

W<sub>1</sub>(cr)

Tungsten (W)

CRYSTAL

Tungsten (W)

T/K	C <sub>p</sub> <sup>a</sup>	Enthalpy Reference Temperature = T <sub>r</sub> = 298.15 K		Standard State Pressure = p <sup>o</sup> = 0.1 MPa		log K <sub>r</sub>
		S <sup>o</sup> / J·K <sup>-1</sup> ·mol <sup>-1</sup>	(G <sup>o</sup> - H <sup>o</sup> (T <sub>r</sub> ))/T	H <sup>o</sup> - H <sup>o</sup> (T <sub>r</sub> ) / kJ·mol <sup>-1</sup>	ΔH <sup>o</sup> / kJ·mol <sup>-1</sup>	
0	0	0	INFINITE	-4.973	0	0
100	16.033	9.612	52.820	-4.321	0	0
200	22.489	23.273	34.839	-2.313	0	0
250	23.686	28.431	33.056	-1.156	0	0
298.15	24.295	32.660	32.660	0	0	0
300	24.313	32.610	32.660	0.045	0	0
350	24.643	34.583	32.957	1.269	0	0
400	24.978	36.893	33.622	2.509	0	0
450	25.314	39.842	34.483	3.760	0	0
500	25.539	43.462	35.456	5.023	0	0
600	25.790	50.163	37.529	7.580	0	0
700	26.229	54.172	39.627	10.181	0	0
800	26.669	57.703	41.670	12.826	0	0
900	27.112	60.870	43.631	15.515	0	0
1000	27.564	63.750	45.501	18.249	0	0
1100	28.017	66.398	47.282	21.028	0	0
1200	28.472	68.855	48.978	23.852	0	0
1300	28.930	71.132	50.596	26.722	0	0
1400	29.393	73.313	52.143	29.639	0	0
1500	29.862	75.357	53.623	32.601	0	0
1600	30.334	77.299	55.042	35.611	0	0
1700	30.807	79.132	56.407	38.668	0	0
1800	31.284	80.927	57.720	41.773	0	0
1900	31.765	82.681	58.986	44.923	0	0
2000	32.254	84.373	60.210	48.126	0	0
2100	32.744	85.858	61.394	51.376	0	0
2200	33.238	87.393	62.541	54.675	0	0
2300	33.736	88.881	63.654	58.024	0	0
2400	34.233	90.328	64.735	61.422	0	0
2500	34.736	91.735	65.787	64.870	0	0
2600	35.246	93.106	66.812	68.366	0	0
2700	35.762	94.453	67.811	71.935	0	0
2800	36.284	95.791	68.786	75.613	0	0
2900	36.812	97.133	69.740	79.440	0	0
3000	37.346	98.490	70.676	83.442	0	0
3100	37.884	99.873	71.595	87.662	0	0
3200	38.426	101.292	72.501	92.132	0	0
3300	38.972	102.752	73.396	96.877	0	0
3400	39.522	104.261	74.281	101.933	0	0
3500	40.076	105.833	75.160	107.563	0	0
3600	40.634	107.499	76.035	113.270	0	0
3680.000	41.196	108.904	76.734	118.383	---	---
3700	41.762	109.265	76.909	119.716	0.191	-0.003
3800	42.332	111.130	77.785	126.712	-31.336	-0.015
3900	42.906	113.089	78.665	134.256	-27.348	-0.025
4000	43.484	115.138	79.551	142.349	-22.811	-0.034
4100	44.066	117.272	80.445	150.992	-17.725	-0.040
4200	44.652	119.487	81.348	160.184	-12.090	-0.045
4300	45.242	121.778	82.261	169.924	-5.905	-0.047
4400	45.836	124.144	83.186	180.214	0.828	-0.048
4500	46.434	126.579	84.123	191.054	8.111	-0.047
4600	47.036	129.082	85.073	202.442	15.943	-0.044
4700	47.642	131.649	86.037	214.379	24.324	-0.039
4800	48.252	134.278	87.014	226.866	33.254	-0.032

PREVIOUS: December 1961

CURRENT: June 1966

$$\Delta H_f^{\circ}(0 \text{ K}) = 0 \text{ kJ}\cdot\text{mol}^{-1}$$

$$\Delta H_f^{\circ}(298.15 \text{ K}) = 0 \text{ kJ}\cdot\text{mol}^{-1}$$

$$\Delta_{\text{sub}}H^{\circ} = 35.40 \pm 10.5 \text{ kJ}\cdot\text{mol}^{-1}$$

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## Enthalpy of Formation

Zero by definition.

