

NBSIR 73-338

(Lop 11773) 2, 1 page

Reference

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CALCULATED LIQUID PHASE THERMODYNAMIC PROPERTIES AND LIQUID-VAPOR EQUILIBRIA FOR FLUORINE-OXYGEN (FLOX) MIXTURES

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NAT'L INST. OF STAND & TECH



A11106 449260

NIST PUBLICATIONS

REFERENCE

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Handwritten: *auth.*

Cryogenics Division
Institute for Basic Standards
National Bureau of Standards
Boulder, Colorado 80302

September 1973

Final Report

Prepared for:

The National Aeronautics and Space Administration
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Cleveland, Ohio 44135

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456
no. 73-338
1973



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* The thermodynamic properties given in these tables are based on the thermochemical BTU = 1054.35J.

CALCULATED LIQUID PHASE THERMODYNAMIC PROPERTIES
AND LIQUID-VAPOR EQUILIBRIA
FOR FLUORINE-OXYGEN (FLOX) MIXTURES*

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ABSTRACT

Liquid phase thermodynamic properties and liquid-vapor equilibria of fluorine-oxygen mixtures, for which no experimental data exist, have been calculated. The results are based on excess properties predicted from the Snider-Herrington equations, with an adjusted combining rule, and the corresponding data for the pure fluids. Mixtures considered are 0.6, 0.7, 0.8, 0.88, and 0.9 weight fraction of fluorine from 55 to 90K up to 70×10^5 Pa. In the compressed liquid, molar volumes, enthalpy, entropy, and constant pressure specific heat were determined. Along the saturation boundary, coexistent vapor compositions and solution vapor pressures were determined as well. Corresponding properties of pure fluorine from experimental data have also been included. Results are tabulated in both British and S.I. units.

Key Words: Calculated properties; enthalpy; entropy; liquid; fluorine; fluorine-oxygen; mixtures; hard-sphere model; liquid-vapor equilibria; specific heat; volume.

1. INTRODUCTION

Liquid mixtures of fluorine and oxygen are potential candidates for oxidizers in chemical rocket propulsion systems. As such, there is a need for thermophysical properties of selected mixtures over a range of temperatures and pressures, for which experimental data do not exist. It is the purpose of this study to provide the "best" calculated mixture properties within the current state-of-the-art.

*This study was carried out at the National Bureau of Standards under the sponsorship of the National Aeronautics and Space Administration.

To predict the thermodynamic properties and phase equilibria of liquid mixtures reliably, it is essential at some point to introduce appropriate numerical data on the components of the mixture and on the nature of the unlike molecule interactions. The latter information is generally obtained by evaluating some experimental data on the fluid mixture, not necessarily on the property of interest, to account for inadequacies of the combining rules. Calculations can then be made on the property of interest, preferably in regions where some data are available to assess the degree to which the real mixture values are reproduced.

Recent work on hard-sphere perturbation models has resulted in at least one set of relatively simple equations for liquid mixtures, i. e., the Snider-Herrington equations [1], which represent the correct temperature dependence of thermophysical excess properties. Staveley [2] showed that results calculated from these equations compare much better with newer data than with some of the data originally used by the authors. This led Miller, Kidnay, and Hiza [3] to apply the Snider-Herrington equations to the infinite dilution Henry's constants for the neon-krypton system with marked success. In these latter calculations, the adjustment to account for the inadequacy of the combining rule for the energy parameters was consistent with the correlation of Hiza and Duncan [4], making it unnecessary to sacrifice experimental data to obtain information on the unlike molecule interactions. Miller [5] also demonstrated that for a number of other systems, the deviation parameters needed to predict mixture properties with the Snider-Herrington equations are consistent with those obtained from independent gas phase data, which for the most part are properly represented in magnitude by the Hiza-Duncan correlation.

Fluorine-oxygen mixtures are particularly intriguing in the above context, since appropriate data are not available to determine the inadequacies of the combining rules nor to determine the reliability of any calculations. However, accurate data for the pure fluids, fluorine [6] and oxygen[7], are available. If one can postulate something about the nature of the unlike molecule interactions, it is possible to obtain liquid fluorine-oxygen mixture properties from the pure fluid properties data and the excess properties determined with the Snider-Herrington equations.

Though the desired information for fluorine-oxygen interactions is not available, previous experience with comparable oxygen systems can be applied to this problem, since oxygen appears to be the responsible molecule in exceptions encountered in the correlation effort [4]. To model the fluorine-oxygen system after other simple oxygen systems, it must be assumed that fluorine is well-behaved, and that the mixtures formed are purely physical solutions.

In this study, the argon-oxygen system was used as the model to estimate an adjustment to the geometric mean combining rule for the characteristic energy parameter of the fluorine-oxygen system. Fluorine-oxygen liquid mixtures containing 0.6, 0.7, 0.8, 0.88, and 0.9 weight fraction of fluorine and pure fluorine were considered in the present calculations from the saturation boundary up to 70×10^5 Pa from 55 to 90K. The properties determined were molar volume, enthalpy, entropy, and constant pressure specific heat. Along the saturation boundary, the coexistent vapor phase compositions and solution vapor pressures were determined as well.

2. CALCULATION METHODS

The equations used in these calculations are well documented; thus, only the specific logic and unique features of the present calculations are included here.

The form of the equation of state of Longuet-Higgins and Widom [8], upon which the mixture equations of Snider and Herrington are based is:

$$Pv/RT = [(1 + \xi + \xi^2)/(1 - \xi)^3] - (a/RTv) \quad (1)$$

where $\xi = \beta/v$ and β is the hard-sphere volume, $(\frac{1}{6} N_A \pi r^3)$. The second term represents a uniform attractive potential field.

Assuming constant "a" and "β" parameters implies that the configurational $C_V = 0$. Recently, Blinowska, Herrington, and Staveley [9] have taken into account the fact that these parameters are not constant over a range of temperatures. To avoid evaluation of temperature and density dependencies in the present calculations, the parameters "a" and "β" were evaluated from the pure fluid properties at each temperature of interest. The "a" parameter was evaluated from the configurational internal energy (the configurational internal energy of an ideal gas is zero); thus, "β", the hard-sphere volume is the only unknown in equation (1).

In applying their equations, Snider and Herrington used the equimolar G^E (excess Gibbs energy) at one temperature to determine the value of the cross parameter, "a₁₂". Sacrificing mixture data in any form in the present calculations was not possible. Thus, a different scheme had to be devised.

In correlating deviations from the geometric mean combining rule for the characteristic energy parameters [4], the following empirical equation was developed:

$$k_{12} = 0.17(I_1 - I_2)^{1/2} \ln(I_1/I_2) \quad (2),$$

where I is the ionization potential in eV and the subscript one denotes the component with the larger value. The parameter k_{12} is defined as:

$$U_{o12} = (1 - k_{12}) (U_{o1} U_{o2})^{1/2} \quad (3)$$

where U_o is the characteristic potential energy parameter. It is anticipated that the deviation parameter k_{12} is always positive.

In comparing predictions of the deviation parameter from equation (2) with observed values for oxygen systems specifically, equation (2) consistently over-predicts the magnitude of k_{12} . This suggests that an effective ionization potential for oxygen, somewhat larger than the observed value of 12.08 eV [10], is necessary to make the predictions for oxygen systems consistent with experiment.

The ionization potential for fluorine is 15.70 - 15.83 eV [11], almost the same as that for argon, 15.79 - 15.94 eV [12]. If one assumes that, in the fluorine-oxygen system, fluorine will be well-behaved and oxygen can be blamed for the unusual behavior noted above, the argon-oxygen system can be used as a model system for the fluorine-oxygen system. This is precisely the assumption made in the present calculations.

In the combining rule used by Miller, Kidnay, and Hiza [3]

$$a_{12} = (1 - k_{12}) (a_{11} a_{22})^{1/2} \left[\frac{(r_1 + r_2)}{2(r_1 r_2)^{1/2}} \right]^3 \quad (4)$$

where "r" is the hard-sphere diameter, a k_{12} of 0.015 was selected for the fluorine-oxygen system by examination of the equimolar G^E from argon-oxygen liquid-vapor equilibria data [4].

Since the liquid-vapor equilibria data for argon-oxygen are well represented by the two-suffix Margules equation [13], this framework

was used here to establish the liquid phase total vapor pressure and the vapor composition of the fluorine-oxygen system. In this sense, the excess Gibbs energy and the activity coefficient of each component, γ_1 and γ_2 , are related to the liquid phase mole fractions, x_1 and x_2 , with a single constant, A, by

$$G^E = A x_1 x_2 \quad (5)$$

$$RT \ln \gamma_1 = A x_2^2 \quad (6)$$

$$RT \ln \gamma_2 = A x_1^2 \quad (7)$$

The standard expression for the activity coefficient of each component

$$RT \ln \gamma_1 = RT \ln (y_1 P / x_1 p_{o1}) + (B_{11} - v_{o1}) (P - p_{o1}) + P y_2^2 \delta_{12} \quad (8)$$

$$RT \ln \gamma_2 = RT \ln (y_2 P / x_2 p_{o2}) + (B_{22} - v_{o2}) (P - p_{o2}) + P y_1^2 \delta_{12} \quad (9)$$

can then be related to equations (6) and (7) to obtain two equations in two unknowns, i.e. the solution vapor pressure, P, and the vapor phase mole fraction, y_1 where ($y_1 + y_2 = 1$). B_{11} , v_{o1} , and p_{o1} , and the corresponding terms with a subscript two, are the pure component second virial coefficient, saturated liquid molar volume, and vapor pressure respectively. δ_{12} is an excess interaction second virial coefficient term defined by

$$\delta_{12} = 2B_{12} - B_{11} - B_{22} \quad (10).$$

The second virial coefficients were calculated using

$$B(T/T_c) = v_c \sum_{i=1}^5 B_i^* (T/T_c)^{(1-i)/4} \quad (11)$$

where B_i^* are the reduced form of the fitted polynomial coefficients B_i of pure fluorine second virial coefficients given by Prydz and Straty [6]. The relationship between the two sets of coefficients is

$$B_i^* = \frac{B_i(T_c)}{v_c} \quad (12).$$

Since the reduced polynomial coefficients for pure oxygen are essentially identical to those for fluorine, i.e. two parameter corresponding states is satisfied, the values of B_i^* for fluorine were used. To obtain B_{12} from equation (11), the following mixing rules were used

$$v_c^{1/3} = v_{c12}^{1/3} = \frac{1}{2} (v_{c1}^{1/3} + v_{c2}^{1/3}) \quad (13)$$

and

$$T_c = T_{c12} = (1 - k_{12}) (T_{c1} T_{c2})^{1/2} \quad (14)$$

with a value of k_{12} equal to 0.015.

