

Crystal Structures and Cell Parameters of Allotropes of the Chemical Elements

The crystal structures of the allotropic forms of the metallic elements are presented in terms of the Pearson symbol, space group, Strukturbericht designation, and prototype of the structure. The temperatures of the phase transformations are listed in degrees Celsius and the pressures are in GPa. A consistent nomenclature is used, whereby all allotropes are labeled by Greek letters. The lattice parameters of the unit cells are given in nanometers (nm) and are considered to be accurate to ± 2 in the last reported digit.

This compilation is restricted to changes in crystal structure that occur as a result of a change in temperature or pressure. Low-temperature structures are included for the diatomic and rare gases, which show many similarities with respect to the metallic elements.

This compilation updates and supersedes previous compilations published in the *Bulletin of Alloy Phase Diagrams*.

Most of the data given below are the same as in the first edition. However, occasionally changes were made when updated information was supplied by a category editor, or by present volume editors. The reader's attention is drawn to the fact that there may differences between values quoted below and similar values given in another table in this edition that has been reproduced from another source. For example, the allotropic transformation temperatures of Mn may differ by as much as 23 °C, etc.

Element phase	Common name	Temperature, °C	Pressure, GPa	Struktur- bericht designation	Prototype	Pearson symbol	Space group	Cell parameters, nm			Comment,c/a, or α or β
								a	b	c	
(Ac)	Ac	25	atm	A1	Cu	<i>cF</i> 4	<i>Fm</i> $\bar{3}$ <i>m</i>	0.5311
(Ag)	Ag	25	atm	A1	Cu	<i>cF</i> 4	<i>Fm</i> $\bar{3}$ <i>m</i>	0.40857
(Al)	α Al	25	atm	A1	Cu	<i>cF</i> 4	<i>Fm</i> $\bar{3}$ <i>m</i>	0.40496
(Al) hp	β Al	25	20.5	A3	Mg	<i>hP</i> 2	<i>P6</i> ₃ / <i>mmc</i>	0.2693	...	0.4398	1.6331
(Am) rt	α Am	25	atm	A3'	Nd	<i>hP</i> 4	<i>P6</i> ₃ / <i>mmc</i>	0.34681	...	1.1241	2 x 1.621
(Am) ht1	β Am	>769	atm	A1	Cu	<i>cF</i> 4	<i>Fm</i> $\bar{3}$ <i>m</i>	0.4894
(Am) ht2	γ Am	>1077	atm	A2	W	<i>cI</i> 2	<i>Im</i> $\bar{3}$ <i>m</i>	?
(Am) hp	δ Am	25	>15	A20	U	<i>oS</i> 4	<i>Cmcm</i>	0.3063	0.5968	0.5169	...

Element phase	Common name	Temperature, °C	Pressure, GPa	Struktur- bericht designation	Prototype	Pearson symbol	Space group	Cell parameters, nm			Comment,c/a, or α or β
								a	b	c	
(Ar)	α Ar	<-189.2	atm	A1	Cu	cF4	Fm $\bar{3}m$	0.53109
(As)	As	25	atm	A7	As	hR6	R $\bar{3}m$	0.41319	$\alpha = 54.12^\circ$
(Au)	Au	25	atm	A1	Cu	cF4	Fm $\bar{3}m$	0.40782
(B) rhom1	β B	25	atm	...	B	hR423	R $\bar{3}m$	1.017	$\alpha = 65.12^\circ$
(Ba)	α Ba	25	atm	A2	W	cI2	I $m\bar{3}m$	0.50227
(Ba) hp1	β Ba	25	>5.33	A3	Mg	hP2	P6 ₃ /mmc	0.3901	...	0.6154	1.5775
(Ba) hp2	γ Ba	25	>23	?	?
(Be) rt	α Be	25	atm	A3	Mg	hP2	P6 ₃ /mmc	0.22859	...	0.35845	1.5681
(Be) ht	β Be	>1270	atm	A2	W	cI2	I $m\bar{3}m$	0.25515
(Be) hp	BeII	25	>28.3	hP*	...	0.4328	...	0.3416	0.7893
(Bi)	α Bi	25	atm	A7	As	hR6	R $\bar{3}m$	0.47460	$\alpha = 57.23^\circ$
(Bi) hp1	β Bi	25	>2.6	...	β Bi	mS4	C12/m1	0.6674	0.6117	0.3304	$\beta = 110.33^\circ$
(Bi) hp2	γ Bi	25	>3.0	mP4	P12 ₁ /m1	0.665	0.420	0.465	$\beta = 85.33^\circ$
(Bi) hp3	δ Bi	25	>4.3	?	?
(Bi) hp4	ζ Bi	25	>9.0	A2	W	cI2	I $m\bar{3}m$	0.3800
(Bk) rt	α Bk	25	atm	A3'	Nd	hP4	P6 ₃ /mmc	0.3416	...	1.1069	2 x 1.620
(Bk) ht	β Bk	>977	atm	A1	Cu	cF4	Fm $\bar{3}m$	0.4997
([Br ₂])	Br	<-7.25	atm	A14	[I ₂]	oS8	Cmce	0.668	0.449	0.874	...
(C) gra	C(graphite)	25	atm	A9	C	hP4	P6 ₃ /mmc	0.24612	...	0.6709	2.7258
(C) dia	C(diamond)	25	>60	A4	C	cF8	Fd $\bar{3}m$	0.35669