## Heat Capacity and Entropy

There have been several investigations of the low temperature heat capacity of tungsten, Lange<sup>1</sup> covered the range 26–91 K. Horowitz and Daulton<sup>2</sup> reported values in the range 1–77 K, while Waite *et al.*<sup>3</sup> worked from 4–15° and DeSorbo<sup>4</sup> from 15–90 K. Zwicker and Schmidt<sup>5</sup> covered the range from 90 to 2521 K. Clusius *et al.*<sup>6</sup> made a thorough study of the heat capacity in the range 12–274 K, and their measurements were adopted leading to S<sup>o</sup>(298.15 K) = 7.806 cal·K<sup>-1</sup>·mol<sup>-1</sup> based on S<sup>o</sup>(12.5) = 0.0088 cal·K<sup>-1</sup>·mol<sup>-1</sup>. This value is in disagreement with the quoted value of 7.83 cal·K<sup>-1</sup>·mol<sup>-1</sup> but agrees exactly with a separate integration by Kirillin *et al.*<sup>7</sup> in the intermediate temperature range the adiabatic heat capacity measurements of Bronson *et al.*<sup>8</sup> from 253 to 773 K joined well with the low temperature measurements and were adopted. The high temperature enthalpies have been measured by several investigators from which were selected the values of Jaeger and Rosenbohm<sup>9</sup> from 273–1800 K; Magnus and Holzmann<sup>10</sup> who covered the range 373 to 1173 K; Hoch and Johnston,<sup>11</sup> who made measurements between 1382 and 2900 K; Kirillin *et al.*<sup>7</sup> who have made several determinations over the range 600 to 3100 K, and have also analysed the above data and have presented smooth functions from 0 to 3500 K. The present table agrees with that of Kirillin *et al.*<sup>7</sup> up to 2700 K, above this temperature the values of C<sub>p</sub> adopted follow those reported by Novikov *et al.*<sup>12</sup> measured by an electric modulation method, up to the melting point. These values rise rapidly above 2700 K and are not inconsistent with the individual measurements of Kirillin *et al.*<sup>7</sup> at their highest temperatures.

## Fusion Data

Langmuir<sup>13</sup> determined the melting point as 3540 K from intrinsic brilliance measurements, this was later corrected to 3655 ± 30 K by Jones *et al.*<sup>14</sup> and Pirani and Alterthum<sup>15</sup> from pyrometer measurements on a black body hole found 3660 ± 60 K. Using the same technique Zalabak<sup>16</sup> reports 3680 K on a low carbon specimen. He reports a decrease of the melting point with increasing carbon content. Rudy and Windisch<sup>17</sup> report a melting point of 3696 ± 20 K. The value adopted is 3680 K ± 20 K. The enthalpy of melting is obtained by assuming an entropy of melting of 2.3 cal·K<sup>-1</sup>·mol<sup>-1</sup> obtained from a comparison of several high melting metals (Fe, Cu, Co, Mg, Al).

## Sublimation Data

Refer to the ideal gas table for details.

## References

- <sup>1</sup>F. Lange, Z. Physik. Chem. 110, 343 (1924).
- <sup>2</sup>M. Horowitz and J. G. Daulton, Phys. Rev. 91, 1099 (1953).
- <sup>3</sup>T. R. Waite, R. S. Craig, and W. E. Wallace, Phys. Rev. 104, 1240 (1956).
- <sup>4</sup>W. DeSorbo, J. Phys. Chem. 62, 965 (1958).
- <sup>5</sup>C. Zwicker and G. Schmidt, Z. Physik. 52, 668 (1928).
- <sup>6</sup>K. Clusius and P. Franzosini, Z. Naturforsch. 14A, 99 (1959).
- <sup>7</sup>V. A. Kirillin, A. E. Sheindlin, V. Ya. Chekhovskoi and V. A. Petrov, Zhur. Fiz. Khim. 37, 2249 (1963).
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- <sup>9</sup>F. M. Jaeger and E. Rosenbohm, Rec. trav. chim. 51, 1 (1932).
- <sup>10</sup>A. Magnus and H. Holzmann, Ann. Physik. Ser. 5, 3, 585 (1929).
- <sup>11</sup>M. Hoch and H. L. Johnston, J. Phys. Chem. 65, 855 (1961).
- <sup>12</sup>I. I. Novikov and P. G. Streikov, Vestnik. Akad. Nauk SSSR 34, 26 (1964).
- <sup>13</sup>J. Langmuir, Phys. Rev. 6, 138 (1915).
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- <sup>15</sup>M. Pirani and H. Alterthum, Z. Elektrochem. 29, 5 (1923).
- <sup>16</sup>C. F. Zalabak, NASA Tech. Note D-761 (1961).
- <sup>17</sup>E. Rudy and S. Windisch, Aerojet-General Tech. Rept. No. AFML-TR-65-2 Part 1, Vol. III (July 1965).

Tungsten (W)

W<sub>1</sub>(cr)

