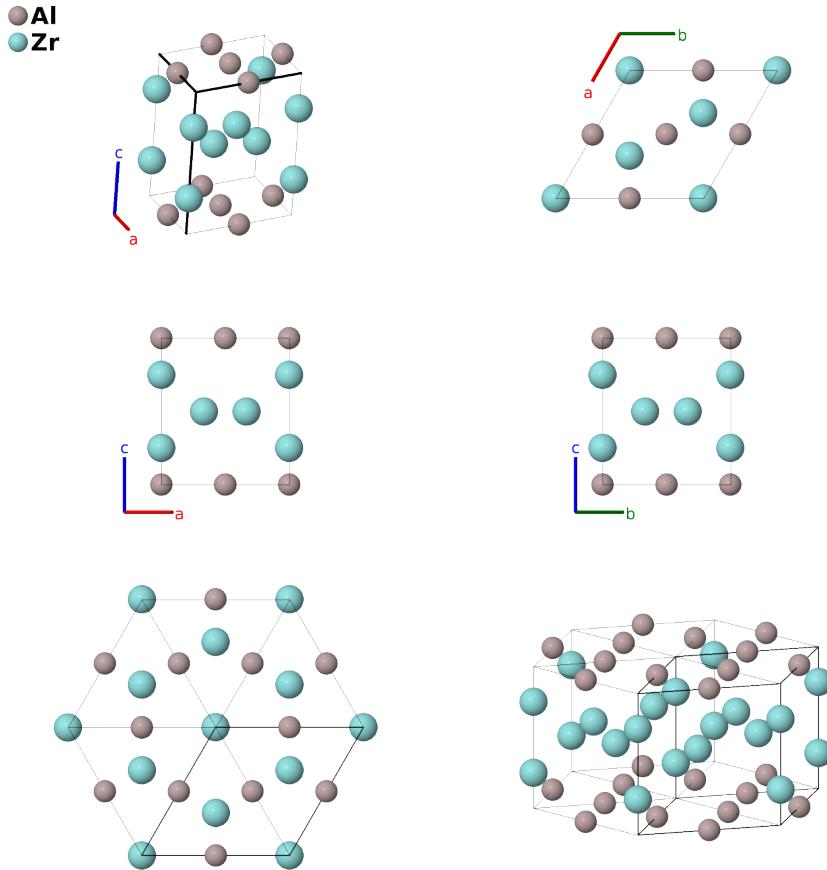


Zr₄Al₃ Structure: A3B4_hP7_191_f_de-001

Cite this page as: H. Eckert, S. Divilov, A. Zettel, M. J. Mehl, D. Hicks, and S. Curtarolo, *The AFLOW Library of Crystallographic Prototypes: Part 4*. In preparation.

<https://aflow.org/p/AMP4>

https://aflow.org/p/A3B4_hP7_191_f_de-001



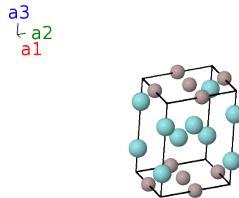
Prototype	Al ₃ Zr ₄
AFLOW prototype label	A3B4_hP7_191_f_de-001
ICSD	150529
Pearson symbol	hP7
Space group number	191
Space group symbol	$P6/mmm$
AFLOW prototype command	<code>aflow --proto=A3B4_hP7_191_f_de-001 --params=a, c/a, z₂</code>

Other compounds with this structure
Hf₄Al₃

- (Wilson, 1960) place this structure in, using modern notation, space group $P\bar{6}m2$ #187, but (Cenzual, 1991) point out that the given atomic coordinates place the structure in the higher symmetry space group $P6/mmm$ #191. We show that structure here.

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	$\frac{1}{3}\mathbf{a}_1 + \frac{2}{3}\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$\frac{1}{2}a\hat{\mathbf{x}} + \frac{\sqrt{3}}{6}a\hat{\mathbf{y}} + \frac{1}{2}c\hat{\mathbf{z}}$	(2d)	Zr I
\mathbf{B}_2	$\frac{2}{3}\mathbf{a}_1 + \frac{1}{3}\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	=	$\frac{1}{2}a\hat{\mathbf{x}} - \frac{\sqrt{3}}{6}a\hat{\mathbf{y}} + \frac{1}{2}c\hat{\mathbf{z}}$	(2d)	Zr I
\mathbf{B}_3	$z_2\mathbf{a}_3$	=	$cz_2\hat{\mathbf{z}}$	(2e)	Zr II
\mathbf{B}_4	$-z_2\mathbf{a}_3$	=	$-cz_2\hat{\mathbf{z}}$	(2e)	Zr II
\mathbf{B}_5	$\frac{1}{2}\mathbf{a}_1$	=	$\frac{1}{4}a\hat{\mathbf{x}} - \frac{\sqrt{3}}{4}a\hat{\mathbf{y}}$	(3f)	Al I
\mathbf{B}_6	$\frac{1}{2}\mathbf{a}_2$	=	$\frac{1}{4}a\hat{\mathbf{x}} + \frac{\sqrt{3}}{4}a\hat{\mathbf{y}}$	(3f)	Al I
\mathbf{B}_7	$\frac{1}{2}\mathbf{a}_1 + \frac{1}{2}\mathbf{a}_2$	=	$\frac{1}{2}a\hat{\mathbf{x}}$	(3f)	Al I

References

[1] C. G. Wilson, D. K. Thomas, and F. J. Spooner, *The crystal structure of Zr₄Al₃*, Acta Cryst. **13**, 56–57 (1960), doi:10.1107/S0365110X60000121.

Found in

[1] K. Cenzual, L. M. Gelato, M. Penzo, and E. Parthé, *Inorganic structure types with revised space groups. I*, Acta Crystallogr. Sect. B **47**, 433–439 (1991), doi:10.1107/S0108768191000903.