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								a	b	c	
(Ca) rt	α Ca	25	atm	A1	Cu	cF4	$Fm\bar{3}m$	0.55884
(Ca) ht	β Ca	>443	atm	A2	W	cI2	$Im\bar{3}m$	0.4480
(Ca) stab	γ Ca	25	>1.5	A3	Mg	hP2	$P6_3/mmc$
(Cd)	Cd	25	atm	A3	Mg	hP2	$P6_3/mmc$	0.29793	...	0.56196	1.8862
(Ce) ht1	α Ce	<-177	atm	A1	Cu	cF4	$Fm\bar{3}m$	0.485
(Ce) rt	β Ce	25	atm	A3'	Nd	hP4	$P6_3/mmc$	0.36810	...	1.1857	2 x 1.611
(Ce) ht1	γ Ce	>61	atm	A1	Cu	cF4	$Fm\bar{3}m$	0.51610
(Ce) ht2	δ Ce	>726	atm	A2	W	cI2	$Im\bar{3}m$	0.412
(Ce) hp	α' Ce	25	>5.4	A20	U	oS4	$Cmcm$	0.3049	0.5998	0.5215	...
(Cf) rt	α Cf	25	atm	A3'	Nd	hP4	$P6_3/mmc$	0.339	...	1.1015	2 x 1.625
(Cf) ht	β Cf	>590	atm	A1	Cu	cF4	$Fm\bar{3}m$?
([Cl ₂])	Cl	<-100.97	atm	A14	[I ₂]	oS8	$Cmce$	0.624	0.448	0.826	...
(Cm) rt	α Cm	25	atm	A3'	Nd	hP4	$P6_3/mmc$	0.3496	...	1.1331	2 x 1.621
(Cm) ht	β Cm	>1277	atm	A1	Cu	cF4	$Fm\bar{3}m$	0.4382
(Co) rt	ε Co	25	atm	A3	Mg	hP2	$P6_3/mmc$	0.25071	...	0.40686	1.6228
(Co) ht	α Co	>422	atm	A1	Cu	cF4	$Fm\bar{3}m$	0.35447
(Cr)	α Cr	25	atm	A2	W	cI2	$Im\bar{3}m$	0.28848
(Cr) hp	α' Cr	25	hp	...	Cr	tI2	$I4/mmm$	0.2882	...	0.2887	1.002
(Cs)	α Cs	25	atm	A2	W	cI2	$Im\bar{3}m$	0.6141
(Cs) hp1	β Cs	25	>2.37	A1	Cu	cF4	$Fm\bar{3}m$	0.6465
(Cs) hp2	β' Cs	25	>4.22	A1	Cu	cF4	$Fm\bar{3}m$	0.5800