**Tungsten (W)**  $A_f = 183.85$  **Tungsten (W)**  $W_r(0)$

T/K	$C_p^*$	$S^*$ $-\{G^*-F(T)\}/T$	$H^*-F(T)$	$\Delta_f G^*$	$\log K_f$
Enthalpy Reference Temperature = $T_r = 298.15$ K					
$\text{J}\cdot\text{K}^{-1}\cdot\text{mol}^{-1}$ $\text{kJ}\cdot\text{mol}^{-1}$					
0	0				
100	24.296	45.678	45.678	0.	46.896
200	24.310	45.828	45.678	0.045	46.896
250	24.342	46.001	45.678	0.169	46.896
300	24.378	46.191	45.678	0.298	46.896
350	24.418	46.397	45.678	0.432	46.896
400	24.462	46.619	45.678	0.571	46.896
450	24.511	46.857	45.678	0.715	46.896
500	24.564	47.112	45.678	0.864	46.896
600	24.631	47.484	45.678	1.018	46.896
700	24.703	47.874	45.678	1.177	46.896
800	24.780	48.281	45.678	1.341	46.896
900	24.862	48.705	45.678	1.510	46.896
1000	24.949	49.146	45.678	1.684	46.896
1100	25.041	49.604	45.678	1.863	46.896
1200	25.138	50.079	45.678	2.047	46.896
1300	25.240	50.571	45.678	2.236	46.896
1400	25.347	51.080	45.678	2.429	46.896
1500	25.459	51.606	45.678	2.626	46.896
1600	25.576	52.149	45.678	2.827	46.896
1700	25.698	52.709	45.678	3.032	46.896
1800	25.825	53.286	45.678	3.241	46.896
1900	25.957	53.880	45.678	3.454	46.896
2000	26.094	54.491	45.678	3.671	46.896
2100	26.236	55.118	45.678	3.892	46.896
2200	26.383	55.762	45.678	4.117	46.896
2300	26.535	56.423	45.678	4.346	46.896
2400	26.692	57.101	45.678	4.579	46.896
2450.000	26.780	57.400	45.678	4.680	46.896
2450.000	26.780	57.400	45.678	4.680	46.896
2500	26.880	57.720	45.678	4.780	46.896
2600	27.080	58.360	45.678	4.980	46.896
2700	27.290	59.020	45.678	5.180	46.896
2800	27.510	59.700	45.678	5.380	46.896
2900	27.740	60.400	45.678	5.580	46.896
3000	27.980	61.120	45.678	5.780	46.896
3100	28.230	61.860	45.678	5.980	46.896
3200	28.490	62.620	45.678	6.180	46.896
3300	28.760	63.400	45.678	6.380	46.896
3400	29.040	64.200	45.678	6.580	46.896
3500	29.330	65.020	45.678	6.780	46.896
3600	29.630	65.860	45.678	6.980	46.896
3680.000	29.780	66.300	45.678	7.100	46.896
3700	29.800	66.350	45.678	7.120	46.896
3800	29.900	66.450	45.678	7.150	46.896
3900	29.990	66.550	45.678	7.180	46.896
4000	30.080	66.650	45.678	7.210	46.896
4100	30.170	66.750	45.678	7.240	46.896
4200	30.260	66.850	45.678	7.270	46.896
4300	30.350	66.950	45.678	7.300	46.896
4400	30.440	67.050	45.678	7.330	46.896
4500	30.530	67.150	45.678	7.360	46.896
4600	30.620	67.250	45.678	7.390	46.896
4700	30.710	67.350	45.678	7.420	46.896
4800	30.800	67.450	45.678	7.450	46.896
4900	30.890	67.550	45.678	7.480	46.896
5000	30.980	67.650	45.678	7.510	46.896
5200	31.160	67.830	45.678	7.550	46.896
5400	31.340	68.010	45.678	7.590	46.896
5600	31.520	68.190	45.678	7.630	46.896
5800	31.700	68.370	45.678	7.670	46.896
6000	31.880	68.550	45.678	7.710	46.896

$S^*(298.15 \text{ K}) = [45.678] \text{ J}\cdot\text{K}^{-1}\cdot\text{mol}^{-1}$   
 $T_{\text{fus}} = 3680 \pm 20 \text{ K}$   
 $\Delta H_f^*(298.15 \text{ K}) = [46.896] \text{ kJ}\cdot\text{mol}^{-1}$   
 $\Delta_{\text{liq}} H^* = 35.40 \pm 10.5 \text{ kJ}\cdot\text{mol}^{-1}$

**Enthalpy of Formation**  
 The enthalpy of formation at 298.15 K was calculated from that of the crystal by adding  $\Delta_{\text{fus}} H^*$  and the difference in molar enthalpy,  $H^*(3680 \text{ K}) - H^*(298.15 \text{ K})$ , between the crystal and liquid.

**Heat Capacity and Entropy**  
 The heat capacity was estimated as  $8.5 \text{ cal}\cdot\text{mol}^{-1}\cdot\text{K}^{-1}$  by analogy with other monatomic metals. The entropy at 298.15 K was calculated in manner analogous to that used for the enthalpy of formation. At 2450 K a glass transition is assumed below which the heat capacity was that of the crystal.

**Fusion Data**  
 Refer to the crystal table for details.

**Vaporization Data**  
 The boiling point and enthalpy of vaporization are calculated from the adopted functions and enthalpy of sublimation in order to maintain proper thermodynamic consistency.

PREVIOUS: December 1961  
 CURRENT: June 1966

**Tungsten (W)**  $W_r(0)$

CRYSTAL-LIQUID

0 to 3680 K crystal  
above 3680 K liquid

Refer to the individual tables for details.

Tungsten (W)

Tungsten (W)

W<sub>1</sub>(cr,l)

