

SbI₃S₂₄ Structure:

A3B24C_hR28_160_b_2b3c_a-001

This structure previously used the label(s) **A3B24C_hR28_160_b_2b3c_a**. Calls to these addresses will be redirected here.

Cite this page as:

- H. Eckert, S. Divilov, M. J. Mehl, D. Hicks, A. C. Zettel, M. Esters, X. Campilongo, and S. Curtarolo, *The AFLOW Library of Crystallographic Prototypes: Part 4*, Comput. Mater. Sci. **240**, 112988 (2024), doi: 10.1016/j.commatsci.2024.112988.
- D. Hicks, M. J. Mehl, M. Esters, C. Oses, O. Levy, G. L. W. Hart, C. Toher, and S. Curtarolo, *The AFLOW Library of Crystallographic Prototypes: Part 3*, Comput. Mater. Sci. **199**, 110450 (2021), doi: 10.1016/j.commatsci.2021.110450.

<https://aflow.org/prototype-encyclopedia/8TMQ>

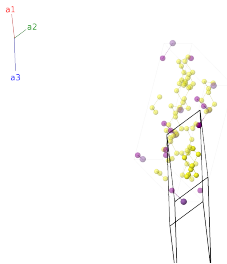
https://aflow.org/prototype-encyclopedia/A3B24C_hR28_160_b_2b3c_a-001

Prototype	I ₃ S ₂₄ Sb
AFLOW prototype label	A3B24C_hR28_160_b_2b3c_a-001
ICSD	14200
CCDC	1595599
Pearson symbol	hR28
Space group number	160
Space group symbol	<i>R</i> 3 <i>m</i>
AFLOW prototype command	<code>aflow --proto=A3B24C_hR28_160_b_2b3c_a-001</code> <code>--params=a, c/a, x₁, x₂, z₂, x₃, z₃, x₄, z₄, x₅, y₅, z₅, x₆, y₆, z₆, x₇, y₇, z₇</code>

- Since there are three S₈ molecules in this structure, (Bjorvatten, 1963) refer to it as SbI₃:3S₈.
- Space group *R*3*m* #160 does not fix the zero of the *z*-axis. Here it is set to coincide with the plane of the iodine atoms.

Rhombohedral primitive vectors

$$\begin{aligned}\mathbf{a}_1 &= \frac{1}{2}a \hat{\mathbf{x}} - \frac{\sqrt{3}}{6}a \hat{\mathbf{y}} + \frac{1}{3}c \hat{\mathbf{z}} \\ \mathbf{a}_2 &= \frac{1}{\sqrt{3}}a \hat{\mathbf{y}} + \frac{1}{3}c \hat{\mathbf{z}} \\ \mathbf{a}_3 &= -\frac{1}{2}a \hat{\mathbf{x}} - \frac{\sqrt{3}}{6}a \hat{\mathbf{y}} + \frac{1}{3}c \hat{\mathbf{z}}\end{aligned}$$



Basis vectors

	Lattice coordinates		Cartesian coordinates	Wyckoff position	Atom type
B₁	=	$x_1 \mathbf{a}_1 + x_1 \mathbf{a}_2 + x_1 \mathbf{a}_3$	=	$cx_1 \hat{\mathbf{z}}$	(1a) Sb I
B₂	=	$x_2 \mathbf{a}_1 + x_2 \mathbf{a}_2 + z_2 \mathbf{a}_3$	=	$\frac{1}{2}a(x_2 - z_2) \hat{\mathbf{x}} + \frac{\sqrt{3}}{6}a(x_2 - z_2) \hat{\mathbf{y}} + \frac{1}{3}c(2x_2 + z_2) \hat{\mathbf{z}}$	(3b) I I

$$\mathbf{B}_{28} = x_7 \mathbf{a}_1 + z_7 \mathbf{a}_2 + y_7 \mathbf{a}_3 = \frac{1}{2}a(x_7 - y_7) \hat{\mathbf{x}} - \frac{\sqrt{3}}{6}a(x_7 + y_7 - 2z_7) \hat{\mathbf{y}} + \frac{1}{3}c(x_7 + y_7 + z_7) \hat{\mathbf{z}} \quad (6c) \quad \text{S V}$$

References

- [1] T. Bjorvatten, O. Hassel, and A. Lindheim, *Crystal Structure of the Addition Compound $\text{SbI}_3 \cdot 3\text{S}_8$* , Acta Chem. Scand. **17**, 689–702 (1963), doi:10.3891/acta.chem.scand.17-0689.

First cited in

- D. Hicks, M. J. Mehl, M. Esters, C. Oses, O. Levy, G. L. W. Hart, C. Toher, and S. Curtarolo, *The AFLOW Library of Crystallographic Prototypes: Part 3*, Comput. Mater. Sci. **199**, 110450 (2021), doi: 10.1016/j.commatsci.2021.110450.

I
S
Sb

