

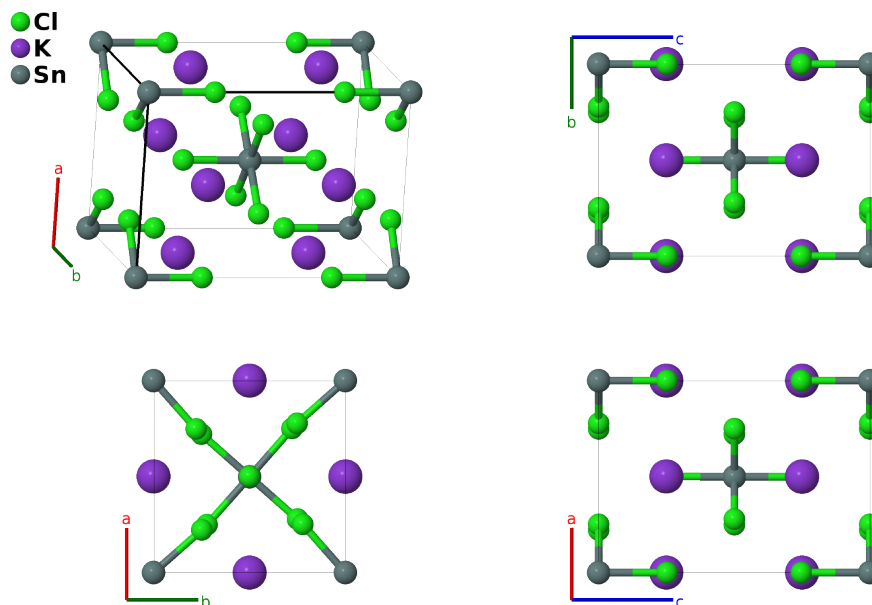
Phase II K_2SnCl_6 Structure: A6B2C_tP18_128_eh_d_a-001

This structure originally had the label A6B2C.tP18.128.eh.d.a. Calls to that address will be redirected here.

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<https://aflow.org/p/AD9M>

https://aflow.org/p/A6B2C_tP18_128_eh_d_a-001

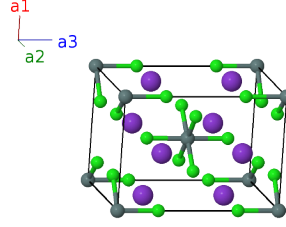


Prototype	$\text{Cl}_6\text{K}_2\text{Sn}$
AFLOW prototype label	A6B2C_tP18_128_eh_d_a-001
ICSD	1669
Pearson symbol	tP18
Space group number	128
Space group symbol	$P4/mnc$
AFLOW prototype command	<code>aflow --proto=A6B2C_tP18_128_eh_d_a-001 --params=a, c/a, z3, x4, y4</code>

- K_2SnCl_6 is found in three forms, depending on the temperature:
 - Below 255K it is in the monoclinic Phase I structure.
 - In the range 255-261K it is in the tetragonal Phase II structure (this structure).
 - Above 261K it transforms to Phase III, which has the K_2PtCl_6 ($J1_1$) structure.
- Data for Phase II was taken at 265K.

Simple Tetragonal primitive vectors

$$\begin{aligned}\mathbf{a}_1 &= a \hat{\mathbf{x}} \\ \mathbf{a}_2 &= a \hat{\mathbf{y}} \\ \mathbf{a}_3 &= c \hat{\mathbf{z}}\end{aligned}$$



