

Gallium (Ga)

$A_r = 69.72$ Gallium (Ga)

Ga (ref)

0 to 302.920 K crystal
 302.92 to 2476.570 K liquid
 above 2476.570 K ideal monatomic gas

Refer to the individual tables for details.

T/K	C_p^o	Enthalpy Reference Temperature = $T_r = 298.15$ K		Standard State Pressure = $p^o = 0.1$ MPa		ΔG^o	log K_r
		S^o - $[C_p^o - (T_r/T)]/T$	$H^o - H^o(T_r)$	S^o - $[C_p^o - (T_r/T)]/T$	ΔH^o		
0	0	0	0	INFINITE	0	0	0
100	18.484	16.085	62.261	0	-5.561	0	0
200	23.815	30.912	43.151	0	-4.618	0	0
298.15	26.064	40.838	0	0	-2.448	0	0
300	26.091	41.000	40.839	0	0.048	0	0
302.920	26.133	41.253	40.842	0	0.124	0	0
302.920	28.493	59.706	40.842	0	3.714	CRYSTAL \leftrightarrow LIQUID	TRANSITION
400	27.154	67.426	46.408	0	8.407	0	0
500	26.819	73.445	51.237	0	11.104	0	0
600	26.694	78.324	55.358	0	13.780	0	0
700	26.568	82.428	58.939	0	16.442	0	0
800	26.568	85.976	62.102	0	19.099	0	0
900	26.568	89.105	64.932	0	21.756	0	0
1000	26.568	91.904	67.492	0	24.413	0	0
1100	26.568	94.436	69.828	0	27.069	0	0
1200	26.568	96.748	71.976	0	29.726	0	0
1300	26.568	98.875	73.965	0	32.383	0	0
1400	26.568	100.844	75.815	0	35.040	0	0
1500	26.568	102.677	77.546	0	37.697	0	0
1600	26.568	104.391	79.171	0	40.354	0	0
1700	26.568	106.092	80.700	0	43.010	0	0
1800	26.568	107.777	82.150	0	45.667	0	0
1900	26.568	109.457	83.524	0	48.324	0	0
2000	26.568	110.320	84.830	0	50.981	0	0
2100	26.568	111.616	86.075	0	53.638	0	0
2200	26.568	112.832	87.264	0	56.295	0	0
2300	26.568	114.033	88.402	0	58.951	0	0
2400	26.568	115.164	89.494	0	61.608	0	0
2476.570	26.568	115.998	90.301	0	63.643	LIQUID \leftrightarrow IDEAL GAS	FUGACITY = 1 bar
2476.570	21.260	220.462	90.301	0	322.353	0	0
2500	21.251	220.662	91.522	0	322.851	0	0
2600	21.215	221.495	96.505	0	324.974	0	0
2700	21.184	222.295	101.189	0	327.094	0	0
2800	21.156	223.065	105.489	0	329.211	0	0
2900	21.131	223.807	109.556	0	331.325	0	0
3000	21.109	224.523	113.377	0	333.437	0	0
3100	21.090	225.214	116.973	0	335.547	0	0
3200	21.073	225.884	120.366	0	337.655	0	0
3300	21.059	226.532	123.574	0	339.762	0	0
3400	21.047	227.160	126.611	0	341.867	0	0
3500	21.036	227.770	129.493	0	343.971	0	0
3600	21.029	228.363	132.231	0	346.075	0	0
3700	21.023	228.939	134.837	0	348.177	0	0
3800	21.020	229.500	137.321	0	350.279	0	0
3900	21.018	230.045	139.691	0	352.381	0	0
4000	21.020	230.578	141.957	0	354.483	0	0
4100	21.024	231.097	144.125	0	356.585	0	0
4200	21.030	231.603	146.202	0	358.688	0	0
4300	21.040	232.098	148.193	0	360.791	0	0
4400	21.052	232.582	150.106	0	362.896	0	0
4500	21.068	233.055	151.944	0	365.002	0	0
4600	21.087	233.519	153.712	0	367.110	0	0
4700	21.109	233.972	155.415	0	369.219	0	0
4800	21.135	234.417	157.056	0	371.332	0	0
4900	21.165	234.853	158.640	0	373.446	0	0
5000	21.198	235.281	160.168	0	375.565	0	0
5100	21.236	235.701	161.645	0	377.686	0	0
5200	21.278	236.114	163.073	0	379.812	0	0
5300	21.324	236.520	164.455	0	381.942	0	0
5400	21.375	236.919	165.793	0	384.077	0	0
5500	21.431	237.312	167.090	0	386.217	0	0
5600	21.491	237.698	168.348	0	388.363	0	0
5700	21.556	238.079	169.568	0	390.515	0	0
5800	21.625	238.455	170.752	0	392.674	0	0
5900	21.700	238.825	171.903	0	394.841	0	0
6000	21.780	239.190	173.021	0	397.015	0	0

PREVIOUS:

