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Selected Values of Chemical Thermodynamic Properties

Tables for the Alkaline Earth Elements
(Elements 92 through 97 in the
Standard Order of Arrangement)

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no. 270-6

1971

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UNITED STATES DEPARTMENT OF COMMERCE
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Selected Values of Chemical Thermodynamic Properties

Tables for the Alkaline Earth Elements
(Elements 92 through 97 in the
Standard Order of Arrangement)

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ABSTRACT

Contains tables of values for the standard heats and Gibbs (free) energies of formation, entropies and enthalpies at 298.15 K and heats of formation at 0 K for compounds of beryllium, magnesium, calcium, strontium, barium, and radium (elements 92-97 in the Standard Order of Arrangement). These tables are a continuation of the comprehensive revision of NBS Circular 500.

Key words: Enthalpy; entropy; Gibbs energy of formation; beryllium compounds; magnesium compounds; calcium compounds; strontium compounds; barium compounds; radium compounds.

PREFACE

This is the sixth of a series of Technical Notes containing the tables of material prepared as a revision of National Bureau of Standards Circular 500, Selected Values of Chemical Thermodynamic Properties, by F. D. Rossini, D. D. Wagman, W. H. Evans, S. Levine and I. Jaffe. This Note contains data for compounds of six elements, numbered 92 through 97 in the Standard Order of Arrangement.

The continued encouragement and support of the Office of Standard Reference Data of the National Bureau of Standards is gratefully acknowledged.

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SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES

INTRODUCTION

Substances and Properties Included in the Tables

The tables contain values where known of the enthalpy and Gibbs energy of formation, enthalpy, entropy and heat capacity at 298.15 K (25°C), and the enthalpy of formation at 0 K, for all inorganic substances and organic molecules containing not more than two carbon atoms. In some instances such as metal-organic compounds, data are given for substances in which each organic radical contains one or two carbon atoms.

No values are given in these tables for metal alloys or other solid solutions, fused salts, or for substances of undefined chemical composition.

Physical States

The physical state of each substance is indicated in the column headed "State" as crystalline solid (c), liquid (liq), glassy or amorphous (amorp), or gaseous (g). Solutions in water are listed as aqueous (aq). For non-aqueous systems the physical state is that normal for the indicated solvent at 298.15 K.

Definition of Symbols

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

Conventions Regarding Pure Substances

The values of the thermodynamic properties of the pure substances given in these tables are for the substances in their standard states (indicated by the superscript ° on the thermodynamic symbol). These standard states are defined as follows:

For a pure solid or liquid, the standard state is the substance in the condensed phase under a pressure of one atmosphere.

For a gas the standard state is the hypothetical ideal gas at unit fugacity, in which state the enthalpy is that of the real gas at the same temperature and at zero pressure.

The values of ΔH_f° and ΔG_f° given in the tables represent the change in the appropriate thermodynamic quantity when one gram-formula weight of the substance in its standard state is formed, isothermally at the indicated temperature, from the elements, each in its appropriate standard reference state. The standard reference state at 25°C for each element except phosphorus has been chosen to be the standard state that is thermodynamically stable at 25°C and at one atmosphere pressure. For phosphorus the standard reference state is the crystalline white form; the more stable forms have not been well characterized thermochemically. The same reference states have been maintained for the elements at 0 K except for the liquid elements bromine and mercury, for which the reference states have been chosen as the stable crystalline forms. The standard reference states are indicated in the tables by the fact that the values of ΔH_f° and ΔG_f° are exactly zero.

The values of $H_{298}^\circ - H_0^\circ$ represent the enthalpy difference for the given substance between 298.15 K and 0 K. If the indicated standard state at 25°C is the gas, the corresponding state at 0 K is the hypothetical ideal gas; if the state at 25°C is solid or liquid, the corresponding state at 0 K is the thermodynamically stable crystalline solid, unless otherwise specifically indicated.

The values of S° represent the virtual or "thermal" entropy of the substance in the standard state at 298.15 K, omitting contributions from nuclear spins. Isotope mixing effects, etc., are also excluded except in the case of the hydrogen-deuterium (^1H - ^2H) system. Where data have been available only for a particular isotope, they have been corrected when possible to the normal isotopic composition.

The values of the enthalpies of formation of gaseous ionic species are computed on the convention that the value of ΔH_f° for the electron is zero. Conversions between 0 and 298.15 K are calculated using the value of $H_{298}^\circ - H_0^\circ = 1.481$ kcal per mole of electrons, and assuming that the values of $H_{298}^\circ - H_0^\circ$ for the ionized and un-ionized molecules are the same.

Conventions Regarding Solutions

Solutions in water are designated as aqueous (aq); other solvents are designated by name or chemical formula. The concentration of the solution is expressed in terms of the number of moles of solvent associated with one mole of the solute. If no concentration is indicated, the solution is assumed to be "dilute".

The standard state for a solute in aqueous solution is taken as the hypothetical ideal solution of unit molality (indicated as "std. state, $m = 1$ "). In this state the partial molal enthalpy and heat capacity of the solute are the same as in the infinitely dilute real solution. For non-aqueous solutions the standard state of the solute is the hypothetical ideal solution of unit mole fraction of solute ("std. state, $x_2 = 1$ ").

The value of ΔH_f° given in the tables for a solute in its standard state is the apparent molal enthalpy of formation of the substance in the infinitely dilute real solution. At this dilution the partial molal enthalpy is equal to the apparent molal quantity. At concentrations other than the standard state, the value of ΔH_f° represents the apparent enthalpy of the reaction of formation of the solution from the elements comprising the solute, each in its standard reference state, and the appropriate total number of moles of solvent. In this representation the value of ΔH_f° for the solvent is not required. The experimental value for a heat of dilution is obtained directly as the difference between the two values of ΔH_f° at the corresponding concentrations.

The values of the thermodynamic properties tabulated for the individual ions in aqueous solution are based on the usual convention that the values of ΔH_f° , ΔG_f° , S° and C_p° for H^+ (aq, std. state, $m = 1$) are zero. The properties of a neutral electrolyte in aqueous solution in the standard state are equal to the algebraic sum of these values for the appropriate kinds and number of individual ions assumed to constitute the molecule of the given electrolyte. When the undissociated species, rather than the sum of the ions, is meant, the notation "undissociated" or "un-ionized" is used. For an ionic species the properties tabulated refer to that undissociated ion. By adopting the above convention with respect to aqueous H^+ , it follows that the thermodynamic relation $\Delta G_f^\circ = \Delta H_f^\circ - T \Delta S_f^\circ$ will not hold for an individual ionic species. However no problem arises when neutral chemical systems are considered.

Unit of Energy and Fundamental Constants

All of the energy values given in these tables are expressed in terms of the thermochemical calorie. This unit, defined as equal to 4.1840 joules, is generally accepted for the presentation of chemical thermodynamic data. Values reported in other units have been converted to calories by means of the conversion factors for molecular energy given in Table A.

The following values of the fundamental physical constants have been used in these calculations:

$$R = \text{gas constant} = 8.3143 \pm 0.0012 \text{ J/deg mol} = 1.98717 \pm 0.00029 \text{ cal/deg mol}$$

$$F = \text{Faraday constant} = 96487.0 \pm 1.6 \text{ coulombs/mol} \\ = 23060.9 \pm 0.4 \text{ cal/volt equivalent}$$

$$Z = Nhc = 11.96258 \pm 0.00107 \text{ J/cm}^{-1} \text{ mol} = 2.85912 \pm 0.00026 \text{ cal/cm}^{-1} \text{ mol}$$

$$c_2 = \text{second radiation constant} = hc/k = 1.43879 \pm 0.00015 \text{ cm deg}$$

$$0^\circ\text{C} = 273.15 \text{ K}$$

These constants are consistent with those given in the Table of General Physical Constants, recommended by the National Academy of Sciences - National Research Council¹. The formula weights in the tables have been calculated for the molecular formula given in the Formula and Description column using the 1961 Table of Relative Atomic Weights based on the atomic mass of ¹²C = 12 exactly².

Internal Consistency of the Tables

All of the values given in these tables have been calculated from the original articles, using consistent values for all subsidiary and auxiliary quantities. The original data were corrected where possible for differences in energy units, molecular weights, temperature scales, etc. Thus we have sought to maintain a uniform scale of energies for all the substances in the tables. In addition the tabulated values of the properties of a substance satisfy all the known physical and thermodynamic relationships among these properties. The quantities ΔH_f° , ΔG_f° , and S° at 298.15 K satisfy the relation:

$$\Delta G_f^\circ = \Delta H_f^\circ - T \Delta S_f^\circ.$$

¹NBS Technical News Bulletin, October 1963.

²A. E. Cameron and E. Wichers, J. Am. Chem. Soc. 84, 4192 (1962).

Furthermore the calculated value of any thermodynamic quantity for a reaction is independent of the path chosen for the evaluation.

In some cases newer data may have become available on certain substances after the values were selected for these tables. Because of the need to maintain the internal consistency of the tables, it is not always possible to incorporate these newer data into the tables without a detailed analysis of the effect of such a change. Unless great care is used, relatively significant errors in calculated values of ΔH° or ΔG° for specific reactions may result from the introduction of such data.

Uncertainties

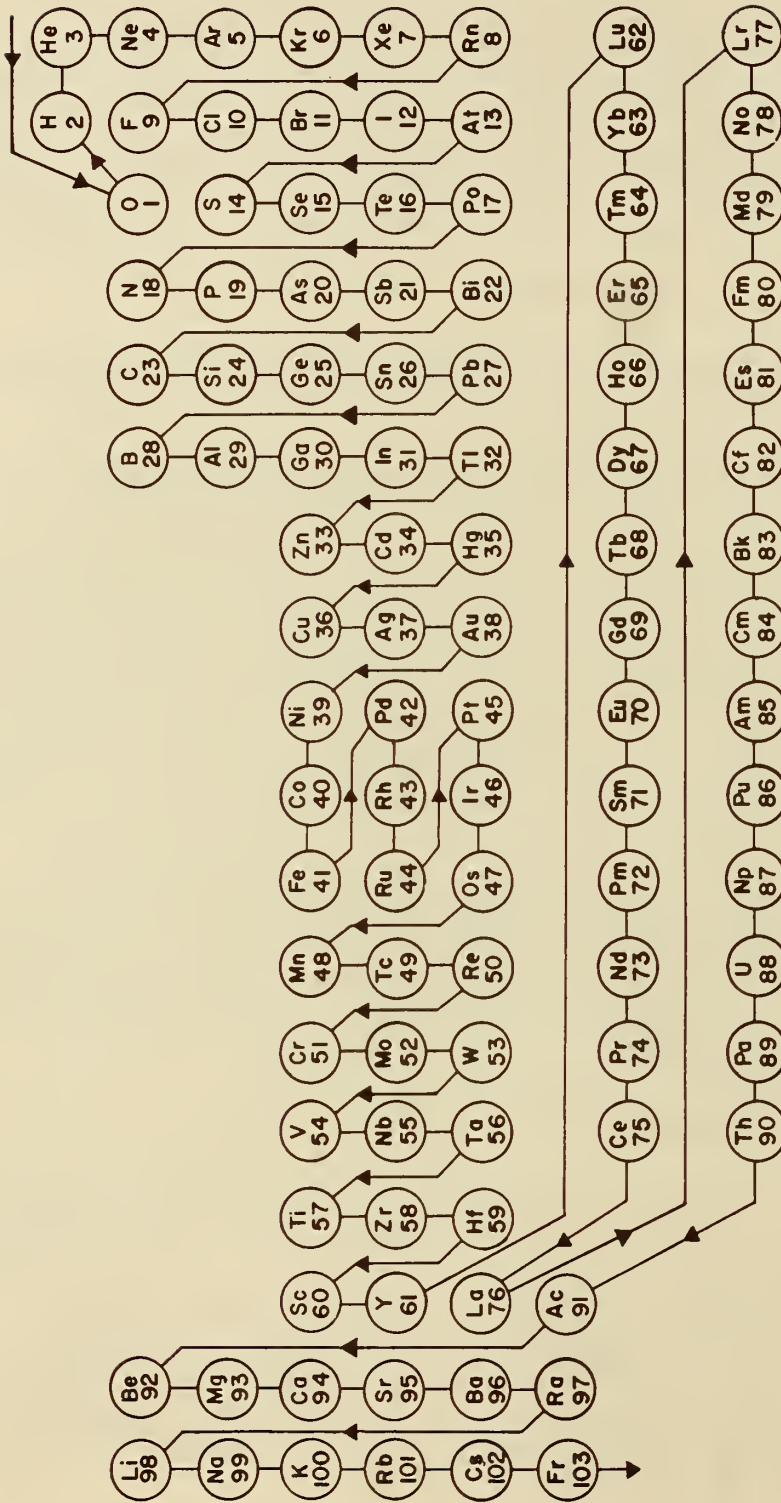
The uncertainty in any value in the tables depends on the uncertainties of all the determinations in the total chain of reactions used to establish the value.

A discussion of the uncertainties will be included in the final publication of these tables in the National Standard Reference Data System. However we have followed certain rules with respect to significant figures to indicate these uncertainties. Values are tabulated in general such that the overall uncertainty lies between 2 and 20 units of the last figure. On the other hand, values are given so that the experimental data from which they are derived may be recovered with an accuracy equal to that of the original quantities. Thus the number of significant figures for any one value in the tables need not represent the absolute accuracy of that value. For solutions of varying composition values are frequently tabulated to more figures to make possible the recovery of enthalpies of solution and dilution. Similarly values of ΔH_f° and $\Delta H_f^\circ_{298.15}$ may be given to different numbers of significant figures. In this instance the quantity with the lesser number of figures is used to represent the uncertainty estimate. The larger number of figures is used for the other quantity to retain the significance of the temperature correction term.

Arrangement of the Tables

The compounds in the tables are entered according to the Standard Order of Arrangement, (see Figure 1), by the principle of latest position. In this scheme, a compound is listed under the element occurring latest in the list; water of hydration is neglected. Within a given element-table will be found all of the compounds of that element with elements occurring earlier in the order; the arrangement within a table follows the same ordering. An exception occurs in the carbon tables (Table 23), which is divided into subgroups consisting of all compounds with one carbon atom, then all with two carbon atoms, etc.

STANDARD ORDER OF ARRANGEMENT



Standard Order of Arrangement of the Elements and Compounds
based on the
Periodic Classification of the Elements

Figure 1

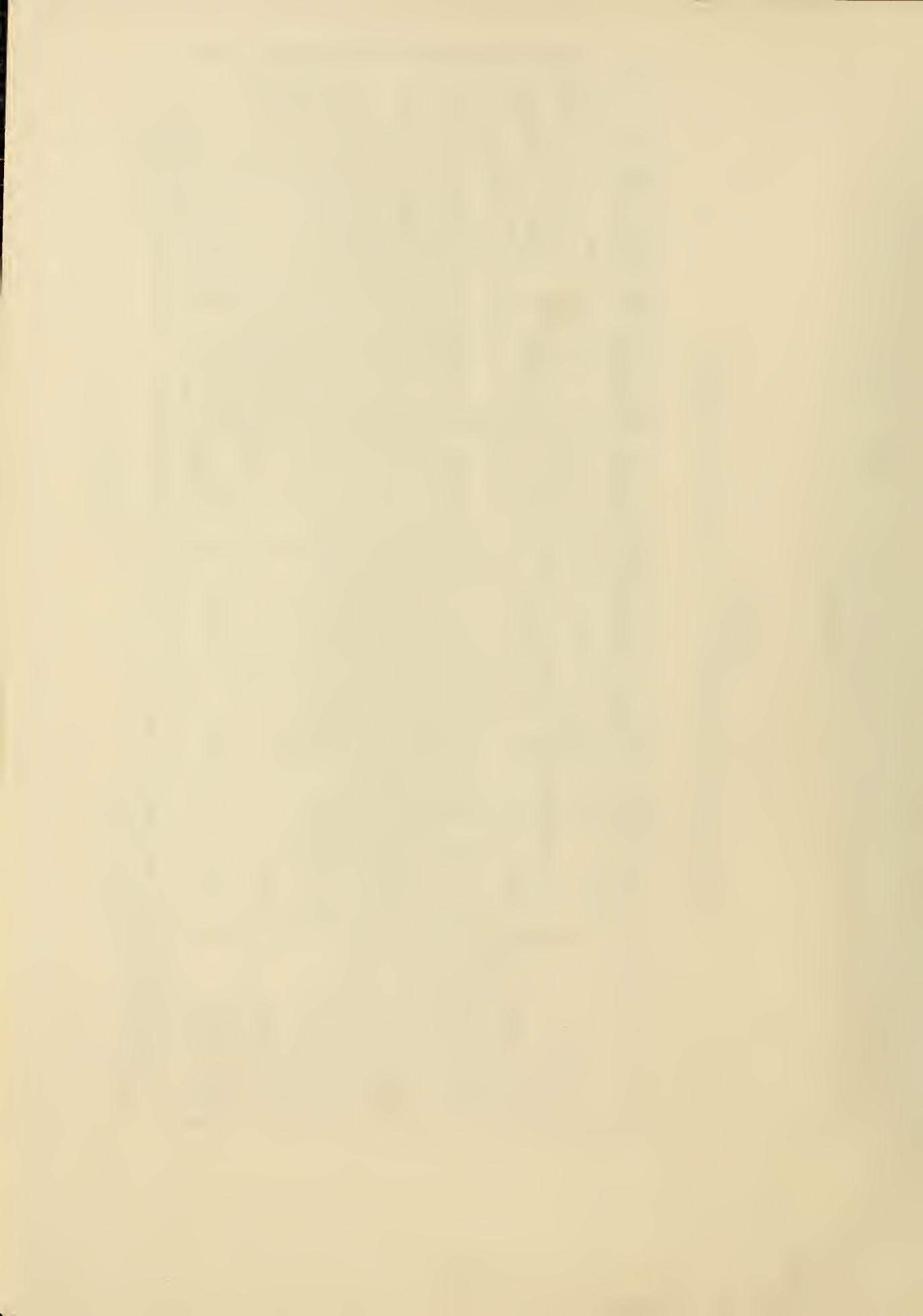
TABLE A
CONVERSION FACTORS FOR UNITS OF MOLECULAR ENERGY

	J/mol	cal/mol	cm ³ atm/mol	kWh/mol	Btu/lb-mol	cm ⁻¹ /molecule	eV/molecule
1 J/mol =	1	2.390057 x 10 ⁻¹	9.86923	2.77778 x 10 ⁻⁷	0.429923	8.35940 x 10 ⁻²	1.036409 x 10 ⁻⁵
1 cal/mol =	<u>4.18400</u>	1	41.2929	1.162222 x 10 ⁻⁶	1.798796	3.49757 x 10 ⁻¹	4.33634 x 10 ⁻⁵
1 cm ³ atm/mol =	<u>0.1013250</u>	2.42173 x 10 ⁻²	1	2.81458 x 10 ⁻⁸	4.35619 x 10 ⁻²	8.47016 x 10 ⁻³	1.050141 x 10 ⁻⁶
1 kWh/mol =	<u>3,600,000</u>	860,421	3.55292 x 10 ⁷	1	1,547,721	300,938	37.3107
1 Btu/lb-mol =	<u>2,32600</u>	5.55927 x 10 ⁻¹	22.9558	6.46111 x 10 ⁻⁷	1	1.944396 x 10 ⁻¹	2.41069 x 10 ⁻⁵
1 cm ⁻¹ /molecule =	11.96258	2.85912	118.0614	3.32294 x 10 ⁻⁶	5.14299	1	1.239812 x 10 ⁻⁴
1 eV/molecule =	<u>96487.0</u>	23060.9	952,252	2.68019 x 10 ⁻²	41482.0	8065.73	1

