

Al₃Ti (*D*0₂₂) Structure:

A3B_tI8_139_ad_b-001

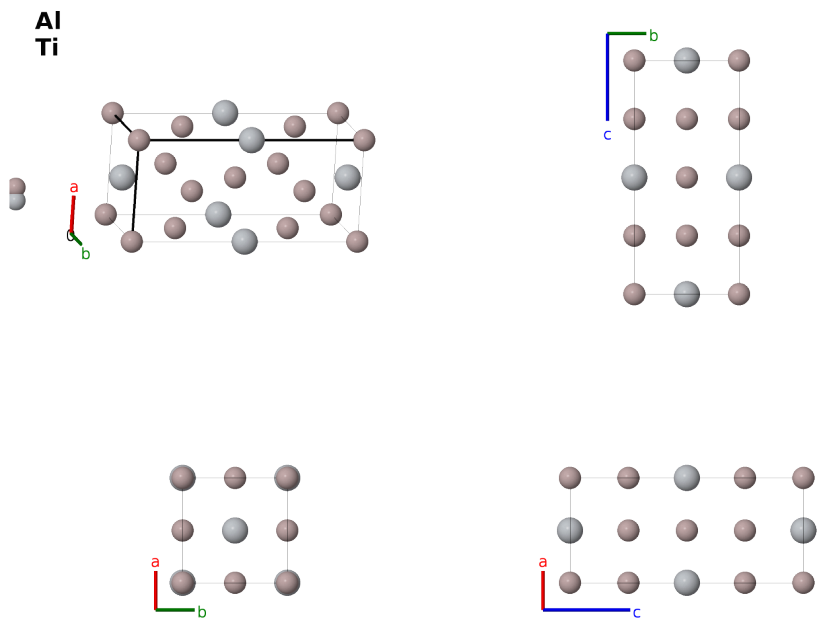
This structure previously used the label(s) **A3B_tI8_139_bd_a**. Calls to these addresses will be redirected here.

Cite this page as:

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<https://aflow.org/prototype-encyclopedia/7W3W>

https://aflow.org/prototype-encyclopedia/A3B_tI8_139_ad_b-001



Prototype	Al ₃ Ti
AFLOW prototype label	A3B_tI8_139_ad_b-001
<i>Strukturbericht</i> designation	<i>D</i> 0 ₂₂
ICSD	58189
CCDC	1622217
Pearson symbol	tI8
Space group number	139
Space group symbol	<i>I</i> 4/ <i>mmm</i>
AFLOW prototype command	<code>aflow --proto=A3B_tI8_139_ad_b-001</code> <code>--params=<i>a</i>, <i>c/a</i></code>

Other compounds with this structure

Al_3Hf , $\text{Al}_{3-x}\text{Fe}_x\text{Mo}$, Al_3Nb , Al_3Sc , Al_3Ta , Al_3V , Ga_3Hf , Ga_3Nb , Ga_3Ti , $\alpha\text{-NbPd}_3$, Ni_3V , Pd_3Nb , Pd_3Ta , Pd_3V , Pt_3V

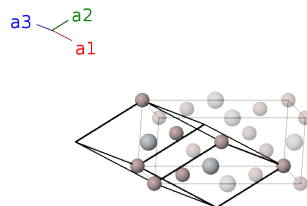
- When $c = 2a$ the atoms are on the sites of a face-centered cubic lattice. When $c/a = 1/\sqrt{2}$, this becomes the cubic $D0_3$ structure.

Body-centered Tetragonal primitive vectors

$$\mathbf{a}_1 = -\frac{1}{2}a\hat{\mathbf{x}} + \frac{1}{2}a\hat{\mathbf{y}} + \frac{1}{2}c\hat{\mathbf{z}}$$

$$\mathbf{a}_2 = \frac{1}{2}a\hat{\mathbf{x}} - \frac{1}{2}a\hat{\mathbf{y}} + \frac{1}{2}c\hat{\mathbf{z}}$$

$$\mathbf{a}_3 = \frac{1}{2}a\hat{\mathbf{x}} + \frac{1}{2}a\hat{\mathbf{y}} - \frac{1}{2}c\hat{\mathbf{z}}$$



Basis vectors

	Lattice coordinates		Cartesian coordinates	Wyckoff position	Atom type
$\mathbf{B}_1$	$=$	0	$=$	0	(2a) Al I
$\mathbf{B}_2$	$=$	$\frac{1}{2}\mathbf{a}_1 + \frac{1}{2}\mathbf{a}_2$	$=$	$\frac{1}{2}c\hat{\mathbf{z}}$	(2b) Ti I
$\mathbf{B}_3$	$=$	$\frac{3}{4}\mathbf{a}_1 + \frac{1}{4}\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	$=$	$\frac{1}{2}a\hat{\mathbf{y}} + \frac{1}{4}c\hat{\mathbf{z}}$	(4d) Al II
$\mathbf{B}_4$	$=$	$\frac{1}{4}\mathbf{a}_1 + \frac{3}{4}\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	$=$	$\frac{1}{2}a\hat{\mathbf{x}} + \frac{1}{4}c\hat{\mathbf{z}}$	(4d) Al II

References

- [1] P. Norby and N. Christensen, *Preparation and Structure of Al_3Ti* , Acta Chem. Scand. pp. 157–159 (1986), doi:10.3891/acta.chem.scand.40a-0157.
- [2] J. Nic, S. Zhang, and D. Mikkola, *Observations on the systematic alloying of Al_3Ti with fourth period elements to yield cubic phases*, Scripta Metallurgica et Materialia **24**, 1099–1104 (1990), doi:10.1016/0956-716X(90)90306-2.

Found in

- [1] W. B. Pearson, *A Handbook of Lattice Spacings and Structures of Metals and Alloys, Volume 2, International Series of Monographs on Metal Physics and Physical Metallurgy*, vol. 8 (Pergamon Press, Oxford, London, Edinburgh, New York, Toronto, Sydney, Paris, Braunschweig, 1967).

First cited in

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