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								a	b	c	
(Cs) hp3	γ Cs	25	>4.27	?
(Cu)	Cu	25	atm	A1	Cu	cF4	Fm $\bar{3}m$	0.36146
(Dy) lt	α 'Dy	<-187	atm	...	U	oS4	Cmcm	0.3595	0.6184	0.5678	...
(Dy) rt	α Dy	25	atm	A3	Mg	hP2	P6 ₃ /mmc	0.35915	...	0.56501	1.5732
(Dy) ht	β Dy	>1381	atm	A2	W	cI2	Im $\bar{3}m$	(0.398)
(Dy) hp	γ Dy	25	>7.5	C19	CdCl ₂	hR9	R $\bar{3}m$	0.3436	...	2.483	4.5 x 1.606
(Er)	Er	25	atm	A3	Mg	hP2	P6 ₃ /mmc	0.35592	...	0.55850	1.5692
(Es)	α Es	25	atm	A3'	Nd	hP4	P6 ₃ /mmc	?
(Es?)	β Es	?	atm	A1	Cu	cF4	Fm $\bar{3}m$?
(Eu)	Eu	25	atm	A2	W	cI2	Im $\bar{3}m$	0.45827
([F ₂]) lt2	α F	<-227.60	atm	...	[F ₂]	mS8	C12/c1	0.550	0.338	0.728	$\beta = 102.17^\circ$
([F ₂]) lt1	β F	<-219.67	atm	...	[F ₂]	cP64	Pm $\bar{3}n$	0.667
(Fe) rt	α Fe	25	atm	A2	W	cI2	Im $\bar{3}m$	0.28665
(Fe) ht	γ Fe	>912	atm	A1	Cu	cF4	Fm $\bar{3}m$	0.36467
(Fe) rt	δ Fe	>1394	atm	A2	W	cI2	Im $\bar{3}m$	0.29315
(Fe) hp	ε Fe	25	>13	A3	Mg	hP2	P6 ₃ /mmc	0.2468	...	0.396	1.603
(Ga)	α Ga	25	atm	A11	Ga	oS8	Cmce	0.45186	0.76570	0.45258	...
(Ga) lt	β Ga	25	>1.2	A6	In	tI2	I4/mmm	0.2808	...	0.4458	1.588
(Ga) hp	γ Ga	-53	>3.0	...	Ga	oS40	Cmcm	1.0593	1.3523	0.5203	...
(Gd) rt	α Gd	25	atm	A3	Mg	hP2	P6 ₃ /mmc	0.36336	...	0.57810	1.5910
(Gd) ht	β Gd	>1235	atm	A2	W	cI2	Im $\bar{3}m$	0.406

Element phase	Common name	Temperature, °C	Pressure, GPa	Struktur-bericht designation	Prototype	Pearson symbol	Space group	Cell parameters, nm			Comment, c/a, or α or β
								a	b	c	
(Gd) hp	γ Gd	25	>3.0	C19	Sm	<i>hR9</i>	$R\bar{3}m$	0.361	...	2.603	4.5 x 1.60
(Ge)	α Ge	25	atm	A4	C	<i>cF8</i>	$Fd\bar{3}m$	0.56574
(Ge) hp1	β Ge	25	>12	A5	Sn	<i>tI4</i>	$I4_1/AMD$	0.4884	...	0.2692	0.551
(Ge) hp2	γ Ge	25	>12 → atm	...	Ge	<i>tP12</i>	$P4_12_12$	0.593	...	0.698	1.18
(Ge) hp3	δ Ge	LT	>12	...	Si	<i>cI16</i>	$Ia\bar{3}$	0.692
([H ₂]) lt1	α H	<-271.9	atm	A1	Cu	<i>cF4</i>	$Fm\bar{3}m$	0.5338
([H ₂])	β H	<-259.34	atm	A3	Mg	<i>hP2</i>	$P6_3/mmc$	0.3776	...	0.6162	1.632
(He) lt2	He3	-269.67	0.163	A3	Mg	<i>hP2</i>	$P6_3/mmc$	0.3501	...	0.5721	1.634
(He) lt1	He4	-269.2	0.129	A3	Mg	<i>hP2</i>	$P6_3/mmc$	0.3470	...	0.5540	1.597
(Hf) rt	α Hf	25	atm	A3	Mg	<i>hP2</i>	$P6_3/mmc$	0.31946	...	0.50510	1.5811
(Hf) ht	β Hf	>1743	atm	A2	W	<i>cI2</i>	$Im\bar{3}m$	0.3610
(Hg)	α Hg	<-38.836	atm	A10	Hg	<i>hR3</i>	$R\bar{3}m$	0.3005	$\alpha = 70.53^\circ$
(Hg) hp	β Hg	<-194	hp	...	Hg	<i>tI2</i>	$I4/mmm$	0.3995	...	0.2825	0.707
(Ho)	α Ho	25	atm	A3	Mg	<i>hP2</i>	$P6_3/mmc$	0.35778	...	0.56178	1.5702
(Ho) hp	β Ho	25	>7.5	C19	Sm	<i>hR9</i>	$R\bar{3}m$	0.334	...	2.45	4.5 x 1.63
([I ₂])	I	25	atm	A14	[I ₂]	<i>oS8</i>	<i>Cmce</i>	0.72697	0.47903	0.97942	...
(In)	In	25	atm	A6	In	<i>tI2</i>	$I4/mmm$	0.3253	...	0.49470	1.5210
(Ir)	Ir	25	atm	A1	Cu	<i>cF4</i>	$Fm\bar{3}m$	0.38392
(K)	K	25	atm	A2	W	<i>cI2</i>	$Im\bar{3}m$	0.5321
(Kr)	Kr	<-157.385	atm	A1	Cu	<i>cF4</i>	$Fm\bar{3}m$	0.5810
(La) rt	α La	25	atm	A3'	Nd	<i>hP4</i>	$P6_3/mmc$	0.37740	...	1.2171	2 x 1.6125