The underlined numbers represent the fundamental values used in deriving this table. The remaining factors were obtained by applying the relationships:

$$n_{ij} = n_{ik} \cdot n_{kj} = 1$$

$$n_{ij} = n_{ik} \cdot n_{kj}$$



TABLES OF SELECTED VALUES OF PROPERTIES

SERIES I

Enthalpy of Formation at 0 K

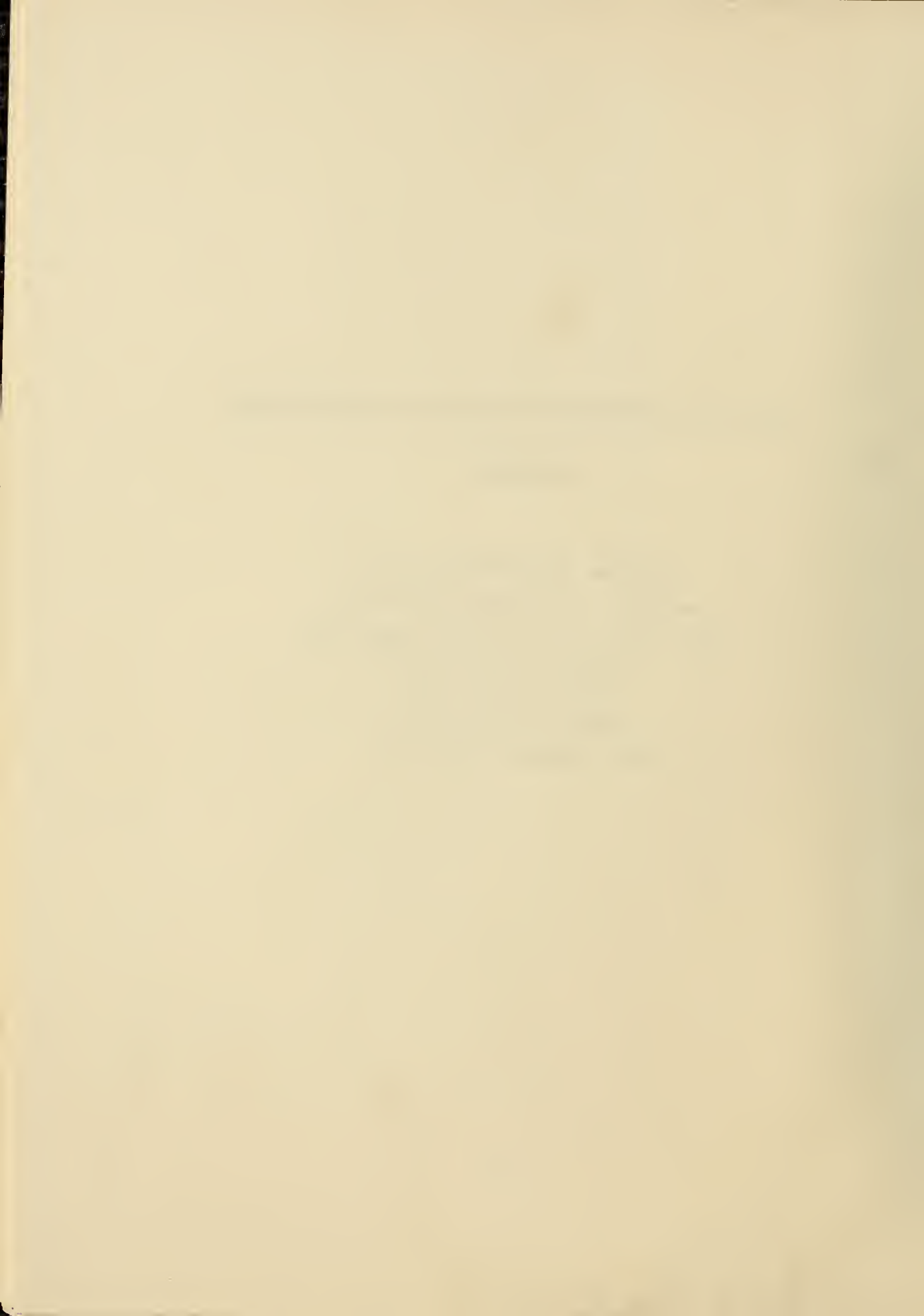
Enthalpy of Formation at 298.15 K

Gibbs Energy of Formation at 298.15 K

Enthalpy at 298.15 K

Entropy at 298.15 K

Heat Capacity at 298.15 K



SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Beryllium

Table 92(1)

Substance	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$		S°	C_p°
						0 K	298.15 K (25°C)		
Be	c	9.0122	0	0	0	0.466	2.27	3.93	
Be ⁺	g		76.49	77.5	68.5	1.481	32.543	4.968	
Be ²⁺	g		291.474	293.965					
Be ³⁺	g		711.426	715.398					
Be ⁴⁺	aq			-91.5	-90.75			-31.0	
Be	g		4260.24	4265.69					
Be	g		9280.86	9287.80					
BeO	c	25.0116	-144.86	-145.7	-138.7	0.669	3.38	6.10	
BeO ₂ ²⁻	g			28.					
Be ₂ O	aq	41.0110		-189.0	-153.0			-38.	
(BeO) ₂	g	34.0238		-20.					
(BeO) ₃	g	50.0232		-103.					
(BeO) ₄	g	75.0348		-260.					
(BeO) ₅	g	100.0464		-391.					
(BeO) ₆	g	125.0580		-518.					
(BeO) ₆	g	150.0696		-651.					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

BeH
92

Table 92(2)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Substance		Beryllium						
Formula and Description	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
			0 K	298.15 K (25°C)				
		kcal/mol						
		cal/deg mol						
BeH	g	10.0202	75.	75.6	68.3	2.062	42.21	6.95
BeH ₂	c			-4.60				
Be(OH) ₂	c	43.0269		-215.7	-194.8		12.4	
	c			-216.5	-195.4		12.	
	amorp			-214.6				
	g			-158.				
	amorp	56.5		-266.9	-430.6			
Be(OH) ₂ •3/4H ₂ O	g	78.0587						
Be ₃ (OH) ₃ ⁺	aq							
	std. state, m = 1							
BeF	g	28.0106	-42.41	-41.8	-48.6	2.082	49.15	7.14
BeF ₂	c	47.0090	-244.85	-245.4	-234.1	2.024	12.75	12.39
	c			-244.7				
	amorp			-244.3				
	g			-189.7				
	aq			-250.0				
	aq			-251.40				
	aq			-251.78				
	aq			-252.38				
	aq			-252.84				
	aq			-253.15				
	in 100 (HF + 1.5 H ₂ O)							
	100 (HF + 2.0 H ₂ O)							
	100 (HF + 3.0 H ₂ O)							
	100 (HF + 4.0 H ₂ O)							
	100 (HF + 5.0 H ₂ O)							

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Beryllium

Table 92(3)

Substance		Formula Weight	State	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
Formula and Description								
BeF ₂	in 100 (HF + 6.0 H ₂ O)		aq	-253.32				
	100 (HF + 7.0 H ₂ O)		aq	-253.40				
	100 (HF + 8.0 H ₂ O)		aq	-253.46				
Be ₂ O ₃		72.0206	g	-292.				
BeCl		44.4652	g	14.	7.	2.12	52.0	7.56
BeCl ₂	α	79.9182	c	-117.40	-106.5	2.863	19.76	15.50
	β		c	-118.57	-107.3	2.729	18.12	14.92
BeCl ₂ ·4H ₂ O	dilute		g	-85.7				
	in 45.8 (HCl + 7.25 H ₂ O)		aq	-171.1				
	38.3 (HCl + 9.72 H ₂ O)		aq	-162.0				
	6.38 (HCl + 10.65 H ₂ O)		aq	-164.3				
Be ₂ Cl ₄	in C ₂ H ₅ OH, ethyl alcohol		aq	-163.6				
		151.9796	c	-155.				
Be ₂ Cl ₄		159.8364	g	-432.2				
			g	-202.				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

BeBr₂

92

Substance		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Beryllium							
Formula and Description	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°	
			0 K	298.15 K (25°C)					
				kcal/mol					
				cal/deg mol					
BeBr ₂	c	168.8302		-84.5					
	aq			-140.2					
BeI ₂	c	262.8210		-46.0					
	aq			-108.5					
BeS	c	41.0762		-56.0					
BeSO ₄	α , tetragonal	105.0738	-285.505	-288.05	-261.44	3.125	18.62	20.48	
	std. state, m = 1			-308.8	-268.72				
	in 15 H ₂ O			-304.77					
	20 H ₂ O			-305.33					
	25 H ₂ O			-305.68					
	50 H ₂ O			-306.36					
	75 H ₂ O			-306.61					
	100 H ₂ O			-306.76					
	200 H ₂ O			-307.06					
	400 H ₂ O			-307.34					
	500 H ₂ O			-307.43					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Beryllium

Table 92(5)

Substance		Formula Weight	State	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
Formula and Description								
BeSO ₄	in 1,000 H ₂ O		aq					
	2,000 H ₂ O		aq	-307.73				
	3,000 H ₂ O		aq	-308.00				
	4,000 H ₂ O		aq	-308.20				
	5,000 H ₂ O		aq	-308.26				
in 100 (HCl + 55.5 H ₂ O)			aq	-308.37				
	200 (HCl + 27.75 H ₂ O)		aq	-295.0				
	300 (HCl + 18.5 H ₂ O)		aq	-294.6				
	400 (HCl + 13.88 H ₂ O)		aq	-294.1				
BeSO ₄ ·H ₂ O		123.0891	c	-364.2				
	BeSO ₄ ·2H ₂ O	141.1045	c	-435.74	-381.99	5.865	39.01	36.63
BeSO ₄ ·4H ₂ O		177.1352	c	-569.682	-497.29	8.306	55.68	51.77
	tetragonal							
BeBr ₂ ·2H ₂ S		236.9901	c	-112.1				
BeI ₂ ·2H ₂ S		330.9809	c	-69.8				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

BeSeO₄
92

Table 92(6)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Beryllium						
Substance	State	Formula Weight	ΔH_f° 0 K	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
Formula and Description								
BeSeO ₄	c	151.970		-212.8				
std. state, m = 1	aq			-234.7				
in 800 H ₂ O	aq			-234.1	-196.2			
BeSeO ₄ ·2H ₂ O	c	188.000		-360.3				
BeSeO ₄ ·4H ₂ O	c	224.031		-505.0				
Be ₃ N ₂	c	55.0500		-140.6				
α, cubic	c			-136.5				
β, hexagonal								
Be(NO ₃) ₂	aq	133.0220		-191.0				
BeCl ₂ ·2NH ₃	c	113.9794		-201.7				
BeCl ₂ ·4NH ₃	c	148.0406		-260.4				
BeCl ₂ ·6NH ₃	c	182.1019		-298.2				
BeCl ₂ ·12NH ₃	c	284.2855		-409.9				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Table 92(7)

Beryllium

Substance	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C° _p
Formula and Description	kcal/mol							
BeBr ₂ •4NH ₃	c	236.9526		-235.1				
BeBr ₂ •6NH ₃	c	271.0138		-274.4				
BeBr ₂ •10NH ₃	c	339.1363		-349.2				
BeI ₂ •4NH ₃	c	330.9434		-207.6				
BeI ₂ •6NH ₃	c	365.0047		-247.2				
BeI ₂ •13NH ₃	c	484.2189		-378.4				
Be ₂ C	c	30.0356		-28.0				
BeCO ₃	c	69.0216		-245.				
Be ₂ SiO ₄	c	110.1080	-510.77	-513.7	-485.8	2.922	15.37	22.84
Be(BO ₂) ₂	g	94.6318		-325.				
Be ₃ P ₂ O ₆	c	144.6550		-750.3				

Be₃P₂O₆
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SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

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BeOAl
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Table 92(8)		Beryllium							
Substance		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity							
Formula and Description	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °	
			0 K	298.15 K	298.15 K	298.15 K			
			kcal/mol					cal/deg mol	
BeOAl	g	51.9931		-5.					
BeO·Al ₂ O ₃ chrysobery1	c	126.9728	-546.32	-549.9	-520.7	3.128	15.84	25.19	
BeO·3Al ₂ O ₃	c	330.8952	-1335.65	-1344.9	-1271.6	8.153	42.0	63.38	
BeMoO ₄	c	168.950		-328.					
NbBe ₂	c	110.9304		-14.6					
NbBe ₈	c	165.0036		-20.					
NbBe ₁₂	c	201.0524		-20.					
UBe ₁₃	c	355.189		-39.					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

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Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Magnesium

Table 93(1)

Substance	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
			0 K	298.15 K (25°C)	cal/deg mol			
Mg	c	24.312	0	0	0	1.195	7.81	5.95
Mg ⁺	g		35.014	35.30				
Mg ²⁺	g		211.333	213.100				
Mg ³⁺	g		558.052	561.299				
Mg ⁴⁺	g			-111.58				
Mg ⁵⁺	g		2406.08	2410.81				
Mg ⁶⁺	g		4927.14	4933.35				
Mg ⁷⁺	g		8184.9	8192.6				
Mg ⁸⁺	g		12486.7	12495.8				
Mg ⁹⁺	g		17674.3	17685.0				
Mg ¹⁰⁺	g		23809.1	23821.2				
Mg ¹¹⁺	g		31372.7	31386.3				
Mg ₂	g	48.624	39847.	39862.				
	g		80472.	80489.				
	g		+68.87	68.59				
	aq				-108.7		-33.0	
	g							4.968

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MgO
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Table 93(2)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity							
		Magnesium							
Substance	Formula and Description	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
				0 K	298.15 K (25°C)				
MgO	macrocrystal (periclase)	c	40.3114	-142.813	-143.81	-136.10	1.235	6.44	8.88
	microcrystal	c		-141.954	-142.92 +4.	-135.27	1.266	6.67	9.00
MgH ₂		g		-16.05	-18.0	-8.6	1.270	7.43	8.45
MgOH ⁺	std. state, m = 1	aq	41.3194			-149.8			
Mg(OH) ₂	precipitate	c	58.3267	-218.402	-220.97	-199.23	2.725	15.10	18.41
	std. state, m = 1	amorp			-220.0				
		g			-134.				
		aq			-221.52			-35.6	
MgF		g	43.3104	-52.89	-53.0	-59.2	2.143	52.79	7.78
MgF ₂		c	62.3088	-267.57	-268.5	-255.8	2.370	13.68	14.72
		g		-172.58	-173.0	-174.6	2.878	61.73	11.62
MgCl		g	59.765	-9.9	-10.	-16.	2.24	55.8	8.32
MgCl ₂		c	95.218	-153.180	-153.28	-141.45	3.288	21.42	17.06
		g			-95.7				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

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Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Table 93(3) Magnesium

Substance	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
Formula and Description			0 K	298.15 K (25°C)				
MgCl ₂	aq							
std. state, m = 1	aq			-191.48	-171.4		-6.0	
in	aq			-186.82				
10 H ₂ O	aq			-187.58				
12 H ₂ O	aq			-188.45				
15 H ₂ O	aq			-189.17				
20 H ₂ O	aq			-189.57				
25 H ₂ O	aq			-190.23				
50 H ₂ O	aq			-190.46				
75 H ₂ O	aq			-190.59				
100 H ₂ O	aq			-190.81				
200 H ₂ O	aq			-190.89				
300 H ₂ O	aq			-190.94				
400 H ₂ O	aq			-190.98				
500 H ₂ O	aq			-191.003				
600 H ₂ O	aq			-191.048				
800 H ₂ O	aq			-191.080				
1,000 H ₂ O	aq			-191.176				
2,000 H ₂ O	aq			-191.269				
5,000 H ₂ O	aq			-191.323				
10,000 H ₂ O	aq			-191.364				
20,000 H ₂ O	aq							

