

Ce₃Ni₆Si₂ Structure:

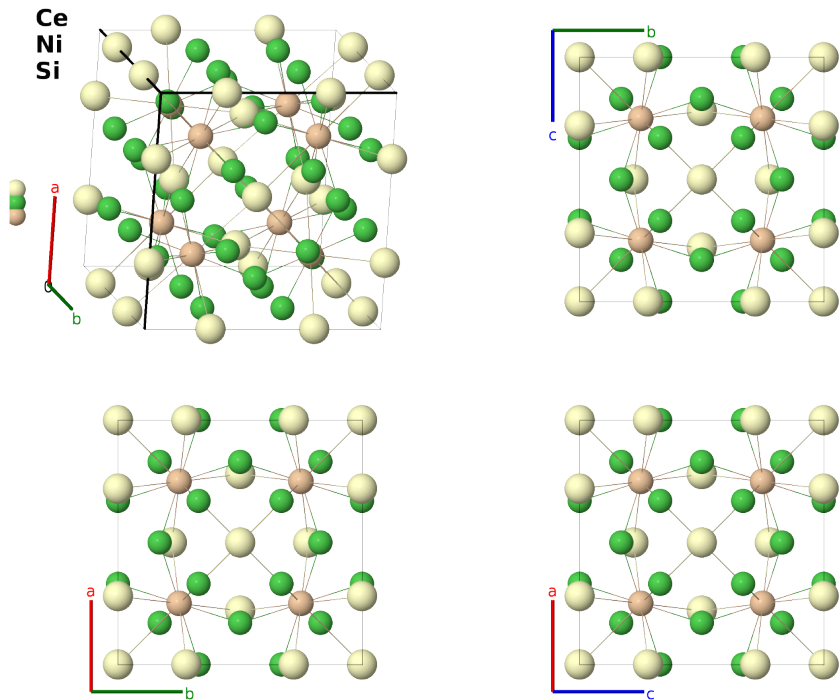
A3B6C2_cI44_229_e_h_c-001

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<https://aflow.org/prototype-encyclopedia/BC5E>

https://aflow.org/prototype-encyclopedia/A3B6C2_cI44_229_e_h_c-001



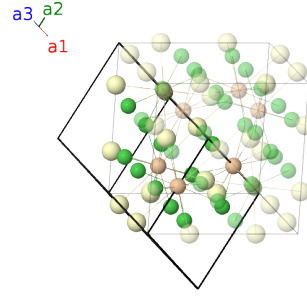
Prototype	Ce ₃ Ni ₆ Si ₂
AFLOW prototype label	A3B6C2_cI44_229_e_h_c-001
ICSD	25622
CCDC	1601272
Pearson symbol	cI44
Space group number	229
Space group symbol	$Im\bar{3}m$
AFLOW prototype command	<code>aflow --proto=A3B6C2_cI44_229_e_h_c-001</code> <code>--params=a, x₂, y₃</code>

Other compounds with this structure

Dy₃Ni₆Si₂, Er₃Ni₆Al₂, Er₃Ni₆Si₂, Eu₃Ni₆Si₂, Gd₃Ni₆Si₂, Ho₃Ni₆Si₂, Lu₃Ni₆Si₂, Nd₃Ni₆Si₂, Pr₃Ni₆Si₂, Sm₃Ni₆Si₂, Tb₃Ni₆Si₂, Tm₃Ni₆Si₂, U₃Ni₆Ge₂, U₃Ni₆Si₂, Yb₃Ni₆Si₂