Element phase	Common name	Temperature, °C	Pressure, GPa	Struktur- bericht designation	Prototype	Pearson symbol	Space group	Cell parameters, nm			Comment,c/a, or α or β
								<i>a</i>	<i>b</i>	<i>c</i>	
(La) ht1	βLa	>310	atm	A1	Cu	<i>cF</i> 4	<i>Fm</i> 3̄ <i>m</i>	0.5303
(La) ht2	γLa	>865	atm	A2	W	<i>cI</i> 2	<i>Im</i> 3̄ <i>m</i>	0.426
(La) hp	β'La	25	>2.0	A1	Cu	<i>cF</i> 4	<i>Fm</i> 3̄ <i>m</i>	0.517
(Li) lt	αLi	<-193	atm	A3	Mg	<i>hP</i> 2	<i>P</i> 6 ₃ /mmc	0.3111	...	0.5093	1.637
(Li) rt	βLi	25	atm	A2	W	<i>cI</i> 2	<i>Im</i> 3̄ <i>m</i>	0.35093
(Lu)	Lu	25	atm	A3	Mg	<i>hP</i> 2	<i>P</i> 6 ₃ /mmc	0.35052	...	0.55494	1.5832
(Mg)	Mg	25	atm	A3	Mg	<i>hP</i> 2	<i>P</i> 6 ₃ /mmc	0.32094	...	0.52107	1.6236
(Mn) rt	αMn	25	atm	A12	Mn	<i>cI</i> 58	<i>I</i> 4̄3 <i>m</i>	0.89126
(Mn) ht1	βMn	>727	atm	A13	Mn	<i>cP</i> 20	<i>P</i> 4 ₁ 32	0.63152
(Mn) ht3	γMn	>1100	atm	A1	Cu	<i>cF</i> 4	<i>Fm</i> 3̄ <i>m</i>	0.3860
(Mn) ht4	δMn	>1138	atm	A2	W	<i>cI</i> 2	<i>Im</i> 3̄ <i>m</i>	0.3080
(Mo)	Mo	25	atm	A2	W	<i>cI</i> 2	<i>Im</i> 3̄ <i>m</i>	0.31470
([N ₂]) lt2	αN	<-237.54	atm	...	[N ₂]	<i>cP</i> 8	<i>P</i> 2 ₁ 3	0.5661
([N ₂]) lt1	βN	<-210.004	atm	...	[N ₂]	<i>hP</i> 24	<i>P</i> 6 ₃ /mmc	0.4050	...	0.6604	1.631
([N ₂]) hp	γN	<-253	>3.3	...	[N ₂]	<i>tP</i> 4	<i>P</i> 4 ₂ /nnm	0.3957	...	0.5109	1.291
(Na) lt	αNa	<-233	atm	A3	Mg	<i>hP</i> 2	<i>P</i> 6 ₃ /mmc	0.3767	...	0.6154	1.634
(Na) rt	βNa	25	atm	A2	W	<i>cI</i> 2	<i>Im</i> 3̄ <i>m</i>	0.42906
(Nb)	Nb	25	atm	A2	W	<i>cI</i> 2	<i>Im</i> 3̄ <i>m</i>	0.33004
(Nd) rt	αNd	25	atm	A3'	Nd	<i>hP</i> 4	<i>P</i> 6 ₃ /mmc	0.36582	...	1.17966	2 x 1.6124
(Nd) ht	βNd	>863	atm	A2	W	<i>cI</i> 2	<i>Im</i> 3̄ <i>m</i>	0.413
(Nd) hp	γNd	25	>5.0	A1	Cu	<i>cF</i> 4	<i>Fm</i> 3̄ <i>m</i>	0.480