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MgCl₂
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Table 93(4)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Magnesium							
Substance	Formula and Description	State	Formula Weight	ΔHf°	ΔHf°	ΔGf°	H ₂₉₈ ^o - H ₀ ^o	S°	C _p ^o
				0 K	298.15 K (25°C)				
MgCl ₂	in 50,000 H ₂ O	aq			-191.404				
	100,000 H ₂ O	aq			-191.425				
	∞ H ₂ O	aq			-191.48				
	in 75.2 HCl + 4190 H ₂ O	aq			-189.84				
Mg Cl ₂ · H ₂ O		c	113.2333		-231.03	-205.98		32.8	27.55
MgCl ₂ · 2H ₂ O		c	131.2487		-305.86	-267.24		43.0	38.05
MgCl ₂ · 4H ₂ O		c	167.2793		-453.87	-388.03		63.1	57.70
MgCl ₂ · 6H ₂ O		c	203.3100		-597.28	-505.49		87.5	75.30
Mg(ClO ₄) ₂		c	223.2132		-135.97	-112.8		54.0	
	std. state, m = 1	aq			-173.40				
	in 18 H ₂ O	aq			-171.549				
	20 H ₂ O	aq			-171.847				
	25 H ₂ O	aq			-172.230				
	30 H ₂ O	aq			-172.426				
	40 H ₂ O	aq			-172.633				
	50 H ₂ O	aq			-172.720				
	75 H ₂ O	aq			-172.815				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

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Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Magnesium

Table 93(5)

Substance	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
			0 K	298.15 K (25°C)				cal/deg mol
Mg(ClO ₄) ₂ in	100 H ₂ O	aq		-172.858				
	200 H ₂ O	aq		-172.924				
	300 H ₂ O	aq		-172.956				
	400 H ₂ O	aq		-172.982				
	500 H ₂ O	aq		-173.003				
	600 H ₂ O	aq		-173.023				
	800 H ₂ O	aq		-173.052				
	1,000 H ₂ O	aq		-173.073				
	2,000 H ₂ O	aq		-173.136				
	5,000 H ₂ O	aq		-173.208				
	10,000 H ₂ O	aq		-173.254				
	20,000 H ₂ O	aq		-173.290				
	50,000 H ₂ O	aq		-173.324				
	100,000 H ₂ O	aq		-173.346				
	∞ H ₂ O	aq		-173.40				

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Mg(ClO₄)₂
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Table 93(6)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Magnesium						
Substance	Formula and Description	State	Formula Weight	ΔHf°	ΔGf°	H ₂₉₈ ° - H ₀ °	S°	C _p °
				0 K	298.15 K (25°C)			
				kcal/mol				
Mg(ClO ₄) ₂	in 2,000 CH ₃ OH							
	in 4,000 C ₂ H ₅ OH							
	∞ C ₂ H ₅ OH							
	in 4,000 n-C ₃ H ₇ OH							
	∞ n-C ₃ H ₇ OH							
	in 4,000 i-C ₃ H ₇ OH							
	∞ i-C ₃ H ₇ OH							
	in 4,000 n-C ₄ H ₉ OH							
	∞ n-C ₄ H ₉ OH							
Mg(ClO ₄) ₂ ·2H ₂ O		c	259.2439	-181.87				
Mg(ClO ₄) ₂ ·4H ₂ O		c	295.2746	-181.2				
Mg(ClO ₄) ₂ ·6H ₂ O		c	331.3052	-181.5	-445.3		124.5	
MgO·MgCl ₂		c	135.5294	-175.5				
MgO·MgCl ₂ ·6H ₂ O		c	243.6214	-176.6				
MgO·MgCl ₂ ·16H ₂ O		c	423.7748	-179.5				
MgOHCl		c	76.7724	-181.1	-174.9		20.0	

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Table 93(7)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Magnesium							
Substance		Formula Weight	State	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°	
Formula and Description									298.15 K (25°C)
		kcal/mol						cal/deg mol	
3Mg(OH) ₂ •MgCl ₂		270.1982	c		-827.9				
3Mg(OH) ₂ •MgCl ₂ •4H ₂ O		342.2596	c		-1123.7				
3Mg(OH) ₂ •MgCl ₂ •7H ₂ O		396.3056	c		-1337.6				
3Mg(OH) ₂ •MgCl ₂ •8H ₂ O		414.3209	c		-1408.8				
5Mg(OH) ₂ •MgCl ₂ •8H ₂ O		530.9744	c		-1847.7				
MgBr ₂		184.130	c		-125.3	-120.4	28.0		
std. state, m = 1			g		-74.0	-158.4	6.4		
in			aq						
100 H ₂ O			aq						
200 H ₂ O			aq						
300 H ₂ O			aq						
400 H ₂ O			aq						
500 H ₂ O			aq						
600 H ₂ O			aq						
800 H ₂ O			aq						
1,000 H ₂ O			aq						
2,000 H ₂ O			aq						
5,000 H ₂ O			aq						

MgBr₂
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MgBr₂
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Table 93(8)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Magnesium								
Substance		Formula	State	Formula Weight	ΔHf°	ΔHf°	ΔGf°	H ₂₉₈ ° - H ₀ °	S°	C _p °
Formula and Description					0 K	298.15 K (25°C)				
						kcal/mol				
						cal/deg mol				
MgBr ₂	in 10,000 H ₂ O		aq			-169.528				
	20,000 H ₂ O		aq			-169.568				
	50,000 H ₂ O		aq			-169.606				
	100,000 H ₂ O		aq			-169.626				
	∞ H ₂ O		aq			-169.68				
	in 81 HBr + 9,000 H ₂ O					-168.4				
MgBr ₂ ·6H ₂ O			c	292.2220		-576.0		-491.4	95.	
MgI ₂			c	278.1208		-87.0		-85.6	31.0	
	std. state, m = 1.		g			-41.				
	in 400 H ₂ O		aq			-137.96		-133.4	20.2	
	800 H ₂ O		aq			-137.49				
	1,000 H ₂ O		aq			-137.57				
	2,000 H ₂ O		aq			-137.59				
	5,000 H ₂ O		aq			-137.67				
	∞ H ₂ O		aq			-137.76				
			aq			-137.96				

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National Bureau of Standards

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Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Magnesium

Table	93(9)	Substance	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
Formula and Description				kcal/mol		cal/deg mol			
		0 K							
MgI ₂		in 81 HI + 9,000 H ₂ O	aq						
MgIO ₃		un-ionized; std. state, m = 1	aq	199.2146	-140.3	1.992			
MgS			c	56.376	-82.7			12.03	10.89
MgSO ₃			c	104.3742	-241.0				
MgSO ₃ ·3H ₂ O			c	158.4202	-461.7				
MgSO ₃ ·6H ₂ O			c	212.4662	-673.4				
MgSO ₄			c	120.3736	-307.1			21.9	23.06
		undissociated; std. state, m = 1	aq		-324.1			-1.7	
		std. state, m = 1	aq		-328.90			-28.2	
		in 20 H ₂ O	aq		-327.12				
		25 H ₂ O	aq		-327.20				
		50 H ₂ O	aq		-327.41				
		75 H ₂ O	aq		-327.52				
		100 H ₂ O	aq		-327.59				
		200 H ₂ O	aq		-327.74				
		300 H ₂ O	aq		-327.82				
		400 H ₂ O	aq		-327.88				

MgSO₄
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MgSO₄
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Table 93(10)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Magnesium							
MgSO ₄	Substance Formula and Description	State	Formula Weight	ΔHf°	ΔHf°	ΔGf°	H ₂₉₈ ° - H ₀ °	S°	C _p °
				0 K	kcal/mol				
MgSO ₄	in 500 H ₂ O	aq			-327.93				
	600 H ₂ O	aq			-327.97				
	800 H ₂ O	aq			-328.04				
	1,000 H ₂ O	aq			-328.08				
	2,000 H ₂ O	aq			-328.202				
	5,000 H ₂ O	aq			-328.358				
	10,000 H ₂ O	aq			-328.467				
	20,000 H ₂ O	aq			-328.581				
	50,000 H ₂ O	aq			-328.707				
	100,000 H ₂ O	aq			-328.774				
	∞ H ₂ O	aq			-328.90				
MgSO ₄	in CH ₃ OH, std. state, m = 1				-310.9	-277.4		+1.	
MgSO ₄	in C ₂ H ₅ OH, std. state, m = 1				-309.3	-276.1		+2.	
MgSO ₄ · H ₂ O		c	138.3889		-382.9	-341.5		30.2	
MgSO ₄ · 2H ₂ O		amorp			-376.4	-335.8		33.0	
MgSO ₄ · 4H ₂ O		c	156.4043		-453.2				
MgSO ₄ · 6H ₂ O		c	192.4350		-596.7				
MgSO ₄ · 7H ₂ O		c	228.4656		-737.8	-629.1		83.2	83.20
MgSO ₄ · 7H ₂ O		c	246.4810		-809.92	-686.4		89.	

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Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Magnesium

Table	93 (11)	Substance	Formula and Description	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
			$Mg_2O_3 \cdot 3H_2O$	c	190.4842		-465.6				
			$Mg_2O_3 \cdot 6H_2O$	c	244.5302		-680.8				
			$3Mg(OH)_2 \cdot MgSO_4 \cdot 8H_2O$	c	439.4765		-1543.3				
			$MgSeO_3$	c	151.270		-215.15				
				amorp			-213.35				
			$MgSeO_3 \cdot 6H_2O$	c	259.362		-647.00				
			$MgSeO_4$	c	167.270		-231.48	-214.2		-20.1	
			std. state, m = 1 in 1,500 H_2O	aq			-254.8				
				aq			-254.31				
			$MgSeO_4 \cdot H_2O$	c	185.285		-309.62				
			$MgSeO_4 \cdot 4H_2O$	c	239.331		-523.40				
			$MgSeO_4 \cdot 6H_2O$	c	275.362		-664.2				
			$MgTe$	c	151.912		-50.				
			Mg_3N_2	c	100.9494		-110.1				

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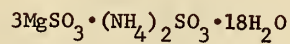
Mg(NO₃)₂
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Table 93(12)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Magnesium								
Substance		Formula	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°	
Formula and Description					0 K	298.15 K (25°C)				
					kcal/mol					cal/deg mol
Mg(NO ₃) ₂	std. state, m = 1		c	148.3218		-188.97	-140.9	39.2	33.92	
	in		aq			-210.70	-161.9	37.0		
	12 H ₂ O		aq			-209.62				
	15 H ₂ O		aq			-209.89				
	20 H ₂ O		aq			-210.07				
	25 H ₂ O		aq			-210.11				
	50 H ₂ O		aq			-210.14				
	75 H ₂ O		aq			-210.15				
	100 H ₂ O		aq			-210.16				
	200 H ₂ O		aq			-210.21				
	300 H ₂ O		aq			-210.25				
	400 H ₂ O		aq			-210.28				
	500 H ₂ O		aq			-210.30				
	600 H ₂ O		aq			-210.317				
	800 H ₂ O		aq			-210.344				
	1,000 H ₂ O		aq			-210.364				
	2,000 H ₂ O		aq			-210.423				
	5,000 H ₂ O		aq			-210.502				
	10,000 H ₂ O		aq			-210.551				
	20,000 H ₂ O		aq			-210.590				

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Table 93 (13)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Magnesium						
Substance		Formula Weight	State	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p
Formula and Description								
				0 K	kcal/mol			cal/deg mol
Mg(NO ₃) ₂	in 50,000 H ₂ O		aq					
	100,000 H ₂ O		aq					
	∞ H ₂ O		aq					
Mg(NO ₃) ₂ · 2H ₂ O		184.3525	c					
Mg(NO ₃) ₂ · 6H ₂ O		256.4138	c					
Mg(NH ₃) ₂ ⁺		58.3732	aq					
MgCl ₂ · NH ₃		112.2486	c					
MgCl ₂ · 2NH ₃		129.2792	c					
Mg(NH ₃) ₂ Cl ₂			aq					
Mg(ClO ₄) ₂ · 6NH ₃		325.3969	c					
MgBr ₂ · NH ₃		201.1606	c					
MgBr ₂ · 2NH ₃		218.1912	c					
MgI ₂ · 2NH ₃		312.1820	c					
Mg(NH ₃) ₂ SO ₄		154.4348	aq					
3MgSO ₃ · (NH ₄) ₂ SO ₃ · 6H ₂ O		537.3540	c					
3MgSO ₃ · (NH ₄) ₂ SO ₃ · 18H ₂ O		753.5381	c					
							108.	

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MgP₂O₇²⁻
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Table 93 (14)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Magnesium						
Substance		Formula Weight	State	ΔHf°	ΔGf°	H ₂₉₈ ° - H ₀ °	S°	C _p °
Formula and Description								
		kcal/mol						
		cal/deg mol						
MgP ₂ O ₇ ²⁻	std. state, m = 1	198.2554	aq	0 K	-577.2	6.467	-19.	
Mg ₂ P ₂ O ₇	α	222.5674	c		-651.5		37.02	42.53
Mg ₃ (PO ₄) ₂		262.8788	c	-896.98	-903.6	7.825	45.22	51.02
MgHPO ₄		120.2914	aq		-419.8			
MgNH ₄ PO ₄ ·6H ₂ O		245.4140	c		-880.0			
MgAs ₄		323.9984	c		-30.			
Mg ₃ As ₂		222.7792	c		-88.8			
Mg ₃ (AsO ₄) ₂		350.7744	c		-739.2			
MgHAsO ₄		164.2392	aq		-325.9			
Mg(H ₂ AsO ₄) ₂		306.1823	aq		-545.5			
MgNH ₄ AsO ₄ ·6H ₂ O		289.3618	c		-792.7			
Mg ₃ Sb ₂	α	316.436	c		-56.			
Mg ₃ Bi ₂	β	490.896	c		-36.8			
MgC ₂		48.3343	c		+20.			
Mg ₂ C ₃		84.6574	c		+17.			

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MgCl₂·6C₂H₅OH
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Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Table 93(15) Magnesium

Substance	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
MgCO ₃	c	84.3214		-261.9	-241.9		15.7	18.05
MgCO ₃ ·3H ₂ O	c	138.3674		-412.6	-412.6			
MgCO ₃ ·5H ₂ O	c	174.3980		-525.7	-525.7			
MgC ₂ O ₄	c	112.3319		-303.3				
std. state, m = 1	aq			-308.8	-269.8		-22.1	
MgC ₂ O ₄ ·2H ₂ O	c	148.3626			-390.4			
Mg(C ₂ O ₄) ₂ ²⁻	aq	200.3518			-436.9			
MgHCO ₃ ⁺	aq				-250.3			
un-ionized; std. state, m = 1	aq	85.3293			-198.7			
Mg(CO ₂ H) ₂	c	114.348		-418.4				
magnesium formate	c	83.3570						
Mg(C ₂ H ₃ O ₂) ₂ ⁺	aq	174.4008		-422.8				
std. state, m = 1	aq							
Mg(C ₂ H ₃ O ₂) ₂	c	174.4008		-561.2				
magnesium glycolate	aq							
Mg(C ₂ H ₃ O ₂) ₂ ·2H ₂ O	c	210.4315						
3MgCO ₃ ·Mg(OH) ₂ ·3H ₂ O	c	365.3368						
MgCl ₂ ·6CH ₃ OH	c	287.4726		-508.				
MgCl ₂ ·6C ₂ H ₅ OH	c	371.6351		-609.	-1100.3			

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MgCN₂
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Table 93(16)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Magnesium								
Substance		Formula Weight	State	ΔHf° 0 K	ΔHf°	ΔGf°	H° ₂₉₈ - H° ₀	S°	C _p °	
Formula and Description										298.15 K (25°C)
		kcal/mol							cal/deg mol	
MgCN ₂	magnesium cyanamide	64.3366	c		-60.3					
Mg(CN) ₂		76.3477	aq		-39.2					
Mg(NH ₂ CH ₂ COO) ⁺ m = 1	un-ionized; std. state,	98.3717	aq			-188.7				
Mg(NO ₃) ₂ ·6CH ₃ OH		340.5764	c		-559.					
Mg(NO ₃) ₂ ·6C ₂ H ₅ OH		424.7389	c		-603.					
Mg ₂ Si		76.710	c		-18.6	-18.0		18.	17.6	
MgSiO ₃	clinoenstatite	100.3962	c	-368.039	-370.22	-349.46	2.895	16.19	19.45	
Mg ₂ SiO ₄	forsterite	140.7076	c	-516.42	-519.6	-491.2	4.129	22.74	28.32	
Mg ₃ Si ₂ O ₅ (OH) ₄	chrysotile	277.1345	c		-1043.4	-965.1		52.9	65.41	
	antigorite		c					53.2	65.47	
Mg ₃ Si ₄ O ₁₀ (OH) ₂	talc	379.2887	c	-1405.57	-1415.5	-1324.8	11.20	62.3	76.9	
Mg ₂ Ce		121.214	c	-25.73	-26.0	-25.3	3.229	20.67	16.62	
Mg ₂ Sn		167.314	c		-19.3					
Mg ₂ Pb		255.814	c		-12.3					
PbI ₂ ·2MgI ₂		1017.240	c		-206.5					

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Table 93(17) Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity
Magnesium