T/K	Enthalpy Reference Temperature = T <sub>r</sub> = 298.15 K		Standard State Pressure = p° = 0.1 MPa		log K <sub>r</sub>
	C <sub>p</sub> <sup>o</sup>	S° - [(C <sub>p</sub> ° - H <sub>f</sub> °(T <sub>r</sub> ))/T]	H° - H <sub>f</sub> °(T)	Δ <sub>f</sub> G°	
		J·K <sup>-1</sup> ·mol <sup>-1</sup>	J·K <sup>-1</sup> ·mol <sup>-1</sup>	kJ·mol <sup>-1</sup>	
0	0	INFINITE	-4.973	0	0
100	16.033	9.612	-4.321	0	0
200	22.489	23.273	-2.313	0	0
250	23.686	28.431	-1.156	0	0
298.15	24.295	32.660	0	0	0
300	24.313	32.810	0.045	0	0
350	24.644	36.583	1.269	0	0
400	24.928	39.893	2.509	0	0
450	25.144	42.842	3.760	0	0
500	25.359	45.502	5.023	0	0
600	25.790	50.163	7.580	0	0
700	26.229	54.172	10.181	0	0
800	26.669	57.703	12.826	0	0
900	27.112	60.870	15.515	0	0
1000	27.564	63.750	18.249	0	0
1100	28.017	66.398	21.028	0	0
1200	28.472	68.855	23.852	0	0
1300	28.930	71.152	26.722	0	0
1400	29.393	73.313	29.639	0	0
1500	29.862	75.357	32.601	0	0
1600	30.334	77.299	35.611	0	0
1700	30.807	79.152	38.668	0	0
1800	31.284	80.927	41.773	0	0
1900	31.765	82.631	44.925	0	0
2000	32.234	84.273	48.126	0	0
2100	32.744	85.858	51.376	0	0
2200	33.238	87.391	54.673	0	0
2300	33.736	88.881	58.024	0	0
2400	34.233	90.328	61.422	0	0
2500	34.736	91.735	64.870	0	0
2600	35.246	93.106	68.366	0	0
2700	35.762	94.453	71.935	0	0
2800	36.284	95.791	75.613	0	0
2900	36.812	97.135	79.440	0	0
3000	41.003	98.490	83.442	0	0
3100	43.430	99.873	87.662	0	0
3200	46.024	101.292	92.132	0	0
3300	48.953	102.752	96.877	0	0
3400	52.300	104.261	101.933	0	0
3500	56.484	105.835	107.363	0	0
3600	61.714	107.499	113.270	0	0
3680.000	66.149	108.904	118.383	CRYSTAL ← → LIQUID	0
3680.000	35.564	118.372	153.780	TRANSITION	0
3700	35.564	118.715	154.491	0	0
3800	35.564	119.995	158.048	0	0
3900	35.564	120.587	161.604	0	0
4000	35.564	121.488	165.161	0	0
4100	35.564	122.566	168.717	0	0
4200	35.564	123.223	172.273	0	0
4300	35.564	124.060	175.830	0	0
4400	35.564	124.877	179.386	0	0
4500	35.564	125.676	182.943	0	0
4600	35.564	126.458	186.499	0	0
4700	35.564	127.223	190.055	0	0
4800	35.564	127.972	193.612	0	0
4900	35.564	128.705	197.168	0	0
5000	35.564	129.423	200.725	0	0
5100	35.564	130.128	204.281	0	0
5200	35.564	130.818	207.837	0	0
5300	35.564	131.496	211.394	0	0
5400	35.564	132.161	214.950	0	0
5500	35.564	132.813	218.507	0	0
5600	35.564	133.454	222.063	0	0
5700	35.564	134.083	225.619	0	0
5800	35.564	134.702	229.176	0	0
5900	35.564	135.310	232.732	0	0
6000	35.564	135.908	236.289	0	0

PREVIOUS:

CURRENT: June 1966

Tungsten (W)

W<sub>1</sub>(cr,l)

Tungsten (W)

Tungsten (W)

IDEAL GAS

W<sub>1</sub>(g)

IP(W,g) = 64400 ± 100 cm<sup>-1</sup>  
 S<sup>o</sup>(298.15 K) = 173.95 ± 0.08 J·K<sup>-1</sup>·mol<sup>-1</sup>

A<sub>r</sub> = 183.85 Tungsten (W)  
 ΔH<sup>o</sup>(0 K) = 849.8 ± 6.3 kJ·mol<sup>-1</sup>  
 ΔH<sup>o</sup>(298.15 K) = 851.0 ± 6.3 kJ·mol<sup>-1</sup>

Electronic Levels and Quantum Weights	State	ε <sub>i</sub> , cm <sup>-1</sup>	g <sub>i</sub>
<sup>3</sup> D <sub>0</sub>	1	0.00	1
<sup>3</sup> D <sub>1</sub>	3	1670.29	3
<sup>3</sup> D <sub>2</sub>	5	3325.53	5
<sup>3</sup> D <sub>3</sub>	7	4830.00	7
<sup>3</sup> D <sub>4</sub>	9	6219.33	9
.	.	.	.
.	.	.	.
.	.	.	.
.	.	.	.
.	.	.	.
IP	11	62154.50	11
.	.	66400.00	.

Enthalpy of Formation

The enthalpy of formation is the enthalpy of sublimation at 298.15 K; this has been obtained by a 2nd and 3rd law analysis of the vapor pressure data of three investigators.

Source	T/K**	Data Points	Δ <sub>sub</sub> H <sup>o</sup> (298.15 K), kcal·mol <sup>-1</sup>	Drift
1	2511-3053	12*	200.9 ± 3.6	201.57 ± 1.6
2	2383-3123	14	214.8 ± 1.4	202.17 ± 2.5
3	2574-3183	10	202.4 ± 2.7	203.35 ± 1.2

\* 1 point rejected due to failure of a statistical test.

\*\* All temperatures are taken from Szwarc<sup>3</sup> who has corrected the earlier measurements.

The values are remarkably good since even at the highest temperature the pressures are less than 10<sup>-6</sup> atm; all workers used the Langmuir technique. Since the data of references 1 and 3 do not drift, it may be assumed that the accommodation coefficient is unity or close to unity. The drift in the second set cannot be eliminated by assuming a non-unity coefficient, nor does there appear to be a constant pressure error. Most probably the drift is due to slight errors in temperature, for example if the readings were 24 K high at the low end and 10 K low at the high end the drift would be eliminated; errors of this magnitude are quite possible. The value adopted for Δ<sub>sub</sub>H<sup>o</sup>(298.15 K) is 203.4 ± 1.5 kcal·mol<sup>-1</sup>.