CURRENT: June 1983 (1 bar)

Gallium (Ga)

Ga (ref)

CRYSTAL

Gallium (Ga)

Gallium (Ga)

Ga₁(cr)

$\Delta_f H^\circ(298.15 \text{ K}) = 40.838 \pm 0.20 \text{ J}\cdot\text{K}^{-1}\cdot\text{mol}^{-1}$
 $\Delta_f H^\circ(0 \text{ K}) = 0 \text{ kJ}\cdot\text{mol}^{-1}$
 $\Delta_{\text{fus}} H^\circ = 5.59 \pm 0.04 \text{ kJ}\cdot\text{mol}^{-1}$
 $T_{\text{fus}} = 302.92 \pm 0.02 \text{ K}$

Enthalpy of Formation

Zero by definition.

Heat Capacity and Entropy

The adopted thermal functions are those recommended by Hultgren *et al.*¹ The heat capacity values were based on the studies of Phillips (0.1–4.2 K),² Seidel and Kiesom (0.35–4.38 K),³ Eichenauer and Schulze (5–22 K),⁴ Adams *et al.* (15–323 K),⁵ and Clusius and Hartick (15–200 K).⁶ Refer to Hultgren *et al.*¹ for details.

Enthalpy of Fusion

Refer to the liquid table for details.

Enthalpy of Sublimation

The enthalpy of sublimation is the enthalpy of formation of the gas, $\Delta_{\text{sub}} H^\circ(298.15 \text{ K}) = \Delta_f H^\circ(\text{g}, 298.15 \text{ K})$.

References

- ¹R. Hultgren, P. D. Desai, *et al.*, "Selected Values of the Thermodynamic Properties of the Elements," American Society for Metals, Metals Park, Ohio (1973).
- ²N. E. Phillips, *Phys. Rev.* **134**, A385 (1964).
- ³G. Seidel and P. H. Kiesom, *Phys. Rev.* **112**, 1083 (1958).
- ⁴W. Eichenauer and M. Schulze, *Z. Naturforsch.* **14a**, 962 (1959).
- ⁵G. B. Adams, H. L. Johnston and E. C. Kerr, *J. Am. Chem. Soc.* **74**, 4784 (1952).
- ⁶K. Clusius and P. Hartick, *Z. Physik. Chem.* **134**, 243 9 (1928).

T/K	Enthalpy Reference Temperature = T _r = 298.15 K		Standard State Pressure = p° = 0.1 MPa	
	C _p ^a	S° - [G° - H°(T _r)]/T	H° - H°(T _r)	Δ _r G°
0	0	INFINITE	0	0
100	18.484	16.085	-5.561	0
200	23.815	30.912	-4.618	0
250	24.907	36.558	-2.948	0
298.15	26.064	40.838	-1.251	0
300	26.091	40.839	0	0
302.920	26.133	41.253	0.048	0
350	27.004	45.091	0.124	CRYSTAL <--> LIQUID ---
400	27.460	48.726	1.376	0.875
450	27.916	51.987	4.122	1.810
500	28.372	54.952	5.529	2.743
				3.672

PREVIOUS:

CURRENT: June 1983

Gallium (Ga)

Ga₁(cr)

Ga₁(l)

A_r = 69.72 Gallium (Ga)

LIQUID

Gallium (Ga)

T/K	C _p ^o	S ^o - (G ^o - H ^o (T))/T	H ^o - H ^o (T)	Δ _{sub} H ^o	Δ _{sub} G ^o	log K _r
0						
100						
200						
250						
298.15	28.559	59.253	0.	5.578	0.088	-0.015
300	28.533	59.253	0.033	5.583	0.054	-0.009
302.920	28.493	59.256	0.126	---	CRYSTAL < -> LIQUID	---
350	27.719	59.598	1.458	0.	0.	0.
400	27.154	60.335	2.829	0.	0.	0.
450	26.981	61.320	4.182	0.	0.	0.
500	26.819	62.393	5.526	0.	0.	0.
600	26.694	64.655	8.202	0.	0.	0.
700	26.568	66.908	10.864	0.	0.	0.
800	26.568	69.075	13.521	0.	0.	0.
900	26.568	71.130	16.177	0.	0.	0.
1000	26.568	73.070	18.834	0.	0.	0.
1100	26.568	74.899	21.491	0.	0.	0.
1200	26.568	76.625	24.148	0.	0.	0.
1300	26.568	78.256	26.805	0.	0.	0.
1400	26.568	79.800	29.462	0.	0.	0.
1500	26.568	81.264	32.118	0.	0.	0.
1600	26.568	82.657	34.775	0.	0.	0.
1700	26.568	83.983	37.432	0.	0.	0.
1800	26.568	85.251	40.089	0.	0.	0.
1900	26.568	86.459	42.746	0.	0.	0.
2000	26.568	87.619	45.403	0.	0.	0.
2100	26.568	88.731	48.060	0.	0.	0.
2200	26.568	89.799	50.716	0.	0.	0.
2300	26.568	90.828	53.373	0.	0.	0.
2400	26.568	91.818	56.030	0.	0.	0.
2476.570	26.568	92.553	58.064	---	FUGACITY = 1 bar	---
2500	26.568	92.774	58.687	-258.586	2.447	-0.051
2600	26.568	93.697	61.344	-258.052	12.878	-0.259
2700	26.568	94.589	64.001	-257.515	23.288	-0.451
2800	26.568	95.453	66.657	-256.975	33.679	-0.628
2900	26.568	96.290	69.314	-256.433	44.050	-0.793
3000	26.568	97.102	71.971	-255.888	54.402	-0.947