Substance	State	Formula Weight	ΔH_f° 0 K	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	s°	C_p°		
									298.15 K (25°C)	
									kcal/mol	
MgB ₂	c	45.934				1.625	8.59	11.59		
MgB ₄	c	67.556				2.270	12.51	16.78		
MgB ₁₂	c	154.044	-52.8	-53.	-52.	4.49	21.4	36.2		
MgO·Al ₂ O ₃	c	142.2726	-549.05	-552.9	-523.1	3.683	19.27	27.77		
Mg ₂ Al ₄ Si ₄ O ₁₈ cordierite	c	584.9692		-2177.	-2055.		97.3	108.1		
MgTi	c	228.682		-12.						
MgZn	c	89.682		-4.0						
MgZn ₂	c	155.052		-7.8				17.7		
Mg ₂ Zn ₁₁	c	767.874		-30.						
MgCu ₂	c	151.392		-8.02	-8.1		24.	17.43		
Mg ₂ Cu	c	112.164		-6.84						
MgCd	c	136.712	-3.814	-3.84	-3.71	2.66	19.76	12.32		
MgCd ₃	c	361.512	-4.662	-4.52	-4.71	5.81	45.56	28.24		
Mg ₃ Cd	c	185.336	-5.244	-5.32	-5.14	5.00	35.20	24.04		
MgHgBr ₄	aq	544.538		-210.4						
Mg(HgBr ₃) ₂	aq	904.946		-249.1						

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National Bureau of Standards

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2MgBr₂•HgBr₂
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Table 93(18)		Magnesium						
Substance		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Formula and Description	State	Formula Weight	ΔHf°	ΔGf°	H° ₂₉₈ - H°	S°	C _p	
			0 K	298.15 K (25°C)				
		kcal/mol						cal/deg mol
2MgBr ₂ •HgBr ₂	aq	728.668		-381.2				
4MgBr ₂ •HgBr ₂	aq	1096.928		-721.2				
MgHg(CN) ₄	aq	328.973		19.4				
Mg[Hg(CN) ₃] ₂	aq	581.599		81.6				
2Hg(CN) ₂ •MgCl ₂	aq	600.469		-57.1				
2Hg(CN) ₂ •MgCl ₂ •6H ₂ O	c	708.561		-476.5				
2Hg(CN) ₂ •MgBr ₂	aq	689.381		-36.0				
2Hg(CN) ₂ •MgBr ₂ •8H ₂ O	c	833.504		-597.9				
2Hg(CN) ₂ •MgI ₂	aq	783.372		-8.4				
2Hg(CN) ₂ •MgI ₂ •8H ₂ O	c	927.495		-574.2			21.20	17.54
MgNi ₂	c	141.732		-13.2	-12.9			
Mg ₂ Ni	c	107.334		-9.5				
MgFe ₂ O ₄	c	200.004		-341.4	-314.8		29.6	34.35

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Table 93 (19)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Magnesium						
Substance	Formula and Description	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
				0 K	298.15 K (25°C)			
$MgFe(CN)_6$	un-ionized; std. state, m=1	aq	236.2661					
$MgFe(CN)_6$	un-ionized; std. state, m=1	aq						
$Mg_2Fe(CN)_6$	std. state, m = 1	aq	260.5781					
	in 1,500 H ₂ O	aq		-114.3	-286.8			
	2,000 H ₂ O	aq		-111.72	-280.0			
	5,000 H ₂ O	aq		-111.90				
	7,500 H ₂ O	aq		-112.34				
	10,000 H ₂ O	aq		-112.49				
	20,000 H ₂ O	aq		-112.59				
	50,000 H ₂ O	aq		-112.80				
	100,000 H ₂ O	aq		-113.10				
	200,000 H ₂ O	aq		-113.39				
	500,000 H ₂ O	aq		-113.65				
	1,000,000 H ₂ O	aq		-113.94				
	∞ H ₂ O	aq		-114.12				
		aq		-114.3				
$MgCrO_4$		c	140.3056	-321.1				
$MgCr_2O_4$		aq		-324.0				
		c	192.3016	-426.3	-398.9		25.34	30.30

MgCr₂O₄
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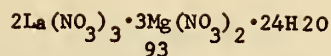
MgMoO₄
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Table 93 (20)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Magnesium						
Substance	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	s°	C_p°
MgMoO ₄	c	184.250		-334.81	-309.69		28.4	26.57
MgWO ₄	g	272.160	-224.	-366.3	-339.6	4.115	24.18	26.14
MgV ₂ O ₆	c	222.1924	-222.	-526.19	-487.43		38.4	39.47
Mg ₂ V ₂ O ₇	g	262.5038		-677.80	-632.24		47.9	48.63
MgTiO ₃	c	120.210		-375.9	-354.8	3.240	17.82	21.96
MgTi ₂ O ₅	c	200.109		-599.8	-565.7	5.428	30.42	35.15
Mg ₂ TiO ₄	c	160.522		-517.5	-489.2	4.502	26.13	30.75
MgY	c	113.217		-6.				
Mg ₅ Y ₂	c	299.370		-21.				
Mg ₁₇ Y ₃	c	680.019		-50.				
MgPr	c	165.219		-8.2				
Mg ₃ Pr	c	213.843		-11.1				
MgCe	c	164.432		-13.				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

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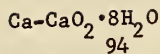
Table 93(21)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Magnesium					
Substance		ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
Formula and Description		0 K	298.15 K (25°C)				
		kcal/mol					
		cal/deg mol					
		Formula Weight					
		State					
Mg ₃ Ce		c	213.056	-17.			
MgLa		c	163.222	-5.7			
Mg ₃ La		c	211.846	-12.8			
2La(NO ₃) ₃ · 3Mg(NO ₃) ₂ · 24H ₂ O		c	1527.183			525.	483.2

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Table 94(1)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Substance		Calcium						
Formula and Description	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
			0 K	298.15 K (25°C)				
		kcal/mol		kcal/mol		cal/deg mol		
Ca	c	40.08	0	0	0	1.364	9.90	6.05
Ca ⁺	g		42.48	42.6				
Ca ²⁺	g		183.45	185.05	34.5	1.481	36.992	4.968
	g		457.21	460.29				
	aq			-129.74	-132.30		-12.7	
	g		1631.19	1635.75				
	g		3178.6	3184.6				
	g		5125.1	5132.7				
	g		7633.7	7642.7				
	g		10579.	10589.				
	g		13974.	13986.				
CaO	c	56.079		-151.79	-144.37		9.50	10.23
	g		+11.					
CaO ₂	c	72.079		-156.0				
CaO ₂ ·8H ₂ O	c	216.202		-718.2				



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Table 94(2)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Calcium						
Substance		Formula Weight	State	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
Formula and Description				0 K	298.15 K (25°C)			
		kcal/mol						
CaH		41.088	g	55.	47.9	2.076	48.19	7.11
CaH ₂		42.096	c	.	-35.2		10.	
CaOH		57.087	g		-48.			
CaOH ⁺	std. state, m = 1		aq		-171.7			
Ca(OH) ₂		74.095	c		-235.68		19.93	20.91
			g		-130.			
	std. state, m = 1		aq		-239.68		-17.8	
Ca(OH) ₂ · H ₂ O		108.109	c		-291.2			
CaF		59.078	g	-64.76	-71.2	2.181	54.8	8.03
CaF ₂		78.077	c		-291.5		16.46	16.02
			g	-186.35	-188.9	3.025	65.55	12.25
	std. state, m = 1		aq		-288.74		-19.3	
CaCl		75.533	g	-23.23	-29.7	2.292	57.70	8.58
CaCl ₂		110.986	c		-190.2		25.0	17.35
			g	-112.76	-114.54	3.613	69.35	14.18
	std. state, m = 1		aq		-209.64		14.3	

 CaCl₂
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CaCl₂
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Table 94(3)		Calcium							
		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity							
CaCl ₂	in	Substance Formula and Description	State	Formula Weight	ΔHf°	ΔGf°	H ₂₉₈ ^o - H ₀ ^o	S°	C _p ^o
					0 K	298.15 K (25°C)			
		10 H ₂ O	aq			-206.20			
		11 H ₂ O	aq			-206.55			
		12 H ₂ O	aq			-206.83			
		13 H ₂ O	aq			-207.06			
		15 H ₂ O	aq			-207.42			
		20 H ₂ O	aq			-207.95			
		25 H ₂ O	aq			-208.19			
		40 H ₂ O	aq			-208.51			
		50 H ₂ O	aq			-208.63			
		75 H ₂ O	aq			-208.769			
		100 H ₂ O	aq			-208.849			
		300 H ₂ O	aq			-209.079			
		400 H ₂ O	aq			-209.144			
		500 H ₂ O	aq			-209.162			
		800 H ₂ O	aq			-209.227			
		1,000 H ₂ O	aq			-209.259			
		2,000 H ₂ O	aq			-209.343			
		5,000 H ₂ O	aq			-209.433			
		10,000 H ₂ O	aq			-209.485			
		20,000 H ₂ O	aq			-209.526			
		50,000 H ₂ O	aq			-209.565			

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

Ca(ClO₄)₂
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Washington, D. C.

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Calcium

Table 94(4)

Substance		Formula Weight	State	ΔHf° 0°K	ΔHf°	ΔGf°	H ₂₉₈ ^o - H ₀ ^o	S°	C _p ^o
Formula and Description									
298.15°K (25°C)									
kcal/mol									
cal/deg mol									
CaCl ₂	in 100,000 H ₂ O		aq		-209.585				
	500,000 H ₂ O		aq		-209.615				
	∞ H ₂ O		aq		-209.64				
CaCl ₂ ·H ₂ O		129.001	c		-265.1				
CaCl ₂ ·2H ₂ O		147.017	c		-335.3				
CaCl ₂ ·4H ₂ O		183.047	c		-480.3				
CaCl ₂ ·6H ₂ O		219.078	c		-623.3				
CaOCl ₂		126.985	c		-178.4				
			aq		-188.9				
CaOCl ₂ ·H ₂ O		145.001	c		-249.1				
Ca(OCl) ₂		142.985	aq		-180.3				
Ca(ClO) ₂		174.984	c		-162.1				
Ca(ClO ₂) ₂		238.981	c		-176.09				
Ca(ClO ₄) ₂			aq		-191.56	-136.42		74.3	
	std. state, m = 1		aq		-187.59				
	in		aq		-188.22				
	7 H ₂ O		aq		-188.72				
	8 H ₂ O		aq		-189.12				
	9 H ₂ O		aq		-189.78				
	10 H ₂ O		aq						
	12 H ₂ O		aq						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

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Ca(ClO₄)₂
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Table 94(5)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity					
Substance		Calcium					
Formula and Description	State	Formula Weight	ΔHf°	ΔGf°	H ₂₉₈ ° - H ₀ °	S°	C _p °
			0 K	298.15 K (25°C)			
Ca(ClO ₄) ₂	in						
15 H ₂ O	aq		-190.35				
20 H ₂ O	aq		-190.75				
25 H ₂ O	aq		-190.92				
30 H ₂ O	aq		-190.99				
40 H ₂ O	aq		-191.08				
50 H ₂ O	aq		-191.14				
75 H ₂ O	aq		-191.21				
100 H ₂ O	aq		-191.20				
200 H ₂ O	aq		-191.13				
300 H ₂ O	aq		-191.17				
500 H ₂ O	aq		-191.18				
1,000 H ₂ O	aq		-191.25				
10,000 H ₂ O	aq		-191.42				
100,000 H ₂ O	aq		-191.51				
∞ H ₂ O	aq		-191.56				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Calcium

Table	94(6)	Substance	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
					kcal/mol				
Formula and Description					0 K	298.15 K (25°C)			
Ca(ClO ₄) ₂		in CH ₃ OH; std. state, m = 1	c	238.981	-204.3				
		in C ₂ H ₅ OH; std. state, m = 5/1			-197.9				
		in n C ₃ H ₇ OH; std. state, m = 3/1			-194.2				
		in i C ₃ H ₇ OH; std. state, m = 1/1			-196.3				
		in n C ₄ H ₉ OH; std. state, m = 4/1			-192.7				
Ca(ClO ₄) ₂ •4H ₂ O		c	311.043	-465.8	-352.97		103.6		
CaCl ₂ •2CaO		c	223.145	-505.					
CaCl ₂ •3CaO		c	279.224	-654.					
CaCl ₂ •3CaO•3H ₂ O		c	333.270	-910.					
CaCl ₂ •3CaO•16H ₂ O		c	567.470	-1832.					
CaClH		c	76.541	-120.5					
CaBr ₂			c	199.898	-163.2	-158.6		31.	
		std. state, m = 1	g		-95.2	-182.00		26.7	

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CaBr₂
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Table 94(7)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Substance		Calcium						
Formula and Description	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °	
			0 K	298.15 K (25°C)				cal/deg mol
CaBr ₂	in 400 H ₂ O	aq		-187.39				
	500 H ₂ O	aq		-187.42				
	800 H ₂ O	aq		-187.47				
	1,000 H ₂ O	aq		-187.494				
	2,000 H ₂ O	aq		-187.566				
	5,000 H ₂ O	aq		-187.644				
	10,000 H ₂ O	aq		-187.693				
	20,000 H ₂ O	aq		-187.729				
	50,000 H ₂ O	aq		-187.767				
	100,000 H ₂ O	aq		-187.786				
	500,000 H ₂ O	aq		-187.815				
	∞ H ₂ O	aq		-187.84				
	CaBr ₂ •6H ₂ O	c	307.990		-514.6		98.	
	Ca (BrO ₃) ₂	c	295.894					
CaBr ₂ •3CaO•16H ₂ O	aq							
	c	656.382						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

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Table 94(8)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Calcium						
Substance		Formula Weight	State	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
Formula and Description								
				kcal/mol		cal/deg mol		
CaBrH		120.997	c		-106.0			
CaI ₂		293.889	c		-127.5	-126.4	34.	
			g		-65.			
	std. state, m = 1		aq		-156.12	-156.96	40.5	
	in 400 H ₂ O		aq		-155.6			
CaI ₂ ·8H ₂ O		438.012	c		-700.2			
Ca(IO ₃) ₂		389.885	c		-239.6	-200.6	55.	
Ca(IO ₃) ₂ ·H ₂ O		407.901	c		-309.1			
Ca(IO ₃) ₂ ·6H ₂ O		497.977	c		-664.6	-542.0	108.	
CaI ₂ ·3CaO·16H ₂ O		750.372	c		-1779.			
Ca ₅ (IO ₆) ₂		646.202	c		-915.			
CaIH		167.992	c		-87.5			
CaS		72.144	c		-115.3	-114.1	13.5	11.33
			g	+32.				
CaSO ₃		120.142	c				24.23	12.92
CaSO ₃ ·2H ₂ O		156.173	c		-418.9	-371.7	44.	42.7

CaSO₃·2H₂O
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National Bureau of Standards

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 CaSO_4
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Table 94(9)		Calcium					
Substance		ΔHf°	ΔHf°	ΔGf°	$\text{H}_{298}^\circ - \text{H}_0^\circ$	S°	C_p°
Formula and Description		0 K	298.15 K (25°C)				
		kcal/mol		cal/deg mol			
		Formula Weight	State				
CaSO_4	insol., anhydrite	136.142	c	-342.76	-315.93	25.5	23.82
	sol., α		c	-340.64	-313.93	25.9	23.95
	sol., β		c	-339.58	-312.87	25.9	23.67
	std. state, $m = 1$		aq	-347.06	-310.27	-7.9	
	in 5,000 H_2O		aq	-346.414			
	7,500 H_2O		aq	-346.518			
	10,000 H_2O		aq	-346.566			
	20,000 H_2O		aq	-346.685			
	50,000 H_2O		aq	-346.827			
	100,000 H_2O		aq	-346.915			
	500,000 H_2O		aq	-347.009			
	1,000,000 H_2O		aq	-347.026			
	$\infty \text{H}_2\text{O}$		aq	-347.06			
$\text{CaSO}_4 \cdot 1/2\text{H}_2\text{O}$	macro; α	145.149	c	-376.85	-343.41	31.2	28.54
	micro; β		c	-376.35	-343.18	32.1	29.69
$\text{CaSO}_4 \cdot 2\text{H}_2\text{O}$	selenite	172.172	c	-483.42	-429.60	46.4	44.46

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Table 94 (10)

Calcium

Substance	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
Formula and Description			0 K	298.15 K (25°C)				
CaS ₂ O ₃ in 35 H ₂ O 100 H ₂ O 200 H ₂ O 400 H ₂ O 1,000 H ₂ O α H ₂ O	aq	152.206		-284.56				
	aq			-284.80				
	aq			-284.95				
	aq			-285.02				
	aq			-285.10				
CaS ₂ O ₆ ·4H ₂ O dithionate	aq			-285.6				
	c	272.266		-697.				
Ca(HS) ₂	aq	106.224		-138.1				
CaSe	c	119.04		-88.0	-86.8		16.	
	c	203.069		-384.9	-341.5			
CaSeO ₃ ·2H ₂ O precipitate	c	183.038		-265.25				
	aq			-273.30				
CaSeO ₄ in 1,600 H ₂ O	c	219.068		-407.9	-355.4		53.	
	c	167.68					19.	
CaTe	c							

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National Bureau of Standards

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Table 94 (11)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Calcium						
Substance		Formula Weight	State	ΔH_f° 0 K	ΔH_f°	$H_{298}^\circ - H_0^\circ$ 298.15 K (25°C)	S°	C _p °
Formula and Description								
Ca(N ₃) ₂	c	124.120			+3.5			
Ca(N ₃) ₂ •0.5H ₂ O	c	133.128			-33.3			
Ca(N ₃) ₂ •1.5H ₂ O	c	151.143			-104.5			
Ca(N ₃) ₂ •4H ₂ O	c	196.182			-278.2			
Ca ₃ N ₂	c	148.253			-103.			
CaN ₂ O ₂ •4H ₂ O	c	172.154			-407.5			
Ca(NO ₂) ₂	c	132.091			-177.2			
	aq				-179.5			
	in 800 H ₂ O				-450.7			
Ca(NO ₂) ₂ •4H ₂ O	c	204.152			-224.28			
Ca(NO ₃) ₂	c	164.090			-228.86		-177.63	46.2
	aq				-228.86		-185.52	57.3
	std. state, m = 1							
	in							
	3 H ₂ O				-228.164			
	4 H ₂ O				-228.793			
	5 H ₂ O				-229.247			
	6 H ₂ O				-229.545			
	7 H ₂ O				-229.758			
	8 H ₂ O				-229.905			
	9 H ₂ O				-229.988			
	10 H ₂ O				-230.037			

