Face-Centered Cubic (Cu, $A1$, fcc) Structure: A_cF4_225_a-001

This structure originally had the label `A_cF4_225_a`. Calls to that address will be redirected here.

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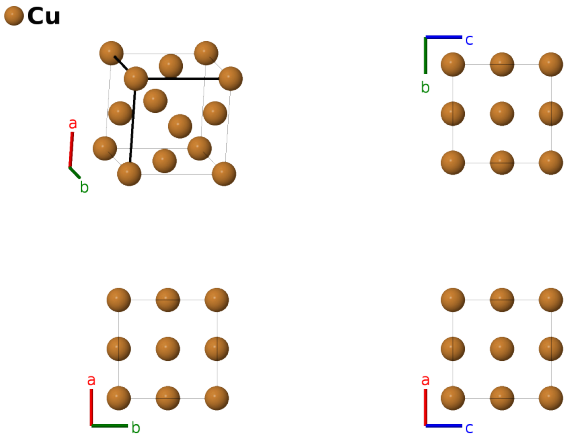
https://aflow.org/p/6EYD

https://aflow.org/p/A_cF4_225_a-001



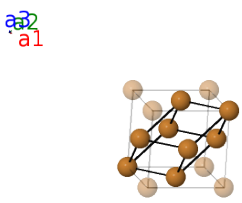
Prototype	Cu
AFLOW prototype label	A_cF4_225_a-001
***Strukturbericht* designation**	$A1$
ICSD	627114
Pearson symbol	cF4
Space group number	225
Space group symbol	$Fm\bar{3}m$
AFLOW prototype command	`aflow --proto=A_cF4_225_a-001 --params=`a

Other compounds with this structure

Al, Ag, Au, Ce, Ir, Ni, Pb, Pd, Pt, Rh, Sr, Tb, Th

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

References

[1] M. E. Straumanis and L. S. Yu, *Lattice parameters, densities, expansion coefficients and perfection of structure of Cu and of Cu-In α phase*, Acta Cryst. **25**, 676–682 (1969), doi:10.1107/S0567739469001549.