

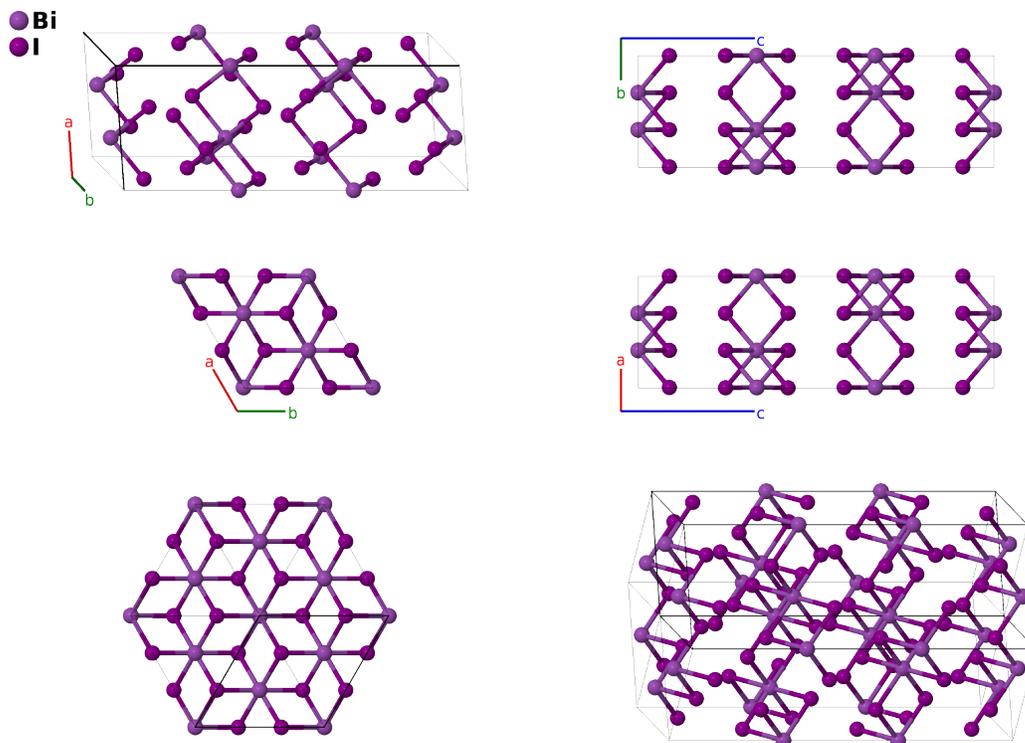
BiI₃ (*D*0₅) Structure: AB3_hR8_148_c_f-001

This structure originally had the label AB3_hR8_148_c_f. Calls to that address will be redirected here.

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<https://aflow.org/p/TZC0>

https://aflow.org/p/AB3_hR8_148_c_f-001



Prototype	BiI ₃
AFLOW prototype label	AB3_hR8_148_c_f-001
<i>Strukturbericht</i> designation	<i>D</i> 0 ₅
ICSD	36182
Pearson symbol	hR8
Space group number	148
Space group symbol	<i>R</i> $\bar{3}$
AFLOW prototype command	<code>aflow --proto=AB3_hR8_148_c_f-001 --params=a, c/a, x₁, x₂, y₂, z₂</code>

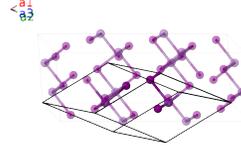
Other compounds with this structure

AsI₃, CrBr₃, CrCl₃, CrI₃, FeCl₃, RhBr₃, SbI₃, VCl₃, VI₃

- Hexagonal settings of this structure can be obtained with the option `--hex`.

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
\mathbf{B}_1	$= x_1 \mathbf{a}_1 + x_1 \mathbf{a}_2 + x_1 \mathbf{a}_3$	$=$	$cx_1 \hat{\mathbf{z}}$	(2c)	Bi I
\mathbf{B}_2	$= -x_1 \mathbf{a}_1 - x_1 \mathbf{a}_2 - x_1 \mathbf{a}_3$	$=$	$-cx_1 \hat{\mathbf{z}}$	(2c)	Bi I
\mathbf{B}_3	$= x_2 \mathbf{a}_1 + y_2 \mathbf{a}_2 + z_2 \mathbf{a}_3$	$=$	$\frac{1}{2}a(x_2 - z_2) \hat{\mathbf{x}} - \frac{\sqrt{3}}{6}a(x_2 - 2y_2 + z_2) \hat{\mathbf{y}} + \frac{1}{3}c(x_2 + y_2 + z_2) \hat{\mathbf{z}}$	(6f)	I I
\mathbf{B}_4	$= z_2 \mathbf{a}_1 + x_2 \mathbf{a}_2 + y_2 \mathbf{a}_3$	$=$	$-\frac{1}{2}a(y_2 - z_2) \hat{\mathbf{x}} + \frac{\sqrt{3}}{6}a(2x_2 - y_2 - z_2) \hat{\mathbf{y}} + \frac{1}{3}c(x_2 + y_2 + z_2) \hat{\mathbf{z}}$	(6f)	I I
\mathbf{B}_5	$= y_2 \mathbf{a}_1 + z_2 \mathbf{a}_2 + x_2 \mathbf{a}_3$	$=$	$-\frac{1}{2}a(x_2 - y_2) \hat{\mathbf{x}} - \frac{\sqrt{3}}{6}a(x_2 + y_2 - 2z_2) \hat{\mathbf{y}} + \frac{1}{3}c(x_2 + y_2 + z_2) \hat{\mathbf{z}}$	(6f)	I I
\mathbf{B}_6	$= -x_2 \mathbf{a}_1 - y_2 \mathbf{a}_2 - z_2 \mathbf{a}_3$	$=$	$-\frac{1}{2}a(x_2 - z_2) \hat{\mathbf{x}} + \frac{\sqrt{3}}{6}a(x_2 - 2y_2 + z_2) \hat{\mathbf{y}} - \frac{1}{3}c(x_2 + y_2 + z_2) \hat{\mathbf{z}}$	(6f)	I I
\mathbf{B}_7	$= -z_2 \mathbf{a}_1 - x_2 \mathbf{a}_2 - y_2 \mathbf{a}_3$	$=$	$\frac{1}{2}a(y_2 - z_2) \hat{\mathbf{x}} - \frac{\sqrt{3}}{6}a(2x_2 - y_2 - z_2) \hat{\mathbf{y}} - \frac{1}{3}c(x_2 + y_2 + z_2) \hat{\mathbf{z}}$	(6f)	I I
\mathbf{B}_8	$= -y_2 \mathbf{a}_1 - z_2 \mathbf{a}_2 - x_2 \mathbf{a}_3$	$=$	$\frac{1}{2}a(x_2 - y_2) \hat{\mathbf{x}} + \frac{\sqrt{3}}{6}a(x_2 + y_2 - 2z_2) \hat{\mathbf{y}} - \frac{1}{3}c(x_2 + y_2 + z_2) \hat{\mathbf{z}}$	(6f)	I I

References

- [1] H. Br, IX. *Die Kristallstruktur der Trijodide von Arsen, Antimon und Wismut*, Z. Kristallogr. **74**, 67–72 (1930), doi:10.1524/zkri.1930.74.1.67.

Found in

- [1] C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).