The value of the constant A in equations (5), (6), and (7) was evaluated at each temperature from the equimolar G^E calculated from the Snider-Herrington equations. All pure component parameters were taken from the experimental data [6, 7].

3. RESULTS

The properties of interest were determined for the selected liquid phase compositions at each temperature in both British and S. I. units. The properties along the saturation boundary and those in the compressed liquid are given in S. I. units in tables A-1 and A-2, respectively. The corresponding properties in British units are given in tables B-1 and B-2. All calculations were done on a molar basis, though some of the properties are given on a weight basis for the convenience of the intended user. Fluorine and oxygen molecular weights of 38.00 and 32.00, respectively, were used in these conversions.

The mixture molar volumes, enthalpies, and entropies are the algebraic sum of the mole fraction average of the pure component properties taken from experiment, the excess property of the mixture calculated from the Snider-Herrington equations, and the ideal mixing contribution. This relationship is simply expressed as

$$M_m = x_1 M_1 + x_2 M_2 + M^E + M^{\text{ideal}} \quad (15)$$

where M is the property of interest, the subscript m denotes mixture, and the superscript E denotes excess. The last term, the ideal mixing contribution, is zero except for entropy and thus Gibbs energy, the ideal entropy of mixing being $-R\sum_i x_i \ln x_i$. In calculating the heat capacity of the mixture, it was assumed that the excess heat capacity is zero. The pure fluorine properties for the compressed liquid listed here have been corrected for a computational error in the original internal energy and enthalpy values given in reference 6; the saturated liquid properties were correct as given. Revised thermophysical properties tables for fluorine have been prepared by G. C. Straty and L. A. Weber and will be published soon. The reference state for the pure component thermodynamic properties is the ideal gas at zero K.

It is worth noting that the saturated liquid mixture densities (more correctly, specific weights), reported graphically by Schmidt [14] from "proportional interpolation" of the pure fluid properties, differ significantly from the values determined in this study. It is clear that Schmidt's values were determined by weight fraction averaging of the weights per unit volume of the pure saturated liquids. Thus, his mixture values are not consistent with the traditional definition of an ideal liquid mixture, i. e. the mole fraction average of the molar volumes. For a 70 weight % fluorine mixture at 140R, the saturation specific volume taken from Schmidt is approximately 2.2% lower than the corresponding value from this study (Table B-1). This difference can be accounted for in three parts (approximately): 0.4% due to the difference in pure fluid properties used, 1.4% due to the two different methods of averaging, and 0.4% due to nonideality considerations accounted for in the present study. If the same pure fluid properties were used, for this mixture, Schmidt's method of averaging results in a negative excess molar volume (equation 15) 1.4% of the ideal while the present calculations result in a positive excess volume 0.4% of the ideal.

The validity of the calculated results given here can only be assessed in the final analysis by comparison with appropriate experimental data, which are not available at the present time. At a minimum, two kinds of data at a single T - x point for the liquid mixture can satisfy this requirement. These are precise values of the solution vapor pressure and the saturated liquid excess molar volume for the equimolar liquid mixture at a single temperature, preferably far removed from the critical temperatures of the pure components. An additional piece of experimental data which is desirable, but not absolutely necessary, is the vapor phase composition in equilibrium with the equimolar liquid solution. Without

this information, it is hoped that these results are a significant improvement over values of the subject properties obtained by simple mole fraction averaging of the pure fluid (molar) properties.

4. ACKNOWLEDGMENTS

Support of this study by the National Aeronautics and Space Administration is gratefully acknowledged. The authors wish to thank G. C. Straty and L. A. Weber for providing computer programs for the pure fluid properties of interest.

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TABLE A-1. THERMODYNAMIC PROPERTIES OF LIQUID FLUORINE-OXYGEN MIXTURES ON THE SATURATION BOUNDARY

TEMPERATURE K	WT. FRACTION F ₂		PRESSURE 10 ⁵ PA	VOLUME CM ³ /MOL	ENTHALPY J/MOL	ENTROPY J/MOL-K	C _P J/MOL-K
	LIQUID	VAPOR					
55.0	0.6000	0.7562	0.00338	23.41	-5963.57	70.86	54.02
	0.7000	0.8146	0.00355	23.19	-5947.42	69.85	54.17
	0.8000	0.8724	0.00372	22.94	-5937.77	68.34	54.32
	0.8800	0.9204	0.00384	22.74	-5935.40	66.66	54.44
	0.9000	0.9329	0.00387	22.68	-5935.63	66.15	54.47
	1.0000	1.0000	0.00401	22.41	-5942.16	62.39	54.63
	60.0	0.6000	0.7524	0.01299	23.84	-5692.30	75.58
0.7000		0.8132	0.01367	23.61	-5675.20	74.59	54.16
0.8000		0.8726	0.01432	23.36	-5664.54	73.10	54.31
0.8800		0.9212	0.01481	23.15	-5661.32	71.43	54.43
0.9000		0.9338	0.01493	23.10	-5661.32	70.92	54.46
1.0000		1.0000	0.01550	22.92	-5666.69	67.18	54.62
65.0		0.6000	0.7468	0.03996	24.28	-5419.60	79.95
	0.7000	0.8100	0.04203	24.05	-5401.21	78.98	54.72
	0.8000	0.8714	0.04401	23.80	-5389.18	77.51	54.96
	0.8800	0.9211	0.04552	23.59	-5384.79	75.86	55.16
	0.9000	0.9337	0.04589	23.53	-5384.56	75.36	55.21
	1.0000	1.0000	0.04769	23.25	-5388.31	71.64	55.46
	70.0	0.6000	0.7492	0.10326	24.75	-5145.50	84.01
0.7000		0.8059	0.10853	24.51	-5125.62	83.07	54.51
0.8000		0.8694	0.11356	24.26	-5112.01	81.62	54.70
0.8800		0.9203	0.11744	24.05	-5105.29	79.99	54.85
0.9000		0.9332	0.11839	24.00	-5105.65	79.49	54.89
1.0000		1.0000	0.12302	23.71	-5107.68	75.80	55.10
75.0		0.6000	0.7333	0.23246	25.25	-4869.73	87.81
	0.7000	0.8014	0.24404	25.01	-4848.26	86.89	54.94
	0.8000	0.8670	0.25514	24.76	-4832.98	85.47	55.18
	0.8800	0.9192	0.26373	24.55	-4825.87	83.85	55.38
	0.9000	0.9323	0.26583	24.49	-4824.88	83.36	55.43
	1.0000	1.0000	0.27612	24.20	-4825.07	79.69	55.69
	80.0	0.6000	0.7257	0.46837	25.78	-4592.60	91.39
0.7000		0.7971	0.49116	25.55	-4569.39	90.49	55.20
0.8000		0.8646	0.51308	25.30	-4552.15	89.09	55.45
0.8800		0.9180	0.53008	25.08	-4543.54	87.49	55.67
0.9000		0.9315	0.53426	25.02	-4542.16	87.01	55.72
1.0000		1.0000	0.55473	24.73	-4540.38	83.36	56.00
85.0		0.6000	0.7208	0.86232	26.35	-4313.57	94.75
	0.7000	0.7932	0.90342	26.12	-4298.44	93.88	55.97
	0.8000	0.8625	0.94315	25.87	-4269.34	92.50	56.31
	0.8800	0.9170	0.97407	25.65	-4259.04	90.93	56.59
	0.9000	0.9307	0.98168	25.60	-4257.09	90.44	56.67
	1.0000	1.0000	1.01909	25.30	-4253.20	86.82	57.03
	90.0	0.6000	0.7154	1.47487	26.97	-4031.98	97.96
0.7000		0.7835	1.54399	26.73	-4004.77	97.11	56.96
0.8000		0.8605	1.61112	26.48	-3983.51	95.76	57.40
0.8800		0.9161	1.66359	26.27	-3971.43	94.20	57.76
0.9000		0.9300	1.67654	26.21	-3969.16	93.72	57.86
1.0000		1.0000	1.74034	25.91	-3962.81	90.12	58.33
95.0		0.6000	0.7101	2.37502	27.64	-3746.61	101.03
	0.7000	0.7850	2.48455	27.41	-3717.20	100.20	57.63
	0.8000	0.8594	2.59143	27.16	-3693.67	98.87	58.10
	0.8800	0.9150	2.67529	26.95	-3679.70	97.33	58.49
	0.9000	0.9291	2.69604	26.89	-3676.96	96.86	58.58
	1.0000	1.0000	2.79849	26.59	-3668.18	93.28	59.09