Heat Capacity and Entropy

The electronic energy levels are taken from those listed by Moore.<sup>4</sup> Levels above 25000 cm<sup>-1</sup> were averaged.

References

- <sup>1</sup>H. A. Jones, I. Langmuir and G. M. J. Mackay, Phys. Rev. 30, 201 (1927).
- <sup>2</sup>C. Zwickler, Physica 5, 249 (1925).
- <sup>3</sup>R. Szwarc, E. R. Plante, and J. J. Diamond, J. Research Nat. Bur. Stand. 69A, 417 (1965).

T/K	C <sub>p</sub> <sup>o</sup>	S <sup>o</sup> - (G <sup>o</sup> - H <sup>o</sup> (T))/T	H <sup>o</sup> - H <sup>o</sup> (T)	ΔH <sup>o</sup>	log K <sub>r</sub>	W <sub>1</sub> (g)
0	0.000	INFINITE	-6.216	849.782	INFINITE	
100	20.786	151.176	-4.138	851.209	-437.231	
200	20.808	165.586	-2.059	852.817	-214.898	
250	20.941	170.280	-1.016	851.166	-170.434	
298.15	21.306	173.955	.000	851.026	808.898	
300	21.326	174.087	.039	808.637	-141.716	
350	22.059	174.425	1.123	850.879	-140.796	
400	23.163	180.438	2.262	850.769	-119.630	
450	24.611	183.247	3.442	850.710	-103.738	
500	26.344	185.977	4.717	850.720	-91.414	
600	30.337	191.077	7.548	850.993	-81.539	
700	34.381	196.061	10.786	851.631	-66.725	
800	37.765	200.883	14.402	852.601	-56.130	
900	40.071	205.476	18.303	853.814	-48.190	
1000	41.235	209.768	22.377	855.154	-42.001	
1100	41.438	213.714	26.518	856.516	-37.041	
1200	40.951	217.303	30.642	857.815	-32.978	
1300	40.037	220.547	34.694	858.997	-29.598	
1400	38.900	223.473	38.642	860.079	-26.711	
1500	37.687	226.115	42.472	860.896	-24.248	
1600	36.488	228.509	46.180	861.595	-22.104	
1700	35.360	230.687	49.772	862.129	-20.230	
1800	34.331	232.678	53.255	862.508	-18.575	
1900	33.417	234.516	56.642	862.742	-17.103	
2000	32.621	236.203	59.943	862.842	-15.785	
2100	31.943	237.778	63.170	862.820	-14.599	
2200	31.379	239.250	66.335	862.686	-13.526	
2300	30.922	240.635	69.449	862.451	-12.551	
2400	30.566	241.943	72.525	862.126	-11.660	
2500	30.304	243.185	75.566	861.721	-10.844	
2600	30.129	244.370	78.586	861.246	-10.094	
2700	30.035	245.505	81.594	860.685	-9.401	
2800	30.013	246.596	84.596	860.008	-8.761	
2900	30.060	247.644	87.589	859.184	-8.166	
3000	30.167	248.671	90.610	858.194	-7.619	
3100	30.331	249.663	93.634	856.998	-7.119	
3200	30.545	250.629	96.678	855.571	-6.665	
3300	30.805	251.575	99.745	853.893	-6.252	
3400	31.106	252.497	102.840	851.933	-5.874	
3500	31.446	253.403	105.967	849.630	-5.545	
3600	31.822	254.294	109.130	846.886	-5.252	
3700	32.230	255.172	112.333	843.627	-5.000	
3800	32.670	256.037	115.578	840.555	-4.781	
3900	33.141	256.892	118.868	837.702	-4.591	
4000	33.642	257.737	122.207	835.072	-4.435	
4100	34.174	258.574	125.597	832.680	-4.308	
4200	34.738	259.404	129.043	830.437	-4.206	
4300	35.335	260.229	132.546	828.342	-4.125	
4400	35.968	261.048	136.111	826.393	-4.061	
4500	36.638	261.864	139.741	824.580	-4.011	
4600	37.348	262.677	143.440	822.896	-3.972	
4700	38.102	263.488	147.212	821.330	-3.943	
4800	38.900	264.298	151.061	819.875	-3.923	
4900	39.746	265.109	154.993	818.511	-3.911	
5000	40.642	265.921	159.012	817.225	-3.905	
5100	41.590	266.735	163.124	816.008	-3.905	
5200	42.591	267.552	167.332	814.850	-3.910	
5300	43.646	268.374	171.644	813.755	-3.919	
5400	44.755	269.200	176.066	812.727	-3.931	
5500	45.916	270.031	180.596	811.765	-3.945	
5600	47.129	270.870	185.248	810.868	-3.961	
5700	48.391	271.715	190.024	810.030	-3.978	
5800	49.698	272.568	194.928	809.250	-3.995	
5900	51.046	273.429	199.965	808.511	-4.011	
6000	52.431	274.298	205.138	807.815	-4.026	

PREVIOUS: June 1966 (1 atm)

CURRENT: June 1966 (1 bar)

Tungsten (W)

W<sub>1</sub>(g)

Tungsten, Ion (W<sup>+</sup>)

IDEAL GAS

Tungsten, Ion (W<sup>+</sup>)

W<sup>+</sup>(g)

IP(W<sup>+</sup>, g) = 142800 ± 4000 cm<sup>-1</sup>  
 S<sup>o</sup>(298.15 K) = 179.74 ± 0.08 J·K<sup>-1</sup>·mol<sup>-1</sup>

