

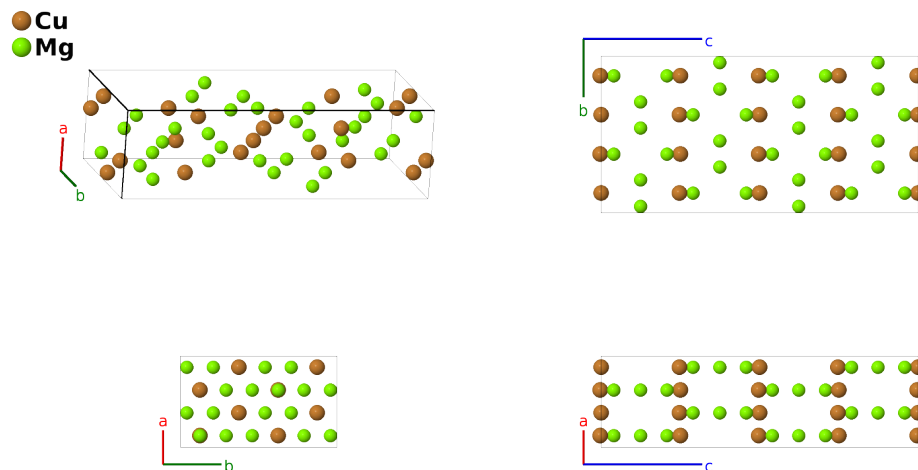
Mg₂Cu (*C_b*) Structure: AB2_oF48_70_e_ef-002

This structure originally had the label AB2_oF48_70_g_fg. Calls to that address will be redirected here.

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<https://aflow.org/p/S336>

https://aflow.org/p/AB2_oF48_70_e_ef-002



Prototype	CuMg ₂
AFLOW prototype label	AB2_oF48_70_e_ef-002
<i>Strukturbericht</i> designation	<i>C_b</i>
ICSD	695334
Pearson symbol	oF48
Space group number	70
Space group symbol	<i>Fddd</i>
AFLOW prototype command	<code>aflow --proto=AB2_oF48_70_e_ef-002 --params=a, b/a, c/a, x₁, x₂, y₃</code>

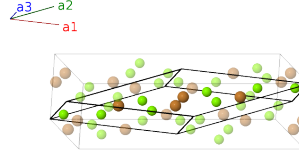
Other compounds with this structure

In₂Ir, NbSn₂, TiBi₂, FeGaGe, SbSnTi

- Mn₂B (*D1_f*) and CuMg₂ (*C_b*) have the same AFLOW prototype label, AB2_oF48_70_e_ef. They are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Face-centered Orthorhombic primitive vectors