Basis vectors

	Lattice coordinates		Cartesian coordinates	Wyckoff position	Atom type
\mathbf{B}_1	0	$=$	0	(2a)	Sn I
\mathbf{B}_2	$\frac{1}{2} \mathbf{a}_1 + \frac{1}{2} \mathbf{a}_2 + \frac{1}{2} \mathbf{a}_3$	$=$	$\frac{1}{2} a \hat{\mathbf{x}} + \frac{1}{2} a \hat{\mathbf{y}} + \frac{1}{2} c \hat{\mathbf{z}}$	(2a)	Sn I
\mathbf{B}_3	$\frac{1}{2} \mathbf{a}_2 + \frac{1}{4} \mathbf{a}_3$	$=$	$\frac{1}{2} a \hat{\mathbf{y}} + \frac{1}{4} c \hat{\mathbf{z}}$	(4d)	K I
\mathbf{B}_4	$\frac{1}{2} \mathbf{a}_1 + \frac{1}{4} \mathbf{a}_3$	$=$	$\frac{1}{2} a \hat{\mathbf{x}} + \frac{1}{4} c \hat{\mathbf{z}}$	(4d)	K I
\mathbf{B}_5	$\frac{1}{2} \mathbf{a}_2 + \frac{3}{4} \mathbf{a}_3$	$=$	$\frac{1}{2} a \hat{\mathbf{y}} + \frac{3}{4} c \hat{\mathbf{z}}$	(4d)	K I
\mathbf{B}_6	$\frac{1}{2} \mathbf{a}_1 + \frac{3}{4} \mathbf{a}_3$	$=$	$\frac{1}{2} a \hat{\mathbf{x}} + \frac{3}{4} c \hat{\mathbf{z}}$	(4d)	K I
\mathbf{B}_7	$z_3 \mathbf{a}_3$	$=$	$cz_3 \hat{\mathbf{z}}$	(4e)	Cl I
\mathbf{B}_8	$\frac{1}{2} \mathbf{a}_1 + \frac{1}{2} \mathbf{a}_2 - (z_3 - \frac{1}{2}) \mathbf{a}_3$	$=$	$\frac{1}{2} a \hat{\mathbf{x}} + \frac{1}{2} a \hat{\mathbf{y}} - c (z_3 - \frac{1}{2}) \hat{\mathbf{z}}$	(4e)	Cl I
\mathbf{B}_9	$-z_3 \mathbf{a}_3$	$=$	$-cz_3 \hat{\mathbf{z}}$	(4e)	Cl I
\mathbf{B}_{10}	$\frac{1}{2} \mathbf{a}_1 + \frac{1}{2} \mathbf{a}_2 + (z_3 + \frac{1}{2}) \mathbf{a}_3$	$=$	$\frac{1}{2} a \hat{\mathbf{x}} + \frac{1}{2} a \hat{\mathbf{y}} + c (z_3 + \frac{1}{2}) \hat{\mathbf{z}}$	(4e)	Cl I
\mathbf{B}_{11}	$x_4 \mathbf{a}_1 + y_4 \mathbf{a}_2$	$=$	$ax_4 \hat{\mathbf{x}} + ay_4 \hat{\mathbf{y}}$	(8h)	Cl II
\mathbf{B}_{12}	$-x_4 \mathbf{a}_1 - y_4 \mathbf{a}_2$	$=$	$-ax_4 \hat{\mathbf{x}} - ay_4 \hat{\mathbf{y}}$	(8h)	Cl II
\mathbf{B}_{13}	$-y_4 \mathbf{a}_1 + x_4 \mathbf{a}_2$	$=$	$-ay_4 \hat{\mathbf{x}} + ax_4 \hat{\mathbf{y}}$	(8h)	Cl II
\mathbf{B}_{14}	$y_4 \mathbf{a}_1 - x_4 \mathbf{a}_2$	$=$	$ay_4 \hat{\mathbf{x}} - ax_4 \hat{\mathbf{y}}$	(8h)	Cl II
\mathbf{B}_{15}	$-(x_4 - \frac{1}{2}) \mathbf{a}_1 + (y_4 + \frac{1}{2}) \mathbf{a}_2 + \frac{1}{2} \mathbf{a}_3$	$=$	$-a (x_4 - \frac{1}{2}) \hat{\mathbf{x}} + a (y_4 + \frac{1}{2}) \hat{\mathbf{y}} + \frac{1}{2} c \hat{\mathbf{z}}$	(8h)	Cl II
\mathbf{B}_{16}	$(x_4 + \frac{1}{2}) \mathbf{a}_1 - (y_4 - \frac{1}{2}) \mathbf{a}_2 + \frac{1}{2} \mathbf{a}_3$	$=$	$a (x_4 + \frac{1}{2}) \hat{\mathbf{x}} - a (y_4 - \frac{1}{2}) \hat{\mathbf{y}} + \frac{1}{2} c \hat{\mathbf{z}}$	(8h)	Cl II
\mathbf{B}_{17}	$(y_4 + \frac{1}{2}) \mathbf{a}_1 + (x_4 + \frac{1}{2}) \mathbf{a}_2 + \frac{1}{2} \mathbf{a}_3$	$=$	$a (y_4 + \frac{1}{2}) \hat{\mathbf{x}} + a (x_4 + \frac{1}{2}) \hat{\mathbf{y}} + \frac{1}{2} c \hat{\mathbf{z}}$	(8h)	Cl II
\mathbf{B}_{18}	$-(y_4 - \frac{1}{2}) \mathbf{a}_1 - (x_4 - \frac{1}{2}) \mathbf{a}_2 + \frac{1}{2} \mathbf{a}_3$	$=$	$-a (y_4 - \frac{1}{2}) \hat{\mathbf{x}} - a (x_4 - \frac{1}{2}) \hat{\mathbf{y}} + \frac{1}{2} c \hat{\mathbf{z}}$	(8h)	Cl II

References

- [1] H. Boysen and A. W. Hewat, *A neutron powder investigation of the structural changes in $K_2\text{SnCl}_6$* , Acta Crystallogr. Sect. B **34**, 1412–1418 (1978), doi:10.1107/S0567740878005816.

Found in

- [1] P. Villars and K. Cenzual, *Pearson's Crystal Data – Crystal Structure Database for Inorganic Compounds* (2013). ASM International.