sub>3</sub>		c	1			
NpF <sub>4</sub>		c	1			
NpCl <sub>3</sub>		c	1			
		aq	1			
NpCl <sub>4</sub>		c	1, 2			
NpCl <sub>5</sub>		c	1			
NpBr <sub>3</sub>		c	1			
NpBr <sub>4</sub>		c	1			
NpI <sub>3</sub>		c	1			
REFERENCES						
1. Calculated						
2. Brewer et al-1						
3. Seaborg, Manning and Katz-1						
4. Westrum-1						

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Table 81  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
U		g	1, 2			
		c			1	3, 4
U <sup>+</sup>		g	2			
U <sup>+++</sup>		aq	5	1		
U <sup>++++</sup>		aq	5	1		
UO <sub>2</sub>		c	6		1	
UO <sub>2</sub> <sup>+</sup>		aq	1	7		
UO <sub>2</sub> <sup>++</sup>		aq	5		5	
UO <sub>3</sub>		c	6, 8		1	
UO <sub>3</sub> ·nH <sub>2</sub> O		c	8, 9			
UO <sub>4</sub> ·2H <sub>2</sub> O		c	10			
U <sub>3</sub> O <sub>8</sub>		c	1, 6			
UH <sub>3</sub>		c	11			
U(OH) <sup>+++</sup>		aq	5	5		
UF <sub>3</sub>		c	1		12	
UF <sub>4</sub>		c	1, 14		13	12
UF <sub>5</sub>		c	1		1	
UF <sub>6</sub>		g	1		13, 15	
		c	15		13	
UO <sub>2</sub> F <sub>2</sub>		c			16	16
UCl <sub>3</sub>		c	1, 5, 17		1	
UCl <sub>4</sub>		c	1, 17		1	
UCl <sub>5</sub>		c	1		1	
UCl <sub>6</sub>		c	1		1	
UO <sub>2</sub> Cl <sub>2</sub>		aq	9			
UBr <sub>3</sub>		c	1		1	
UBr <sub>4</sub>		c	1		1	
UO <sub>2</sub> Br <sub>2</sub>		aq	9			
UI <sub>3</sub>		c	1		1	
UI <sub>4</sub>		c	1		1	
UCl <sub>3</sub> I		c	1		1	
UBr <sub>3</sub> I		c	1			
UO <sub>2</sub> SO <sub>4</sub>		aq	9	12		
UO <sub>2</sub> SO <sub>4</sub> ·3H <sub>2</sub> O		c	9, 19, 21	19, 20		
U(SO <sub>4</sub> ) <sub>2</sub>		c	18			
UN		c	1		1	
U <sub>2</sub> N <sub>3</sub>		c	1		1	
UO <sub>2</sub> (NO <sub>3</sub> ) <sub>2</sub>		c	8		12	
		aq	12		12	
UO <sub>2</sub> (NO <sub>3</sub> ) <sub>2</sub> ·nH <sub>2</sub> O		c	8, 22	23	12	

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Table 81 (Continued)  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
UC <sub>2</sub>	uranyl acetate	c	1, 24, 25		12	
UO <sub>2</sub> (C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>2</sub>		aq	9			
UO <sub>2</sub> (C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>2</sub> ·2H <sub>2</sub> O		c	9			
UO <sub>2</sub> (C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>2</sub> ·NH <sub>4</sub> C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ·6H <sub>2</sub> O		c	9			
UO <sub>2</sub> CrO <sub>4</sub>		aq	12			
UO <sub>2</sub> CrO <sub>4</sub> ·5 1/2H <sub>2</sub> O		c	9			
REFERENCES						
1. Brewer et al-1			19. de Coninck-2			
2. Sitterly-1			20. Secoy-1			
3. Moore and Kelley-2			21. Colani-1			
4. Ginnings and Corruccini-2			22. Coulter, Pitzer, and Latimer-1			
5. Fontana-1			23. Latimer-1			
6. Mixer-14			24. Heuse and Otto-2			
7. Kraus and Nelson-1			25. Kelley-23			
8. de Forcrand-1						
9. Aloy-1						
10. Pissarjewsky-3						
11. Spedding, Newton, Warf, Johnson, Nottorf, Johns, and Daane-1						
12. Calculated						
13. Brickwedde, Hoge, and Scott-1						
14. Domange and Wohlhuter-1						
15. Weinstock and Crist-1						
16. Wacker and Cheney-1						
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Table 83  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Th		g			1	
		c			3	4
Th <sup>++++</sup>		aq	2, 1			
ThO <sub>2</sub>		c	7, 8, 9, 10			11
ThH <sub>4</sub>		c	12			
Th(OH) <sub>4</sub>		c	2			
ThF <sub>4</sub>		c	14			
ThCl <sub>4</sub>		c	14			
		aq	7, 8			
ThCl <sub>4</sub> ·nH <sub>2</sub> O		c	8			
ThOCl <sub>2</sub>		c	8			
Th(OH)Cl <sub>3</sub> ·H <sub>2</sub> O		c	8			
ThBr <sub>4</sub>		c	14			
		aq	8			
ThBr <sub>4</sub> ·nH <sub>2</sub> O		c	8			
ThOBr <sub>2</sub>		c	8			
ThI <sub>4</sub>		c	8, 14			
ThOI <sub>2</sub>		c	8			
ThOI <sub>2</sub> ·3 1/2H <sub>2</sub> O		c	8			
Th(OH)I <sub>3</sub> ·10H <sub>2</sub> O		c	8			
Th <sub>2</sub> S <sub>3</sub>		c	14			
ThOSO <sub>4</sub>		c	13			
Th(SO <sub>4</sub> ) <sub>2</sub>		c	16			4
		aq	1, 15			
Th(SO <sub>4</sub> ) <sub>2</sub> ·nH <sub>2</sub> O		c	15			
Th <sub>3</sub> N <sub>4</sub>		c	10, 17			18
Th(NO <sub>3</sub> ) <sub>4</sub>		aq	19			
Th(NO <sub>3</sub> ) <sub>4</sub> ·NH <sub>4</sub> NO <sub>3</sub> ·nH <sub>2</sub> O		c	20			
5Th(NO <sub>3</sub> ) <sub>4</sub> ·4NH <sub>4</sub> NO <sub>3</sub> ·nH <sub>2</sub> O		c	20			
ThCl <sub>4</sub> ·NH <sub>4</sub> Cl		c	8, 21			
ThCl <sub>4</sub> ·2NH <sub>4</sub> Cl·10H <sub>2</sub> O		c	8, 21			
ThCl <sub>4</sub> ·nNH <sub>3</sub>		c	8, 21			
[Th(NH <sub>3</sub> ) <sub>6</sub> ]Cl <sub>4</sub>		c	8, 21			
[Th(NH <sub>3</sub> ) <sub>6</sub> ]Cl <sub>4</sub> ·nNH <sub>3</sub>		c	8, 21			
ThC <sub>2</sub>		c	5, 6			



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Table 83 (Continued)  
SPECIFIC REFERENCES

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18. Sato-19
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Table 85  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Be		g	3	3	1	1
		c			2	2
Be <sup>n+</sup>		g	4			
Be <sup>++</sup>		aq	1			
BeO		g	1, 7, 8, 9		5, 6	5
		c	10, 11, 12, 13, 14, 16, 26		2, 17	2
BeO <sub>2</sub> <sup>--</sup>		aq	18, 19			
BeH		g	20		1, 6, 20	1, 6, 20
Be(OH) <sub>2</sub>		c	21, 18, 19			
BeO·Be(OH) <sub>2</sub>		c	22, 23			
BeF <sub>2</sub>		aq	18, 19, 24, 25, 26, 27			
BeCl <sub>2</sub>		c	14			
		aq	18, 19, 29, 30			
	in C <sub>2</sub> H <sub>5</sub> l		31			
BeCl <sub>2</sub> ·4H <sub>2</sub> O		c	29			
BeBr <sub>2</sub>		c	29, 30			
		aq	1			
BeI <sub>2</sub>		c	29, 30			
		aq	1			
BeS		c	32			
BeSO <sub>4</sub>		c	18, 19, 33, 34			
		aq	22, 35			
BeSO <sub>4</sub> ·nH <sub>2</sub> O		c	18, 19, 36, 22, 33			
BeSO <sub>4</sub> ·4BeO		c	34, 37			
BeBr <sub>2</sub> ·2H <sub>2</sub> S		c	38			
BeI <sub>2</sub> ·2H <sub>2</sub> S		c	38			
Be <sub>3</sub> N <sub>2</sub>		c	12, 16			
Be(NO <sub>3</sub> ) <sub>2</sub>		aq	22, 18			
BeCl <sub>2</sub> ·nNH <sub>3</sub>		c	29, 30			
BeBr <sub>2</sub> ·nNH <sub>3</sub>		c	29, 30			
BeI <sub>2</sub> ·nNH <sub>3</sub>		c	29, 30		17	17
Be <sub>2</sub> SiO <sub>4</sub>		c	17			
BeMoO <sub>4</sub>		c	15			
Be(AlO <sub>2</sub> ) <sub>2</sub>		c				28

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Table 85 (Continued)  
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16. Neumann, Kröger, and Haebler-2
17. Kelley-11
18. Matignon and Marchal-2
19. Matignon and Marchal-4
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22. Thomsen-16
23. Latimer-1
24. Mulert-1
25. Petersen-5
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28. Nilson and Pettersson-2
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Table 86  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Mg		g	2,3,4,5		1	1
		c			6	6
Mg <sup>n+</sup>		g	7			
Mg <sup>++</sup>		aq	1	1		
MgO		c	8,9,10,11,12		6,13,14	6,13,14
MgO <sub>2</sub>		c	15			
MgH		g	17		1,16,17	
Mg(OH) <sub>2</sub>		c	20,21,22,23,24,14, 25,26	18,19,1	14	14
MgF		g	17		1,16	1,16
MgF <sub>2</sub>		c	22,27,28		29	29
MgCl		g	17			
MgCl <sub>2</sub>		c	8,30,31		32	32
		aq	20,8,14,23,24,33, 34,35,36,37,38,39, 40,41,42,	1		
MgCl <sub>2</sub> ·nH <sub>2</sub> O		c	8,43,20,44,45,46, 47,48,50,51		32,48	32,48
MgO·MgCl <sub>2</sub>		c	52			
MgO·MgCl <sub>2</sub> ·nH <sub>2</sub> O		c	53			
Mg(ClO <sub>4</sub> ) <sub>2</sub>		c	1,53			
		aq	1,53			
Mg(ClO <sub>4</sub> ) <sub>2</sub> ·nH <sub>2</sub> O		c	1,53			
Mg(OH)Cl		c	54			
MgBr <sub>2</sub>		c	1,55			
		aq	55	1,56		
MgBr <sub>2</sub> ·6H <sub>2</sub> O		c	56	1		
MgI <sub>2</sub>		c	57,58			
		aq	1,59,60	1		
MgS		c	60,61,62			
MgSO <sub>3</sub>		c	63			
MgCO <sub>3</sub> ·nH <sub>2</sub> O		c	63			
Mg·O <sub>2</sub>		c	20,42,64,65		66	
		aq	20,67,68	1		
Mg·O <sub>2</sub> ·nH <sub>2</sub> O		c	64,69,70,71,72			
MgSe		c	73,74,75,41			75
MgTe		c	76,77,78,79,80,81	1	81	81

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Table 86 (Continued)  
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Formula	Description	State	$\Delta H_f^\circ$	$\Delta G_f^\circ$	$S^\circ$	$\log K$
$\text{Mg}(\text{NH}_3)_2^{++}$		aq	1, 21			
$\text{MgCl}_2 \cdot n\text{NH}_3$		c	87			
$\text{MgCl}_2 \cdot 2\text{NH}_3$		aq	21			
$\text{MgBr}_2 \cdot n\text{NH}_3$		c	87			
$\text{MgI}_2 \cdot 2\text{NH}_3$		c	87			
$\text{Mg}(\text{NH}_3)_2\text{SO}_4$		aq	21			
$3\text{MgSO}_3 \cdot (\text{NH}_4)_2\text{SO}_3 \cdot n\text{H}_2\text{O}$		c	88			
$\text{Mg}_3(\text{PO}_4)_2$		c	89			
$\text{MgHPO}_4$		aq	90			
$\text{Mg}(\text{NH}_4)_2\text{PO}_4 \cdot 6\text{H}_2\text{O}$		c	91			
$\text{Mg}_3(\text{AsO}_4)_2$		c	92			
$\text{MgHAsO}_4$		aq	92			
$\text{Mg}(\text{H}_2\text{AsO}_4)_2$		aq	92			
$\text{Mg}(\text{NH}_4)\text{AsO}_4 \cdot 6\text{H}_2\text{O}$		c	92			
$\text{Mg}_3\text{Sb}_2$		c	93			
$\text{Mg}_3\text{Bi}_2$		c	94, 95, 96			
$\text{MgCO}_3$		c	97, 98, 100		95	77
$\text{Mg}(\text{C}_2\text{H}_3\text{O}_3)_2$	magnesium glycolate	c	101			
		aq	20, 101			
$\text{Mg}(\text{C}_2\text{H}_3\text{O}_3)_2 \cdot 2\text{H}_2\text{O}$		c	101, 20			
$\text{MgCl}_2 \cdot 6\text{CH}_3\text{OH}$		c	102			
$\text{MgCl}_2 \cdot 6\text{C}_2\text{H}_5\text{OH}$		c	102			
$\text{MgCN}_2$	magnesium cyanamide	c	103, 104			
$\text{Mg}(\text{CN})_2$		aq	105			
$\text{Mg}(\text{NO}_3)_2 \cdot 6\text{CH}_3\text{OH}$		c	102			
$\text{Mg}(\text{NO}_3)_2 \cdot 6\text{C}_2\text{H}_5\text{OH}$		c	102			
$\text{Mg}_2\text{Si}$		c	106			
$\text{MgSiO}_3$		c	22		107	108
$\text{Mg}_2\text{SiO}_4$	forsterite	c	22		107	108
$\text{Mg}_2\text{Sn}$		c	93, 103, 107, 110			
$\text{Mg}_2\text{P}_6$		c	95			
$2\text{MgI}_2 \cdot \text{PbI}_2$		c	111			
$\text{MgTi}$		c	110			
$\text{MgZn}_2$		c	113			
$\text{MgCd}$		c	113			
$\text{MgBr}_2 \cdot n\text{H}_2\text{O}$		aq	114			
$\text{Mg}(\text{OH})_2 \cdot n\text{H}_2\text{O}$		aq	115			
$\text{Mg}(\text{OH})_2 \cdot n\text{H}_2\text{O}$		aq	115			
$2\text{Hg}(\text{OH})_2 \cdot n\text{H}_2\text{O}$		c	115			
$\text{MgCrO}_4$		c	116			
		aq	116			

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Table 86 (Continued)  
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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
MgCr <sub>2</sub> O <sub>4</sub>		c			117	117
MgTiO <sub>3</sub>		c			118	118
Mg <sub>4</sub> Al <sub>3</sub>		c	96,113			
MgPr		c	119			
Mg <sub>3</sub> Pr		c	119			
MgLa		c	40			
Mg <sub>3</sub> La		c	40			
2La(NO <sub>3</sub> ) <sub>2</sub> · 3Mg(NO <sub>3</sub> ) <sub>2</sub> · 24H <sub>2</sub> O		c	120		120	120
MgCe		c	110			
MgCe <sub>3</sub>		c	110			

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17. Gaydon-1	37. Tammann and Ohler-1
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20. Thomsen-16	40. Canneri and Rossi-1

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Table B6 (Continued)  
SPECIFIC REFERENCES

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| 46. Sano-22                         | 88. Hartog-1                             |
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| 64. Pickering-3                     | 106. Kubaschewski and Villa-1            |
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| 69. Carpenter and Jette-1           | 111. Mosnier-1                           |
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| 72. Biltz-6                         | 114. Varet-6                             |
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Distance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$\text{CaS}_2\text{O}_3$		aq	6, 64			
$\text{CaS}_2\text{O}_3 \cdot \text{H}_2\text{O}$		c		45, 73		
$\text{CaSe}$		c	74		6	
$\text{CaTe}$		c			6	
$\text{CaN}_6$	calcium azide	c	75			
$\text{Ca}_3\text{N}_2$		c	76, 77, 78		6	79
$\text{CaN}_2\text{O}_2 \cdot 4\text{H}_2\text{O}$		c	80			
$\text{Ca}(\text{NO}_2)_2$		c	81			
		aq	6			
$\text{Ca}(\text{NO}_3)_2$		c	15		82	82
		aq	41, 47, 45, 51, 83, 84	6		
	in $\text{C}_2\text{H}_5\text{OH}$		85			
$\text{Ca}(\text{NO}_3)_2 \cdot \text{nH}_2\text{O}$		c	15, 41, 83, 86	6, 86	6	
$\text{Ca}(\text{NH}_2)_2$		c	46			
$\text{Ca}(\text{NO}_3)_2 \cdot \text{Ca}(\text{OH})_2$		c	88			
$\text{Ca}(\text{NC}_3)_2 \cdot \text{Ca}(\text{OH})_2 \cdot 2\frac{1}{2}\text{H}_2\text{O}$		c	88			
$\text{CaCl}_2 \cdot \text{nNH}_3$		c	89, 90, 91, 92, 93			
$\text{CaBr}_2 \cdot \text{nNH}_3$		c	89, 92			
$\text{CaI}_2 \cdot \text{nNH}_3$		c	89			
$\text{Ca}_3\text{F}_2$		c	94			
$\text{Ca}_3(\text{PO}_4)_2$		c	8, 95, 97, 98		96	8, 96
$\text{CaHPO}_4$		c	95, 97		6	
$\text{CaH}_2\text{O}_4 \cdot 2\text{H}_2\text{O}$		c	97, 99	37		
$\text{Ca}(\text{H}_2\text{IO}_4)_2$		c	95, 97			
$\text{Ca}(\text{H}_2\text{IO}_4)_2 \cdot \text{H}_2\text{O}$		c	95, 97			
$\text{Ca}(\text{H}_2\text{IO}_4)_2 \cdot 2\text{H}_2\text{O} \cdot \text{nNH}_3$		c	100			
$\text{Ca}_3(\text{AsO}_4)_2$		c	101			
$\text{CaH}_2\text{O}_4$		aq	101			
$\text{Ca}(\text{H}_2\text{IO}_4)_2 \cdot \text{H}_2\text{O}$		c	101	37		
$\text{Ca}(\text{H}_2\text{IO}_4)_2 \cdot 2\text{H}_2\text{O}$		aq	101			
$\text{Ca}_2\text{F}_2$		c	102			
$\text{Ca}_2\text{Cl}_2$		c	102			
$\text{Ca}_2\text{O}_2$		c	52, 104, 105, 106		107	107
$\text{CaCO}_3$	calcium carbonate	c	108, 109, 110, 111, 114	108, 112, 114	113	115
	aragonite	c	111, 112		113	115
$\text{CaCO}_3 \cdot \text{nNH}_3$		c	116, 117, 118, 119	117, 118, 119	120	120
$\text{Ca}(\text{HCO}_3)_2$	calcium formate	c	120			
		aq	120			
$\text{Ca}(\text{HCO}_3)_2$		aq	121	109, 122		

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$\text{Ca}(\text{C}_2\text{H}_3\text{O}_2)_2$	calcium acetate	c	120			
		aq	120			
$\text{Ca}(\text{C}_2\text{H}_3\text{O}_2)_2 \cdot \text{H}_2\text{O}$		c	120			
$\text{Ca}(\text{C}_2\text{H}_3\text{O}_3)_2$	calcium glycollate	c	122			
		aq	122			
$\text{Ca}(\text{C}_2\text{H}_3\text{O}_3)_2 \cdot n\text{H}_2\text{O}$		c	122			
$\text{Ca}(\text{C}_2\text{H}_5\text{O})_2$	calcium ethylate	c	123			
$\text{Ca}(\text{C}_2\text{H}_5\text{O})_2 \cdot 2\text{C}_2\text{H}_5\text{OH}$		c	123			
$3\text{CaO} \cdot 4\text{C}_2\text{H}_5\text{OH}$		c	124			
$\text{CaCl}_2 \cdot n\text{C}_2\text{H}_5\text{OH}$		c	125			
$\text{CaCN}_2$	calcium cyanamide	c	126, 127, 128			
$\text{Ca}(\text{CN})_2$		c	129			
		aq	130			
$\text{Ca}(\text{NO}_3)_2 \cdot 2\text{CH}_3\text{OH}$		c	131			
$3\text{CaO} \cdot \text{Ca}(\text{CN})_2 \cdot 15\text{H}_2\text{O}$		c	130			
$\text{CaSi}_2$		c	132, 133, 134			
$\text{Ca}_2\text{Si}$		c	134			
$\text{Ca}_2\text{Si}_2$		c	132, 133, 134			
$\text{CaSiO}_3$		c	34, 112, 135, 136, 137		8, 138, 140	8
$\text{Ca}_2\text{SiO}_4$		c	110, 135, 141, 142			
$\text{Ca}_3\text{SiO}_5$		c	141, 143			
$\text{CaSn}$		c	134			
$\text{CaSn}_3$		c	134, 144			
$\text{Ca}_2\text{Sn}$		c	134			
$\text{CaPb}$		c	102			
$\text{CaPb}_3$		c	102			
$\text{Ca}_2\text{Pb}$		c	102			
$\text{CaI}_2 \cdot 2\text{PbI}_2$		c	146			
$\text{CaI}_2 \cdot 2\text{PbI}_2 \cdot 7\text{H}_2\text{O}$		c	146			
$\text{CaTl}$		c	102			
$\text{CaZn}$		c	147			
$\text{CaZn}_2$		c	147			
$\text{CaZn}_5$		c	147, 148			
$\text{CaZn}_{13}$		c	147, 148, 149			
$\text{Ca}_4\text{Zn}$		c	148			
$\text{CaCd}_3$		c	148			
$\text{CaHgBr}_4$		aq	150			
$\text{CaHg}_2\text{Br}_6$		aq	150			
$\text{CaHg}(\text{CN})_4$		aq	151			
$\text{CaHg}_2(\text{CN})_6$		aq	151			
$\text{CaX}_2 \cdot 2\text{Hg}(\text{CN})_2$		aq	151			

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Formula	Description	State				
$\text{CaX}_2 \cdot 2\text{Hg}(\text{CN})_2 \cdot n\text{H}_2\text{O}$	inyoite	c	151	37		
$\text{CaAg}(\text{CN})_3$		aq	150			
$\text{CaAg}_2(\text{CN})_4$		aq	150			
$\text{CaO} \cdot \text{Fe}_2\text{O}_3$		c	152			
$\text{CaH}_2\text{Fe}(\text{CN})_6$		aq	153			
$\text{Ca}_2\text{Fe}(\text{CN})_6$		aq	37, 153			
$\text{Ca}_2\text{Fe}(\text{CN})_6 \cdot 11\text{H}_2\text{O}$		c	153			
$\text{CaCrO}_4$		c	37			
		aq	154			
$\text{CaWO}_4$		c	155			
$\text{CaTiO}_3$		c			156	156
$\text{CaO} \cdot n\text{B}_2\text{O}_3$		c	17, 157		158	158
$\text{CaO} \cdot 2\text{B}_2\text{O}_3$		gls	157, 159			
$2\text{CaO} \cdot n\text{B}_2\text{O}_3$		c	17, 157, 160		158	158
$2\text{CaO} \cdot 3\text{B}_2\text{O}_3 \cdot 13\text{H}_2\text{O}$		c	160			
$3\text{CaO} \cdot \text{B}_2\text{O}_3$		c	139, 157		158	158
$\text{CaAl}_3$		c	145, 148			
$\text{CaO} \cdot \text{Al}_2\text{O}_3$		gls	161			
$2\text{CaO} \cdot \text{Al}_2\text{O}_3$		c	162			
		gls	161			
$2\text{CaO} \cdot \text{Al}_2\text{O}_3 \cdot 5\text{H}_2\text{O}$		c	162			
$3\text{CaO} \cdot \text{Al}_2\text{O}_3$	heulandite	c	143, 162, 163			
		gls	161			
$3\text{CaO} \cdot \text{Al}_2\text{O}_3 \cdot 6\text{H}_2\text{O}$		c	163			
$4\text{CaO} \cdot \text{Al}_2\text{O}_3$		c	162			
$12\text{CaO} \cdot 7\text{Al}_2\text{O}_3$		c	164			
		gls	164			
$3\text{CaCl}_2 \cdot 4\text{AlCl}_3$		c	44			
$\text{CaO} \cdot \text{Al}_2\text{O}_3 \cdot 6\text{SiO}_2$		c	103			
$3\text{CaO} \cdot \text{Al}_2\text{O}_3 \cdot 2\text{SiO}_2$		c	139			
$4\text{CaO} \cdot \text{Al}_2\text{O}_3 \cdot \text{Fe}_2\text{O}_3$		c	145			
$\text{CaMg}_2$		c	13			
$\text{CaCl}_2 \cdot 2\text{MgCl}_2 \cdot 2\text{H}_2\text{O}$		c	166			
$\text{CaCO}_3 \cdot \text{MgCO}_3$		c	165, 167			

Table 87 (Continued)  
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| 5. Ruff and Hartmann-1                | 47. Lange and Streeck-1                   |
| 6. Calculated                         | 48. Nernst and Orthmann-2                 |
| 7. Clusius and Vaughan-1              | 49. Plake-1                               |
| 8. Kelley-24                          | 50. Richards and Dole-1                   |
| 9. Moore-1                            | 51. Thomsen-16                            |
| 10. Latimer, Schutz, and Hicks, Jr.-2 | 52. de Forcrand-22                        |
| 11. Guntz and Bassett Jr.-1           | 53. Ochi-2,4                              |
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| 13. Biltz and Hohorst-1               | 55. Andre-2                               |
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| 15. Young-1                           | 57. Tassilly-1                            |
| 16. Taylor and Wells-1                | 58. Varet-2                               |
| 17. Torgeson and Shomate-1            | 59. de Forcrand-58                        |
| 18. Schwiete and Pranschke-1          | 60. von Wartenberg-13                     |
| 19. Centnerszwer and Blumenthal-4     | 61. Sabatier-1                            |
| 20. de Forcrand-31                    | 62. Anderson-5                            |
| 21. Bergius-1                         | 63. Kelley and Moore-4                    |
| 22. Grundström-1                      | 64. Bichowsky-3                           |
| 23. Herzberg-9                        | 65. Newmann and Wells-1                   |
| 24. Gaydon-1                          | 66. Kelley, Southard, and Anderson-1      |
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| 26. Sabatier-2                        | 68. Trautz and Pakschwer-1                |
| 27. Schwiete and Hey-1                | 69. Zawadski-1                            |
| 28. Moissan-14,17                     | 70. Lange and Monheim-2                   |
| 29. Washburn et al-1                  | 71. Latimer, Hicks Jr., and Schutz-2      |
| 30. Berthelot-21                      | 72. Volzbenski-1                          |
| 31. Hellwege-1                        | 73. Zil'berman and Ivanov-1               |
| 32. Johnson-5                         | 74. Fabre-3                               |
| 33. Guntz-1                           | 75. Wöhler and Martin-2                   |
| 34. Torgeson and Sahama-1             | 76. Guntz and Bassett-2                   |
| 35. Kohlrausch-2                      | 77. Moser and Herzner-1                   |
| 36. Todd-1                            | 78. Franck and Bodea-1                    |
| 37. Seidell-1                         | 79. Sato-13                               |
| 38. Hellwege-1                        | 80. Berthelot-106                         |
| 39. Parker-1                          | 81. Dode'-2                               |
| 40. Favre and Silbermann-3            | 82. Shomate and Kelley-1                  |
| 41. Pickering-9                       | 83. Ewing, Rogers, Miller, and McGovern-1 |
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97. Matignon and Séon-1	139. Tschernobaeff and Wologdine-1
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109. Wells and Taylor-1	151. Varet-2
110. von Gronow and Elsner-1	152. Marchal-7
111. Backstrom-1,5	153. Joannis-8
112. Roth and Chall-1	154. von Wartenberg, Reusch, and Saran-1
113. Frear and Johnston-1	155. Tammann and Westerholt-1
114. Bichowsky and Rossini-1	156. Shomate-4
115. Anderson-8	157. Griveau-1
116. Berthelot-11	158. Kelley, Todd, and Shomate-1
117. Kohlrausch-2	159. King, Torgeson, and Cook-1
118. Aumeras-1	160. Kurnakov, Nikolaev and Chelishcheva-1
119. Pedersen-2	161. Tschernobaeff-1
120. Berthelot-9	162. Cirilli-2
121. Randall-3	163. Thorvaldson, Brown, and Peaker-2
122. de Forcrand-3	164. Roth and Wolf-3
123. de Forcrand-4	165. Backstrom-2
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Formula	Description	State				
Sr	amalgam	g	2, 3		1	1
		c	4	4	1	5
			6			
Sr <sup>+</sup>		g	7			
Sr <sup>++</sup>		g	7			
		aq	8, 9, 10, 11, 12, 13 33	14, 15, 16 17, 18, 19	1	
SrO		c	20, 21, 22		23	23
SrO <sub>2</sub>		c	24, 25, 26, 27, 28			
SrO <sub>2</sub> ·8H <sub>2</sub> O		c	28			
Sr <sub>2</sub> O		c	29			
SrH		g	30	31, 32, 1	31, 32, 1	
SrH <sub>2</sub>		c	9, 34			
Sr(OH) <sub>2</sub>		c	11, 35, 36, 37, 38, 39 21			
		aq	35, 28, 11, 40	40	40	
Sr(OH) <sub>2</sub> ·H <sub>2</sub> O		c	21, 41, 42			
Sr(OH) <sub>2</sub> ·8H <sub>2</sub> O		c	21, 11, 35, 41, 42			
SrF		g	30			
SrF <sub>2</sub>		c	43, 39			
SrCl		g	30			
SrCl <sub>2</sub>		c	44, 45, 46, 47, 29, 48 11, 49		1	5
		aq	9, 10, 11, 49, 50	40	40	
SrCl <sub>2</sub> ·nH <sub>2</sub> O		c	11, 44, 51, 47, 46, 53 54, 55			5
SrOCl <sub>2</sub>		aq	56			
Sr(OCl) <sub>2</sub>		aq	56			
SrCl <sub>2</sub> ·SrO·nH <sub>2</sub> O		c	57, 58			
SrBr <sub>2</sub>		c	11, 44, 59, 60			5
		aq	40, 50	40	40	
SrBr <sub>2</sub> ·nH <sub>2</sub> O		c	51, 11, 44			5
SrBr <sub>2</sub> ·SrO·nH <sub>2</sub> O		c	59, 60			
SrI <sub>2</sub>		c	59, 61, 44			5
		aq	40, 1	40	40	
SrI <sub>2</sub> ·nH <sub>2</sub> O		c	44, 62, 59, 51			5
SrS		g	63, 30			
		c	64			
SrSO <sub>4</sub>		c	11, 65, 66, 67, 68	68		
		aq	68, 40	68, 40	40	
SrS <sub>2</sub> O <sub>6</sub>		aq	40			
SrS <sub>2</sub> O <sub>6</sub> ·4H <sub>2</sub> O		c	11			
SrI <sub>2</sub> ·nSO <sub>2</sub>		c	69			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
SrSe		c	70			
Sr(N <sub>3</sub> ) <sub>2</sub>		c	71,72			
Sr <sub>3</sub> N <sub>2</sub>		c	9			
SrN <sub>2</sub> O <sub>2</sub> ·5H <sub>2</sub> O		c	73			
Sr(NO <sub>2</sub> ) <sub>2</sub>		c	74,75			
		aq	74,75,40			
Sr(NO <sub>3</sub> ) <sub>2</sub>		c	11,76,47,77			5
		aq	76,11,40,50,78	40	40	
Sr(NO <sub>3</sub> ) <sub>2</sub> ·4H <sub>2</sub> O		c	76,47,77,11			
Sr(NH <sub>2</sub> ) <sub>2</sub>		c	9			
Sr(NH <sub>3</sub> ) <sub>6</sub>		c	79			
SrCl <sub>2</sub> ·nNH <sub>3</sub>		c	80,81,82			
SrBr <sub>2</sub> ·nNH <sub>3</sub>		c	80,81			
SrI <sub>2</sub> ·nNH <sub>3</sub>		c	80,81			
(NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub> ·SrSO <sub>4</sub>		c	83			
Sr <sub>3</sub> (PO <sub>4</sub> ) <sub>2</sub>		c	84			
SrHPO <sub>4</sub>		c	84,85			
Sr(H <sub>2</sub> PO <sub>4</sub> ) <sub>2</sub> ·H <sub>2</sub> O		c	84			
Sr <sub>3</sub> (AsO <sub>4</sub> ) <sub>2</sub>		c	86			
SrHASO <sub>4</sub>		aq	86			
Sr(H <sub>2</sub> AsO <sub>4</sub> ) <sub>2</sub>		aq	86			
SrCO <sub>3</sub>		c	11,87,14,15,8,40	14,15,17,18	19	19
Sr(C <sub>2</sub> O <sub>4</sub> )		aq	66	68	1	
Sr(C <sub>2</sub> O <sub>4</sub> )·H <sub>2</sub> O		c		68		
Sr(C <sub>2</sub> O <sub>4</sub> )·2 1/2 H <sub>2</sub> O		c	66		1	
Sr(CHO <sub>2</sub> ) <sub>2</sub>	strontium formate	c	76,46			
		aq	40			
Sr(CHO <sub>2</sub> ) <sub>2</sub> ·2H <sub>2</sub> O		c	76,46			
Sr(HCO <sub>3</sub> ) <sub>2</sub>		aq	88,40	88,40	40	
Sr(C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>2</sub>	strontium acetate	c	66			
		aq	66,40			
Sr(C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ) <sub>2</sub> ·1/2H <sub>2</sub> O		c	66			
Sr(C <sub>2</sub> H <sub>3</sub> O <sub>3</sub> ) <sub>2</sub>	strontium glycollate	c	89			
		aq	89			
SrBr <sub>2</sub> ·1/2C <sub>2</sub> H <sub>5</sub> OH		c	90			
Sr(CN) <sub>2</sub>		aq	91			
Sr(CN) <sub>2</sub> ·4H <sub>2</sub> O		c	91			
SrSi		c	92			
SrSi <sub>2</sub>		c	92			
SrSiO <sub>3</sub>		c	93			
		gls	94			
Sr <sub>2</sub> SiO <sub>4</sub>		c	93			

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Formula	Description	State				
$\text{SrI}_2 \cdot 2\text{PbI}_2$		c	95			
$\text{SrI}_2 \cdot 2\text{PbI}_2 \cdot 7 \text{H}_2\text{O}$		c	95			
$\text{SrHgBr}_4$		aq	96			
$\text{SrHg}_2\text{Br}_6$		aq	96			
$\text{Sr}_2\text{HgBr}_6$		aq	96			
$\text{SrHg}(\text{CN})_4$		aq	97			
$\text{SrHg}_2(\text{CN})_6$		aq	97			
$\text{SrX}_2 \cdot 2\text{Hg}(\text{CN})_2$		aq	97			
$\text{SrX}_2 \cdot 2\text{Hg}(\text{CN})_2 \cdot n\text{H}_2\text{O}$		c	97			
$\text{SrAg}(\text{CN})_3$		aq	98			
$\text{SrAg}_2(\text{CN})_4$		aq	98			
$\text{SrNi}(\text{CN})_4$		aq	99			
$\text{Sr}_3[\text{FeCO}(\text{CN})_5]_2$		c	100,101			
		aq	100,101			
$\text{Sr}_3[\text{FeCO}(\text{CN})_5]_2 \cdot 4\text{H}_2\text{O}$		c	100,101			
$\text{SrWO}_4$		c	102			
$\text{SrAlCl}_5$		c	45			
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5. Kelley-25			25. Blumenthal-2			
6. Smith and Braley-1			26. Holtermann and Laffitte-1			
7. Bacher and Goudsmit-1			27. Holtermann-1			
8. Kapustinskiĭ and Dezideryeva-1			28. de Forcrand-29			
9. Guntz and Benoit-1			29. Benoit-1			
10. Guntz and Raederer-1			30. Gaydon-1			
11. Thomsen-16			31. Humphreys and Fredrickson-1			
12. Lange and Streeck-1			32. Watson and Fredrickson-1			
13. Lehtonen-1			33. Guntz and Benoit-2			
14. Kapustinskiĭ and Stakhanova-1			34. Guntz-17			
15. Kelley-20			35. Berthelot-21			
16. Kapustinskiĭ and Dezideryeva-1			36. Tamaru and Shiomi-1			
17. Townley, Whitney, and Felsing-1			37. Tamaru and Shiomi-2			
18. Hogge and Johnston-1			38. Sano-19			
19. Anderson-8			39. Guntz-1			
20. de Forcrand-49			40. Calculated from the ions			



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| 51. Trip-1                                | 93. Nacken-3                           |
| 52. Sabatier-2                            | 94. Tschernobaeff-1                    |
| 53. Bell-1                                | 95. Mosnier-1                          |
| 54. Schumb-1                              | 96. Varet-6                            |
| 55. Saxter and Lansing-1                  | 97. Varet-2                            |
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| 58. Andre-3                               | 100. Muller-8                          |
| 59. Tassilly-1                            | 101. Muller-9                          |
| 60. Tassilly-5                            | 102. Tammann and Westerholt-1          |
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| 62. Tassilly-4                            |  |
| 63. Mathur-1                              |  |
| 64. Sabatier-1                            |  |
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| 70. Fabre-1                               |  |
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| 72. Wöhler and Martin-2                   |  |
| 73. Berthelot-106                         |  |
| 74. Dode-1                                |  |
| 75. Dode-2                                |  |
| 76. Berthelot-8                           |  |
| 77. Ewing, Rogers, Miller, and McGovern-1 |  |
| 78. Lange and Streeck-2                   |  |
| 79. Siltz and Hüttig-3                    |  |
| 80. Siltz-7                               |  |
| 81. Hüttig-2                              |  |
| 82. Suffington-1                          |  |