Ca(N₃)₂
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SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Table 94 (12)

Calcium

Substance		Formula Weight	State	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
Formula and Description	0 K							
Ca(NO ₃) ₂	11 H ₂ O		aq					
	12 H ₂ O		aq					
	13 H ₂ O		aq					
	15 H ₂ O		aq					
	20 H ₂ O		aq					
	25 H ₂ O		aq					
	40 H ₂ O		aq					
	50 H ₂ O		aq					
	75 H ₂ O		aq					
	100 H ₂ O		aq					
	300 H ₂ O		aq					
	400 H ₂ O		aq					
	500 H ₂ O		aq					
	800 H ₂ O		aq					
	1,000 H ₂ O		aq					
	2,000 H ₂ O		aq					
5,000 H ₂ O		aq						
10,000 H ₂ O		aq						
20,000 H ₂ O		aq						
50,000 H ₂ O		aq						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

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Table 94(13)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Substance		Calcium						
Formula and Description	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
			0 K	298.15 K (25°C)				
Ca(NO ₃) ₂ in 100,000 H ₂ O 500,000 H ₂ O ∞ H ₂ O	aq			-228.809				
	aq			-228.836				
	aq			-228.86				
in 100 C ₂ H ₅ OH in ∞ C ₂ H ₅ OH	c	200.120		-368.25	-293.82		64.4	
	c	218.136		-439.3	-351.8		76.3	
Ca(NO ₃) ₂ •4H ₂ O	c	236.151		-509.64	-409.53		89.7	
Ca(NH ₂) ₂	c	72.125		-91.9				
Ca(N ₃) ₂ •2N ₂ H ₄	c	156.165		+12.6				
Ca(N ₃) ₂ •2N ₂ H ₄	c	188.211		+4.0				
Ca(NO ₃) ₂ •Ca(OH) ₂	c	238.185		-461.5				
Ca(NO ₃) ₂ •Ca(OH) ₂ •2½H ₂ O	c	283.223		-640.9				
CaCl ₂ •NH ₃	c	128.017		-217.5				
CaCl ₂ •2NH ₃	c	145.047		-242.8				
CaCl ₂ •4NH ₃	c	179.108		-285.2				
CaCl ₂ •8NH ₃	c	247.231		-368.9				
Ca(ClO ₄) ₂ •6NH ₃		341.165		-315.7				

 Ca(NO₃)₂
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SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

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Table 94(14)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Calcium						
Substance	Formula and Description	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
				0 K	298.15 K (25°C)			
				kcal/mol				
	CaBr ₂ ·NH ₃	c	216.929	-191.5				
	CaBr ₂ ·2NH ₃	c	233.959	-219.0				
	CaBr ₂ ·6NH ₃	c	302.082	-309.1				
	CaBr ₂ ·8NH ₃	c	336.143	-350.7				
	CaI ₂ ·NH ₃	c	310.919	-157.4				
	CaI ₂ ·2NH ₃	c	327.950	-186.6				
	CaI ₂ ·6NH ₃	c	396.072	-284.9				
	CaI ₂ ·8NH ₃	c	430.134	-324.3				
	Ca ₃ P ₂	c	182.188	-121.				
	Ca(PO ₃) ₂	c	198.024		5.715		35.05	34.68
	β							
	glassy	amorp		-587.0	-748.6	7.430	45.23	44.89
	β	c	254.103	-792.88				
	Ca ₂ P ₂ O ₇	c	310.183					
	β, low temp. form	c		-984.9	-928.5		56.4	54.45
	Ca ₃ (PO ₄) ₂	c		-982.3	-926.3		57.58	55.35
	α, high temp. form	c						
	std. state, m = 1	aq		-999.8	-883.9		-144.	

Ca₃(PO₄)₂
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SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

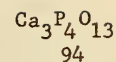


Table 94 (15)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Calcium						
Substance		Formula Weight	State	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
Formula and Description				0 K	298.15 K (25°C)			
$\text{Ca}_3\text{P}_4\text{O}_{13}$	glassy	452.127	amorp	-1374.				
CaHPO_4	std. state, m = 1	136.059	c	-430.299	-401.83	4.455	26.62	26.30
$\text{CaHPO}_4 \cdot 2\text{H}_2\text{O}$		172.090	c	-559.496	-392.64		-20.7	
$\text{Ca}(\text{H}_2\text{PO}_4)_2$	std. state, m = 1	170.057	c	-418.9	-515.00	7.490	45.28	47.10
	in 50 H_2O		aq	-423.1				
	100 H_2O		aq	-419.8				
	200 H_2O		aq	-420.6				
	300 H_2O		aq	-421.4				
	400 H_2O		aq	-421.8				
	600 H_2O		aq	-422.1				
	800 H_2O		aq	-422.4				
	1,000 H_2O		aq	-422.6				
$\text{Ca}(\text{H}_2\text{PO}_4)_2$	std. state, m = 1	234.055	c	-742.04	-672.64		30.5	
$\text{Ca}(\text{H}_2\text{PO}_4)_2 \cdot \text{H}_2\text{O}$			aq	-749.38				
$\text{Ca}_8\text{H}_2(\text{PO}_4)_6 \cdot 5\text{H}_2\text{O}$		252.070	c	-797.013	-730.98	9.950	62.1	61.86
		982.561	c		-2931.			

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Table 94 (16)		Calcium						
Substance		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Formula and Description	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
			0 K	298.15 K (25°C)				
		kcal/mol		cal/deg mol				
$Ca_{10}(PO_4)_6(OH)_2$ hydroxyapatite std. state, m = 1	c	1004.643	-3201.4	-3221.	-3030.	30.71	186.6	184.0
$Ca_{10}(PO_4)_6F_2$ fluorapatite	aq			-3239.1	-2859.2		-450.	
$Ca(H_2PO_4)_2 \cdot H_2O \cdot NH_3$	c	1008.625	-3267.0	-3285.	-3103.	30.34	185.4	179.7
$Ca(H_2PO_4)_2 \cdot H_2O \cdot NH_3$	c	269.101		-37.5				
$Ca(H_2PO_4)_2 \cdot H_2O \cdot 2NH_3$	c	286.131		-857.6				
$Ca(H_2PO_4)_2 \cdot H_2O \cdot 4NH_3$	c	320.192		-894.8				
$Ca_3(AsO_4)_2$ hydrated precipitate	c	398.078		-788.4	-732.1		54.	
$CaHAsO_4$	aq	180.007		-799.				
$Ca(H_2AsO_4)_2$	aq	321.950		-345.6				
Ca_3Sb_2	c	363.74		-563.6				
Ca_3B_2	c	538.200		-174.				
CaC_2	c	64.102	-15.14	-107.		2.711	16.72	14.99
$CaCO_3$ calcite	c			-14.3	-15.5			
$CaCO_3$ aragonite	c	100.089		-288.46	-269.80		22.2	19.57
$CaCO_3$ std. state, m = 1	aq			-288.51	-269.55		21.2	19.42
				-291.58	-258.47		-26.3	

 $CaCO_3$
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CaC₂O₄
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Table 94 (17)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Substance		Calcium						
Formula and Description	State	Formula Weight	ΔHf°	ΔGf°	H ₂₉₈ ° - H ₀ °	S°	C _p °	
			0 K	298.15 K (25°C)				cal/deg mol
CaC ₂ O ₄	c	128.100	-325.2	-293.37		-1.8		
std. state, m = 1	aq		-326.9					
CaC ₂ O ₄ · H ₂ O	c	146.115	-400.30	-361.85		37.4	36.52	
Ca(HCOO) ₂	c	130.116	-331.4					
in 400 H ₂ O	aq		-332.5					
Ca(C ₂ H ₃ O ₂) ⁺	aq	99.125	-353.6	-222.1				
Ca(C ₂ H ₃ O ₂) ₂	c	158.170	-362.06	-308.88		28.7		
std. state, m = 1	aq		-360.1					
in 100 H ₂ O	aq		-360.7					
200 H ₂ O	aq		-360.9					
300 H ₂ O	aq		-361.1					
400 H ₂ O	aq		-361.4					
1,000 H ₂ O	aq		-361.7					
3,000 H ₂ O	aq		-361.8					
5,000 H ₂ O	aq		-361.9					
10,000 H ₂ O	aq		-362.0					
50,000 H ₂ O	aq		-362.06					
∞ H ₂ O	aq		-423.6					
Ca(C ₂ H ₃ O ₂) ₂ · H ₂ O	c	176.185						

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National Bureau of Standards

Washington, D. C.

Table 94 (18)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Substance		Calcium						
Formula and Description	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	s°	c _p °	
								298.15 K (25°C)
		0 K	kcal/mol					cal/deg mol
Ca(C ₂ H ₃ O ₃) ₂	c	190.169	-443.6					
Ca(C ₂ H ₃ O ₃) ₂ •3H ₂ O	aq	244.215	-442.0					
Ca(C ₂ H ₃ O ₃) ₂ •5H ₂ O	c	280.246	-654.0					
Ca(C ₂ H ₅ O) ₂	c	130.203	-791.4					
Ca(C ₂ H ₅ O) ₂ •2C ₂ H ₅ OH	c	222.342	-227.0					
3CaO•4C ₂ H ₅ OH	c	352.516	-368.2					
CaCl ₂ •3C ₂ H ₅ OH	c	249.195	-775.9					
CaCl ₂ •4C ₂ H ₅ OH	c	295.264	-399.9					
CaCN ₂	c	80.105	-467.5					
Ca(CN) ₂	c	92.116	-83.8					
3CaO•Ca(CN) ₂ •15H ₂ O	aq	530.584	-44.1					
Ca(NO ₃) ₂ •2CH ₃ OH	c	228.175	-56.9					
			-1606.					
			-343.0					

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CaSi
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Table 94 (19)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Substance		Calcium					S°	C _p °
Formula and Description	State	Formula Weight	ΔHf°	ΔHf°	ΔGf°	H ₂₉₈ ° - H ₀ °		
			0 K	298.15 K (25°C)				
		kcal/mol						
		cal/deg mol						
CaSi	c	68.166		-36.				
CaSi ₂	c	96.252		-36.				
Ca ₂ Si	c	108.246		-50.				
CaO·SiO ₂	c	116.164		-390.76	-370.39		19.58	20.38
wollastonite	c			-389.2	-369.2		20.88	20.67
pseudowollastonite	c			-382.65				
glassy	amorp							
β	c	172.244	-548.95	-551.5	-524.1	5.098	30.53	30.78
γ	c	228.323	-551.25	-554.0	-526.1	4.898	28.87	30.27
3CaO·SiO ₂	c	288.408	-942.23	-700.1	-665.4		40.3	41.08
rankinite	c			-946.7	-899.0	8.424	50.38	51.24
3CaO·2SiO ₂ ·CaF ₂	c	366.485		-1251.				
cuspidine	c							
CaO·GeO ₂	c	160.668		-308.7				
2CaO·GeO ₂	c	216.748		-476.8				
3CaO·GeO ₂	c	272.827		-623.5				
CaSn	c	158.770		-38.				
Ca ₂ Sn	c	198.850		-75.				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

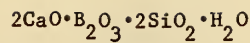
Washington, D. C.

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Calcium

Table 94(20)

Substance	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°		
								298.15 K (25°C)	
								cal/deg mol	
CaPb	c	247.27	-29.						
CaPb ₃	c	661.65	-30.						
Ca ₂ Pb	c	287.35	-51.						
CaI ₂ ·2PbI ₂	c	1215.886	-222.9						
CaI ₂ ·2PbI ₂ ·7H ₂ O	c	1341.994	-712.4						
CaO·B ₂ O ₃	c	125.700	-485.41	-459.87		25.06	24.85		
CaO·2B ₂ O ₃	c	195.320	-803.12	-756.96		32.2	37.75		
glassy	amorp		-790.48						
2CaO·B ₂ O ₃	c	181.779	-653.54	-620.62		34.68	35.16		
2CaO·3B ₂ O ₃	c	321.019	-1290.						
2CaO·3B ₂ O ₃ ·13H ₂ O Inyoite	c	555.219	-2221.						
3CaO·B ₂ O ₃	c	237.858	-819.57	-779.14		43.9	44.90		
Ca(BF ₄) ₂	c	213.689	-930.						
2CaO·B ₂ O ₃ ·2SiO ₂	c	301.949	-1100.						
2CaO·B ₂ O ₃ ·2SiO ₂ ·H ₂ O	c	319.964	-1170.						



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Table 94 (21)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Calcium						
Substance		ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°		
Formula and Description		0 K	298.15 K (25°C)					
		kcal/mol						
		cal/deg mol						
Formula and Description	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°	
CaAl ₂	c	94.043	-52.					
CaAl ₄	c	148.006	-51.					
CaO·Al ₂ O ₃	c	158.041	-556.0	-527.9	4.569	27.30	28.87	
CaO·2Al ₂ O ₃	amorp		-550.					
CaO·Al ₂ O ₃	c	260.002	-944.98	-901.2	7.286	42.50	48.00	
2CaO·Al ₂ O ₃	c	214.120	-707					
	amorp		-701.					
2CaO·Al ₂ O ₃ ·5H ₂ O	c	304.197	-1081.					
3CaO·Al ₂ O ₃	c	270.199	-853.21	-815.4	8.216	49.2	50.16	
	amorp		-849.					
3CaO·Al ₂ O ₃ ·6H ₂ O	c	378.291	-1326.					
4CaO·Al ₂ O ₃	c	326.279	-1008.					
12CaO·7Al ₂ O ₃	c	1386.681	-4620.1	-4414.	41.98	250.2	259.5	
	amorp		-4562.					
3CaCl ₂ ·4AlCl ₃	c	866.320	-1265.					

CaAl₂
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Table 94 (22)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Calcium						
Substance		Formula Weight	State	ΔH_f° 0 K	ΔG_f°	$F_{298}^\circ - H_0^\circ$	S°	C_p°
Formula and Description	kcal/mol							
		298.15 K (25°C)						
		cal/deg mol						
$3CaO \cdot Al_2O_3 \cdot CaSO_4 \cdot 12H_2O$	c	622.525		-2098.				
$3CaO \cdot Al_2O_3 \cdot 3CaSO_4 \cdot 31H_2O$	c	1237.100		-4121.				
$3CaO \cdot Al_2O_3 \cdot 3CaSO_4 \cdot 32H_2O$	c	1255.115		-4192.				
$3CaO \cdot Al_2O_3 \cdot CaCO_3 \cdot 10.68H_2O$	c	562.693		-1954.				
$3CaO \cdot Al_2O_3 \cdot 3CaCO_3 \cdot 30H_2O$	c	1110.928		-3876.				
$CaO \cdot Al_2O_3 \cdot 2SiO_2$ anorthite, triclinic	c	278.210		-1009.2	-955.5		48.4	50.46
anorthite, hexagonal	c			-1004.3	-949.8		45.8	49.76
glassy	amorp			-991.8				
$CaO \cdot Al_2O_3 \cdot 2SiO_2 \cdot 2H_2O$ lawsonite	c	314.241		-1161.2	-1076.8		56.8	68.16
$CaO \cdot Al_2O_3 \cdot 6SiO_2$ heulandite	c	518.549		-1907.				
$2CaO \cdot Al_2O_3 \cdot SiO_2$ gehlenite	c	274.205		-952.6	-904.3		47.4	49.10
glassy	amorp			-940.7				
$2CaO \cdot 2Al_2O_3 \cdot 8SiO_2 \cdot 7H_2O$ leonhardite	c	922.867		-3396.6	-3146.0		220.4	228.0
$3CaO \cdot Al_2O_3 \cdot 2SiO_2$	c	390.369		-1329.				
$CaAlGaSi_2O_8$	c	320.949		-935.6				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

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Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Calcium

CaTl
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Table 94 (23)

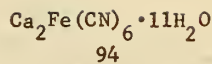
Substance	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
Formula and Description			0 K				
CaTl	c	244.45	-39.				
CaHgBr ₄	aq	560.306	-228.1				
CaHg ₂ Br ₆	aq	920.714	-266.3				
CaHg(CN) ₄ in 700 H ₂ O	aq	344.741	-2.0				
CaHg ₂ (CN) ₆ in 1,000 H ₂ O	aq	597.367	+65.8				
CaCl ₂ •2Hg(CN) ₂ in 1,000 H ₂ O	aq	616.237	-74.5				
CaCl ₂ •2Hg(CN) ₂ •6H ₂ O	c	724.329	-498.0				
CaBr ₂ •2Hg(CN) ₂ in 1,000 H ₂ O	aq	705.149	-53.6				
CaBr ₂ •2Hg(CN) ₂ •7H ₂ O	c	831.257	-550.7				
CaI ₂ •2Hg(CN) ₂ in 1,000 H ₂ O	aq	799.140	-26.1				
CaI ₂ •2Hg(CN) ₂ •7H ₂ O	c	925.248	-525.8				
CaAg(CN) ₃ in 1,200 H ₂ O	aq	226.004	-29.3				
CaAg ₂ (CN) ₄ in 550 H ₂ O	aq	359.891	-0.9				

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Table 94 (24)

Calcium

Substance	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p
Formula and Description			kcal/mol		cal/deg mol		
CaO·Fe ₂ O ₃	c	215.772	-361.787	-337.67	6.076	34.74	36.71
2CaO·Fe ₂ O ₃	c	271.851	-508.804	-478.44	7.564	45.12	46.19
CaFe(CN) ₆ ⁻	aq	252.034		+38.1			
CaFe(CN) ₆ ²⁻	aq			+28.6			
Ca ₂ Fe(CN) ₆	aq	292.114		-98.51		-2.7	
undissoc.; std. state, m = 1	aq			-105.6			
undissoc.; std. state, m = 1	aq			-149.27			
in 1,500 H ₂ O	aq			-149.22			
2,000 H ₂ O	aq			-149.28			
5,000 H ₂ O	aq			-149.37			
7,500 H ₂ O	aq			-149.44			
10,000 H ₂ O	aq			-149.59			
20,000 H ₂ O	aq			-149.76			
50,000 H ₂ O	aq			-149.86			
100,000 H ₂ O	aq			-150.17			
500,000 H ₂ O	aq			-150.37			
1,000,000 H ₂ O	aq			-150.6			
∞ H ₂ O	aq						



