

Cu₂Fe[CN]₆ Structure: A12B2C_cF60_196_h_ac_b-001

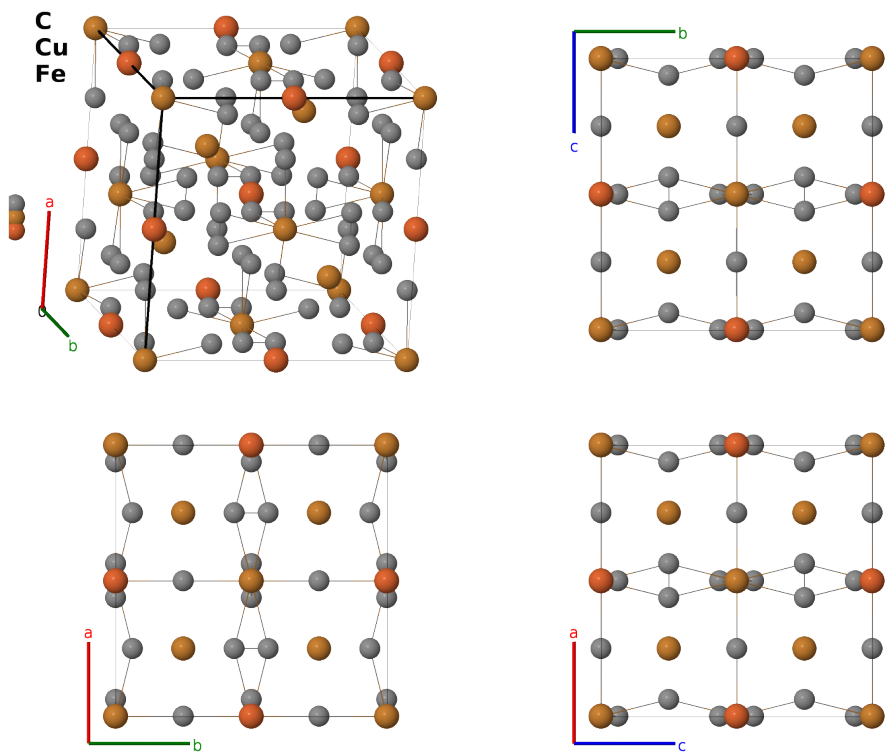
This structure previously used the label(s) **A12B2C_cF60_196_h_bc_a**. Calls to these addresses will be redirected here.

Cite this page as:

- H. Eckert, S. Divilov, M. J. Mehl, D. Hicks, A. C. Zettl, M. Esters, X. Campilongo, and S. Curtarolo, *The AFLOW Library of Crystallographic Prototypes: Part 4*, Comput. Mater. Sci. **240**, 112988 (2024), doi: 10.1016/j.commatsci.2024.112988.
- D. Hicks, M. J. Mehl, M. Esters, C. Oses, O. Levy, G. L. W. Hart, C. Toher, and S. Curtarolo, *The AFLOW Library of Crystallographic Prototypes: Part 3*, Comput. Mater. Sci. **199**, 110450 (2021), doi: 10.1016/j.commatsci.2021.110450.

<https://aflow.org/prototype-encyclopedia/DJ7F>

https://aflow.org/prototype-encyclopedia/A12B2C_cF60_196_h_ac_b-001



Prototype	C ₆ Cu ₂ FeN ₆
AFLOW prototype label	A12B2C_cF60_196_h_ac_b-001
Pearson symbol	cF60
Space group number	196
Space group symbol	<i>F</i> 23
AFLOW prototype command	<code>aflow --proto=A12B2C_cF60_196_h_ac_b-001</code> <code>--params=<i>a</i>, <i>x</i>₄, <i>y</i>₄, <i>z</i>₄</code>

- The sites we have labeled as carbon are actually a 50-50 mixture of carbon and nitrogen.