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Formula	Description	State				
Ba		g	3,4,5		2	1
		c			2	1
Ba <sup>n+</sup>		g	6,7			
Ba <sup>++</sup>		aq	2			
BaO		g	9,10,11,22,27		2,8	2,8
		c			12,13	13
BaO <sub>2</sub>		c	14,15,16,30			
BaO <sub>2</sub> ·nH <sub>2</sub> O		c	14,17			
Ba <sub>2</sub> O		c	18			
BaH		g	19		2,8	2,8
BaH <sub>2</sub>		c	20,21			
Ba(OH) <sub>2</sub>		c	10,16,22,23		2	
		aq	10,11,17,24,25,26,27,28,29			
Ba(OH) <sub>2</sub> ·nH <sub>2</sub> O		c	10,11,22,27,29			
BaO <sub>2</sub> ·H <sub>2</sub> O <sub>2</sub>		c	17,30			
BaF		g	19			
BaF <sub>2</sub>		c	31,32		33	13
		aq	31,32			
BaCl		g	19			
		c	18			
BaCl <sub>2</sub>		c	11,18,34,35,36,37,38,39,40		41	1
		aq	20,21,42,43,44,45			
BaCl <sub>2</sub> ·nH <sub>2</sub> O		c	11,37,38,39,46,47,48,49,50,51	2		1,13,41
BaOCl <sub>2</sub>		aq	51			
Ba(OC1) <sub>2</sub>		aq	52			
Ba(ClO <sub>2</sub> ) <sub>2</sub>		c	11,53			
Ba(ClO <sub>3</sub> ) <sub>2</sub>		c	11,54			
		aq	11,54			
Ba(ClO <sub>3</sub> ) <sub>2</sub> ·H <sub>2</sub> O		c	11,54			1
Ba(ClO <sub>4</sub> ) <sub>2</sub>		c	53			
		aq	11,55			
Ba(ClO <sub>4</sub> ) <sub>2</sub> ·3H <sub>2</sub> O		c	53			
BaCl <sub>2</sub> ·BaO·nH <sub>2</sub> O		c	56,57			
BaBr <sub>2</sub>		c	11,58	2		
		aq	48			
BaBr <sub>2</sub> ·nH <sub>2</sub> O		c	11,48			
BaOBr <sub>2</sub>		aq	59			
Ba(BrO <sub>3</sub> ) <sub>2</sub> ·H <sub>2</sub> O		c			1,60	1,60
BaBr <sub>2</sub> ·BaO·nH <sub>2</sub> O		c	56			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
BaI <sub>2</sub>		c	56		2	
		aq	2			
BaI <sub>2</sub> ·nH <sub>2</sub> O		c	11,48			
Ba(IO <sub>3</sub> ) <sub>2</sub>		aq	2,61			
Ba(IO <sub>3</sub> ) <sub>2</sub> ·H <sub>2</sub> O		c	2,61			
BaI <sub>2</sub> ·8aO·nH <sub>2</sub> O		c	56			
BaS		g	19			
		c	62			
		aq	11,63			
BaSO <sub>3</sub>		c	64			
BaSO <sub>4</sub>		c	11,65,66,67,68	67	68	68
		aq	11,65,66,67,68			
BaS <sub>2</sub> O <sub>3</sub>		c				1
BaS <sub>2</sub> O <sub>6</sub>		aq	11			
BaS <sub>2</sub> O <sub>6</sub> ·2H <sub>2</sub> O		c	11			
BaS <sub>2</sub> O <sub>8</sub>		aq	69			
BaS <sub>2</sub> O <sub>8</sub> ·4H <sub>2</sub> O		c	69			
BaS <sub>4</sub> O <sub>6</sub>		aq	2.			
BaS <sub>4</sub> O <sub>6</sub> ·2H <sub>2</sub> O		c	70			
Ba(HS) <sub>2</sub>		aq	11			
Ba(HSO <sub>3</sub> ) <sub>2</sub>		aq	64			
BaSO <sub>4</sub> ·H <sub>2</sub> SO <sub>4</sub>		c	71			
BaSO <sub>4</sub> ·2H <sub>2</sub> SO <sub>4</sub> ·H <sub>2</sub> O		c	71			
BaI <sub>2</sub> ·nSO <sub>2</sub>		c	72			
BaSe		c	73,74			
BaSeO <sub>4</sub>		c	73,74			
Ba(N <sub>3</sub> ) <sub>2</sub>		c	75			
		aq	75			
Ba <sub>3</sub> N <sub>2</sub>		c	21			
Ba(NO <sub>2</sub> ) <sub>2</sub>		c	53,76,77,78			
		aq	53,76,77,78			
Ba(NO <sub>2</sub> ) <sub>2</sub> ·H <sub>2</sub> O		c	53,78			
Ba(NO <sub>3</sub> ) <sub>2</sub>		c	40	79	80	80
		aq	11,41,63,81,82,83,84,85			
BaNH		c	21			
Ba(NH <sub>2</sub> ) <sub>2</sub>		c	21			
Ba(NH <sub>3</sub> ) <sub>6</sub>		c	86,87			
BaCl <sub>2</sub> ·8NH <sub>3</sub>		c	87,88,89			
BaBr <sub>2</sub> ·nNH <sub>3</sub>		c	87,88,89			
BaI <sub>2</sub> ·nNH <sub>3</sub>		c	88			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$Ba_3(PO_4)_2$		c	90, 92, 93			
		aq	90, 92, 93			
$BaHPO_4$		c	90, 92, 93			
$8a(H_2PO_2)_2$		aq	11			
$Ba(H_2PO_2)_2 \cdot H_2O$		c	11			
$Ba(H_2PO_4)_2$		c	90, 92, 93			
$Ba_3(AsO_4)_2$		c	94, 95			
$BaHAsO_4 \cdot H_2O$		c	94, 95			
$Ba(H_2AsO_4)_2 \cdot 2H_2O$		c	94, 95			
$8a_3Sb_2$		c	96			
$Ba_3Bi_2$		c	96			
$BaCO_3$	witherite	c	11, 97, 100, 101	98, 99	12, 13	12, 13
		aq	2	9B, 99		
$8aC_2O_4 \cdot 1/2H_2O$	barium oxalate	c	102			
$8aC_2O_4 \cdot nH_2O$		c	2, 102			
$BaO \cdot BaCO_3$		c	97, 103			
$Ba(CHO_2)_2$	barium formate	c	104			
		aq	39, 104			
$Ba(HCO_3)_2$		aq	105		2	
$Ba(C_2H_3O_2)_2$	barium acetate	c	11, 83			
		aq	11, 83			
$Ba(C_2H_3O_2)_2 \cdot 3H_2O$		c	11, 83			
$Ba(C_2H_3O_3)_2$	barium glycollate	c	64			
		aq	64			
$Ba(OC_2H_5)_2$	barium ethylate	c	106			
$3BaO \cdot 4CH_3OH$		c	106			
$3BaO \cdot 4C_2H_5OH$		c	106			
$Ba(HSO_3)_2 \cdot C_2H_2O_2$	barium bisulfite-glyoxal	aq	64			
$8a(HSO_3)_2 \cdot C_2H_2O_2 \cdot 2 \frac{1}{2}H_2O$		c	64			
$Ba(OSO_3C_2H_5)_2$	barium ethylsulfate	aq	11			
$Ba(OSO_3C_2H_5)_2 \cdot 2H_2O$		c	11			
$BaCN_2$	barium cyanamide	c	107			
$Ba(CN)_2$		c	107, 108			
		aq	107, 10B			
$Ba(CN)_2 \cdot nH_2O$		c	107, 10B			
$Ba(CNO)_2$	barium cyanate	c	109			
		aq	109			
$BaSi_3$		c	110			
$Ba_2Si_2$		c	110			
$8aSiO_3$		c	111, 112, 113, 114			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Ba <sub>2</sub> SiO <sub>4</sub>		c	111,113			
BaSiF <sub>6</sub>		c	115			
BaSn <sub>3</sub>		c	96			
Ba <sub>2</sub> Sn		c	96			
BaPb		c	96			
BaPb <sub>3</sub>		c	96			
Ba <sub>2</sub> Pb		c	96			
BaI <sub>2</sub> ·2PbI <sub>2</sub>		c	116			
BaI <sub>2</sub> ·2PbI <sub>2</sub> ·7H <sub>2</sub> O		c	116			
BaHgBr <sub>4</sub>		aq	117			
BaHg <sub>2</sub> Br <sub>6</sub>		aq	117			
Ba <sub>2</sub> HgBr <sub>6</sub>		aq	117			
Ba <sub>4</sub> HgBr <sub>10</sub>		aq	117			
BaHg(CN) <sub>4</sub>		aq	5B			
BaHg <sub>2</sub> (CN) <sub>6</sub>		aq	5B			
BaCl <sub>2</sub> ·2Hg(CN) <sub>2</sub>		aq	58			
BaCl <sub>2</sub> ·2Hg(CN) <sub>2</sub> ·5H <sub>2</sub> O		c	5B			
BaBr <sub>2</sub> ·2Hg(CN) <sub>2</sub>		aq	5B			
BaBr <sub>2</sub> ·2Hg(CN) <sub>2</sub> ·7H <sub>2</sub> O		c	5B			
BaI <sub>2</sub> ·2Hg(CN) <sub>2</sub>		aq	58			
BaI <sub>2</sub> ·2Hg(CN) <sub>2</sub> ·6H <sub>2</sub> O		c	58			
BaAg(CN) <sub>3</sub>		aq	117			
BaAg <sub>2</sub> (CN) <sub>4</sub>		aq	117			
BaPtCl <sub>6</sub>		c	118,119			
		aq	118,119			
BaPtCl <sub>6</sub> ·6H <sub>2</sub> O		c	118,119			
BaOsCl <sub>6</sub>		c	120			
BaPdCl <sub>4</sub>		c	120			
Ba <sub>3</sub> (RhCl <sub>6</sub> ) <sub>2</sub>		c	121			
BaNi(CN) <sub>4</sub>		aq	122			
Ba <sub>2</sub> Fe(CN) <sub>6</sub>		aq	123			
Ba <sub>2</sub> Fe(CN) <sub>6</sub> ·6H <sub>2</sub> O		c	123			
BaH <sub>2</sub> Fe(CN) <sub>6</sub>		aq	123			
Ba <sub>3</sub> [FeCO(CN) <sub>5</sub> ] <sub>2</sub>		c	124			
		aq	124			
Ba <sub>3</sub> [FeCO(CN) <sub>5</sub> ] <sub>2</sub> ·11H <sub>2</sub> O		c	124			
BaMnO <sub>4</sub>		c	125			
BaCrO <sub>4</sub>		c	126			
BaMoO <sub>4</sub>		c	127			
BaWO <sub>4</sub>		c	127			
BaCl <sub>2</sub> ·2AlCl <sub>3</sub>		c	36			
3BaCl <sub>2</sub> ·4AlCl <sub>3</sub>		c	36			

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| 2. Calculated                    | 44. Schwartz and Coblaus-1           |
| 3. Kelley-19                     | 45. Richards and Dole-1              |
| 4. Brewer et al-1                | 46. Baxter and Cooper-1              |
| 5. Rudbery and Lempert-1         | 47. Bell-1                           |
| 6. Bacher and Goudsmit-1         | 4B. Trip-1                           |
| 7. Rasmussen-5                   | 49. Khomjakov-1                      |
| B. Herzberg-9                    | 50. Perreu-2                         |
| 9. de Forcrand-13,49,5B          | 51. Neumann and Muller-2             |
| 10. Berthelot-21                 | 52. Bruni and Levi-1                 |
| 11. Thomsen-16                   | 53. Berthelot-9                      |
| 12. Anderson-1                   | 54. Berthelot-39                     |
| 13. Kelley-24                    | 55. Berthelot-79                     |
| 14. Berthelot-28                 | 56. Tassilly-1                       |
| 15. Biltz-15                     | 57. Andre-1,3                        |
| 16. Le Chatelier-6               | 5B. Varet-2                          |
| 17. de Forcrand-27,2B            | 59. Berthelot-45                     |
| 18. Benoit-1                     | 60. Greensfelder and Latimer-1       |
| 19. Gaydon-1                     | 61. Chroustchhoff-1                  |
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| 22. de Forcrand-52               | 64. de Forcrand-3                    |
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| 24. Steinwehr-1                  | 66. Melchner-1                       |
| 25. Anderson and Noyes-1         | 67. Seidell-1                        |
| 26. Guntz-16                     | 68. Latimer, Hicks Jr., and Schutz-2 |
| 27. de Forcrand-12               | 69. Berthelot-109                    |
| 2B. Tassilly-5                   | 70. Portillo-1                       |
| 29. Tamaru and Shiomi-1,2        | 71. Volchovskii-1                    |
| 30. Berthelot-60                 | 72. Ephraim and Kornblum-1           |
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| 36. Baud-1                       | 7B. Berthelot-26                     |
| 37. Schottky-1                   | 79. Latimer-1                        |
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| 42. Guntz and Benoit-2           | 84. Stackelberg-1                    |

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| 89. Gillespie and Lurie-1          |                               |
| 90. Berthelot-102                  |                               |
| 91. Hart and Partington-1          |                               |
| 92. Berthelot and Louguinine-3     |                               |
| 93. Berthelot-52                   |                               |
| 94. Blarez-2                       |                               |
| 95. Joly-5                         |                               |
| 96. Kubaschewski and Villa-1       |                               |
| 97. Finkelstein-1                  |                               |
| 98. Hogges and Johnston-1          |                               |
| 99. Townley, Whitney and Felsing-1 |                               |
| 100. Roth-6                        |                               |
| 101. Berthelot-8                   |                               |
| 102. Berthelot-9                   |                               |
| 103. Dutoit-1                      |                               |
| 104. Berthelot-29                  |                               |
| 105. Randall and Tamele-1          |                               |
| 106. de Forcrand-2,19,20           |                               |
| 107. Frank and Bank-1              |                               |
| 108. Joannis-1                     |                               |
| 109. Lemoult-1                     |                               |
| 110. Wöhler and Schuff-1           |                               |
| 111. Nacken-3                      |                               |
| 112. Tschernobaeff-1               |                               |
| 113. Grube and Trucksess-1         |                               |
| 114. Jander-1                      |                               |
| 115. Hantke-1                      |                               |
| 116. Mosnier-1                     |                               |
| 117. Varet-6                       |                               |
| 118. Gire-3                        |                               |
| 119. Gire-2                        |                               |
| 120. Puche-1                       |                               |
| 121. Gire and Puche-1              |                               |
| 122. Varet-4                       |                               |
| 123. Joannis-8                     |                               |
| 124. Muller-9                      |                               |
| 125. Kapustinskiĭ and Bayushkina-1 |                               |
| 126. Chroustschoff and Martinoff-1 |                               |

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Ra		g	2		1	1
		c			1	
Ra <sup>n+</sup>		g	3,4,5			
Ra <sup>++</sup>		aq	1	1		
RaO		c	1,6			
RaCl <sub>2</sub>		c			1	
RaCl <sub>2</sub> ·2H <sub>2</sub> O		c		1,7	1	
RaSO <sub>4</sub>		c		1,7	1	
Ra(NO <sub>3</sub> ) <sub>2</sub>		c		1,7	1	
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1. Calculated 2. Brewer et al-1 3. Rasmussen-2 4. Russell-8 5. Rasmussen-3 6. Roth and Schwartz-1 7. Seidell-1						



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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Li		g	2, 3	2, 3	1	1
		c			4, 5	4, 5
Li <sup>n+</sup>		g	1, 6			
Li <sup>+</sup>		aq	7, 9, 10	8, 1, 9, 10, 11		
Li <sub>2</sub>		g	12, 13, 14, 15		2	1
Li <sub>2</sub> O		c	17, 18			
Li <sub>2</sub> O <sub>2</sub>		c	19			
		aq	19			
LiH		g	12		1, 20, 21	1, 20, 21
		c	22, 23, 24		5	5
LiOH		c	10		1	
		aq	25, 26, 27, 23, 29, 30	10		
LiOH·H <sub>2</sub> O		c	10, 31		1	
Li <sub>2</sub> O <sub>2</sub> ·H <sub>2</sub> O <sub>2</sub> ·3H <sub>2</sub> O		c	19			
LiF		c	32, 9		33	33
		aq	34, 35	1		
LiHF <sub>2</sub>		aq	1, 35			
LiCl		g	37		1, 36	1, 36
		c	38, 39, 40, 41, 42, 43, 44, 45, 46			47
		aq	40, 48, 49, 50, 51, 52, 53	1		
	in CH <sub>3</sub> OH		52, 39, 54			
	in C <sub>2</sub> H <sub>5</sub> OH		52, 54, 44			
LiCl·nH <sub>2</sub> O		c	46, 55, 56	8, 17		47
LiClO		aq	57			
LiBr		g	37		1, 36	1, 36
		c	46, 58			47
		aq	49, 58, 60	1		
LiBr·nH <sub>2</sub> O		c	46, 56			47
LiI		g	37		1, 36	1, 36
		c	46, 61, 62, 63			47
		aq	49		1	
LiI·nH <sub>2</sub> O		c	64, 46			47
Li <sub>2</sub> SO <sub>4</sub>		c	65, 41, 66, 67			
		aq	68, 65, 69	1		
Li <sub>2</sub> SO <sub>4</sub> ·H <sub>2</sub> O		c	66, 65, 41			
LiI·nSO <sub>2</sub>		c	70			
Li <sub>2</sub> Se		c	71			
		aq	71			
Li <sub>2</sub> Se·9H <sub>2</sub> O		c	71			
Li <sub>3</sub> N		c	72, 73			74

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$		
Formula	Description	State						
LiNO <sub>2</sub>	in C <sub>2</sub> H <sub>5</sub> OH	c	76	1				
LiNO <sub>3</sub>		c	44, 41, 42					
		aq	50					
			44					
LiNO <sub>3</sub> ·3H <sub>2</sub> O		c	77	1, 93				
LiNH <sub>2</sub>		c	78, 79					
Li(NH <sub>3</sub> ) <sub>4</sub>		liq	80, 81					
Li <sub>2</sub> NH		c	78					
LiCl·nNH <sub>3</sub>		c	45, 82, 83, 84, 85, 86					
LiBr·nNH <sub>3</sub>		c	45, 87, 84, 86					
LiI·nNH <sub>3</sub>		c	86, 84, 88					
Li <sub>3</sub> Sb <sub>2</sub>	in CH <sub>3</sub> OH in C <sub>2</sub> H <sub>5</sub> OH	c	89	1, 8	7	7		
Li <sub>3</sub> Bi		c	90					
Li <sub>2</sub> C <sub>2</sub>		c	91					
Li <sub>2</sub> CO <sub>3</sub>		c	7, 92					
		aq	92, 7, 94					
LiHCO <sub>3</sub>		aq	94, 1					
LiOCH <sub>3</sub>			95					
LiOC <sub>2</sub> H <sub>5</sub>			95					
LiCN		aq	1, 96					
LiCl·nCH <sub>3</sub> NH <sub>2</sub>		c	45, 98, 97	1				
LiCl·nC <sub>2</sub> H <sub>5</sub> NH <sub>2</sub>	c	45						
LiCl·n(CH <sub>3</sub> ) <sub>2</sub> NH	c	97						
LiBr·n(CH <sub>3</sub> ) <sub>2</sub> NH	c	97						
LiBr·nCH <sub>3</sub> NH <sub>2</sub>	c	97						
LiI·n(CH <sub>3</sub> ) <sub>2</sub> NH	c	97						
LiI·nCH <sub>3</sub> NH <sub>2</sub>	c	97						
Li <sub>2</sub> SiO <sub>3</sub>	gls	99						
Li <sub>2</sub> SiO <sub>6</sub>	c	100						
	aq	100, 1						
LiSn	c	89						
LiSn <sub>2</sub>	c	101						
Li <sub>2</sub> Sn	c	101						
Li <sub>4</sub> Sn	c	89						
Li <sub>5</sub> Sn <sub>2</sub>	c	101						
Li <sub>7</sub> Sn <sub>2</sub>	c	89						
LiPb	c	90						
Li <sub>3</sub> Pb	c	101						
Li <sub>4</sub> Pb	c	101						
Li <sub>5</sub> Pb <sub>2</sub>	c	101						
Li <sub>7</sub> Pb <sub>2</sub>	c	90						

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
• 2LiI•PbI <sub>2</sub>		c	61			
2LiI•PbI <sub>2</sub> •4H <sub>2</sub> O		c	61			
LiTl		c	101			
LiHg		c	69			
LiHg <sub>2</sub>		c	69			
LiHg <sub>3</sub>		c	69			
LiHg <sub>99</sub>		liq	102			
nLiBr•HgBr <sub>2</sub>		aq	103			
nLiX•Hg(CN) <sub>2</sub>		aq	104			
LiX•Hg(CN) <sub>2</sub> •nH <sub>2</sub> O		aq	104			
LiCN•Hg(CN) <sub>2</sub>		aq	104			
2LiCN•Hg(CN) <sub>2</sub>		aq	104			
LiReO <sub>4</sub>		c	105			
LiReO <sub>4</sub> •nH <sub>2</sub> O		c	105			
LiCl•ThCl <sub>4</sub> •8H <sub>2</sub> O		c	75			
2LiCl•ThCl <sub>4</sub>		c	75			
LiBH <sub>4</sub>		c	16			16
LiAlH <sub>4</sub>		c	16			16

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| <ol style="list-style-type: none"> <li>1. Calculated</li> <li>2. Gordon-4</li> <li>3. Hartman and Schneider-1</li> <li>4. Simon and Swain-1</li> <li>5. Kelley-24</li> <li>6. Sitterly-1</li> <li>7. Brown and Latimer-1</li> <li>8. Seidell-1</li> <li>9. Payne-1</li> <li>10. Ueda-1</li> <li>11. Robinson and Stokes-2</li> <li>12. Gaydon-1</li> <li>13. Loomis and Nusbaum-2</li> <li>14. Lewis-6</li> <li>15. Bartlett and Furry-1</li> <li>16. Davis, Mason, and Stegeman-1</li> <li>17. de Forcrand-47</li> <li>18. Beketoff-2</li> <li>19. de Forcrand-32</li> <li>20. Crawford and Jorgesen-1</li> </ol> | <ol style="list-style-type: none"> <li>21. Herzberg-9</li> <li>22. Gibb, Jr.-1</li> <li>23. Moers-1</li> <li>24. Kapustinskiĭ, Shamovskii, and Bayushkina-1</li> <li>25. Rossini-4</li> <li>26. Rossini-6</li> <li>27. Gucker and Schminke-1</li> <li>28. Bichowsky and Rossini-1</li> <li>29. Guntz-13</li> <li>30. Zukowsky-1</li> <li>31. de Forcrand-45</li> <li>32. de Forcrand-13</li> <li>33. Clusius-4</li> <li>34. Lange and Leighton-1</li> <li>35. Petersen-4</li> <li>36. Stevenson-2</li> <li>37. Desai-2</li> <li>38. Lange and Martin-3</li> <li>39. Askew, Bullock, Smith, Tinkler, Gatty, and Wolfenden-1</li> </ol> |
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| 44. Pickering-9                    | 86. Hart and Partington-1      |
| 45. Bonnefoi-1                     | 87. Bonnefoi-4                 |
| 46. Slonim and Hüttig-1            | 88. Ephraim-5                  |
| 47. Kelley-25                      | 89. Kubaschewski and Seith-1   |
| 48. Muller-1                       | 90. Seith and Kubaschewski-1   |
| 49. Richards and Rowe-3            | 91. Guntz-15                   |
| 50. Richards and Rowe-2            | 92. de Forcrand-50             |
| 51. Dunnington and Hoggard-1       | 93. Kelley-20                  |
| 52. Lemoine-2                      | 94. Muller-1                   |
| 53. Tucker-2                       | 95. de Forcrand-17             |
| 54. Slansky-1                      | 96. Varet-3                    |
| 55. Bogorodskii-2                  | 97. Simon and Glauner-1        |
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| 59. Bodisko-2                      | 101. Weibke and Kubaschewski-1 |
| 60. Miscenko-1                     | 102. Lewis and Keyes-2         |
| 61. Mosnier-1                      | 103. Varet-6                   |
| 62. Bodisko-1                      | 104. Varet-2                   |
| 63. Beketoff-8                     | 105. Smith and Long-1          |
| 64. Hüttig and Pohle-1             |                                |
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| 66. Pickering-6                    |                                |
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Formula	Description	State				
Na		g	2,3,4,5,6,7,8,10		1,12	1,12
		c			9	9
Na <sup>n+</sup>		g	12			
Na <sup>+</sup>		aq	13		9	
Na <sub>2</sub>		g	2		2	
NaO <sub>2</sub>		c	14			
Na <sub>2</sub> O		c	15,17,18,19		16	10
Na <sub>2</sub> O <sub>2</sub>		c	17,18,20,21,22			
NaH		g	23,25,26		1,24	1,24
		c	27,28,29,30			
NaOH		c	31,32,33,34			10
		aq	1,34,29,35,36,37,38,39,40	1		
NaOH·H <sub>2</sub> O		c	33,45	42,43,44		
NaHO <sub>2</sub>		aq	46			
NaF		g	23			
		c	47,48,49		49	
		aq	34,48	1		
NaHF <sub>2</sub>		c	48,51			
		aq	34			
NaCl		g	1,23,24			
		c	54,55,56,57,58,59,60,61,62,63,64,65,66,67	42,44	52,53	52,53
		aq	50,39,68,69,70,71,72,73	1		
	in CH <sub>3</sub> OH		62,66,74			
NaClO		aq	34,75,76			
NaClO <sub>2</sub>		c	77			
		aq	1,77			
NaClO <sub>3</sub>		c	78,79			
		aq	34			
NaClO <sub>4</sub>		c	66,63,78,80			10
		aq	81,1	1,81		
	in CH <sub>3</sub> OH		66			
	in C <sub>2</sub> H <sub>5</sub> OH		66			
NaBr		g	23			
		c	59,82,66			10
		aq	54,72,82,59,83,84,85,1	1		
	in CH <sub>3</sub> OH		62,66			
	in C <sub>2</sub> H <sub>5</sub> OH		62,66			
NaBr·2H <sub>2</sub> O		c	78,34,86,87			

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Formula	Description	State				
NaBrO		aq	34, 88			
NaI		g	23			
		c	66, 34, 78, 89, 90, 59			10
		aq	72	1		
	in CH <sub>3</sub> OH		66, 62			
	in C <sub>2</sub> H <sub>5</sub> OH		66, 89			
NaI·2H <sub>2</sub> O		c	78, 34, 91			
NaIO <sub>3</sub>		aq	1			
Na <sub>2</sub> S		c	92, 93, 94, 95, 96			
		aq	93, 97, 34			
Na <sub>2</sub> S·nH <sub>2</sub> O		c	93			
Na <sub>2</sub> S <sub>2</sub>		aq	96, 93			
Na <sub>2</sub> S <sub>3</sub>		aq	96, 93			
Na <sub>2</sub> S <sub>4</sub>		c	96, 93			
		aq	93, 96			
Na <sub>2</sub> SO <sub>3</sub>		c	99		100	100
		aq	1, 99, 34, 101			
Na <sub>2</sub> SO <sub>3</sub> ·7H <sub>2</sub> O		c	102, 103			
Na <sub>2</sub> SO <sub>4</sub>		c	34, 104, 105, 106, 107, 108, 109, 67, 110, 111, 112		105	105
		aq	34, 113, 114, 115, 116, 117, 118, 119, 120		1	
Na <sub>2</sub> SO <sub>4</sub> ·10H <sub>2</sub> O		c	122, 123, 105, 87, 111, 67, 112, 78, 34, 124	125, 126, 1	105	105
Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub>		c	127, 128			10
		aq	129, 34, 130, 131			
Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> ·5H <sub>2</sub> O		c	130, 34, 129, 132			
Na <sub>2</sub> S <sub>2</sub> O <sub>5</sub>		c	54, 99			
		aq	54			
Na <sub>2</sub> S <sub>2</sub> O <sub>6</sub>		c	1, 34			
		aq	1, 34			
Na <sub>2</sub> S <sub>2</sub> O <sub>6</sub> ·2H <sub>2</sub> O		c	34			
Na <sub>2</sub> S <sub>3</sub> O <sub>6</sub>		aq	130			
Na <sub>2</sub> S <sub>3</sub> O <sub>6</sub> ·3H <sub>2</sub> O		c	130			
Na <sub>2</sub> S <sub>4</sub> O <sub>6</sub>		aq	34, 130, 132			
Na <sub>2</sub> S <sub>4</sub> O <sub>6</sub> ·2H <sub>2</sub> O		c	130			
NaHS		c	93, 133			
		aq	93, 98, 34, 97			
NaHS·2H <sub>2</sub> O		c	93			
NaHSO <sub>3</sub>		aq	34, 99, 134			

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Formula	Description	State				
NaHSO <sub>4</sub>	sodium azide	c	78, 34			
		aq	135, 34			
NaHSO <sub>4</sub> ·H <sub>2</sub> O		c	136			
NaI·nSO <sub>2</sub>		c	137, 138			
Na <sub>2</sub> Se		c	1, 139			
		aq	1, 139			
Na <sub>2</sub> Se·nH <sub>2</sub> O		c	139			
Na <sub>2</sub> SeO <sub>3</sub>		aq	34			
Na <sub>2</sub> SeO <sub>4</sub>		c	140			
		aq	34, 141, 142			
NaHSe		c	133			
		aq	139			
NaHSeO <sub>3</sub>		aq	34			
NaHSeO <sub>4</sub>		aq	34, 141			
Na <sub>2</sub> Te		c	96			
Na <sub>2</sub> Te <sub>2</sub>		c	96			
Na <sub>2</sub> TeO <sub>4</sub>		c	140			
NaN <sub>3</sub>		c				10
NaNO <sub>2</sub>		c	143, 144, 145			153
		aq	143, 145, 146			
NaNO <sub>3</sub>		c	78, 123, 34, 67, 54, 167, 148, 149, 150, 151, 152, 154, 159			
		aq	39, 117, 34, 118, 149, 84, 151, 155, 156			153
NaNH <sub>2</sub>		c	157, 158			10
NaCl·5NH <sub>3</sub>		c	160, 161			
NaBr·nNH <sub>3</sub>		c	160			
NaI·nNH <sub>3</sub>		c	160, 161			10
Na <sub>2</sub> SO <sub>4</sub> ·(NH <sub>4</sub> ) <sub>2</sub> SO <sub>4</sub> ·H <sub>2</sub> O		c	163			
NaPO <sub>3</sub>		c	164			10
		aq	34, 164, 1			
Na <sub>3</sub> PO <sub>4</sub>		c	165			
		aq	167, 34			10
Na <sub>3</sub> PO <sub>4</sub> ·12H <sub>2</sub> O		c	166			
Na <sub>4</sub> P <sub>2</sub> O <sub>7</sub>		c	34			
		aq	34, 164, 54			10
Na <sub>4</sub> P <sub>2</sub> O <sub>7</sub> ·10H <sub>2</sub> O		c	168			
NaH <sub>2</sub> PO <sub>3</sub>		c	169			
		aq	34			10
NaH <sub>2</sub> PO <sub>3</sub> ·2 1/2H <sub>2</sub> O		c	169			
NaH <sub>2</sub> PO <sub>4</sub>		aq	167, 34			

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Formula	Description	State				
$\text{NaH}_3\text{P}_2\text{O}_7$		c	164			
		aq	34, 164			
$\text{NaH}_3\text{P}_2\text{O}_7 \cdot \text{H}_2\text{O}$		c	164			
$\text{Na}_2\text{HPO}_3$		c	169			
		aq	34			
$\text{Na}_2\text{HPO}_3 \cdot 5\text{H}_2\text{O}$		c	169			
$\text{Na}_2\text{HPO}_4$		c	34, 170, 238			
		aq	34, 167			
$\text{Na}_2\text{HPO}_4 \cdot n\text{H}_2\text{O}$		c	34, 170, 171, 238			
$\text{Na}_2\text{H}_2\text{P}_2\text{O}_5$		c	169			
		aq	169			
$\text{Na}_2\text{H}_2\text{P}_2\text{O}_7$		c	164			
		aq	34, 164			
$\text{Na}_2\text{H}_2\text{P}_2\text{O}_7 \cdot 6\text{H}_2\text{O}$		c	164			
$\text{Na}_3\text{HP}_2\text{O}_7$		c	164			
		aq	164, 34, 1			
$\text{Na}_3\text{HP}_2\text{O}_7 \cdot n\text{H}_2\text{O}$		c	164			
$\text{NaNH}_4\text{HPO}_4$		aq	167			
$\text{NaNH}_4\text{HPO}_4 \cdot 4\text{H}_2\text{O}$		c	34			
$\text{Na}_3\text{AsO}_4$		c	165			
		aq	34			
$\text{Na}_3\text{AsO}_4 \cdot 12\text{H}_2\text{O}$		c	166			
$\text{NaH}_2\text{AsO}_3$		aq	34			
$\text{NaH}_2\text{AsO}_4$		aq	34			
$\text{Na}_2\text{HASO}_4$		aq	34			
$\text{Na}_3\text{Sb}$		c	172, 174			
$\text{Na}_3\text{SbO}_4$		c	165			
$\text{Na}_3\text{SbS}_3$		aq	175			
$\text{Na}_3\text{Bi}$		c	173, 174			
$\text{Na}_3\text{BiO}_4$		c	165			
$\text{Na}_2\text{C}_2$	sodium carbide	c	176, 177		185	185
$\text{Na}_2\text{CO}_3$		c	78, 108, 159, 34, 178, 179, 180, 181, 182, 183, 184			
		aq	180, 134, 117, 186, 187			
$\text{Na}_2\text{CO}_3 \cdot n\text{H}_2\text{O}$		c	34, 183, 159, 182, 188, 189			
$\text{Na}_2\text{C}_2\text{O}_4$	sodium oxalate	c	78			190
		aq	191, 34			
$\text{NaHC}_2$	sodium formate	c	177			
$\text{NaCHC}_2$		c	78			
		aq	34, 78, 191, 192			



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Formula	Description	State				
$\text{NaCHO}_2 \cdot n\text{H}_2\text{O}$		c	193			
$\text{NaHCO}_3$	sodium bicarbonate	c	78, 188, 178, 194		185	185
		aq	186, 34, 117			
$\text{NaCH}_3\text{O}$	in $\text{CH}_3\text{OH}$		195			
$\text{NaHC}_2\text{O}_4$	sodium acid oxalate	c	78			
		aq	34, 191			
$\text{NaHC}_2\text{O}_4 \cdot \text{H}_2\text{O}$		c	78			
$\text{NaC}_2\text{H}_3\text{O}_2$	sodium acetate	c	34, 191, 182			
		aq	34, 191, 1, 196, 197, 198			
	in $\text{C}_2\text{H}_5\text{OH}$		89			
$\text{NaC}_2\text{H}_3\text{O}_2 \cdot 3\text{H}_2\text{O}$		c	194, 34, 182, 78, 199			
$\text{NaC}_2\text{H}_3\text{O}_3$	sodium glycollate	c	200			
		aq	200			
$\text{NaC}_2\text{H}_3\text{O}_3 \cdot 1/2\text{H}_2\text{O}$		c	200			
$\text{NaC}_2\text{H}_5\text{O}$	in $\text{C}_2\text{H}_5\text{OH}$		195			
$\text{NaC}_2\text{H}_5\text{O}_2$	monosodium glycol	c	201			
$\text{NaC}_2\text{H}_5\text{O}_2 \cdot \text{CH}_3\text{OH}$	monosodium glycol methanol	c	201			
$\text{NaC}_2\text{H}_3\text{O}_3 \cdot \text{HC}_2\text{H}_3\text{O}_3$	sodium acid glycollate	c	200			
$\text{NaC}_2\text{H}_5\text{O}_2 \cdot \text{C}_2\text{H}_5\text{OH}$	monosodium glycol ethanol	c	201			
$\text{NaC}_2\text{H}_5\text{O}_2 \cdot \text{C}_2\text{H}_6\text{O}_2$	monosodium glycol glycol	c	201			
$\text{Na}_2\text{C}_2\text{H}_2\text{O}_3$	disodium glycollate	c	200			
		aq	200			
$\text{Na}_2\text{C}_2\text{H}_2\text{O}_3 \cdot 2\text{H}_2\text{O}$		c	200			
$\text{Na}_2\text{CO}_3 \cdot \text{NaHCO}_3 \cdot 2\text{H}_2\text{O}$	trona	c	178			
$\text{NaCCl}_3\text{CO}_2$	sodium trichloroacetate	aq	34, 202			
$\text{NaCH}_2\text{ClCO}_2$	sodium chloroacetate	aq	34, 202			
$\text{NaCHCl}_2\text{CO}_2$	sodium dichloroacetate	aq	34			
$\text{NaI} \cdot 3\text{CH}_3\text{OH}$		c	203			
$\text{NaC}_2\text{H}_5\text{OSO}_3$	sodium ethylsulfate	aq	34			
$2\text{NaHSO}_3 \cdot \text{C}_2\text{H}_2\text{O}_2 \cdot \text{H}_2\text{O}$	sodium bisulfite glyoxal	c	200			
		aq	200			
$\text{NaCN}$		c	204			
		aq	34			
$\text{NaCN} \cdot n\text{H}_2\text{O}$		c	204			
$\text{NaCNO}$	sodium cyanate	c	205			
		aq	1, 205			
$\text{NaCN}_2\text{H}$	sodium cyanamid	aq	205			
$\text{NaCO}_2\text{NH}_2$	sodium carbamate	c	206			

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Formula	Description	State				
NaCNS	sodium thiocyanate	c	207, 208			
		aq	204, 208			
	in C <sub>2</sub> H <sub>5</sub> OH		208			
Na <sub>2</sub> SiO <sub>3</sub>		c	209, 210, 211, 212		213	213
		gls	214, 215, 216			
Na <sub>2</sub> SiO <sub>3</sub> ·nH <sub>2</sub> O		c	217			
Na <sub>2</sub> Si <sub>2</sub> O <sub>5</sub>		c			213	213
Na <sub>4</sub> SiO <sub>4</sub>		c			213	213
Na <sub>2</sub> SiF <sub>6</sub>		c	218, 219, 220			
		aq	34, 219			
NaHSiF <sub>6</sub>		aq	34			
NaSn		c	172, 174, 221, 222			
NaSn <sub>2</sub>		c	172, 174, 221, 222			
Na <sub>2</sub> Sn		c	172, 174, 221, 222			
Na <sub>4</sub> Sn		c	172, 174, 221, 222			
Na <sub>4</sub> Sn <sub>3</sub>		c	172, 174, 221, 222			
Na <sub>2</sub> SnO <sub>3</sub>		c	223			
Na <sub>4</sub> SnO <sub>4</sub>		aq	34			
NaPb		c	173, 174			
Na <sub>2</sub> Pb		c	173, 174			
Na <sub>2</sub> Pb <sub>5</sub>		c	173, 174			
Na <sub>4</sub> Pb		c	173, 174			
Na <sub>5</sub> Pb <sub>2</sub>		c	173, 174			
Na <sub>2</sub> PbO <sub>3</sub>		c	224			
2NaI·PbI <sub>2</sub>		c	225			
2NaI·PbI <sub>2</sub> ·nH <sub>2</sub> O		c	225			
2Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> ·PbS <sub>2</sub> O <sub>3</sub>		c	226			
		aq	226			
Na <sub>2</sub> ZnO <sub>2</sub>		c	227			
Na <sub>2</sub> Zn(SO <sub>4</sub> ) <sub>2</sub>		c	228			
Na <sub>2</sub> Zn(SO <sub>4</sub> ) <sub>2</sub> ·4H <sub>2</sub> O		c	228			
NaCd <sub>2</sub>		c	173, 174, 229, 230			
NaCd <sub>6</sub>		c	173, 174, 229, 230			
Na <sub>2</sub> CdO <sub>2</sub>		c	231			
NaHg		c	174, 222, 172, 232, 233, 234			
NaHg <sub>2</sub>		c	174, 222, 172, 232, 233, 234			
NaHg <sub>4</sub>		c	172, 222, 174, 232, 233, 234			
NaHg <sub>27.5</sub>		liq	233, 234, 235			
NaHg <sub>50</sub>		liq	233, 234, 235			

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Formula	Description	State				
NaHg <sub>200</sub>		liq	233, 234, 235			
NaHg <sub>500</sub>		liq	233, 234, 235			
Na <sub>3</sub> Hg		c	174, 222, 172, 232, 233, 234			
Na <sub>3</sub> Hg <sub>2</sub>		c	174, 222, 172, 232, 233, 234			
Na <sub>5</sub> Hg <sub>2</sub>		c	174, 222, 172, 232, 233, 234			
Na <sub>7</sub> Hg <sub>8</sub>		c	174, 222, 172, 232, 233, 234			
Na <sub>2</sub> HgBr <sub>4</sub>		aq	236			
Na <sub>2</sub> HgS <sub>2</sub>		c	237			
NaHg(CN) <sub>3</sub>		aq	239			
Na <sub>2</sub> Hg(CN) <sub>4</sub>		aq	239			
NaCl·Hg(CN) <sub>2</sub>		aq	239			
NaCl·Hg(CN) <sub>2</sub> · 1 1/4H <sub>2</sub> O		c	239			
2NaCl·Hg(CN) <sub>2</sub>		aq	239			
NaBr·Hg(CN) <sub>2</sub>		aq	239			
NaBr·Hg(CN) <sub>2</sub> ·2H <sub>2</sub> O		c	239			
2NaBr·Hg(CN) <sub>2</sub>		aq	239			
NaI·Hg(CN) <sub>2</sub>		aq	239			
NaI·Hg(CN) <sub>2</sub> ·2H <sub>2</sub> O		c	239			
2NaI·Hg(CN) <sub>2</sub>		aq	239			
Na <sub>2</sub> O·CuO <sub>2</sub>		c	240			
Na <sub>2</sub> CO <sub>3</sub> ·CuCO <sub>3</sub>		c	241			
Na <sub>2</sub> CO <sub>3</sub> ·CuCO <sub>3</sub> ·3H <sub>2</sub> O		c	241			
Na <sub>2</sub> O·Ag <sub>2</sub> O <sub>2</sub>		c	240			
Na <sub>3</sub> Ag(S <sub>2</sub> O <sub>3</sub> ) <sub>2</sub>		aq	226			
NaAg(CN) <sub>2</sub>		aq	242			
Na <sub>2</sub> Ag(CN) <sub>3</sub>		aq	242			
Na <sub>2</sub> PtCl <sub>4</sub>		aq	34, 245			
Na <sub>2</sub> PtCl <sub>6</sub>		c	34, 243, 244			
		aq	243, 244			
Na <sub>2</sub> PtCl <sub>6</sub> ·nH <sub>2</sub> O		c	34			
Na <sub>2</sub> PtBr <sub>6</sub>		c	34			
		aq	34			
Na <sub>2</sub> PtBr <sub>6</sub> ·6H <sub>2</sub> O		c	34			
Na <sub>2</sub> PtI <sub>6</sub>		aq	246			
Na <sub>2</sub> IrCl <sub>6</sub>		c	247			
Na <sub>3</sub> IrCl <sub>6</sub>		c	247			
Na <sub>2</sub> OsCl <sub>6</sub>		c	247			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
NaReO <sub>4</sub>		c	248			
		aq	248, 1			
Na <sub>3</sub> RhCl <sub>6</sub>		c	243			
		aq	243, 247			
Na <sub>3</sub> RhCl <sub>6</sub> · 12H <sub>2</sub> O		c	243			
Na <sub>2</sub> Ni(CN) <sub>4</sub>		aq	242			
Na <sub>2</sub> CoO <sub>3</sub>		c	227			
Na <sub>2</sub> Fe <sub>2</sub> O <sub>4</sub>		c	249, 250, 251			
Na <sub>3</sub> FeCO(CN) <sub>5</sub>		c	252, 256			
		aq	252, 256			
Na <sub>3</sub> FeCO(CN) <sub>5</sub> · 7H <sub>2</sub> O		c	252, 256			
Na <sub>2</sub> MnO <sub>4</sub>		c	227			
Na <sub>2</sub> SO <sub>4</sub> · MnSO <sub>4</sub>		c	253			
Na <sub>2</sub> SO <sub>4</sub> · MnSO <sub>4</sub> · 2H <sub>2</sub> O		c	253			
Na <sub>2</sub> CrO <sub>4</sub>		c	255, 257, 258			
		aq	34, 255			
Na <sub>2</sub> CrO <sub>4</sub> · 4H <sub>2</sub> O		c	257, 258			
Na <sub>2</sub> Cr <sub>2</sub> O <sub>7</sub>		aq	34, 254			
Na <sub>2</sub> MoO <sub>4</sub>		c	261			
		aq	259, 260			
Na <sub>2</sub> WO <sub>4</sub>		c	255			
		aq	260			
NaVO <sub>5</sub>		aq	262, 263			
Na <sub>3</sub> VO <sub>4</sub>		c	264			
Na <sub>3</sub> V <sub>3</sub> O <sub>9</sub>		aq	262			
Na <sub>2</sub> TiO <sub>3</sub>		c			265	265
Na <sub>2</sub> Ti <sub>2</sub> O <sub>5</sub>		c			265	265
Na <sub>2</sub> Ti <sub>3</sub> O <sub>7</sub>		c			265	265
NaBO <sub>2</sub>		c	255			10
		aq	34, 117, 267			
NaBO <sub>3</sub>		aq	268			
NaBO <sub>3</sub> · 4H <sub>2</sub> O		c	268			
Na <sub>2</sub> B <sub>4</sub> O <sub>7</sub>		c	269			10
		aq	34, 117, 270			
Na <sub>2</sub> B <sub>4</sub> O <sub>7</sub> · nH <sub>2</sub> O		c	269			10
NaBH <sub>4</sub>	sodium borohydride	c	272		271	271
NaAlO <sub>2</sub>		c	255, 273			
Na <sub>3</sub> AlF <sub>6</sub>		c	274			10
Na <sub>3</sub> AlF <sub>6</sub> · 3 1/2H <sub>2</sub> O		c	274			
NaCl · AlCl <sub>3</sub>		c	274			
3NaCl · nAlCl <sub>3</sub>		c	274			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$\text{NaCl} \cdot \text{AlCl}_3 \cdot 6\text{NH}_3$		c	274			
$\text{Na}_2\text{O} \cdot \text{Al}_2\text{O}_3 \cdot 3\text{SiO}_2$	natrolite	c	275			
$\text{Na}_2\text{O} \cdot \text{Al}_2\text{O}_3 \cdot 4\text{SiO}_2$	dehydrated analcite	c	275			
$\text{Na}_2\text{UO}_4$		c	264			
$\text{Na}_2\text{U}_2\text{O}_7 \cdot 1 \frac{1}{2}\text{H}_2\text{O}$		c	260			
$(\text{Na}_2\text{O}_2)_2\text{UO}_4$		aq	260			
$(\text{Na}_2\text{O}_2)_2 \cdot \text{UO}_4 \cdot 9\text{H}_2\text{O}$		c	260			
$2\text{NaCl} \cdot \text{ThCl}_4$		c	276			
$2\text{NaCl} \cdot \text{ThCl}_4 \cdot 10\text{H}_2\text{O}$		c	276			
$\text{Na}_2\text{SO}_4 \cdot \text{CaSO}_4$		c	277			
$2\text{Na}_2\text{SO}_4 \cdot \text{CaSO}_4 \cdot 2\text{H}_2\text{O}$		c	277			
$\text{NaSrPO}_4 \cdot 9\text{H}_2\text{O}$		c	278, 279, 41			
$\text{NaSrAsO}_4 \cdot 9\text{H}_2\text{O}$		c	279			
$\text{NaBaPO}_4 \cdot 9\text{H}_2\text{O}$		c	278, 279, 41			
$\text{NaBaAsO}_4 \cdot 9\text{H}_2\text{O}$		c	279			
$\text{NaLiICl}$		c	121			

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| 6. Rodebush and Walters-1  | 20. de Forcrand-26                |
| 7. Rodebush and Henry-1    | 21. Blumenthal-1                  |
| 8. Loomis and Nusbaum-4    | 22. Centnerszwer and Blumenthal-1 |
| 9. Kelley-23               | 23. Gaydon-1                      |
| 10. Kelley-24              | 24. Herzberg-9                    |
| 11. Kelley-9               | 25. Dufay-1                       |
| 12. Moore-1                | 26. Hori-6                        |
| 13. Smith and Taylor-1     | 27. Sollers and Crenshaw-2        |
| 14. Gilles et al-1         | 28. Keyes-1                       |