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Enthalpy and Gibbs Energy of Formation, Entropy and Heat Capacity

Table 94(25)

Calcium

Substance	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
$\text{Ca}_2\text{Fe}(\text{CN})_6 \cdot 11\text{H}_2\text{O}$	c	490.283	-903.7				
$\text{Ca}_3[\text{Fe}(\text{CN})_6]_2$ std. state, m = 1 in	aq	544.148	-120.6	-48.3		91.1	
	1,500 H ₂ O	aq	-118.91				
	2,000 H ₂ O	aq	-118.96				
	5,000 H ₂ O	aq	-119.13				
	7,500 H ₂ O	aq	-119.22				
	10,000 H ₂ O	aq	-119.28				
	20,000 H ₂ O	aq	-119.43				
	50,000 H ₂ O	aq	-119.65				
	100,000 H ₂ O	aq	-119.82				
	500,000 H ₂ O	aq	-120.20				
1,000,000 H ₂ O	aq	-120.32					
∞ H ₂ O	aq	-120.6					
$\text{CaH}_2\text{Fe}(\text{CN})_6$	aq	254.050	-20.2				
CaFeSiO_4	c	188.011	-456.				
CaCrO_4	aq	156.074	-340.9				

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Calcium

Table 94(26)

Substance	State	Formula Weight	298.15 K (25°C)				S°	C _p °
			ΔH _f ° 0 K	ΔH _f °	ΔG _f °	H ₂₉₈ ° - H ₀ °		
Formula and Description		kcal/mol		cal/deg mol				
CaMoO ₃	c	184.018		-296.				
CaMoO ₄	c	200.018		-368.4	-342.9	29.3	27.32	
	g		-197.					
	aq			-368.2		-6.2		
CaWO ₃	g	271.928	-113.					
CaWO ₄	c	287.928	-391.272	-393.20	-367.71	30.21	27.28	
	g		-225.					
	aq			-386.8				
				-556.71	-518.57	42.8	39.86	
CaO·V ₂ O ₅	c	237.960						
2CaO·V ₂ O ₅	c	294.040		-736.94	-691.49	52.7	50.08	
3CaO·V ₂ O ₅	c	350.119		-902.95	-851.12	65.7	60.29	
CaO·TiO ₂	c	135.978		-396.9	-376.5	22.38	23.34	
3CaO·2TiO ₂	c	328.036		-944.2	-896.6	56.1	57.20	
CaTiSiO ₅	c	196.063		-622.2	-588.4	30.88	33.21	
CaZrO ₃	c	197.298		-422.3	-401.8	23.92	23.88	

CaZrO₃
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SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

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Table 94 (27)		Calcium					
Substance		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity					
Formula and Description	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
			0 K	298.15 K (25°C)			
		kcal/mol					
CaHfO ₃	c	266.568	-433.0				
CaMg ₂	c	88.704	-9.6	-9.4	3.720	24.92	18.09
CaCl ₂ ·2MgCl ₂ ·2H ₂ O	c	337.453	-704.9				
CaMgC ₂ O ₆	c	184.411	-556.0	-517.1	6.210	37.09	37.65
CaO·MgO·SiO ₂	c	156.476	-540.89				
CaO·MgO·2SiO ₂	c	216.560	-766.3	-724.7		34.16	39.80
2CaO·MgO·2SiO ₂	amorp		-744.				
ankermanite	c	272.640	-926.67	-879.53		50.0	50.67
glassy	amorp		-918.9				
3CaO·MgO·2SiO ₂	c	328.719	-1091.7	-1037.4		60.5	60.29
mervinite	c		-2954.	-2780.	23.34	131.2	156.7
2CaO·5MgO·8SiO ₂ ·H ₂ O	c	812.410					
tremolite							

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

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Sr-SrO₂·8H₂O
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Table 95(1)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Strontium							
Substance		ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°			
Formula and Description		0 K	298.15 K (25°C)						
			kcal/mol					cal/deg mol	
	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°		
Sr	c	87.62	0	0		12.5	6.3		
Sr ⁺	g			39.3					
Sr ²⁺	g			172.11					
Sr	g			427.96					
	aq			-130.45	-133.71				
	g			1435.					
	g			2751.					
	g			4403.					
	g			6549.					
	g			8995.					
	g			11816.					
SrO	c	103.619		-141.5	-134.3	13.0	10.76		
SrO ₂	g			-2.					
	c	119.619		-151.4					
SrO ₂ ·8H ₂ O	c	263.742		-722.6					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

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Sr₂O
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Table 95(2)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Strontium						
Substance		Formula Weight	State	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
Formula and Description								
		kcal/mol						
Sr ₂ O		191.239	c					
SrH		88.628	g	-154.7				
SrH ₂		89.636	c	-56.	45.	2.080	50.80	7.17
SrOH		104.627	g	+52.				
SrOH ⁺			c	-43.1				
Sr(OH) ₂	std. state, m = 1	121.635	aq	-41.	-172.4			
			c	-229.2				
			g	-135.				
	in 800 H ₂ O		aq	-240.1				
Sr(OH) ₂ ·H ₂ O		139.650	c	-302.3				
Sr(OH) ₂ ·8H ₂ O		265.757	c	-801.2				
SrF		106.618	g	-69.0	-62.9	2.219	57.31	8.27
SrF ₂		125.617	c	-290.7	-278.4	3.125	19.63	16.73
			g	-182.7	-185.3	3.192	69.54	12.66

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National Bureau of Standards

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Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Strontium

Table 95(3)

Substance		State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°	
Formula and Description										298.15 K (25°C)
		kcal/mol							cal/deg mol	
SrCl		g	123.073							
SrCl ₂	α	c	158.526							
	std. state, m = 1	g								
	in	aq								
	200 H ₂ O	aq		-30.2	-36.5	2.338	60.2	8.73		
	300 H ₂ O	aq		-198.1	-186.7	3.880	27.45	18.07		
	400 H ₂ O	aq		-116.1	-118.6	3.454	74.26	13.33		
	500 H ₂ O	aq		-210.35	-196.45		19.2			
	800 H ₂ O	aq		-209.78						
	1,000 H ₂ O	aq		-209.83						
	2,000 H ₂ O	aq		-209.871						
	5,000 H ₂ O	aq		-209.900						
	10,000 H ₂ O	aq		-209.955						
	20,000 H ₂ O	aq		-209.977						
	50,000 H ₂ O	aq		-210.057						
	100,000 H ₂ O	aq		-210.144						
	500,000 H ₂ O	aq		-210.195						
	1,000,000 H ₂ O	aq		-210.236						
	∞ H ₂ O	aq		-210.274						
		aq		-210.295						
		aq		-210.325						
		aq		-210.332						
		aq		-210.35						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

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SrCl₂

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Table 95(4)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Substance		Strontium						
Formula and Description	State	Formula Weight	ΔHf°	ΔHf°	ΔGf°	H ₂₉₈ ° - H ₀ °	S°	C _p °
			0 K	298.15 K (25°C)				
SrCl ₂	in 28 (HCl + 59.5 H ₂ O)	aq		-209.2				
	37 (HCl + 44.6 H ₂ O)	aq		-209.2				
	56 (HCl + 29.8 H ₂ O)	aq		-208.8				
	115 (HCl + 14.4 H ₂ O)	aq		-207.7				
	726 (HCl + 12.8 H ₂ O)	aq		-207.2				
SrCl ₂ ·H ₂ O	c	176.541		-271.7			41.	28.7
SrCl ₂ ·2H ₂ O	c	194.557		-343.7			52.	38.3
SrCl ₂ ·6H ₂ O	c	266.618		-627.1			93.4	
Sr(OCl) ₂	aq	190.525		-180.5				
Sr(ClO ₂) ₂	aq	222.524		-162.				
Sr(ClO ₄) ₂	c	286.521		-182.31				
	std. state, m = 1			-192.27			79.2	

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Sr(ClO₄)₂

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Table 95(5)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Strontium					
Substance	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
Formula and Description		kcal/mol					
Sr(ClO ₄) ₂	in		0 K				
	13 H ₂ O	aq		-191.85			
	15 H ₂ O	aq		-192.03			
	20 H ₂ O	aq		-192.21			
	25 H ₂ O	aq		-192.28			
	30 H ₂ O	aq		-192.30			
	40 H ₂ O	aq		-192.252			
	50 H ₂ O	aq		-192.204			
	75 H ₂ O	aq		-192.111			
	100 H ₂ O	aq		-192.052			
	150 H ₂ O	aq		-191.996			
	200 H ₂ O	aq		-191.975			
	300 H ₂ O	aq		-191.960			
400 H ₂ O	aq		-191.958				
500 H ₂ O	aq		-191.962				
800 H ₂ O	aq		-191.978				
1,000 H ₂ O	aq		-191.992				
2,000 H ₂ O	aq		-192.033				
5,000 H ₂ O	aq		-192.092				
10,000 H ₂ O	aq		-192.132				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

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Table 95(6)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Strontium					
Substance		ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°	
Formula and Description		0 K	298.15 K (25°C)				
		Formula Weight	kcal/mol				
		State	cal/deg mol				
Sr(ClO ₄) ₂	in 20,000 H ₂ O	aq	-192.165				
	50,000 H ₂ O	aq	-192.199				
	100,000 H ₂ O	aq	-192.218				
	500,000 H ₂ O	aq	-192.245				
	1,000,000 H ₂ O	aq	-192.252				
	∞ H ₂ O	aq	-192.27				
in CH ₃ OH							
	C ₂ H ₅ OH		-206.9				
	n-C ₃ H ₇ OH		-198.0				
	i-C ₃ H ₇ OH		-195.7				
	n-C ₄ H ₉ OH		-197.31				
Sr(ClO ₄) ₂ ·2H ₂ O		c	-194.7				
Sr(ClO ₄) ₂ ·4H ₂ O		c	-328.5				
SrCl ₂ ·SrO·H ₂ O		c	-469.0				
SrCl ₂ ·SrO·9H ₂ O		c	-437.2				
		c	-1006.6				

 Sr(ClO₄)₂
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SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

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Table 95(7) Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity
Strontium

Substance	Formula and Description	State	Formula Weight	ΔH_f^0	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °	
										0 K
kcal/mol										
SrClH		c	124.081		-125.5					
SrFCl		c	142.071		-250.0					
SrBr ₂		c	247.438		-171.5	-166.6	4.261	32.29	18.01	
	std. state, m = 1	g			-98.	-106.	3.8	77.3	14.5	
	in	aq			-188.55	-183.41		31.6		
	400 H ₂ O	aq			-188.14					
	500 H ₂ O	aq			-188.16					
	800 H ₂ O	aq			-188.21					
	1,000 H ₂ O	aq			-188.225					
	2,000 H ₂ O	aq			-188.286					
	5,000 H ₂ O	aq			-188.359					
	10,000 H ₂ O	aq			-188.405					
	20,000 H ₂ O	aq			-188.442					
	50,000 H ₂ O	aq			-188.477					
	100,000 H ₂ O	aq			-188.497					
500,000 H ₂ O	aq			-188.525						
1,000,000 H ₂ O	aq			-188.532						
∞ H ₂ O	aq			-188.55						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

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Table 95(8)		Strontium						
Substance		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity					SrBr ₂ ·H ₂ O	
Formula and Description	State	Formula Weight	ΔHf°	ΔGf°	H ₂₉₈ ° - H ₀ °	S°	C _p °	
			0 K	298.15 K (25 °C)				cal/deg mol
SrBr ₂ ·H ₂ O	c	265.453						
SrBr ₂ ·6H ₂ O	c	355.530						
Sr(BrO ₃) ₂ ·H ₂ O	c	361.450	-246.5	-228.1		43.	28.9	
SrBr ₂ ·SrO·3H ₂ O	c	405.103	-605.0	-519.7		97.	82.1	
SrBr ₂ ·SrO·9H ₂ O	c	513.195	-264.	-189.1		67.		
SrBrH	c	168.537	-553.0					
SrI ₂	c	341.429	-985.0					
	g		-111.0				19.5	
	aq		-133.4					
			-65.			45.4		
	std. state, m = 1		-156.83	-158.37			28.5	
SrI ₂ ·H ₂ O	c	359.444	-211.9				39.1	
SrI ₂ ·2H ₂ O	c	377.459	-282.6				84.9	
SrI ₂ ·6H ₂ O	c	449.521	-570.9					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity
Strontium

Table 95(9)

Substance	Formula and Description	State	Formula Weight	ΔH_f°		ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
				0 K	298.15 K (25°C)				
				kcal/mol					
Sr(IO ₃) ₂		c	437.425		-243.6	-204.4		56.	
Sr(IO ₃) ₂ ·H ₂ O		c	455.441		-313.2	-260.4		66.	
Sr(IO ₃) ₂ ·6H ₂ O		c	545.517		-666.8	-543.7		109.	
SrIH		c	215.532		-92.2				
SrS		c	119.684		-108.3	-107.2		16.3	11.64
		g			26.				
SrSO ₃		c	167.682		-281.3				
SrSO ₄		c	183.682		-347.3	-320.5		28.	
	precipitate	c			-346.5				
	std. state, m = 1	aq			-347.77	-311.68			
SrS ₂ O ₆ ·4H ₂ O		c	319.806		-699.1				
SrI ₂ ·2SO ₂		c	469.554		-297.4				
SrI ₂ ·4SO ₂		c	597.680		-460.8				

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SrSe
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Table 95(10)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Strontium						
Substance		Formula Weight	State	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
Formula and Description								
		kcal/mol						
		cal/deg mol						
SrSe		166.58	c		-92.2			
SrSeO ₃		214.578	c		-250.4			
SrSeO ₄		230.578	c		-273.1			
Sr(N ₃) ₂		171.660	c		+2.1			
Sr ₃ N ₂		290.873	c		-93.5			
SrN ₂ O ₂ ·5H ₂ O		237.709	c		-476.1			
Sr(NO ₂) ₂		179.631	c		-182.2			
	in 800 H ₂ O		aq		-180.5			
		197.646	c		-254.2			
Sr(NO ₂) ₂ ·H ₂ O		211.630	c		-233.80	6.854	46.50	35.83
Sr(NO ₃) ₂			aq		-229.57		62.2	
	std. state, m = 1		aq		-231.713			
	in 20 H ₂ O		aq		-231.435			
	25 H ₂ O		aq		-231.210			
	30 H ₂ O		aq		-230.857			
	40 H ₂ O		aq		-230.595			
	50 H ₂ O		aq		-230.180			
	75 H ₂ O		aq					

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

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Table 95(11)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Strontium						
Substance		Formula Weight	State	ΔHf° 0 K	ΔGf°	H ₂₉₈ ° - H ₀ °	S°	C _p °
Formula and Description								
		kcal/mol				cal/deg mol		
Sr(NO ₃) ₂	in		aq					
		100 H ₂ O		aq	-229.966			
		150 H ₂ O		aq	-229.751			
		200 H ₂ O		aq	-229.590			
		300 H ₂ O		aq	-229.533			
		400 H ₂ O		aq	-229.486			
		500 H ₂ O		aq	-229.454			
		800 H ₂ O		aq	-229.414			
		1,000 H ₂ O		aq	-229.405			
		2,000 H ₂ O		aq	-229.396			
		5,000 H ₂ O		aq	-229.424			
		10,000 H ₂ O		aq	-229.451			
		20,000 H ₂ O		aq	-229.476			
		50,000 H ₂ O		aq	-229.505			
		100,000 H ₂ O		aq	-229.521			
	500,000 H ₂ O		aq	-229.546				
	1,000,000 H ₂ O		aq	-229.553				
	∞ H ₂ O		aq	-229.57				
	Sr(NO ₃) ₂ ·4H ₂ O	283.691	c		-413.65		88.2	

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

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Sr(NH₂)₂
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Table 95 (12)		Strontium						
Substance		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Formula and Description	State	Formula Weight	ΔHf°	ΔHf°	ΔGf°	H ₂₉₈ ° - H ₀ °	S°	C _p °
			0 K	298.15 K (25°C)				
				kcal/mol				
Sr(NH ₂) ₂	c	119.665		-83.1				
Sr(NH ₂) ₆	c	189.804		-123.7				
SrCl ₂ ·NH ₃	c	175.557		-220.6				
SrCl ₂ ·8NH ₃	c	294.771		-367.1				
Sr(ClO ₄) ₂ ·2NH ₃	c	320.582		-226.2				
Sr(ClO ₄) ₂ ·6NH ₃	c	388.705		-314.5				
Sr(ClO ₄) ₂ ·7NH ₃	c	405.735		-334.6				
SrBr ₂ ·NH ₃	c	264.469		-199.3				
SrBr ₂ ·2NH ₃	c	281.499		-223.1				
SrBr ₂ ·8NH ₃	c	383.683		-354.9				
SrI ₂ ·NH ₃	c	358.459		-162.7				
SrI ₂ ·2NH ₃	c	375.490		-189.2				
SrI ₂ ·6NH ₃	c	443.612		-283.5				
SrI ₂ ·8NH ₃	c	477.674		-328.0				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

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Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Strontium