Δ<sub>f</sub>H<sup>o</sup>(0 K) = 1620.2 ± 7 kJ·mol<sup>-1</sup>  
 Δ<sub>f</sub>H<sup>o</sup>(298.15 K) = [1627.623] kJ·mol<sup>-1</sup>

Electronic Levels and Quantum Weights	g <sub>i</sub>
State	ε <sub>i</sub> , cm <sup>-1</sup>
<sup>5</sup> D <sub>3/2</sub>	0.00
<sup>5</sup> D <sub>5/2</sub>	1518.78
<sup>5</sup> D <sub>7/2</sub>	3172.52
<sup>5</sup> D <sub>9/2</sub>	4716.32
<sup>5</sup> D <sub>9/2</sub>	6147.16
.	.
.	.
<sup>4</sup> F <sub>6/2</sub>	64516.37
IP	[142800]

Enthalpy of Formation

Δ<sub>f</sub>H<sup>o</sup>(W<sup>+</sup>, g, 0 K) is calculated from Δ<sub>f</sub>H<sup>o</sup>(W, g, 0 K) using the spectroscopic value of IP(W) = 64400 ± 100 cm<sup>-1</sup> (770.40 ± 1.20 kJ·mol<sup>-1</sup>) from Moore.<sup>2</sup> The ionization limit is converted from cm<sup>-1</sup> to kJ·mol<sup>-1</sup> using the factor, 1 cm<sup>-1</sup> = 0.01196266 kJ·mol<sup>-1</sup>, which is derived from the 1973 CODATA fundamental constants.<sup>3</sup> Rosenstock *et al.*<sup>4</sup> and Levin and Lias<sup>5</sup> have summarized additional ionization and appearance potential data.

Δ<sub>f</sub>H<sup>o</sup>(W<sup>+</sup>, g, 298.15 K) is calculated from Δ<sub>f</sub>H<sup>o</sup>(W, g, 0 K) by using IP(W) with JANAF<sup>1</sup> enthalpies, H<sup>o</sup>(0 K) - H<sup>o</sup>(298.15 K), for W(g), W<sup>+</sup>(g), and e<sup>-</sup> (ref). Δ<sub>f</sub>H<sup>o</sup>(W → W<sup>+</sup> + e<sup>-</sup>, 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.*<sup>4</sup> Δ<sub>f</sub>H<sup>o</sup>(298.15 K) should be changed by -6.197 kJ·mol<sup>-1</sup> 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,<sup>2</sup> is incomplete because many theoretically predicted levels have not been observed. Although we have listed only the ground, the first four excited states, the highest observed excited state, and the ionization potential for W<sup>+</sup>(g), all levels listed by Moore<sup>2</sup> as well as estimated levels, are used in the calculation. The observed levels are too numerous to list completely. The calculations indicate that for W<sup>+</sup>(g), the thermodynamic functions are independent of the estimated missing levels (for n = 6) and the cut-off procedure; the Gibbs energy function showing variations of 0.8% at this temperature. The reported uncertainty is S<sup>o</sup>(298.15 K) is due to uncertainties in the relative ionic mass, the fundamental constants, and the low lying atomic energy levels. Extension of these calculations above 6000 K may require consideration of the higher excited states (n > 6), and use of different fill and cut-off procedures.<sup>1</sup>

References

- JANAF Thermochemical Tables: W(g), 6-30-66; e<sup>-</sup> (ref), 3-31-82.
- C. E. Moore, U. S. Nat. Bur. Stand., NSRDS-NBS-34, 8 pp. (1970).
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- H. M. Rosenstock, K. Draxl *et al.*, J. Phys. Chem. Ref. Data, 6, Supp. 1, 783 pp. (1977).
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- C. E. Moore, U. S. Nat. Bur. Stand., NSRDS-NBS-35, Volume III, (1970) [Reprint of NBS Circular 467, Volume III, 1958].
- J. R. Downey, Jr., The Dow Chemical Company, AFOSR-TR-78-0960, Contract No. F44620-75-1-0048, (1978).