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32. Berthelot-19	73. Gillespie, Lambert, and Gibson-1
33. de Forcrand-36	74. Moss and Wolfenden-1
34. Thomsen-16	75. Berthelot-155
35. Joannis-2	76. Neumann and Muller-2
36. Rengade-1	77. Fontana and Latimer-1
37. Fricke-1	78. Berthelot-9
38. Bertetti and McCabe-1	79. Bronsted-11
39. Rossini-4	80. Berthelot-79
40. Sturtevant-7	81. Harned and Owen-1
41. Berthelot-102	82. Wallace-1
42. Robinson and Stokes-2	83. Lange and Robinson-1
43. Akerlof and Kegeles-1	84. van Deventer and van de Stadt-1
44. Seidell-1	85. Dunnington and Hoggard-1
45. Pickering-13	86. Dingemans-2
46. Joyner-1	87. Bell-1
47. Schmitz and Schumacher-4	88. Berthelot-46
48. Guntz-1	89. Pickering-9
49. Ivett and De Vries-1	90. Varet-1
50. Roth-2	91. Dingemans-1
51. de Forcrand-61	92. Courtois-1
52. Nernst-8	93. Sabatier-1
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56. Allmand and Pollock-1	97. Berthelot-13
57. Randall and Bisson-1	98. Zeumer and Roth-3
58. Sandonnini-20	99. de Forcrand-4
59. Wüst and Lange-2	100. Kelley and Moore-4
60. Lippett, Johnson, and Maass-2	101. Ramstetter and Hantke-1
61. Cohen and Kooy-1	102. Tarasenkoy-1
62. Slansky-1	103. Arie-1
63. Lange and Martin-3	104. Ruff and Friedrich-1
64. Becker and Roth-7	105. Pitzer and Coulter-1
65. Richardson and Wells-1	106. Berthelot-8
66. Askew, Bullock, Smith, Tinkler, Gatty, and Wolfenden-1	107. Tilden-1
67. Voskresenskaya and Ponomareva-1	108. Berthelot and Ilosvay-1
68. Pitzer-3	109. Mixer-6
69. Kegeles-1	110. Voskresenskaya and Ponomareva-2
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| 115. Lange and Streeck-2       | 157. Juza, Fasold, and Haeberle-1    |
| 116. Richards and Rowe-4       | 158. de Forcrand-25                  |
| 117. Berthelot-4               | 159. Tilden-3                        |
| 118. Mathews and Germann-4     | 160. Biltz and Hansen-1              |
| 119. Recoura-2                 | 161. Joannis-7                       |
| 120. Berthelot-7               | 162. Hart and Partington-1           |
| 121. Beketoff-8                | 163. Matignon and Meyer-1            |
| 122. Higuchi-1                 | 164. Giran-1                         |
| 123. Tilden-3                  | 165. Mixter-9                        |
| 124. Favre and Valson-3        | 166. Joly-5                          |
| 125. Richards and Yngve-1      | 167. Berthelot and Louguinine-3      |
| 126. Stokes-4                  | 168. Perreu-16                       |
| 127. Berthelot-93              | 169. Amat-1                          |
| 128. Fogh-1                    | 170. Pfaundler-2                     |
| 129. Muller-15                 | 171. Perreu-17                       |
| 130. Berthelot-104             | 172. Kubaschewski and Seith-1        |
| 131. Bichowsky-3               | 173. Seith and Kubaschewski-1        |
| 132. Zimmermann and Latimer-1  | 174. Weibke and Kubaschewski-1       |
| 133. Teichert and Klemm-1      | 175. Berthelot-100                   |
| 134. Lindner-1                 | 176. de Forcrand-24                  |
| 135. Berthelot-6               | 177. Matignon-20                     |
| 136. Favre-8                   | 178. Torgeson-1                      |
| 137. Ephraim and Kornblum-1    | 179. Matsui, Kambara, and Miyamura-1 |
| 138. de Forcrand and Taboury-1 | 180. Swallow and Alty-1              |
| 139. Fabre-1                   | 181. Ostwald-1                       |
| 140. Mixter-10                 | 182. Pickering-8                     |
| 141. Metzner-1                 | 183. Donnan and Hope-1               |
| 142. Metzner-2                 | 184. Mixter-5                        |
| 143. Swietoslawski-6           | 185. Anderson-7                      |
| 144. Matignon and Marchal-3    | 186. Muller-1                        |
| 145. Dode-1                    | 187. Laksonen-1                      |
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| 147. Winkelmann-1              | 189. Perreu-1                        |
| 148. Scholz-1                  | 190. Cherbov and Chernyak-1          |
| 149. von Stackelberg-1         | 191. Berthelot-7                     |
| 150. Varali-Thevenet-1         | 192. Roth and Eymann-1               |
| 151. Mondain-Monval-3          | 193. Takagi and Oomi-1               |
| 152. Haigh-1                   | 194. Perreu-13                       |
| 153. Southard and Nelson-1     | 195. de Forcrand-17                  |

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203. Lloyd, Brown, Bunnel, and Jones-1	245. Miller and Terrey-1
204. Joannis-1	246. Pigeon-1
205. Lemoult-1	247. Puche-1
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208. Partington and Soper-1	250. Matsui and Kingo-1
209. Ruff and Greiger-1	251. Terashkevich and Vishnevskii-1
210. Matignon-14	252. Muller-9
211. Mulert-1	253. Graham-2
212. Roth and Richter-1	254. Morges-1
213. Kelley-10	255. Mixter-6
214. Mixter-5	256. Muller-8
215. Tschernabaeff-1	257. Berthelot-52
216. Tammann-3	258. Morgan and Benson-1
217. Lange and von Stockelburg-1	259. Pechard-1
218. Caillot-1	260. Pissarjewsky-3
219. Hantke-1	261. Mixter-10
220. Truchot-3	262. Matignon-15
221. Biltz and Holverscheid-1	263. Ruff and Friedrich-1
222. Biltz and Meyer-2	264. Mixter-14
223. Mixter-7	265. Shomate-7
224. Mixter-8	266. Berthelot-12
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Formula	Description	State				
K		g	1	1		1
		c			2	2
$K^{n+}$		g	3			
$K^+$		aq	6, 7	4, 5, 8, 9, 10		
$K_2$		g	11, 12, 13, 14		15	
$K_2O$		c	7, 16, 17			
$K_2O_2$		c	18			
$K_2O_3$		c	18, 19			
$K_2O_4$		c	18, 20			
KH		g	21		2	
		c	22, 23, 24			
KOH		c	25, 26, 27, 28			
		aq	29, 30		31	
$KOH \cdot nH_2O$		c	25, 32			
KF		c	33, 34, 35, 36, 37		37	37
		aq	33, 35, 38, 39			
$KF \cdot nH_2O$		c	33, 35, 38, 40		31	
$KHF_2$		c	37	37		37
		aq	35			
$KF \cdot nHF$		c	41			
KCl		g	14		31, 42	31, 42
		c	27, 43, 44, 45, 46, 47, 48, 49		50, 51	2
		aq	28, 52, 53, 54, 55, 56, 57		31	
KClO		aq	58, 59, 60			
KClO <sub>3</sub>		c	61, 62, 64		63	63
		aq	26, 28, 65, 66	9		
KClO <sub>4</sub>		c	62, 68		67	67
		aq	65, 66, 67, 69, 70	9		
KBr		g			31, 42	31, 42
		c	28, 34, 71, 72, 73, 74, 75		71	2
		aq	52, 53, 73, 76		31	
	in CH <sub>3</sub> OH		43, 77			
KBr <sub>3</sub>		aq	31, 78			
KBr <sub>5</sub>		aq	31			
KBrO		aq	28, 79			
KBrO <sub>3</sub>		c	28, 34, 66, 80		80	80
		aq	28, 31		9	
KI		g			42	42
		c	28, 73, 81, 82, 83, 84, 85		71	71

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Formula	Description	State				
KI		aq	53, 55, 73		31	
	in CH <sub>3</sub> OH		43, 77, 85			
	in C <sub>2</sub> H <sub>5</sub> OH		77, 85			
	in CH <sub>3</sub> COCH <sub>3</sub>		85			
	in CH <sub>3</sub> CN		85			
KI <sub>3</sub>		c	86			
		aq	78, 86, 87, 88			
KIO <sub>3</sub>		c	28, 31, 66, 89		80	80
		aq	28, 89, 90, 91		31	
KIO <sub>4</sub>		aq	28			
KIO <sub>3</sub> ·HIO <sub>3</sub>		c	89			
		aq	28			
KH <sub>4</sub> IO <sub>6</sub>		aq	28			
K <sub>2</sub> H <sub>3</sub> IO <sub>6</sub>		aq	28			
KI·KCl	fresh melt	c	92			
K <sub>2</sub> S		c	93, 94, 95			
		aq	95			
K <sub>2</sub> S·nH <sub>2</sub> O		c	95			
K <sub>2</sub> S <sub>4</sub>		c	95			
		aq	95			
K <sub>2</sub> S <sub>4</sub> ·nH <sub>2</sub> O		c	95			
K <sub>2</sub> SO <sub>3</sub>		c	96, 97			
		aq	96, 98			
K <sub>2</sub> SO <sub>4</sub>		c	28, 100, 101, 102, 107, 110	103, 104	99	99
		aq	28, 102, 105, 106, 108			
K <sub>2</sub> S <sub>2</sub> O <sub>3</sub>		aq	31			
K <sub>2</sub> S <sub>2</sub> O <sub>5</sub>		c	109			
		aq	109			
K <sub>2</sub> S <sub>2</sub> O <sub>5</sub> ·1 1/2H <sub>2</sub> O		c	109			
K <sub>2</sub> S <sub>2</sub> O <sub>6</sub>		c	28, 97			
		aq	28, 111			
K <sub>2</sub> S <sub>2</sub> O <sub>8</sub>		c	112			
		aq	31			
K <sub>2</sub> S <sub>4</sub> O <sub>6</sub>		c	28, 97			
		aq	31			
K <sub>2</sub> S <sub>5</sub> O <sub>6</sub>		c				
		aq	113			
K <sub>2</sub> S <sub>5</sub> O <sub>6</sub> ·1 1/2H <sub>2</sub> O		c	114			
KHS		c	95, 115			
		aq	95			

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Substance			$\Delta H_f^\circ$	$\Delta P_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
KHS•1/4H <sub>2</sub> O		c	95			
KHSO <sub>3</sub>		aq	116			
KHSO <sub>4</sub>		c	26, 28, 117			
		aq	28, 118			
KI•4SO <sub>2</sub>		c	119			
K <sub>2</sub> Se		c	120			
		aq	120			
K <sub>2</sub> Se•nH <sub>2</sub> O		c	120			
K <sub>2</sub> SeO <sub>4</sub>		aq	121, 122			
KHSe		c	115			
		aq	120			
KHSeO <sub>4</sub>		aq	121, 122			
K <sub>2</sub> TeO <sub>3</sub>		aq	121			
K <sub>2</sub> TeO <sub>4</sub>		aq	121			
KNO <sub>2</sub>		c	123			
		aq	124			
KNO <sub>3</sub>		c	29, 45, 125, 126, 127, 128, 129, 130, 131		132	132
		aq	28, 45, 133, 134		31	31
KNH <sub>2</sub>		c	135			
KBr•4NH <sub>3</sub>		c	136			
KI•nNH <sub>3</sub>		c	136			
KH <sub>2</sub> PO <sub>4</sub>		c	137, 138, 139			
		aq	137, 138			
KH <sub>2</sub> AsO <sub>4</sub>		c	139		140	140
		aq	31			
K <sub>2</sub> CO <sub>3</sub>		c	26, 28, 81, 141			
		aq	134, 142			
K <sub>2</sub> CO <sub>3</sub> •nH <sub>2</sub> O		c	26, 28			
K <sub>2</sub> C <sub>2</sub> O <sub>4</sub>	potassium oxalate	c	26			
		aq	26, 106, 143			
K <sub>2</sub> C <sub>2</sub> O <sub>4</sub> •H <sub>2</sub> O		c	26, 28, 138			
KCHO <sub>2</sub>	potassium formate	c	26			
		aq	28, 144			
KHCO <sub>3</sub>		c	26, 145, 146			
		aq	134, 142			
KOCH <sub>3</sub>	in CH <sub>3</sub> OH		147			
KC <sub>2</sub> H <sub>3</sub> O <sub>2</sub>	potassium acetate	c	26, 28			
		aq	26, 28, 143			
KC <sub>2</sub> H <sub>3</sub> O <sub>3</sub>	potassium glycollate	aq	148			
KOC <sub>2</sub> H <sub>5</sub>	in C <sub>2</sub> H <sub>5</sub> OH		147, 149			

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Formula	Description	State				
$5(K_2CO_3 \cdot 1 \frac{1}{2}H_2O) \cdot 4KHCO_3$	potassium oxamate	c	145	166, 167		
KCN		c	26, 28, 34			
		aq	150			
KCNO		c	26			
		aq	151			
$KCO_2CONH_2$		aq	152			
KCNS		c	153, 154, 155			
		aq	153, 154, 155			
$KCNS \cdot nSO_2$		c	119			
$K_2SiF_6$		c	156, 157, 158			
$K_2SnCl_4 \cdot H_2O$		c	28			
$K_2SnCl_6$		c	28			
		aq	28			
$KCl \cdot PbCl_2 \cdot 1/3H_2O$		c	159			
$KCl \cdot 2PbCl_2$		c	159			
$2KI \cdot PbI_2$		c	160			
$2KI \cdot PbI_2 \cdot 2H_2O$		c	160			
$4KI \cdot 3PbI_2$		c	160			
$4KI \cdot 3PbI_2 \cdot 6H_2O$		c	160			
$K_2SO_4 \cdot PbSO_4$		c	161, 162			
$K_2SO_4 \cdot ZnSO_4$		c	28			
$K_2SO_4 \cdot ZnSO_4 \cdot nH_2O$		c	28, 163			
$2KCN \cdot Zn(CN)_2$		c	164			
KHg		c	165			
KHg <sub>2</sub>		c	165			
KHg <sub>12</sub>		c	16 <sup>F</sup>			
KHg <sub>50</sub>		liq	165, 166, 168, 169			
KHg <sub>100</sub>		liq	165, 166, 168, 169			
KHg <sub>200</sub>		liq	165, 166, 168, 169			
KHg <sub>1000</sub>		liq	165, 166, 168, 169			
$KHgCl_3$		c	170			
		aq	170			
$KHgCl_3 \cdot H_2O$		c	170			
$K_2HgCl_4$		c	170			
		aq	28, 170			
$K_2HgCl_4 \cdot H_2O$		c	28, 170			
$4KCl \cdot 3HgCl_2$		c	170			
$4KCl \cdot 3HgCl_2 \cdot 3H_2O$		c	170			
KHgBr <sub>3</sub>		c	170			
		aq	170, 171			

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Formula	Description	State				
KHgBr <sub>3</sub> ·H <sub>2</sub> O		c	170			
KHgI <sub>3</sub>		c	170			
KHgI <sub>3</sub> ·H <sub>2</sub> O		c	170			
K <sub>2</sub> HgI <sub>4</sub>		c	28			
		aq	28, 172			
KHg(CN) <sub>3</sub>		aq	170, 173			
K <sub>2</sub> Hg(CN) <sub>4</sub>		c	26			
		aq	170, 173			
KCN·2Hg(CN) <sub>2</sub>		aq	170			
KCl·Hg(CN) <sub>2</sub>		c	172			
KCl·Hg(CN) <sub>2</sub> ·H <sub>2</sub> O		c	172			
KBr·Hg(CN) <sub>2</sub>		c	170			
		aq	170, 173			
KBr·Hg(CN) <sub>2</sub> ·1 1/2H <sub>2</sub> O		c	170, 173			
KI·Hg(CN) <sub>2</sub>		c	170			
		aq	170			
KI·Hg(CN) <sub>2</sub> ·1/4H <sub>2</sub> O		c	170			
KCl·CuCl <sub>2</sub>		c	174			
2KCl·CuCl		c	159			
2KCl·CuCl <sub>2</sub>		c	174			
2KCl·CuCl <sub>2</sub> ·2H <sub>2</sub> O		c	175			
K <sub>2</sub> Cu(SO <sub>4</sub> ) <sub>2</sub>		c	28, 176			
K <sub>2</sub> Cu(SO <sub>4</sub> ) <sub>2</sub> ·H <sub>2</sub> O		c	28			
K <sub>2</sub> Cu(SO <sub>4</sub> ) <sub>2</sub> ·6H <sub>2</sub> O		c	28, 177, 178			
K <sub>2</sub> Cu(CO <sub>3</sub> ) <sub>2</sub>		c	179			
KCl·AgCl		c	180			
KBr·AgBr		c	180			
3KBr·AgBr·1/2H <sub>2</sub> O		c	180			
KI·AgI		c	180			
KI·AgI·1/4H <sub>2</sub> O		c	180			
2KI·AgI·1/2H <sub>2</sub> O		c	180			
3KI·AgI		c	180			
3KI·AgI·1/2H <sub>2</sub> O		c	180			
3KI·2AgI·H <sub>2</sub> O		c	172			
KAg(CN) <sub>2</sub>		c	26, 181			
		aq	164, 171			
K <sub>2</sub> Ag(CN) <sub>2</sub>		aq	171			
KAu(CN) <sub>2</sub>		aq	182			
K <sub>2</sub> PtCl <sub>4</sub>		c	28			
		aq	28			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$K_2PtCl_6$		c	28, 183			
		aq	182, 183		184	184
$K_2PtBr_4$		c	184			
		aq	28			
$K_2PtBr_6$		c	28			
		aq	28			
$K_2IrCl_6$		c	183			
		aq	183			
$K_3IrCl_6$		c	183			
		aq	183			
$K_2OsCl_6$		c	185			
$KReO_4$		c	186			
		aq	186			
$K_2PdCl_4$		c	28, 187			
		aq	28, 187			
$K_2PdCl_6$		c	28, 185			
		aq	31			
$K_2PdBr_4$		c	187			
		aq	187			
$K_3RhCl_6$		c	185			
$K_2Ni(CN)_4$		aq	171			
$K_2CO_3 \cdot CoCO_3$		c	188			
$K_2CO_3 \cdot CoCO_3 \cdot 4H_2O$		c	188			
$KFe(SO_4)_2$		aq	28			
$K_2Fe(SO_4)_2$		aq	189			
$K_2Fe(SO_4)_2 \cdot nH_2O$		c	163, 190			
$K_3Fe(CN)_6$		c	154			
		aq	154			
$K_4Fe(CN)_6$		c	26, 191			191
		aq	154, 192, 193			
$K_4Fe(CN)_6 \cdot 3H_2O$		c	26, 191, 194			
$KH_2Fe(CN)_6$		aq	154			
$KH_3Fe(CN)_6$		aq	195			
$K_2HFe(CN)_6$		aq	154			
$K_2H_2Fe(CN)_6$		aq	154, 195			
$K_3HFe(CN)_6$		aq	195			
$K_3FeCO(CN)_5$		c	196			
		aq	196			
$K_3FeCO(CN)_5 \cdot 3 \frac{1}{2}H_2O$		c	196			
$KMnO_4$		c	197, 198		199	199
		aq	28, 197, 198			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$K_2Mn(SO_4)_2$		c	28			
$K_2Mn(SO_4)_2 \cdot nH_2O$		c	28			
$K_2CrO_4$		c	198, 200			
		aq	198, 200, 201, 202, 203			
$K_2Cr_2O_7$		c	28, 198, 204			
		aq	198, 202, 204, 205			
$K_2Cr_2O_7 \cdot CrO_3$		c	139			
$KCl \cdot CrO_3$		c	200			
$KCr(SO_4)_2$		c	28, 203			
		aq	206			
$KCr(SO_4)_2 \cdot nH_2O$		c	28, 206			
$KNH_4CrO_4$		c	202			
		aq	202			
$K_2MoO_4$		aq	207			
$KVO_3$		aq	208			
$KVO_4$		c	208			
		aq	208			
$KVO_5$		aq	31			
$3KF \cdot AlF_3$		c	209			
		aq	209			
$3KF \cdot AlF_3 \cdot 3 \frac{1}{2}H_2O$		c	209			
$KCl \cdot AlCl_3$		c	209			
$3KCl \cdot AlCl_3$		c	209			
$3KCl \cdot 2AlCl_3$		c	209			
$KAl(SO_4)_2$		c	206, 210		211	211
		aq	28			
$KAl(SO_4)_2 \cdot nH_2O$		c	28, 206		211	211
$KCl \cdot AlCl_3 \cdot 6NH_3$		c	209			
$K_2O \cdot Al_2O_3 \cdot 4SiO_2$	leucite	c	212, 213			
		gls	213			
$K_2O \cdot Al_2O_3 \cdot 6SiO_2$		c	212, 213			
		gls	212, 213			
$2KCl \cdot UO_2Cl_2 \cdot 2H_2O$		c	214			
$KCl \cdot ThCl_4 \cdot 9H_2O$		c	215			
$2KCl \cdot ThCl_4$		c	215			
$KCl \cdot MgCl_2$		c	81			
$KCl \cdot MgCl_2 \cdot 6H_2O$	carnallite	c	81, 216			
$2KCl \cdot MgCl_2$	fresh melt	c	81			
$4KCl \cdot MgCl_2$	fresh melt	c	81			

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Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$K_2SO_4 \cdot MgSO_4$	fresh melt	c	217, 218			
		c	81, 217			
$K_2SO_4 \cdot MgSO_4 \cdot nH_2O$	langbeinite	c	28, 163, 190, 217, 218			
$K_2SO_4 \cdot 2MgSO_4$		c	216			
$KCl \cdot MgSO_4$	hainite	c	216			
$2KCl \cdot CaCl_2$	fresh melt	c	81			
$K_2SO_4 \cdot CaSO_4 \cdot H_2O$		c	219			
$K_2SO_4 \cdot 5CaSO_4 \cdot H_2O$		c	219			
$KCaFeCO(CN)_5 \cdot 5H_2O$		c	220			
$K_2SO_4 \cdot SrSO_4$		c	219			
$KI \cdot LiCl$	fresh melt	c	221			
$KNa$		g	222			
		liq	6, 223			
$KNa_2$		liq	6, 223			
$K_2Na$		liq	6, 223			
$K_3Na$		liq	6, 223			
$KCl \cdot NaCl$	fresh melt	c	223			
$KI \cdot NaCl$	fresh melt	c	223			
$KNa_2PO_4$		aq	224			
$3KCNS \cdot NaCNS$		c	155			
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2. Kelley-24			16. Rengade-3			
3. Moore-1			17. Beketoff-2			
4. Lewis and Randall-7			18. Centnerszwer and Blumenthal-1			
5. Armbruster and Crenshaw-1			19. de Forcrand-67			
6. Joannis-2			20. de Forcrand-66			
7. Rengade-1			21. Almy and Beiler-1			
8. Robinson and Stokes-2			22. Keyes-1			
9. Seidell-1			23. Sollers and Crenshaw-1			
10. Hill and Ricci-1			24. Moutier-1			
11. Lewis-6			25. de Forcrand-35			
12. Loomis-2			26. Berthelot-9			
13. Loomis and Nusbaum-3			27. Voskresenskaya and Ponomareva-1			
14. Gaydon-1			28. Thomsen-16			



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| 30. Roth and Zeumer-3                             | 72. Brönsted-5  |
| 31. Calculated                                    | 73. Wüst and Lange-2  |
| 32. Berthelot-19                                  | 74. Popov, Bundel, and Choller-1                              |
| 33. Lange and Eichler-1                           | 75. Ishikawa and Tachiki-1                                    |
| 34. Lange and Martin-3                            | 76. Hammerschmid and Robinson-1                               |
| 35. Guntz-1                                       | 77. Askew, Bullock, Smith, Tinkler, Gatty, and<br>Wolfenden-1 |
| 36. de Forcrand-13                                | 78. Pickering-9   |
| 37. Westrum Jr. and Pitzer-1                      | 79. Berthelot-45  |
| 38. de Forcrand-59                                | 80. Ahlberg and Latimer-1                                     |
| 39. Lange and Messner-2                           | 81. Berthelot and Ilsovay-1                                   |
| 40. Bell-1  | 82. Scholz-1  |
| 41. Guntz-4                                       | 83. Hieber and Mühlbauer-1                                    |
| 42. Herzberg-9                                    | 84. Ishikawa and Tachiki-2                                    |
| 43. Slansky-1                                     | 85. Walden-4  |
| 44. Fineman and Wallace-1                         | 86. Berthelot-63  |
| 45. Lange and Monheim-2                           | 87. von Wartenburg and Klinkott-1                             |
| 46. Wüst and Lange-2                              | 88. Bertram and Roth-1  |
| 47. Miscenko-1                                    | 89. Berthelot-46  |
| 48. Roth and Eymann-1                             | 90. Matsuyama-2   |
| 49. Popov, Chomyakov, Feodossiev, and Shirokich-1 | 91. Berthelot-42  |
| 50. Clusius-4                                     | 92. Beketoff-8  |
| 51. Southard and Nelson-1                         | 93. Rengade and Costeanu-1                                    |
| 52. Rossini-5                                     | 94. Favre and Silbermann-3                                    |
| 53. Richards and Rowe-3                           | 95. Sabatier-1  |
| 54. Berthelot-85                                  | 96. Berthelot-92  |
| 55. Berthelot-18                                  | 97. Martin and Metz-1   |
| 56. Muller-11                                     | 98. Roth and Zeumer-2   |
| 57. Gillespie, Lambert, and Gibson-1              | 99. Moore and Kelley-1  |
| 58. Ramstetter and Hantke-1                       | 100. Pickering-2  |
| 59. Neumann and Muller-2                          | 101. Cohen and Kooy-1   |
| 60. Berthelot-24                                  | 102. Mischenko and Pronima                                    |
| 61. Wöhler and Schuff-1                           | 103. Shibata, Oda, and Furukawa-1                             |
| 62. Hofmann and Marin-1                           | 104. Kelley-22  |
| 63. Latimer, Schutz, and Hicks Jr.-3              | 105. Lange and Streeck-2                                      |
| 64. Brönsted-11                                   | 106. Plake-1  |
| 65. Andauer and Lange-1                           | 107. Pickering-8  |
| 66. Stackelberg-1                                 | 108. Berthelot-4  |
| 67. Latimer and Ahlberg-1                         | 109. Berthelot-94   |
| 68. Berthelot and Vielle-4                        | 110. Pickering-6  |
| 69. Berthelot-79                                  | 111. Zeumer and Roth-1  |
| 70. Noyes and Sammet-1                            |   |

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113. Berthelot-105	155. Wrzesnewsky-1
114. Berthelot-104	156. Caillot-1
115. Teichert and Klemm-1	157. Truchot-3
116. Berthelot-92	15B. Hantke-1
117. Favre-B	159. Bronsted-11
118. Berthelot-6	160. Berthelot-2
119. Ephraim and Kornblum-1	161. Bronsted-10
120. Fabre-1	162. Barre-1
121. Metzner-1	163. Caven and Ferguson-1
122. Metzner-2	164. Berthelot-117
123. Dode-2	165. Berthelot-56
124. Dode-1	166. Armbuster and Crenshaw-1
125. Haigh-1	167. Bent and Gilfillan-1
126. Roth and Eymann-1	16B. Lewis and Keyes-1
127. Roth-26	169. Smith and Ball-1
128. Voskresenskaya and Ponomereva-1	170. Berthelot-83
129. Hieber and Feder-1	171. Varet-6
130. Nacken-3	172. Berthelot-86
131. Hieber and Muhlbauer-1	173. Varet-2
132. Southard and Nelson-1	174. Agostini-1
133. Rossini-5	175. Bouzat and Chauvenet-2
134. Berthelot-4	176. Pickering-4
135. Juza, Fasold, and Haeberle-1	177. Pickering-6
136. Biltz and Hansen-1	178. Caven and Ferguson-2
137. Chomyakov, Yavorovskaya, and Schirokich-1	179. Pickering-12
138. Perreu-16	180. Berthelot-88
139. Graham-2	1B1. Berthelot-89
140. Stephenson and Zettlemoyer-1	182. Biltz and Wien-1
141. Ostwald-1	183. Gire-1,2
142. Muller-1	184. Coulter, Pitzer, and Latimer-1
143. Berthelot and Guntz-1	185. Puche-1
144. Berthelot-7	1B6. Roth and Becker-4
145. de Forcrand-54	187. Joannis-5
146. Caven and Land-2	18B. de Carli-1
147. de Forcrand-17	1B9. Berthelot-24
14B. de Forcrand-3	190. Graham-2
149. Deventer and Reicher-2	191. Schottky-1
150. Berthelot-25, B5, 116	192. Muller-2
151. Berthelot-111	193. Chretien and Guinchant-1
152. Matignon-1	194. Sano-15
153. Partington and Soper-1	195. Muller-2

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197. Roth and Becker-4
198. Perreu-16
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200. Morges-1
201. Roth, Schwartz, and Buckner-1
202. Sabatier-5
203. Berthelot-96
204. Stackelberg-1
205. Berthelot-95
206. Kraus, Fricke, and Querengasser-1
207. Pechard-1
208. Pissarjewsky-5
209. Baud-1
210. Young-2
211. Shomate-2
212. Mulert-1
213. Tammann-3
214. Aloy-1
215. Chauvenet-2
216. Richardson and Wells-1
217. Pickering-4
218. Pickering-6
219. Barre-1
220. Lecocq-1
221. Beketoff-8
222. Loomis and Arvin-1
223. Kawakami-3
224. Berthelot and Louguinine-3

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Formula	Description	State				
Rb		g	2,3,4		1	1
		c			4,5	4,5
Rb <sup>+</sup>		aq	1	6,7	5	
Rb <sup>n+</sup>		g	9,10,11			
Rb <sub>2</sub>		g	12,13			
Rb <sub>2</sub> O		c	14,15,16			
Rb <sub>2</sub> O <sub>2</sub>		c	17,18			
Rb <sub>2</sub> O <sub>3</sub>		c	17,18			
Rb <sub>2</sub> O <sub>4</sub>		c	17,18			
RbH		g	13			
RbOH		c	19,20			
		aq	15,14,20,1	1		
RbOH·nH <sub>2</sub> O		c	19,20			
RbF		c	21,22		23	
		aq	1,24	1		
RbF·nH <sub>2</sub> O		c	21			
RbHF <sub>2</sub>		c	25			
		aq	1			
RbCl		c	22,26,27,28			23
		aq	7	1		
	in CH <sub>3</sub> OH		29			
RbClO <sub>3</sub>		c	31	1	5	5
		aq	1,30	1,30		
RbClO <sub>4</sub>		c	31		32	
		aq	1	1,30		
RbBr		c	27,34,22		33	33
		aq	1	30		
RbI		c	22,27,36	30,35	33	33
		aq	1	1		
Rb <sub>2</sub> S		c	37,38			
		aq	37			
Rb <sub>2</sub> SO <sub>4</sub>		c	26			
		aq	26,39	1		
RbHS		c	40			
		aq	1			
RbHSO <sub>4</sub>		c	41			
		aq	26			
RbI·4SO <sub>2</sub>		c	42,43			
RbHSe		c	40			
		aq	1			

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Formula	Description	State				
RbNO <sub>3</sub>	rubidium thiocyanate	c	44	1		23
		aq	1			
RbNH <sub>2</sub>		c	45			
RbBr·3NH <sub>3</sub>		c	46			
RbI·6NH <sub>3</sub>		c	46			
Rb <sub>2</sub> CO <sub>3</sub>		c	47,48			
		aq	48			
Rb <sub>2</sub> CO <sub>3</sub> ·nH <sub>2</sub> O		c	47,48			
RbHCO <sub>3</sub>		c	50,51			
		aq	1,49			
3Rb <sub>2</sub> CO <sub>3</sub> ·2RbHCO <sub>3</sub> · 4 1/2H <sub>2</sub> O		c	52			
RbCNS		c	42			
		aq	1			
RbCNS·1/2SO <sub>2</sub>		c	42			
Rb <sub>2</sub> SiF <sub>6</sub>		c	53			
2RbI·PbI <sub>2</sub>		c	36			
2RbI·PbI <sub>2</sub> ·4H <sub>2</sub> O		c	36			
Rb <sub>2</sub> CuCl <sub>4</sub>		c	54			
Rb <sub>2</sub> CuCl <sub>4</sub> ·2H <sub>2</sub> O		c	54			
Rb <sub>2</sub> IrCl <sub>6</sub>		c	55			
RbReO <sub>4</sub>		c	56			
		aq	1			
RbAl(SO <sub>4</sub> ) <sub>2</sub>		c	57			
RbAl(SO <sub>4</sub> ) <sub>2</sub> ·nH <sub>2</sub> O		c	57			
2RbCl·ThCl <sub>4</sub>		c	58			
2RbCl·ThCl <sub>4</sub> ·9H <sub>2</sub> O		c	58			
4RbCl·ThCl <sub>4</sub>		c	58			

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1. Calculated	11. Tombouliau-1
2. Hackspill-2	12. Tsi-Ze and San-Tsiang-1
3. Scott-1	13. Gaydon-1
4. Kelley-19	14. Beketoff-3
5. Kelley-24	15. Rengade-1
6. Lewis and Argo-1	16. Rengade-3
7. Harned and Owen-1	17. Centnerszwer and Blumenthal-1
8. Lautie-1	18. de Forcrand-67
9. Bacher and Goudsmit-1	19. de Forcrand-45
10. Laporte, Miller and Sawyer-1	20. de Forcrand-56

Table 94 (Continued)  
SPECIFIC REFERENCES

REFERENCES	
21. de Forcrand-60	
22. de Forcrand-58	
23. Kelley-25	
24. Lange and Monheim-2	
25. de Forcrand-61	
26. de Forcrand-46	
27. de Forcrand-13	
28. Zemczuzny and Rambach-1	
29. Slansky-1	
30. Seidell-1	
31. Pitzer-4	
32. Latimer, Pitzer, and Smith-1	
33. Clusius, Goldman, and Perlick-1	
34. Lange and Martin-3	
35. Robinson and Stokes-2	
36. Mosnier-1	
37. Rengade and Costeanu-2	
38. Rengade and Costeanu-1	
39. Lange and Streeck-2	
40. Teichert and Klemm-1	
41. de Forcrand-46	
42. Ephraim and Kornblum-1	
43. de Forcrand and Taboury-1	
44. Haigh-1	
45. Juza, Fasold, and Haeberle-1	
46. Biltz and Hanson-1	
47. de Forcrand-50	
48. de Forcrand-53	
49. de Forcrand-54	
50. Caven and Land-2	
51. Caven and Ferguson-1	
52. de Forcrand-55	
53. Caillot-1	
54. Bouzat and Chauvenet-2	
55. Puche-1	
56. Smith and Long-1	
57. Kraus, Fricke and Querengässer-1	
58. Chauvenet-2	

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Table 95  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
Cs		g	1	4	1	1
		c				
Cs <sup>n+</sup>		g	3		1	2
Cs <sup>+</sup>		aq	5			
Cs <sub>2</sub>		g	6, 7			
Cs <sub>2</sub> O		c	8, 9			
Cs <sub>2</sub> O <sub>2</sub>		c	10			
Cs <sub>2</sub> O <sub>3</sub>		c	10			
Cs <sub>2</sub> O <sub>4</sub>		c	10, 11			
CsH		g	6		12, 13	5
CsOH		c	14, 15	5		
		aq	8, 16		5	
CsOH·H <sub>2</sub> O		c	15			
CsF		c	17			
		aq	17		5	
CsF·nH <sub>2</sub> O		c	17			
CsHF <sub>2</sub>		c	18			
		aq	18			
CsCl		c	17, 19, 20, 21			
		aq	22, 23	5		
CsClO <sub>4</sub>		c	25	26, 27	24	24
		aq	25	26, 27		
CsBr		c	17, 21, 28	5		29
		aq	17, 21, 28	5		
CsI		c	17, 21, 30	5		29
		aq	17, 21	26		
Cs <sub>2</sub> S		c	5	5		
		aq	31			
Cs <sub>2</sub> SO <sub>4</sub>		c	20			
		aq	20, 32			
CsHS		c	5, 33			
		aq	5, 33			
CsHSO <sub>4</sub>		c	20			
		aq	20			
CsI·4SO <sub>2</sub>		c	34, 35			
CsHSe		c	33			
		aq	33			
CsNO <sub>3</sub>		c	19			
		aq	23			
CsNH <sub>2</sub>		c	36			

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Table 95 (Continued)  
SPECIFIC REFERENCES

Substance			$\Delta H_f^\circ$	$\Delta F_f^\circ$	$S^\circ$	$C_p^\circ$
Formula	Description	State				
$\text{Cs}_2\text{CO}_3$		c	37			
		aq	37,38			
$\text{Cs}_2\text{CO}_3 \cdot 3 \frac{1}{2}\text{H}_2\text{O}$		c	37,38			
$\text{CsHCO}_3$		c	39			
		aq	39,40,41			
$5\text{Cs}_2\text{CO}_3 \cdot 2\text{CsHCO}_3 \cdot \text{nH}_2\text{O}$		c	42			
$\text{Cs}_2\text{SiF}_6$		c	43			
$\text{Cs}_2\text{CuCl}_4$		c	44			
$\text{Cs}_2\text{CuCl}_4 \cdot 2\text{H}_2\text{O}$		c	44			
$\text{CsReO}_4$		c	45			
		aq	45			
$\text{CsAl}(\text{SO}_4)_2$		c	46			
$\text{CsAl}(\text{SO}_4)_2 \cdot \text{nH}_2\text{O}$		c	46	47		47
$2\text{CsCl} \cdot \text{ThCl}_4$		c	48			
$2\text{CsCl} \cdot \text{ThCl}_4 \cdot 8\text{H}_2\text{O}$		c	48			
$4\text{CsCl} \cdot \text{ThCl}_4$		c	48			
$\text{CsLiICl}$		c	30			
$\text{CsNaICl}$		c	30			
$\text{CsKICl}$		c	30			

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|-----------------------------------|----------------------------------|
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| 2. Kelley-24                      | 16. de Forcrand-56               |
| 3. Bacher and Goudsmit-1          | 17. de Forcrand-58               |
| 4. Bent, Forbes, and Forziati-1   | 18. de Forcrand-61               |
| 5. Calculated                     | 19. Haigh-1                      |
| 6. Gaydon-1                       | 20. de Forcrand-46               |
| 7. Loomis and Kusch-1             | 21. de Forcrand-13               |
| 8. Rengade-1                      | 22. Lange and Messner-2          |
| 9. Beketoff-6                     | 23. Richards and Rowe-2          |
| 10. Centnerszwer and Blumenthal-1 | 24. Pitzer, Smith, and Latimer-1 |
| 11. de Forcrand-57                | 25. Pitzer-4                     |
| 12. Herzberg-9                    | 26. Seidell-1                    |
| 13. Almy and Rassweiler-1         | 27. Latimer-1                    |
| 14. Beketoff-4                    | 28. Lange and Martin-3           |



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Table 95 (Continued)  
SPECIFIC REFERENCES

REFERENCES

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30. Beketoff-8
31. Rengade and Costeanu-1
32. Lange and Streeck-2
33. Teichert and Klemm-1
34. Ephraim and Kornblum-1
35. de Forcrand and Taboury-1
36. Juza, Fasold, and Haeberle-1
37. de Forcrand-50
38. de Forcrand-53
39. de Forcrand-54
40. Caven and Land-2
41. Caven and Ferguson-1
42. de Forcrand-55
43. Caillot-1
44. Bouzat and Chauvenet-2
45. Smith and Long-1
46. Kraus, Fricke, and Querengässer-1
47. Latimer and Greensfelder-1
48. Chauvenet-2



V. SPECIFIC REFERENCES FOR THE TABLES OF  
PROPERTIES OF SERIES II

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 1  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
O <sub>2</sub>		Transition	1,7,9	1	1
		Fusion	1,6,7,8,9	1	1
		Vaporization	2	1,3	1,5
O <sub>3</sub>		Vaporization	4	4	
REFERENCES					
1. Giauque and Johnston-1 2. Stimson-1 3. Frank and Clusius-2 4. Riesenfeld and Beja-1 5. Wagman, Kilpatrick, Taylor, Pitzer, and Rossini-1 6. Lisman and Keesom-1 7. Aoyama and Kanda-4 8. Henning and Otto-1 9. Clusius-6					

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Table 2 SPECIFIC REFERENCES					
Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$H_2$		Fusion	1	1	1
		Vaporization	1	1	
$H_2O$		Fusion	6	2,3	3,4,5
		Vaporization	6	4,7	4,7
$H_2O_2$		Fusion	8	8	
		Vaporization	8	8	

REFERENCES

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| 1. Simon and Lange-1<br>2. Osborne-1<br>3. Giauque and Stout-1<br>4. Osborne, Stimson, and Ginnings-2,3<br>5. Brown, Bornes, and Maass-1<br>6. Stimson-1 | 7. Wagman, Kilpatrick, Taylor, Pitzer, and Rossini-1<br>8. Maass and Hiebert-1 |
|--|--|

Table 2a

$1H_2$		Fusion	1	1	1
		Vaporization	1	1	

REFERENCES

1. Calculated

Table 2b

$2H_2$		Fusion	1,2	1,2	1,2
		Vaporization	1,2,3		
$1H_2H$		Fusion	1	4	4
		Vaporization	1	4	4
$2H_2O$		Fusion	5,6	5,6,7,8,9	5,7,8
		Vaporization	10	10	
$1H_2HO$		Vaporization	11	11	

REFERENCES

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|--|--|
| 1. Woolley, Scott, and Brickwedde-1<br>2. Clusius and Bartholomé-2<br>3. Lewis and Hanson-1<br>4. Calculated<br>5. Long and Kemp-1<br>6. Redlich and Zentner-1,2 | 7. Brown, Barnes and Maass-1<br>8. Bartholome and Clusius-1<br>9. Jacobs-2<br>10. Rossini, Knowlton, and Johnston-1<br>11. Wahl and Urey-1 |
|--|--|

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Table 3					
SPECIFIC REFERENCES					
Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
He		Fusion	1,2,3,4,5	1,2,3,4,5	1,6
		Transition	1,6	1,6	
		Vaporization	1	1	
REFERENCES					
1. Keesom-2		6. Keesom, Weber and Norgaard-1			
2. Keesom and Keesom-2					
3. Keesom, Weber, and Schmidt-1					
4. Kaishev and Simon-1					
5. Simon and Steckel-1					
Table 4					
Ne		Fusion	1,2	1,2	1,3,4
		Vaporization	3	3	1,3,4
REFERENCES					
1. Clusius-8					
2. Henning and Otto-1					
3. Keesom and Haantjes-3					
4. Clusius-5					
Table 5					
A		Fusion	1	1	1,3,4
		Vaporization	2	2	1,3,4
REFERENCES					
1. Clusius-8					
2. Frank and Clusius-2					
3. Eucken and Hauch-1					
4. Rice-1					

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Table 6  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Kr		Fusion	1	1	1,2
		Vaporization	2	2	1,2