ASM Alloy Phase Diagram Database™

Element phase	Common name	Temperature, °C	Pressure, GPa	Struktur-bericht designation	Prototype	Pearson symbol	Space group	Cell parameters, nm			Comment, c/a, or α or β
								a	b	c	
(Ne)	Ne	<-248.587	atm	A1	Cu	cF4	Fm $\bar{3}m$	0.4462
(Ni)	Ni	25	atm	A1	Cu	cF4	Fm $\bar{3}m$	0.35240
(Np) rt	α Np	25	atm	A_c	Np	<i>o</i> P8	Pnma	0.6663	0.4723	0.4887	...
(Np) ht1	β Np	>280	atm	A_d	Np	<i>t</i> P4	P4/nmm	0.4883	...	0.3389	0.694
(Np) ht2	γ Np	>576	atm	A2	W	<i>c</i> I2	Im $\bar{3}m$	0.352
([O ₂]) lt3	α O	<-249.283	atm	...	[O ₂]	<i>m</i> S4	C12/m1	0.5403	0.3429	0.5086	$\beta = 132.53^\circ$
([O ₂])	β O	<-229.349	atm	...	[O ₂]	<i>h</i> R6	R $\bar{3}m$	0.4210	$\alpha = 46.27^\circ$
([O ₂]) lt1	γ O	<-218.789	atm	...	[O ₂]	<i>c</i> P64	Pm $\bar{3}n$	0.683
(Os)	Os	25	atm	A3	Mg	<i>h</i> P2	P6 ₃ /mmc	0.27341	...	0.43198	1.5800
(P) whi rt	α P(white)	25	atm	...	Mn	<i>c</i> I58	I $\bar{4}3m$	0.718
(P) bla	P(black)	25	atm	A17	P	<i>o</i> S8	Cmce	0.33136	1.0478	0.43763	...
(Pa) rt	α Pa	25	atm	A_a	Pa	<i>t</i> I2	I4/mmm	0.3921	...	0.3235	0.825
(Pa) ht	β Pa	>1170	atm	A2	W	<i>c</i> I2	Im $\bar{3}m$	0.381
(Pb)	α Pb	25	atm	A1	Cu	cF4	Fm $\bar{3}m$	0.49502
(Pb) hp	β Pb	25	>10.3	A3	Mg	<i>h</i> P2	P6 ₃ /mmc	0.3265	...	0.5387	1.650
(Pd)	Pd	25	atm	A1	Cu	cF4	Fm $\bar{3}m$	0.38903
(Pm) rt	α Pm	25	atm	A3'	Nd	<i>h</i> P4	P6 ₃ /mmc	0.365	...	1.165	2 x 1.60
(Po) rt	α Po	25	atm	A_h	Po	<i>c</i> P1	Pm $\bar{3}m$	0.3366
(Po) ht	β Po	>54	atm	A_i	Po	<i>h</i> R3	R3m	0.3373	$\alpha = 98.08^\circ$
(Pr) rt	α Pr	25	atm	A3'	Nd	<i>h</i> P4	P6 ₃ /mmc	0.36721	...	1.18326	2 x 1.6111
(Pr) ht	β Pr	>795	atm	A2	W	<i>c</i> I2	Im $\bar{3}m$	0.413