Table 95(13)	Substance	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
				0 K	298.15 K (25°C)				
Formula and Description				kcal/mol					
	$Sr(NH_4)_2(SO_4)_2$	c	315.820		-630.8				
	Sr_3P_2	c	324.808		-152.				
	$Sr_3(PO_4)_2$	c	452.803		-985.4				
	$SrHPO_4$	c	183.599		-435.4	-403.6		29.	
	$Sr(H_2PO_4)_2$	c	281.595		-749.2				
	Sr_3As_2	c	412.703		-147.3				
	$Sr_3(AsO_4)_2$, hydrated precipitate	c	540.698		-792.8 -803.	-736.2		61.	
	$SrHASO_4$	aq	227.547		-345.7				
	$Sr(H_2AsO_4)_2$	aq	369.490		-563.2				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

SrSb
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Table 95(14)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Strontium						
Substance	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°	
			0 K	298.15 K (25°C)				
Formula and Description			kcal/mol				cal/deg mol	
SrSb	c	209.37	-46.					
Sr ₂ Sb	c	296.99	-77.					
Sr ₃ Sb ₂	c	506.36	-135.					
SrBi	c	296.600	-43.					
Sr ₂ Bi	c	384.220	-75.					
Sr ₃ Bi ₂	c	680.820	-127.					
SrC ₂	c	111.642	-18.					
SrCO ₃	c	147.629	-291.6	-272.5		23.2	19.46	
strontianite			-292.29	-259.88		-21.4		
std. state, m = 1	aq		-327.6					
SrC ₂ O ₄	c	175.640	-327.6					
std. state, m = 1	aq			-294.8		+2.1		

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

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Table 95(15)		Strontium					
Substance		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity					
Formula and Description	State	Formula Weight	ΔHf°	ΔGf°	H ₂₉₈ ^o - H ₀ ^o	S°	C _p ^o
			298.15 K (25°C)				
		kcal/mol					
		cal/deg mol					
Sr(OH) ₂	c	177.656	-333.0				
Sr(OH) ₂ ·2H ₂ O	c	213.687	-475.8				
Sr(HCO ₃) ₂	aq	209.655	-460.8				
SrC ₂ H ₃ O ₄ ⁺	aq	146.665		-223.5			
Sr(C ₂ H ₃ O ₂) ₂	c	205.710	-355.5				
	aq		-361.4				
Sr(C ₂ H ₃ O ₂) ₂ ·0.5H ₂ O	c	214.718	-390.0				
Sr(C ₂ H ₃ O ₂) ₂	c	237.709	-441.4				
	aq		-441.2				
SrBr ₂ ·1/2C ₂ H ₅ OH	c	270.473	-205.				
SrCN ₂	c	127.645	-72.5				
Sr(CN) ₂	aq	139.656	-57.0				
Sr(CN) ₂ ·4H ₂ O	c	211.717	-333.7				

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National Bureau of Standards

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SrSi
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Table 95(16)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Strontium						
Substance		State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
Formula and Description								
				kcal/mol				
				cal/deg mol				
SrSi		c	115.706	-133.				
SrSi ₂		c	143.792	-188.				
Sr ₂ Si		c	203.326	-116.				
SrSiO ₃		c	163.704	-390.5	-370.4		23.1	21.16
Sr ₂ SiO ₄	glassy	amorp	267.324	-380.				
Sr ₂ Ce		c	247.83	-550.8	-523.7		36.6	32.09
Sr ₂ Sn		c	293.93	-75.				
Sr ₂ Pb		c	382.43	-83.				
SrI ₂ ·2PbI ₂		c	1263.426	-80.				
SrI ₂ ·2PbI ₂ ·7H ₂ O		c	1389.534	-222.4				
				-714.0				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Strontium

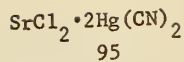
Table 95(17)

Substance	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
Formula and Description			0 K	kcal/mol		cal/deg mol		
Sr(BF ₄) ₂	c	261.229		-921.				
SrO·Al ₂ O ₃	c	205.581		-555.6				
3SrO·Al ₂ O ₃	c	412.819		-847.2				
4SrO·Al ₂ O ₃ α	c	516.439		-978.2				
β	c			-976.7				
3SrCl ₂ ·4AlCl ₃	c	1008.940		-1284.4				
2SrO·Al ₂ O ₃ ·SiO ₂	c	369.285		-941.1				
SrBr ₂ ·HgBr ₂	aq	607.846		-229.2				
SrBr ₂ ·2HgBr ₂	aq	968.254		-268.0				
2SrBr ₂ ·HgBr ₂	aq	855.284		-419.6				
SrHg(CN) ₄	aq	392.281		-3.9				
Sr[Hg(CN) ₃] ₂	aq	644.907		+61.8				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

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Table 95(18)		Strontium						
Substance		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Formula and Description	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°	
			0 K	298.15 K (25°C)				cal/deg mol
$\text{SrCl}_2 \cdot 2\text{Hg}(\text{CN})_2$	aq	663.777		-77.3				
$\text{SrCl}_2 \cdot 2\text{Hg}(\text{CN})_2 \cdot 6\text{H}_2\text{O}$	c	771.869		-503.7				
$\text{SrBr}_2 \cdot 2\text{Hg}(\text{CN})_2$	aq	752.689		-56.4				
$\text{SrBr}_2 \cdot 2\text{Hg}(\text{CN})_2 \cdot 6\text{H}_2\text{O}$	c	860.781		-485.7				
$\text{SrI}_2 \cdot 2\text{Hg}(\text{CN})_2$	aq	846.680		-29.1				
$\text{SrI}_2 \cdot 2\text{Hg}(\text{CN})_2 \cdot 7\text{H}_2\text{O}$	c	972.788		-528.0				
$\text{SrAg}(\text{CN})_3$	aq	273.544		-39.3				
$\text{Sr}[\text{Ag}(\text{CN})_2]_2$	aq	407.431		-0.4				
$\text{SrNi}(\text{CN})_4$	aq	250.401		-39.3				
$2\text{SrO} \cdot \text{Fe}_2\text{O}_3$	c	366.931		-510.9				
$3\text{SrO} \cdot \text{Fe}_2\text{O}_3$	c	470.550		-686.3				
$7\text{SrO} \cdot 5\text{Fe}_2\text{O}_3$	c	1523.797		-2038.				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Table 95(19)

Strontium

Substance	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
Formula and Description			kcal/mol					
SrFe(CN) ₆ ⁻	aq	299.574			+36.7			
Sr ₂ Fe(CN) ₆	aq	387.194			-101.33		+7.1	
in	aq			-152.0				
std. state, m = 1	aq			-150.04				
std. state, m = 1	aq			-150.08				
1,500 H ₂ O	aq			-150.21				
2,000 H ₂ O	aq			-150.29				
5,000 H ₂ O	aq			-150.36				
7,500 H ₂ O	aq			-150.57				
10,000 H ₂ O	aq			-150.84				
20,000 H ₂ O	aq			-151.03				
50,000 H ₂ O	aq			-151.80				
100,000 H ₂ O	aq			-151.49				
200,000 H ₂ O	aq			-151.73				
500,000 H ₂ O	aq			-152.0				
1,000,000 H ₂ O	aq							
∞ H ₂ O	aq							

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Strontium

Table 95(20)

Substance	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
Formula and Description			0 K			298.15 K (25°C)		
Sr ₃ [Fe(CN) ₆] ₂ std. state, m = 1 in	aq	686.768						
	1,500 H ₂ O			-122.8	-52.5		105.8	
	2,000 H ₂ O			-121.16				
	5,000 H ₂ O			-121.22				
	7,500 H ₂ O			-121.43				
	10,000 H ₂ O			-121.54				
	20,000 H ₂ O			-121.62				
	50,000 H ₂ O			-121.79				
	100,000 H ₂ O			-121.99				
	200,000 H ₂ O			-122.12				
	500,000 H ₂ O			-122.25				
	1,000,000 H ₂ O			-122.43				
	∞ H ₂ O			-122.55				
Sr ₃ [FeCO(CN) ₅] ₂	aq			-122.8				
	c	690.754		-258.7				
Sr ₃ [FeCO(CN) ₅] ₂ •4H ₂ O	aq			-300.3				
	c	762.815		-540.0				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Table 95(21)

Strontium

Substance	Formula and Description	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
SrMoO ₃		c	231.558	-306.				
SrMoO ₄		g	247.558	-130.				
		c		-370.				
		g		-235.				
SrWO ₃		g	319.468	-140.				
SrWO ₄		c	335.468	-391.9	-366.		33.	
		g		-245.				
SrTiO ₃		c	183.518	-399.71	-379.64		26.0	23.51
Sr ₂ TiO ₄		c	287.138	-546.7	-520.7		38.0	34.34
SrZrO ₃		c	226.838	-422.4	-402.2		27.5	24.71
SrHfO ₃		c	314.108	-433.7				
SrMg ₂	β phase	c	136.244	-5.1				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

Table 96(1)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Substance		Barium						
Formula and Description	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
			0 K	298.15 K (25°C)				
Ba	c	137.34	0	0	0	1.65	15.0	6.71
Ba ⁺	g		43.2	43.	35.	1.481	40.663	4.968
Ba ²⁺	g		163.39	164.67				
Ba	g		394.09	396.86				
	aq			-128.50	-134.02		+2.3	
BaO	c	153.339	-132.00	-132.3	-125.5	2.385	16.83	11.42
	g			-28.				7.9
BaO ₂	c	169.339		-151.6				16.0
BaO ₂ ·H ₂ O	c	187.354		-222.3				
BaO ₂ ·8H ₂ O	c	313.462		-718.6				
Ba ₂ O	c	290.679		-147.1				
	g			-72.				
Ba ₂ O ₂	g	306.679		-135.				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Barium

Table 96(2)

Substance	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
			0 K	298.15 K (25°C)				
BaH	g	138.348	53.6	53.	47.	2.082	52.29	7.19
BaH ₂	c	139.356		-42.7				
BaOH	g	154.347		-52.				
BaOH ⁺	aq				-174.6			
Ba(OH) ₂	c	171.355		-225.8				
	g			-140.				
	aq			-237.9				
	c	189.370		-298.4				
Ba(OH) ₂ ·H ₂ O	c	225.401		-442.0				
Ba(OH) ₂ ·3H ₂ O	c	315.477		-798.8			102.	
Ba(OH) ₂ ·8H ₂ O	c			-208.5				
BaO ₂ ·H ₂ O ₂	c	203.354						

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

BaF
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Table 96(3)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Barium						
Substance		ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°	
Formula and Description		0 K	298.15 K (25°C)					
		kcal/mol						
		cal/deg mol						
State	Formula Weight							
BaF	g	156.338	-77.53	-78.0	-83.8	2.232	58.8	8.33
BaF ₂	c	175.337	-288.19	-288.5	-276.5	3.452	23.03	17.02
	g		-195.0	-195.5	-198.0	3.262	71.91	12.85
	aq			-287.50	-267.30		-4.3	
BaCl	c	172.793		-110.	-46.	2.36	61.8	8.80
	g			-40.				
BaCl ₂	c	208.246	-205.35	-205.2	-193.7	3.993	29.56	17.96
	g		-125.37	-125.7	-128.5	3.511	77.72	13.43
	aq			-208.40	-196.76		29.3	
	aq			-207.747				
	aq			-207.754				
	aq			-207.762				
	aq			-207.783				
	aq			-207.803				
	aq			-207.836				
	aq			-207.863				
	std. state, m = 1							
	in							
	30 H ₂ O'							
	40 H ₂ O							
	50 H ₂ O							
	75 H ₂ O							
	100 H ₂ O							
	150 H ₂ O							
	200 H ₂ O							

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

Washington, D. C.

National Bureau of Standards

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Table 96(4)

Barium

BaCl ₂	Substance	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
				0 K	kcal/mol				
	in	aq							
	300 H ₂ O	aq			-207.905				
	400 H ₂ O	aq			-207.938				
	500 H ₂ O	aq			-207.963				
	800 H ₂ O	aq			-208.012				
	1,000 H ₂ O	aq			-208.036				
	2,000 H ₂ O	aq			-208.110				
	5,000 H ₂ O	aq			-208.197				
	10,000 H ₂ O	aq			-208.249				
	20,000 H ₂ O	aq			-208.288				
	50,000 H ₂ O	aq			-208.326				
	100,000 H ₂ O	aq			-208.346				
	500,000 H ₂ O	aq			-208.375				
	1,000,000 H ₂ O	aq			-208.382				
	∞ H ₂ O	aq			-208.40				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

BaCl₂
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Table 96(5)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity					
Substance		Barium					
Formula and Description	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
			0 K	298.15 K (25°C)			
BaCl ₂ in 4.0 mola1 HCl 3.0 mola1 HCl 2.0 mola1 HCl 1.0 mola1 HCl 0.1 mola1 HCl	aq		-205.1				
	aq		-205.7				
	aq		-206.4				
	aq		-207.3				
	aq		-208.3				
BaCl ₂ •H ₂ O	c	226.261	-277.4	-252.32		39.9	
BaCl ₂ •2H ₂ O	c	244.277	-348.98	-309.86		48.5	38.71
Ba(OCl) ₂	aq	240.245	-179.3				
Ba(ClO) ₂	c	272.244	-162.6				
Ba(ClO ₂) ₂ •3.5H ₂ O	c	335.297	-407.9				
Ba(ClO ₃) ₂	c	304.242	-182.3				
Ba(ClO ₃) ₂ in 400 H ₂ O	aq		-176.2				
Ba(ClO ₃) ₂ •H ₂ O	c	322.258	-255.5				50.6
Ba(ClO ₄) ₂	c	336.241	-191.2				

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

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Table 96(6)		Barium						
Substance		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Formula and Description	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
			0 K	298.15 K (25°C)				
		kcal/mol						
		cal/deg mol						
Ba(ClO ₄) ₂ in 900 H ₂ O in ethanol, ∞ diln in propanol, ∞ diln in i-propanol, ∞ diln in n-butanol, ∞ diln	aq							
		390.287	-190.2					
		415.631	-196.6					
		451.662	-195.4					
		505.708	-196.0					
BaCl ₂ ·3H ₂ O	c		-194.8				94.	
BaCl ₂ ·BaO·3H ₂ O	c		-404.3	-303.7				
BaCl ₂ ·BaO·5H ₂ O	c		-578.0					
BaCl ₂ ·BaO·8H ₂ O	c		-727.0					
BaClH	c	173.801	-128.1					
BaClF	c	191.791	-254.4					
BaBr ₂	c	297.158	-181.0	-176.1		35.		
	g		-101.4	-113.	3.9	79.		14.7

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

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Table 96 (7)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Substance		Barium		ΔHf°	ΔGf°	H ₂₉₈ ^o - H ₀ ^o	S°	C _p ^o
Formula and Description	State	Formula Weight	0 K					
				kcal/mol		cal/deg mol		
BaBr ₂	std. state, m = 1							
	in	aq					41.7	
	400 H ₂ O	aq		-186.60	-183.72			
	500 H ₂ O	aq		-186.20				
	800 H ₂ O	aq		-186.22				
	1,000 H ₂ O	aq		-186.26				
	2,000 H ₂ O	aq		-186.280				
	5,000 H ₂ O	aq		-186.339				
	10,000 H ₂ O	aq		-186.408				
	20,000 H ₂ O	aq		-186.453				
	50,000 H ₂ O	aq		-186.490				
	100,000 H ₂ O	aq		-186.526				
	500,000 H ₂ O	aq		-186.546				
	1,000,000 H ₂ O	aq		-186.575				
	∞ H ₂ O	aq		-186.582				
BaBr ₂ ·H ₂ O		c	315.173					
BaBr ₂ ·2H ₂ O		c	333.189		-294.1			54.

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National Bureau of Standards

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Table 96(8)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Barium						
Substance	Formula and Description	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
Ba(BrO) ₂		aq	329.157	-174.5				
Ba(BrO ₃) ₂		c	393.154	-179.89	-138.0		58.	
	in 100,000 H ₂ O	aq		-168.45				
	50,000 H ₂ O	aq		-168.61				
	20,000 H ₂ O	aq		-169.17				
Ba(BrO ₃) ₂ ·H ₂ O		c	411.170	-252.1	-197.09		69.9	52.90
BaBr ₂ ·BaO·2H ₂ O		c	486.528	-486.0				
BaBr ₂ ·BaO·5H ₂ O		c	540.574	-705.4				
BaBrH		c	218.257	-116.4				
BaI ₂		c	391.149	-143.9				
		g		-77.2			82.	14.7
	std. state, m = 1	aq		-154.88	-158.68	4.0	55.5	

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Table 96(9)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Substance		Barium						
Formula and Description	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
			0 K	298.15 K (25°C)				
		kcal/mol						
BaI ₂ •H ₂ O	c	409.164	-219.7					
BaI ₂ •2H ₂ O	c	427.179	-290.8					
BaI ₂ •2.5H ₂ O	c	436.187	-325.9					
BaI ₂ •7H ₂ O	c	517.256	-639.7					
Ba(IO ₃) ₂	c	487.145	-243.31		-206.7	8.84	59.6	44.8
std. state, m = 1	aq				-195.2		58.9	
Ba(IO ₃) ₂ •H ₂ O	c	505.161	-316.0		-263.9		71.	
BaI ₂ •BaO•2H ₂ O	c	580.519	-455.3					
BaI ₂ •BaO•9H ₂ O	c	706.626	-955.0					
BaIH	c	265.252	-98.1					
BaS	c	169.404	-110.		-109.		18.7	11.80
	g		+12.					
	aq		-117.9					
Ba ₂ S	g	338.808	-90.					