T/K	Enthalpy Reference Temperature = T <sub>r</sub> = 298.15 K		Standard State Pressure = p <sup>o</sup> = 0.1 MPa		log K <sub>i</sub>
	C <sub>p</sub> <sup>o</sup>	S <sup>o</sup> - [G <sup>o</sup> - H <sup>o</sup> (T <sub>r</sub> )]/T	H <sup>o</sup> - H <sup>o</sup> (T <sub>r</sub> )	Δ <sub>f</sub> G <sup>o</sup>	
0	.000	INFINITE	-6.221	1620.178	-276.374
100	20.786	156.939	1.225	1627.623	-274.616
200	20.822	171.350	1.125	1627.656	-274.616
250	20.989	176.011	1.067	1627.656	-274.616
298.15	21.372	179.737	.000	1627.656	-274.616
300	21.391	179.869	.040	1627.656	-274.616
350	21.043	183.213	1.125	1628.556	-234.120
400	22.891	186.210	2.247	1629.479	-203.730
450	23.844	188.961	3.415	1630.434	-180.081
500	24.828	191.524	4.632	1631.428	-161.150
600	26.695	196.219	7.210	1633.527	-132.727
700	28.294	200.458	9.962	1635.757	-112.398
800	29.616	204.325	12.860	1638.088	-97.130
900	30.703	207.878	15.877	1640.495	-85.238
1000	31.603	211.161	18.994	1642.957	-75.710
1100	32.347	214.209	22.193	1645.455	-67.902
1200	32.951	217.051	25.469	1647.975	-61.333
1300	33.441	219.709	28.780	1650.505	-55.864
1400	33.813	222.201	32.145	1653.041	-51.153
1500	34.082	224.544	35.559	1655.582	-47.008
1600	34.261	226.750	38.957	1658.029	-43.403
1700	34.362	228.830	42.389	1660.482	-40.216
1800	34.401	230.795	45.827	1662.895	-37.380
1900	34.393	232.655	49.267	1665.261	-34.838
2000	34.351	234.419	52.705	1667.576	-32.472
2100	34.288	236.093	56.137	1669.837	-30.247
2200	34.218	237.683	59.562	1672.042	-28.148
2300	34.148	239.206	62.980	1674.190	-26.856
2400	34.078	240.658	66.392	1676.282	-25.271
2500	34.041	242.048	69.798	1678.319	-23.810
2600	34.013	243.383	73.201	1680.304	-22.461
2700	34.008	244.666	76.602	1682.214	-21.210
2800	34.026	245.900	80.003	1684.016	-20.047
2900	34.057	247.098	83.405	1685.672	-18.951
3000	34.130	248.254	86.817	1687.159	-17.951
3100	34.216	249.375	90.234	1688.435	-17.003
3200	34.320	250.462	93.661	1689.470	-16.113
3300	34.442	251.520	97.099	1690.241	-15.277
3400	34.578	252.551	100.550	1690.715	-14.490
3500	34.725	253.555	104.015	1690.838	-13.748
3600	34.882	254.535	107.495	1690.480	-13.047
3700	35.044	255.493	110.991	1689.634	-12.387
3800	35.209	256.430	114.504	1688.309	-11.772
3900	35.375	257.347	118.033	1686.507	-11.188
4000	35.538	258.244	121.579	1684.232	-10.632
4100	35.698	259.124	125.141	1681.486	-10.103
4200	35.851	259.986	128.718	1678.272	-9.598
4300	35.996	260.831	132.311	1674.604	-9.116
4400	36.132	261.660	135.917	1670.486	-8.655
4500	36.257	262.474	139.537	1665.923	-8.215
4600	36.370	263.272	143.168	1660.910	-7.792
4700	36.471	264.055	146.810	1655.450	-7.388
4800	36.559	264.824	150.462	1649.549	-6.997
4900	36.634	265.579	154.122	1643.201	-6.627
5000	36.695	266.319	157.788	1636.419	-6.268
5100	36.743	267.047	161.460	1629.203	-5.923
5200	36.778	267.760	165.136	1621.562	-5.591
5300	36.800	268.461	168.815	1613.503	-5.271
5400	36.811	269.149	172.496	1605.031	-4.963
5500	36.810	269.825	176.177	1596.149	-4.665
5600	36.799	270.488	179.858	1586.852	-4.378
5700	36.779	271.139	183.537	1577.140	-4.100
5800	36.749	271.778	187.213	1567.002	-3.832
5900	36.713	272.406	190.886	1556.438	-3.572
6000	36.669	273.023	194.557	1545.455	-3.321

PREVIOUS:

CURRENT: March 1984 (1 bar)

Tungsten, Ion (W<sup>+</sup>)

W<sup>+</sup>(g)

Tungsten, Ion (W<sup>-</sup>) IDEAL GAS W<sub>r</sub>(g)

EA(W, g) = 0.815 ± 0.008 eV M<sub>r</sub> = 183.85055 Tungsten, Ion (W<sup>-</sup>)  
 S°(298.15 K) = 188.781 ± 0.08 J·K<sup>-1</sup>·mol<sup>-1</sup> ΔH<sup>o</sup>(0 K) = 771.147 ± 2.0 kJ·mol<sup>-1</sup> ΔH<sup>o</sup>(298.15 K) = [766.174] kJ·mol<sup>-1</sup>

Electronic Level and Quantum Weight	
State	g
<sup>6</sup> S <sub>3/2</sub>	6

**Enthalpy of Formation**  
 Δ<sub>f</sub>H<sup>o</sup>(W<sup>-</sup>, g, 0 K) is calculated from Δ<sub>f</sub>H<sup>o</sup>(W, g, 0 K)<sup>1</sup> using the adopted electron affinity of EA(W) = 0.815 ± 0.008 eV (78.635 ± 0.772 kJ·mol<sup>-1</sup>). This value, recommended by Hotop and Lineberger,<sup>2</sup> is based on a laser photodetachment electron spectrometry study.<sup>3</sup> Additional information on W<sup>-</sup>(g) may be obtained in the critical discussions of Hotop and Lineberger,<sup>2,4</sup> Rosenstock *et al.*,<sup>5</sup> and Massey.<sup>6</sup> Δ<sub>f</sub>H<sup>o</sup>(W<sup>-</sup>, g, 298.15 K) is obtained from Δ<sub>f</sub>H<sup>o</sup>(W, g, 0 K) by using EA(W) with JANAF<sup>1</sup> enthalpies, H<sup>o</sup>(0 K)–H<sup>o</sup>(298.15 K), for W<sup>-</sup>(g), W(g), and e<sup>-</sup>(ref). Δ<sub>f</sub>H<sup>o</sup>(W<sup>-</sup> → W + e<sup>-</sup>, 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.*<sup>5</sup> Δ<sub>f</sub>H<sup>o</sup>(298.15 K) should be changed by +6.197 kJ·mol<sup>-1</sup> 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 W<sup>-</sup>(g) is given by Hotop and Lineberger<sup>2,4</sup> and Rosenstock *et al.*<sup>5</sup> Lacking any experimental evidence as to the stability of any excited states, we assume that no stable excited states exist.