$S^{\circ}(298.15\text{ K}) = [59.253] \text{ J}\cdot\text{K}^{-1}\cdot\text{mol}^{-1}$
 $T = 302.92 \pm 0.02\text{ K}$
 $\Delta_{\text{sub}}H^{\circ}(298.15\text{ K}) = [5.578] \text{ kJ}\cdot\text{mol}^{-1}$
 $\Delta_{\text{sub}}H^{\circ} = 5.59 \pm 0.04 \text{ kJ}\cdot\text{mol}^{-1}$

Enthalpy of Formation
 The enthalpy of formation, $\Delta_f H^{\circ}$ (l, 298.15 K), is calculated from that of the crystal by adding the enthalpy of fusion and the difference in enthalpy, $H^{\circ}(302.92\text{ K}) - H^{\circ}(298.15\text{ K})$, between the crystal and liquid.

Heat Capacity and Entropy
 The heat capacity of the liquid is based on the recommendations of Hultgren *et al.*¹ These values are based on the enthalpy data of Kochehova and Rezukhina (315–712 K),² and the heat capacity data of Chen and Turnbull (245–611 K).³

Enthalpy of Fusion
 The adopted value for the enthalpy of fusion is derived from the direct calorimetric measurement of Roth *et al.*⁴
 The temperature of fusion, $T_{\text{fus}} = 302.92 \pm 0.2\text{ K}$ (corrected to IPTS–68), was derived from the study of Boedtker *et al.*⁵

Enthalpy of Vaporization
 The temperature at which $\Delta_{\text{sub}}G^{\circ} = 0$ (1 bar) for the process $\text{Ga(l)} \rightarrow \text{Ga(g)}$ is the adopted T_{vap} (1 bar). The difference in the enthalpies of formation of the ideal gas and liquid at T_{vap} is $\Delta_{\text{sub}}H^{\circ}$. Refer to the ideal gas table for additional information.

- References**
¹R. Hultgren, P. D. Desai, *et al.*, "Selected Values of the Thermodynamic Properties of the Elements," American Society for Metals, Metals Parks, Ohio (1973).
²H. M. Kochehova and T. N. Rezukhina, Proc. 4th All-Union Conf. on Semiconductor Materials (Eng. transl.), 26 (1961).
³H. S. Chen and S. Turnbull, Acta Met. 16, 369 (1968).
⁴W. A. Roth, I. Meyer and H. Zeumar, Z. Anorg. Chem. 214, 309 (1933); Z. Anorg. Chem. 216, 303 (1934).
⁵O. A. Boedtker, R. C. LaForce, *et al.*, Trans. Faraday Soc. 61, 665 (1965).

PREVIOUS:

CURRENT: June 1983

Ga₁(l)

Gallium (Ga)

CRYSTAL-LIQUID

0 to 302.92 K crystal
above 302.92 K liquid

Refer to the individual tables for details.

Gallium (Ga)

A₁ = 69.72 Gallium (Ga)

Ga₁(cr,l)

