

substance: (transition metal)-V-VI compounds**property: crystallographic and physical data for pyrite-type derivatives**

(RT values). Seebeck coefficient from [62H].

Compound	<i>a</i> [Å]	<i>d_X</i> [g cm ⁻³]	Ref.	<i>E_g</i> [eV]	Ref.	10 ⁶ · <i>χ_m</i> [cm ³ mol ⁻¹]	Ref.	<i>S</i> [μV K ⁻¹]
CoPS (tetr.)	5.414 (<i>c</i> = 5.430)	5.08	74N	0.4	59H			
	5.422 (h?)		63H					
CoAsS (r) (C _{2v} ⁵ -Pca2 ₁) pseudocell*)	5.582		65G	0.75	59H			- 50
	5.578	6.35	63H	0.60	76A			
	5.572		65K					
CoSbS (p) (T ⁴ -P2 ₁ 3)	5.844	7.08	75H			0	75H	
CoPSe (p) (T _h ⁶ -Pa3)	5.63	6.29	75H					
CoAsSe (p) (T _h ⁶ -Pa3 or T ⁴ -P2 ₁ 3)	5.76**)		75H					
			63H					
RhPS	5.640	6.14	63H					+ 10
RhAsS	5.780	7.22	63H					+ 50
RhSbS (ordered)	6.027	7.79	63H					- 300
RhBiS (T ⁴ -P2 ₁ 3)	6.138	9.88	63H					- 350
RhPSe (ordered)	5.795	7.26	63H			- 15	62H	
RhAsSe	5.934	8.16	63H					+ 100
RhSbSe (ordered)	6.176	8.56	63H					+ 50
RhBiSe (ordered)	6.283	10.47	63H					- 250
RhAsTe	6.165	8.66	63H					- 100
RhSbTe	6.392	8.96	63H	> 0.15	80H	- 70	62H	- 50
RhBiTe	6.504	10.61	63H					- 30
IrPS	5.650	9.40	63H					
IrAsS	5.791	10.23	63H			- 66	62H	+ 10
IrSbS (ordered)	6.036	10.45	63H					- 1000
IrBiS (ordered)	6.143	12.41	63H			- 78	62H	- 70
IrPSe (ordered)	5.798	10.30	63H					
IrAsSe	5.940	10.97	63H					+ 10
IrSbSe (ordered)	6.184	11.04	63H					
IrBiSe (ordered)	6.290	12.82	63H					- 50
IrAsTe	6.164	11.19	63H					+ 60
IrSbTe (ordered)	6.397	11.20	63H					
IrBiTe	6.500	12.79	63H					- 50

*) (h): annealing at 550°C completely disorders As and S [65G]. Most mineral specimens are partially ordered [65G, 82B]. **) Stabilized by impurities?

References:

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