TABLE A-2. THERMODYNAMIC PROPERTIES OF FLUOPIPINE-OXYGEN MIXTURES IN THE COMPRESSED LIQUID REGION

WT. FRACTION F2 =0.60

PRESSURE 10 ⁵ PA	TEMPERATURE K	VOLUME CM ³ /MOL	ENTHALPY J/MOL	ENTROPY J/MOL-K	C _P J/MOL-K	TEMPERATURE K	VOLUME CM ³ /MOL	ENTHALPY J/MOL	ENTROPY J/MOL-K	C _P J/MOL-K
0.50000	55.0	23.41	-5362.54	70.86	54.37	55.0	23.41	-5961.58	70.85	54.36
	60.0	23.44	-5591.34	75.58	54.78	60.0	23.44	-5673.62	75.50	54.71
	65.0	24.25	-5419.33	79.94	55.79	65.0	24.26	-5401.07	79.86	55.82
	70.0	25.25	-5145.38	87.81	58.27	70.0	25.22	-5128.70	87.91	55.70
5.00000	55.0	23.82	-5954.06	70.82	54.30	55.0	23.81	-5944.62	70.78	54.23
	60.0	24.27	-5683.02	75.90	54.75	60.0	24.26	-5673.62	75.50	54.71
	65.0	25.24	-5137.10	83.96	55.81	65.0	25.22	-5128.70	83.91	55.82
	70.0	26.35	-4584.94	91.34	58.32	70.0	26.31	-4576.41	91.28	55.36
20.00000	55.0	26.92	-4306.36	94.71	58.65	55.0	26.92	-4293.81	94.65	58.60
	60.0	27.62	-3744.33	100.99	57.22	60.0	27.59	-3736.20	100.92	57.11
	65.0	33.79	-5325.74	70.70	54.10	65.0	33.77	-5307.02	70.52	54.06
	70.0	24.09	-5655.08	79.41	55.65	70.0	24.21	-5636.30	79.33	55.80
40.00000	55.0	25.18	-4933.14	87.61	54.68	55.0	25.17	-4917.92	87.52	54.65
	60.0	26.26	-4559.37	91.17	55.33	60.0	26.22	-4544.69	91.07	55.30
	65.0	27.52	-4282.31	97.78	56.08	65.0	27.45	-4268.04	97.65	56.04
	70.0	37.93	-5325.74	70.70	54.10	70.0	37.93	-5307.02	70.52	54.06
60.00000	55.0	23.75	-5945.15	70.54	53.97	55.0	23.73	-5926.60	70.46	53.93
	60.0	24.19	-5675.31	75.20	54.50	60.0	24.16	-5656.14	75.11	54.46
	65.0	25.12	-5129.38	83.63	55.62	65.0	25.09	-5109.68	83.54	55.58
	70.0	26.18	-4579.19	90.46	56.30	70.0	26.13	-4559.92	90.36	56.22
20.00000	55.0	26.73	-4247.35	94.31	56.69	55.0	26.71	-4228.74	94.20	56.65
	60.0	27.39	-3792.57	100.51	55.30	60.0	27.34	-3767.68	100.38	55.26
	65.0	33.75	-5325.74	70.54	53.97	65.0	33.73	-5307.02	70.46	53.93
	70.0	24.09	-5675.31	75.20	54.50	70.0	24.09	-5656.14	75.11	54.46
60.00000	55.0	23.71	-5945.15	70.54	53.97	55.0	23.69	-5926.60	70.46	53.93
	60.0	24.58	-5675.31	75.43	54.49	60.0	24.55	-5656.14	75.35	54.46
	65.0	25.55	-5129.38	83.42	55.55	65.0	25.52	-5109.68	83.35	55.51
	70.0	26.59	-4579.19	90.76	56.16	70.0	26.56	-4559.92	90.68	56.10
20.00000	55.0	26.92	-4306.36	94.71	58.65	55.0	26.91	-4293.81	94.65	58.60
	60.0	27.62	-3744.33	100.25	57.14	60.0	27.61	-3736.20	100.13	57.08
	65.0	33.79	-5325.74	70.70	54.10	65.0	33.78	-5307.02	70.52	54.06
	70.0	24.09	-5675.31	75.20	54.50	70.0	24.08	-5656.14	75.11	54.46

TABLE A-2. THERMODYNAMIC PROPERTIES OF FLUORINE-OXYGEN MIXTURES IN THE COMPRESSED LIQUID REGION.

WT. FRACTION F2 = 0.70

PRESSURE 10 ⁵ PA	TEMPERATURE K	VOLUME CM ³ /MOL	ENTHALPY J/MOL	ENTROPY J/MOL-K	C _P J/MOL-K	10 ⁵ PA	TEMPERATURE K	VOLUME CM ³ /MOL	ENTHALPY J/MOL	ENTROPY J/MOL-K	C _P J/MOL-K
0.50000	55.0	23.19	-53.46	69.85	54.57	1.01325	55.0	23.13	-59.45	69.84	54.57
	60.0	23.61	-55.74	74.59	55.07		60.0	23.61	-56.73	74.58	55.06
	70.0	24.05	-54.00	78.98	56.26		70.0	24.05	-53.99	78.98	56.26
	75.0	24.51	-51.23	83.05	54.98		75.0	24.51	-51.24	83.05	54.98
5.00000	55.0	25.01	-48.47	86.89	55.61	10.00000	55.0	25.01	-45.68	86.88	55.65
	60.0	25.55	-43.21	90.34	55.99		60.0	25.55	-42.57	90.37	55.98
	70.0	26.11	-41.01	93.84	55.99		70.0	26.11	-39.92	93.77	55.99
	75.0	26.79	-37.14	97.38	57.70		75.0	26.79	-37.07	97.00	57.58
20.00000	55.0	27.39	-33.77	100.16	57.70	30.00000	55.0	27.39	-31.78	100.09	57.58
	60.0	28.04	-31.66	102.92	56.26		60.0	28.04	-29.92	102.92	56.26
	70.0	28.71	-29.92	105.67	54.98		70.0	28.71	-28.41	105.67	54.98
	75.0	29.40	-28.41	108.42	53.69		75.0	29.40	-27.19	108.42	53.69
40.00000	55.0	29.99	-27.19	111.17	52.38	50.00000	55.0	29.99	-26.23	111.17	52.38
	60.0	30.70	-26.23	113.92	51.07		60.0	30.70	-24.93	113.92	51.07
	70.0	31.43	-24.93	116.67	49.76		70.0	31.43	-23.86	116.67	49.76
	75.0	32.18	-23.86	119.42	48.45		75.0	32.18	-23.00	119.42	48.45
60.00000	55.0	32.85	-23.00	122.17	47.14	70.00000	55.0	32.85	-22.23	122.17	47.14
	60.0	33.61	-22.23	124.92	45.83		60.0	33.61	-21.61	124.92	45.83
	70.0	34.39	-21.61	127.67	44.52		70.0	34.39	-21.10	127.67	44.52
	75.0	35.18	-21.10	130.42	43.21		75.0	35.18	-20.70	130.42	43.21

TABLE A-2. THERMODYNAMIC PROPERTIES OF FLUORINE-OXYGEN MIXTURES IN THE COMPRESSED LIQUID REGION

WT. FRACTION F2 = 0.80

PRESSURE 10 ⁵ PA	TEMPERATURE K	VOLUME CM ³ /MOL	ENTHALPY J/MOL	ENTROPY J/MOL-K	C _P J/MOL-K	PRESSURE 10 ⁵ PA	TEMPERATURE K	VOLUME CM ³ /MOL	ENTHALPY J/MOL	ENTROPY J/MOL-K	C _P J/MOL-K
0.50000	55.0	22.34	-536.73	68.34	54.79	1.01325	55.0	22.94	-535.79	68.33	54.78
	60.0	23.36	-537.55	77.51	55.76		60.0	23.34	-564.24	73.02	54.62
	65.0	23.30	-538.34	81.50	56.75		65.0	23.73	-537.93	77.43	55.81
	70.0	24.26	-541.36	85.47	57.25		70.0	24.23	-509.51	81.52	55.21
5.00000	75.0	24.76	-4332.64	85.47	57.96	10.00000	75.0	25.25	-451.52	85.99	55.10
	80.0						80.0	25.63	-425.51	88.90	55.37
	85.0						85.0	26.83	-397.33	92.65	55.76
	90.0						90.0	27.11	-366.88	95.65	55.10
20.00000	95.0					95.0					
	55.0	23.93	-532.47	68.31	54.71	30.00000	55.0	22.92	-591.3	68.27	54.62
	60.0	23.79	-535.33	73.06	55.33		60.0	23.34	-564.24	73.02	55.81
	65.0	24.25	-537.35	77.47	56.25		65.0	24.23	-509.51	81.52	55.21
70.0	24.75	-544.83	81.52	57.00	70.0		25.25	-451.52	85.99	55.37	
40.00000	75.0	25.28	-454.83	85.99	57.96	50.00000	75.0	25.63	-425.51	88.90	55.76
	80.0	25.78	-423.77	89.46	58.37		80.0	26.83	-397.33	92.65	55.10
	85.0	26.39	-333.56	92.52	58.87		85.0	26.99	-365.45	95.65	55.37
	90.0	27.05	-337.77	95.52	59.83		90.0	27.11	-336.88	98.49	55.76
60.00000	95.0					95.0					
	55.0	23.92	-530.72	68.19	54.44	70.00000	55.0	22.89	-588.46	68.11	54.27
	60.0	23.76	-532.15	72.34	55.22		60.0	23.33	-560.95	72.25	55.83
	65.0	24.21	-535.73	77.43	56.24		65.0	24.13	-533.57	77.33	55.23
70.0	24.70	-547.93	81.52	57.00	70.0		24.67	-505.79	81.57	55.72	
80.00000	75.0	25.22	-479.19	85.99	57.96	80.00000	75.0	25.14	-478.29	85.17	57.05
	80.0	25.78	-423.39	89.46	58.37		80.0	25.74	-450.31	88.17	57.05
	85.0	26.39	-333.56	92.52	58.87		85.0	26.34	-422.31	92.17	57.05
	90.0	27.05	-337.77	95.52	59.83		90.0	26.99	-394.45	95.65	57.36
20.00000	95.0					95.0					
	55.0	23.97	-536.73	68.03	54.10	50.00000	55.0	22.85	-584.5	67.95	54.27
	60.0	23.76	-537.39	72.77	55.14		60.0	23.26	-557.46	72.69	55.10
	65.0	24.15	-541.36	77.17	56.24		65.0	23.69	-529.96	77.08	55.92
70.0	24.54	-546.36	81.50	57.00	70.0		24.13	-502.47	81.15	55.18	
40.00000	75.0	24.99	-479.19	85.99	57.96	50.00000	75.0	22.5	-474.74	84.56	57.80
	80.0	25.49	-423.39	89.46	58.37		80.0	22.5	-446.79	88.56	57.00
	85.0	25.99	-333.56	92.52	58.87		85.0	22.5	-419.12	91.95	57.97
	90.0	26.49	-337.77	95.52	59.83		90.0	22.5	-391.45	95.16	57.30
60.00000	95.0					95.0					
	55.0	23.97	-535.19	67.60	53.94	70.00000	55.0	22.85	-533.6	67.51	53.94
	60.0	23.66	-536.03	72.00	54.94		60.0	23.21	-506.92	72.00	54.94
	65.0	24.17	-540.35	76.00	56.16		65.0	23.63	-479.19	76.00	56.16
70.0	24.70	-479.19	80.00	57.00	70.0		24.04	-451.52	80.00	57.00	
80.00000	75.0	25.21	-423.39	84.46	57.96	70.00000	75.0	22.5	-423.39	88.56	57.96
	80.0	25.71	-371.95	88.44	58.91		80.0	22.5	-394.45	91.95	57.84
	85.0	26.18	-319.53	91.84	59.83		85.0	22.5	-365.45	95.16	57.30
	90.0	26.61	-311.11	95.09	60.61		90.0	22.5	-336.88	98.49	57.36