T/K	C _p ^a	Enthalpy Reference Temperature = T _r = 298.15 K		Standard State Pressure = p° = 0.1 MPa		log K _f
		S° - (G° - H°(T _r))/T	H° - H°(T _r)	ΔH°	ΔG°	
		J·K ⁻¹ ·mol ⁻¹	kJ·mol ⁻¹	kJ·mol ⁻¹	kJ·mol ⁻¹	
0	0.	INFINITE	-5.561	0.	0.	0.
100	18.484	16.085	-4.618	0.	0.	0.
200	23.815	30.912	-2.448	0.	0.	0.
250	24.907	36.338	-1.231	0.	0.	0.
298.15	26.064	40.838	0.	0.	0.	0.
300	26.091	41.000	0.048	0.	0.	0.
302.920	26.133	41.253	0.174	0.	0.	0.
302.920	28.493	59.706	5.714	CRYSTAL → LIQUID	CRYSTAL → LIQUID	TRANSITION
350	27.719	63.764	7.037	0.	0.	0.
400	27.154	67.476	8.407	0.	0.	0.
450	26.961	70.613	9.760	0.	0.	0.
500	26.819	73.445	11.104	0.	0.	0.
600	26.694	78.324	13.780	0.	0.	0.
700	26.568	82.428	16.442	0.	0.	0.
800	26.568	85.976	19.099	0.	0.	0.
900	26.568	89.105	21.756	0.	0.	0.
1000	26.568	91.904	24.413	0.	0.	0.
1100	26.568	94.436	27.069	0.	0.	0.
1200	26.568	96.748	29.726	0.	0.	0.
1300	26.568	98.875	32.383	0.	0.	0.
1400	26.568	100.844	35.040	0.	0.	0.
1500	26.568	102.677	37.697	0.	0.	0.
1600	26.568	104.391	40.354	0.	0.	0.
1700	26.568	106.002	43.010	0.	0.	0.
1800	26.568	107.521	45.667	0.	0.	0.
1900	26.568	108.957	48.324	0.	0.	0.
2000	26.568	110.320	50.981	0.	0.	0.
2100	26.568	111.616	53.638	0.	0.	0.
2200	26.568	112.852	56.295	0.	0.	0.
2300	26.568	114.033	58.951	0.	0.	0.
2400	26.568	115.164	61.608	0.	0.	0.
2476.570	26.568	115.998	63.643	---	---	---
				---	---	---
2500	26.568	116.249	64.265	-258.586	2.447	-0.051
2600	26.568	117.291	66.922	-258.052	12.878	-0.259
2700	26.568	118.293	69.579	-257.515	23.288	-0.451
2800	26.568	119.260	72.236	-256.975	33.679	-0.628
2900	26.568	120.192	74.892	-256.433	44.050	-0.793
3000	26.568	121.093	77.549	-255.888	54.402	-0.947

PREVIOUS:

CURRENT: June 1983

Gallium (Ga)

Ga₁(cr,l)

Gallium (Ga)

IP(Ga, g) = 48380 ± 5 cm⁻¹
 S°(298.15 K) = 169.042 ± 0.01 J·K⁻¹·mol⁻¹

IDEAL GAS

A_r = 51.996 Gallium (Ga)

Δ_fH°(0 K) = 270.97 ± 2.1 kJ·mol⁻¹
 Δ_fH°(298.15 K) = 271.96 ± 2.1 kJ·mol⁻¹

Electronic Levels and Quantum Weights	g _i
State	
¹ P _{1/2}	0
³ P ₂	2
³ P _{1/2}	4
³ P ₀	2
¹ P _{1/2}	2
¹ P _{3/2}	4

Enthalpy of Formation

The enthalpy of formation of Ga(g), as recommended by Huligren *et al.*,¹ was derived from the third law analysis of seven vaporization studies and a spectroscopic calculation. The studies are: Matern *et al.* (1081–1213 K),² Macur *et al.* (1196–1473 K),³ Alcock *et al.* (1173–1437 K),⁴ Munir and Searcy (1174–1603 K),⁵ Cochran and Foster (1179–1383 K),⁶ Speiser and Johnston (1230–1518 K),⁷ Hartick (1198–1391 K),⁸ and the spectroscopic calculation by Gurvich.⁹

Heat Capacity and Entropy

The thermal functions were calculated using the atomic levels as given by Moore.¹⁰ The resulting thermal functions are those recommended by Huligren *et al.*¹

References

- ¹R. Huligren, P. D. Desai, *et al.*, "Selected Values of the Thermodynamic Properties of the Elements," American Society for Metals, Metals Park, Ohio, (1973).
- ²G. Matern, Yu. A. Sapozhnikov, *et al.*, *Zvest. Akad. Nauk SSSR, Metall.*, (3) 210 (1969).
- ³G. J. Macur, R. K. Edwards, and P. G. Wahibick, *J. Phys. Chem.*, 70, 2956 (1966).
- ⁴C. B. Alcock, J. B. Cornish, and P. Grievson, "Thermodynamics," Proc. Symp. Vienna, Vol. 1, 211–30, 367–8 (1965).
- ⁵Z. A. Munir and A. W. Searcy, *J. Electrochem. Soc.*, 111, 1170 (1964).
- ⁶C. N. Cochran and L. M. Foster, *J. Electrochem. Soc.*, 109, 144 (1962).
- ⁷R. Speiser and H. L. Johnston, *J. Am. Chem. Soc.*, 75, 1469 (1953).
- ⁸P. Hartick, *Z. Physik. Chem.*, 134, 1 (1960).
- ⁹L. V. Gurvich, *Zh. Fiz. Khim.*, 34, 1690 (1960).
- ¹⁰C. E. Moore, *U. S. Nat. Bur. Stand., NSRDS-NBS-35*, Vol. II, 1971 [Reprint of NBS Circ. 497, (1952)].