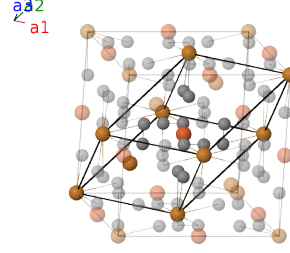
- This structure is taken from (Villars, 2013), repeated in (Villars, 2023). Our print versions of (Rigamonti, 1937) gives a structure consistent with the K_2PtCl_6 ($J1_1$) prototype.

Face-centered Cubic primitive vectors

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

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

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



Basis vectors

	Lattice coordinates		Cartesian coordinates	Wyckoff position	Atom type
$\mathbf{B}_1$	$=$	0	$=$	0	(4a) Cu I
$\mathbf{B}_2$	$=$	$\frac{1}{2} \mathbf{a}_1 + \frac{1}{2} \mathbf{a}_2 + \frac{1}{2} \mathbf{a}_3$	$=$	$\frac{1}{2}a \hat{\mathbf{x}} + \frac{1}{2}a \hat{\mathbf{y}} + \frac{1}{2}a \hat{\mathbf{z}}$	(4b) Fe I
$\mathbf{B}_3$	$=$	$\frac{1}{4} \mathbf{a}_1 + \frac{1}{4} \mathbf{a}_2 + \frac{1}{4} \mathbf{a}_3$	$=$	$\frac{1}{4}a \hat{\mathbf{x}} + \frac{1}{4}a \hat{\mathbf{y}} + \frac{1}{4}a \hat{\mathbf{z}}$	(4c) Cu II
$\mathbf{B}_4$	$=$	$(-x_4 + y_4 + z_4) \mathbf{a}_1 +$ $(x_4 - y_4 + z_4) \mathbf{a}_2 +$ $(x_4 + y_4 - z_4) \mathbf{a}_3$	$=$	$ax_4 \hat{\mathbf{x}} + ay_4 \hat{\mathbf{y}} + az_4 \hat{\mathbf{z}}$	(48h) C I
$\mathbf{B}_5$	$=$	$(x_4 - y_4 + z_4) \mathbf{a}_1 +$ $(-x_4 + y_4 + z_4) \mathbf{a}_2 -$ $(x_4 + y_4 + z_4) \mathbf{a}_3$	$=$	$-ax_4 \hat{\mathbf{x}} - ay_4 \hat{\mathbf{y}} + az_4 \hat{\mathbf{z}}$	(48h) C I
$\mathbf{B}_6$	$=$	$(x_4 + y_4 - z_4) \mathbf{a}_1 -$ $(x_4 + y_4 + z_4) \mathbf{a}_2 +$ $(-x_4 + y_4 + z_4) \mathbf{a}_3$	$=$	$-ax_4 \hat{\mathbf{x}} + ay_4 \hat{\mathbf{y}} - az_4 \hat{\mathbf{z}}$	(48h) C I
$\mathbf{B}_7$	$=$	$-(x_4 + y_4 + z_4) \mathbf{a}_1 +$ $(x_4 + y_4 - z_4) \mathbf{a}_2 +$ $(x_4 - y_4 + z_4) \mathbf{a}_3$	$=$	$ax_4 \hat{\mathbf{x}} - ay_4 \hat{\mathbf{y}} - az_4 \hat{\mathbf{z}}$	(48h) C I
$\mathbf{B}_8$	$=$	$(x_4 + y_4 - z_4) \mathbf{a}_1 +$ $(-x_4 + y_4 + z_4) \mathbf{a}_2 +$ $(x_4 - y_4 + z_4) \mathbf{a}_3$	$=$	$az_4 \hat{\mathbf{x}} + ax_4 \hat{\mathbf{y}} + ay_4 \hat{\mathbf{z}}$	(48h) C I
$\mathbf{B}_9$	$=$	$-(x_4 + y_4 + z_4) \mathbf{a}_1 +$ $(x_4 - y_4 + z_4) \mathbf{a}_2 +$ $(-x_4 + y_4 + z_4) \mathbf{a}_3$	$=$	$az_4 \hat{\mathbf{x}} - ax_4 \hat{\mathbf{y}} - ay_4 \hat{\mathbf{z}}$	(48h) C I
$\mathbf{B}_{10}$	$=$	$(-x_4 + y_4 + z_4) \mathbf{a}_1 +$ $(x_4 + y_4 - z_4) \mathbf{a}_2 -$ $(x_4 + y_4 + z_4) \mathbf{a}_3$	$=$	$-az_4 \hat{\mathbf{x}} - ax_4 \hat{\mathbf{y}} + ay_4 \hat{\mathbf{z}}$	(48h) C I
$\mathbf{B}_{11}$	$=$	$(x_4 - y_4 + z_4) \mathbf{a}_1 -$ $(x_4 + y_4 + z_4) \mathbf{a}_2 +$ $(x_4 + y_4 - z_4) \mathbf{a}_3$	$=$	$-az_4 \hat{\mathbf{x}} + ax_4 \hat{\mathbf{y}} - ay_4 \hat{\mathbf{z}}$	(48h) C I
$\mathbf{B}_{12}$	$=$	$(x_4 - y_4 + z_4) \mathbf{a}_1 +$ $(x_4 + y_4 - z_4) \mathbf{a}_2 +$ $(-x_4 + y_4 + z_4) \mathbf{a}_3$	$=$	$ay_4 \hat{\mathbf{x}} + az_4 \hat{\mathbf{y}} + ax_4 \hat{\mathbf{z}}$	(48h) C I