TABLE A-2. THERMODYNAMIC PROPERTIES OF FLUORINE-OXYGEN MIXTURES IN THE COMPRESSED LIQUID REGION

WT. FRACTION F2 = 0.88

PRESSURE	TEMPERATURE	VOLUME	ENTHALPY	ENTROPY	C _P	PRESSURE	TEMPERATURE	VOLUME	ENTHALPY	ENTROPY	C _P
10 ⁵ PA	K	CM ³ /MOL	J/MOL	J/MOL-K	J/MOL-K	10 ⁵ PA	K	CM ³ /MOL	J/MOL	J/MOL-K	J/MOL-K
	55.0	22.74	-53.34	66.65	54.96		55.0	22.74	-53.34	66.65	54.96
	60.0	23.13	-53.59	71.42	55.61		60.0	23.13	-53.59	71.42	55.61
0.50000	65.0	23.53	-53.85	75.86	57.17	1.01325	65.0	23.53	-53.85	75.86	57.17
	70.0	24.05	-54.09	79.97	58.47		70.0	24.05	-54.09	79.97	58.47
	75.0	24.55	-54.25	83.85	59.25		80.0	25.09	-54.42	87.48	56.25
	55.0	22.73	-59.26	66.62	54.88		55.0	22.72	-59.19	66.58	54.78
	60.0	23.14	-59.52	71.39	55.57		60.0	23.13	-59.43	71.38	55.52
	65.0	23.58	-59.79	75.82	57.19		65.0	23.57	-59.67	75.78	57.04
5.00000	70.0	24.03	-60.03	79.93	58.48	10.00000	70.0	24.02	-60.00	79.88	58.48
	75.0	24.54	-60.28	83.44	59.63		80.0	24.52	-60.28	83.75	56.42
	80.0	25.06	-60.54	87.00	60.89		85.0	25.16	-60.51	87.00	59.05
	85.0	25.52	-60.71	90.00	61.64		90.0	25.62	-60.71	90.00	59.49
	90.0	26.03	-60.89	92.70	62.66		95.0	26.03	-60.89	92.70	59.49
	55.0	22.70	-59.99	66.50	54.59		55.0	22.72	-59.80	66.42	54.40
	60.0	23.11	-60.26	71.27	55.46		60.0	23.10	-60.27	71.18	55.26
	65.0	23.55	-60.54	75.69	57.26		65.0	23.52	-60.70	75.60	57.47
	70.0	24.00	-60.82	79.79	58.48		70.0	24.45	-61.01	79.70	56.01
20.00000	75.0	24.49	-61.11	83.65	59.69	30.00000	75.0	24.97	-61.31	83.55	56.36
	80.0	25.01	-61.41	87.28	60.97		80.0	25.52	-61.57	87.18	56.32
	85.0	25.56	-61.70	90.71	62.33		85.0	26.11	-61.79	90.60	56.32
	90.0	26.17	-61.99	93.97	63.78		90.0	26.70	-61.99	93.84	56.32
	95.0	26.83	-62.28	97.08	65.20		95.0	27.30	-62.28	96.95	57.73
	55.0	22.67	-59.59	66.35	54.28		55.0	22.65	-59.40	66.26	54.14
	60.0	23.07	-59.87	71.10	55.38		60.0	23.05	-59.71	71.02	55.26
	65.0	23.49	-60.16	75.52	56.81		65.0	23.47	-60.03	75.43	56.43
	70.0	23.94	-60.45	79.61	58.46		70.0	23.92	-60.35	79.36	57.73
40.00000	75.0	24.42	-60.74	83.46	59.11	50.00000	75.0	24.91	-61.06	83.36	56.17
	80.0	24.94	-61.03	87.07	60.49		80.0	25.44	-61.34	87.00	56.38
	85.0	25.47	-61.32	90.49	61.97		85.0	26.00	-61.61	90.38	56.38
	90.0	26.07	-61.61	93.72	63.54		90.0	26.60	-61.89	93.68	57.74
	95.0	26.70	-61.92	96.81	65.17		95.0	27.20	-62.16	96.68	59.14
	60.0	23.03	-59.27	70.95	54.39		60.0	23.01	-59.08	70.86	54.26
	65.0	23.43	-59.56	75.43	55.82		65.0	23.41	-59.37	75.37	55.59
	70.0	23.86	-59.85	79.56	57.41		70.0	23.83	-59.66	79.47	57.09
	75.0	24.30	-60.14	83.26	58.95		75.0	24.25	-60.00	83.16	58.24
	80.0	24.76	-60.43	86.66	60.23		80.0	24.70	-60.35	86.76	59.16
60.00000	85.0	25.23	-60.72	90.27	61.59	70.00000	85.0	25.20	-60.71	90.16	59.35
	90.0	25.71	-61.01	93.48	63.04		90.0	25.71	-61.01	93.16	59.42
	95.0	26.21	-61.30	96.55	64.57		95.0	26.25	-61.30	96.15	59.42

TABLE A-2. THERMODYNAMIC PROPERTIES OF FLUORINE-OXYGEN MIXTURES IN THE COMPRESSED LIQUID REGION

WT. FRACTION F2 = 1.00

WT. FRACTION F2	PRESSURE 10 ⁵ PA	TEMPERATURE K	VOLUME CM ³ /MOL	ENTHALPY J/MOL	ENTROPY J/MOL-K	C _P J/MOL-K	PRESSURE 10 ⁵ PA	TEMPERATURE K	VOLUME CM ³ /MOL	ENTHALPY J/MOL	ENTROPY J/MOL-K	C _P J/MOL-K
1.00	1.00	55.0	23.41	-5341.30	62.39	55.24	1.00	55.0	23.41	-5947.36	62.38	55.23
1.00	1.00	60.0	23.32	-5365.31	61.18	55.29	1.00	60.0	23.32	-5968.90	61.18	55.28
1.00	1.00	70.0	23.25	-5386.36	71.64	57.78	1.00	70.0	23.25	-5386.97	71.64	57.78
1.00	1.00	75.0	24.21	-5107.36	75.71	55.82	1.00	75.0	24.21	-4823.82	79.68	55.82
1.00	1.00	75.0	24.21	-4324.68	79.69	56.71	1.00	75.0	24.21	-4533.64	83.35	56.68
1.00	5.00	55.0	22.41	-5333.20	62.35	55.14	5.00	55.0	22.41	-5924.18	62.31	55.03
1.00	5.00	60.0	22.32	-5357.21	61.14	55.19	5.00	60.0	22.32	-5648.94	61.16	55.90
1.00	5.00	70.0	23.24	-5379.17	71.60	57.83	5.00	70.0	23.24	-5378.59	71.56	57.84
1.00	5.00	75.0	24.19	-5110.16	75.74	55.82	5.00	75.0	24.19	-5091.42	75.59	55.94
1.00	5.00	80.0	24.71	-4333.18	79.31	56.77	5.00	80.0	24.71	-4408.77	83.26	56.77
1.00	5.00	85.0	25.28	-4237.11	83.78	57.19	5.00	85.0	25.28	-4225.07	86.77	57.17
1.00	5.00	90.0	25.99	-3359.49	89.08	58.23	5.00	90.0	25.99	-3953.42	90.14	58.23
1.00	5.00	95.0	26.97	-3265.35	93.25	59.23	5.00	95.0	26.97	-3265.82	93.17	59.21
1.00	20.00	55.0	22.77	-5306.17	62.23	54.83	20.00	55.0	22.77	-5889.14	62.15	54.83
1.00	20.00	60.0	22.78	-5331.35	61.02	55.89	20.00	60.0	22.78	-5611.36	61.03	55.78
1.00	20.00	70.0	23.56	-5074.22	71.47	57.14	20.00	70.0	23.56	-5335.41	71.38	57.14
1.00	20.00	75.0	24.14	-4792.27	73.60	55.85	20.00	75.0	24.14	-5055.82	73.51	55.85
1.00	20.00	80.0	24.66	-4508.18	79.15	56.82	20.00	80.0	24.66	-4779.59	79.04	56.82
1.00	20.00	85.0	25.21	-4232.11	86.60	57.18	20.00	85.0	25.21	-4420.53	86.49	57.18
1.00	20.00	90.0	25.84	-3345.67	89.89	57.73	20.00	90.0	25.84	-3923.22	89.76	57.73
1.00	20.00	95.0	26.44	-3245.47	93.03	58.85	20.00	95.0	26.44	-3243.41	92.90	58.83
1.00	40.00	55.0	23.74	-5270.24	62.08	54.40	40.00	55.0	23.74	-5852.64	61.97	54.40
1.00	40.00	60.0	23.74	-5295.44	60.85	55.46	40.00	60.0	23.74	-5572.70	60.87	55.29
1.00	40.00	70.0	24.51	-5037.20	71.30	56.22	40.00	70.0	24.51	-5300.15	71.31	56.22
1.00	40.00	75.0	24.98	-4759.10	73.42	55.99	40.00	75.0	24.98	-5020.20	73.20	55.99
1.00	40.00	80.0	25.43	-4476.00	79.38	56.88	40.00	80.0	25.43	-4746.93	79.27	56.88
1.00	40.00	85.0	25.93	-4192.99	86.64	56.99	40.00	85.0	25.93	-4472.82	86.62	56.99
1.00	40.00	90.0	26.43	-3307.89	89.76	58.11	40.00	90.0	26.43	-3893.81	89.53	58.11
1.00	40.00	95.0	27.03	-3208.35	92.76	59.21	40.00	95.0	27.03	-3602.47	92.63	59.19
1.00	60.00	55.0	22.30	-5334.66	61.91	54.41	60.00	55.0	22.30	-5814.45	61.83	54.41
1.00	60.00	60.0	22.31	-5359.67	60.69	55.48	60.00	60.0	22.31	-5540.16	60.71	55.31
1.00	60.00	70.0	23.15	-5081.27	71.13	56.23	60.00	70.0	23.15	-5266.25	71.03	56.23
1.00	60.00	75.0	24.02	-4803.18	73.27	55.97	60.00	75.0	24.02	-4790.42	73.07	55.97
1.00	60.00	80.0	24.54	-4520.08	79.16	56.82	60.00	80.0	24.54	-4424.84	79.03	56.82
1.00	60.00	85.0	25.05	-4237.01	86.60	57.18	60.00	85.0	25.05	-4149.79	86.66	57.18
1.00	60.00	90.0	25.56	-3352.93	89.76	58.21	60.00	90.0	25.56	-3881.89	89.57	58.19
1.00	60.00	95.0	26.07	-3253.39	92.76	59.24	60.00	95.0	26.07	-3591.96	92.67	59.21

