Erratum to: Theoretical Calculations of the Surface Tension of Liquid Transition Metals

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THE printer regrets an error in Table 1 of this paper.
The correct Table 1 is as follows:

Table I. Parameters Needed for Calculating Surface Tension of Metals. Calculated and Experimental Surface Tension Values of Some d-Block Metals

Metal	Atomic Weight	Melting Point (K/°C)	Density at Melting Point (g/cm ³)	Molar Volume (cm ³ /mol)	Sublimation Energy (kJ/mol)	f	ϕ ($\times 10^{-16}$)	Calculated Surface Tension (mJ/m ²)	Experimental Surface Tension (mJ/m ²)
Ti	47.86	1941/1668	4.11	11.64	470	0.287	7.86	1380	1500
Zr	91.224	2128/1855	5.8	15.73	609	0.287	9.61	1461	1480
Fe	55.845	1811/1538	6.98	8.0	418	0.287	6.12	1560	1830
Co	58.933	1768/1495	7.75	7.60	428	0.287	5.92	1668	1830
Ni	58.693	1728/1455	7.81	7.52	430	0.287	5.87	1810	1838
Pd	106.42	1828/1555	10.38	10.25	377	0.287	7.22	1205	1475
Pt	195.078	2041/1768	19.77	9.87	566	0.287	7.042	1855	1860
Cu	63.546	1357/1084	8.02	7.92	338	0.287	6.1	1200	1310
Ag	107.868	1235/962	9.32	11.57	285	0.287	7.82	841	910
Au	196.967	1337/1064	17.31	11.38	368	0.287	7.74	1100	1149
Zn	65.39	693/420	6.57	9.95	130	0.240	7.08	854	789
Cd	112.411	594/321	7.996	14.05	112	0.240	8.91	584	637
Hg	200.59	234/-39	13.534	14.82	61	0.210	9.24	428	438

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