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



Basis vectors

	Lattice coordinates	Cartesian coordinates	Wyckoff position	Atom type
\mathbf{B}_1	$= -(x_1 - \frac{1}{4})\mathbf{a}_1 + x_1\mathbf{a}_2 + x_1\mathbf{a}_3 =$	$ax_1\hat{\mathbf{x}} + \frac{1}{8}b\hat{\mathbf{y}} + \frac{1}{8}c\hat{\mathbf{z}}$	(16e)	Cu I
\mathbf{B}_2	$= x_1\mathbf{a}_1 - (x_1 - \frac{1}{4})\mathbf{a}_2 - (x_1 - \frac{1}{4})\mathbf{a}_3 =$	$-a(x_1 - \frac{1}{4})\hat{\mathbf{x}} + \frac{1}{8}b\hat{\mathbf{y}} + \frac{1}{8}c\hat{\mathbf{z}}$	(16e)	Cu I
\mathbf{B}_3	$= (x_1 + \frac{3}{4})\mathbf{a}_1 - x_1\mathbf{a}_2 - x_1\mathbf{a}_3 =$	$-ax_1\hat{\mathbf{x}} + \frac{3}{8}b\hat{\mathbf{y}} + \frac{3}{8}c\hat{\mathbf{z}}$	(16e)	Cu I
\mathbf{B}_4	$= -x_1\mathbf{a}_1 + (x_1 + \frac{3}{4})\mathbf{a}_2 + (x_1 + \frac{3}{4})\mathbf{a}_3 =$	$a(x_1 + \frac{3}{4})\hat{\mathbf{x}} + \frac{3}{8}b\hat{\mathbf{y}} + \frac{3}{8}c\hat{\mathbf{z}}$	(16e)	Cu I
\mathbf{B}_5	$= -(x_2 - \frac{1}{4})\mathbf{a}_1 + x_2\mathbf{a}_2 + x_2\mathbf{a}_3 =$	$ax_2\hat{\mathbf{x}} + \frac{1}{8}b\hat{\mathbf{y}} + \frac{1}{8}c\hat{\mathbf{z}}$	(16e)	Mg I
\mathbf{B}_6	$= x_2\mathbf{a}_1 - (x_2 - \frac{1}{4})\mathbf{a}_2 - (x_2 - \frac{1}{4})\mathbf{a}_3 =$	$-a(x_2 - \frac{1}{4})\hat{\mathbf{x}} + \frac{1}{8}b\hat{\mathbf{y}} + \frac{1}{8}c\hat{\mathbf{z}}$	(16e)	Mg I
\mathbf{B}_7	$= (x_2 + \frac{3}{4})\mathbf{a}_1 - x_2\mathbf{a}_2 - x_2\mathbf{a}_3 =$	$-ax_2\hat{\mathbf{x}} + \frac{3}{8}b\hat{\mathbf{y}} + \frac{3}{8}c\hat{\mathbf{z}}$	(16e)	Mg I
\mathbf{B}_8	$= -x_2\mathbf{a}_1 + (x_2 + \frac{3}{4})\mathbf{a}_2 + (x_2 + \frac{3}{4})\mathbf{a}_3 =$	$a(x_2 + \frac{3}{4})\hat{\mathbf{x}} + \frac{3}{8}b\hat{\mathbf{y}} + \frac{3}{8}c\hat{\mathbf{z}}$	(16e)	Mg I
\mathbf{B}_9	$= y_3\mathbf{a}_1 - (y_3 - \frac{1}{4})\mathbf{a}_2 + y_3\mathbf{a}_3 =$	$\frac{1}{8}a\hat{\mathbf{x}} + by_3\hat{\mathbf{y}} + \frac{1}{8}c\hat{\mathbf{z}}$	(16f)	Mg II
\mathbf{B}_{10}	$= -(y_3 - \frac{1}{4})\mathbf{a}_1 + y_3\mathbf{a}_2 - (y_3 - \frac{1}{4})\mathbf{a}_3 =$	$\frac{1}{8}a\hat{\mathbf{x}} - b(y_3 - \frac{1}{4})\hat{\mathbf{y}} + \frac{1}{8}c\hat{\mathbf{z}}$	(16f)	Mg II
\mathbf{B}_{11}	$= -y_3\mathbf{a}_1 + (y_3 + \frac{3}{4})\mathbf{a}_2 - y_3\mathbf{a}_3 =$	$\frac{3}{8}a\hat{\mathbf{x}} - by_3\hat{\mathbf{y}} + \frac{3}{8}c\hat{\mathbf{z}}$	(16f)	Mg II
\mathbf{B}_{12}	$= (y_3 + \frac{3}{4})\mathbf{a}_1 - y_3\mathbf{a}_2 + (y_3 + \frac{3}{4})\mathbf{a}_3 =$	$\frac{3}{8}a\hat{\mathbf{x}} + b(y_3 + \frac{3}{4})\hat{\mathbf{y}} + \frac{3}{8}c\hat{\mathbf{z}}$	(16f)	Mg II

References

- [1] F. Gingl, P. Selvam, and K. Yvon, *Structure refinement of Mg_2Cu and a comparison of the Mg_2Cu , Mg_2Ni and Al_2Cu structure types*, Acta Crystallogr. Sect. B **49**, 201–203 (1993), doi:10.1107/S0108768192008723.

Found in

- [1] V. Vreshch, *Crystal Structure of $CuMg_2$* (2018). Crystallography online.com.