TABLE B-1. THERMODYNAMIC PROPERTIES OF LIQUID FLUORINE-OXYGEN MIXTURES ON THE SATURATION BOUNDARY

TEMPERATURE R	WT. FRACTION F ₂		PRESSURE PSIA	VOLUME FT ³ /LB	ENTHALPY BTU/LB	ENTROPY BTU/LB/R	C _P BTU/LB/R
	LIQUID	VAPOR					
100.0	0.6000	0.7552	0.0577	0.01063	-72.215	0.4828	0.3661
	0.7000	0.3139	0.0606	0.01034	-70.761	0.4677	0.3603
	0.8000	0.3720	0.0634	0.01005	-69.301	0.4495	0.3556
	0.8800	0.3203	0.0656	0.00982	-68.359	0.4322	0.3514
	0.9000	0.3328	0.0660	0.00976	-68.111	0.4274	0.3503
	1.0000	1.0000	0.0685	0.00946	-66.930	0.3959	0.3451
110.0	0.6000	0.7514	0.2461	0.01084	-68.543	0.5178	0.3633
	0.7000	0.3127	0.2590	0.01055	-67.140	0.5022	0.3583
	0.8000	0.3725	0.2712	0.01026	-65.820	0.4836	0.3525
	0.8800	0.3213	0.2805	0.01002	-64.828	0.4659	0.3481
	0.9000	0.3338	0.2828	0.00996	-64.590	0.4609	0.3470
	1.0000	1.0000	0.2937	0.00966	-63.458	0.4290	0.3413
120.0	0.6000	0.7447	0.8092	0.01107	-64.850	0.5499	0.3667
	0.7000	0.3088	0.8511	0.01078	-63.494	0.5339	0.3615
	0.8000	0.3708	0.8910	0.01047	-62.219	0.5149	0.3559
	0.8800	0.3208	0.9216	0.01023	-61.264	0.4969	0.3522
	0.9000	0.3335	0.9291	0.01017	-61.034	0.4919	0.3511
	1.0000	1.0000	0.9655	0.00986	-59.948	0.4595	0.3453
130.0	0.6000	0.7371	2.1809	0.01132	-61.135	0.5797	0.3683
	0.7000	0.3039	2.2912	0.01101	-59.822	0.5633	0.3633
	0.8000	0.3683	2.3966	0.01071	-58.592	0.5440	0.3588
	0.8800	0.3198	2.4780	0.01046	-57.672	0.5257	0.3543
	0.9000	0.3328	2.4980	0.01040	-57.451	0.5206	0.3539
	1.0000	1.0000	2.5952	0.01008	-56.408	0.4878	0.3483
140.0	0.6000	0.7236	5.0370	0.01157	-57.397	0.6073	0.3707
	0.7000	0.7990	5.2846	0.01127	-56.126	0.5907	0.3657
	0.8000	0.3656	5.5225	0.01096	-54.939	0.5710	0.3609
	0.8800	0.3135	5.7067	0.01071	-54.053	0.5524	0.3563
	0.9000	0.3319	5.7519	0.01064	-53.841	0.5473	0.3560
	1.0000	1.0000	5.9731	0.01032	-52.840	0.5142	0.3511
150.0	0.6000	0.7227	10.2958	0.01185	-53.634	0.6332	0.3740
	0.7000	0.7945	10.7897	0.01154	-52.406	0.6163	0.3694
	0.8000	0.3631	11.2663	0.01123	-51.259	0.5964	0.3643
	0.8800	0.3174	11.6369	0.01097	-50.404	0.5776	0.3611
	0.9000	0.3309	11.7281	0.01091	-50.200	0.5723	0.3601
	1.0000	1.0000	12.1755	0.01058	-49.240	0.5390	0.3555
160.0	0.6000	0.7166	19.0933	0.01216	-49.837	0.6576	0.3803
	0.7000	0.7904	19.9913	0.01184	-48.648	0.6404	0.3750
	0.8000	0.3609	20.8626	0.01152	-47.541	0.6203	0.3731
	0.8800	0.3153	21.5430	0.01126	-46.718	0.6013	0.3699
	0.9000	0.3301	21.7108	0.01120	-46.522	0.5960	0.3691
	1.0000	1.0000	22.5372	0.01086	-45.600	0.5623	0.3652
170.0	0.6000	0.7107	32.7573	0.01249	-45.987	0.6808	0.3863
	0.7000	0.7864	34.2707	0.01217	-44.837	0.6634	0.3825
	0.8000	0.3536	35.7466	0.01184	-43.767	0.6430	0.3789
	0.8800	0.3152	36.9040	0.01158	-42.975	0.6238	0.3759
	0.9000	0.3292	37.1904	0.01151	-42.787	0.6184	0.3752
	1.0000	1.0000	38.6038	0.01118	-41.903	0.5845	0.3715
180.0	0.6000	0.7043	52.7445	0.01286	-42.067	0.7030	0.3923
	0.7000	0.7824	55.1428	0.01253	-40.953	0.6854	0.3892
	0.8000	0.3553	57.4945	0.01220	-39.921	0.6647	0.3857
	0.8800	0.3140	59.3467	0.01194	-39.158	0.6453	0.3829
	0.9000	0.3283	59.8061	0.01187	-38.977	0.6399	0.3822
	1.0000	1.0000	62.0800	0.01152	-38.131	0.6058	0.3787

TABLE B-2. THERMODYNAMIC PROPERTIES OF FLUORINE-OXYGEN MIXTURES IN THE COMPRESSED LIQUID REGION

WT. FRACTION F2 = 0.60

PRESSURE	TEMPERATURE	VOLUME	ENTHALPY	ENTROPY	C _P	PSIA	R	FT ³ /LB	BTU/LB	BTU/LB/R	ENTROPY	C _P
PSIA	R	FT ³ /LB	BTU/LB	BTU/LB/R	BTU/LB/R	PSIA	R	FT ³ /LB	BTU/LB	BTU/LB/R	ENTROPY	C _P
8.000	100.0	0.01063	-72.203	0.4828	0.3595	14.696	100.0	0.01053	-72.192	0.4828	0.3594	
	110.0	0.01084	-63.532	0.5178	0.3700		110.0	0.01094	-68.520	0.5177	0.3700	
	120.0	0.01107	-64.845	0.5499	0.3726		120.0	0.01127	-64.305	0.5499	0.3726	
	130.0	0.01132	-61.121	0.5797	0.3752		130.0	0.01157	-61.185	0.5797	0.3752	
140.0	0.01157	-57.355	0.6073	0.3779	140.0	0.01185	-57.355	0.6073	0.3779			
30.000	100.0	0.01063	-72.169	0.4827	0.3593	50.000	100.0	0.01053	-72.137	0.4826	0.3592	
	110.0	0.01084	-63.497	0.5177	0.3700		110.0	0.01094	-68.467	0.5176	0.3700	
	120.0	0.01107	-64.807	0.5498	0.3727		120.0	0.01131	-64.776	0.5495	0.3727	
	130.0	0.01131	-61.092	0.5795	0.3753		130.0	0.01157	-61.333	0.5795	0.3753	
140.0	0.01157	-57.363	0.6072	0.3779	140.0	0.01185	-57.581	0.6071	0.3779			
150.0	0.01185	-53.603	0.6331	0.3805	150.0	0.01215	-53.841	0.6330	0.3805			
160.0	0.01215	-49.827	0.6576	0.3831	160.0	0.01248	-49.801	0.6575	0.3831			
170.0	0.01248	-46.041	0.6804	0.3857	170.0	0.01282	-45.537	0.6807	0.3857			
180.0	0.01282	-42.255	0.7027	0.3883	180.0	0.01317	-41.917	0.7019	0.3883			
100.000	100.0	0.01062	-72.058	0.4824	0.3588	200.000	100.0	0.01052	-71.900	0.4820	0.3582	
	110.0	0.01084	-63.389	0.5174	0.3700		110.0	0.01093	-68.253	0.5170	0.3700	
	120.0	0.01106	-64.699	0.5495	0.3729		120.0	0.01130	-64.585	0.5491	0.3722	
	130.0	0.01131	-60.988	0.5793	0.3759		130.0	0.01155	-60.849	0.5788	0.3753	
140.0	0.01156	-57.261	0.6069	0.3789	140.0	0.01183	-57.171	0.6064	0.3789			
150.0	0.01184	-53.513	0.6327	0.3818	150.0	0.01213	-53.371	0.6322	0.3818			
160.0	0.01214	-49.735	0.6572	0.3848	160.0	0.01245	-49.605	0.6566	0.3848			
170.0	0.01248	-45.912	0.6804	0.3878	170.0	0.01282	-45.791	0.6797	0.3878			
180.0	0.01284	-42.084	0.7027	0.3908	180.0	0.01322	-41.917	0.7019	0.3908			
300.000	100.0	0.01061	-71.747	0.4817	0.3576	400.000	100.0	0.01050	-71.584	0.4813	0.3570	
	110.0	0.01085	-63.077	0.5166	0.3700		110.0	0.01092	-67.922	0.5162	0.3700	
	120.0	0.01105	-64.386	0.5486	0.3725		120.0	0.01134	-64.251	0.5483	0.3725	
	130.0	0.01129	-60.675	0.5783	0.3750		130.0	0.01177	-60.540	0.5779	0.3750	
140.0	0.01154	-56.911	0.6059	0.3776	140.0	0.01153	-56.825	0.6054	0.3750			
150.0	0.01181	-53.124	0.6317	0.3802	150.0	0.01130	-53.092	0.6312	0.3802			
160.0	0.01211	-49.272	0.6560	0.3828	160.0	0.01210	-49.341	0.6555	0.3811			
170.0	0.01244	-45.463	0.6791	0.3854	170.0	0.01247	-45.545	0.6785	0.3854			
180.0	0.01279	-41.614	0.7012	0.3880	180.0	0.01287	-41.705	0.6988	0.3880			
500.000	100.0	0.01060	-71.435	0.4809	0.3563	1000.000	100.0	0.01057	-70.837	0.4791	0.3562	
	110.0	0.01081	-62.764	0.5158	0.3700		110.0	0.01098	-67.179	0.5153	0.3700	
	120.0	0.01103	-64.073	0.5479	0.3725		120.0	0.01139	-63.520	0.5459	0.3725	
	130.0	0.01127	-60.362	0.5776	0.3750		130.0	0.01182	-60.000	0.5753	0.3750	
140.0	0.01152	-56.611	0.6052	0.3776	140.0	0.01147	-56.490	0.6027	0.3750			
150.0	0.01176	-52.825	0.6307	0.3802	150.0	0.01120	-52.944	0.6282	0.3802			
160.0	0.01200	-49.019	0.6549	0.3828	160.0	0.01200	-49.440	0.6549	0.3828			
170.0	0.01225	-45.183	0.6789	0.3854	170.0	0.01283	-45.779	0.6792	0.3854			
180.0	0.01275	-41.315	0.7012	0.3880	180.0	0.01323	-41.916	0.7016	0.3880			