REFERENCES

1. Clusius-8
2. Clusius, Kruis, and Konnertz-1

Table 7

Xe		Fusion	1	1	1
		Vaporization	1,2	1,2	1

REFERENCES

1. Clusius and Riccoboni-1
2. Clusius and Wiegand-1

Table 8

Rn		Fusion	1	2	
		Vaporization	1	2	

REFERENCES

1. Washburn et. al.-1
2. Unpublished Calculations

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Table 9  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
F <sub>2</sub>		Fusion	1	1, 2, 7	1, 5
		Vaporization	3, 4	3, 4, 7	1, 5, 6
F <sub>2</sub> O		Vaporization	8	8, 18	
HF		Fusion	9	9	9, 13, 14, 15
		Vaporization	10, 11, 12, 17, 18	10, 11, 16, 17	9, 14, 15, 13
REFERENCES					
1. Kanda-1 2. Aoyama and Kanda-5 3. Claussen-1 4. Cady and Hildebrand-1 5. Kelley-18 6. Murphy and Vance-1 7. Henglein-1 8. Ruff and Menzel-4 9. Dahmlos and Jung-1 10. Fredenhagen-2 11. Simons and Bouknight-1 12. Claussen and Hildebrand-1 13. Clusius, Hiller, and Vaughen-1 14. Pranschke and Schwiete-1		15. Roth, Pahlke, Bertram and Borger-1 16. Kolosovskii-1 17. Simons-1 18. Kelley-19			



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Table 10  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Cl <sub>2</sub>  ClO <sub>2</sub> Cl <sub>2</sub> O Cl <sub>2</sub> O <sub>7</sub> HCl  HCl·2H <sub>2</sub> O ClF ClF <sub>3</sub> ClO <sub>2</sub> F		Fusion	1, 2, 5, 6, 7, 8, 9	1, 2, 5, 6, 7, 8, 9	1, 4
		Vaporization	1, 2, 3, 32	1, 2, 3, 32	1, 4, 12, 13
		Vaporization	3, 10	3, 10	
		Vaporization	3	3	
		Vaporization	11	11	
		Transition	14	14	14
		Fusion	7, 14, 26	7, 14, 26	14
		Vaporization	6, 14, 16, 17, 18, 19, 20, 21, 22, 23, 24, 25, 32	6, 14, 16, 17, 18, 19, 20, 21, 22, 23, 24, 25, 32	14
		Fusion	15, 28, 27	15, 28, 27	
		Vaporization	29, 30	29, 30	
		Vaporization	30	30	
		Vaporization	31	31	

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|---------------------------------------|--------------------------------|
| 1. Giauque and Powell-1               | 20. Ansdell-1                  |
| 2. Kelley-24                          | 21. McIntosh and Steele-1      |
| 3. Goodeve-1                          | 22. Briner-4                   |
| 4. Brucksch, Jr, and Ziegler          | 23. Drozdowski and Pietrgale-1 |
| 5. Estreicher and Staniewski-1        | 24. Cardoso and Germann-1      |
| 6. Estreicher and Schneer-1           | 25. Henning and Stock-1        |
| 7. Eucken and Karwat-1                | 26. Eucken and Donath-1        |
| 8. Harteck-2                          | 27. Roozeboom-4                |
| 9. Rideal-1                           | 28. Roozeboom-3                |
| 10. King and Partington-1             | 29. Ruff and Laass-1           |
| 11. Goodeve and Powney-1              | 30. Ruff and Krug-1            |
| 12. Kelley-18                         | 31. Schmitz and Schumacher-1   |
| 13. Kelley-25                         | 32. Kelley-19                  |
| 14. Giauque and Wiebe-1               |                                |
| 15. Berthelot-47                      |                                |
| 16. Elliott and McIntosh-1            |                                |
| 17. Steele, McIntosh, and Archibald-1 |                                |
| 18. Tsurita-1                         |                                |
| 19. Faraday-1                         |                                |

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Table 11  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Br <sub>2</sub>		Fusion	1,2,3,4,5,6,7,8,10	1,2,3,4,5,6,7,8,10,11	9,11,12
		Vaporization	13,14,15,17,18,19,20,21,22,23,24,25,3,26,27,28,29	13,14,16,17,18,19,20,21,22,23,24,25,3,26,27,28,29	
HBr		Fusion	30,31,32,33,34,35	30,31,32,33,34,35	11,30
		Vaporization	30,31,33,34,36,37,38	30,31,33,34,36,37,38	
HBr•H <sub>2</sub> O		Fusion	39	39	
BrF <sub>5</sub>		Fusion	10,40	10,40	
		Vaporization	11,40	11,40	

REFERENCES

1. Regnault-2	21. Regnault-7
2. Broune-1	22. Thomsen-16
3. Isnardi-1	23. Scheffer and Voogd-1
4. Ramsay and Young-1	24. Bouzat and Leluan-1
5. Cuthbertson and Cuthbertson-1	25. Cuthbertson and Cuthbertson-1
6. Henglein-2	26. Ramsay and Young-1
7. Henglein, Rosenberg, and Muchlinski-1	27. Roozeboom-3
8. Rideal-1	28. Roozeboom-4
9. Kelley-18	29. Lewis and Randall-7
10. Kelley-21	30. Giaque and Wiebe-2
11. Unpublished Calculations	31. Eucken and Karwat-1
12. Latimer and Hoenshel-1	32. Henglein-3
13. Smits and Cannegieter-1	33. Maass and McIntosh-1
14. Smits and Cannegieter-2	34. Steele, McIntosh, and Archibald-1
15. Thorpe-1	35. Drozdowski and Pietrzak-1
16. Fleischman-1	36. Elliott and McIntosh-1
17. Andrews-5	37. Estreicher and Schnee-1
18. Andrews-9	38. Steele and Bagster-1
19. Berthelot and Ogier-6	39. Roozeboom-6
20. LeChatelier-4	40. Ruff and Menzel-2

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Table 12  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$I_2$		Fusion	1,2,3,4,5,6,7	1,2,3,4,5,6,7	16
		Sublimation	8,9,10,11,12,13,14,15	8,9,10,11,12,13,14,15	
HI		Fusion	16,17	16,17	
		Vaporization	16	16	
$IF_5$		Fusion	2,18	2,18	
		Vaporization	2,18	2,18	
$IF_7$		Sublimation	2,19	2,19	
ICl		Fusion	20,21,22,23,24	20,21,22,23,24	
		Vaporization	20,25,26,27	20,25,26,27	

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|------------------------------------|-------------------------|
| 1. Frederick and Hildebrand-1      | 20. Cornog and Bauer-1  |
| 2. Unpublished Calculations        | 21. Tanatar-6           |
| 3. Ramsay and Young-1              | 22. Stortenbeker-1      |
| 4. Stelzner and Niederschulte-1    | 23. Berthelot-63        |
| 5. Rassow-1                        | 24. Berthelot-144       |
| 6. Rideal-1                        | 25. Thorpe-1            |
| 7. Kelley-21                       | 26. McMorris and Yost-2 |
| 8. Giauque-3                       | 27. Cornog and Karges-1 |
| 9. Baxter and Grose-1              |                         |
| 10. Baxter, Hickey and Holmes-1    |                         |
| 11. Ramsay and Young-1             |                         |
| 12. Haber and Kerschbaum-1         |                         |
| 13. Wohl-3                         |                         |
| 14. Dewar-1                        |                         |
| 15. Nernst-1                       |                         |
| 16. Giauque and Wiebe-3            |                         |
| 17. Bates, Halford, and Anderson-1 |                         |
| 18. Ruff and Braida-3              |                         |
| 19. Ruff and Keim-1                |                         |

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Table 14  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
S		Transition	1,2,3,4,5,6,7,8,9,10,11	1,2,3,4,5,6,7,8,9,10,11	1,2,3,12
		Sublimation	3,13,14,15	3,13,14,15	
		Fusion	1,2,3,9,11,12,13,16,17,18	1,2,3,9,11,12,13,16,17,18	1,2,3,12,13
		Transition		8,10,19	
		Vaporization	3,20,21,22,23,24,25,26,27,28,29,30,31,32,33,34	3,20,21,22,23,24,25,26,27,28,29,30,31,32,33,34	
SO <sub>2</sub>		Fusion	21,35,36,37,38,39,40	21,35,36,37,38,39,40	2,3,35,36,41,42,43
		Vaporization	35,36,42,43,44,45	35,36,42,43,44,45	35,43
SO <sub>3</sub>		Fusion	3,37,38,39,40,46,47,48	3,37,38,39,40,46,47,48	
		Sublimation	3,37,38,39,40,46,47,48	3,37,38,39,40,46,47,48	
		Vaporization	3,46,47,48,49,50,51,52,53,54,55,56,57,58,59,60,61,62,63,64,65,66,67,68,69,70,71	3,46,47,48,49,50,51,52,53,54,55,56,57,58,59,60,61,62,63,64,65,66,67,68,69,70,71	
S <sub>2</sub> O <sub>7</sub>		Fusion	72		
H <sub>2</sub> S		Transition	73,74,75,76	73,74,75,76	2,73,74,75
		Fusion	73,74,75	73,74,75	2,73,74,75
		Vaporization	40,51,73,74,77,78,79,80,81,82,83	40,51,73,74,77,78,79,80,81,82,83	2,73,74,77
H <sub>2</sub> S <sub>2</sub>		Fusion	84,85	3,84	3,84
		Vaporization	84,85	3,84	3,84
H <sub>2</sub> S <sub>3</sub>		Fusion	86,87		
		Vaporization	87		
H <sub>2</sub> SO <sub>4</sub>		Fusion	88,89,90,91,92,93	88,89,90,91,92,93	88
H <sub>2</sub> SO <sub>4</sub> ·H <sub>2</sub> O		Fusion	88,89,90,91,92,93,94	88,89,90,91,92,93,94	88
H <sub>2</sub> SO <sub>4</sub> ·2H <sub>2</sub> O		Fusion	90		
H <sub>2</sub> SO <sub>4</sub> ·4H <sub>2</sub> O		Fusion	94		
H <sub>2</sub> SO <sub>5</sub>		Fusion	95		
H <sub>2</sub> S <sub>2</sub> O <sub>7</sub>		Fusion	96	96	96
H <sub>2</sub> S <sub>2</sub> O <sub>8</sub>		Fusion	95		
SF <sub>4</sub>		Fusion	97		
		Vaporization	97	97	
SF <sub>6</sub>		Transition	98	98	98
		Sublimation	99,100	99,100	
		Fusion	98,99,100,101	98,99,100,101	98
		Vaporization	99,102	99,102	

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Table 14 (Continued)  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
S <sub>2</sub> F <sub>2</sub>		Fusion	103		
		Vaporization	103		
S <sub>2</sub> F <sub>10</sub>		Fusion	104		
		Vaporization	104	104	
SOF <sub>2</sub>		Fusion	105	105	
		Vaporization	105		
SO <sub>2</sub> F <sub>2</sub>		Fusion	106,107		
		Vaporization	106,107		
HSO <sub>3</sub> F		Vaporization	108		
S <sub>2</sub> Cl <sub>2</sub>		Fusion	109		
		Vaporization	110,111,112,113	110,111,112,113	
SOCl <sub>2</sub>		Fusion	114		
		Vaporization	112,115,116	112,115,116	
SO <sub>2</sub> Cl <sub>2</sub>		Fusion	117		
		Vaporization	117	117	
S <sub>2</sub> O <sub>5</sub> Cl <sub>2</sub>		Fusion	118		
		Vaporization	112,115,116,119	112,115,116,119	
SOFCl		Fusion	120		
		Vaporization	120	120	
SO <sub>2</sub> FCl		Fusion	121		
		Vaporization	121	121	
S <sub>2</sub> Br <sub>2</sub>		Fusion	122		
		Vaporization	122		
SOBr <sub>2</sub>		Fusion	114		
		Vaporization	114	114	

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Table 14 (Continued)  
SPECIFIC REFERENCES

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39. Burrell and Robertson-1	81. Regnault-6
40. Steele and Bagster-1	82. Maass and McIntosh-2
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50. Eucken-7	92. Berthelot-6
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Table 14 (Continued)  
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Table 14a  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta G_p$
Formula	Description				
$^2\text{H}_2\text{S}$  $^1\text{H}^2\text{H}_3$		Transition	1	1	1
		Fusion	1	1	1
		Transition	2	2	2
		Fusion	2	2	2

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Table 15  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Se		Transition	1	1	1
		Fusion	1,2	1	1
		Vaporization	3,4	3,4	
SeO <sub>2</sub>		Sublimation	5,6,7	5,6,7	
H <sub>2</sub> Se		Transition	8	8	8
		Fusion	8	8	8
		Vaporization	8,9,10	9,10	
H <sub>2</sub> SeO <sub>4</sub>		Fusion	11,12	11	
H <sub>2</sub> SeO <sub>4</sub> ·H <sub>2</sub> O		Fusion	11,12	11	
H <sub>2</sub> SeO <sub>4</sub> ·4H <sub>2</sub> O		Fusion	12		
SeF <sub>4</sub>		Fusion	13		
		Vaporization	13		
		Sublimation	14,15	14,15	
SeF <sub>6</sub>		Fusion	14,15	16	
		Vaporization	14,15	14,15	
		Fusion	17		
SeOF <sub>2</sub>		Vaporization	17		
		Sublimation	18		
SeCl <sub>4</sub>		Fusion	18		
		Fusion	19	19	
SeOCl <sub>2</sub>		Vaporization	19	19	
		Fusion	20		
Se(OH) <sub>3</sub> ·ClO <sub>4</sub>		Fusion	21		
SeOBr <sub>2</sub>		Fusion	21		
		Vaporization	21		

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Table 15a  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$2H_2Se$		Transition	1	1	1
		Fusion	1	1	1
$1H2HSe$		Transition	1,2	1,2	1,2
		Fusion	2	2	2
REFERENCES					
1. Kruis and Clusius-1 2. Kruis-1					

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Table 16  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Te		Transition	1,2	1,2	1,2,3,4
		Fusion	1,2,3,4,5,6	1,2,3,4	
		Vaporization	7,8,9,10,11	7,8,9,10	
H <sub>2</sub> Te		Fusion	12,13	12	
		Vaporization	12,13,14	12,14	
TeF <sub>6</sub>		Transition	15	15	
		Sublimation	15,16	15,16	
		Fusion	15,16	15,16	
		Vaporization	16	16	
TeCl <sub>4</sub>		Fusion	17,18	17,18	17
		Vaporization	18	18	

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Table 18  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
N <sub>2</sub>		Transition	1	1	
		Fusion	1,2,3,4,5,6,7,8	1	
		Vaporization	1	1	
NO		Fusion	9	9	9
		Vaporization	9,10	9,10	
N <sub>2</sub> O		Fusion	11,12	11,12	11,13
		Vaporization	11,12	11,12	
N <sub>2</sub> O <sub>3</sub>		Fusion	14		
		Vaporization	14,15	14,15	
N <sub>2</sub> O <sub>4</sub>		Fusion	16	16	16
		Vaporization	16	16	
N <sub>2</sub> O <sub>5</sub>		Sublimation	17	17	
NH <sub>3</sub>		Fusion	18	15,18	
		Vaporization	18,19	15,18,19	
NH <sub>3</sub> · 1/2 H <sub>2</sub> O		Fusion	20		
NH <sub>3</sub> · H <sub>2</sub> O		Fusion	20		
N <sub>2</sub> H <sub>4</sub>		Fusion	21,22,23		
		Vaporization	21,22	22	
N <sub>2</sub> H <sub>4</sub> · H <sub>2</sub> O		Fusion	22		
		Vaporization	22		
HN <sub>3</sub>		Fusion	24		
		Vaporization	24,25,26	26	
NH <sub>4</sub> N <sub>3</sub>		Sublimation	27	27	
		Fusion	28		
N <sub>2</sub> H <sub>4</sub> HN <sub>3</sub>		Fusion	28		
HNO <sub>3</sub>		Fusion	29	29	29
		Vaporization	30	30	
HNO <sub>3</sub> · H <sub>2</sub> O		Fusion	29	29	29
HNO <sub>3</sub> · 3H <sub>2</sub> O		Fusion	29	29	29
NH <sub>2</sub> OH		Fusion	31		
		Vaporization	31		
NH <sub>4</sub> NO <sub>3</sub>		Transition	32,33,34,35	32,33,34,35	
		Fusion	32,33,36	32,33,36	
N <sub>2</sub> H <sub>4</sub> · HNO <sub>2</sub>		Fusion	37		
N <sub>2</sub> H <sub>4</sub> · HNO <sub>3</sub>	metastable	Fusion	38		
	stable	Fusion	38,39,40		
N <sub>2</sub> H <sub>4</sub> · 2HNO <sub>3</sub>		Fusion	39		
NH <sub>4</sub> NO <sub>3</sub> · 2HNO <sub>3</sub>		Fusion	41		
NF <sub>3</sub>		Transition	42		
		Fusion	42		
		Vaporization	43	43	

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Table 18 (Continued)  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
NOF		Fusion	44		
		Vaporization	44	44	
NO <sub>2</sub> F		Fusion	44		
		Vaporization	44	44	
NO <sub>3</sub> F		Fusion	45		
		Vaporization	45		
NH <sub>2</sub> F		Sublimation	46		
NHF <sub>2</sub>		Fusion	46		
		Vaporization	46		
NH <sub>4</sub> F•HF		Fusion	47		
N <sub>2</sub> H <sub>4</sub> •2HF		Fusion	48		
NCI <sub>3</sub>		Vaporization	49		
NOC1		Fusion	50		
		Vaporization	50	50	
NO <sub>2</sub> Cl		Fusion	51		
		Vaporization	51	51	
NH <sub>2</sub> Cl		Fusion	52		
NH <sub>4</sub> Cl		Transition	53, 54, 55, 56, 57, 58, 59, 60	53, 54, 55, 56, 57, 58, 59, 60	
		Fusion	61		
NH <sub>4</sub> Cl•3NH <sub>3</sub>		Fusion	62		
N <sub>2</sub> H <sub>4</sub> •HCl		Fusion	63	63	
N <sub>2</sub> H <sub>4</sub> •2HCl		Fusion	64		
NH <sub>4</sub> ClO <sub>4</sub>		Transition	65		
N <sub>3</sub> 8r		Fusion	66		
NO8r		Fusion	67		
NO8r <sub>3</sub>		Fusion	67		
NH <sub>4</sub> 8r		Transition	59	59	
		Fusion	61		
NH <sub>4</sub> 8r•3NH <sub>3</sub>		Fusion	62		
N <sub>2</sub> H <sub>4</sub> •H8r		Fusion	48		
N <sub>2</sub> H <sub>4</sub> •2H8r		Fusion	48		
NH <sub>4</sub> I		Transition	55, 59, 68	55, 59, 68	
		Fusion	61		
NH <sub>4</sub> I•3NH <sub>3</sub>		Fusion	62		
NH <sub>4</sub> I•4NH <sub>3</sub>		Fusion	62		
N <sub>2</sub> H <sub>4</sub> HI		Fusion	48		
N <sub>2</sub> H <sub>4</sub> •2HI		Fusion	48		
NH <sub>4</sub> Cl <sub>2</sub> I		Fusion	69		
NH <sub>4</sub> Cl <sub>4</sub> I		Fusion	69		
NH <sub>4</sub> I8r <sub>2</sub>		Fusion	69		
N <sub>2</sub> S <sub>5</sub>		Fusion	70		

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Table 1B (Continued)  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$N_4S_4$		Fusion	71,72,73	76	
$N_2O_3 \cdot 2SO_3$		Fusion	74		
		Vaporization	74		
$NO_2SO_2OH$		Fusion	75		
$NH_4HSO_4$		Transition	59		
		Fusion	76		
$N_2H_4 \cdot H_2SO_4$		Fusion	64		
$(NH_4)_2SO_4$		Fusion	77		
$(H_2NSO_2)_2NH$		Fusion	78		
$(N_2H_4)_2H_2SO_4$		Fusion	37		
$NOSO_3F$		Fusion	79		

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Table 19  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta G_p$
Formula	Description				
P		Transition	1,2,3,4,5	1,2,3,4,5,6	
		Fusion	1,2,3,4,5	1,2,3,4,5	
		Vaporization	4,5	4,5	
P <sub>4</sub> O <sub>6</sub>		Fusion	7		
		Vaporization	7	7	
P <sub>4</sub> O <sub>10</sub>		Sublimation	8,9,10,11	8,9,10,11	
		Fusion	8,9,10,11,12	8,9,10,11,12	
		Vaporization	12	12	
PH <sub>3</sub>		Transition	13,14	13,14	
		Fusion	13,14	13,14	
		Vaporization	13,14	13,14	
P <sub>2</sub> H <sub>4</sub>		Vaporization	15		
H <sub>3</sub> PO <sub>2</sub>		Fusion	16,17	17	
H <sub>3</sub> PO <sub>3</sub>		Fusion	17	17	
H <sub>3</sub> PO <sub>4</sub>		Fusion	18,19	17	
H <sub>3</sub> PO <sub>4</sub> ·1/2 H <sub>2</sub> O		Fusion	18,19,20	20	
H <sub>4</sub> P <sub>2</sub> O <sub>6</sub>		Fusion	21		
H <sub>4</sub> P <sub>2</sub> O <sub>6</sub> ·2H <sub>2</sub> O		Fusion	21	21	
H <sub>4</sub> P <sub>2</sub> O <sub>7</sub>		Fusion	22	22	
H <sub>4</sub> P <sub>2</sub> O <sub>7</sub> ·1 1/2H <sub>2</sub> O		Fusion	22	22	
H <sub>6</sub> P <sub>4</sub> O <sub>13</sub>		Fusion	23		
PF <sub>3</sub>		Fusion	24,25,26		
		Vaporization	24,26	24,26	
PF <sub>5</sub>		Fusion	27	27	
		Sublimation	27	27	
		Vaporization	27	27	
POF <sub>3</sub>		Vaporization	25,26,28,29	26,29	
		Sublimation	26	26	
		Fusion	26,29	26,29	
HPO <sub>2</sub> F <sub>2</sub>		Vaporization	26		
PCl <sub>3</sub>		Fusion	30		
		Vaporization	30,31	31	
PCl <sub>5</sub>		Sublimation	32	32	
		Fusion	32		
P <sub>2</sub> Cl <sub>4</sub>		Fusion	33		
POCl <sub>3</sub>		Fusion	26,34,35,36	36	
		Vaporization	26,31,35,37,38,39	31,39	
PH <sub>4</sub> Cl		Transition	40		
		Fusion	40		
PF <sub>2</sub> Cl		Fusion	41		
		Vaporization	41	41	

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Table 19 (Continued)  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
PFCl <sub>2</sub>		Fusion	41		
		Vaporization	41	41	
POF <sub>2</sub> Cl		Fusion	29		
		Vaporization	29	29	
POFCl <sub>2</sub>		Fusion	29		
		Vaporization	29	29	
PBr <sub>3</sub>		Fusion	30		
		Vaporization	30,42,43	42	
PBr <sub>5</sub>		Sublimation	44	44	
POBr <sub>3</sub>		Fusion	45		
		Vaporization	38,46	46	
PF <sub>2</sub> Br		Fusion	47		
		Vaporization	47	47	
PFBr <sub>2</sub>		Fusion	47		
		Vaporization	47	47	
PF <sub>3</sub> Br <sub>2</sub>		Fusion	48		
		Fusion	49		
POF <sub>2</sub> Br		Vaporization	49	49	
		Fusion	49		
POFBr <sub>2</sub>		Vaporization	49	49	
		Fusion	49		
POCl <sub>2</sub> Br		Fusion	34		
POClBr <sub>2</sub>		Fusion	34		
POClFBr		Vaporization	50		
PI <sub>3</sub>		Fusion	38,51,52		
P <sub>2</sub> I <sub>4</sub>		Fusion	51		
P <sub>2</sub> S <sub>3</sub>		Fusion	6		
P <sub>4</sub> S <sub>3</sub>		Fusion	53		
		Vaporization	53		
P <sub>4</sub> S <sub>7</sub>		Fusion	54,55		
P <sub>4</sub> S <sub>10</sub>		Fusion	56,57		
P <sub>4</sub> O <sub>6</sub> S <sub>4</sub>		Fusion	58		
		Vaporization	58		
PSF <sub>3</sub>		Fusion	25,59		
		Vaporization	59		
PSCl <sub>3</sub>		Fusion	59		
		Vaporization	34,37		
PSF <sub>2</sub> Cl		Fusion	59		
		Vaporization	59	59	
PSFCl <sub>2</sub>		Fusion	59		
		Vaporization	59	59	
PSBr <sub>3</sub>		Fusion	60		



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Table 19 (Continued)  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
PSF <sub>2</sub> Br		Fusion	60	61	
		Vaporization	61		
PSFBr <sub>2</sub>		Fusion	61	61	
		Vaporization	61		
PSCl <sub>2</sub> Br		Fusion	34		
		Vaporization	34		
PSClBr <sub>2</sub>		Fusion	34		
		Vaporization	34		
NH <sub>2</sub> OH·H <sub>3</sub> PO <sub>2</sub>		Fusion	62		
NH <sub>4</sub> H <sub>2</sub> PO <sub>3</sub>		Fusion	62, 66		
N <sub>2</sub> H <sub>4</sub> ·H <sub>3</sub> PO <sub>3</sub>		Fusion	62		
NH <sub>4</sub> NH <sub>2</sub> HPO <sub>3</sub>		Fusion	62		
N <sub>2</sub> H <sub>4</sub> ·H <sub>3</sub> PO <sub>4</sub>		Fusion	62		
N <sub>2</sub> H <sub>4</sub> ·H <sub>4</sub> P <sub>2</sub> O <sub>6</sub>		Fusion	62		
N <sub>2</sub> H <sub>4</sub> ·(H <sub>3</sub> PO <sub>3</sub> ) <sub>2</sub>		Fusion	62		
(NH <sub>4</sub> ) <sub>2</sub> H <sub>2</sub> P <sub>2</sub> O <sub>6</sub>		Fusion	62		
N <sub>3</sub> P <sub>3</sub> F <sub>8</sub> H <sub>2</sub> ·2H <sub>2</sub> O		Fusion	64		
NH <sub>4</sub> HPO <sub>3</sub> F		Fusion	65		
NH <sub>4</sub> PO <sub>2</sub> F <sub>2</sub>		Fusion	65		
(PNCI <sub>2</sub> ) <sub>3</sub>		Fusion	66, 67		
(PNCI <sub>2</sub> ) <sub>4</sub>		Vaporization	66, 67, 68	66, 68	
		Fusion	67		
(PNCI <sub>2</sub> ) <sub>5</sub>		Vaporization	67, 68	67, 68	
		Fusion	67		
(PNCI <sub>3</sub> ) <sub>6</sub>		Vaporization	67		
		Fusion	67		
N <sub>4</sub> P <sub>4</sub> Cl <sub>2</sub> F <sub>6</sub>		Vaporization	67		
		Fusion	67		
PNBr <sub>2</sub>		Vaporization	67		
		Fusion	69		

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Table 20  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
As		Sublimation	1,2,3,4,5,6,7	1,2,3,4,5,6,7	
		Fusion	2,8,9,10,11		
		Vaporization	1,2,3,4,5,6,7,9,10,11,12		
As <sub>4</sub> O <sub>6</sub>		Sublimation	13,14,15,16,17,18,19,20,21	13,14,15,16,17,18,19,20,21	
		Fusion	13,14,15,16,17,18,19,20,21	13,14,15,16,17,18,19,20,21	
		Vaporization	15,16,18,20,22	15,16,18,20,22	
AsH <sub>3</sub>		Fusion	23	23	
		Vaporization	23	23	
H <sub>3</sub> AsO <sub>4</sub> · 1/2 H <sub>2</sub> O		Fusion	24		
AsF <sub>3</sub>		Fusion	25	25	
		Vaporization	25,26	25	
AsF <sub>5</sub>		Fusion	27	27	
		Vaporization	27		
AsCl <sub>3</sub>		Fusion	28,29	28,29	
		Vaporization	26,29,30,31,32,33,34,35,36,37	29,30,32,33,34,35,36,37	
AsBr <sub>3</sub>		Fusion	30,36,37,38,39,40	30,38	
		Vaporization	30,37,39	30,41	
AsI <sub>2</sub>		Fusion	42		
AsI <sub>3</sub>		Fusion	30,43,44,45,46		
		Vaporization	46	46	
As <sub>2</sub> S <sub>2</sub>		Transition	47		
		Fusion	47,48		
		Vaporization	47		
As <sub>2</sub> S <sub>3</sub>		Transition	47		
		Fusion	47		
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Table 21  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Sb		Transition	1,2,3,4		
		Fusion	10	5,6,7,8,9,11	
		Vaporization	12,13		
Sb <sub>4</sub> O <sub>6</sub>		Transition	14	14	
		Fusion	14	14	
		Vaporization	14	14	
SbH <sub>3</sub>		Fusion	15		
		Vaporization	15		
SbF <sub>3</sub>		Fusion	16,17		
		Vaporization	18		
SbF <sub>5</sub>		Fusion	19		
		Vaporization	20,21		
2SbF <sub>3</sub> ·SbF <sub>5</sub>		Vaporization	20		
SbCl <sub>3</sub>		Fusion	22,23,24,25,26,27	22,26,27	
		Vaporization	20,23,25,28,29,30,31,32,33	30,31,32,33	
SbCl <sub>5</sub>		Fusion	24,34,35,36,37	34,35,37	
		Vaporization	30,34	30,34	
SbCl <sub>3</sub> F <sub>2</sub>		Fusion	38		
SbBr <sub>3</sub>		Fusion	22,23,24,25,26,27	22,25,26,27	
		Vaporization	23		
SbI <sub>3</sub>		Fusion	23,39,40		
Sb <sub>4</sub> S <sub>6</sub>		Fusion	41,42,43,44	42,44	
Sb <sub>2</sub> Se <sub>3</sub>		Fusion	45		
Sb <sub>2</sub> Te <sub>3</sub>		Fusion	46		
3NH <sub>4</sub> Cl·2SbCl <sub>3</sub>		Fusion	25		

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Table 22  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Bi		Fusion	1, 2, 3, 4, 5, 6, 7, 8, 9, 10, 11, 12, 13, 14, 15, 16, 17, 18, 19, 20, 21	1, 11, 13, 14, 15, 16, 17, 18, 19, 20, 21	
		Vaporization	22, 23, 24		
Bi <sub>2</sub> O <sub>3</sub>		Fusion	21, 25	21, 25	
BiF <sub>3</sub>		Fusion	26		
BiF <sub>5</sub>		Sublimation	26		
BiCl <sub>3</sub>		Fusion	7, 21, 27, 28, 29	21, 27, 28, 29	
		Vaporization	24	24	
BiCl <sub>4</sub>		Fusion	27		
BiBr <sub>3</sub>		Transition	30		
		Fusion	7, 27, 28, 29, 30, 31		
		Vaporization	32	32	
BiS		Fusion	33		
BiSe		Transition	34		
Bi <sub>2</sub> Se <sub>3</sub>		Fusion	10, 34		
Bi <sub>2</sub> Te <sub>3</sub>		Fusion	2, 4, 8, 12		
Bi <sub>2</sub> Te <sub>3</sub> ·Bi <sub>2</sub> S <sub>3</sub>		Fusion	8, 35		

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Table 23  
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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
C	graphite	Sublimation	1	1	
CO	carbon monoxide	Transition	2,3	2,3	2,3
		Fusion	2,3	2,3	2,3
		Vaporization	2,4	2,4	
CO <sub>2</sub>	carbon dioxide	Sublimation	4,5	4,5	
		Fusion	6	6	
CH <sub>4</sub>	methane	Fusion	7,130	7,130	
		Vaporization	7,130	7,130	
CH <sub>2</sub> O	formaldehyde	Fusion	8		
		Vaporization	8	8	
CH <sub>2</sub> O <sub>2</sub>	formic acid	Fusion	9,10	9,10	9
		Vaporization	10	10	
CH <sub>3</sub> OH	methanol	Transition	11	11	
		Fusion	11	11	11
		Vaporization	12,13,14	12,13,14	
CH <sub>4</sub> O <sub>2</sub>	methyl hydrogen peroxide	Vaporization	15	15	
CF <sub>4</sub>	tetrafluoromethane	Transition	16	16	
		Fusion	16,17	16,17	
		Vaporization	17	17	
COF <sub>2</sub>	carbonyl fluoride	Fusion	18		
		Vaporization	18	18	
CH <sub>3</sub> F	fluoromethane	Vaporization	19	19	
CH <sub>2</sub> F <sub>2</sub>	difluoromethane	Vaporization	20		
CHF <sub>3</sub>	trifluoromethane	Fusion	21		
		Vaporization	21,22	21,22	
CCl <sub>4</sub>	tetrachloromethane	Transition	23,24,57,26,27,28	23,24,57,27,28	
		Fusion	23,24,57,27,28,29,30,31,32	23,24,57,27,28,29,30,31,32	
		Vaporization	14,33,34	14,33,34	
COCl <sub>2</sub>	carbonyl chloride	Fusion	35	35	
		Vaporization	35	35	
CH <sub>3</sub> Cl	chloromethane	Fusion	36	36	36
		Vaporization	34,36	34,36	
CH <sub>2</sub> Cl <sub>2</sub>	dichloromethane	Fusion	37,38	37,38	
		Vaporization	14,39,40	14	
CHCl <sub>3</sub>	trichloromethane	Fusion	37,41,42,43	37,42	
		Vaporization	14,34,39,40,41,44	14,34	
CF <sub>3</sub> Cl	trifluorochloromethane	Fusion	45		
		Vaporization	46,47,48	46,47,48	



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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
CF <sub>2</sub> Cl <sub>2</sub>	difluorodichloromethane	Fusion	49,50		
		Vaporization	47,48,50,51	47,48,50	
CFCl <sub>3</sub>	fluorotrichloromethane	Fusion	52	52	
		Vaporization	40,51,52,53	40,51,52,53	52,54
COFCl	carbonylfluorochloride	Fusion	55		
		Vaporization	55	55	
CH <sub>2</sub> FCl	fluorochloromethane	Fusion			
		Vaporization	40,55	40,55	
CHF <sub>2</sub> Cl	difluorochloromethane	Fusion	25		
		Vaporization	56,25	56,25	
CHFC1 <sub>2</sub>	fluorodichloromethane	Fusion	25		
		Vaporization	56,25	56,25	
CBr <sub>4</sub>	tetrabromomethane	Transition	58,59	58,59	
		Fusion	59,60,61	59,60	
		Vaporization	62	62	
COBr <sub>2</sub>	carbonyl bromide	Vaporization	63	63	
CH <sub>3</sub> Br	bromomethane	Transition	64	64	
		Fusion	64	64	
		Vaporization	64	64	
CH <sub>2</sub> Br <sub>2</sub>	dibromomethane	Fusion	39,65,66,67	39,66,67	
		Vaporization	39,66,	39,66	
CHBr <sub>3</sub>	tribromomethane	Fusion	68,69,70	69,70	
		Vaporization	39,71,72,73	71,72,73	
CF <sub>3</sub> Br	trifluorobromomethane	Fusion			
		Vaporization	74		
CBr <sub>2</sub> F <sub>2</sub>	difluorodibromomethane	Fusion			
		Vaporization	75		
CBr <sub>3</sub> F	fluorotribromomethane	Fusion			
		Vaporization	75		
CHF <sub>2</sub> Br	difluorobromomethane	Fusion			
		Vaporization	76		
CHBr <sub>2</sub>	fluorodibromomethane	Fusion			
		Vaporization	77		
CI <sub>4</sub>	tetraiodomethane	Fusion	39		
CH <sub>3</sub> I	iodomethane	Fusion	78		
		Vaporization	39,66,79	39,66,79	
CH <sub>2</sub> I <sub>2</sub>	diiodomethane	Fusion	65,80	80	
		Vaporization	81		
CHI <sub>3</sub>	triiodomethane	Fusion	37	37	
CHF <sub>2</sub> I	difluoroiodomethane	Fusion	21		
		Vaporization	21	21	

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
CHFI <sub>2</sub>	fluorodiodomethane	Fusion	21		
		Vaporization	21	21	
CS <sub>2</sub>	carbon disulfide	Fusion	82,83,84	83	
		Vaporization	14,84,85	14,84,85	
COS	carbonoxysulfide	Fusion	86,87	86,87	86
		Vaporization	86	86	
CH <sub>4</sub> S	methanethiol	Transition	88	88	
		Fusion	88	88	88
		Vaporization	88	88	
H <sub>2</sub> CS <sub>3</sub>	trithiocarbonic acid	Fusion	89		
CSe <sub>2</sub>	carbon diselenide	Fusion	90		
		Vaporization	90	90	
COSe	carbonoxyselenide	Fusion	91		
		Vaporization	91	91	
CSSe	carbon sulfoselenide	Fusion	92		
		Vaporization	92	92	
CSTe	carbon sulfotelluride	Fusion	93		
C(NO <sub>2</sub> ) <sub>4</sub>	tetranitromethane	Fusion	25		
		Vaporization	94	94	
HCN	hydrogen cyanide	Transition	95	95	
		Fusion	95	95	95
		Vaporization	95	95	
CH <sub>5</sub> N	methylamine	Fusion	96	96	
		Vaporization	96	96	
CH <sub>2</sub> N <sub>2</sub>	cyanamide	Fusion	97	97	
HNCO	hydrogen isocyanide	Fusion	98		
		Vaporization	99	99	
CH <sub>3</sub> ON	formamide	Fusion	100,101,102	101	
CH <sub>3</sub> O <sub>2</sub> N	nitromethane	Fusion	103	103	
		Vaporization	14,103,104,105,106	103	
CH <sub>3</sub> O <sub>2</sub> N	methyl nitrite	Fusion			
		Vaporization	107	107	
CH <sub>3</sub> O <sub>3</sub> N	methyl nitrate	Fusion			
		Vaporization	107	107	
CH <sub>4</sub> ON <sub>2</sub>	urea	Fusion	108,109	108,109	
CFN	cyanogen fluoride	Sublimation	110	110	
COF <sub>3</sub> N	N,N-difluorofluoroformamide	Fusion	111		
		Vaporization	111	111	
COF <sub>3</sub> N	trifluoronitrosomethane	Fusion	111		
		Vaporization	111	111	

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
CClN	cyanogen chloride	Fusion	112, 113	112	
		Vaporization	42	112	
CO <sub>2</sub> Cl <sub>3</sub> N	trichloronitromethane	Fusion			
		Vaporization	114, 115	114, 115	
CH <sub>3</sub> ClN	methylamine hydrochloride	Transition	116	116	
CBrN	cyanogen bromide	Sublimation	114	114	
		Fusion	117, 118		
		Vaporization	118		
CIN	cyanogen iodide	Sublimation	119, 120	119, 120	
		Fusion	118		
HCNS	hydrogen thiocyanate	Fusion	121		
NH <sub>4</sub> SCN	ammonium thiocyanate	Transition	58	58	
COF <sub>2</sub> NP	phosphorus difluoroiso- cyanate	Fusion	122		
		Vaporization	122	122	
CF <sub>2</sub> SNP	phosphorus difluoroiso- thiocyanate	Vaporization	112	112	
CCl <sub>2</sub> SNP	phosphorus dichlorothio- cyanate	Fusion	113		
		Vaporization	113	113	
CH <sub>3</sub> As	methyl arsine	Fusion	124		
		Vaporization	124	124	
CH <sub>3</sub> F <sub>2</sub> As	methyl difluoroarsine	Fusion	125		
		Vaporization	125	125	
CH <sub>3</sub> Cl <sub>2</sub> As	methyl dichloroarsine	Fusion	126		
		Vaporization	126	126	
C <sub>2</sub> H <sub>2</sub>	ethyne (acetylene)	Sublimation	127, 130	127	
		Fusion	127	127	
		Vaporization	127	127	
C <sub>2</sub> H <sub>4</sub>	ethene (ethylene)	Fusion	128, 129, 130	128, 129, 130	
		Vaporization	7	7	
C <sub>2</sub> H <sub>6</sub>	ethane	Fusion	131, 130	131, 130	131
		Vaporization	130	130	130
C <sub>2</sub> H <sub>2</sub> O	ketene	Fusion	132		
		Vaporization	132		
C <sub>2</sub> H <sub>2</sub> O <sub>2</sub>	glyoxal	Fusion	133		
		Vaporization	133	133	
C <sub>2</sub> H <sub>2</sub> O <sub>4</sub>	oxalic acid	Sublimation	134	134	
C <sub>2</sub> H <sub>4</sub> O	acetaldehyde	Fusion	135, 137	135, 137	
		Vaporization	136, 138, 139	136, 138, 139	139
C <sub>2</sub> H <sub>4</sub> O·H <sub>2</sub> O	acetaldehyde monohydrate	Fusion	140	140	140
C <sub>2</sub> H <sub>4</sub> O	ethylene oxide	Fusion	141	141	141
		Vaporization	141, 142	141, 142	142

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$C_2H_4O_2$	acetic acid	Fusion	143, 144, 145, 146	144, 145	147
		Transition	144	144	
		Vaporization	143, 144	148, 149	
$C_2H_4O_2$	methyl formate	Fusion	137, 150	150	
		Vaporization	14, 137	14	
$C_2H_4O_2$	hydroxyacetaldehyde	Fusion	151		
$C_2H_4O_3$	hydroxyacetic acid (glycollic acid)	Fusion	152	152	
$C_2H_6O$	ethanol	Fusion	153	153	
		Vaporization	155, 156	155	
$C_2H_6O$	dimethyl ether	Fusion	157	157	157
		Vaporization	157	157	157
$C_2H_6O_2$	1,2-ethanediol (ethylene glycol)	Fusion	158		
		Vaporization	159, 160, 161, 162		
$C_2F_4$	tetrafluoroethane	Fusion	163		
		Vaporization	163	163	
$C_2F_6$	hexafluoroethane	Transition	164	164	164
		Fusion	163, 164	163, 164	164
		Vaporization	164	164	
$C_2OF_4$	trifluoroacetyl fluoride	Vaporization	165		
$C_2H_3F$	fluoroethene	Fusion	25		
		Vaporization	25	25	
$C_2H_5F$	fluoroethane	Vaporization	166, 167		
$C_2H_4F_2$	1,1-difluoroethane	Fusion	25		
		Vaporization	25	25	
$C_2HF_3$	trifluoroethene	Vaporization	168		
$C_2H_3F_3$	1,1,1-trifluoroethane	Transition	169		169
		Fusion	169	169	
		Vaporization	169	169	
$C_2H_3OF$	acetyl fluoride	Vaporization	170, 171	170	
$C_2H_3O_2F$	fluoroacetic acid	Fusion	172		
$C_2H_5OF$	2-fluoroethanol	Fusion	172, 173		
		Vaporization	172, 173, 174	174	
$C_2H_2O_2F_2$	difluoroacetic acid	Fusion	172		
		Vaporization	172, 175		
$C_2H_4OF_2$	2,2-difluoroethanol	Fusion	172, 176, 177		
		Vaporization	176, 177		
$C_2HO_2F_3$	trifluoroacetic acid	Fusion	165		
		Vaporization	165		
$C_2H_3OF_3$	2,2,2-trifluoroethanol	Fusion	178		
		Vaporization	178		