- References**  
<sup>1</sup>JANAF Thermochemical Tables: W(g), 6–30–66; e<sup>-</sup>(ref), 3–31–82.  
<sup>2</sup>H. Hotop and W. C. Lineberger, *J. Phys. Chem. Ref. Data*, **14**, 731 (1985).  
<sup>3</sup>C. S. Feigler, R. R. Cordermann, S. V. Bobashev and W. C. Lineberger, *J. Chem. Phys.* **74**, 1580 (1981).  
<sup>4</sup>H. Hotop and W. C. Lineberger, *J. Phys. Chem. Ref. Data* **4**, 539 (1975).  
<sup>5</sup>H. M. Rosenstock, K. Draxl *et al.*, *J. Phys. Chem. Ref. Data* **6**, Supp. 1, 783 pp. (1977).  
<sup>6</sup>H. S. W. Massey, "Negative Ions", 3rd ed., Cambridge University Press, Cambridge, 741 pp. (1976).

T/K	Enthalpy Reference Temperature = T <sub>r</sub> = 298.15 K		Standard State Pressure = P <sup>o</sup> = 0.1 MPa		log K <sub>r</sub>
	C <sub>p</sub> <sup>o</sup>	S <sup>o</sup> – (G <sup>o</sup> – H <sup>o</sup> (T <sub>r</sub> ))/T	H <sup>o</sup> – H <sup>o</sup> (T <sub>r</sub> )/T	Δ <sub>f</sub> H <sup>o</sup>	
0	0	INFINITE	–6.197	771.147	–171.171
100	20.786	166.073	–0.7261	766.174	–126.344
200	20.786	180.481	–2.040	766.129	–107.302
250	20.786	185.119	–1.001	766.129	–91.043
298.15	20.786	188.781	0	766.129	–83.971
300	20.786	188.909	0.038	766.129	–83.971
350	20.786	192.113	0.778	766.129	–73.129
400	20.786	194.889	2.117	766.129	–68.011
450	20.786	197.337	3.156	766.129	–60.420
500	20.786	199.527	4.196	766.129	–54.838
600	20.786	203.317	6.274	766.129	–49.987
700	20.786	206.588	8.535	766.129	–46.452
800	20.786	209.297	10.431	766.129	–43.629
900	20.786	211.745	12.510	766.129	–41.292
1000	20.786	213.935	14.589	766.129	–39.632
1100	20.786	215.916	16.667	766.129	–38.087
1200	20.786	217.725	18.746	766.129	–37.144
1300	20.786	219.388	20.824	766.129	–36.663
1400	20.786	220.929	22.903	766.129	–36.451
1500	20.786	222.363	24.982	766.129	–36.451
1600	20.786	223.704	27.060	766.129	–36.663
1700	20.786	224.965	29.139	766.129	–37.144
1800	20.786	226.153	31.217	766.129	–37.807
1900	20.786	227.277	33.296	766.129	–38.629
2000	20.786	228.343	35.375	766.129	–39.600
2100	20.786	229.357	37.453	766.129	–40.717
2200	20.786	230.324	39.532	766.129	–41.989
2300	20.786	231.248	41.611	766.129	–43.414
2400	20.786	232.132	43.689	766.129	–44.992
2500	20.786	232.981	45.768	766.129	–46.720
2600	20.786	233.796	47.846	766.129	–48.600
2700	20.786	234.581	49.925	766.129	–50.639
2800	20.786	235.337	52.004	766.129	–52.843
2900	20.786	236.066	54.082	766.129	–55.214
3000	20.786	236.771	56.161	766.129	–57.757
3100	20.786	237.452	58.239	766.129	–60.474
3200	20.786	238.112	60.318	766.129	–63.373
3300	20.786	238.752	62.397	766.129	–66.453
3400	20.786	239.372	64.475	766.129	–69.720
3500	20.786	239.975	66.554	766.129	–73.181
3600	20.786	240.560	68.632	766.129	–76.843
3700	20.786	241.130	70.711	766.129	–80.606
3800	20.786	241.684	72.790	766.129	–84.474
3900	20.786	242.224	74.868	766.129	–88.449
4000	20.786	242.750	76.947	766.129	–92.529
4100	20.786	243.264	79.025	766.129	–96.714
4200	20.786	243.765	81.104	766.129	–101.006
4300	20.786	244.254	83.183	766.129	–105.406
4400	20.786	244.732	85.262	766.129	–110.006
4500	20.786	245.199	87.341	766.129	–114.806
4600	20.786	245.656	89.420	766.129	–119.806
4700	20.786	246.103	91.499	766.129	–125.006
4800	20.786	246.540	93.578	766.129	–130.406
4900	20.786	246.969	95.657	766.129	–136.006
5000	20.786	247.389	97.736	766.129	–141.806
5100	20.786	247.800	99.815	766.129	–147.806
5200	20.786	248.204	101.894	766.129	–154.006
5300	20.786	248.600	103.969	766.129	–160.406
5400	20.786	248.988	106.047	766.129	–167.006
5500	20.786	249.370	108.126	766.129	–173.806
5600	20.786	249.744	110.204	766.129	–180.806
5700	20.786	250.112	112.283	766.129	–188.006
5800	20.786	250.474	114.362	766.129	–195.406
5900	20.786	250.829	116.440	766.129	–203.006
6000	20.786	251.178	118.519	766.129	–210.806

PREVIOUS: CURRENT: March 1984 (1 bar)