TABLE B-2. THERMODYNAMIC PROPERTIES OF FLUORINE-OXYGEN MIXTURES IN THE COMPRESSED LIQUID REGION

WT. FRACTION F2 = 0.70

TEMPERATURE	TEMPERATURE	VOLUME	ENTHALPY	ENTROPY	C _P	TEMPERATURE	VOLUME	ENTHALPY	ENTROPY	C _P
PSIA	R	FT ³ /LB	BTU/LB	BTU/LB/R	BTU/LB/R	R	FT ³ /LB	BTU/LB	BTU/LB/R	BTU/LB/R
8.000	100.0	0.010334	-70.743	0.4677	0.3647	100.0	0.010334	-70.739	0.4676	0.3647
	110.0	0.01055	-67.130	0.5022	0.3653	110.0	0.01055	-67.120	0.5021	0.3653
	120.0	0.01078	-63.484	0.5338	0.3667	120.0	0.01078	-63.474	0.5337	0.3667
	130.0	0.01101	-59.781	0.5632	0.3683	130.0	0.01101	-59.803	0.5632	0.3683
	140.0	0.01127	-56.095	0.5906	0.3707	140.0	0.01127	-56.119	0.5905	0.3707
	150.0	0.01154	-52.382	0.6162	0.3739	150.0	0.01154	-52.403	0.6161	0.3739
	160.0	0.01184	-48.633	0.6404	0.3754	160.0	0.01184	-48.614	0.6403	0.3754
30.000	100.0	0.01334	-70.716	0.4676	0.3645	100.0	0.01334	-70.685	0.4675	0.3644
	110.0	0.01355	-67.096	0.5021	0.3653	110.0	0.01355	-67.066	0.5020	0.3654
	120.0	0.01378	-63.451	0.5338	0.3667	120.0	0.01378	-63.421	0.5338	0.3668
	130.0	0.01401	-59.781	0.5632	0.3683	130.0	0.01401	-59.752	0.5631	0.3685
	140.0	0.01427	-56.095	0.5906	0.3707	140.0	0.01427	-56.066	0.5905	0.3685
	150.0	0.01454	-52.382	0.6162	0.3739	150.0	0.01454	-52.355	0.6161	0.3709
	160.0	0.01484	-48.633	0.6404	0.3754	160.0	0.01484	-48.614	0.6403	0.3752
100.000	100.0	0.0134	-70.693	0.4674	0.3640	100.0	0.0133	-70.455	0.4670	0.3635
	110.0	0.01377	-67.047	0.5028	0.3657	110.0	0.01354	-66.839	0.5014	0.3655
	120.0	0.01401	-63.373	0.5339	0.3680	120.0	0.01378	-63.197	0.5325	0.3691
	130.0	0.01426	-59.679	0.5633	0.3705	130.0	0.01399	-59.534	0.5898	0.3685
	140.0	0.01453	-55.962	0.6159	0.3740	140.0	0.01452	-52.152	0.6153	0.3710
	150.0	0.01483	-52.286	0.6400	0.3746	150.0	0.01481	-48.425	0.6395	0.3735
	160.0	0.01512	-48.551	0.6685	0.3829	160.0	0.01514	-44.650	0.6624	0.3797
	170.0	0.01542	-44.814	0.6880	0.3899	170.0	0.01549	-40.811	0.6844	0.3884
300.000	100.0	0.00533	-70.302	0.4666	0.3656	100.0	0.01033	-70.148	0.4667	0.3659
	110.0	0.00575	-66.643	0.5017	0.3674	110.0	0.01054	-66.591	0.5027	0.3671
	120.0	0.00619	-62.983	0.5361	0.3699	120.0	0.01078	-62.941	0.5327	0.3680
	130.0	0.00663	-59.317	0.5695	0.3726	130.0	0.01101	-59.277	0.5689	0.3696
	140.0	0.00707	-55.651	0.6029	0.3751	140.0	0.01127	-55.561	0.6383	0.3710
	150.0	0.00751	-51.985	0.6363	0.3776	150.0	0.01154	-51.840	0.6670	0.3782
	160.0	0.00795	-48.319	0.6696	0.3801	160.0	0.01184	-48.113	0.6957	0.3843
	170.0	0.00839	-44.653	0.7028	0.3826	170.0	0.01216	-44.386	0.7240	0.3910
500.000	100.0	0.0031	-63.994	0.4659	0.3512	100.0	0.01029	-63.228	0.4641	0.3589
	110.0	0.00352	-60.328	0.5033	0.3575	110.0	0.01070	-60.470	0.4983	0.3655
	120.0	0.00397	-56.662	0.5407	0.3638	120.0	0.01112	-57.712	0.5300	0.3686
	130.0	0.00442	-52.996	0.5781	0.3701	130.0	0.01154	-54.954	0.5617	0.3678
	140.0	0.00487	-49.330	0.6155	0.3764	140.0	0.01199	-52.196	0.6151	0.3688
	150.0	0.00532	-45.664	0.6529	0.3827	150.0	0.01249	-49.438	0.6574	0.3739
	160.0	0.00577	-41.998	0.6903	0.3890	160.0	0.01299	-46.680	0.6916	0.3779
	170.0	0.00622	-38.332	0.7277	0.3953	170.0	0.01349	-43.922	0.7259	0.3819

TABLE B-2. THERMODYNAMIC PROPERTIES OF F-URINE-OXYGEN MIXTURES IN THE COMPRESSED LIQUID REGION

WT. FRACTION F2 = 0.80

PRESSURE	TEMPERATURE	VOLUME	ENTHALPY	ENTROPY	C_P	PSIA	TEMPERATURE	VOLUME	ENTHALPY	ENTROPY	C_P
PSIA	R	FT ³ /LB	BTU/LB	BTU/LB/R	BTU/LB/R	PSIA	R	FT ³ /LB	BTU/LB	BTU/LB/R	BTU/LB/R
6.000	100.0	0.01005	-63.373	0.4496	0.3500	14.696	100.0	0.01025	-69.369	0.4495	0.3600
	110.0	0.01027	-62.810	0.4836	0.3607		110.0	0.01025	-65.800	0.4836	0.3607
	120.0	0.01071	-62.212	0.5149	0.3644		120.0	0.01071	-62.200	0.5149	0.3644
	130.0	0.01071	-58.582	0.5540	0.3633		130.0	0.01071	-58.573	0.5540	0.3633
	140.0	0.01096	-54.933	0.5740	0.3639		140.0	0.01096	-54.929	0.5740	0.3639
30.000	100.0	0.01005	-63.347	0.4495	0.3598	50.000	100.0	0.01005	-69.317	0.4494	0.3596
	110.0	0.01026	-62.777	0.4835	0.3607		110.0	0.01025	-65.748	0.4834	0.3607
	120.0	0.01071	-62.151	0.5148	0.3643		120.0	0.01071	-62.149	0.5147	0.3643
	130.0	0.01095	-58.503	0.5540	0.3639		130.0	0.01095	-58.523	0.5540	0.3639
	140.0	0.01123	-54.837	0.5740	0.3665		140.0	0.01122	-54.881	0.5709	0.3665
100.000	100.0	0.01005	-63.347	0.4495	0.3598	200.000	100.0	0.01005	-69.288	0.4492	0.3592
	110.0	0.01026	-62.777	0.4835	0.3607		110.0	0.01025	-65.720	0.4834	0.3607
	120.0	0.01071	-62.151	0.5148	0.3643		120.0	0.01071	-62.149	0.5147	0.3643
	130.0	0.01095	-58.503	0.5540	0.3639		130.0	0.01095	-58.523	0.5540	0.3639
	140.0	0.01123	-54.837	0.5740	0.3665		140.0	0.01122	-54.881	0.5709	0.3665
300.000	100.0	0.01004	-63.347	0.4492	0.3577	400.000	100.0	0.01004	-69.259	0.4488	0.3585
	110.0	0.01025	-62.787	0.4832	0.3584		110.0	0.01024	-65.693	0.4828	0.3585
	120.0	0.01069	-62.171	0.5147	0.3637		120.0	0.01069	-62.132	0.5142	0.3636
	130.0	0.01093	-58.533	0.5540	0.3643		130.0	0.01093	-58.516	0.5542	0.3643
	140.0	0.01119	-54.882	0.5740	0.3668		140.0	0.01121	-54.913	0.5693	0.3669
500.000	100.0	0.01004	-63.347	0.4492	0.3577	600.000	100.0	0.01004	-69.224	0.4488	0.3585
	110.0	0.01025	-62.787	0.4832	0.3584		110.0	0.01024	-65.658	0.4828	0.3585
	120.0	0.01069	-62.171	0.5147	0.3637		120.0	0.01069	-62.132	0.5142	0.3636
	130.0	0.01093	-58.533	0.5540	0.3643		130.0	0.01093	-58.516	0.5542	0.3643
	140.0	0.01119	-54.882	0.5740	0.3668		140.0	0.01121	-54.913	0.5693	0.3669