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(Pr) hp	γ Pr	25	>4.0	A1	Cu	cF4	Fm $\bar{3}$ m	0.488
(Pt)	Pt	25	atm	A1	Cu	cF4	Fm $\bar{3}$ m	0.39236
(Pu) rt	α Pu	25	atm	...	Pu	mP16	P12 ₁ /m1	0.6183	0.4822	1.0963	$\beta = 101.97^\circ$
(Pu) ht1	β Pu	>125	atm	...	Pu	mS34	C12/m1	0.9284	1.0463	0.7859	$\beta = 92.13^\circ$
(Pu) ht2	γ Pu	>215	atm	...	Pu	oF8	Fddd	0.31587	0.57682	1.0162	...
(Pu) ht3	δ Pu	>320	atm	A1	Cu	cF4	Fm $\bar{3}$ m	0.46371
(Pu) ht4	δ 'Pu	>463	atm	A6	In	tI2	I4/mmm	0.33261	...	0.44630	1.3418
(Pu) ht5	ε Pu	>483	atm	A2	W	cI2	Im $\bar{3}$ m	0.36343
(Ra)	Ra	25	atm	A2	W	cI2	Im $\bar{3}$ m	0.5148
(Rb)	α Rb	25	atm	A2	W	cI2	Im $\bar{3}$ m	0.5705
(Rb) hp1	β Rb	25	>1.08	?
(Rb) hp2	γ Rb	25	>2.05	?
(Re)	Re	25	atm	A3	Mg	hP2	P6 ₃ /mmc	0.27609	...	0.4458	1.6145
(Rh)	Rh	25	atm	A1	Cu	cF4	Fm $\bar{3}$ m	0.38032
(Ru)	Ru	25	atm	A3	Mg	hP2	P6 ₃ /mmc	0.27058	...	0.42816	1.5824
(S) 8 α rt	α S	25	atm	A16	S	oF128	Fddd	1.0464	1.28660	2.44860	...
(S) 8 β ht	β S	>95.5	atm	...	S	mP64	P2 ₁ /c1	1.102	1.096	1.090	$\beta = 96.7^\circ$
(Sb)	α Sb	25	atm	A7	As	hR6	R $\bar{3}$ m	0.45067	$\alpha = 57.11^\circ$
(Sb) hp1	β Sb	25	>5.0	A _h	Po	cP1	Pm $\bar{3}$ m	0.2992
(Sb) hp2	δ Sb	25	>14.0	mP3	?	0.556	0.404	0.422	$\beta = 86.0^\circ$
(Sc) rt	α Sc	25	atm	A3	Mg	hP2	P6 ₃ /mmc	0.33088	...	0.52680	1.5921



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(Sc) ht	β Sc	>1337	atm	A2	W	cI2	$I\bar{m}\bar{3}m$	0.373
(Se)	γ Se	25	atm	A8	Se	hP3	P3 ₁ 21	0.43659	...	0.49537	1.1346
(Si)	α Si	25	atm	A4	C	cF8	$Fd\bar{3}m$	0.54306
(Si) hp1	β Si	25	>9.5	A5	Sn	tI4	$I4_1/AMD$	0.4686	...	0.2585	0.552
(Si) hp2	γ Si	25	>16.0	...	Si	cI16	$Ia\bar{3}$	0.6636
(Si) hp3	δ Si	25	>16 → atm	A3'	Nd	hP4	P6 ₃ /mmc	0.380	...	0.628	1.653
(Sm) rt	α Sm	25	atm	C19	Sm	hR9	$R\bar{3}m$	0.36290	...	2.6207	4.5 x 1.6048
(Sm) ht1	β Sm	>734	atm	A3	Mg	hP2	P6 ₃ /mmc	0.36630	...	0.58448	1.5956
(Sm) ht2	γ Sm	>922	atm	A2	W	cI2	$I\bar{m}\bar{3}m$?
(Sm) hp	δ Sm	25	>4.0	A3'	La	hP4	P6 ₃ /mmc	0.3618	...	1.166	2 x 1.611
(Sn) lt	α Sn	<13	atm	A4	C	cF8	$Fd\bar{3}m$	0.64892
(Sn) rt	β Sn	25	atm	A5	Sn	tI4	$I4_1/AMD$	0.58318	...	0.31818	0.5456
(Sn) hp	γ Sn	25	>9.0	...	Pa	tI2	$I4/mmm$	0.370	...	0.337	0.91
(Sr) rt	α Sr	25	atm	A1	Cu	cF4	$Fm\bar{3}m$	0.6084
(Sr) ht	β Sr	>547	atm	A2	W	cI2	$I\bar{m}\bar{3}m$	0.487
(Sr) hp	β' Sr	25	>3.5	A2	W	cI2	$I\bar{m}\bar{3}m$	0.4437
(Ta)	Ta	25	atm	A2	W	cI2	$I\bar{m}\bar{3}m$	0.33030
(Tb) lt	α Tb	<-53	atm	...	Dy	oS4	Cmcm	0.3605	0.6244	0.5706	...
(Tb) rt	α' Tb	25	atm	A3	Mg	hP2	P6 ₃ /mmc	0.36055	...	0.56966	1.5800
(Tb) ht	β Tb	>1289	atm	A2	W	cI2	$I\bar{m}\bar{3}m$	(0.402)
(Tb) hp	γ Tb	25	>6.0	C19	Sm	hR9	$R\bar{3}m$	0.341	...	2.45	4.5 x 1.60