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$C_2Cl_2$	dichloroethyne	Vaporization	179		
$C_2Cl_4$	tetrachloroethene	Fusion	37,137,143	37	
		Vaporization	14,25,137,143,180	14,25,180	
$C_2Cl_6$	hexachloroethane	Transition	181,182	181	
		Sublimation	181	181	
$C_2O_2Cl_4$	trichloromethyl chloroformate	Vaporization	183	183	
$C_2H_3Cl$	chloroethene	Fusion	184		
		Vaporization	184	184	
$C_2H_5Cl$	chloroethane	Fusion	185,186	185,186	185
		Vaporization	34,185,187	34,185,187	
$C_2H_2Cl_2$	1,1-dichloroethene	Fusion	25		
		Vaporization	25	25	
$C_2H_2Cl_2$	cis-1,2-dichloroethene	Fusion	25		
		Vaporization	188	188	
$C_2H_2Cl_2$	trans-1,2-dichloroethene	Fusion	25		
		Vaporization	188	188	
$C_2H_4Cl_2$	1,1-dichloroethane	Fusion	25,37	25,37	
		Vaporization	25,189,190	25,189,190	
$C_2H_4Cl_2$	1,2-dichloroethane	Fusion	191,192,193	191,192	192
		Vaporization	14,193,194	14,194	
$C_2HCl_3$	trichloroethene	Fusion	180,137		
		Vaporization	14,180,195	14,180,195	
$C_2H_3Cl_3$	1,1,1-trichloroethane	Transition	196	197	
		Fusion	37,196	37,196	
		Vaporization	25,37,190,196	196	
$C_2H_3Cl_3$	1,1,2-trichloroethane	Fusion	25,137		
		Vaporization	25	25	
$C_2H_2Cl_4$	1,1,1,2-tetrachloroethane	Fusion	25		
		Vaporization	25,190	25,190	
$C_2H_2Cl_4$	1,1,2,2-tetrachloroethane	Fusion	180,197		
		Vaporization	14,180,190,198	14,180,190,198	
$C_2HCl_5$	pentachloroethane	Fusion	37,137	37	
		Vaporization	180,190,193,198	180,190,198	
$C_2H_3OCl$	acetyl chloride	Fusion	199		
		Vaporization	200	200	
$C_2H_3OCl$	chloroacetaldehyde	Vaporization	201		
$C_2H_3O_2Cl$	chloroacetic acid	Fusion	202,203,204	202,203,204	202
		Vaporization	205	205	
$C_2H_5OCl$	2-chloroethanol	Fusion	193,197		
		Vaporization	14,193,197	14	

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
C <sub>2</sub> H <sub>5</sub> OC1	chloromethyl, methyl ether	Vaporization	206,207		
C <sub>2</sub> H <sub>2</sub> OC1 <sub>2</sub>	chloroacetyl chloride	Vaporization	208		
C <sub>2</sub> H <sub>2</sub> OC1 <sub>2</sub>	dichloroacetaldehyde	Vaporization	209		
C <sub>2</sub> H <sub>2</sub> O <sub>2</sub> C1 <sub>2</sub>	dichloroacetic acid	Fusion	203	203	203
		Vaporization	205	205	
C <sub>2</sub> H <sub>4</sub> O <sub>2</sub> C1 <sub>2</sub>	dichloroacetaldehyde monohydrate	Fusion	210		
C <sub>2</sub> HOCl <sub>3</sub>	trichloroacetaldehyde (chloral)	Fusion	25	25	
		Vaporization	25	25	
C <sub>2</sub> HO <sub>2</sub> C1 <sub>3</sub>	trichloroacetic acid	Fusion	203,211	203,211	203
		Vaporization	25	25	
C <sub>2</sub> H <sub>3</sub> OC1 <sub>3</sub>	2,2,2-trichloroethanol	Fusion	212		
		Vaporization	212		
C <sub>2</sub> H <sub>3</sub> O <sub>2</sub> C1 <sub>3</sub>	trichloroacetaldehyde monohydrate	Fusion	213,214	214	214
C <sub>2</sub> F <sub>3</sub> Cl	trifluorochloroethene	Fusion	215		
		Vaporization	25,215	25	
C <sub>2</sub> F <sub>5</sub> Cl	pentafluorochloroethane	Vaporization	216		
C <sub>2</sub> F <sub>2</sub> Cl <sub>2</sub>	1,1-difluoro-2,2-dichloroethene	Fusion	217		
		Vaporization	217		
C <sub>2</sub> F <sub>2</sub> Cl <sub>2</sub>	cis-1,2-difluoro-1,2-dichloroethene	Fusion	216		
		Vaporization	216		
C <sub>2</sub> F <sub>2</sub> Cl <sub>2</sub>	trans-1,2-difluoro-1,2-dichloroethene	Fusion	216		
		Vaporization	216		
C <sub>2</sub> F <sub>4</sub> Cl <sub>2</sub>	1,1,1,2-tetrafluoro-2,2-dichloroethane	Vaporization	216		
C <sub>2</sub> F <sub>4</sub> Cl <sub>2</sub>	1,1,2,2-tetrafluoro-1,2-dichloroethane	Fusion	25		
		Vaporization	25,218,219	25,218,219	
C <sub>2</sub> FC1 <sub>3</sub>	fluorotrichloroethene	Fusion	215		
		Vaporization	215		
C <sub>2</sub> F <sub>3</sub> Cl <sub>3</sub>	1,1,2-trifluoro-1,2,2-trichloroethane	Fusion	25		
		Vaporization	25,220,221	25,220,221	221
C <sub>2</sub> F <sub>3</sub> Cl <sub>3</sub>	1,1,1-trifluoro-2,2,2-trichloroethane	Fusion	217		
		Vaporization	217		
C <sub>2</sub> F <sub>2</sub> Cl <sub>4</sub>	1,2-difluoro-1,1,2,2-tetrachloroethane	Fusion	25,220		
		Vaporization	25,220	25,220	
C <sub>2</sub> F <sub>2</sub> Cl <sub>4</sub>	1,1-difluoro-1,2,2,2-tetrachloroethane	Fusion	217		
		Vaporization	216,217		
C <sub>2</sub> FC1 <sub>5</sub>	fluoropentachloroethane	Fusion	216,222		
		Vaporization	216,222		
C <sub>2</sub> OF <sub>3</sub> Cl	trifluoroacetyl chloride	Fusion	223		
		Vaporization	223	223	

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$C_2OF_2Cl_2$	difluorochloroacetyl chloride	Vaporization	224		
$C_2H_3F_2Cl$	1,1-difluoro-1-chloroethane	Fusion	219	219	219
		Vaporization	225	225	
$C_2H_3F_2Cl$	1,1-difluoro-2-chloroethane	Vaporization	176		
$C_2H_2F_3Cl$	1,1,2-trifluoro-2-chloroethane	Vaporization	166, 226		
$C_2HFCl_2$	2-fluoro-1,1-dichloroethene	Vaporization	166, 226		
$C_2H_2F_2Cl_2$	1,1-difluoro-2,2-dichloroethane	Vaporization	166, 226		
$C_2H_2FCl_3$	1-fluoro-1,2,2-trichloroethane	Vaporization	166, 226		
$C_2HFCl_4$	2-fluoro-1,1,1,2-tetrachloroethane	Vaporization	226		
$C_2H_2OFCl$	fluoroacetyl chloride	Vaporization	227	227	
$C_2H_2OFCl$	chloroacetyl fluoride	Vaporization	227	227	
$C_2HO_2F_2Cl$	difluoroacetyl chloride	Vaporization	175		
$C_2HO_2F_2Cl$	difluorochloroacetic acid	Fusion	228		
		Vaporization	228		
$C_2HOFC_2Cl_2$	dichloroacetyl fluoride	Vaporization	226, 229		
$C_2Br_4$	tetrabromoethene	Sublimation	230		
		Fusion	231		
$C_2H_3Br$	bromoethene	Fusion	137		
		Vaporization	232, 233		
$C_2H_5Br$	bromoethane	Fusion	67, 234, 235, 236, 237	67, 234, 235	
		Vaporization	234, 236	234	238
$C_2H_2Br_2$	1,1-dibromoethene	Vaporization	231		
$C_2H_2Br_2$	cis-1,2-dibromoethene	Fusion	239		
		Vaporization	239		
$C_2H_2Br_2$	trans-1,2-dibromoethene	Fusion	239		
		Vaporization	239		
$C_2H_4Br_2$	1,1-dibromoethane	Vaporization	240		
$C_2H_4Br_2$	1,2-dibromoethane	Transition	241	241	241
		Fusion	241	241	241
		Vaporization	14, 43	14	
$C_2H_3Br_3$	1,1,2-tribromoethane	Fusion	242		
		Vaporization	25, 240	25	
$C_2H_2Br_4$	1,1,1,2-tetrabromoethane	Vaporization	243		
$C_2H_2Br_4$	1,1,2,2-tetrabromoethane	Fusion	224, 244	244	
		Vaporization	240, 245	240, 245	
$C_2HBr_5$	pentabromoethane	Fusion	246		
$C_2H_3OBr$	acetyl bromide	Fusion	199		
		Vaporization	247	247	

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$C_2H_3OBr$	bromoacetaldehyde	Vaporization	248		
$C_2H_3O_2Br$	bromoacetic acid	Fusion	249, 250		
		Vaporization	251		
$C_2H_5OBr$	2-bromoethanol	Vaporization	252		
$C_2H_5OBr$	bromoethyl methyl ether	Vaporization	253		
$C_2H_2OBr_2$	bromoacetyl bromide	Vaporization	254		
$C_2H_2O_2Br_2$	dibromoacetic acid	Fusion	250, 255		
		Vaporization	255		
$C_2H_4OBr_2$	bis(bromomethyl)ether	Vaporization	256	256	
$C_2HOBr_3$	tribromoacetaldehyde (bromal)	Vaporization	25	25	
$C_2HO_2Br_3$	tribromoacetic acid	Fusion	250, 257		
$C_2H_3OBr_3$	2,2,2-tribromoethanol	Fusion	258		
$C_2H_3O_2Br_3$	tribromoacetaldehyde mono-hydrate	Fusion	259, 260		
$C_2F_3Br$	trifluorobromoethene	Vaporization	168		
$C_2F_5Br$	pentafluorobromoethane	Vaporization	261		
$C_2F_4Br_2$	1,1,2,2-tetrafluoro-1,2-dibromoethane	Fusion	216, 262		
		Vaporization	262	262	
$C_2FBr_3$	fluorotribromoethene	Vaporization	263		
$C_2F_2Br_4$	1,1-difluoro-1,2,2,2-tetrabromoethane	Fusion	243		
		Vaporization	243		
$C_2OF_3Br$	trifluoroacetyl bromide	Fusion	264		
		Vaporization	264	264	
$C_2H_2FBr$	1-fluoro-1-bromoethene	Vaporization	166		
$C_2H_4FBr$	1-fluoro-2-bromoethane	Vaporization	243		
$C_2HF_2Br$	1,1-difluoro-2-bromoethene	Vaporization	243		
$C_2H_3F_2Br$	1,1-difluoro-2-bromoethane	Vaporization	166, 177, 265, 266		
$C_2H_2F_3Br$	1,1,2-trifluoro-1-bromoethane	Vaporization	243		
$C_2H_2F_3Br$	1,1,1-trifluoro-2-bromoethane	Vaporization	178		
$C_2H_3FBr_2$	2-fluoro-1,1-dibromoethane	Vaporization	266		
$C_2H_3FBr_2$	1-fluoro-1,2-dibromoethane	Fusion	265		
		Vaporization	265		
$C_2H_2F_2Br_2$	1,1-difluoro-2,2-dibromoethane	Vaporization	265		
$C_2H_2F_2Br_2$	1,1-difluoro-1,2-dibromoethane	Vaporization	267		
$C_2HF_3Br_2$	1,1,2-trifluoro-1,2-dibromoethane	Vaporization	168		
$C_2H_2FBr_3$	1-fluoro-1,2,2-tribromoethane	Vaporization	267		
$C_2H_2FBr_3$	1-fluoro-1,1,2-tribromoethane	Vaporization	243, 266		



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Formula	Description				
$C_2HF_2Br_3$	1,1-difluoro-1,2,2-tri-bromoethane	Vaporization	243		
$C_2HBr_4$	1-fluoro-1,1,2,2-tetrabromoethane	Vaporization	243		
$C_2H_2O_2FBr$	fluorobromoacetic acid	Fusion	226		
		Vaporization	226		
$C_2HO_2FBr_2$	fluorobromoacetyl bromide	Vaporization	226		
$C_2Cl_2Br_2$	1,2-dichloro-1,2-dibromoethene	Fusion	179, 268		
		Vaporization	239, 26B		
$C_2Cl_4Br_2$	1,1,2,2-tetrachloro-1,2-dibromoethane	Sublimation	269		
$C_2OCl_3Br$	trichloroacetyl bromide	Vaporization	25, 20B	25	
$C_2OCl_2Br_2$	chlorodibromoacetyl chloride	Vaporization	270		
$C_2H_2ClBr$	cis-1-chloro-2-bromoethene	Fusion	239, 271		
		Vaporization	239, 271		
$C_2H_2ClBr$	trans-1-chloro-2-bromoethene	Fusion	239, 271		
		Vaporization	239, 271		
$C_2H_4ClBr$	1-chloro-1-bromoethane	Fusion	25		
		Vaporization	25, 190	25, 190	
$C_2H_4ClBr$	1-chloro-2-bromoethane	Fusion	25, 191	25, 191	191
		Vaporization	25, 190	25, 190	
$C_2HCl_2Br$	1,1-dichloro-2-bromethene	Fusion	272		
		Vaporization	272		
$C_2HCl_2Br$	cis-1,2-dichloro-1-bromoethene	Fusion	273, 274		
		Vaporization	271, 273		
$C_2HCl_2Br$	trans-1,2-dichloro-1-bromoethane	Fusion	271		
$C_2H_3Cl_2Br$	1,1-dichloro-2-bromethane	Vaporization	25		
$C_2H_2Cl_3Br$	1,2,2-trichloro-1-bromoethane	Fusion	239		
		Vaporization	239		
$C_2H_2Cl_2Br_2$	1,1-dichloro-1,2-dibromoethane	Fusion	197, 274, 275		
		Vaporization	275		
$C_2H_2Cl_2Br_2$	1,2-dichloro-1,2-dibromoethane	Fusion	137, 276		
		Vaporization	137, 276, 277	277	
$C_2HCl_3Br_2$	1,1,2-trichloro-1,2-dibromoethane	Fusion	137		
$C_2HCl_3Br_2$	1,1,1-trichloro-2,2-dibromoethane	Fusion	278		
$C_2H_2ClBr_3$	2-chloro-1,1,2-tribromoethane	Fusion	239, 276, 278		
		Vaporization	239, 276		
$C_2HCl_2Br_3$	1,2-dichloro-1,1,2-tribromoethane	Fusion	239, 26B, 278		
		Vaporization	239, 26B		
$C_2HCl_2Br_2$	1,1-dichloro-1,2,2-tribromoethane	Fusion	272, 275		
		Vaporization	272, 275		

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$C_2H_2OC1Br$	bromoacetyl chloride	Vaporization	279		
$C_2F_3Cl_8r_2$	1,1,2-trifluoro-2-chloro-1,2-dibromoethane	Vaporization	216		
$C_2F_2Cl_2Br_2$	1,1-difluoro-2,2-dichloro-1,2-dibromoethane	Fusion	216, 217		
		Vaporization	216		
$C_2F_2Cl_2Br_2$	1,2-difluoro-1,2-dichloro-1,2-dibromoethane	Fusion	216, 217		
		Vaporization	216, 217		
$C_2FCl_3Br_2$	1-fluoro-1,2,2-trichloro-1,2-dibromoethane	Fusion	216		
$C_2OF_2ClBr$	fluorochlorobromoacetyl-fluoride	Vaporization	270		
$C_2HFC1_2Br_2$	2-fluoro-1,1-dichloro-1,2-dibromoethane	Vaporization	226		
$C_2HOFC1Br$	fluorobromoacetyl chloride	Vaporization	226		
$C_2HO_2FC1Br$	fluorochlorobromoacetic acid	Fusion	270		
		Vaporization	270		
$(CI_2)_2$	tetraiodoethene	Fusion	280		
$C_2H_5I$	iodoethane	Fusion	68		
		Vaporization	66, 281	66, 281	
$C_2H_2I_2$	1,1-diiodoethene	Fusion	282		
$C_2H_2I_2$	cis-1,2-diiodoethene	Fusion	283		
		Vaporization	271		
$C_2H_2I_2$	trans-1,2-diiodoethene	Sublimation	284	284	
		Fusion	271, 283		
		Vaporization	271		
$(CH_2I)_2$	1,2-diiodoethane	Sublimation	285	285	
		Fusion	285, 286		
$C_2H_3OI$	acetyl iodide	Vaporization	287		
$C_2H_3O_2I$	iodoacetic acid	Fusion	288		
$C_2H_5OI$	2-iodoethanol	Vaporization	289		
$C_2H_5OI$	iodomethyl methyl ether	Vaporization	207		
$C_2H_3F_2I$	1,1-difluoro-2-iodoethane	Vaporization	267		
$C_2H_2F_3I$	1,1,1-trifluoro-2-iodoethane	Vaporization	290		
$C_2H_2O_2FI$	fluoriodoacetic acid	Fusion	226		
$C_2OC1_3I$	trichloroacetyl iodide	Vaporization	208		
$C_2H_2ClI$	cis-1-chloro-2-iodoethene	Fusion	271		
		Vaporization	271		
$C_2H_2ClI$	trans-1-chloro-2-iodoethene	Fusion	271		
		Vaporization	271		
$C_2H_2OC1I$	iodoacetyl chloride	Vaporization	288		
$C_2H_6S$	ethanethiol	Fusion	199		
		Vaporization	291, 292	291, 292	

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$C_2H_6S$	dimethyl sulfide	Fusion	293	293	293
		Vaporization	291, 293	291, 293	
$C_2H_6OS$	dimethyl sulfoxide	Fusion	294		295
		Vaporization	295	295	
$C_2H_6O_2S$	dimethyl sulfone	Fusion	294, 296		
		Vaporization	294		
$C_2H_6O_3S$	dimethyl sulfite	Vaporization	294		
$C_2H_6O_4S$	dimethyl sulfate	Fusion	199, 278		
		Vaporization	183, 199, 278		
$C_2H_6O_4S$	ethylsulfuric acid	Fusion	297		
$C_2H_6O_4S \cdot H_2O$	ethylsulfuric acid mono-hydrate	Fusion	297		
$C_2H_5O_3FS$	ethyl fluorosulfonate	Vaporization	174	174	
$C_2H_5O_3ClS$	ethylchlorosulfonate	Vaporization	183		
$C_2H_4Br_2S$	bis(bromomethyl)sulfide	Vaporization	183	183	
$C_2H_6Se$	dimethylselenide	Vaporization	298		
$C_2H_6O_4Se$	dimethyl selenate	Vaporization	298		
$C_2H_6Te$	dimethyl telluride	Vaporization	299		
$C_2H_6F_2Te$	dimethyltellurium difluoride	Fusion	300		
$C_2H_6Cl_2Te$	dimethyltellurium chloride	Fusion	299		
$C_2H_6Br_2Te$	dimethyltellurium dibromide	Fusion	299		
$C_2H_6I_2Te$	dimethyltellurium diiodide	Fusion	299, 301		
$C_2N_2$	cyanogen	Fusion	302	302	302
		Vaporization	302, 303, 304		
$C_2H_3N$	acetonitrile	Fusion	37, 143, 305	37, 305	
		Vaporization	143, 305, 306	305, 306	
$C_2H_7N$	ethylamine	Fusion	307		34
		Vaporization	34, 307	34, 307	
$C_2H_7N$	dimethylamine	Fusion	308	308	308
		Vaporization	308	308	
$C_2H_8N_2$	1,2-ethanediamine (ethylenediamine)	Fusion	309	309	
		Vaporization	309	309	
$C_2H_8N_2$	1,1-dimethylhydrazine	Vaporization	310		
$C_2H_8N_2$	1,2-dimethylhydrazine	Vaporization	310		
$C_2H_5ON$	acetamide	Fusion	311, 312, 313, 314	311, 313, 314	314
		Transition	314	314	
$C_2H_5O_2N$	nitroethane	Fusion	25		
		Vaporization	315	315	
$C_2H_5NO_2$	ethyl nitrite	Vaporization	315	315	
$C_2H_5NO_3$	ethyl nitrate	Fusion	330		
		Vaporization	315	315	

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$C_2H_7ON$	2-aminoethanol	Fusion	316		
		Vaporization	316		
$C_2H_4O_3N_2$	nitroacetamide	Fusion	317		
$C_2H_4O_6N_2$	glycol dinitrate	Fusion	318	318	
		Vaporization	319		
$C_2H_6O_2N_2$	ethylnitramine	Fusion	320		
$C_2F_3N$	trifluoroacetamide	Vaporization	165		
$C_2F_6N_2$	hexafluoroazomethane	Fusion	321		
		Vaporization	321	321	
$C_2H_2FN$	fluoroacetamide	Vaporization	174	174	
$C_2H_5F_2N$	2,2-difluoroethylamine	Vaporization	322		
$C_2HF_6N$	bis(trifluoromethyl)amine	Vaporization	321	321	
$C_2H_2OFN$	fluoroacetamide	Fusion	172, 175, 266, 322		
$C_2H_4O_2FN$	2-fluoroethylnitrite	Vaporization	174	174	
$C_2H_3OF_2N$	difluoroacetamide	Fusion	172, 266, 322		
		Vaporization	172, 266, 322		
$C_2H_2OF_3N$	trifluoroacetamide	Fusion	165		
		Vaporization	165		
$C_2H_4O_2F_2N_2$	2,2-difluoroethylnitramine	Fusion	172, 320		
$C_2H_2ClN$	chloroacetonitrile	Vaporization	323		
$C_2H_6ClN$	N-chlorodimethylamine	Vaporization	324		
$C_2H_4OClN$	chloroacetamide	Fusion	325		
$C_2H_4OClN$	N-chloroacetamide	Fusion	326		
$C_2H_2OF_2ClN$	difluorochloroacetamide	Fusion	224		
		Vaporization	224		
$C_2H_4OBrN$	bromoacetamide	Fusion	327		
$C_2H_4OBrN$	N-bromoacetamide	Fusion	326		
$C_2H_2OBr_3N$	tribromoacetamide	Fusion	328		
$C_2H_2OFClBrN$	fluorochlorobromoacetamide	Fusion	270		
$C_2H_3SN$	methyl thiocyanate	Fusion	199		
		Vaporization	25	25	
$C_2H_3SN$	methyl isothiocyanate	Fusion	25		
		Vaporization	25	25	
$C_2H_6O_6TeN_2$	dimethyltellurium dinitrate	Fusion	299		
$C_2H_7P$	dimethylphosphine	Vaporization	329	329	
$C_2H_6O_3FP$	dimethylfluorophosphonate	Vaporization	174	174	
$C_2O_2FN_2P$	phosphorous fluorodiisocyanate	Fusion	330		
		Vaporization	330	330	
$C_2O_2ClN_2P$	phosphorous chlorodiisocyanate	Fusion	330		
		Vaporization	330	330	

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$C_2H_7As$	dimethylarsine	Vaporization	331		
$C_2H_7As$	ethylarsine	Vaporization	332		
$C_2H_5OAs$	ethylarsenic oxide	Vaporization	183	183	
$C_2H_5F_2As$	ethyldifluoroarsine	Fusion	125		
		Vaporization	125	125	
$C_2H_6ClAs$	dimethylchloroarsine	Vaporization	333		
$C_2H_5Cl_2As$	ethyldichloroarsine	Fusion	183		
		Vaporization	334	334	
$C_2H_2Cl_3As$	2-chlorovinylldichloroarsine	Vaporization	174	174	
$C_2H_2OClAs$	2-chlorovinylarsine oxide	Fusion	335		
$C_2H_4O_3ClAs$	2-chlorovinylarsonic acid	Fusion	335		
$C_2H_2F_2ClAs$	2-chlorovinylldifluoroarsine	Vaporization	125		
$C_2H_6BrAs$	dimethylbromoarsine	Vaporization	333		
$C_2H_5Br_2As$	ethyldibromoarsine	Vaporization	174	174	
$C_2H_4O_3BrAs$	2-bromovinylarsonic acid	Fusion	335		
$C_2H_2ClBr_2As$	2-chlorovinylldibromoarsine	Vaporization	335		
$C_2H_6IAs$	dimethyliodoarsine	Vaporization	333		
$C_2H_2ClI_2As$	2-chlorovinylldiiodoarsine	Fusion	335		

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14. Mathews-3	30. Bond and Beach-1
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16. Eucken and Schröder-1	32. Kireev and Skvorshova-1
	33. Pitzer-8

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| 39. Biltz, Sapper, and Wünnenberg-1                                      | 80. Beckmann-5                          |
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| 58. Bridgman-11  | 99. Linhard-1                           |
| 59. Frederick and Hildebrand-3   | 100. Lowry and Culter-1                 |
| 60. Kelley-21  | 101. Getman-5                           |
| 61. Pirsch-1   | 102. Timmermans and Hennaut-Roland-2    |
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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Si		Fusion	1,2,3	1,2,3	9
SiO <sub>2</sub>		Transition	4,5,6,7,8	4,5,6,7	
		Fusion	4,5,6,7,8	4,5,6,7	
SiH <sub>4</sub>	silane	Transition	9	9	
		Fusion	9	9	
		Vaporization	10,11,12,13	10,11,12,13	
Si <sub>2</sub> H <sub>6</sub>	disilane	Fusion	10		
		Vaporization	10,12,13	10,12,13	
Si <sub>3</sub> H <sub>8</sub>	trisilane	Fusion	10		
		Vaporization	10,13	10,13	
Si <sub>4</sub> H <sub>10</sub>	tetrasilane	Fusion	14		
		Vaporization	10,13,14	10,13,14	
(SiH <sub>3</sub> ) <sub>2</sub> O	disiloxane	Fusion	15		
		Vaporization	15	15	
SiF <sub>4</sub>	tetrafluorosilane	Sublimation	16,17,18,19,20	16,17,18,19,20	
		Fusion	18	18	
		Vaporization	16,17,18,19,20	16,17,18,19,20	
Si <sub>2</sub> F <sub>6</sub>	hexafluorodisilane	Sublimation	21	21	
		Fusion	21	21	
		Vaporization	21	21	
Si <sub>2</sub> OF <sub>6</sub>	hexafluorodisiloxane	Fusion	22		
		Vaporization	22	22	
SiH <sub>3</sub> F	fluorosilane	Vaporization	23	23	
SiH <sub>2</sub> F <sub>2</sub>	difluorosilane	Fusion	23	23	
		Vaporization	23	23	
SiHF <sub>3</sub>	trifluorosilane	Fusion	23,24		
		Vaporization	23,24	23,24	
SiCl <sub>4</sub>	tetrachlorosilane	Fusion	15,25,26,27	15,26,27	
		Vaporization	15,28,29	15,28,29	
Si <sub>2</sub> Cl <sub>6</sub>	hexachlorodisilane	Fusion	30		
		Vaporization	30	30	
Si <sub>3</sub> Cl <sub>8</sub>	octachlorotrisilane	Vaporization	30,31	30	
Si <sub>2</sub> OC <sub>2</sub> Cl <sub>3</sub>	hexachlorodisiloxane	Fusion	15		
		Vaporization	15	15	
SiH <sub>3</sub> Cl	chlorosilane	Fusion	32		
		Vaporization	32	32	
SiH <sub>2</sub> Cl <sub>2</sub>	dichlorosilane	Fusion	32		
		Vaporization	32	32	
SiHCl <sub>3</sub>	trichlorosilane	Fusion	24,33		
		Vaporization	24,33		

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
SiF <sub>3</sub> Cl	trifluorochlorosilane	Fusion	21,34		
		Vaporization	21,34	21,34	
SiF <sub>2</sub> Cl <sub>2</sub>	difluorodichlorosilane	Fusion	34		
		Vaporization	21,34	21,34	
SiFCl <sub>3</sub>	fluorotrichlorosilane	Fusion	34		
		Vaporization	34	34	
Si <sub>2</sub> O <sub>2</sub> F <sub>4</sub> Cl <sub>2</sub>	1,1,1,3-tetrafluoro-3,3-dichlorodisiloxane	Fusion	22		
		Vaporization	22	22	
Si <sub>2</sub> O <sub>2</sub> F <sub>3</sub> Cl <sub>3</sub>	1,1,1-Trifluoro-3,3,3-trichlorodisiloxane	Fusion	22		
		Vaporization	22	22	
SiHFC1 <sub>2</sub>	fluorodichlorosilane	Fusion	24		
		Vaporization	24	24	
SiBr <sub>4</sub>	tetrabromosilane	Fusion	35,36	35,36	
		Vaporization	35,36,37	35,36,37	
Si <sub>2</sub> Br <sub>6</sub>	hexabromodisilane	Fusion	38		
		Vaporization	3B		
Si <sub>3</sub> Br <sub>8</sub>	octabromotrisilane	Fusion	38		
SiH <sub>3</sub> Br	bromosilane	Fusion	39		
		Vaporization	39	39	
SiH <sub>2</sub> Br <sub>2</sub>	dibromosilane	Fusion	39		
		Vaporization	39	39	
SiHBr <sub>3</sub>	tribromosilane	Fusion	40		
		Vaporization	40	40	
Si <sub>2</sub> H <sub>5</sub> Br	bromodisilane	Fusion	41		
		Vaporization	41	41	
SiF <sub>3</sub> Br	trifluorobromosilane	Fusion	42		
		Vaporization	42	42	
SiF <sub>2</sub> Br <sub>2</sub>	difluorodibromosilane	Fusion	42		
		Vaporization	42	42	
SiFBr <sub>3</sub>	fluorotribromosilane	Fusion	42		
		Vaporization	42	42	
SiCl <sub>3</sub> Br	trichlorobromosilane	Vaporization	43		
SiCl <sub>2</sub> Br <sub>2</sub>	dichlorodibromosilane	Vaporization	43		
SiClBr <sub>3</sub>	chlorotribromosilane	Fusion	43		
		Vaporization	43		
SiFCl <sub>2</sub> Br	fluoredichlorobromosilane	Fusion	44		
		Vaporization	44	44	
SiFClBr <sub>3</sub>	fluoro-chlorodibromosilane	Fusion	44		
		Vaporization	44	44	
SiI <sub>4</sub>	tetraiodosilane	Fusion	37,45,46		
		Vaporization	45		

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta G_p$
Formula	Description				
$\text{Si}_2\text{I}_6$	hexoiododisilane	Fusion	47		
$\text{SiH}_3\text{I}$	iodosilane	Fusion	46		
		Vaporization	46	46	
$\text{SiH}_2\text{I}_2$	diiodosilane	Fusion	46		
		Vaporization	46	46	
$\text{SiHI}_3$	triiodosilane	Fusion	48		
		Vaporization	48	48	
$\text{SiCl}_3\text{I}$	trichloroiodosilane	Vaporization	43		
$\text{SiCl}_2\text{I}_2$	dichlorodiiodosilane	Vaporization	43		
$\text{SiClI}_3$	chlorotriiodosilane	Fusion	43		
		Vaporization	43		
$\text{SiBr}_3\text{I}$	tribromoiodosilane	Fusion	43		
		Vaporization	43		
$\text{SiBr}_2\text{I}_2$	dibromodiiodosilane	Fusion	43		
		Vaporization	43		
$\text{SiBrI}_3$	bromotriiodosilane	Fusion	43		
		Vaporization	43		
$\text{SiS}_2$		Fusion	49		
$\text{SiSCl}_2$		Fusion	50		
$\text{SiSBr}_2$		Fusion	51		
$(\text{SiH}_3)_3\text{N}$	trisilylamine	Fusion	52		
		Vaporization	52	52	
$3\text{Bi}_2\text{O}_3 \cdot \text{SiO}_2$		Fusion	53		
$38\text{i}_2\text{O}_3 \cdot 2\text{SiO}_2$		Fusion	53		
$8\text{i}_2\text{O}_3 \cdot 8\text{SiO}_2$		Fusion	53		
$\text{SiH}_6\text{C}$	methylsilane	Fusion	32		
		Vaporization	32	32	
$\text{SiH}_8\text{C}_2$	ethylsilane	Vaporization	54	54	
$\text{SiH}_8\text{C}_2$	dimethylsilane	Fusion	32		
		Vaporization	32	32	
$\text{SiH}_{10}\text{C}_3$	n-propylsilane	Vaporization	54	54	
$\text{SiH}_{10}\text{C}_3$	trimethylsilane	Vaporization	55		
$\text{SiH}_{12}\text{C}_4$	diethylsilane	Vaporization	55	54	
$\text{SiH}_{12}\text{C}_4$	tetramethylsilane	Fusion	56	56	56
		Vaporization	56	56	
$\text{SiH}_{14}\text{C}_5$	trimethylethylsilane	Vaporization	57, 58	57	
$\text{SiH}_{18}\text{C}_7$	methyltriethylsilane	Vaporization	57	57	
$\text{SiH}_{20}\text{C}_8$	tetramethylsilane	Vaporization	57	57	
$\text{Si}_2\text{H}_{18}\text{C}_6$	hexamethyldisilane	Fusion	59		
		Vaporization	59	59	
$\text{SiH}_{10}\text{OC}_3$	trimethylsilanol	Vaporization	60		

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$\text{SiH}_{14}\text{OC}_5$	ethoxytrimethylsilane	Vaporization	61	61	
$\text{SiH}_{16}\text{O}_2\text{C}_6$	diethoxydimethylsilane	Vaporization	61	61	
$\text{SiH}_{18}\text{O}_3\text{C}_7$	triethoxymethylsilane	Vaporization	61	61	
$\text{SiH}_{20}\text{O}_4\text{C}_8$	tetraethoxysilane	Fusion	62	62	
		Vaporization	62	62	
$\text{Si}_2\text{H}_{18}\text{OC}_6$	hexamethyldisiloxane	Vaporization	63	63	
$\text{Si}_2\text{O}_3\text{H}_{22}\text{C}_8$	1,3-diethoxytetramethyl- disiloxane	Vaporization	62	62	
$\text{Si}_2\text{H}_{30}\text{O}_6\text{C}_{12}$	hexaethoxydisilane	Vaporization	64	64	
$\text{Si}_3\text{H}_{24}\text{O}_2\text{C}_8$	octamethyltrisiloxane	Vaporization	63	63	
$\text{Si}_3\text{H}_{28}\text{O}_4\text{C}_{10}$	1,5-diethoxyhexamethyl- trisiloxane	Vaporization	61	61	
$\text{Si}_4\text{H}_{24}\text{O}_4\text{C}_8$	octamethylcyclotetrasiloxane	Vaporization	63	63	
$\text{Si}_4\text{H}_{30}\text{O}_3\text{C}_{10}$	decamethyltetrasiloxane	Vaporization	63	63	
$\text{SiH}_3\text{F}_3\text{C}$	methyltrifluorosilane	Fusion	65		
		Vaporization	65	65	
$\text{SiH}_6\text{F}_2\text{C}_2$	dimethyldifluorosilane	Fusion	66		
		Vaporization	66	66	
$\text{SiH}_5\text{F}_3\text{C}_2$	ethyltrifluorosilane	Fusion	67		
		Vaporization	67	67	
$\text{SiH}_9\text{FC}_3$	trimethylfluorosilane	Fusion	66		
		Vaporization	66	66	
$\text{SiH}_{10}\text{F}_2\text{C}_4$	diethyldifluorosilane	Fusion	23		
		Vaporization	23	23	
$\text{SiH}_5\text{ClC}$	methylchlorosilane	Fusion	32		
		Vaporization	32	32	
$\text{SiH}_4\text{Cl}_2\text{C}$	methyldichlorosilane	Fusion	32		
		Vaporization	32	32	
$\text{SiH}_3\text{Cl}_3\text{C}$	methyltrichlorosilane	Fusion	65		
		Vaporization	65	65	
$\text{SiH}_6\text{Cl}_2\text{C}_2$	dimethyldichlorosilane	Fusion	66		
		Vaporization	66	66	
$\text{SiH}_5\text{Cl}_3\text{C}_2$	ethyltrichlorosilane	Fusion	67		
		Vaporization	67	67	
$\text{SiHgClC}_3$	trimethylchlorosilane	Fusion	66		
		Vaporization	66	66	
$\text{SiH}_{10}\text{Cl}_2\text{C}_4$	diethyldichlorosilane	Vaporization	61	61	
$\text{SiH}_{15}\text{ClC}_6$	triethylchlorosilane	Vaporization	61	61	
$\text{SiH}_5\text{OCl}_3\text{C}_2$	ethoxytrichlorosilane	Vaporization	61	61	
$\text{SiH}_8\text{OCl}_2\text{C}_3$	ethoxymethyldichlorosilane	Vaporization	61	61	
$\text{Si}_2\text{H}_{12}\text{OCl}_2\text{C}_4$	tetramethyl-1,3-dichloro- disilane	Fusion	68		
		Vaporization	68	68	

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$\text{SiH}_3\text{FCl}_2\text{C}$	methylfluorodichlorosilane	Fusion	65		
		Vaporization	65	65	
$\text{SiH}_6\text{FClC}_2$	dimethylfluorochlorosilane	Fusion	66		
		Vaporization	66	66	
$\text{SiH}_5\text{F}_2\text{ClC}_2$	ethyldifluorochlorosilane	Vaporization	67	67	
$\text{SiH}_5\text{F}_2\text{Cl}_2\text{C}_2$	ethylfluorodichlorosilane	Vaporization	67	67	
$\text{SiO}_4\text{N}_4\text{C}_4$	silicon tetracyanate	Fusion	69		
		Vaporization	69	69	
$\text{SiO}_4\text{N}_4\text{C}_4$	silicon tetraisocyanate	Fusion	69		
		Vaporization	69	69	
$\text{Si}_2\text{H}_9\text{NC}$	2-methyldisilazane	Vaporization	70,71	71	
$\text{Si}_2\text{H}_{11}\text{NC}_2$	2-ethyldisilane	Fusion	71		
		Vaporization	70,71	71	
$\text{SiH}_9\text{ONC}_4$	trimethylsilicon isocyanate	Fusion	72		
		Vaporization	72	72	
$\text{SiH}_6\text{O}_2\text{N}_2\text{C}_4$	dimethylsilicon diisocyanate	Fusion	72,73		
		Vaporization	72,73	72,73	
$\text{SiH}_3\text{O}_3\text{N}_3\text{C}_4$	methylsilicon triisocyanate	Fusion	72		
		Vaporization	72	72	
$\text{SiH}_5\text{O}_4\text{N}_3\text{C}_5$	ethoxysilicon triisocyanate	Vaporization	73	73	
$\text{SiH}_{10}\text{O}_4\text{N}_2\text{C}_6$	diethoxysilicon diisocyanate	Vaporization	73	73	
$\text{SiH}_{15}\text{O}_4\text{NC}_7$	triethoxysilicon isocyanate	Vaporization	73	73	
$\text{SiOF}_3\text{NC}$	trifluorosilicon isocyanate	Vaporization	74		
$\text{SiO}_2\text{F}_2\text{N}_2\text{C}_2$	difluorosilicon diisocyanate	Fusion	74		
		Vaporization	74	74	
$\text{SiO}_3\text{FN}_3\text{C}_3$	fluorosilicon triisocyanate	Fusion	74		
		Vaporization	74	74	
$\text{SiS}_4\text{N}_4\text{C}_4$	silicon tetraisothiocyanate	Fusion	75		
		Vaporization	75	75	
$\text{SiH}_9\text{SNC}_4$	trimethylsilicon isothiocyanate	Fusion	75		
		Vaporization	75	75	
$\text{SiH}_6\text{S}_2\text{N}_2\text{C}_4$	silicon diisothiocyanate	Fusion	75		
		Vaporization	75	75	
$\text{SiH}_3\text{S}_3\text{N}_3\text{C}_4$	methylsilicon triisothiocyanate	Fusion	75		
		Vaporization	75	75	
$\text{SiH}_9\text{O}_3\text{SNC}_4$	trimethoxysilicon isothiocyanate	Vaporization	75	75	

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| 11. Finholt, Bond, Wilzbach, and Schlesinger-1 | 53. Otin-1   |
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| 17. Moissan-2                                  | 59. Brockway and Davidson-1  |
| 18. Patnode and Papish-1                       | 60. Sauer-1  |
| 19. Forbes and Anderson-2                      | 61. Stull-3  |
| 20. Fisher and Weidemann-1                     | 62. Solara and Moles-1   |
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| 22. Booth and Osten-1                          | 64. Martin-1   |
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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Ge		Fusion	1,2	1,2	
GeO		Sublimation	1		
GeO <sub>2</sub>		Transition	1		
GeH <sub>4</sub>	germane	Transition	5	5	
		Fusion	2,5,6	5	
		Vaporization	5,8	5	
Ge <sub>2</sub> H <sub>6</sub>	digermane	Fusion	2,9		
		Vaporization	8	8	
Ge <sub>3</sub> H <sub>8</sub>	trigermane	Fusion	9		
		Vaporization	9	9	
GeF <sub>4</sub>		Sublimation	10	10	
		Fusion	2,10		
GeCl <sub>4</sub>		Fusion	2,7,11		
		Vaporization	2,11,12,13,14	2,13,14	
GeHCl <sub>3</sub>		Fusion	15		
		Vaporization	2,15	2,15	
GeF <sub>3</sub> Cl		Fusion	11		
		Vaporization	11		
GeF <sub>2</sub> Cl <sub>2</sub>		Fusion	11		
		Vaporization	11	11	
GeFCl <sub>3</sub>		Fusion	11		
		Vaporization	11	11	
GeBr <sub>4</sub>		Fusion	2,7,16		
		Vaporization	2,16	2,16	
GeI <sub>4</sub>		Fusion	2,7		
GeS		Fusion	3		
GeSe		Fusion	17		
GeSe <sub>2</sub>		Fusion	17		
Ge(CH <sub>3</sub> ) <sub>4</sub>	tetramethylgermane	Fusion	2,7		
		Vaporization	2,7	2,7	
Ge(C <sub>2</sub> H <sub>5</sub> ) <sub>4</sub>	tetraethylgermane	Fusion	2		
		Vaporization	2		
GeH <sub>6</sub> OC <sub>2</sub>	dimethylgermanium oxide	Fusion	18		
		Vaporization	18		
GeH <sub>5</sub> F <sub>3</sub> C <sub>2</sub>	ethylgermanium trifluoride	Fusion	19		
		Vaporization	19		
GeH <sub>3</sub> Cl <sub>3</sub> C	methylgermanium trichloride	Vaporization	20		
GeH <sub>6</sub> Cl <sub>2</sub> C <sub>2</sub>	dimethylgermanium dichloride	Fusion	20		
		Vaporization	20		
GeH <sub>5</sub> Cl <sub>3</sub> C <sub>2</sub>	ethylgermanium trichloride	Vaporization	19		



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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$\text{GeH}_5\text{Br}_3\text{C}_2$	ethylgermanium tribromide	Vaporization	19	21	
$\text{GeH}_5\text{I}_3\text{C}_2$	ethylgermanium triiodide	Fusion	19		
		Vaporization	19		
$\text{GeH}_6\text{SC}_2$	dimethylgermanium sulfide	Fusion	18		
		Vaporization	18		
$\text{GeO}_4\text{N}_4\text{C}_4$	germanium tetraisocyanate	Fusion	21		
		Vaporization	21		