TABLE B-2. THERMODYNAMIC PROPERTIES OF FLUORINE-OXYGEN MIXTURES IN THE COMPRESSED LIQUID REGION

WT. FRACTION F2 = 0.88

PRESSURE	TEMPERATURE	VOLUME	ENTHALPY	ENTROPY	C _P	PSIA	R	FT ³ /LB	BTU/LB	BTU/LB/R	ENTROPY	ENTHALPY	ENTROPY	C _P	
PSIA	R	FT ³ /LB	BTU/LB	BTU/LB/R	BTU/LB/R	PSIA	R	FT ³ /LB	BTU/LB	BTU/LB/R	ENTROPY	ENTHALPY	ENTROPY	BTU/LB/R	
8.000	100.0	0.00982	-63.343	0.4323	0.3562	14.696	100.0	0.00932	-68.338	0.4322	0.4322	-68.338	0.4322	0.3562	
	110.0	0.01052	-64.811	0.4499	0.3567		110.0	0.01001	-64.754	0.4321	0.4321	0.4321	-64.754	0.4321	0.3571
	130.0	0.01123	-67.253	0.4669	0.3570		130.0	0.01023	-67.195	0.4310	0.4310	0.4310	-67.195	0.4310	0.3598
	140.0	0.01046	-57.662	0.5252	0.3503		150.0	0.01070	-57.604	0.4255	0.4255	0.4255	-57.604	0.4255	0.3604
30.000	100.0	0.01071	-54.052	0.5525	0.3508	50.000	100.0	0.01037	-50.359	0.4273	0.4273	-50.359	0.4273	0.3575	
	110.0	0.01126	-45.711	0.6012	0.3579		120.0	0.01135	-46.687	0.4211	0.4211	0.4211	-46.687	0.4211	0.3677
	130.0	0.00982	-64.787	0.4658	0.3560		150.0	0.01158	-42.364	0.4237	0.4237	0.4237	-42.364	0.4237	0.3771
	140.0	0.01023	-61.223	0.4968	0.3570		160.0	0.01001	-61.223	0.4321	0.4321	0.4321	-61.223	0.4321	0.3558
100.000	100.0	0.00981	-68.214	0.4319	0.3554	200.000	100.0	0.00931	-68.069	0.4315	0.4315	-68.069	0.4315	0.3546	
	110.0	0.01021	-61.125	0.4966	0.3613		110.0	0.01022	-64.543	0.4651	0.4651	0.4651	-64.543	0.4651	0.3573
	130.0	0.01045	-53.923	0.5253	0.3605		130.0	0.01044	-50.383	0.4961	0.4961	0.4961	-50.383	0.4961	0.3618
	140.0	0.01036	-45.293	0.5772	0.3532		140.0	0.01059	-50.796	0.4516	0.4516	0.4516	-50.796	0.4516	0.3602
300.000	100.0	0.01152	-42.911	0.6008	0.3766	400.000	100.0	0.01123	-46.510	0.4766	0.4766	-46.510	0.4766	0.3634	
	110.0	0.01152	-33.124	0.6450	0.3840		120.0	0.01155	-42.802	0.4827	0.4827	0.4827	-42.802	0.4827	0.3659
	130.0	0.00981	-67.353	0.4312	0.3575		140.0	0.01190	-39.029	0.4644	0.4644	0.4644	-39.029	0.4644	0.3824
	140.0	0.01044	-51.261	0.5245	0.3622		150.0	0.01030	-67.777	0.4308	0.4308	0.4308	-67.777	0.4308	0.3599
500.000	100.0	0.01094	-50.039	0.5997	0.3637	1000.000	100.0	0.00976	-66.406	0.4287	0.4287	-66.406	0.4287	0.3494	
	110.0	0.01153	-42.692	0.6437	0.3808		110.0	0.01036	-62.392	0.4622	0.4622	0.4622	-62.392	0.4622	0.3578
	130.0	0.01066	-43.773	0.5732	0.3620		130.0	0.01078	-56.739	0.4515	0.4515	0.4515	-56.739	0.4515	0.3618
	140.0	0.01091	-42.149	0.5986	0.3620		140.0	0.01051	-52.134	0.4480	0.4480	0.4480	-52.134	0.4480	0.3616
150.0	0.01149	-42.149	0.5986	0.3620	150.0	0.01112	-45.534	0.4561	0.4561	0.4561	-45.534	0.4561	0.3613		
160.0	0.01183	-34.733	0.6423	0.3766	160.0	0.01111	-41.181	0.4581	0.4581	0.4581	-41.181	0.4581	0.3610		
170.0	0.01183	-34.733	0.6423	0.3766	170.0	0.01172	-38.821	0.4329	0.4329	0.4329	-38.821	0.4329	0.3715		

TABLE B-2. THERMODYNAMIC PROPERTIES OF F-UOPRINE-OXYGEN MIXTURES IN THE COMPRESSED LIQUID REGION

WT. FRACTION F2 = 0.90

PRESSURE	TEMPERATURE	VOLUME	ENTHALPY	ENTROPY	C_P	PSIA	TEMPERATURE	VOLUME	ENTHALPY	ENTROPY	C_P
PSIA	R	FT ³ /LB	BTU/LB	BTU/LB/R	BTU/LB/R		P	FT ³ /LB	BTU/LB	BTU/LB/R	BTU/LB/R
8.000	100.0	0.00976	-68.090	0.4273	0.3552		100.0	0.00976	-68.090	0.4273	0.3552
	110.0	0.00999	-64.571	0.4260	0.3561		110.0	0.00999	-64.571	0.4260	0.3561
	120.0	0.01017	-61.023	0.4919	0.3599	14.696	120.0	0.01017	-61.023	0.4919	0.3599
	130.0	0.01040	-57.441	0.5206	0.3588		130.0	0.01040	-57.441	0.5206	0.3588
30.000	140.0	0.01064	-53.833	0.5473	0.3594		140.0	0.01064	-53.833	0.5473	0.3594
	150.0	0.01021	-50.180	0.5722	0.3622	50.000	150.0	0.01021	-50.180	0.5722	0.3622
	160.0	0.01120	-45.514	0.5959	0.3671		160.0	0.01120	-45.514	0.5959	0.3671
	170.0	0.00996	-68.039	0.4273	0.3549		170.0	0.00996	-68.039	0.4273	0.3549
100.000	110.0	0.00996	-64.543	0.4608	0.3561		110.0	0.00996	-64.543	0.4608	0.3561
	120.0	0.01017	-60.994	0.4918	0.3599		120.0	0.01017	-60.994	0.4918	0.3599
	130.0	0.01040	-57.411	0.5202	0.3589	50.000	130.0	0.01040	-57.411	0.5202	0.3589
	140.0	0.01063	-53.713	0.5469	0.3594		140.0	0.01063	-53.713	0.5469	0.3594
300.000	150.0	0.01090	-50.083	0.5719	0.3622	200.000	150.0	0.01090	-50.083	0.5719	0.3622
	160.0	0.01118	-46.432	0.5955	0.3662		160.0	0.01118	-46.432	0.5955	0.3662
	170.0	0.01150	-42.723	0.6180	0.3759		170.0	0.01150	-42.723	0.6180	0.3759
	180.0	0.01166	-39.043	0.6396	0.3833		180.0	0.01166	-39.043	0.6396	0.3833
500.000	100.0	0.00975	-67.822	0.4271	0.3544		100.0	0.00975	-67.822	0.4271	0.3544
	110.0	0.00995	-64.443	0.4605	0.3562		110.0	0.00995	-64.443	0.4605	0.3562
	120.0	0.01016	-60.896	0.4915	0.3594	200.000	120.0	0.01016	-60.896	0.4915	0.3594
	130.0	0.01039	-57.315	0.5202	0.3591		130.0	0.01039	-57.315	0.5202	0.3591
1000.000	140.0	0.01063	-53.713	0.5469	0.3594	200.000	140.0	0.01063	-53.713	0.5469	0.3594
	150.0	0.01090	-50.083	0.5719	0.3622		150.0	0.01090	-50.083	0.5719	0.3622
	160.0	0.01118	-46.432	0.5955	0.3662	400.000	160.0	0.01118	-46.432	0.5955	0.3662
	170.0	0.01150	-42.723	0.6180	0.3759		170.0	0.01150	-42.723	0.6180	0.3759
1000.000	180.0	0.01166	-39.043	0.6396	0.3833		180.0	0.01166	-39.043	0.6396	0.3833
	100.0	0.00974	-67.677	0.4263	0.3527		100.0	0.00974	-67.677	0.4263	0.3527
	110.0	0.00994	-64.163	0.4596	0.3564		110.0	0.00994	-64.163	0.4596	0.3564
	120.0	0.01015	-60.615	0.4907	0.3594	400.000	120.0	0.01015	-60.615	0.4907	0.3594
1000.000	130.0	0.01038	-57.042	0.5194	0.3595		130.0	0.01038	-57.042	0.5194	0.3595
	140.0	0.01061	-53.452	0.5459	0.3601	400.000	140.0	0.01061	-53.452	0.5459	0.3601
	150.0	0.01087	-49.833	0.5709	0.3629		150.0	0.01087	-49.833	0.5709	0.3629
	160.0	0.01115	-46.219	0.5944	0.3624		160.0	0.01115	-46.219	0.5944	0.3624
1000.000	170.0	0.01147	-42.505	0.6168	0.3713		170.0	0.01147	-42.505	0.6168	0.3713
	180.0	0.01161	-38.797	0.6383	0.3801		180.0	0.01161	-38.797	0.6383	0.3801
	100.0	0.00973	-67.383	0.4256	0.3511		100.0	0.00973	-67.383	0.4256	0.3511
	110.0	0.00994	-63.873	0.4591	0.3547		110.0	0.00994	-63.873	0.4591	0.3547
1000.000	120.0	0.01016	-60.323	0.4901	0.3580		120.0	0.01016	-60.323	0.4901	0.3580
	130.0	0.01036	-56.767	0.5185	0.3597	1000.000	130.0	0.01036	-56.767	0.5185	0.3597
	140.0	0.01059	-53.197	0.5450	0.3614		140.0	0.01059	-53.197	0.5450	0.3614
	150.0	0.01085	-49.563	0.5693	0.3611		150.0	0.01085	-49.563	0.5693	0.3611
1000.000	160.0	0.01112	-45.884	0.5933	0.3699		160.0	0.01112	-45.884	0.5933	0.3699
	170.0	0.01143	-42.204	0.6157	0.3770		170.0	0.01143	-42.204	0.6157	0.3770
	180.0	0.01176	-38.513	0.6369	0.3847		180.0	0.01176	-38.513	0.6369	0.3847
	100.0	0.00971	-66.666	0.4232	0.3497		100.0	0.00971	-66.666	0.4232	0.3497
1000.000	110.0	0.00990	-63.124	0.4567	0.3520		110.0	0.00990	-63.124	0.4567	0.3520
	120.0	0.01010	-59.580	0.4881	0.3560		120.0	0.01010	-59.580	0.4881	0.3560
	130.0	0.01032	-56.037	0.5164	0.3599		130.0	0.01032	-56.037	0.5164	0.3599
	140.0	0.01055	-52.423	0.5429	0.3613		140.0	0.01055	-52.423	0.5429	0.3613
1000.000	150.0	0.01079	-48.745	0.5670	0.3605		150.0	0.01079	-48.745	0.5670	0.3605
	160.0	0.01113	-45.019	0.5897	0.3663		160.0	0.01113	-45.019	0.5897	0.3663
	170.0	0.01146	-41.247	0.6128	0.3730		170.0	0.01146	-41.247	0.6128	0.3730
	180.0	0.01179	-37.417	0.6338	0.3798		180.0	0.01179	-37.417	0.6338	0.3798