BaI₂•H₂O
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Table 96(10)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity					
Substance		Barium					C _p ^o
Formula and Description	State	Formula Weight	ΔH _f ^o	ΔG _f ^o	H ₂₉₈ ^o - H ₀ ^o	S ^o	
			kcal/mol		298.15 K (25°C)		cal/deg mol
BaSO ₃	c	217.402	-281.9				
BaSO ₄	c	233.402	-352.1	-325.6		31.6	24.32
precipitated	c		-350.5				
std. state, m = 1	aq		-345.82	-311.99		7.1	40.7
BaS ₂ O ₃	c	249.466	-415.4				
BaS ₂ O ₆	aq	297.464	-558.6				
BaS ₂ O ₆ ·2H ₂ O	c	333.495	-453.7				
BaS ₂ O ₈	aq	329.463	-738.4				
BaS ₂ O ₈ ·4H ₂ O	c	401.525	-564.3				
BaS ₄ O ₆ ·2H ₂ O	c	397.623					
Ba(HS) ₂	aq	203.484	-134.3				

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Table 96(11)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity Barium						
Substance		Formula Weight	State	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
Formula and Description								
				0 K				
Ba(HSO ₃) ₂	aq	299.480						
BaSO ₄ ·H ₂ SO ₄	c	331.479		-430.0				
BaSO ₄ ·2H ₂ SO ₄ ·H ₂ O	c	447.572		-548.6				
				-819.8				
BaI ₂ ·2SO ₂	c	519.274		-308.5				
BaI ₂ ·4SO ₂	c	647.400		-470.2				
BaSe	c	216.30		-89.				
BaSeO ₃	c	264.298		-248.7	-231.4		40.	
BaSeO ₄	c	280.298		-274.0	-249.7		42.	
BaN ₂	c	165.353		-41.				
Ba(N ₃) ₂ ·H ₂ O	c	239.396		-73.7	-25.1		45.	
Ba ₂ N	c	288.687		-52.0				
Ba ₃ N ₂	c	440.033		-86.8				

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Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Table 96(12)

Barium

Substance	Formula	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
BaNO ₃ ⁺	std. state, m = 1	aq	199.345	-175.8	-159.35		41.	
Ba(NO ₂) ₂		c	229.351	-183.6				
		aq		-178.5				
Ba(NO ₂) ₂ ·H ₂ O		c	247.366	-254.8				
Ba(NO ₃) ₂		c	261.350	-237.11	-190.42		51.1	36.18
	std. state, m = 1	aq		-227.62	-187.24		72.3	
	in	aq		-228.474				
	200 H ₂ O	aq		-228.160				
	300 H ₂ O	aq		-228.005				
	400 H ₂ O	aq		-227.914				
	500 H ₂ O	aq		-227.770				
	800 H ₂ O	aq		-227.719				
	1,000 H ₂ O	aq		-227.617				
	2,000 H ₂ O	aq		-227.556				
	5,000 H ₂ O	aq						

Ba(NO₃)₂
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Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Barium

Table 96(13)

Substance	State	Formula Weight	ΔH_{f0}°	ΔH_f°	ΔG_f°	$H_{298}^{\circ} - H_0^{\circ}$	S ^o	C _p ^o
Formula and Description		kcal/mol						
Ba(NO ₃) ₂	in	10,000 H ₂ O		-227.545				
		20,000 H ₂ O		-227.549				
		50,000 H ₂ O		-227.563				
		100,000 H ₂ O		-227.575				
		500,000 H ₂ O		-227.597				
		1,000,000 H ₂ O		-227.603				
		∞ H ₂ O		-227.62				
	in 500 NH ₃ (liq)		-252.42					
BaNH	c	152.355		-53.4				
Ba(NH ₂) ₂	c	169.385		-78.5				
Ba(NH ₃) ₆	c	239.524		-124.0				

Ba(NO₃)₂
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National Bureau of Standards

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Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Barium

Table 96(14)

Substance	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	s°	C _p °
			0 K				
Ba ₂ NO ₃	c	324.140		-147.3			
BaCl ₂ ·8NH ₃	c	344.491		-364.5			
Ba(ClO ₄) ₂ ·2NH ₃	c	370.302		-233.0			
Ba(ClO ₄) ₂ ·6NH ₃	c	438.425		-314.0			
Ba ₂ NBr	c	368.596		-137.0			
BaBr ₂ ·NH ₃	c	314.189		-203.9			
BaBr ₂ ·2NH ₃	c	331.219		-225.5			
BaBr ₂ ·4NH ₃	c	365.280		-267.9			
BaBr ₂ ·8NH ₃	c	433.403		-351.8			
Ba ₆ Ni ₉	c	1980.186		-701.			

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National Bureau of Standards

Washington, D. C.

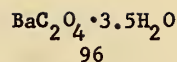
Table 96(15)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Substance		Barium						
Formula and Description	State	Formula Weight	ΔHf°	ΔGf°	H ₂₉₈ ° - H ₀ °	S°	C _p °	
			0 K	298.15 K (25°C)				
		kcal/mol		cal/deg mol				
BaI ₂ ·2NH ₃	c	425.210	-192.9					
BaI ₂ ·4NH ₃	c	459.271	-237.5					
BaI ₂ ·6NH ₃	c	493.332	-281.7					
BaI ₂ ·8NH ₃	c	527.394	-325.1					
BaI ₂ ·9NH ₃	c	544.424	-346.1					
BaI ₂ ·10NH ₃	c	561.455	-364.7					
Ba ₃ (PO ₄) ₂	colloidal	601.963	-978.					
BaHPO ₄	c	233.319	-433.7					
Ba(H ₂ PO ₄) ₂	c	267.317	-421.2					
	in 400 H ₂ O		-421.4					
	800 H ₂ O		-421.6					
	1,200 H ₂ O		-421.8					
	∞ H ₂ O		-421.9					
Ba(H ₂ PO ₄) ₂ ·H ₂ O	c	285.332	-490.5					
Ba(H ₂ PO ₄) ₂	c	331.315	-747.					

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Table 96(16)		Barium						
		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
		Substance	Formula Weight	State	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°
0 K					298.15 K (25°C)			
Formula and Description			kcal/mol					
$Ba_3(AsO_4)_2$	hydrated precipitate	689.858		-819.				
$BaHAsO_4 \cdot H_2O$	c	295.283		-412.6				
$Ba(H_2AsO_4)_2 \cdot 2H_2O$	c	455.241		-696.9				
BaC_2	c	161.362		-18.				
$BaCO_3$	c	197.349		-290.7	-271.9		26.8	20.40
	witherite			-290.34	-260.19		-11.3	
	std. state, m = 1			-327.1				
BaC_2O_4	c	225.360		-365.4				
$BaC_2O_4 \cdot 0.5H_2O$	c	234.368		-471.1				
$BaC_2O_4 \cdot 2H_2O$	c	261.391		-576.6				
$BaC_2O_4 \cdot 3.5H_2O$	c	288.414						



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Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Barium

Table 96(17)

Substance	State	Formula Weight	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
Formula and Description			kcal/mol				
Ba(CHO ₂) ₂	c	227.376	-333.6				
barium formate in 440 H ₂ O	aq		-331.6				
Ba(HCO ₃) ₂	aq	259.375	-459.28	-414.54		45.9	
Ba(C ₂ H ₃ O ₂) ₂	c	255.430	-354.8				
std. state, m = 1	aq		-360.82	-310.60		43.7	
in 500 H ₂ O	aq		-360.2				
Ba(C ₂ H ₃ O ₂) ₂ ·3H ₂ O	c	309.476	-566.2				
Ba(C ₂ H ₃ O ₂) ₃ /2	c	287.429	-444.7				
barium glycolate	aq		-439.6				
Ba(C ₂ H ₅ O) ₂	c	227.463	-218.6				
3BaO·4CH ₃ OH	c	588.188	-712.2				
3BaO·4C ₂ H ₅ OH	c	644.296	-735.2				
BaCS ₃	c	245.543	-127.				

Ba(CHO₂)₂
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Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Table 96(18)

Barium

Substance	State	Formula Weight	ΔH_f°	ΔH_f°	$H_{298}^\circ - H_0^\circ$	s°	C _p °
Formula and Description			kcal/mol				
Ba(HSO ₃) ₂ ·C ₂ H ₂ O ₂	aq	357.517		-522.6			
Ba(HSO ₃) ₂ ·C ₂ H ₂ O ₂ ·2.5H ₂ O glyoxal barium bisulfite	c	402.556		-701.6			
Ba(C ₂ H ₅ SO ₄) ₂ in 400 H ₂ O	aq	387.588		-547.3			
Ba(C ₂ H ₅ SO ₄) ₂ ·2H ₂ O	c	423.618		-688.4			
BaCN ₂	c	177.365		-63.6			
Ba(CN) ₂	c	189.376		-52.2			
Ba(CN) ₂ ·H ₂ O	aq	207.391		-55.0			
Ba(CN) ₂ ·2H ₂ O	c	225.406		-124.4			
Ba(CNO) ₂	c	221.374		-195.4			
				-213.			

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BaSi₃
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Table 96(19)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity							
Substance		Barium					S°	C _p °	
Formula and Description	State	Formula Weight	ΔHf°	ΔHf°	ΔGf°	H ₂₉₈ ° - H ₀ °			
				298.15 K (25°C)					
				kcal/mol					
				cal/deg mol					
BaSi ₃	c	221.598		-458.					
Ba ₂ Si ₂	c	330.852		-400.					
BaO•SiO ₂	c	213.424		-388.05	-368.13		26.2	21.51	
	amorp			-376.					
BaO•2SiO ₂	c	273.509		-609.0	-576.2		36.6	32.05	
2BaO•SiO ₂	c	366.764		-546.8	-519.8		42.1	32.24	
2BaO•3SiO ₂	c	486.933		-1000.2	-947.2		61.7	53.68	
BaSiF ₆	c	279.416		-705.6	-667.8		39.		
BaO•GeO ₂	c	257.928		-295.8					
2BaO•GeO ₂	c	411.268		-446.3					
3BaO•GeO ₂	c	564.607		-589.7					

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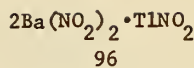
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Table 96(20)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Substance		Barium						
Formula and Description	State	Formula Weight	ΔHf°	ΔHf°	ΔGf°	H ₂₉₈ ° - H ₀ °	S°	C _p °
			0 K	298.15 K (25°C)				
BaI ₂ •2PbI ₂	c	1313.146		-229.3				
BaI ₂ •2PbI ₂ •7H ₂ O	c	1439.254		-716.8				
BaAl ₄	c	245.266		-28.				
BaO•Al ₂ O ₃	c	255.301		-556.				
BaO•Al ₂ O ₃ •H ₂ O	c	273.316		-633.3				
BaO•Al ₂ O ₃ •2H ₂ O	c	291.331		-705.0				
BaO•Al ₂ O ₃ •4H ₂ O	c	327.362		-850.6				
BaO•Al ₂ O ₃ •7H ₂ O	c	381.408		-1061.5				
2BaO•Al ₂ O ₃ •5H ₂ O	c	498.717		-1084.8				
3BaO•Al ₂ O ₃	c	561.979		-841.				
7BaO•6Al ₂ O ₃ •36H ₂ O	c	2333.695		-6058.				
BaCl ₂ •Al ₂ Cl ₆	c	474.927		-547.1				
3BaCl ₂ •2Al ₂ Cl ₆	c	1158.100		-1303.4				

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Table 96(21)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Substance		Barium						
Formula and Description	State	Formula Weight	ΔHf°	ΔGf°	$\text{H}_{298}^\circ - \text{H}_0^\circ$	S°	C_p°	
			kcal/mol		298.15 K (25°C)		cal/deg mol	
$2\text{Ba}(\text{NO}_2)_2 \cdot \text{TlNO}_2$	c	709.078	-381.9					
BaHgBr_4	aq	657.566	-227.5					
$\text{Ba}(\text{HgBr}_3)_2$	aq	1017.974	-266.2					
Ba_2HgBr_6	aq	954.724	-415.7					
$\text{Ba}_4\text{HgBr}_{10}$	aq	1549.040	-788.3					
$\text{BaHg}(\text{CN})_4$	aq	442.001	-0.9					
$\text{Ba}[\text{Hg}(\text{CN})_3]_2$	aq	694.627	+63.8					
$\text{BaCl}_2 \cdot 2\text{Hg}(\text{CN})_2$	aq	713.497	-75.3					
$\text{BaCl}_2 \cdot 2\text{Hg}_2(\text{CN})_2 \cdot 5\text{H}_2\text{O}$	c	803.574	-434.9					
$\text{BaBr}_2 \cdot 2\text{Hg}(\text{CN})_2$	aq	802.409	-54.4					
$\text{BaBr}_2 \cdot 2\text{Hg}(\text{CN})_2 \cdot 7\text{H}_2\text{O}$	c	928.517	-554.2					

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Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Barium

Table 96(22)

Substance	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
Formula and Description	kcal/mol							
cal/deg mol								
2Hg(CN) ₂ ·BaI ₂	aq	896.400		-27.0				
2Hg(CN) ₂ ·BaI ₂ ·6H ₂ O	c	1004.492		-460.0				
BaAg(CN) ₃	aq	323.264		-28.6				
Ba[Ag(CN) ₂] ₂	aq	457.151		+1.9				
Ba[Ni(CN) ₄]	aq	300.121		-34.1				
Ba ₂ Fe(CN) ₆	aq	486.634		-145.7				
Ba ₂ Fe(CN) ₆ ·6H ₂ O	c	594.726		-567.0				
Ba ₃ [Fe(CN) ₆] ₂ std. state, m = 1	aq	835.928		-116.9	-53.5		136.1	
in	aq			-115.03				
1,500 H ₂ O	aq			-115.07				
2,000 H ₂ O	aq			-115.24				
5,000 H ₂ O	aq			-115.35				
7,500 H ₂ O	aq			-115.43				
10,000 H ₂ O	aq							

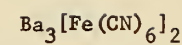
SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

Enthalpy and Gibbs Energy of Formation: Entropy and Heat Capacity

Barium



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Table	96(23)	Substance		Formula Weight	State	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
		Formula and Description	298.15 K (25°C)								
						kcal/mol					
						cal/deg mol					
$\text{Ba}_3[\text{Fe}(\text{CN})_6]_2$	in	20,000 H_2O	aq			0 K					
		50,000 H_2O	aq				-115.66				
		100,000 H_2O	aq				-116.00				
		200,000 H_2O	aq				-116.18				
		500,000 H_2O	aq				-116.31				
		1,000,000 H_2O	aq				-116.51				
		$\infty \text{H}_2\text{O}$	aq								
$\text{Ba}_3[\text{FeCO}(\text{CN})_5]_2$			c	839.914							
$\text{Ba}_3[\text{FeCO}(\text{CN})_5]_2 \cdot 11\text{H}_2\text{O}$			c	1038.082							
$\text{BaH}_2\text{Fe}(\text{CN})_6$			aq	351.310							
BaPdCl_4			c	385.552							
$\text{Ba}_3(\text{RhCl}_6)_2$			c	1043.266							

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

National Bureau of Standards

Washington, D. C.

Table 96 (24)		Barium						
Substance		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Formula and Description	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C_p°
			0 K	298.15 K (25°C)				
		kcal/mol						
		cal/deg mol						
BaRuO ₄ · H ₂ O	c	320.423		-312.7				
BaPtCl ₆	c	545.148		-283.2				
BaPtCl ₆ · 6H ₂ O	aq	653.240		-292.5				
BaOsCl ₆	c	653.240		-703.5				
BaOsCl ₆	c	540.258		-269.6				
BaMnO ₄	c	256.276			-267.5			
Ba (ReO ₄) ₂ · 4H ₂ O	c	709.797		-805.	-697.5		90.	
BaCrO ₄	c	253.334		-345.6	-321.53		37.9	
BaMoO ₃	c	281.278		-295.				
BaMoO ₄	c	297.278		-370.	-344.1		33.	33.6

BaMoO₄
96

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

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BaWO₄
96

Table 96(25)		Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity						
Substance		Barium						
Formula and Description	State	Formula Weight	ΔH_f^0	ΔH_f^0	ΔG_f^0	$H_{298}^0 - H_0^0$	S^0	C_p^0
			0 K	298.15 K (25°C)				
		kcal/mol				cal/deg.mol		
BaWO ₄	c	385.188	-407.					
Ba ₃ WO ₆	c	691.866	-752.					
BaTiO ₃	c	233.238	-396.7		-375.8		25.8	24.49
Ba ₂ TiO ₄	c	386.578	-536.1		-509.8		47.0	36.48
BaZrO ₃	c	276.558	-425.3		-405.0		29.8	24.31
BaHfO ₃	c	363.828	-437.9					
BaMg ₂	c	185.964	-1.5					
BaCa(CO ₃) ₂	c	297.439			-543.2			
barytoalcite	c				-543.0			
BaSrTiO ₄	c	336.858	-544.0		-518.1		45.8	34.95

SELECTED VALUES OF CHEMICAL THERMODYNAMIC PROPERTIES - SERIES I

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National Bureau of Standards

Enthalpy and Gibbs Energy of Formation; Entropy and Heat Capacity

Table 97(1)

Radium

Substance	State	Formula Weight	ΔH_f°	ΔH_f°	ΔG_f°	$H_{298}^\circ - H_0^\circ$	S°	C _p °
Formula and Description			0 K	298.15 K (25°C)				
Ra	c	226.025	0	0	0	17.	4.97	
Ra ⁺	g			38.	31.	42.15		
Ra ²⁺	g			161.22				
Ra	g			396.70				
std. state, m = 1	aq			-126.1	-134.2	13.		
RaO	c	242.0244		-125.				
RaCl ₂	c	296.931		-206.0	-196.9	32.		
std. state, m = 1	aq			-350.	-311.4	40.		
RaCl ₂ ·2H ₂ O	c	332.9617		-245.4	-207.6	51.		
Ra(IO ₃) ₂	c	575.8302		-351.6	-326.4	65.		
RaSO ₄	c	322.0866		-343.4	-312.2	33.		
std. state, m = 1	aq			-237.	-190.3	18.		
Ra(NO ₃) ₂	c	350.0348		-225.2	-187.4	53.		
std. state, m = 1	aq					83.		

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