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Sn	stannane	Transition	1,2,3,4,5	1,2,3,4,5	31
		Fusion	6,7	7	
		Vaporization	8,9,10		
SnO <sub>2</sub>		Transition	11,12	11,12	
SnH <sub>4</sub>		Fusion	13		
		Vaporization	13	13	
SnF <sub>4</sub>		Vaporization	14		
SnCl <sub>2</sub>		Fusion	15,16	15,16	
		Vaporization	17,18	17,18	
SnCl <sub>4</sub>		Fusion	15,19,20,21,22	15,19	
		Vaporization	20,22,23,24,25	23	
SnBr <sub>2</sub>		Fusion	15,16	15	
		Vaporization	18	18	
SnBr <sub>4</sub>		Transition	21		
		Fusion	7,15,16,21,22	7,15	
		Vaporization	18,22	18	
SnCl <sub>3</sub> Br		Fusion	26		
SnCl <sub>2</sub> Br <sub>2</sub>		Fusion	26		
SnCl <sub>3</sub> Br <sub>3</sub>		Fusion	26		
SnI <sub>2</sub>		Fusion	27,28,29		
		Vaporization	18,29	18	
SnI <sub>4</sub>		Fusion	27,29,30,31	30,31	
		Vaporization	29,31	31	
SnS		Fusion	32,33		
SnSe		Fusion	33		
SnTe		Fusion	33		
SnH <sub>6</sub> C	methylstannane	Vaporization	34	34	
SnH <sub>8</sub> C <sub>2</sub>	dimethylstannane	Vaporization	34	34	
SnH <sub>10</sub> C <sub>3</sub>	trimethylstannane	Vaporization	34	34	
SnH <sub>12</sub> C <sub>4</sub>	tetramethylstannane	Vaporization	34,35	34,35	
SnH <sub>14</sub> C <sub>5</sub>	ethyltrimethylstannane	Vaporization	35	35	
SnH <sub>16</sub> C <sub>6</sub>	n-propyltrimethylstannane	Vaporization	35	35	
SnH <sub>20</sub> C <sub>8</sub>	tetraethylstannane	Fusion	36		
		Vaporization	37		
SnH <sub>3</sub> Cl <sub>3</sub> C	methyltin trichloride	Fusion	38		
SnH <sub>6</sub> Cl <sub>2</sub> C <sub>2</sub>	dimethyltin dichloride	Fusion	39		
SnH <sub>9</sub> ClC <sub>3</sub>	trimethyltin chloride	Fusion	40		
SnH <sub>15</sub> ClC <sub>6</sub>	triethyltin chloride	Fusion	41		
		Vaporization	41		
SnH <sub>3</sub> Br <sub>3</sub> C	methyltin tribromide	Fusion	38,42		
SnH <sub>6</sub> Br <sub>2</sub> C <sub>2</sub>	dimethyltin dibromide	Fusion	38,39		

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Table 26 (Continued)  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$\text{SnH}_9\text{BrC}_3$	trimethyltin bromide	Fusion	38		
$\text{SnH}_{15}\text{BrC}_6$	trimethyltin bromide	Fusion	36		
		Vaporization	36		
$\text{SnH}_3\text{I}_3\text{C}$	methyltin triiodide	Fusion	38,41		
$\text{SnH}_6\text{I}_2\text{C}_2$	dimethyltin diiodide	Fusion	38		
$\text{SnH}_9\text{IC}_3$	trimethyltin iodide	Fusion	36		
		Vaporization	38		
$\text{SnH}_{15}\text{IC}_6$	triethyltin iodide	Fusion	36		
		Vaporization	36		

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Table 27  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Pb		Fusion	1,2,3	1,2,3	
		Vaporization	2,4,5,6,7	2,4,6,7	
PbO		Transition	8,9		
		Fusion	1,10,11	1	
		Vaporization	4	4	
PbF <sub>2</sub>		Fusion	1,12,13,18	1	
		Vaporization	4,14	4,14	
PbF <sub>4</sub>		Fusion	15		
PbCl <sub>2</sub>		Fusion	1,12,16,17,18, 19,20,21,24	1,16,17	
		Vaporization	2,4,14,22,23	2,4,14,22,23	
PbCl <sub>2</sub> •2PbO		Fusion	21		
PbCl <sub>2</sub> •4PbO		Fusion	21		
PbFCl		Fusion	13		
PbBr <sub>2</sub>		Fusion	1,2,13,16,25	1,2,13,16	
		Vaporization	4,14,1,17,22,23	4,14,17,22, 23	
PbBr <sub>2</sub> •2PbO		Fusion	25		
PbBrF		Fusion	13		
PbI <sub>2</sub>		Fusion	1,2,26,27	1,2,26	
		Vaporization	4,22,28	4,22,28	
PbS		Sublimation	29	29	
		Fusion	1,30	1	
PbSO <sub>4</sub>		Transition	31	31	
		Fusion	1	1	
PbSO <sub>4</sub> •PbO		Fusion	32		
PbSO <sub>4</sub> •2PbO		Fusion	32		
PbSe		Fusion	33		
PbTe		Fusion	34		
Pb(PO <sub>3</sub> ) <sub>2</sub>		Fusion	24		
Pb(PO <sub>4</sub> ) <sub>2</sub>		Fusion	12		
Pb•Pb <sub>3</sub> (PO <sub>4</sub> ) <sub>2</sub>		Fusion	18		
5PbO•Pb <sub>3</sub> (PO <sub>4</sub> ) <sub>2</sub>		Fusion	18		
3Pb <sub>3</sub> (PO <sub>4</sub> ) <sub>2</sub> •PbF <sub>2</sub>		Fusion	12		
3Pb <sub>3</sub> (PO <sub>4</sub> ) <sub>2</sub> •PbCl <sub>2</sub>		Fusion	18		
Pb <sub>3</sub> (AsO <sub>4</sub> ) <sub>2</sub>		Fusion	18		
5PbO•Pb <sub>3</sub> (AsO <sub>4</sub> ) <sub>2</sub>		Fusion	18		
3Pb <sub>3</sub> (AsO <sub>4</sub> ) <sub>2</sub> •PbF <sub>2</sub>		Fusion	18		
3Pb <sub>3</sub> (AsO <sub>4</sub> ) <sub>2</sub> •PbCl <sub>2</sub>		Fusion	18		
PbO•Sb <sub>2</sub> O <sub>3</sub>		Fusion	35		
2PbO•Bi <sub>2</sub> O <sub>3</sub>		Fusion	36		

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Table 27  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$2\text{PbO} \cdot 3\text{Bi}_2\text{O}_3$	tetramethyllead	Fusion	36		
$\text{PbH}_{12}\text{C}_4$		Fusion	37		
		Vaporization	38	38	
$\text{PbH}_{14}\text{C}_5$	ethyltrimethyllead	Vaporization	39	39	
$\text{PbH}_{16}\text{C}_6$	diethyldimethyllead	Vaporization	39	39	
$\text{PbH}_{18}\text{C}_7$	triethylmethyllead	Vaporization	39	39	
$\text{PbH}_{20}\text{C}_8$	tetraethyllead	Fusion	40		
		Vaporization	41	41	
$\text{PbO} \cdot \text{SiO}_2$		Fusion	10		
$2\text{PbO} \cdot \text{SiO}_2$		Fusion	10		
$4\text{PbO} \cdot \text{SiO}_2$		Transition	10		
$3\text{PbS} \cdot 2\text{SiS}_2$		Fusion	30		

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12. Amadori-18	34. Kimura-4
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18. Amadori-14	40. Buckler and Norrish-1
19. Hackmeister-1	41. Thomson-1
20. Korreng-1	
21. Ruer-1	
22. Niwa-3	

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Table 28  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Ga		Transition	1	1	
		Fusion	1,2,3,4,5	1,2	
		Vaporization	6	6	
Ga <sub>2</sub> O <sub>3</sub>		Fusion	7		
GaF <sub>3</sub>		Vaporization	8,9		
GaCl <sub>2</sub>		Fusion	10		
Ga <sub>2</sub> Cl <sub>6</sub>		Sublimation	10,11	10,11	
		Fusion	10,11	10,11	
		Vaporization	10,11	10,11	
Ga <sub>2</sub> Br <sub>6</sub>		Fusion	12		
		Vaporization	11,12	11	
Ga <sub>2</sub> I <sub>6</sub>		Fusion	12		
		Vaporization	11,12	11	
GaS		Fusion	13,14		
Ga <sub>2</sub> S <sub>3</sub>		Fusion	13,14		
GaSe		Fusion	13		
GaTe		Fusion	13		
Ga <sub>2</sub> Te <sub>3</sub>		Fusion	13		
GaCl <sub>3</sub> ·NH <sub>3</sub>		Fusion	15		
GaBr <sub>3</sub> ·NH <sub>3</sub>		Fusion	15		
GaI <sub>3</sub> ·NH <sub>3</sub>		Fusion	15		
Ga(CH <sub>3</sub> ) <sub>3</sub>		Fusion	16,17		
		Vaporization	17	17	
Ga(C <sub>2</sub> H <sub>5</sub> ) <sub>3</sub>		Fusion	18		
		Vaporization	18	18	
Ga(CH <sub>3</sub> ) <sub>3</sub> ·NH <sub>3</sub>		Fusion	17		
Ga(CH <sub>3</sub> ) <sub>3</sub> N(CH <sub>3</sub> ) <sub>3</sub>		Fusion	16	16	
		Vaporization	16	16	
Ga(CH <sub>3</sub> ) <sub>3</sub> ·N(C <sub>2</sub> H <sub>5</sub> ) <sub>3</sub>		Sublimation	16	16	
		Fusion	16	16	
		Vaporization	16	16	
Ga(CH <sub>3</sub> ) <sub>2</sub> Cl·NH <sub>3</sub>		Fusion	17		
Ga(CH <sub>3</sub> ) <sub>2</sub> Cl·2NH <sub>3</sub>		Fusion	17		

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Table 28(Continued)  
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Table 29  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
In		Fusion	1,2,3	1	
		Vaporization			
InF <sub>3</sub>		Fusion	5,6		
InCl		Transition	7		
		Fusion	7,8		
		Vaporization	8,9	8,9	
InCl <sub>2</sub>		Fusion	8,10		
		Vaporization	8,9	8,9	
InCl <sub>3</sub>		Sublimation	8	8	
		Fusion	7		
InBr		Fusion	8		
		Vaporization	8,9	8,9	
InBr <sub>2</sub>		Fusion	8		
		Vaporization	8,9	8,9	
InBr <sub>3</sub>		Sublimation	8	8	
		Fusion	7,8		
InI		Fusion	8,11		
		Vaporization	8,9	8,9	
InI <sub>3</sub>		Fusion	7		
InS		Fusion	12,13		
In <sub>2</sub> S		Fusion	13		
In <sub>2</sub> S <sub>3</sub>		Fusion	12,13		
InSe		Fusion	12		
In <sub>2</sub> Se <sub>3</sub>		Fusion	12		
InTe		Fusion	12		
In <sub>2</sub> Te <sub>3</sub>		Fusion	12		
In(CH <sub>3</sub> ) <sub>3</sub>		Sublimation	14	14	
		Fusion	14	14	
		Vaporization	14	14	
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Table 30  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Tl		Transition	1,2,3,4,5,6	1,2,5	5
		Fusion	4,5,6,7,8,9	5,7	
		Vaporization	10,11,12,13	11,12,13	
Tl <sub>2</sub> O		Fusion	14		
		Vaporization	15		
Tl <sub>2</sub> O <sub>3</sub>		Fusion	16		
TlF		Fusion	17		
		Vaporization	17		
TlF <sub>3</sub>		Fusion	18,19		
TlCl		Sublimation	20,21	20,21	
		Fusion	22,23,24,25,26,27	22,27	
		Vaporization	20	20	
TlCl <sub>3</sub> ·4H <sub>2</sub> O		Fusion	28		
TlClO <sub>4</sub>		Transition	29,30		
		Fusion	31		
TlBr		Sublimation	20,21,32	20,21,32	
		Fusion	22	22	
		Vaporization	20,21	20,21	
TlI		Transition	33,34,35		
		Sublimation	20	20	
		Fusion	20,21,32	20,21,32	
		Vaporization	20	20	
Tl <sub>2</sub> S		Fusion	36,37,38	38	
Tl <sub>2</sub> S <sub>3</sub>		Fusion	39		
Tl <sub>2</sub> S <sub>5</sub>		Fusion	39		
Tl <sub>2</sub> SO <sub>4</sub>		Fusion	23,38	38	
Tl <sub>2</sub> Se		Fusion	36		
Tl <sub>2</sub> Se <sub>3</sub>		Transition	40		
Tl <sub>2</sub> Te <sub>3</sub>		Fusion	41		
TlNO <sub>2</sub>		Fusion	42		
TlNO <sub>3</sub>		Transition	43,44	43,44	
		Fusion	33,38,43,45,46	38	
		Vaporization	47		
TlPO <sub>3</sub>		Fusion	48		
TlH <sub>2</sub> PO <sub>2</sub>		Fusion	42		
TlH <sub>2</sub> PO <sub>3</sub>		Fusion	49		
TlH <sub>2</sub> PO <sub>4</sub>		Fusion	48		
TlAsS <sub>2</sub>		Fusion	37		
Tl <sub>3</sub> Bi <sub>5</sub>		Fusion	50		
(TlCl) <sub>3</sub> ·BiCl <sub>3</sub>		Fusion	51		
Tl <sub>2</sub> Co <sub>3</sub>		Transition	52		

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Table 30 (Continued)  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$Tl_2CO_3$		Fusion	23,38	38	
$Tl(C_2H_5)_3$		Fusion	53		
$TlCHO_2$	thallous formate	Fusion	54		
$TlC_2H_3O_2$	thallous acetate	Fusion	54		
$TlC_2F_3O_2$	thallous trifluoroacetate	Fusion	55		
		Vaporization	55		
$Tl_2CO_3 \cdot TlNO_3$		Fusion	46		
$Tl_2CO_3 \cdot 2TlNO_3$		Fusion	46		
$TlCl \cdot SnCl_2$		Fusion	56		
$3TlCl \cdot SnCl_2$		Fusion	56		
$TlCl \cdot 2PbCl_2$		Fusion	56		
$3TlCl \cdot PbCl_2$		Fusion	56		

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Table 31  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Zn		Fusion	1,2,3,4	1,2,3	
		Vaporization	2,5,6	2,5,6	
ZnO		Fusion	7		
ZnF <sub>2</sub>		Fusion	8		
		Vaporization	9	9	
ZnCl <sub>2</sub>		Fusion	10,11,12,13	12,13	
		Vaporization	12,13	12,13	
ZnCl <sub>2</sub> ·1 1/2 H <sub>2</sub> O		Fusion	14		
ZnCl <sub>2</sub> ·2H <sub>2</sub> O		Fusion	14		
ZnCl <sub>2</sub> ·2 1/2 H <sub>2</sub> O		Fusion	14		
ZnCl <sub>2</sub> ·3H <sub>2</sub> O		Fusion	14		
ZnBr <sub>2</sub>		Sublimation	13	13	
		Fusion	14,15		
		Vaporization	17		
ZnBr <sub>2</sub> ·2H <sub>2</sub> O		Fusion	14		
ZnBr <sub>2</sub> ·3H <sub>2</sub> O		Fusion	14		
ZnI <sub>2</sub>		Sublimation	13	13	
		Fusion	14,15,18		
ZnI <sub>2</sub> ·2H <sub>2</sub> O		Fusion	14		
ZnS		Sublimation	19	19	
		Transition	20		
ZnTe		Fusion	21		
Zn(NO <sub>3</sub> ) <sub>2</sub> ·H <sub>2</sub> O		Fusion	22		
Zn(NO <sub>3</sub> ) <sub>2</sub> ·2H <sub>2</sub> O		Fusion	22		
Zn(NO <sub>3</sub> ) <sub>2</sub> ·4H <sub>2</sub> O		Fusion	22		
Zn(NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O		Fusion	22		
2ZnCl <sub>2</sub> ·NH <sub>4</sub> Cl		Fusion	23		
ZnAs <sub>2</sub>		Fusion	24		
ZnSb		Fusion	25		
Zn <sub>3</sub> Sb <sub>2</sub>		Fusion	25,26,27	26	
Zn(CH <sub>3</sub> ) <sub>2</sub>		Vaporization	28	28	
Zn(C <sub>2</sub> H <sub>5</sub> ) <sub>2</sub>		Vaporization	28	28	
ZnBr <sub>2</sub> ·2CH <sub>3</sub> OH		Fusion	29		
ZnSiO <sub>3</sub>		Fusion	30		
Zn <sub>2</sub> SiO <sub>4</sub>		Fusion	7		
2TlCl·ZnCl <sub>2</sub>		Fusion	11		
TlCl·2ZnCl <sub>2</sub>		Fusion	11		

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9. Ruff and Boucher-1
10. Sandonnini-10
11. Korreng-1
12. Tarasenkova and Skulkova-1
13. Niwa-3
14. Dietz-1
15. Carnelley-2
16. Graetz-1
17. Meyer and Freyer-1
18. Hampe-1
19. Veselovskii-1
20. Allen and Crenshaw-2
21. Kobayashi-4
22. Ewing, McGovern, and Mathews-1
23. Hachmeister-1
24. Heike-1
25. Mönkemeyer-2
26. Dewitt and Seltz-1
27. Zemczuzny-3
28. Thompson and Linnett-2
29. Menschutkin-5
30. Stein-3

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

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Table 32  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Cd		Sublimation	1, 2, 3, 4, 5, 6, 7, 8, 9, 10, 11, 12, 13, 14, 15, 16, 17, 18, 19, 20	1	
		Fusion	1, 2, 3, 4, 5, 6, 7, 8, 9, 10, 11, 12, 13, 14, 15, 16, 17, 18, 19, 20	1, 6, 7, 17, 18, 19, 20	
		Vaporization	1, 4, 9, 11, 21, 22, 23, 24, 25, 26, 27, 28, 29, 30, 31, 32, 33	1, 9, 24, 26, 27, 28, 29, 30, 31, 32, 33	
CdO		Sublimation	1, 34, 35	1, 34, 35	
CdF <sub>2</sub>		Fusion	1, 36, 37	1	
		Vaporization	38	38	
CdCl <sub>2</sub>		Sublimation	39, 40, 41, 42	39, 40, 41, 42	
		Fusion	1, 43, 44, 45, 46, 47, 48, 49, 50, 51, 52, 53, 54, 55	1, 54	
		Vaporization	39, 40, 41, 42, 46, 54	39, 40, 41, 42	
CdBr <sub>2</sub>		Sublimation	42	42	
		Fusion	1, 45, 54, 55	1, 54	
		Vaporization	54		
CdI <sub>2</sub>		Sublimation	37, 42, 44, 45, 51, 52	42, 44	
		Fusion	37, 42, 44, 45, 51, 52	42, 44	
CdS		Sublimation	57, 58, 59	58, 59	
CdSO <sub>4</sub>		Fusion	51		
CdTe		Fusion	60		
Cd(NO <sub>3</sub> ) <sub>2</sub> ·4H <sub>2</sub> O		Fusion	61, 62	61	
2CdCl <sub>2</sub> ·NH <sub>4</sub> Cl		Fusion	46		
CdSb		Fusion	63	63	
Cd <sub>3</sub> Sb <sub>2</sub>		Fusion	15, 14		
Cd(HCOO) <sub>2</sub> ·3NH <sub>3</sub>	cadmium formate triammine	Fusion	64		
CdSiO <sub>3</sub>		Fusion	65		
TlCl·CdCl <sub>2</sub>		Fusion	48, 50		

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| 5. Kobayashi-3                | 11. Mott-1                     |
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14. Kurnakov and Konstantinov-2	56. Schmidt and Walter-1
15. Treitschke-1	57. Biltz-3
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17. Umino-3	59. Pogorelyi-1
18. Person-1	60. Kobayashi-3
19. Richards-1	61. Riesenfeld and Milchsack-2
20. Roos-2	62. Hasselblatt-1
21. von Leitgebel-1	63. Kubaschewski-1
22. Kordes and Raoz-1	64. Grossmann and Jaeger-1
23. Braune-1	65. Klooster-3
24. Richnow-1	
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27. Jenkins-3	
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30. Fogler and Rodebush-1	
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37. Carnelley-2	
38. Ruff and Boucher-1	
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42. Niwa-3	
43. Menge-1	
44. Bergman-2	
45. Hermann-2	
46. Hachmeister-1	
47. Sandonnini and Scarpa-1	
48. Sandonnini-12	
49. Aten-2	
50. Korreng-1	
51. Ruff and Plato-1	
52. Etard-1	
53. Helfenstein-1	
54. Weber-2	

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Table 33  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Hg		Sublimation	1,2,3,4,5	1,2,3,5	
		Fusion	1,2,3,4	1,2,3	
		Vaporization	5,7,8	5	
HgF <sub>2</sub>		Sublimation	9	9	
		Fusion	9		
HgCl <sub>2</sub>		Sublimation	1,10,11,12,13	1	
		Fusion	1,10,11,12,13	1	
		Vaporization	5,13,14	5	
HgBr <sub>2</sub>		Sublimation	5,11,13,15	3,5	
		Fusion	5,11,13,15	5	
		Vaporization	3,5,13,14	3,5	
HgI <sub>2</sub>		Transition	1,3,16,17,18,19	1,3	
		Sublimation	1,13,20	1	
		Fusion	1,13,20	1	
		Vaporization	5,13	5	
HgS		Transition	3,21,22,23	3,21,22	
HgI <sub>2</sub> ·2N <sub>2</sub> H <sub>5</sub> I·H <sub>2</sub> O		Fusion	24		
Hg(CH <sub>3</sub> ) <sub>2</sub>		Vaporization	25,26	25,26	
Hg(C <sub>2</sub> H <sub>5</sub> ) <sub>2</sub>		Vaporization	25	25	
Hg <sub>2</sub> Te		Fusion	27		
HgCl <sub>2</sub> ·TeCl <sub>4</sub>		Fusion	28		
HgCl <sub>2</sub> ·2TeNO <sub>3</sub>		Fusion	13		

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| 4. Endo-1                            | 18. Bridgman-11                       |
| 5. Kelley-19                         | 19. Shibata and Miwa-1                |
| 6. Kordes and Raoz-1                 | 20. Smits-1                           |
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| 9. Ruff and Bahlan-1                 | 23. Rinse-3                           |
| 10. Jonker-1                         | 24. Ferratine-1                       |
| 11. Carnelley-2                      | 25. Thompson-2                        |
| 12. Sandonnini-10,12                 | 26. Ratman-2                          |
| 13. Bergman-1                        | 27. Kurnakov and Puschin-1            |
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Table 34  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Cu		Fusion	1,2,3,4,5,6,7,8,9,10	1,2,3	
		Vaporization	11	11	
Cu <sub>2</sub> O		Fusion	1,12,13,14,15	1,12	
CuCl		Fusion	1,16	1	
CuBr		Fusion	17		
CuI		Fusion	16		
CuI <sub>2</sub>		Transition	1		
Cu <sub>2</sub> S		Transition	1	1	
		Fusion	1	1	
CuSO <sub>4</sub> ·5H <sub>2</sub> O		Transition	19		
Cu <sub>2</sub> Se		Transition	1	1	1
		Fusion	20		
Cu <sub>2</sub> Te		Fusion	21		
Cu <sub>4</sub> Te <sub>3</sub>		Transition	21		
		Fusion	21		
Cu(NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O		Fusion	22		
Cu <sub>3</sub> P		Fusion	23		
Cu <sub>3</sub> As		Fusion	6		
Cu <sub>5</sub> As <sub>2</sub>		Fusion	6		
Cu <sub>3</sub> Sb		Fusion	24,25		
Cu <sub>5</sub> Sb <sub>2</sub>		Fusion	26		
CuCN		Fusion	1		
Cu(CH <sub>3</sub> COO) <sub>2</sub> ·2NH <sub>3</sub>	cupric acetate diammine	Fusion	27		
Cu(CH <sub>2</sub> ClCOO) <sub>2</sub> ·2NH <sub>3</sub>	cupric chloroacetate diammine	Fusion	27		
Cu(CCl <sub>3</sub> COO) <sub>2</sub> ·5NH <sub>3</sub>	cupric trichloroacetate pentammine	Fusion	27		
CuBr <sub>4</sub> (CH <sub>2</sub> NH <sub>3</sub> ) <sub>2</sub>	ethylenediammonium cupric bromide	Fusion	27		
(C <sub>2</sub> H <sub>5</sub> ) <sub>3</sub> PO·CuCl <sub>2</sub>	triethylphosphine oxide cupric chloride	Fusion	28		
Cu <sub>5</sub> Si		Fusion	29		
4CuS·SiS <sub>2</sub>		Fusion	30		
Cu <sub>2</sub> Ga		Fusion	31		
Cu <sub>2</sub> Cd <sub>3</sub>		Fusion	32		



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Table 34 (Continued)  
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29. Arrhenius and Westgren-1
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32. Kubaschewski-1

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Table 35  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Ag		Fusion	1,2	1	
		Vaporization	3	3	
AgF		Fusion	4		
AgCl		Fusion	1,5,6,7,8	1,5	
		Vaporization	9,10	9,10	
AgClO <sub>3</sub>		Fusion	11		
AgBr		Transition	12		
		Fusion	1,13,14	1	
		Vaporization	15	15	
AgI		Transition	8,16,17,18	16,17,18	
		Fusion	1,19	1	
		Vaporization	3	3	
Ag <sub>2</sub> H <sub>3</sub> IO <sub>6</sub>		Transition	20,21	20	
Ag <sub>2</sub> S		Transition	22,23,24	23	
		Fusion	1,22	1	
Ag <sub>2</sub> SO <sub>4</sub>		Transition	25,26	25	
		Fusion	1,27,28	1	
Ag <sub>2</sub> SO <sub>4</sub> ·H <sub>2</sub> SO <sub>4</sub>		Transition	29		
Ag <sub>2</sub> Se		Transition	25	25	
AgTe		Transition	30		
Ag <sub>2</sub> Te		Transition	30,31		
		Fusion	30,32,34		
Ag <sub>3</sub> Te <sub>2</sub>		Transition	32		
AgNO <sub>3</sub>		Transition	25	25	25
		Fusion	1,34	1	1
AgPO <sub>3</sub>		Fusion	13		
Ag <sub>3</sub> PO <sub>4</sub>		Fusion	13		
Ag <sub>4</sub> P <sub>2</sub> O <sub>7</sub>		Fusion	13		
Ag <sub>2</sub> S·As <sub>2</sub> S <sub>3</sub>		Fusion	22		
3Ag <sub>2</sub> S·As <sub>2</sub> S <sub>3</sub>		Fusion	22		
Ag <sub>2</sub> S·Sb <sub>2</sub> S <sub>3</sub>		Fusion	22		
2Ag <sub>2</sub> S·Sb <sub>2</sub> S <sub>3</sub>		Fusion	22		
3Ag <sub>2</sub> S·Sb <sub>2</sub> S <sub>3</sub>		Fusion	22		
AgCN		Fusion	1	1	
Ag <sub>2</sub> HgI <sub>4</sub>		Transition	35	35	
AgNO <sub>3</sub> ·HgI <sub>2</sub>		Transition	11		
2AgNO <sub>3</sub> ·HgI <sub>2</sub>		Fusion	11		
Ag <sub>3</sub> Cu <sub>2</sub>		Fusion	36		

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Table 35 (Continued)  
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13. Carnelley-2
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17. Cohen and Joss-1
18. Mallard and LeChatelier-1
19. Germann and Metz-1
20. Stephenson and Adams-2
21. Stephenson-2
22. Jaeger and van Klooster-2
23. Kapustinskiĭ and Veselovskii-2
24. Sandonnini-9
25. Kelley-18
26. Nacken-1
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28. Marchal-5
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31. Pellini-1
32. Kracek and Ksanda-1
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35. Ketelaar-1
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Table 36  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Au		Fusion	1,2,3,4,5,6,7,8	1,2,7	1
		Vaporization	9,10,11,12,13,14	9,10,11,12,13,14	9,10,11,12
AuTe <sub>2</sub>		Fusion	15		
AuSb <sub>2</sub>		Transition	1,16	1,16	1
Au <sub>2</sub> Si		Fusion	17		
AuSn		Fusion	18,19	18	18
AuGa		Fusion	20		
AuGa <sub>2</sub>		Fusion	20		
AuZn		Fusion	18	18	18
AuCd		Fusion	18	18	18
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Table 37  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Pt		Fusion	1, 2, 3, 4, 5, 6, 7, 8	1	
		Vaporization	9, 10		
PtTe <sub>2</sub>		Fusion	11		
PtSb		Transition	12		
PtSb <sub>2</sub>		Fusion	12		
PtCl <sub>2</sub> ·CO		Fusion	13, 14		
PtCl <sub>2</sub> ·2CO		Fusion	13, 14		
PtCl <sub>4</sub> ·3CO		Fusion	13, 14		
PtBr <sub>2</sub> ·CO		Fusion	14		
PtI <sub>2</sub> ·CO		Fusion	15		
PtSi		Fusion	16		
Pt <sub>2</sub> Si		Transition	16		
		Fusion	16		

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12. Nemilow and Woronow-1	
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14. Pullinger-1	

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Table 38  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta G_p$
Formula	Description				
Ir		Fusion	1	3	
		Vaporization	2		
IrF <sub>6</sub>		Vaporization	3		
Ir <sub>2</sub> Cl <sub>5</sub> ·4(C <sub>2</sub> H <sub>5</sub> ) <sub>2</sub> S		Fusion	4		

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Table 39

Os		Fusion	1	3,4	
		Vaporization	1,2		
OsO <sub>4</sub>		Fusion	3,4		
		Vaporization	5		
OsF <sub>6</sub>		Vaporization	6	6	
OsF <sub>8</sub>		Fusion	6		
		Vaporization	6		
Os(CO) <sub>3</sub> Cl <sub>2</sub>		Fusion	7		

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Table 40  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Re		Fusion	1,2		
ReO <sub>3</sub>		Fusion	3		
ReO <sub>4</sub>		Fusion	4		
Re <sub>2</sub> O <sub>7</sub>		Fusion	4,5,6,7	5,7	
		Vaporization	5,7	5,7	
ReF <sub>4</sub>		Fusion	8		
ReF <sub>6</sub>		Fusion	9	9	
		Vaporization	9	9	
ReOF <sub>4</sub>		Sublimation	8	8	
		Fusion	8	8	
ReO <sub>3</sub> Cl		Fusion	10		
		Vaporization	10		
ReO <sub>2</sub> Cl <sub>3</sub>		Fusion	11		
ReOCl <sub>4</sub>		Fusion	10		
TlReO <sub>4</sub>		Fusion	12		
AgReO <sub>4</sub>		Fusion	12		
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Table 41  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Pd		Fusion	1,2,3,4,5,6,7,8,9,10,11	1,2	
		Vaporization	12,13,14		
PdCl <sub>2</sub>		Fusion	15	15	
PdS		Fusion	16		
PdSb		Fusion	17		
Pd <sub>3</sub> Sb		Transition	10,17,18	10,18	
		Fusion	17		

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| 5. Fairchild, Hoover and Peters-1 | 16. Weibke, Laar and Meisel-1 |
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| 7. Roeser and Wensel-1            | 18. Poppema-1                 |
| 8. Day and Sosman-1               |                               |
| 9. Waidner and Burgess-1          |                               |
| 10. Jaeger and Poppema-1          |                               |
| 11. Holborn and Valentiner-1      |                               |

Table 42

Rh		Fusion	1,2,3		
		Vaporization	4,5		

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Table 43  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Ru		Transition	1,2	1,2	
		Fusion	3		
		Vaporization	4		
RuO <sub>4</sub>		Fusion	5,6,7		
RuF <sub>5</sub>		Fusion	8		
		Vaporization	8		
Ru(CO) <sub>5</sub>		Fusion	9		
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Table 45  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Ni	nickel ethylenediammonium thiocyanate	Transition	1,2,3,4,5,6	1,2	1,2
		Fusion	1,2,7,8,9,10,11,12,13	1,2,9,13	1,2
		Vaporization	15,16,17	14,15,16	
NiO		Sublimation	18,19,20,21,22	18	
		Fusion	18,19,20,21,22		
NiCl <sub>2</sub>		Sublimation	16,23	16	
		Fusion	23		
NiBr <sub>2</sub>		Fusion	23		
NiI <sub>2</sub>		Fusion	23		
NiS		Fusion	24		
Ni <sub>2</sub> S		Fusion	25,26	25	
Ni <sub>3</sub> S <sub>2</sub>		Fusion	25,26,27	25	
NiSO <sub>4</sub> ·6H <sub>2</sub> O		Transition	28		
Ni(NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O		Fusion	29		
Ni(NO <sub>3</sub> ) <sub>2</sub> ·6NH <sub>3</sub>		Transition	30	30	
Ni <sub>2</sub> P		Fusion	31		
Ni <sub>5</sub> P <sub>2</sub>		Fusion	31		
NiSb		Fusion	32		
Ni <sub>5</sub> Sb <sub>2</sub>		Fusion	32		
Ni(CO) <sub>4</sub>		Fusion	33,34		
		Vaporization	16	1,16	
Ni(SCN) <sub>4</sub> ·C <sub>2</sub> H <sub>4</sub> (NH <sub>3</sub> ) <sub>2</sub>		Fusion	35		
NiSi		Fusion	36		
Ni <sub>2</sub> Si		Fusion	36		
NiZn <sub>3</sub>		Fusion	37,38,39		

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Table 45(Continued)  
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Table 46  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Co	cobalt ethylenediammonium thiocyanate	Transition	1,2,3,4	1,2,3,4	
		Fusion	5,6,7,8	5	
		Vaporization	9,10		
CoO		Fusion	11,12,13,14,15		
CoCl <sub>2</sub>		Fusion	5,16,17,18	5	
		Vaporization	19,20	19,20	
CoBr <sub>2</sub>		Fusion	21		
CoI <sub>2</sub>		Fusion	22,29		
CoS		Fusion	15		
Co(NO <sub>3</sub> ) <sub>2</sub> ·3H <sub>2</sub> O		Fusion	23		
Co(NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O		Fusion	23,24	24	
CoI <sub>3</sub> ·6NH <sub>3</sub>		Transition	25	25	
Co <sub>2</sub> P		Fusion	26		
CoSb		Fusion	8		
Co(SCN) <sub>4</sub> ·C <sub>2</sub> H <sub>4</sub> (NH <sub>3</sub> ) <sub>2</sub>		Fusion	27		
CoSi		Fusion	8		
CoSi <sub>2</sub>		Fusion	8		
CoSi <sub>3</sub>		Fusion	8		
Co <sub>2</sub> Si		Fusion	8		
Co <sub>2</sub> Sn		Fusion	8,28		

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Table 47  
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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Fe	wüstite	Transition	1,2,3,4,5,6	1,2,3,4,5,6	8
		Sublimation	1,5,7	1,5,7	
		Fusion	1,8,9,10,11,12,13	1,8	
		Vaporization	5,7,14	5,7,14	
Fe <sub>0.95</sub> O		Fusion	9	9	
Fe <sub>2</sub> O <sub>3</sub>		Transition	15,16		
Fe <sub>3</sub> O <sub>4</sub>		Fusion	17	17	
FeF <sub>2</sub>		Fusion	18	18	
FeCl <sub>2</sub>		Fusion	19,20,21,22		
		Vaporization	7	7	
FeCl <sub>3</sub>		Sublimation	7,23,24,25	7,23,24,25	
		Fusion	8,26	8	
		Vaporization	7	7	
FeCl <sub>3</sub> ·2H <sub>2</sub> O		Fusion	27		
FeCl <sub>3</sub> ·2 1/2 H <sub>2</sub> O		Fusion	27		
FeCl <sub>3</sub> ·3 1/2 H <sub>2</sub> O		Fusion	27		
FeCl <sub>3</sub> ·6H <sub>2</sub> O		Fusion	27		
FeCl <sub>3</sub> ·HCl·2H <sub>2</sub> O		Fusion	27		
FeCl <sub>3</sub> ·HCl·4H <sub>2</sub> O		Fusion	27		
FeCl <sub>3</sub> ·HCl·6H <sub>2</sub> O		Fusion	27		
FeBr <sub>2</sub>		Fusion	28		
FeI <sub>2</sub>		Fusion	29		
FeI <sub>2</sub> ·4H <sub>2</sub> O		Fusion	30		
FeI <sub>2</sub> ·9H <sub>2</sub> O		Fusion	30		
FeS		Transition	6,31,32	6	6
FeSO <sub>4</sub> ·7H <sub>2</sub> O		Fusion	33		
Fe(NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O		Fusion	34,35		
Fe(NO <sub>3</sub> ) <sub>3</sub> ·9H <sub>2</sub> O		Fusion	36		
FeCl <sub>3</sub> ·NH <sub>4</sub> Cl		Fusion	26		
NH <sub>4</sub> Fe(SO <sub>4</sub> ) <sub>2</sub> ·12H <sub>2</sub> O		Fusion	37		
Fe <sub>2</sub> P		Fusion	38		
Fe <sub>2</sub> O <sub>3</sub> ·P <sub>2</sub> O <sub>5</sub>		Fusion	39		
FeAs		Fusion	40		
Fe <sub>2</sub> As		Fusion	40		
Fe <sub>3</sub> Sb <sub>2</sub>		Fusion	41		
Fe(CO) <sub>5</sub>		Fusion	42,43,44,45	42,44	
		Vaporization	45,46,47	45,46,47	
Fe(CO) <sub>4</sub> H <sub>2</sub>		Fusion	48		
		Vaporization	48	48	
Fe(CO) <sub>2</sub> (NO) <sub>2</sub>		Sublimation	49	49	
		Fusion	49	49	

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Table 47 (Continued)  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$\text{Fe}(\text{CO})_2(\text{NO})_2$		Vaporization	49	49	
$\text{FeCl}_3 \cdot \text{CH}_3\text{NH}_3\text{Cl}$	methyammonium ferric chloride	Fusion			
$\text{FeCl}_3 \cdot (\text{CH}_3)_2\text{NH}_2\text{Cl}$	dimethylammionium ferric chloride	Fusion	50		
$\text{FeCl}_3 \cdot \text{C}_2\text{H}_5\text{NH}_3\text{Cl}$	ethylammonium ferric chloride	Fusion	50		
$\text{FeSi}$		Fusion	38,51		
$\text{FeO} \cdot \text{SiO}_2$	$\text{FeSiO}_3$	Fusion	52		
$2\text{FeO} \cdot \text{SiO}_2$	$\text{Fe}_2\text{SiO}_4$	Fusion	52		
$4\text{FeO} \cdot 3\text{SiO}_2$	$\text{Fe}_4\text{Si}_3\text{O}_{10}$	Fusion	52		
$\text{Fe}_2\text{O}_3 \cdot \text{PbO}$		Fusion	53		
$\text{FeGa}_2$		Fusion	54		
$\text{FeCl}_3 \cdot 2\text{TlCl}$		Fusion	55		
$\text{Fe}(\text{SO}_4)_2 \cdot 11 \cdot 12\text{H}_2\text{O}$		Fusion	37		
$\text{Fe}_2\text{O}_3 \cdot \text{ZnO}$		Fusion	53		
$\text{Fe}_2\text{O}_3 \cdot \text{CdO}$		Fusion	53		
$\text{Fe}_2\text{O}_3 \cdot \text{CuO}$		Fusion	53		
$2\text{FeCl}_3 \cdot \text{CuCl}_2$		Fusion	56		
$\text{Fe}_2\text{O}_3 \cdot \text{NiO}_3$		Fusion	53		
$2\text{FeS} \cdot \text{Ni}_2\text{S}$		Fusion	57		
$\text{Fe}_2\text{O}_3 \cdot \text{CoO}$		Fusion	53		
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Table 48  
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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Mn	manganese ethylenediammonium thiocyanate	Transition	1,2,3,4,5,6	1,2,3	
		Fusion	1,2,3,7	1	
		Vaporization	1,8,9,10,11	1,10	
MnO		Fusion	12		
Mn <sub>2</sub> O <sub>3</sub>		Transition	13		
Mn <sub>3</sub> O <sub>4</sub>		Transition	14	14	
		Fusion	15,16,17		
MnF <sub>2</sub>		Fusion	18		
MnCl <sub>2</sub>		Fusion	19,20,21,22,23,24	19	
		Vaporization	25	25	
MnBr <sub>2</sub>		Fusion	26,27		
MnI <sub>2</sub>		Fusion	28		
MnS		Fusion	29		
MnSO <sub>4</sub>		Fusion	30,31		
Mn(NO <sub>3</sub> ) <sub>2</sub> ·3H <sub>2</sub> O		Fusion	32,33	32	
Mn(NO <sub>3</sub> ) <sub>2</sub> ·4H <sub>2</sub> O		Fusion	34		
Mn(NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O		Fusion	33,34,35,36	35	
Mn <sub>5</sub> P <sub>2</sub>		Fusion	37		
Mn <sub>2</sub> Sb		Fusion	38		
Mn <sub>3</sub> C		Transition	39	39	
Mn(SCN) <sub>2</sub> · C <sub>2</sub> H <sub>4</sub> (NH <sub>3</sub> ) <sub>2</sub> (SCN) <sub>2</sub>		Fusion	40		
MnSiO <sub>3</sub>		Transition	41		
		Fusion	41,42	42	
Mn <sub>2</sub> SiO <sub>4</sub>		Fusion	41,43		
MnAu		Fusion	44		
MnPd		Fusion	7		
MnFe <sub>2</sub> O <sub>4</sub>		Fusion	45		

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Table 49  
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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Cr		Fusion	1,2,3,4,5,10	1,10	
Cr <sub>2</sub> O <sub>3</sub>		Transition	6,7		
		Fusion	8,9		
CrCl <sub>2</sub>		Sublimation	11,12	11,12	
		Fusion	11,12	11,12	
CrO <sub>2</sub> Cl <sub>2</sub>		Vaporization	13,14	13,14	
Cr <sub>8</sub> r <sub>2</sub>		Fusion	15		
CrI <sub>2</sub>		Fusion	16		
NH <sub>4</sub> Cr(SO <sub>4</sub> ) <sub>2</sub> ·12H <sub>2</sub> O		Fusion	17		
CrSb		Fusion	18		
Cr(CO) <sub>6</sub>		Sublimation	19,20	19,20	
CrSi		Fusion	21		
CrSi <sub>2</sub>		Fusion	21		
PbCrO <sub>4</sub>		Transition	22		
PbCrO <sub>4</sub> ·PbO		Fusion	22		
TlCr(SO <sub>4</sub> ) <sub>2</sub> ·12H <sub>2</sub> O		Fusion	17		
Pd <sub>2</sub> Cr <sub>3</sub>		Fusion	2		

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Table 50  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Mo		Fusion	1,2		
MoO <sub>3</sub>		Sublimation	3	3	
		Fusion	3,4,5,6		
		Vaporization	3,7	3,7	
MoF <sub>6</sub>		Sublimation	8,9,10	8,10	
		Fusion	8,9,10	8,10	
		Vaporization	8,9,10	8,10	
MoOF <sub>4</sub>		Fusion	9		
MoCl <sub>5</sub>		Fusion	9,11,12,13		
		Vaporization	9,11,12,13		
Bi <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>		Fusion	14		
MoC		Fusion	15,16		
Mo <sub>2</sub> C		Fusion	15,16		
Mo(CO) <sub>6</sub>		Sublimation	17	17	
PbMoO <sub>4</sub>		Fusion	6		
PbMoO <sub>4</sub> ·PbO		Fusion	6		
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Table 51  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
W		Fusion	1,2,3,4,5,6,7		
WO <sub>3</sub>		Sublimation	8	8	
		Fusion	9		
WF <sub>6</sub>		Transition	10,11	10,11	
		Fusion	10,11	10,11	
WOF <sub>4</sub>		Fusion	12		
		Vaporization	12		
WCl <sub>5</sub>		Fusion	13,14		
		Vaporization	13,14		
WCl <sub>6</sub>		Transition	15	15	
		Sublimation	15	15	
		Fusion	15	15	
		Vaporization	15	15	
WOCl <sub>4</sub>		Sublimation	12,14,16,17,18	16	
		Fusion	16	16	
		Vaporization	14,16	16	
WBr <sub>5</sub>		Fusion	14		
		Vaporization	14		
WOBr <sub>4</sub>		Fusion	14		
		Vaporization	14		
WCl <sub>6</sub> ·3WBr <sub>6</sub>		Fusion	19		
Bi <sub>2</sub> (WO <sub>4</sub> ) <sub>3</sub>		Fusion	20		
WC		Fusion	21,22		
W <sub>2</sub> C		Fusion	22		
W(CO) <sub>6</sub>		Sublimation	23	23	
PbWO <sub>4</sub>		Transition	9,20		
		Fusion	9,20		
PbWO <sub>4</sub> ·PbO		Fusion	9		
W <sub>2</sub> Re		Fusion	7		