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(Te)	α Te	25	atm	A8	Se	hP3	P3 ₁ 21	0.44566	...	0.59264	1.3298
(Te) hp1	β Te	25	>2.0	A7	As	hR6	R ₃ m	0.469	$\alpha = 53.30^\circ$
(Te) hp2	γ Te	25	>7.0	A_i	Po	hR3	R ₃ m	0.3002	$\alpha = 103.3^\circ$
(Th) rt	α Th	25	atm	A1	Cu	cF4	Fm ₃ m	0.50842
(Th) ht	β Th	>1360	atm	A2	W	cI2	I _m ₃ m	0.411
(Ti) rt	α Ti	25	atm	A3	Mg	hP2	P6 ₃ /mmc	0.29506	...	0.46835	1.5873
(Ti) ht	β Ti	>882	atm	A2	W	cI2	I _m ₃ m	0.33065
(Ti) hp1	ω Ti	25	hp \rightarrow atm	...	Ti	hP3	P6/mmm	0.4625	...	0.2813	0.6082
(Tl) rt	α Tl	25	atm	A3	Mg	hP2	P6 ₃ /mmc	0.34566	...	0.55248	1.5983
(Tl) ht	β Tl	>230	atm	A2	W	cI2	I _m ₃ m	0.3879
(Tl) hp	γ Tl	25	hp	A1	Cu	cF4	Fm ₃ m	?
(Tm)	Tm	25	atm	A3	Mg	hP2	P6 ₃ /mmc	0.35375	...	0.55540	1.5700
(U) rt	α U	25	atm	A20	U	oS4	Cmcm	0.28537	0.58695	0.49548	...
(U) ht1	β U	>668	atm	A_b	U	tP30	P4 ₂ /nnm	1.0759	...	0.5656	0.526
(U) ht2	γ U	>776	atm	A2	W	cI2	I _m ₃ m	0.3524
(V)	V	25	atm	A2	W	cI2	I _m ₃ m	0.30240
(W)	W	25	atm	A2	W	cI2	I _m ₃ m	0.31652
(Xe)	Xe	<-111.758	atm	A1	Cu	cF4	Fm ₃ m	0.6350
(Y) rt	α Y	25	atm	A3	Mg	hP2	P6 ₃ /mmc	0.36482	...	0.57318	1.5711
(Y) ht	β Y	>1478	atm	A2	W	cI2	I _m ₃ m	(0.407)
(Yb) stab	α Yb	<-3	atm	A3	Mg	hP2	P6 ₃ /mmc	0.38799	...	0.63859	1.6459



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(Yb) rt	β Yb	25	atm	A1	Cu	<i>cF</i> 4	<i>Fm</i> $\bar{3}$ <i>m</i>	0.54848
(Yb) ht	γ Yb	>795	atm	A2	W	<i>cI</i> 2	<i>Im</i> $\bar{3}$ <i>m</i>	0.444
(Zn)	α Zr	25	atm	A3	Mg	<i>hP</i> 2	<i>P6</i> ₃ / <i>mmc</i>	0.32316	...	0.51475	1.5929
(Zr) ht	β Zr	>863	atm	A2	W	<i>cI</i> 2	<i>Im</i> 3 <i>m</i>	0.36090
(Zr) rt	ω Zr	25	hp \rightarrow atm	...	ω Ti	<i>hP</i> 2	<i>P6</i> / <i>mmm</i>	0.5036	...	0.3109	0.617

Note: Values in parentheses are estimated.