Body-centered Cubic primitive vectors

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



Basis vectors

	Lattice coordinates		Cartesian coordinates	Wyckoff position	Atom type
$\mathbf{B}_1$	$= \frac{1}{2}\mathbf{a}_1 + \frac{1}{2}\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	$=$	$\frac{1}{4}a\hat{\mathbf{x}} + \frac{1}{4}a\hat{\mathbf{y}} + \frac{1}{4}a\hat{\mathbf{z}}$	(8c)	Si I
$\mathbf{B}_2$	$= \frac{1}{2}\mathbf{a}_3$	$=$	$\frac{1}{4}a\hat{\mathbf{x}} + \frac{1}{4}a\hat{\mathbf{y}} - \frac{1}{4}a\hat{\mathbf{z}}$	(8c)	Si I
$\mathbf{B}_3$	$= \frac{1}{2}\mathbf{a}_2$	$=$	$\frac{1}{4}a\hat{\mathbf{x}} - \frac{1}{4}a\hat{\mathbf{y}} + \frac{1}{4}a\hat{\mathbf{z}}$	(8c)	Si I
$\mathbf{B}_4$	$= \frac{1}{2}\mathbf{a}_1$	$=$	$-\frac{1}{4}a\hat{\mathbf{x}} + \frac{1}{4}a\hat{\mathbf{y}} + \frac{1}{4}a\hat{\mathbf{z}}$	(8c)	Si I
$\mathbf{B}_5$	$= x_2\mathbf{a}_2 + x_2\mathbf{a}_3$	$=$	$ax_2\hat{\mathbf{x}}$	(12e)	Ce I
$\mathbf{B}_6$	$= -x_2\mathbf{a}_2 - x_2\mathbf{a}_3$	$=$	$-ax_2\hat{\mathbf{x}}$	(12e)	Ce I
$\mathbf{B}_7$	$= x_2\mathbf{a}_1 + x_2\mathbf{a}_3$	$=$	$ax_2\hat{\mathbf{y}}$	(12e)	Ce I
$\mathbf{B}_8$	$= -x_2\mathbf{a}_1 - x_2\mathbf{a}_3$	$=$	$-ax_2\hat{\mathbf{y}}$	(12e)	Ce I
$\mathbf{B}_9$	$= x_2\mathbf{a}_1 + x_2\mathbf{a}_2$	$=$	$ax_2\hat{\mathbf{z}}$	(12e)	Ce I
$\mathbf{B}_{10}$	$= -x_2\mathbf{a}_1 - x_2\mathbf{a}_2$	$=$	$-ax_2\hat{\mathbf{z}}$	(12e)	Ce I
$\mathbf{B}_{11}$	$= 2y_3\mathbf{a}_1 + y_3\mathbf{a}_2 + y_3\mathbf{a}_3$	$=$	$ay_3\hat{\mathbf{y}} + ay_3\hat{\mathbf{z}}$	(24h)	Ni I
$\mathbf{B}_{12}$	$= y_3\mathbf{a}_2 - y_3\mathbf{a}_3$	$=$	$-ay_3\hat{\mathbf{y}} + ay_3\hat{\mathbf{z}}$	(24h)	Ni I
$\mathbf{B}_{13}$	$= -y_3\mathbf{a}_2 + y_3\mathbf{a}_3$	$=$	$ay_3\hat{\mathbf{y}} - ay_3\hat{\mathbf{z}}$	(24h)	Ni I
$\mathbf{B}_{14}$	$= -2y_3\mathbf{a}_1 - y_3\mathbf{a}_2 - y_3\mathbf{a}_3$	$=$	$-ay_3\hat{\mathbf{y}} - ay_3\hat{\mathbf{z}}$	(24h)	Ni I
$\mathbf{B}_{15}$	$= y_3\mathbf{a}_1 + 2y_3\mathbf{a}_2 + y_3\mathbf{a}_3$	$=$	$ay_3\hat{\mathbf{x}} + ay_3\hat{\mathbf{z}}$	(24h)	Ni I
$\mathbf{B}_{16}$	$= -y_3\mathbf{a}_1 + y_3\mathbf{a}_3$	$=$	$ay_3\hat{\mathbf{x}} - ay_3\hat{\mathbf{z}}$	(24h)	Ni I
$\mathbf{B}_{17}$	$= y_3\mathbf{a}_1 - y_3\mathbf{a}_3$	$=$	$-ay_3\hat{\mathbf{x}} + ay_3\hat{\mathbf{z}}$	(24h)	Ni I
$\mathbf{B}_{18}$	$= -y_3\mathbf{a}_1 - 2y_3\mathbf{a}_2 - y_3\mathbf{a}_3$	$=$	$-ay_3\hat{\mathbf{x}} - ay_3\hat{\mathbf{z}}$	(24h)	Ni I
$\mathbf{B}_{19}$	$= y_3\mathbf{a}_1 + y_3\mathbf{a}_2 + 2y_3\mathbf{a}_3$	$=$	$ay_3\hat{\mathbf{x}} + ay_3\hat{\mathbf{y}}$	(24h)	Ni I
$\mathbf{B}_{20}$	$= y_3\mathbf{a}_1 - y_3\mathbf{a}_2$	$=$	$-ay_3\hat{\mathbf{x}} + ay_3\hat{\mathbf{y}}$	(24h)	Ni I
$\mathbf{B}_{21}$	$= -y_3\mathbf{a}_1 + y_3\mathbf{a}_2$	$=$	$ay_3\hat{\mathbf{x}} - ay_3\hat{\mathbf{y}}$	(24h)	Ni I
$\mathbf{B}_{22}$	$= -y_3\mathbf{a}_1 - y_3\mathbf{a}_2 - 2y_3\mathbf{a}_3$	$=$	$-ay_3\hat{\mathbf{x}} - ay_3\hat{\mathbf{y}}$	(24h)	Ni I

References

- [1] E. I. Hladyschewskij, P. I. Krypiakewytsh, and O. I. Bodak, *Die Kristallstruktur von $Ce_3Ni_6Si_2$ und verwandten Verbindungen*, Z. Anorganische und Allgemeine Chemie **344**, 95–101 (1966), doi:10.1002/zaac.19663440113.

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