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Table 51 (Continued)  
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Table 52  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
V		Fusion	1,2,3,4		
V <sub>2</sub> O <sub>3</sub>		Fusion	5		
V <sub>2</sub> O <sub>4</sub>		Transition	6	6	6
		Fusion	5,6	6	6
V <sub>2</sub> O <sub>5</sub>		Fusion	6	6	6
VF <sub>5</sub>		Sublimation	7		
VCl <sub>4</sub>		Fusion	8,9		
		Vaporization	7,8,9,10,11	8	
VOCl <sub>3</sub>		Fusion	12		
		Vaporization	10,12,13,14,15	12	
VN		Fusion	16,17		
NH <sub>4</sub> V(SO <sub>4</sub> ) <sub>2</sub> ·12H <sub>2</sub> O		Fusion	18		
VC		Fusion	19		
VS <sub>2</sub>		Fusion	20		
2PbO·V <sub>2</sub> O <sub>5</sub>		Fusion	21		
3PbO·V <sub>2</sub> O <sub>5</sub>		Fusion	21		
8PbO·V <sub>2</sub> O <sub>5</sub>		Fusion	21		
3Pb <sub>3</sub> (VO <sub>4</sub> ) <sub>2</sub> ·PbF <sub>2</sub>		Fusion	22		
3Pb <sub>3</sub> (VO <sub>4</sub> ) <sub>2</sub> ·PbCl <sub>2</sub>		Transition	23		
		Fusion	22,23		
TlV(SO <sub>4</sub> ) <sub>2</sub> ·12H <sub>2</sub> O		Fusion	18		
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9. Biltz and Keuncke-1			23. Eissner-1		
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Table 53  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Cb		Sublimation	1,2	2	
		Fusion	2		
		Fusion	3		
		Fusion	4,5		
		Fusion	6,7		
		Vaporization	6,7		
		Sublimation	8,9,10,11		
		Fusion	8,9,10,11		
		Vaporization	8,9,10,12		
		Fusion	13		
		Fusion	14,15		
		Fusion	16		

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Table 54  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Ta		Sublimation	1,2,3,4	1	
		Fusion	1,2,3,4		
Ta <sub>2</sub> O <sub>5</sub>		Fusion	4,5		
TaF <sub>5</sub>		Fusion	6	6	
		Vaporization	6		
TaCl <sub>5</sub>		Sublimation	7,8,9,10	7,8	
		Fusion	7,8,9,10	7,8	
		Vaporization	7,8,9,11	7,8	
TaBr <sub>5</sub>		Sublimation	12,13	12	
		Fusion	12,13	12	
		Vaporization	12,13	12	
TaI <sub>5</sub>		Fusion	14		
TaN		Fusion	15,16	15	
TaC		Fusion	15,17,18		
TaFe		Fusion	19		

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Table 55  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Ti		Transition	1,2,3		
		Sublimation	4	4	
		Fusion	5,6,7,8		
TiO		Transition	9	9	
TiO <sub>2</sub>		Transition	10,11,12,13	10,11	
		Fusion	14,15,16		
Ti <sub>2</sub> O <sub>3</sub>		Transition	9	9	
Ti <sub>3</sub> O <sub>5</sub>		Transition	9	9	
TiF <sub>4</sub>		Vaporization	17		
TiCl <sub>4</sub>		Fusion	18,19,20,21,22,23,24,	18,19,20	
		Vaporization	25,26,27,28,29	25,26	
		Transition	30		
TiBr <sub>4</sub>		Fusion	18,22,30,31,32,33	18	
		Vaporization	22,27,30,33		
		Fusion	22		
TiI <sub>4</sub>		Vaporization	34	34	
		Fusion	35		
TiCl <sub>4</sub> ·SCl <sub>4</sub>		Fusion	36,37		
TiN		Fusion	38		
TiCl <sub>4</sub> ·POCl <sub>3</sub>		Fusion	38		
		Vaporization	38		
TiC		Fusion	37,39		
TiFe <sub>2</sub>		Fusion	40		
FeTiO <sub>3</sub>		Fusion	41	41	
Fe <sub>2</sub> TiO <sub>4</sub>		Fusion	42		
MnTiO <sub>3</sub>		Fusion	43		

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Table 55(Continued)  
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Table 56  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Zr		Transition	1,2	1	
		Fusion	3		
ZrO		Transition	4		
ZrO <sub>2</sub>		Transition	6,7		
		Fusion	3,5,8,9,10	5	
ZrCl <sub>4</sub>		Sublimation	11,12	11,12	
		Fusion	12		
ZrBr <sub>4</sub>		Sublimation	11,12	11,12	
		Fusion	12		
ZrI <sub>4</sub>		Sublimation	11,12	11,12	
		Fusion	12		
ZrN		Fusion	13,14		
Zr(NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O		Fusion	15		
2ZrCl <sub>4</sub> ·PCl <sub>5</sub>		Fusion	3		
		Vaporization	3,17		
2ZrCl <sub>4</sub> ·POCl <sub>3</sub>		Vaporization	3,17		
ZrC		Fusion	18,19		
ZrSiO <sub>4</sub>		Fusion	20,21		
ZrCu <sub>3</sub>		Fusion	22		
ZrAu <sub>3</sub>		Fusion	22		
Zr <sub>2</sub> Fe <sub>3</sub>		Fusion	16		
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11. Kelley-19					
12. Rahlfs and Fischer-1					
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Table 57  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Hf		Transition	1		
		Fusion	1,2		
HfO <sub>2</sub>		Fusion	3,4		
HfCl <sub>4</sub>		Sublimation	5	5	
		Fusion	5		
HfBr <sub>4</sub>		Sublimation	5	5	
		Fusion	5		
HfC		Fusion	6,7		
REFERENCES					
1. Zwikker-3 2. de Boer and Fast-1 3. Clausing-1 4. Henning-5 5. Fischer, Gewehr, and Wingchen-1 6. Agte and Alterthum-1 7. Moers-2					

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Table 58  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
B		Fusion	1		
B <sub>2</sub> O <sub>3</sub>		Fusion	2,3	2	
		Vaporization	4	4	
B <sub>2</sub> H <sub>6</sub>	diborane	Fusion	5,6	6	
		Vaporization	6	6	
B <sub>4</sub> H <sub>10</sub>	tetraborane	Fusion	5		
		Vaporization	7	7	
B <sub>5</sub> H <sub>9</sub>	pentaborane	Fusion	5		
		Vaporization	7	7	
B <sub>5</sub> H <sub>11</sub>		Fusion	8		
		Vaporization	8	8	
B <sub>6</sub> H <sub>10</sub>	hexaborane	Fusion	5		
B <sub>10</sub> H <sub>14</sub>	decaborane	Sublimation	7	7	
		Fusion	9,10	10	
		Vaporization	7	7	
HBO <sub>2</sub>		Fusion	3		
BF <sub>3</sub>		Sublimation	11,12,13,14	11,12,13,14	
		Fusion	15	15	
		Vaporization	11,12,13,14	11,12,13,14	
BF <sub>3</sub> ·H <sub>2</sub> O		Fusion	16		
BF <sub>3</sub> ·2H <sub>2</sub> O		Fusion	17		
		Vaporization	17		
BCl <sub>3</sub>		Fusion	18		
		Vaporization	7	7	
B <sub>2</sub> H <sub>5</sub> Cl		Fusion	19		
BBr <sub>3</sub>		Fusion	1		
		Vaporization	7	7	
B <sub>2</sub> H <sub>5</sub> Br		Fusion	20		
		Vaporization	7	7	
BI <sub>3</sub>		Fusion	21		
		Vaporization	21		
B <sub>2</sub> S <sub>3</sub>		Fusion	22		
S <sub>2</sub> S <sub>5</sub>		Fusion	23		
8(HSO <sub>4</sub> ) <sub>3</sub>		Fusion	24		
SCl <sub>4</sub> ·BCl <sub>3</sub>		Fusion	22		
B <sub>3</sub> N <sub>3</sub> H <sub>6</sub>	hexahydro-s-triazatri-borine, borazole	Fusion	25		
		Vaporization	26,27	26,27	
B(CH <sub>3</sub> ) <sub>3</sub>	trimethylborine	Fusion	28		
		Vaporization	28	28	
(CH <sub>3</sub> ) <sub>2</sub> B <sub>2</sub> H <sub>4</sub>	unsym. dimethyldiborane	Fusion	29		
		Vaporization	29	29	

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Table 58 (Continued)  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$(CH_3)_3B_2H_3$	1,1,2-trimethyldiborane	Fusion	29		
		Vaporization	29	29	
$(CH_3)_4B_2H_2$	sym. tetramethyldiborane	Fusion	29		
		Vaporization	29	29	
$(C_2H_5)_2B_2H_4$	unsym. diethyldiborane	Vaporization	30	30	
$(C_2H_5)_4B_2H_2$	sym. tetraethyldiborane	Fusion	30		
$(CH_3O)_2BH$	dimethoxyborine	Fusion	31		
		Vaporization	31	31	
$B_3O_3(CH_3)_3$	trimethylboroxole	Fusion	32		
		Vaporization	32	32	
$B(CH_3)F_2$	methyldifluoroborine	Fusion	32		
		Vaporization	32	32	
$B(CH_3)_2F$	dimethylfluoroborine	Fusion	32		
		Vaporization	32	32	
$B(CH_3)Cl_2$	methyldichloroborine	Fusion	25		
		Vaporization	25	25	
$B(CH_3)_2Cl$	dimethylchloroborine	Vaporization	25	25	
$B(C_2H_5)_2Cl$	diethylchloroborine	Fusion	25		
		Vaporization	25	25	
$B(CH_3)Br_2$	methyldibromoborine	Fusion	25		
		Vaporization	25	25	
$B(CH_3)_2Br$	dimethylbromoborine	Fusion	25		
		Vaporization	25	25	
$B(C_2H_5)_2Br$	diethylbromoborine	Fusion	25		
		Vaporization	25	25	
$B(CH_3)_2I$	dimethyliodoborine	Fusion	25		
		Vaporization	25	25	
$BH_2N(CH_3)_2$	dimethylaminoborine	Fusion	25		
$B(CH_3)_2NH_2$	aminomethylborine	Fusion	25		
$B(CH_3)_2NH(CH_3)$	methylaminodimethylborine	Vaporization	33	33	
$BH_3 \cdot NH(CH_3)_2$	dimethylamine borine	Fusion	25		
$B(CH_3)_3 \cdot NH_2CH_3$	methylamine trimethylborine	Fusion	33		
$BH[N(CH_3)_2]_2$	bis(dimethylamino)borine	Fusion	25		
$B(CH_3)_3 \cdot NH(CH_3)_2$	dimethylamine trimethylborine	Fusion	25		
$B(CH_3)_2N(C_2H_5)_2$	diethylaminodimethylborine	Vaporization	25	25	
$B(C_2H_5)_2N(CH_3)_2$	dimethylaminodiethylborine	Vaporization	25	25	
$B[N(CH_3)_2]_3$	tris(dimethylamino)-borine	Fusion	34		
		Vaporization	34	34	
$[B(CH_3)_2NH(CH_3)]_2$	bis(methylaminodimethylborine)	Vaporization	33	33	
$[B(N(CH_3)_2)_3]_2$		Vaporization	25		

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$(CH_3)_2B_3N_3H_5$	1-methylborazole (N-methylborazole)	Vaporization	30,35	30,35	
$(CH_3)_2B_3N_3H_5$	2-methylborazole (B-methylborazole)	Fusion	30		
		Vaporization	30,35	30,35	
$(CH_3)_2B_3N_3H_4$	1,3-dimethylborazole	Vaporization	35	35	
$(CH_3)_2B_3N_3H_4$	2,4-dimethylborazole	Fusion	30		
		Vaporization	30	30	
$(CH_3)_2B_3N_3H_4$	1,2-dimethylborazole	Vaporization	35	35	
$(CH_3)_3B_3N_3H_3$	1,3,5-trimethylborazole	Fusion	25		
		Vaporization	25,35	35	
$(CH_3)_3B_3N_3H_3$	2,4,6-trimethylborazole	Fusion	25		
		Vaporization	25,30	30	
$(CH_3)_3B_3N_3H_3$	1,2,4-Trimethylborazole	Vaporization	35	35	
$(CH_3)_4B_3N_3H_2$	1,2,4,6-tetramethylborazole	Vaporization	35	35	
$(CH_3)_6B_3N_3$	hexamethylborazole	Fusion	25		
		Vaporization	33	33	
$(CH_3)_3B_3O_3 \cdot NH_2(CH_3)$	trimethylborazol methylamine	Fusion	25		
$BCl_2N(CH_3)_2$	dimethylaminodichloroborine	Fusion	25		
		Vaporization	34	34	
$BCl[N(CH_3)_2]_2$	bis(dimethylamino)-chloroborine	Fusion	34		
		Vaporization	34	34	
$[BCl_2N(CH_3)_2]_2$	bis(dimethylaminodichloroborine)	Fusion	34		
$BBr_2N(CH_3)_2$	dimethylaminodibromoborine	Fusion	25		
$[BBr_2N(CH_3)_2]_2$	bis(dimethylaminodibromoborine)	Fusion	25		
$PbO \cdot B_2O_3$		Fusion	36		
$PbO \cdot 2B_2O_3$		Fusion	36		
$PbO \cdot 3B_2O_3$		Fusion	36		
$2PbO \cdot 5B_2O_3$		Fusion	36		
$3ZnO \cdot 2B_2O_3$		Fusion	37		
$2CdO \cdot 3B_2O_3$		Fusion	36		
$Ni_2B$		Fusion	38		
$Ni_3B_2$		Transition	38		
		Fusion	38		
$Fe_2B$		Fusion	39		
$MnO \cdot B_2O_3$		Fusion	36		
$MnO \cdot 2B_2O_3$		Fusion	36		
$MnO \cdot 3B_2O_3$		Fusion	36		
$ZrB$		Fusion	40		
$Zr(BH_4)_4$	zirconium borohydride	Sublimation	41	41	

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Zr(BH <sub>4</sub> ) <sub>4</sub>	zirconium borohydride	Fusion	41	41	
		Vaporization	41	41	
HfB		Fusion	40		
Hf(BH <sub>4</sub> ) <sub>4</sub>		Sublimation	41	41	
		Fusion	41	41	
		Vaporization	41	41	

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Table 59  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Al		Fusion	1, 2, 3, 4, 5, 6, 7, 8	1, 2, 3, 4, 6, 8	
		Vaporization	9, 10, 11, 12, 13	9, 10, 11, 12, 13	
Al <sub>2</sub> O <sub>3</sub>		Fusion	3, 14, 15, 16, 17, 18, 19, 20, 21	3	
AlF <sub>3</sub>		Sublimation	22, 23, 24	22, 23, 24	
Al <sub>2</sub> Cl <sub>6</sub>		Sublimation	25	25	
		Fusion	3, 26, 27	3, 26	
Al <sub>2</sub> Br <sub>6</sub>		Fusion	26, 27, 28, 29, 30	26	
		Vaporization	25, 31	25, 31	
Al <sub>2</sub> I <sub>6</sub>		Fusion	26, 27, 32	26	
		Vaporization	25, 33	25, 33	
AlBr <sub>3</sub> ·H <sub>2</sub> S		Fusion	34		
Al <sub>3</sub> Se <sub>4</sub>		Fusion	35		
Al <sub>2</sub> Te <sub>3</sub>		Fusion	36		
AlCl <sub>3</sub> ·NH <sub>3</sub>		Fusion	32		
		Vaporization	37	37	
AlBr <sub>3</sub> ·NH <sub>3</sub>		Fusion	32		
NH <sub>4</sub> Al(SO <sub>4</sub> ) <sub>2</sub> ·12H <sub>2</sub> O		Transition	38	38	
		Fusion	39		
AlCl <sub>3</sub> ·PH <sub>3</sub>		Fusion	32		
AlCl <sub>3</sub> ·POCl <sub>3</sub>		Fusion	41		
AlBr <sub>3</sub> ·PH <sub>3</sub>		Fusion	32		
AlI <sub>3</sub> ·PH <sub>3</sub>		Fusion	32		
AlSb		Fusion	41		
AlBr <sub>3</sub> ·SbBr <sub>3</sub>		Fusion	29, 42		
AlBr <sub>3</sub> ·BiBr <sub>3</sub>		Fusion	29, 42		
Al(CH <sub>3</sub> ) <sub>3</sub>	trimethylaluminum	Fusion	43		
		Vaporization	44, 45	44, 45	
Al(C <sub>2</sub> H <sub>5</sub> ) <sub>3</sub>	triethylaluminum	Vaporization	45	45	
(CH <sub>3</sub> ) <sub>2</sub> O·Al(CH <sub>3</sub> ) <sub>3</sub>	trimethylaluminum di-methylether	Fusion	46		
		Vaporization	46	46	
[(CH <sub>3</sub> ) <sub>2</sub> AlO(CH <sub>3</sub> )] <sub>3</sub>	tris(dimethylaluminum methoxide)	Fusion	46		
		Vaporization	46	46	
[CH <sub>3</sub> AlCl <sub>2</sub> ] <sub>2</sub>	bis(methylaluminum dichloride)	Fusion	47		
[(CH <sub>3</sub> ) <sub>2</sub> AlCl] <sub>2</sub>	bis(dimethylaluminum chloride)	Fusion	46		
		Vaporization	46	46	
(CH <sub>3</sub> ) <sub>2</sub> O·Al(CH <sub>3</sub> ) <sub>2</sub> Cl	dimethylaluminum chloride dimethylether	Vaporization	46	46	
(C <sub>2</sub> H <sub>5</sub> ) <sub>2</sub> O·AlCl <sub>3</sub>	aluminum chloride diethylether	Fusion	34		
[(CH <sub>3</sub> ) <sub>2</sub> AlBr] <sub>2</sub>	bis(dimethylaluminum bromide)	Fusion	46		
		Vaporization	46	46	

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$(C_2H_5)_2O \cdot AlBr_3$	aluminum bromide diethylether	Fusion	34		
$(CH_3)_2S \cdot Al(CH_3)_3$	trimethylaluminum dimethyl sulfide	Fusion	46		
		Vaporization	46	46	
$[(CH_3)_2AlSCH_3]_2$	bis(methylmercaptodimethyl aluminum)	Fusion	46		
		Vaporization	46	46	
$AlH_3 \cdot N(CH_3)_3$	aluminumhydride trimethylamine	Fusion	47		
$Al(CH_3)_3 \cdot NH(CH_3)_2$	trimethylaluminum dimethylamine	Fusion	46		
		Vaporization	46	46	
$Al(CH_3)_3 \cdot N(CH_3)_3$	trimethylaluminum trimethylamine	Sublimation	46,47	46	
		Vaporization	46,47	46	
$AlH_3 \cdot 2N(CH_3)_3$	aluminum hydride di(trimethylamine)	Fusion	47		
$[(CH_3)_2AlN(CH_3)_2]_2$	bis(dimethylaminodimethylaluminum)	Sublimation	46	46	
		Fusion	46		
$AlCl_3 \cdot C_2H_5NH_2$	aluminum chloride ethylamine	Fusion	34		
$Al(CH_3)_2Cl \cdot (CH_3)_3N$	dimethylaluminum chloride trimethylamine	Sublimation	46	46	
		Fusion	46		
$(CH_3)_2PH \cdot Al(CH_3)_3$	trimethylaluminum dimethyl phosphine	Vaporization	46		
$(CH_3)_3P \cdot Al(CH_3)_3$	trimethylaluminum trimethylphosphine	Sublimation	46	46	
		Vaporization		46	
$[(CH_3)_2AlP(CH_3)_2]_3$	tris(dimethylphosphinodimethylaluminum)	Sublimation	46	46	
$Al_2O_3 \cdot SiO_2$		Fusion	58		
$Al_2Cl_6 \cdot SnCl_2$		Fusion	42		
$Al_2Cl_6 \cdot 2SnCl_2$		Fusion	42		
$Al_2Br_6 \cdot SnBr_2$		Fusion	42		
$Al_2Br_6 \cdot 2SnBr_2$		Fusion	42		
$Al_2Br_6 \cdot PbBr_2$		Fusion	42		
$AlGa$		Fusion	49		
$AlGa_2$		Fusion	49		
$Al_2Ga$		Fusion	49		
$Al_2Cl_6 \cdot 2TlCl$		Fusion	42		
$Al_2Br_6 \cdot TlBr$		Fusion	42		
$Al_2Br_6 \cdot 2TlBr$		Fusion	42		
$AlTl(SO_4)_2 \cdot 12H_2O$		Fusion	39		
$Al_2O_3 \cdot ZnO$		Fusion	50		
$Al_2Br_6 \cdot ZnBr_2$		Fusion	42		
$Al_2Br_6 \cdot CdBr_2$		Fusion	42		
$Al_2Br_6 \cdot HgBr_2$		Fusion	42		
$Al_2Br_6 \cdot Hg_2Br_2$		Fusion	42		

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Table 59 (Continued)					
SPECIFIC REFERENCES					
Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
AlCu		Fusion	51		
AlCu <sub>3</sub>		Fusion	51,52,53		
Al <sub>2</sub> Cu		Fusion	51		
Al <sub>2</sub> Cl <sub>6</sub> ·Cu <sub>2</sub> Cl <sub>2</sub>		Fusion	42		
AlAg <sub>2</sub>		Fusion	51,54		
AlAg <sub>3</sub>		Fusion	51,55		
Al <sub>2</sub> Br <sub>6</sub> ·2HgBr		Fusion	42		
AlAu <sub>2</sub>		Fusion	56		
Al <sub>2</sub> Au		Fusion	56		
NiAl		Fusion	52		
Al <sub>2</sub> O <sub>3</sub> ·NiO		Fusion	57		
CoAl		Fusion	52		
Al <sub>2</sub> O <sub>3</sub> ·CoO		Fusion	57		
FeAl <sub>3</sub>		Fusion	58,59		
Al <sub>3</sub> Ti		Fusion	60		
Al <sub>4</sub> Ti		Fusion	61		
Al <sub>2</sub> O <sub>3</sub> ·TiO <sub>2</sub>		Fusion	62		
Al <sub>2</sub> O <sub>3</sub> ·2TiO <sub>2</sub>		Fusion	57		
Al(BH <sub>4</sub> ) <sub>3</sub>	aluminum borohydride	Fusion	63		
		Vaporization	63	63	
Al(BH <sub>4</sub> ) <sub>3</sub> ·N(CH <sub>3</sub> ) <sub>3</sub>	aluminum borohydride trimethylamine	Fusion	63		
REFERENCES					
1. Richnow-1			15. Rankin and Merwin-1		
2. Awbery and Griffiths-1			16. von Wartenberg, Linde and Junge-1		
3. Kelley-21			17. Kanolt-2		
4. Kelley-25			18. Ruff, Seiferheld and Suda-1		
5. Roeser and Wensel-3			19. Ruff-6		
6. Bichowsky and Rossini-1			20. Weigel and Kaysser-1		
7. Vogel-4			21. Ruff and Schmidt-1		
B. Awbery-1			22. Ruff and Boucher-1		
9. Baur and Brunner-1			23. Olbrich-1		
10. Farkas-1			24. Naryushkin-1		
11. von Wartenberg-22			25. Keller-19		
12. Greenwood-2			26. Fischer-2		
13. Brewer et al-1			27. Blitz and Keunecke-1		
14. Bunting-3,5			28. Sugden-2		

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30. Puschin and Makuts-1
31. Zhuravlev-1
32. Klemm and Tanke-1
33. Moles and Vian-1
34. Nespital-1
35. Chikashige and Aoki-1
36. Chikashige and Nose-1
37. Klemm, Clausen, and Jacobi-1
38. Shomate-3
39. Locke-1
40. Casselman-1
41. Campbell and Mathews-1
42. Kendall, Crittenden, and Miller-1
43. Long and Norrish-1
44. Laubengayer and Gilliam-1
45. Bamford, Levi, and Newitt-1
46. Davidson and Brown-1
47. Klemm-1
48. Rankin and Wright-1
49. Puschin and Stojic-1
50. Bunting-4
51. Schübel-1
52. Gwyer-1
53. Carpenter and Edwards-1
54. Petrenko-1
55. Tischtschenko-1
56. Heycock and Neville-4
57. von Wartenberg and Reusch-1
58. Osawa-1
59. Biltz and Hasse-1
60. Manchot and Leber-1
61. Erckelens-1
62. Bunting-5
63. Schlesinger, Sanderson, and Burg-1

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Table 60  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Sc		Fusion	1		
		Vaporization	2		
ScCl <sub>3</sub>		Sublimation	3,4,5	3	
		Fusion	3,4,5		
ScBr <sub>3</sub>		Sublimation	3	3	
		Fusion	3		
ScI <sub>3</sub>		Sublimation	3	3	
		Fusion	3		
ScN		Fusion	6		

REFERENCES

1. Fischer, Brünger and Grieneisen-1
2. Richardson-6
3. Fischer, Gewehr and Wingchen-1
4. Biltz and Klemm-1
5. Voigt and Biltz-1
6. Friederich and Sittig-1

Table 61

Y		Fusion	1		
		Vaporization	2		
Y <sub>2</sub> O <sub>3</sub>		Vaporization	3		
YCl <sub>3</sub>		Fusion	4,5,6,7		
YBr <sub>3</sub>		Fusion	4,7		
YI <sub>3</sub>		Fusion	4,7		
Y <sub>2</sub> C <sub>3</sub>		Vaporization	3		
Y <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>		Fusion	8		

REFERENCES

1. Thompson, Holton, and Kremers-1
2. Richardson-6
3. Mott-1
4. Jantsch, Jawurek, Skalla, and Gawalowski-1
5. Klemm and Biltz-1
6. Kleinheksel and Kremers-1
7. Jantsch and Wein-1
8. Zambonini-1

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Table 62  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
LuCl <sub>3</sub>		Fusion	1,2		
LuI <sub>3</sub>		Fusion	1,2		

REFERENCES

1. Jantsch, Skalla, and Grubitsch-1
2. Jantsch and Wein-1

Table 63

YbCl <sub>3</sub>		Fusion	1,2		
YbBr <sub>3</sub>		Fusion	2		

REFERENCES

1. Jantsch, Skalla and Jawurek-1
2. Jantsch and Wein-1

Table 64

Tm		Vaporization	1		
TmCl <sub>3</sub>		Fusion	2,3,4		
TmI <sub>3</sub>		Fusion	4		

REFERENCES

1. Richardson-1
2. Kleinheksel and Kremers-1
3. Jantsch, Skalla, and Grubitsch-1
4. Jantsch and Wein-1

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Table 65  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
ErCl <sub>3</sub>		Fusion	1,2		
ErBr <sub>3</sub>		Fusion	1,2		
ErI <sub>3</sub>		Fusion	1,2		

REFERENCES

1. Jantsch and Wein-1
2. Jantsch, Jawurek, Skalla, and Gawalowski-1

Table 66

HoCl <sub>3</sub>		Fusion	1,2,3		
HoBr <sub>3</sub>		Fusion	1,3		
HoI <sub>3</sub>		Fusion	1,3		

REFERENCES

1. Jantsch, Jawurek, and Skalla-1
2. Kleinheksel and Kremers-1
3. Jantsch and Wein-1

Table 67

DyBr <sub>3</sub>		Fusion	1,2		
DyI <sub>3</sub>		Fusion	1,2		

REFERENCES

1. Jantsch, Jawurek, Skalla and Gawalowski-1
2. Jantsch and Wein-1

Table 68

TbCl <sub>3</sub>		Fusion	1		
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REFERENCES

1. Bourion-2

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Table 69  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
GdCl <sub>3</sub>		Fusion	1,2,3		
GdBr <sub>3</sub>		Fusion	1,2		
GdI <sub>3</sub>		Fusion	1,2		

REFERENCES

1. Jantsch, Jawurek, Skalla, and Gawalowski-1
2. Jantsch and Wein-1
3. Bourion-2

Table 70

EuCl <sub>3</sub>		Fusion	1,2		
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REFERENCES

1. Jantsch, Grubitsch, Hoffman and Alber-1
2. Jantsch and Wein-1

Table 71

Sm		Fusion	1		
SmCl <sub>2</sub>		Fusion	2		
SmCl <sub>3</sub>		Fusion	3,4,5,6,7,8		
SmBr <sub>3</sub>		Fusion	5,6		
SmI <sub>3</sub>		Fusion	4,6		
Sm <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>		Fusion	9		

REFERENCES

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Table 73  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Nd		Transition	1,2		
		Fusion	3		
NdF <sub>3</sub>		Fusion	4		
NdCl <sub>3</sub>		Fusion	5,6,7,8,9		
NdCl <sub>3</sub> ·6H <sub>2</sub> O		Fusion	10		
NdBr <sub>3</sub>		Fusion	9		
NdI <sub>3</sub>		Fusion	5,9		
Nd(MoO <sub>4</sub> ) <sub>3</sub>		Fusion	11		

REFERENCES

1. Jaeger, Bottema and Rosenbohm-3,4
2. Trombe-2
3. Muthmann and Weiss-1
4. von Wartenberg-2
5. Jantsch, Grubitsch, Hoffmann, and Alber-1
6. Kleinheksel and Kremers-1
7. Matignon-6,21
8. Bourion-2
9. Jantsch and Wein-1
10. Matignon-5
11. Zambonini-1

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Table 74  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta G_p$
Formula	Description				
Pr		Fusion	1,2	1	
PrF <sub>3</sub>		Fusion	3		
PrCl <sub>3</sub>		Fusion	4,5,6		
PrCl <sub>3</sub> ·7H <sub>2</sub> O		Fusion	6		
PrBr <sub>3</sub>		Fusion	7		
PrI <sub>3</sub>		Fusion	7,8		
PrSn <sub>3</sub>		Fusion	9		
Pr <sub>2</sub> Sn		Fusion	9		
Pr <sub>2</sub> Sn <sub>3</sub>		Fusion	9		
PrPb		Fusion	9		
PrPb <sub>3</sub>		Fusion	9		
Pr <sub>2</sub> Pb		Fusion	9		
PrGa <sub>2</sub>		Fusion	10		
PrTl		Fusion	9		
PrTl <sub>3</sub>		Fusion	9		
PrCu <sub>2</sub>		Fusion	9		
PrCu <sub>6</sub>		Fusion	9		
PrAg		Fusion	9		
PrAg <sub>3</sub>		Fusion	9		
PrAu		Fusion	9		
PrAu <sub>2</sub>		Fusion	9,11		
PrAu <sub>3</sub>		Fusion	9		
PrAu <sub>4</sub>		Fusion	11		
PrNi		Fusion	12		
PrNi <sub>5</sub>		Fusion	12		
Pr <sub>3</sub> Ni		Fusion	12		
Pr <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>		Fusion	13		
PrAl <sub>2</sub>		Fusion	9		

REFERENCES

1. Kelley-21	11. Rossi-1
2. Canneri and Rossi-3	12. Vogel and Fülling-1
3. von Wartenberg-2	13. Zambonini-4
4. Kleinheksel and Kremers-1	
5. Matignon-21	
6. Matignon-6	
7. Jantsch and Wein-1	
8. Jantsch, Grubitsch, Hoffmann, and Alber-1	
9. Rolla, Jandelli, Canneri, and Vogel-1	
10. Jandelli-1	

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Table 75  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Ce		Transition	1,2		
		Fusion	3,4,5,6	3	
Ce <sub>2</sub> O <sub>3</sub>		Fusion	7		
CeF <sub>3</sub>		Fusion	8		
CeCl <sub>3</sub>		Fusion	9,10	10	
CeBr <sub>3</sub>		Fusion	11		
CeI <sub>3</sub>		Fusion	11,12		
CeBi		Fusion	13		
Ce <sub>4</sub> Bi <sub>3</sub>		Fusion	14		
CeSn <sub>2</sub>		Fusion	15		
CeSn <sub>3</sub>		Fusion	4		
Ce <sub>2</sub> Sn		Fusion	4,15		
Ce <sub>2</sub> Sn <sub>3</sub>		Fusion	4,15		
CePb <sub>3</sub>		Fusion	4,13		
Ce <sub>2</sub> Pb		Fusion	4		
CeTi		Fusion	4		
CeTi <sub>3</sub>		Fusion	4		
CeZn <sub>9</sub>		Fusion	16		
CeCu <sub>2</sub>		Fusion	4		
CeCu <sub>6</sub>		Fusion	4		
CeAg		Fusion	4		
CeAg <sub>3</sub>		Fusion	4		
CeAu		Fusion	4		
CeAu <sub>2</sub>		Fusion	4		
CeAu <sub>3</sub>		Fusion	4		
CeNi		Fusion	17		
CeNi <sub>5</sub>		Fusion	17		
Ce <sub>3</sub> Ni		Fusion	17		
Ce <sub>3</sub> Co		Fusion	17		
Ce <sub>2</sub> O <sub>3</sub> •Cr <sub>2</sub> O <sub>3</sub>		Fusion	18		
Ce <sub>2</sub> (MoO <sub>4</sub> ) <sub>3</sub>		Fusion	19		
Ce <sub>2</sub> (WO <sub>4</sub> ) <sub>3</sub>		Fusion	19		
CeAl <sub>2</sub>		Fusion	4,20		
CeAl <sub>4</sub>		Transition	21		
Ce <sub>3</sub> Al		Fusion	21		

Table 75 (Continued)  
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2. Jaeger, Bottema, and Rosenbohm-1
3. Kelley-21
4. Rolla, Jandelli, Canneri, and Vogel-1
5. Cubiciotti-1
6. Vogel-9
7. Friederich and Sittig-3
8. von Wartenberg-11
9. Kleinheksel and Kremers-1
10. Brewer et al-1
11. Jantsch and Wein-1
12. Jantsch, Grubitsch, Hoffmann, and Alber-1
13. Iandelli-1
14. Vogel-6
15. Vogel-3
16. Schramm-1
17. Vogel and Füllung-1
18. von Wartenberg and Eckhardt-1
19. Zambonini-4
20. Vogel-4
21. Notwozny-1

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Table 76  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
La		Transition	1,2		
		Fusion	3,4,5		
La <sub>2</sub> O <sub>3</sub>		Fusion	6		
LaCl <sub>3</sub>		Fusion	7,8,9,10		
LaBr <sub>3</sub>		Fusion	8		
LaI <sub>3</sub>		Fusion	8,10		
La(NO <sub>3</sub> ) <sub>3</sub> ·6H <sub>2</sub> O		Transition	11		
		Fusion	12		
LaSn <sub>3</sub>		Fusion	5		
La <sub>2</sub> Sn		Fusion	5,13		
La <sub>2</sub> Sn <sub>3</sub>		Fusion	5,13		
LaPb		Fusion	5,13		
LaPb <sub>3</sub>		Fusion	5		
La <sub>2</sub> Pb		Fusion	5,13		
LaTl		Fusion	5,13,14		
LaTl <sub>3</sub>		Fusion	5,13		
LaCu <sub>2</sub>		Fusion	5,13		
LaCu <sub>6</sub>		Fusion	5		
LaZn <sub>9</sub>		Fusion	15		
LaAg		Fusion	5,13		
LaAg <sub>3</sub>		Fusion	5,13		
LaAu		Fusion	5,13		
LaAu <sub>2</sub>		Fusion	5,13		
LaAu <sub>3</sub>		Fusion	5,13		
LaNi		Fusion	16		
LaNi <sub>5</sub>		Fusion	16		
La <sub>3</sub> Ni		Fusion	16		
La(MoO <sub>4</sub> ) <sub>3</sub>		Fusion	17		
LaAl <sub>2</sub>		Fusion	5,18		
LaAl <sub>4</sub>		Transition	19		

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Table 79  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
PuF <sub>3</sub>		Sublimation	1,2	1,2	
		Vaporization	1,2	1,2	
PuCl <sub>3</sub>		Sublimation	1,2,3	1,2,3	
		Vaporization	1,2,3	1,2,3	
PuBr <sub>3</sub>		Sublimation	1,2	1,2	
		Vaporization	1,2	1,2	

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1. Brewer et al-1
2. Seaborg, Katz, and Manning-1
3. Weinstock-1

Table 80

NpF <sub>6</sub>		Fusion	1		
NpCl <sub>3</sub>		Fusion	1		
NpCl <sub>4</sub>		Fusion	1		
NpBr <sub>4</sub>		Fusion	1		
NpI <sub>3</sub>		Fusion	1		

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Table 81  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H^\circ$	$\Delta C_p$
Formula	Description				
U		Transition	1,2	1,2	
		Fusion	1		
UF <sub>4</sub>		Sublimation	1	1	
		Vaporization	1	1	
UF <sub>6</sub>		Sublimation	1,3,4	1,3	
		Fusion	4	4	
		Vaporization	1,4	1,4	
UCl <sub>3</sub>		Sublimation	1	1	
		Vaporization	1	1	
UCl <sub>4</sub>		Sublimation	1	1	
		Vaporization	1	1	
UBr <sub>3</sub>		Sublimation	1	1	
		Fusion	1	1	
		Vaporization	1	1	
UBr <sub>4</sub>		Sublimation	1	1	
		Vaporization	1	1	
UC <sub>2</sub>		Fusion	1	1	
U <sub>2</sub> C <sub>3</sub>		Fusion	1	1	
(UO <sub>2</sub> )Cl <sub>4</sub> (CH <sub>2</sub> NH <sub>3</sub> ) <sub>2</sub>	ethylene diammonium uranyl chloride	Fusion	5		
UO <sub>2</sub> (NO <sub>3</sub> ) <sub>4</sub> (CH <sub>2</sub> NH <sub>3</sub> ) <sub>2</sub>	diethylene diammonium uranyl nitrate	Fusion	5		
U(BH <sub>4</sub> ) <sub>4</sub>	uranium borohydride	Vaporization	6	6	
U(BH <sub>4</sub> ) <sub>3</sub> (BH <sub>3</sub> CH <sub>3</sub> )		Vaporization	6	6	
U(BH <sub>3</sub> CH <sub>3</sub> ) <sub>4</sub>		Vaporization	6	6	

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4. Brickwedde, Hoge and Scott-1
5. Grossmann and Schück-1
6. Schlesinger and Brown-1

Table 82

PaCl <sub>5</sub>		Fusion	1		
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REFERENCES

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Table 83  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Th		Transition	1		
		Fusion	2,3		
ThO <sub>2</sub>		Fusion	4		
ThCl <sub>4</sub>		Sublimation	5	5	
		Vaporization	5	5	
ThBr <sub>4</sub>		Sublimation	5	5	
		Vaporization	5	5	
ThI <sub>4</sub>		Fusion	5		
		Vaporization	5	5	
Th(BH <sub>4</sub> ) <sub>4</sub>		Sublimation	6	6	
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Table 85  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Be		Sublimation	1,2,3,4	1,2,3,4	
		Fusion	1,2,3,4	1,2,3,4	
BeO		Sublimation	5	5	
		Fusion	6		
BeCl <sub>2</sub>		Sublimation	7,8,9	7,8,9	
		Vaporization	7,8,9	7,8,9	
BeBr <sub>2</sub>		Sublimation	7,8	7,8	
		Fusion	7,10		
BeI <sub>2</sub>		Sublimation	7,8	7,8	
		Fusion	7		
Be <sub>4</sub> O(CH <sub>3</sub> CO <sub>2</sub> ) <sub>6</sub>		Sublimation	11	11	
		Transition	11	11	
		Fusion	11	11	
BeCl <sub>2</sub> ·2(C <sub>2</sub> H <sub>5</sub> ) <sub>2</sub> O		Fusion	12		
BeBr <sub>2</sub> ·2(C <sub>2</sub> H <sub>5</sub> ) <sub>2</sub> O		Fusion	12		
BeB <sub>2</sub> H <sub>8</sub>		Sublimation	13	13	
BeB <sub>2</sub> H <sub>8</sub> ·N(CH <sub>3</sub> ) <sub>3</sub>		Vaporization	13	13	
		Fusion	13		
3BeO·2ZrO <sub>2</sub>		Fusion	14		
BeO·Al <sub>2</sub> O <sub>3</sub>		Fusion	15		

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Table 86  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Mg		Fusion	1, 2	3	
		Vaporization	4, 5, 6, 7, 8	8	
MgO		Fusion	9	3	
MgF <sub>2</sub>		Fusion	10, 11	10	
		Vaporization	8, 12	8, 12	
MgCl <sub>2</sub>		Fusion	13, 14, 15, 16, 17	13, 18	
		Vaporization	19, 20, 21	19, 20	
MgCl <sub>2</sub> ·6H <sub>2</sub> O		Fusion	23, 24	23, 25, 22	
MgBr <sub>2</sub>		Fusion	26, 27	3	
MgBr <sub>2</sub> ·6H <sub>2</sub> O		Fusion	28		
MgSO <sub>4</sub>		Fusion	3		
Mg <sub>3</sub> N <sub>2</sub>		Transition	29		
Mg(NO <sub>3</sub> ) <sub>2</sub> ·2H <sub>2</sub> O		Fusion	30, 31, 32		
Mg(NO <sub>3</sub> ) <sub>2</sub> ·6H <sub>2</sub> O		Fusion	69, 30, 31, 32, 34	69	
Mg <sub>2</sub> P <sub>2</sub> O <sub>7</sub>		Transition	35		
		Fusion	36		
Mg <sub>3</sub> (PO <sub>4</sub> ) <sub>2</sub>		Fusion	3		
Mg <sub>3</sub> As <sub>2</sub>		Fusion	37		
Mg <sub>3</sub> Sb <sub>2</sub>		Transition	38		
3Mg(NO <sub>3</sub> ) <sub>2</sub> ·2Bi·24H <sub>2</sub> O		Fusion	39		
Mg(CH <sub>3</sub> COO) <sub>2</sub> ·4H <sub>2</sub> O		Fusion	40		
MgBr <sub>2</sub> ·6HCOOH		Fusion	41, 42		
MgBr <sub>2</sub> ·6CH <sub>3</sub> OH		Fusion	41		
MgBr <sub>2</sub> ·6CH <sub>3</sub> COOH		Fusion	41, 42		
MgBr <sub>2</sub> ·6C <sub>2</sub> H <sub>5</sub> OH		Fusion	41, 42		
MgI <sub>2</sub> ·6CH <sub>3</sub> OH		Fusion	41		
MgI <sub>2</sub> ·6CH <sub>3</sub> COOH		Fusion	41		
MgI <sub>2</sub> ·6C <sub>2</sub> H <sub>5</sub> OH		Fusion	41		
MgBr <sub>2</sub> ·3CH <sub>3</sub> CN		Fusion	43		
MgBr <sub>2</sub> ·6CH <sub>3</sub> CONH <sub>2</sub>		Fusion	43		
MgI <sub>2</sub> ·6CH <sub>3</sub> CONH <sub>2</sub>		Fusion	43		
Mg <sub>2</sub> Si		Fusion	44		
MgSiO <sub>3</sub>		Fusion	45, 46	3	
Mg <sub>2</sub> SiO <sub>4</sub>		Fusion	47		
Mg <sub>2</sub> Ge		Fusion	48		
Mg <sub>5</sub> Tl		Fusion	49		
MgZn <sub>2</sub>		Fusion	3, 50, 51	3	
MgCu <sub>2</sub>		Fusion	1, 51		
Mg <sub>2</sub> Cu		Fusion	1		
MgCuSb		Fusion	52		
MgAg		Fusion	51		