$$\mathbf{B}_{13} = \begin{pmatrix} -x_4 + y_4 + z_4 \\ x_4 + y_4 + z_4 \\ x_4 - y_4 + z_4 \end{pmatrix} \begin{pmatrix} \mathbf{a}_1 \\ \mathbf{a}_2 \\ \mathbf{a}_3 \end{pmatrix} = -ay_4 \hat{\mathbf{x}} + az_4 \hat{\mathbf{y}} - ax_4 \hat{\mathbf{z}} \quad (48h) \quad \text{C I}$$

$$\mathbf{B}_{14} = \begin{pmatrix} -(x_4 + y_4 + z_4) \\ -x_4 + y_4 + z_4 \\ x_4 + y_4 - z_4 \end{pmatrix} \begin{pmatrix} \mathbf{a}_1 \\ \mathbf{a}_2 \\ \mathbf{a}_3 \end{pmatrix} = ay_4 \hat{\mathbf{x}} - az_4 \hat{\mathbf{y}} - ax_4 \hat{\mathbf{z}} \quad (48h) \quad \text{C I}$$

$$\mathbf{B}_{15} = \begin{pmatrix} x_4 + y_4 - z_4 \\ x_4 - y_4 + z_4 \\ x_4 + y_4 + z_4 \end{pmatrix} \begin{pmatrix} \mathbf{a}_1 \\ \mathbf{a}_2 \\ \mathbf{a}_3 \end{pmatrix} = -ay_4 \hat{\mathbf{x}} - az_4 \hat{\mathbf{y}} + ax_4 \hat{\mathbf{z}} \quad (48h) \quad \text{C I}$$

References

- [1] R. Rigamonti, *Structure of Cupriferrocyanides I. Copper Ferrocyanide and Potassium Copper Ferrocyanide*, Gazz. Chim. Ital. **67**, 137–146 (1937).

Found in

- [1] P. Villars and K. Cenzual, *Pearson's Crystal Data – Crystal Structure Database for Inorganic Compounds* (2013). ASM International.
- [2] P. Villars, *PAULING FILE in: Inorganic Solid Phases, SpringerMaterials (online database)*, Springer, Heidelberg (ed.) (2023). Cu₂Fe(CN)₆ (Cu₂Fe[CN]₆) Crystal Structure, sd₁902947.

First cited in

D. Hicks, M. J. Mehl, E. Gossett, C. Toher, O. Levy, R. M. Hanson, G. Hart, and S. Curtarolo, *The AFLOW Library of Crystallographic Prototypes: Part 2*, Comput. Mater. Sci. **161**, S1 (2019). doi: 10.1016/j.commatsci.2018.10.043