Ga₁(g)

T/K	C _p ^o	S° - (G° - H°(T _r))/T	H° - H°(T _r)/T	Δ _f H°	ΔG°	log K _r
		J·K ⁻¹ ·mol ⁻¹		kJ·mol ⁻¹		
		Standard State Pressure = p° = 0.1 MPa				
0	0	INFINITE	INFINITE	270.970	270.970	INFINITE
100	20.802	144.848	189.570	259.229	259.229	-135.407
200	22.311	159.555	171.266	246.337	246.337	-64.337
300	23.914	164.703	169.452	239.913	239.913	-50.177
298.15	23.914	164.703	169.452	239.913	239.913	-50.177
350	25.304	169.042	169.042	233.736	233.736	-40.950
400	26.430	170.189	169.043	227.999	227.999	-30.656
450	27.174	170.957	170.957	222.966	222.966	-20.072
500	27.691	171.300	171.300	218.523	218.523	-9.085
600	28.589	172.809	172.809	211.030	211.030	75.192
700	29.489	174.247	174.247	206.245	206.245	122.102
800	30.389	175.616	175.616	202.626	202.626	174.466
900	31.289	176.923	176.923	200.000	200.000	226.626
1000	32.189	178.171	178.171	198.000	198.000	278.626
1100	33.089	179.369	179.369	196.500	196.500	330.626
1200	33.989	180.517	180.517	195.200	195.200	382.626
1300	34.889	181.615	181.615	194.000	194.000	434.626
1400	35.789	182.663	182.663	192.900	192.900	486.626
1500	36.689	183.661	183.661	191.900	191.900	538.626
1600	37.589	184.609	184.609	191.000	191.000	590.626
1700	38.489	185.507	185.507	190.200	190.200	642.626
1800	39.389	186.355	186.355	189.500	189.500	694.626
1900	40.289	187.153	187.153	188.900	188.900	746.626
2000	41.189	187.901	187.901	188.400	188.400	798.626
2100	42.089	188.600	188.600	188.000	188.000	850.626
2200	42.989	189.259	189.259	187.700	187.700	902.626
2300	43.889	189.878	189.878	187.500	187.500	954.626
2400	44.789	190.457	190.457	187.400	187.400	1006.626
2476.570	45.689	191.000	191.000	187.400	187.400	1058.626
2500	46.589	191.517	191.517	187.400	187.400	1110.626
2600	47.489	192.000	192.000	187.400	187.400	1162.626
2700	48.389	192.449	192.449	187.400	187.400	1214.626
2800	49.289	192.857	192.857	187.400	187.400	1266.626
2900	50.189	193.224	193.224	187.400	187.400	1318.626
3000	51.089	193.550	193.550	187.400	187.400	1370.626
3100	51.989	193.835	193.835	187.400	187.400	1422.626
3200	52.889	194.080	194.080	187.400	187.400	1474.626
3300	53.789	194.285	194.285	187.400	187.400	1526.626
3400	54.689	194.450	194.450	187.400	187.400	1578.626
3500	55.589	194.575	194.575	187.400	187.400	1630.626
3600	56.489	194.660	194.660	187.400	187.400	1682.626
3700	57.389	194.705	194.705	187.400	187.400	1734.626
3800	58.289	194.710	194.710	187.400	187.400	1786.626
3900	59.189	194.675	194.675	187.400	187.400	1838.626
4000	60.089	194.600	194.600	187.400	187.400	1890.626
4100	60.989	194.485	194.485	187.400	187.400	1942.626
4200	61.889	194.330	194.330	187.400	187.400	1994.626
4300	62.789	194.135	194.135	187.400	187.400	2046.626
4400	63.689	193.900	193.900	187.400	187.400	2098.626
4500	64.589	193.625	193.625	187.400	187.400	2150.626
4600	65.489	193.310	193.310	187.400	187.400	2202.626
4700	66.389	192.955	192.955	187.400	187.400	2254.626
4800	67.289	192.560	192.560	187.400	187.400	2306.626
4900	68.189	192.125	192.125	187.400	187.400	2358.626
5000	69.089	191.650	191.650	187.400	187.400	2410.626
5100	69.989	191.135	191.135	187.400	187.400	2462.626
5200	70.889	190.580	190.580	187.400	187.400	2514.626
5300	71.789	189.985	189.985	187.400	187.400	2566.626
5400	72.689	189.350	189.350	187.400	187.400	2618.626
5500	73.589	188.675	188.675	187.400	187.400	2670.626
5600	74.489	187.960	187.960	187.400	187.400	2722.626
5700	75.389	187.205	187.205	187.400	187.400	2774.626
5800	76.289	186.410	186.410	187.400	187.400	2826.626
5900	77.189	185.575	185.575	187.400	187.400	2878.626
6000	78.089	184.700	184.700	187.400	187.400	2930.626

PREVIOUS:

CURRENT: June 1983 (1 bar)

Gallium (Ga)

Ga₁(g)

IDEAL GAS

Gallium, Ion (Ga⁺)

