

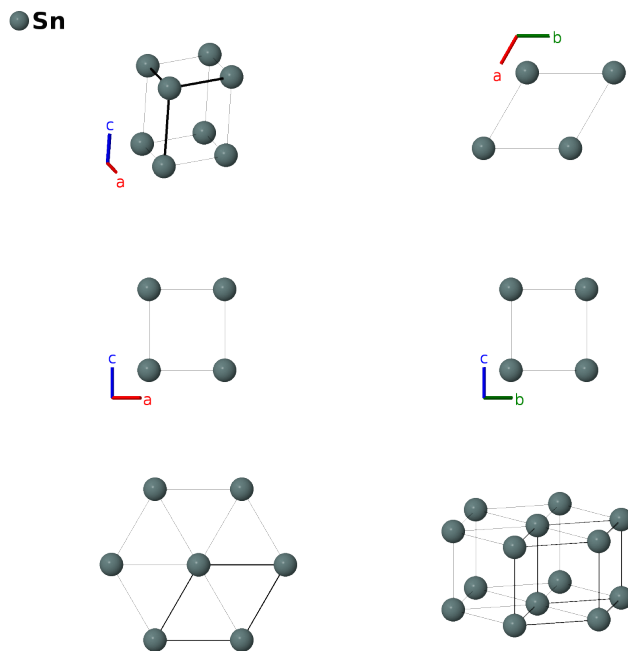
Simple Hexagonal ($\text{HgSn}_{6-10} A_f$) Structure: A_hP1_191_a-001

This structure originally had the label A_hP1_191_a. Calls to that address will be redirected here.

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<https://afLOW.org/p/3V7H>

https://afLOW.org/p/A_hP1_191_a-001



Prototype	HgSn_{6-10}
AFLOW prototype label	A_hP1_191_a-001
<i>Strukturbericht</i> designation	A_f
ICSD	639211
Pearson symbol	hP1
Space group number	191
Space group symbol	$P6/mmm$
AFLOW prototype command	<code>afLOW --proto=A_hP1_191_a-001 --params=a, c/a</code>

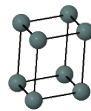
Other compounds with this structure

Si (metastable), disordered phases of BiIn, CdSn_{19} , In_7Sb_3 , InSb

- Unlike the simple cubic lattice, there are no elements which take this structure as the ground state. There is a metastable silicon phase with this structure. The prototype state is a mercury-tin alloy. Thus the atom type “M” represents an average of Hg and Sn atoms.

Hexagonal primitive vectors

$$\begin{aligned}\mathbf{a}_1 &= \frac{1}{2}a \hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a \hat{\mathbf{y}} \\ \mathbf{a}_2 &= \frac{1}{2}a \hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a \hat{\mathbf{y}} \\ \mathbf{a}_3 &= c \hat{\mathbf{z}}\end{aligned}$$



Basis vectors

	Lattice coordinates		Cartesian coordinates	Wyckoff position	Atom type
\mathbf{B}_1	=	0	=	0	(1a) Sn I

References

- [1] G. V. Raynor and J. A. Lee, *The tin-rich intermediate phases in the alloys of tin with cadmium, indium and mercury*, *Acta Metall.* **2**, 616–620 (1954), doi:10.1016/0001-6160(54)90197-2.

Found in

- [1] P. Villars and L. Calvert, *Pearson's Handbook of Crystallographic Data for Intermetallic Phases* (ASM International, Materials Park, OH, 1991), 2nd edn.