TABLE B-2. THERMODYNAMIC PROPERTIES OF FLUORINE-OXYGEN MIXTURES IN THE COMPRESSED LIQUID REGION

WT. FRACTION F2 =1.00

TEMPERATURE	R	VOLUME	ENTHALPY	ENTROPY	TEMPERATURE	VOLUME	ENTHALPY	ENTROPY	C _p
PSIA	FT ³ /LB	BTU/LB/R	BTU/LB/R	BTU/LB/R	PSIA	FT ³ /LB	BTU/LB/R	BTU/LB/R	BTU/LB/R
8.000	100.0	0.009346	-65.920	0.3959	100.0	0.009346	-66.311	0.3958	0.3505
	110.0	0.009666	-63.443	0.4290	110.0	0.009666	-63.440	0.4289	0.3515
	120.0	0.009886	-59.939	0.4595	120.0	0.009886	-59.930	0.4595	0.3557
	130.0	0.01009	-56.397	0.4878	130.0	0.01032	-56.388	0.4877	0.3551
	140.0	0.01032	-52.838	0.5142	140.0	0.01032	-52.830	0.5142	0.3549
	150.0	0.01058	-49.261	0.5389	150.0	0.01058	-49.252	0.5388	0.3578
	160.0	0.01086	-45.592	0.5623	160.0	0.01117	-45.570	0.5622	0.3627
	170.0	0.01116	-41.841	0.5841	170.0	0.01117	-41.832	0.5845	0.3729
	180.0	0.01151	-38.093	0.6055	180.0	0.01149	-38.080	0.6049	0.3784
30.000	100.0	0.00946	-65.792	0.3956	100.0	0.00946	-66.552	0.3952	0.3487
	110.0	0.00955	-63.321	0.4286	110.0	0.00955	-63.183	0.4282	0.3519
	120.0	0.00986	-59.813	0.4592	120.0	0.00986	-59.677	0.4587	0.3569
	130.0	0.01008	-56.275	0.4875	130.0	0.01007	-56.143	0.4870	0.3550
	140.0	0.01032	-52.720	0.5138	140.0	0.01030	-52.592	0.5134	0.3555
	150.0	0.01057	-49.154	0.5385	150.0	0.01056	-49.011	0.5380	0.3584
	160.0	0.01085	-45.581	0.5619	160.0	0.01074	-45.400	0.5614	0.3608
	170.0	0.01116	-41.941	0.5841	170.0	0.01115	-41.738	0.5836	0.3681
	180.0	0.01151	-38.293	0.6055	180.0	0.01149	-38.080	0.6049	0.3784
100.000	100.0	0.00945	-65.511	0.3948	100.0	0.00944	-66.371	0.3945	0.3469
	110.0	0.00964	-63.041	0.4279	110.0	0.00944	-63.307	0.4275	0.3524
	120.0	0.00984	-59.541	0.4583	120.0	0.00944	-59.392	0.4581	0.3544
	130.0	0.01006	-56.010	0.4866	130.0	0.01035	-55.378	0.4862	0.3559
	140.0	0.01029	-52.463	0.5129	140.0	0.01039	-51.335	0.5124	0.3559
	150.0	0.01055	-48.919	0.5376	150.0	0.01073	-47.266	0.5371	0.3575
	160.0	0.01082	-45.382	0.5608	160.0	0.01111	-43.170	0.5603	0.3668
	170.0	0.01113	-41.832	0.5829	170.0	0.01115	-41.526	0.5824	0.3751
	180.0	0.01147	-38.281	0.6042	180.0	0.01145	-38.624	0.6035	0.3791
300.000	100.0	0.00944	-65.231	0.3941	100.0	0.00941	-65.529	0.3934	0.3430
	110.0	0.00963	-62.761	0.4277	110.0	0.00941	-62.777	0.4273	0.3487
	120.0	0.00983	-59.265	0.4577	120.0	0.00941	-59.575	0.4568	0.3532
	130.0	0.01006	-55.742	0.4856	130.0	0.01030	-55.384	0.4848	0.3561
	140.0	0.01026	-52.203	0.5120	140.0	0.01033	-51.562	0.5133	0.3559
	150.0	0.01049	-48.633	0.5367	150.0	0.01073	-44.718	0.5374	0.3574
	160.0	0.01076	-45.053	0.5597	160.0	0.01113	-41.463	0.5592	0.3654
	170.0	0.01109	-41.473	0.5818	170.0	0.01142	-38.240	0.5790	0.3673
	180.0	0.01142	-37.893	0.6029	180.0	0.01142	-37.893	0.5997	0.3733

U.S. DEPT. OF COMM. BIBLIOGRAPHIC DATA SHEET	1. PUBLICATION OR REPORT NO. NBSIR 73-338	2. Gov't Accession No.	3. Recipient's Accession No.
4. TITLE AND SUBTITLE Calculated Liquid Phase Thermodynamic Properties and Liquid-Vapor Equilibria for Fluorine-Oxygen (FLOX) Mixtures		5. Publication Date September 1973	6. Performing Organization Code
7. AUTHOR(S) W. R. Parrish and M. J. Hiza		8. Performing Organization	
9. PERFORMING ORGANIZATION NAME AND ADDRESS NATIONAL BUREAU OF STANDARDS, Boulder Laboratories DEPARTMENT OF COMMERCE Boulder, Colorado 80302		10. Project/Task/Work Unit No. 2750547	11. Contract/Grant No. Purchase Req. C-32369-C
12. Sponsoring Organization Name and Address NASA Lewis Research Center 21000 Brookpark Rd. Cleveland, Ohio 44135		13. Type of Report & Period Covered Final Nov '72 -- Sept '73	14. Sponsoring Agency Code
15. SUPPLEMENTARY NOTES			
<p>16. ABSTRACT (A 200-word or less factual summary of most significant information. If document includes a significant bibliography or literature survey, mention it here.)</p> <p>Liquid phase thermodynamic properties and liquid-vapor equilibria of fluorine-oxygen mixtures, for which no experimental data exist, have been calculated. The results are based on excess properties predicted from the Snider-Herrington equations, with an adjusted combining rule, and the corresponding data for the pure fluids. Mixtures considered are 0.6, 0.7, 0.8, 0.88, and 0.9 weight fraction of fluorine from 55 to 90 K up to 70×10^5 Pa. In the compressed liquid, molar volumes, enthalpy, entropy, and constant pressure specific heat were determined. Along the saturation boundary, coexistent vapor compositions and solution vapor pressures were determined as well. Corresponding properties of pure fluorine from experimental data have also been included. Results are tabulated in both British and S.I. units.</p>			
17. KEY WORDS (Alphabetical order, separated by semicolons) Calculated thermophysical properties; compressed liquid phase, fluorine-oxygen mixtures, hard-sphere model; liquid-vapor equilibria.			
18. AVAILABILITY STATEMENT <input checked="" type="checkbox"/> UNLIMITED. <input type="checkbox"/> FOR OFFICIAL DISTRIBUTION. DO NOT RELEASE TO NTIS.		19. SECURITY CLASS (THIS REPORT) UNCLASSIFIED	21. NO. OF PAGES
		20. SECURITY CLASS (THIS PAGE) UNCLASSIFIED	22. Price