IP(Ga⁺, g) = 165458 ± 5 cm⁻¹
 $\Delta_f H^\circ(0 \text{ K}) = 849.8 \pm 4.0 \text{ kJ}\cdot\text{mol}^{-1}$
 $\Delta_f H^\circ(298.15 \text{ K}) = 161.790 \pm 0.04 \text{ kJ}\cdot\text{mol}^{-1}$

Electronic Levels and Quantum Weights	g _i
State	
¹ S ₀	0
³ P ₀	47370
³ P ₁	47816
³ P ₂	48750

Enthalpy of Formation

$\Delta_f H^\circ(\text{Ga}^+, \text{g}, 0 \text{ K})$ is calculated from $\Delta_f H^\circ(\text{Ga}, \text{g}, 0 \text{ K})$ using the spectroscopic value of IP(Ga) = 48387.63 ± 0.1 cm⁻¹ (578.837 ± 0.001 kJ·mol⁻¹) from Moore.² The ionization limit is converted from cm⁻¹ to kJ·mol⁻¹ using the factor, 1 cm⁻¹ = 0.01196266 kJ·mol⁻¹, which is derived from the 1973 CODATA fundamental constants.³ Rosenstock *et al.*⁴ and Levin and Lias⁵ have summarized additional ionization and appearance potential data.

$\Delta_f H^\circ(\text{Ga}^+, \text{g}, 298.15 \text{ K})$ is calculated from $\Delta_f H^\circ(\text{Ga}, \text{g}, 0 \text{ K})$ by using IP(Ga) with JANAF¹ enthalpies, $H^\circ(0 \text{ K}) - H^\circ(298.15 \text{ K})$, for Ga(g), Ga⁺(g), and e⁻(ref). $\Delta_f H^\circ(\text{Ga} \rightarrow \text{Ga}^+ + \text{e}^-, 298.15 \text{ K})$ differs from a room temperature threshold energy due to inclusion of these enthalpies and to threshold effects discussed by Rosenstock *et al.*⁴ $\Delta_f H^\circ(298.15 \text{ K})$ should be changed by -6.197 kJ mol⁻¹, if it is to be used in the ion convention that excludes the enthalpy of the electron.

Heat Capacity and Entropy

The information on electronic energy levels and quantum weights, given by Moore,^{2,6} is incomplete because many theoretically predicted levels have not been observed. Our calculations indicate that any reasonable method of filling in these missing levels and cutting off the summation in the partition function⁷ has no effect on the thermodynamic functions to 6000 K. This is a result of the high energy of all levels other than the ground state and the first three excited states; the next excited state is 70700 cm⁻¹ above the ground state. Since inclusion of these excited states has no effect on the thermodynamic functions (to 6000 K), we list only the ground state and the first three excited states. The position of the ³P₀ levels are uncertain by 249 cm⁻¹. The reported uncertainty in $S^\circ(298.15 \text{ K})$ is due to uncertainties in the relative ionic mass and the fundamental constants. Extension of these calculations above 6000 K may require consideration of the excited states and use of different fill and cutoff procedures.⁷

References

- ¹JANAF Thermochemical Tables: Ga(g); e⁻(ref), 3-31-82.
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Gallium, Ion (Ga⁺)Ga⁺(g)