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Table 86 (Continued)  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta G_p$
Formula	Description				
MgNi <sub>2</sub>	boracite	Fusion	51		
MgO·Fe <sub>2</sub> O <sub>3</sub>		Fusion	53,54		
MgO·2TiO <sub>2</sub>		Fusion	55		
2MgO·TiO <sub>2</sub>		Fusion	55		
MgO·ZrO <sub>2</sub>		Fusion	56		
MgO·B <sub>2</sub> O <sub>3</sub>		Fusion	57		
2MgO·B <sub>2</sub> O <sub>3</sub>		Fusion	57		
3MgO·B <sub>2</sub> O <sub>3</sub>		Fusion	57		
MgCl <sub>2</sub> ·5MgO·7B <sub>2</sub> O <sub>3</sub>		Transition	58		
Mg <sub>2</sub> Al <sub>3</sub>		Fusion	59		
Mg <sub>4</sub> Al <sub>3</sub>		Fusion	60		
MgO·Al <sub>2</sub> O <sub>3</sub>		Fusion	61,62,63		
MgPr		Fusion	64		
Mg <sub>3</sub> Pr		Fusion	64		
MgCe		Fusion	65		
Mg <sub>3</sub> Ce		Fusion	64,65,66		
Mg <sub>4</sub> Ce		Fusion	64,65		
MgLa		Fusion	67		
Mg <sub>3</sub> La		Fusion	64,65		
MgO·La <sub>2</sub> O <sub>3</sub>		Fusion	68		

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Table 87  
SPECIFIC REFERENCES

Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Ca		Transition	1,2,3	12	2
		Sublimation	3,4,5	3,4,5	
		Fusion	1,6,7,8	8	
CaO		Fusion	9,10,11	9,10,11	
CaF <sub>2</sub>		Transition	2,12	2,12	2
		Fusion	2,12,13	2,12	2
		Vaporization	12	12	
CaCl <sub>2</sub>		Fusion	2,15,16,17,18,19	2,15	2
		Sublimation	20	20	
CaCl <sub>2</sub> ·6H <sub>2</sub> O		Fusion	59		
4CaCl <sub>2</sub> ·CaO		Fusion	17		
CaBr <sub>2</sub> ·6H <sub>2</sub> O		Fusion	60		
CaSO <sub>4</sub>		Transition	21		
		Fusion	8,21	8	
Ca <sub>3</sub> N <sub>2</sub>		Fusion	7		
Ca(NO <sub>3</sub> ) <sub>2</sub>		Fusion	8,22	8	
Ca(NO <sub>3</sub> ) <sub>2</sub> ·NH <sub>2</sub> O		Fusion	23		
CaO·P <sub>2</sub> O <sub>5</sub>		Fusion	24,25		
2CaO·P <sub>2</sub> O <sub>5</sub>		Fusion	24		
3CaO·P <sub>2</sub> O <sub>5</sub>		Transition	2	2	2
		Fusion	24		
		Fusion	25		
4CaO·P <sub>2</sub> O <sub>5</sub>		Fusion	26		
Ca <sub>3</sub> (AsO <sub>4</sub> ) <sub>2</sub>		Fusion	27		
Ca <sub>3</sub> Bi <sub>2</sub>		Fusion	27		
CaC <sub>2</sub>		Transition	2,28	2,28	2
CaC <sub>2</sub> ·CaO		Fusion	29		
CaCl <sub>2</sub> ·3CH <sub>3</sub> OH		Fusion	30		
CaCl <sub>2</sub> ·3C <sub>2</sub> H <sub>5</sub> OH		Fusion	30		
CaSi <sub>2</sub>		Fusion	31		
CaSiO <sub>3</sub>		Transition	32		
		Fusion	32,33,34,35,36,37		
Ca <sub>2</sub> SiO <sub>4</sub>		Transition	2,32,36,38,39	2	2
		Fusion	32,37		
CaSn <sub>3</sub>		Fusion	40		
CaPb <sub>3</sub>		Fusion	27,40,41		
Ca <sub>2</sub> Pb		Fusion	41		
CaTl		Fusion	27,41		
CaCl <sub>2</sub> ·TlCl		Fusion	16		
CaZn <sub>10</sub>		Fusion	40,42		
Ca <sub>2</sub> Zn <sub>3</sub>		Fusion	40,42		
CaCd <sub>3</sub>		Fusion	42		

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
CaCu <sub>4</sub>		Fusion			
CaAg		Fusion	41,43		
CaAg <sub>2</sub>		Fusion	43		
CaAg <sub>3</sub>		Fusion	41,43		
CaAg <sub>4</sub>		Fusion	43		
CaAu <sub>2</sub>		Fusion	44		
CaAu <sub>4</sub>		Fusion	44		
2CaO·3Fe <sub>2</sub> O <sub>3</sub>		Fusion	45		
CaO·9FeO		Fusion	46		
CaO·FeO·SiO <sub>2</sub>		Fusion	47		
CaCrO <sub>4</sub>		Fusion	48		
CaTiO <sub>3</sub>		Transition	2,49	2,49	2
CaO·TiO <sub>2</sub> ·SiO <sub>2</sub>		Fusion	50		
CaO·ZrO <sub>2</sub>		Fusion	51		
CaO·B <sub>2</sub> O <sub>3</sub>		Fusion	52,53	52	52
CaO·2B <sub>2</sub> O <sub>3</sub>		Fusion	2,52,53	2,52	2
2CaO·B <sub>2</sub> O <sub>3</sub>		Transition	2,52	2,52	2
		Fusion	52,53	52	2
3CaO·B <sub>2</sub> O <sub>3</sub>		Fusion	52,53	52	2
5CaO·B <sub>2</sub> O <sub>3</sub> ·SiO <sub>2</sub>		Fusion	54		
CaAl <sub>2</sub>		Fusion	55		
CaAl <sub>3</sub>		Fusion	56		
CaO·Al <sub>2</sub> O <sub>3</sub>		Fusion	32,36,37		
12CaO·7Al <sub>2</sub> O <sub>3</sub>		Fusion	36,37		
CaBr <sub>2</sub> ·AlBr <sub>3</sub>		Fusion	57		
CaO·Al <sub>2</sub> O <sub>3</sub> ·2SiO <sub>2</sub>		Fusion	36,37	8	8
2CaO·Al <sub>2</sub> O <sub>3</sub> ·SiO <sub>2</sub>		Fusion	36,37		
CaMg <sub>2</sub>		Fusion	41		
CaMg(SiO <sub>3</sub> ) <sub>2</sub>		Fusion	58		
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Table 88  
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Substance		Type of Process	Temperature °	$\Delta H$	$\Delta C_p$
Formula	Description				
Sr		Fusion	2,3,4,5	1	
		Vaporization	6	6	
SrO		Fusion	7		
SrF <sub>2</sub>		Fusion	1	1	
SrCl <sub>2</sub>		Fusion	8,9,10,11,12	1	
4SrCl <sub>2</sub> ·SrO		Fusion	8		
SrCl <sub>2</sub> ·SrF <sub>2</sub>		Fusion	13		
SrBr <sub>2</sub>		Fusion	14	1	
SrI <sub>2</sub>		Fusion	15		
SrSO <sub>4</sub>		Transition	16		
SrSO <sub>4</sub>		Fusion	16,17		
Sr(NO <sub>2</sub> ) <sub>2</sub>		Transition	18		
Sr(NO <sub>3</sub> ) <sub>2</sub>		Fusion	15		
Sr <sub>3</sub> (AsO <sub>4</sub> ) <sub>2</sub>		Fusion	19		
Sr <sub>3</sub> (PO <sub>4</sub> ) <sub>2</sub>		Fusion	1	1	
SrCO <sub>3</sub>		Transition	20		
		Fusion	21		
SrO·SiO <sub>2</sub>		Fusion	22,23		
2SrO·SiO <sub>2</sub>		Fusion	24		
SrPb <sub>3</sub>		Fusion	25		
SrAg		Fusion	2		
SrAg <sub>4</sub>		Fusion	2		
Sr <sub>3</sub> Ag <sub>2</sub>		Fusion	2		
Sr <sub>3</sub> Ag <sub>5</sub>		Fusion	2		
SrO·B <sub>2</sub> O <sub>3</sub>		Fusion	26		
SrO·2B <sub>2</sub> O <sub>3</sub>		Fusion	26		
2SrO·B <sub>2</sub> O <sub>3</sub>		Fusion	26		
SrO·Al <sub>2</sub> O <sub>3</sub>		Fusion	27		
SrMg <sub>2</sub>		Fusion	28		
SrMg <sub>9</sub>		Fusion	28		



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Table 88 (Continued)  
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2. Weibke-3
3. Hartmann and May-1
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6. Kelley-19
7. Schumacher-1
8. Neumann, Kröger, and Jüttner-1
9. Korreng-1
10. Sandonnini-10
11. Ferrari and Inganni-2
12. Sandonnini-5
13. Winter-2
14. Kellner-1
15. Carnelley-2
16. Grahmann-1
17. Amadori-3
18. Bureau-2
19. Guérin-1
20. Zimens-1
21. Boeke-4
22. Eskola-1
23. Jaeger and van Klooster-1
24. Stein-3
25. Iandelli-1
26. Guertler-2
27. von Wartenberg and Reusch-1
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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Ba		Transition	1		
		Sublimation	2,3	2	
		Fusion	1,3,4,5		
		Vaporization	6	6	
BaO		Sublimation	6,7	6,7	
		Fusion	8		
		Fusion	9	9	
Ba(OH) <sub>2</sub>		Fusion	10,11,12,12	10	
BaF <sub>2</sub>		Vaporization	14	14	
		Transition	16,17		
BaCl <sub>2</sub>		Fusion	10,13,16,17,18,19	10,19	
		Vaporization	15	15	
Ba(ClO <sub>4</sub> ) <sub>2</sub>		Transition	20		
		Fusion	10,13,21,22	10	
BaI <sub>2</sub>		Fusion	13		
BaSO <sub>4</sub>		Fusion	10,23	10	
		Transition	24		
Ba(NO <sub>3</sub> ) <sub>2</sub>		Fusion	10,22,25	10	
Ba <sub>3</sub> (PO <sub>4</sub> ) <sub>2</sub>		Fusion	10	10	
Ba <sub>3</sub> (AsO <sub>4</sub> ) <sub>2</sub>		Fusion	26		
BaCO <sub>3</sub>		Transition	27,28	27	
BaSiO <sub>3</sub>		Fusion	29,30		
BaO·2SiO <sub>2</sub>		Fusion	29,31		
BaAg <sub>4</sub>		Fusion	4		
Ba <sub>2</sub> Ag <sub>3</sub>		Fusion	4		
BaO·Fe <sub>2</sub> O <sub>3</sub>		Fusion	32		
Ba(MoO <sub>4</sub> )		Fusion	33		
Ba <sub>2</sub> V <sub>2</sub> O <sub>7</sub>		Fusion	22		
BaTiO <sub>3</sub>		Transition	34		
BaO·ZrO <sub>2</sub>		Fusion	35		
BaO·B <sub>2</sub> O <sub>3</sub>		Fusion	36		
BaO·2B <sub>2</sub> O <sub>3</sub>		Fusion	37		
BaO·3B <sub>2</sub> O <sub>3</sub>		Fusion	37		
BaO·4B <sub>2</sub> O <sub>3</sub>		Fusion	37		
2BaO·B <sub>2</sub> O <sub>3</sub>		Fusion	36		
3BaO·B <sub>2</sub> O <sub>3</sub>		Fusion	36		
BaO·Al <sub>2</sub> O <sub>3</sub>		Fusion	38		
BaMg <sub>2</sub>		Fusion	39		
BaMg <sub>9</sub>		Fusion	39		
BaBeF <sub>4</sub>		Fusion	40		
BaBe <sub>2</sub> F <sub>6</sub>		Fusion	40		

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15. Maier-1
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17. Sandonnini-1
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19. Gire and Puche-1
20. Vorländer and Kaascht-1
21. Kellner-1
22. Carnelley-2
23. Calcagni-1
24. Grahmann-1
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27. Lashchenko-1
28. Zimens-2
29. Eskola-1
30. Jaeger and von Klooster-1
31. Bowen-9
32. van Arkel, Verwey, and Van Brugger-1
33. Liebhafsky, Rochow, and Winslow-1
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35. von Wartenberg and Werth-1
36. Guertler-2
37. de Carli-4
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39. Klemm and Denkelacker-1
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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Ra		Fusion	1		
RaCl <sub>3</sub>		Fusion	2		
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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Li		Transition	1		
		Fusion	2,3,4,5,6	2,3	
		Vaporization	7		
LiOH		Fusion	8		
LiF		Fusion	2,8,9,10,11,12	2	
		Vaporization	13	13	
LiCl		Sublimation	14	14	
		Fusion	2,8,9,11,15,16,17,18,19	2	
		Vaporization	13	13	
LiClO <sub>3</sub>		Transition	20		
		Fusion	20,21		
LiClO <sub>3</sub> ·3H <sub>2</sub> O		Fusion	20		
LiClO <sub>4</sub>		Fusion	22		
LiBr		Fusion	2,9,16,23,24	2	
		Vaporization	13	13	
LiI		Fusion	8,9,16,23		
		Vaporization	13	13	
Li <sub>2</sub> SO <sub>4</sub>		Transition	18,25,26,27,28,29	25	
		Fusion	2,18,26,28,29,30	2	
LiNO <sub>3</sub>		Fusion	11,31,32,33,34,35,36,37	31	32
		Fusion	38,39	38	
Li <sub>3</sub> PO <sub>4</sub>		Fusion	40		
Li <sub>3</sub> Bi		Fusion	41		
Li <sub>2</sub> CO <sub>3</sub>		Fusion	11,18,28,42		
LiC <sub>2</sub> H <sub>3</sub> O <sub>2</sub>		Fusion	43,44		
LiC <sub>2</sub> H <sub>3</sub> O <sub>2</sub> ·2H <sub>2</sub> O		Fusion	44		
Li <sub>2</sub> SiO <sub>3</sub>		Fusion	2,45	2	
Li <sub>4</sub> SiO <sub>4</sub>		Fusion	2,45,46	2	
Li <sub>2</sub> GeO <sub>3</sub>		Fusion	47		
Li <sub>4</sub> GeO <sub>4</sub>		Fusion	47		
LiReO <sub>4</sub>		Fusion	48		
LiCl·CoCl <sub>2</sub>		Fusion	49		
Li <sub>2</sub> MoO <sub>4</sub>		Fusion	2,50	2	
Li <sub>2</sub> WO <sub>4</sub>		Fusion	2,50,51	2	
Li <sub>2</sub> W <sub>2</sub> O <sub>7</sub>		Transition	51		
		Fusion	50,51		
LiVO <sub>3</sub>		Fusion	9		
2LiSiO <sub>4</sub> ·3ZrSiO <sub>4</sub>		Fusion	52		
LiBO <sub>2</sub>		Fusion	2,11,45	2	

Table 91 (Continued)  
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Substance		Type of Process	Temperature	$\Delta H$	$\Delta G_p$
Formula	Description				
$\text{Li}_2\text{O} \cdot \text{B}_2\text{O}_3$		Fusion	53		
$\text{Li}_2\text{O} \cdot 2\text{B}_2\text{O}_3$		Fusion	53		
$\text{Li}_2\text{O} \cdot 3\text{B}_2\text{O}_3$		Fusion	53		
$\text{Li}_2\text{O} \cdot 4\text{B}_2\text{O}_3$		Fusion	53		
$\text{Li}_2\text{O} \cdot 5\text{B}_2\text{O}_3$		Fusion	53		
$3\text{LiF} \cdot \text{AlF}_3$		Fusion	54,55		
$\text{LiCl} \cdot \text{AlCl}_3$		Fusion	24		
$\text{LiBr} \cdot \text{AlBr}_3$		Fusion	24		
$\text{LiBr} \cdot 7\text{AlBr}_3$		Fusion	24		
$\text{Li}_2\text{BeF}_4$		Fusion	56,57		
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2. Kelley-21			31. Goodwin and Kalmus-1		
3. Binayendra-1			32. Kelley-25		
4. Bidwell-2			33. Wagner-3		
5. Losano-3			34. Puschin and Radoicic-1		
6. Bridgman-13			35. Harkins and Clark-1		
7. Gordon-4			36. Briscoe, Evans, and Robinson-1		
8. Scarpa-3			37. Lehrman and Breslau-1		
9. Schmitz-Dumont and Schmitz-1			38. Morgan and Benson-2		
10. Puschin and Baskow-1			39. Donnan and Burt-1		
11. Jaeger-3			40. Carnelley-2		
12. Tacchini-1			41. Iandelli-1		
13. Kelley-19			42. Amadori-3		
14. Niwa-1			43. Davidson and McAllister-1		
15. Sandonnini-4,11,18			44. Sidgwick and Gentle-1		
16. Ginnings and Phipps-1			45. van Klooster-2,3		
17. Hackmeister-1,2			46. Biltz and Lemke-1		
18. Hüttner and Tammann-1			47. Schwarz-2		
19. Richards and Meldrum-1			48. Smith and Long-1		
20. Kraus and Burgess-1			49. Ferrari and Baroni-1		
21. Bruhl-3			50. Hoerman-1		
22. Richards and Willard-1			51. van Liempt-8		
23. Sandonnini and Scarpa-4			52. Schwarz and Haacke-1		
24. Kendall, Crittenden, and Miller-1			53. Mazzetti and DeCarli-1		
25. Hare-1			54. Fedotieff and Timofeeff-1		
26. Nacken-1			55. Puschin and Baskow-1		
27. Ewles-1			56. Goldschmidt-3		
28. Amadori-8			57. Thilo and Lehmann-1		
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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Na		Fusion	1, 2, 3, 4, 5, 6, 7	1	1
		Vaporization	8, 9		
Na <sub>2</sub> O <sub>2</sub>		Fusion	10		
NaOH		Transition	11, 12, 13	11	
		Fusion	11, 12, 14	11, 12, 14	
NaOH·H <sub>2</sub> O		Fusion	15		
NaOH·3 1/2H <sub>2</sub> O		Fusion	15		
NaF		Sublimation	16, 17, 18, 19, 20, 21, 22, 23, 24, 25	16	
		Fusion	1, 14, 16, 17, 18, 19, 20, 21, 22, 23, 24, 25	1, 14	
		Vaporization	19	19	
NaCl		Sublimation	27, 28	27, 28	
		Fusion	1, 21, 23, 29, 30, 31, 32, 33, 34, 35, 36	1	1
		Vaporization	26	26	
NaClO <sub>3</sub>		Fusion	1, 37	1, 37	1
NaClO <sub>4</sub>		Transition	38, 39		
NaBr		Sublimation	16, 28	16, 28	
		Fusion	14, 21, 22, 23, 33, 34, 40, 41, 42	14	1
		Vaporization	26	26	
NaI		Fusion	14, 22, 23, 34, 40, 41, 43, 44	14	
		Vaporization	26	26	
Na <sub>2</sub> S		Fusion	45, 46, 47	46, 47	
Na <sub>2</sub> S <sub>2</sub>		Fusion	47, 48		
Na <sub>2</sub> S <sub>3</sub>		Fusion	47		
Na <sub>2</sub> S <sub>4</sub>		Fusion	45, 47, 49		
Na <sub>2</sub> S <sub>5</sub>		Fusion	47, 48		
Na <sub>4</sub> S <sub>3</sub>		Fusion	47		
Na <sub>4</sub> S <sub>5</sub>		Fusion	47		
Na <sub>4</sub> S <sub>7</sub>		Fusion	47		
Na <sub>4</sub> S <sub>9</sub>		Fusion	47		
Na <sub>2</sub> SO <sub>4</sub>		Transition	49, 50, 51, 52, 53, 54		
		Fusion	23, 34, 49, 52, 53, 54, 55		
Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> ·5H <sub>2</sub> O		Fusion	56, 57, 58	56, 57	
NaSH		Transition	50	50	
NaHSO <sub>4</sub>		Fusion	59		
NaSeH		Transition	50		
Na <sub>2</sub> Te		Transition	60		
NaNO <sub>2</sub>		Transition	61		

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
NaNO <sub>3</sub>		Transition	62		
		Fusion	1,42,37,63,64,65,66	1	1
NaNH <sub>2</sub>		Fusion	67,68,69		
NaPO <sub>3</sub>		Fusion	23,70		
Na <sub>4</sub> P <sub>2</sub> O <sub>7</sub>		Fusion	14,17	14	
NaH <sub>2</sub> PO <sub>3</sub>		Fusion	71		
NaBi		Fusion	72		
Na <sub>3</sub> Bi		Fusion	72,73		
Na <sub>2</sub> CO <sub>3</sub>		Transition	74		
		Fusion	14,74,75,76	14	
NaO <sub>2</sub> CH	sodium formate	Fusion	77,78		
NaC <sub>2</sub> H <sub>3</sub> O <sub>2</sub>	sodium acetate	Transition	79		
NaCN		Transition	80		
		Fusion	14	14	
		Vaporization	26	26	
NaCNS		Fusion	14	14	
Na <sub>2</sub> SiO <sub>3</sub>		Fusion	1,81,82	1	1
Na <sub>2</sub> SiO <sub>5</sub>		Fusion	14,81,82	14	
Na <sub>2</sub> GeO <sub>3</sub>		Fusion	83,84		
Na <sub>2</sub> Ge <sub>2</sub> O <sub>5</sub>		Fusion	84		
Na <sub>2</sub> Ge <sub>4</sub> O <sub>9</sub>		Fusion	84		
NaSn		Transition	72		
		Fusion	72		
Na <sub>2</sub> Sn		Fusion	72		
NaPb		Fusion	72,85,86		
Na <sub>2</sub> Pb		Fusion	72,86		
Na <sub>2</sub> Pb <sub>5</sub>		Fusion	72,86		
Na <sub>4</sub> Pb		Fusion	72,86,87		
Na <sub>5</sub> Pb <sub>2</sub>		Fusion	85		
NaCd <sub>2</sub>		Fusion	86		
NaCd <sub>5</sub>		Fusion	86		
NaReO <sub>4</sub>		Fusion	87		
Na <sub>2</sub> O•Fe <sub>2</sub> O <sub>3</sub>		Fusion	88		
Na <sub>2</sub> CrO <sub>4</sub>		Transition	89		
		Fusion	89		
Na <sub>2</sub> CrO <sub>4</sub> •10H <sub>2</sub> O		Fusion	90		
Na <sub>2</sub> Cr <sub>2</sub> O <sub>7</sub>		Fusion	91		
Na <sub>2</sub> MoO <sub>4</sub>		Transition	30	30	
		Fusion	14,23	14	
Na <sub>2</sub> WO <sub>4</sub>		Transition	92	92	
		Fusion	14,23,92	14,92	



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Substance		Type of Process	Temperature	$\Delta H$	$\Delta G_p$
Formula	Description				
$\text{Na}_2\text{TiO}_3$		Transition	1		
		Fusion	1,93	1	1
$\text{Na}_2\text{Ti}_2\text{O}_5$		Fusion	93		
$\text{Na}_2\text{Ti}_3\text{O}_7$		Fusion	93		
$\text{NaBO}_2$		Fusion	23		
$\text{Na}_2\text{O} \cdot \text{B}_2\text{O}_3$		Fusion	94		
$\text{Na}_2\text{O} \cdot \text{B}_2\text{O}_3 \cdot 8\text{H}_2\text{O}$		Fusion	94		
$\text{Na}_2\text{O} \cdot 2\text{B}_2\text{O}_3$		Fusion	95,96		
$\text{Na}_2\text{O} \cdot 4\text{B}_2\text{O}_3$		Fusion	95,96		
$2\text{NaO} \cdot \text{B}_2\text{O}_3$		Fusion	95		
$\text{Na}_3\text{AlF}_6$		Transition	1	1	1
		Fusion	1	1	1
$\text{NaAlSiO}_4$		Transition	97		
		Fusion	97		
$\text{Na}_2\text{SO}_4 \cdot \text{MgSO}_4$		Transition	98		
		Fusion	99		

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7. Grube and Schmidt-1	27. Zimm and Mayer-1
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11. von Hevesy-1	31. Korreng-1
12. Seward-2	32. Menge-1
13. Ewles-1	33. Sandonnini and Scarpa-3
14. Kelley-21	34. Hüttner and Tammann-1
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16. Niwa-1,6	36. Ray and Dayal-1
17. Carnelley-2	37. Goodwin and Kalmus-1
18. Moissan-3	38. Vorländer and Kaascht-1
19. Fedotieff and Iljinsky-1	39. Herrmann and Ilge-1
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45. Pearson and Robinson-2	87. Smith and Long-1
46. Friedrich-7	88. Knick and Kohlmeier-1
47. Friedrich-9	89. Hartford-1
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49. Perrier and Bellanca-1	91. Lehrman, Selditch, and Skell-1
50. Teichert-1	92. Goranson and Kracek-2
51. Calcagni and Marotta-3	93. Washburn and Bunting-1
52. Boeke-1	94. Cole, Scholes, and Amberg-1
53. Nacken-1	95. Morey and Merwin-1
54. Calgani-1	96. Ponomareff-1
55. Roberts-2	97. Bowen-10
56. Sturley-1	98. Speranskaya-1
57. Müller-15	99. Druzhinin-1
58. Tilden-2	
59. Ishikawa, Masuda, and Hagesawa-1	
60. Pellini and Quercigh-3	
61. Jaffray-1	
62. Kracek-6	
63. Puschin and Radoicic-1	
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65. Harkins and Clark-1	
66. Bergman-2	
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71. Partridge, Hicks, and Smith-1	
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75. Makarov and Schul'gina-1	
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78. Kendall and Adler-1	
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81. Morey and Bowen-1	
82. Kracek-5	

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
K		Fusion	1, 2, 3, 4, 5, 6, 7	1, 2	
		Vaporization	8		
KO <sub>2</sub>		Transition	9		
K <sub>2</sub> O <sub>2</sub>		Fusion	10		
K <sub>2</sub> O <sub>3</sub>		Fusion	10		
KOH		Transition	11, 12, 13	11	
		Fusion	11, 13, 14	11, 13, 14	
		Vaporization	19	19	
KOH·H <sub>2</sub> O		Fusion	16		
KF		Fusion	2, 17, 18, 19, 20, 21	2	2
		Vaporization	15	15	
KF·4H <sub>2</sub> O		Fusion	22		
KHF <sub>2</sub>		Transition	23	23	
		Fusion	23, 24	23	
KF·2HF		Fusion	24		
KF·3HF		Fusion	24		
KF·4HF		Fusion	24		
2KF·5HF		Fusion	24		
KCl		Sublimation	25, 26	25, 26	
		Fusion	2, 20, 21, 27, 28, 29, 30, 31, 32, 33, 34, 35, 36, 37	2, 27	2
		Vaporization	15, 38, 39	15, 38	
KClO <sub>4</sub>		Transition	40, 41, 42	40	
KBr		Sublimation	25, 26, 43, 44	25, 26, 43, 44	
		Fusion	15, 20, 25, 26, 30, 32, 33, 35, 36, 37, 43, 44, 45, 46, 47	15, 25, 26, 43, 44	
		Vaporization	15	15	
KI		Sublimation	26, 43, 44	26, 43, 44	
		Fusion	21, 30, 32, 33, 35, 36, 37, 48, 49, 50		
		Vaporization	15	15	
KI <sub>3</sub>		Fusion	51		
KIO <sub>3</sub>		Fusion	37		
KIBr <sub>2</sub>		Fusion	52		
K <sub>2</sub> S		Transition	53	53	
K <sub>2</sub> S <sub>5</sub>		Fusion	54		
K <sub>2</sub> SO <sub>4</sub>		Transition	27, 28, 35, 55, 56, 57, 58, 59, 60, 61	27, 56, 57	55
		Fusion	27, 28, 56, 57, 62, 63	56, 57	55
K <sub>2</sub> SO <sub>4</sub> ·3K <sub>2</sub> S		Fusion	62		
KSH		Transition	64	64	

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
KHSO <sub>4</sub>		Transition	65	65	
		Fusion	66		
K <sub>2</sub> SO <sub>4</sub> •3H <sub>2</sub> SO <sub>4</sub>		Fusion	66		
KSeH		Transition	64	64	
KN <sub>3</sub>		Fusion	67		
KNO <sub>2</sub>		Fusion	68,69		
KNO <sub>3</sub>		Transition	2,70,71,72,73	2,70,71	2
		Fusion	14,21,72,73,74,75,76	14,74	2
KNH <sub>2</sub>		Fusion	77,78		
KNO <sub>3</sub> •2HNO <sub>3</sub>		Fusion	79		
KPO <sub>3</sub>		Transition	28		
		Fusion	14,21,80	14	
K <sub>3</sub> PO <sub>4</sub>		Fusion	81	81	
K <sub>4</sub> P <sub>2</sub> O <sub>7</sub>		Transition	81		
		Fusion	14,28,81	14	
KH <sub>2</sub> PO <sub>4</sub>		Transition	82	82	
K <sub>2</sub> HPO <sub>4</sub> •6H <sub>2</sub> O		Fusion	83		
KPO <sub>3</sub> •KF		Transition	81		
		Fusion	81		
KH <sub>2</sub> AsO <sub>4</sub>		Transition	84	84	
KSb		Fusion	85		
K <sub>3</sub> Sb		Fusion	85		
KBi <sub>2</sub>		Fusion	86		
K <sub>2</sub> Bi		Fusion	87		
K <sub>3</sub> Bi		Fusion	87,88		
K <sub>3</sub> Bi <sub>2</sub>		Fusion	87		
K <sub>2</sub> CO <sub>3</sub>		Transition	35,89		
		Fusion	2,35,37,89,90,91	2	
KCHO <sub>2</sub>	potassium formate	Fusion	92		
KC <sub>2</sub> H <sub>3</sub> O <sub>2</sub>	potassium acetate	Fusion	93,94		
KCHO <sub>2</sub> •HCO <sub>2</sub> H		Fusion	92		
HC <sub>2</sub> H <sub>3</sub> O <sub>2</sub> •HC <sub>2</sub> H <sub>3</sub> O <sub>2</sub>		Fusion	93		
KCN		Transition	95	95	
		Fusion	14,49	14	
KCNS		Transition	96,97	96	
		Fusion	14,96,97,98,99,100	14,96,98	
K <sub>2</sub> SiO <sub>3</sub>		Fusion	101		
K <sub>2</sub> O•2SiO <sub>2</sub>		Fusion	101		
K <sub>2</sub> O•4SiO <sub>2</sub>		Transition	102	102	
		Fusion	101,102,103	102	

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
$K_2O \cdot GeO_2$		Fusion	104		
$K_2O \cdot 4GeO_2$		Fusion	104		
$2KF \cdot GeF_4$		Fusion	105		
$KPbI_3$		Fusion	48		
$K_2SO_4 \cdot 2PbSO_4$		Transition	106		
		Fusion	106		
KTl		Fusion	107		
KCd		Fusion	7		
$KReO_4$		Sublimation	108, 109, 110	109	
		Fusion	108, 109, 110	109	
$K_2SO_4 \cdot 2CoSO_4$		Fusion	111		
$KFe(SO_4)_2 \cdot 12H_2O$		Fusion	112		
$KCl \cdot MnCl_2$		Fusion	113		
$K_2SO_4 \cdot 2MnSO_4$		Fusion	114		
$K_2CrO_4$		Transition	27, 61, 115, 116, 117	27	
		Fusion	14, 61, 115, 116, 117	14	
$K_2Cr_2O_7$		Transition	118		
		Fusion	2, 21, 57, 61, 115, 116, 117, 118, 119, 120	2, 119	2
$K_2MoO_4$		Transition	59, 121, 61, 117, 122		
		Fusion	14, 21, 35, 59, 61, 117, 121, 122	14	
$K_2Mo_2O_7$		Fusion	61, 117		
$K_2O \cdot 3MoO_3$		Fusion	121		
$K_2WO_4$		Transition	59, 117, 121, 122, 123		
		Fusion	14, 21, 35, 59, 121, 123	14	
$K_2W_2O_7$		Fusion	61, 117		
$KVO_3$		Fusion	20, 124		
$KBO_2$		Fusion	21, 80	15	
$K_2O \cdot B_2O_3$		Fusion	125, 126		
$K_2O \cdot 2B_2O_3$		Fusion	125		
$K_2O \cdot 4B_2O_3$		Fusion	125		
$KBF_4$		Transition	127		
		Fusion	128		
$KBO_2 \cdot KPO_3$		Fusion	80		
$3KF \cdot AlF_3$		Fusion	17		
$KBr \cdot AlBr_3$		Fusion	129		
$KBr \cdot 2AlBr_3$		Fusion	129		
$KAl(SO_4)_2 \cdot 12H_2O$		Transition	130	130	
		Fusion	112, 131, 132	132	

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta G_p$
Formula	Description				
$KAlSi_2O_6$	lincite	Transition	133	2	
		Fusion	134		
$K_2SO_4 \cdot BeSO_4$		Fusion	106		
$KF \cdot MgF_2$		Fusion	135		
$KCl \cdot MgCl_2$		Fusion	2,136,137		
$2KCl \cdot MgCl_2$		Fusion	137		
$K_2SO_4 \cdot 2MgSO_4$		Fusion	106,138		
$KCl \cdot CaCl_2$		Fusion	136		
$K_2SO_4 \cdot 2CaSO_4$		Transition	136		
		Fusion	136		
$KCaPO_4$		Transition	139		
$K_2O \cdot CaO \cdot SiO_2$		Fusion	140		
$K_2O \cdot 3CaO \cdot 6SiO_2$		Fusion	140		
$2K_2O \cdot CaO \cdot 3SiO_2$		Fusion	140		
$2K_2O \cdot CaO \cdot 6SiO_2$		Fusion	140		
$4K_2O \cdot CaO \cdot 10SiO_2$		Fusion	140		
$KBr \cdot 2SrBr_2$		Fusion	141		
$2KBr \cdot SrBr_2$		Fusion	141		
$K_2SO_4 \cdot 2SrSO_4$		Transition	106		
$2KBr \cdot BaBr_2$		Fusion	141		
$K_2SO_4 \cdot Li_2SO_4$		Fusion	138		
$K_2MoO_4 \cdot Li_2MoO_4$		Transition	121		
		Fusion	121		
$K_2WO_4 \cdot Li_2WO_4$		Fusion	123		

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$
Formula	Description				
Rb		Transition	1		
		Sublimation	2	2	
		Fusion	1,3,4,5,6	3	
		Vaporization	2	2	
RbOH		Transition	7	7	
		Fusion	7,8	7,8	
RbF		Fusion	3,9	3,9	
		Vaporization	2	2	
RbF·HF		Fusion	10		
RbF·2HF		Fusion	10		
RbCl		Fusion	3,11,12,13		
		Vaporization	2,14	2,14	
RbBr		Fusion	3,14,15	3,14	
		Vaporization	2	2	
RbI		Fusion	3,15	3	
		Vaporization	2	2	
Rb <sub>2</sub> S <sub>2</sub>		Fusion	16		
Rb <sub>2</sub> S <sub>3</sub>		Fusion	16		
Rb <sub>2</sub> S <sub>5</sub>		Fusion	16		
Rb <sub>2</sub> S <sub>6</sub>		Fusion	16		
Rb <sub>2</sub> SO <sub>4</sub>		Transition	17,18		
		Fusion	15		
RbSH		Transition	19	19	
RbSeH		Transition	19		
RbN <sub>3</sub>		Fusion	20		
RbNO <sub>3</sub>		Transition	21,22		
		Fusion	3,15,21,22	3	
RbNO <sub>3</sub> ·HNO <sub>3</sub>		Fusion	23		
Rb <sub>2</sub> CO <sub>3</sub>		Fusion	24		
RbCHO <sub>2</sub>	rubidium formate	Fusion	25		
RbC <sub>2</sub> H <sub>3</sub> O <sub>2</sub>	rubidium acetate	Fusion	25		
RbCNS		Fusion	26		
RbBF <sub>4</sub>		Fusion	27		
3Rb·AlF <sub>3</sub>		Fusion	28		
RbNO <sub>3</sub> ·LiNO <sub>3</sub>		Fusion	22		
RbOH·2NaOH		Fusion	7		
2RbOH·KOH		Fusion	7		

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Substance		Type of Process	Temperature	$\Delta H$	$\Delta C_p$	
Formula	Description					
Cs		Sublimation	1	1		
		Fusion	2,3,4,5,6	2,3		
		Vaporization	1	1		
Cs <sub>2</sub> O		Fusion	7			
Cs <sub>7</sub> O		Fusion	7			
CsOH		Transition	8			
CsF		Fusion	8	8		
		Fusion	2,9,10	2,9		
CsF•HF		Fusion	11			
CsF•2HF		Fusion	11			
CsF•3HF		Fusion	11			
CsF•6HF		Fusion	11			
CsCl		Transition	12	12		
		Fusion	2,10,13,14	2		
		Vaporization	1	1		
CsBr		Fusion	1,10,15	1,15		
		Vaporization	1	1		
CsBrO <sub>3</sub>		Fusion	16			
CsI		Fusion	10			
		Vaporization	1	1		
Cs <sub>2</sub> S <sub>2</sub>		Fusion	17			
Cs <sub>2</sub> S <sub>3</sub>		Fusion	17			
Cs <sub>2</sub> S <sub>5</sub>		Fusion	17			
Cs <sub>2</sub> S <sub>6</sub>		Fusion	17			
Cs <sub>2</sub> SO <sub>4</sub>		Transition	18			
		Fusion	10			
CsN <sub>3</sub>		Fusion	19			
CsNH <sub>2</sub>		Fusion	20,21			
CsNO <sub>3</sub>		Transition	22,23			
	Fusion	1,10,22	1			
CsNO <sub>3</sub> •4H <sub>2</sub> O		Fusion	24			
CsNO <sub>3</sub> •HNO <sub>3</sub>		Fusion	25			
CsHCO <sub>2</sub>		Fusion	26			
CsHCO <sub>2</sub> •H <sub>2</sub> O	cesium formate	Transition	26			
		Fusion	26			
CsC <sub>2</sub> H <sub>3</sub> O <sub>2</sub>	cesium acetate	Fusion	26			
CsCl•2CuCl		Fusion	27			
		Fusion	28			
Cs <sub>2</sub> TiF <sub>6</sub>		Fusion	28			
CsBF <sub>4</sub>		Fusion	29			
3CsF•AlF <sub>3</sub>		Fusion	30			

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## VII. INDEX



As indicated in the Introduction, the formula of the given substance, written according to the standard arrangement, determines, first, the element under which the compound will be listed and, second, the position the given compound will occupy in that section.

The tables of values and of specific references for compounds occurring under given elements may be located from the following numbers referring to pages in this book.

Element	Tables of values		Specific references		Element	Tables of values		Specific references	
	Series I page	Series II page	Series I page	Series II page		Series I page	Series II page	Series I page	Series II page
Actinium					Molybdenum	293	707	928	1, 113
Aluminum	318	731	941	1, 127	Neodymium	344	752	957	1, 135
Americium					Neon	14	543	827	1, 028
Antimony	90	582	852	1, 051	Neptunium	356	764	963	1, 140
Argon	15	544	827	1, 028	Nickel	245	690	910	1, 104
Arsenic	85	579	850	1, 049	Nitrogen	52	561	843	1, 041
Astatine					Osmium	238	684	905	1, 100
Barium	413	787	984	1, 152	Oxygen	8	538	824	1, 026
Beryllium	368	770	968	1, 143	Palladium	210	687	907	1, 102
Bismuth	95	584	854	1, 053	Phosphorus	72	570	847	1, 045
Boron	313	721	939	1, 123	Platinum	234	682	902	1, 099
Bromine	27	550	832	1, 032	Plutonium	355	763	962	1, 140
Cadmium	190	670	885	1, 091	Polonium	51		842	
Calcium	386	778	971	1, 147	Potassium	483	803	1, 007	1, 161
Carbon	98	585	855	1, 054	Praseodymium	346	753	958	1, 136
Cerium	348	756	959	1, 137	Promethium	343		956	
Cesium	528	819	1, 021	1, 169	Protactinium		768		1, 141
Chlorine	21	549	830	1, 031	Radium	430	791	990	1, 154
Chromium	284	705	925	1, 112	Radon	18	547	828	1, 029
Cobalt	252	692	913	1, 106	Rhenium	239	685	906	1, 101
Columbium	302	714	932	1, 117	Rhodium	242	688	908	1, 102
(Niobium)					Rubidium	520	815	1, 018	1, 167
Copper	207	674	891	1, 094	Ruthenium	243	689	909	1, 103
Curium					Samarium	341	751	955	1, 134
Dysprosium	337	747	951	1, 133	Scandium	328	740	944	1, 131
Erbium	335	745	949	1, 133	Selenium	46	557	839	1, 038
Europium	340	750	954	1, 134	Silicon	147	629	869	1, 072
Fluorine	19	548	829	1, 030	Silver	222	677	896	1, 096
Francium					Sodium	447	796	995	1, 157
Gadolinium	339	749	953	1, 134	Strontium	402	784	980	1, 150
Gallium	166	657	877	1, 084	Sulfur	35	552	836	1, 034
Germanium	151	645	871	1, 078	Tantalum	303	715	933	1, 118
Gold	230	681	900	1, 098	Technetium	244		909	
Hafnium	312	720	938	1, 122	Tellurium	49	560	841	1, 040
Helium	13	542	827	1, 028	Terbium	338	748	952	1, 133
Holmium	336	746	950	1, 133	Thallium	173	662	879	1, 087
Hydrogen	9	539	825	1, 027	Thorium	363	769	966	1, 142
Indium	170	660	878	1, 086	Thulium	334	744	948	1, 132
Iodine	31	551	834	1, 033	Tin	154	649	872	1, 080
Iridium	237	683	904	1, 100	Titanium	304	716	934	1, 119
Iron	261	694	917	1, 107	Tungsten	296	709	930	1, 114
Krypton	16	545	828	1, 029	Uranium	357	765	964	1, 141
Lanthanum	351	760	961	1, 139	Vanadium	299	712	931	1, 116
Lead	158	653	874	1, 082	Xenon	17	546	828	1, 029
Lithium	431	792	991	1, 155	Ytterbium	333	743	947	1, 132
Lutetium	332	742	946	1, 132	Yttrium	330	741	945	1, 131
Magnesium	373	772	970	1, 144	Zinc	177	667	881	1, 089
Manganese	273	702	921	1, 110	Zirconium	308	718	936	1, 121
Mercury	199	672	888	1, 093					