T/K	C _p ^a	Enthalpy Reference Temperature = T _r = 298.15 K		H ^o - H ^o (T _r)/T	Standard State Pressure = p ^o = 0.1 MPa		log K _r
		S ^o	-[G ^o - H ^o (T _r)]/T		Δ _r H ^o	Δ _r G ^o	
0	0	0	INFINITE	-6.197	849.815		
100	20.786	139.083	180.270	-4.119			
200	20.786	153.491	163.691	-2.040			
250	20.786	158.129	162.132	-1.001			
298.15	20.786	161.790	161.790	0	856.648		814.332
300	20.786	161.919	161.790	0.038	856.677		814.069
350	20.786	165.123	162.044	1.078	851.767		807.783
400	20.786	167.898	162.606	2.117	852.475		801.451
450	20.786	170.347	163.333	3.156	853.201		795.285
500	20.786	172.537	164.145	4.196	853.935		788.577
600	20.786	176.326	165.869	6.274	855.417		775.307
700	20.786	179.351	167.598	8.353	856.912		761.849
800	20.786	182.506	169.267	10.431	858.415		748.152
900	20.786	184.754	170.834	12.510	859.913		734.280
1000	20.786	186.944	172.356	14.589	861.413		720.240
1100	20.786	188.926	173.774	16.667	862.914		706.050
1200	20.786	190.734	175.113	18.746	864.414		691.723
1300	20.786	192.398	176.379	20.824	865.914		677.272
1400	20.786	193.938	177.579	22.903	867.415		662.705
1500	20.786	195.372	178.718	24.982	868.915		648.030
1600	20.786	196.714	179.801	27.060	870.415		633.255
1700	20.786	197.974	180.834	29.139	871.916		618.387
1800	20.786	199.162	181.819	31.217	873.416		603.430
1900	20.786	200.286	182.762	33.295	874.917		588.390
2000	20.786	201.352	183.665	35.373	876.417		573.271
2100	20.786	202.366	184.532	37.453	877.917		558.077
2200	20.786	203.333	185.364	39.532	879.418		542.811
2300	20.786	204.257	186.166	41.610	880.918		527.477
2400	20.786	205.142	186.938	43.689	882.418		512.078
2500	20.786	205.991	187.683	45.768	883.915		496.664
2600	20.786	206.806	188.403	47.846	885.415		481.233
2700	20.786	207.590	189.103	49.925	886.914		465.784
2800	20.786	208.346	189.773	52.004	888.414		450.329
2900	20.786	209.076	190.427	54.082	889.914		434.868
3000	20.786	209.780	191.060	56.161	891.415		419.402
3100	20.786	210.462	191.675	58.239	892.916		403.937
3200	20.786	211.122	192.272	60.318	894.416		388.472
3300	20.786	211.761	192.853	62.397	895.916		373.007
3400	20.786	212.382	193.419	64.475	897.416		357.542
3500	20.786	212.984	193.969	66.554	898.916		342.077
3600	20.786	213.570	194.505	68.632	899.916		326.612
3700	20.786	214.140	195.028	70.711	900.916		311.147
3800	20.786	214.694	195.539	72.790	901.916		295.682
3900	20.786	215.234	196.037	74.868	902.916		280.217
4000	20.787	215.760	196.523	76.947	903.916		264.752
4100	20.787	216.273	196.999	79.026	904.916		249.287
4200	20.787	216.774	197.464	81.104	905.916		233.822
4300	20.788	217.263	197.919	83.183	906.916		218.357
4400	20.789	217.741	198.364	85.262	907.916		202.892
4500	20.790	218.209	198.799	87.341	908.916		187.427
4600	20.791	218.665	199.226	89.420	909.916		171.962
4700	20.792	219.113	199.645	91.499	910.916		156.497
4800	20.794	219.550	200.055	93.578	911.916		141.032
4900	20.797	219.979	200.457	95.658	912.916		125.567
5000	20.799	220.399	200.852	97.738	913.916		110.102
5100	20.803	220.811	201.239	99.818	914.916		94.637
5200	20.807	221.215	201.619	101.898	915.916		79.172
5300	20.812	221.612	201.993	103.979	916.916		63.707
5400	20.818	222.001	202.360	106.061	917.916		48.242
5500	20.825	222.383	202.720	108.143	918.916		32.777
5600	20.833	222.758	203.075	110.226	919.916		17.312
5700	20.843	223.127	203.423	112.310	920.916		1.847
5800	20.854	223.489	203.766	114.394	921.916		-13.118
5900	20.866	223.846	204.104	116.480	922.916		-28.193
6000	20.880	224.197	204.436	118.568	923.916		-43.268

PREVIOUS:

CURRENT: June 1983 (1 bar)

Gallium, Ion (Ga⁺)Ga⁺(g)

Gallium, Ion (Ga⁺)

EA(Ga, g) = 0.30 ± 0.15 eV
 S°(298.15 K) = [175.801 ± 0.1] J·K⁻¹·mol⁻¹

IDEAL GAS

Electronic Levels and Quantum State	ε _l , cm ⁻¹	Weights g _l
² P ₀	0	1
² P ₁	220	3
² P ₂	580	5

Enthalpy of Formation

Δ_fH°(Ga⁺, g, 0 K) is calculated from Δ_fH°(Ga, g, 0 K)¹ using the adopted electron affinity of EA(Ga) = 0.30 ± 0.15 eV (28.945 ± 14.473 kJ·mol⁻¹). This value, recommended by Hotop and Lineberger,² is based on a photodetachment threshold (using conventional light sources) and semiempirical extrapolation.^{3, 4, 5} Additional information on Ga⁺(g) may be obtained in the critical discussions of Hotop and Lineberger, Rosenstock *et al.*,⁶ and Massey.⁷

Δ_fH°(Ga⁺, g, 298.15 K) is obtained from Δ_fH°(Ga, g, 0 K) by using EA(Ga) with JANAF¹ enthalpies, H°(0 K) - H°(298.15 K), for Ga⁺(g), Ga(g), and e⁻(ref). Δ_fH°(Ga⁺ → Ga + e⁻, 298.15 K) differs from a room-temperature threshold energy due to inclusion of these enthalpies and to threshold effects discussed by Rosenstock *et al.*,⁶ Δ_fH°(298.15 K) should be changed by +6.197 kJ·mol⁻¹ if it is to be used in the ion convention that excludes the enthalpy of the electron.

Heat Capacity and Entropy

The ground state electronic configuration for Ga⁺(g) is given by Hotop and Lineberger,² Rosenstock *et al.*,⁶ and Massey.⁷ The fine-structure separation has been calculated via an isoelectronic extrapolation of ratios of fine structure separations, and a quadratic isoelectronic extrapolation^{4, 5} and is that recommended by Hotop and Lineberger.²

A comparison of the isoelectronic sequences - Ga⁺(g), Ge⁺(g), As⁺(g), Se⁺(g) - would suggest that a state, ¹D₂, may exist at low wave numbers. However, this is most likely a metastable state lying at an energy greater than the electron affinity, as in the case of Al⁺(g). As discussed by Hotop and Lineberger,^{2, 5} Rosenstock *et al.*,⁶ and Massey,⁷ no stable excited states have been observed. Thus, we assume no stable excited states exist.

References

- JANAF Thermochemical Tables: Ga(g); e⁻(ref), 3-31-82.
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Gallium, Ion (Ga⁺)

Ga⁺(g)

T/K	Enthalpy Reference Temperature = T _r = 298.15 K		Standard State Pressure = P° = 0.1 MPa		log K _r
	C _p ^o	S° - (G° - H°(T _r))/T	H° - H°(T _r)	Δ _f H°	
0	0	0	INFINITE	242.025	
100	29.645	143.110	-5.865		
200	30.076	164.238	-5.884		
250	28.867	170.824	-2.834		
298.15	27.618	175.801	-1.360		
300	27.572	175.972	0	238.531	204.547
350	26.407	180.132	0.051	238.496	204.336
400	25.437	183.593	1.400	231.817	199.788
450	24.656	186.542	2.695	230.702	195.070
500	24.034	189.107	3.947	229.562	190.684
600	23.155	190.872	5.163	228.395	186.427
700	22.612	196.926	7.519	225.996	178.257
800	22.287	200.976	9.805	223.542	170.404
900	22.106	202.539	12.048	221.050	163.086
1000	22.021	204.863	14.268	218.533	155.991
1100	21.998	206.961	16.473	216.003	149.177
1200	22.012	208.875	18.674	213.469	142.617
1300	22.049	210.638	20.874	210.933	136.288
1400	22.097	212.274	23.077	208.401	130.171
1500	22.149	213.801	25.284	205.873	124.248
1600	22.199	215.232	27.497	203.350	118.505
1700	22.246	216.579	29.714	200.832	112.931
1800	22.287	217.852	31.936	198.318	107.515
1900	22.323	219.058	34.163	195.810	102.246
2000	22.353	220.203	36.394	193.305	97.116
2100	22.377	221.295	38.627	190.803	92.118
2200	22.395	222.336	40.864	188.304	87.245
2300	22.408	223.332	43.103	185.807	82.491
2400	22.417	224.286	45.348	183.316	77.851
2500	22.422	225.201	47.594	180.818	73.319
2600	22.423	226.080	49.828	178.321	68.826
2700	22.422	226.927	52.054	175.824	64.381
2800	22.417	227.742	54.311	173.327	59.936
2900	22.410	228.528	56.553	170.830	55.491
3000	22.402	229.288	58.794	168.333	51.046
3100	22.391	230.022	61.035	165.836	46.601
3200	22.379	230.733	63.274	163.339	42.156
3300	22.366	231.422	65.513	160.842	37.711
3400	22.351	232.089	67.750	158.345	33.266
3500	22.336	232.737	69.986	155.848	28.821
3600	22.320	233.366	72.220	153.351	24.376
3700	22.303	233.977	74.453	150.854	19.931
3800	22.286	234.572	76.684	148.357	15.486
3900	22.268	235.150	78.914	145.860	11.041
4000	22.250	235.714	81.141	143.363	6.596
4100	22.232	236.263	83.367	140.866	2.151
4200	22.214	236.799	85.591	138.369	-2.294
4300	22.195	237.321	87.814	135.872	-6.839
4400	22.176	237.831	90.034	133.375	-11.384
4500	22.158	238.329	92.253	130.878	-15.929
4600	22.139	238.816	94.469	128.381	-20.474
4700	22.120	239.292	96.684	125.884	-25.019
4800	22.102	239.757	98.897	123.387	-29.564
4900	22.083	240.213	101.108	120.890	-34.109
5000	22.065	240.659	103.318	118.393	-38.654
5100	22.047	241.096	105.525	115.896	-43.199
5200	22.029	241.524	107.731	113.399	-47.744
5300	22.011	241.943	109.934	110.902	-52.289
5400	21.993	242.354	112.136	108.405	-56.834
5500	21.976	242.758	114.336	105.908	-61.379
5600	21.958	243.154	116.535	103.411	-65.924
5700	21.941	243.542	118.732	100.914	-70.469
5800	21.924	243.924	120.927	98.417	-75.014
5900	21.908	244.298	123.120	95.920	-79.559
6000	21.891	244.666	125.311	93.423	-84.104
			223.416	90.926	-88.649

PREVIOUS.

CURRENT: June 1983 (1 bar)

Gallium, Ion (Ga⁺)

Ga⁺(g)