

**Name and formula**

Reference code: 00-001-1241  
PDF index name: Copper  
Empirical formula: Cu  
Chemical formula: Cu

**Crystallographic parameters**

Crystal system: Cubic  
Space group: Fm-3m  
Space group number: 225  
  
a (Å): 3,6077  
b (Å): 3,6077  
c (Å): 3,6077  
Alpha (°): 90,0000  
Beta (°): 90,0000  
Gamma (°): 90,0000  
  
Measured density (g/cm<sup>3</sup>): 8,95  
Volume of cell (10<sup>6</sup> pm<sup>3</sup>): 46,96  
Z: 4,00  
  
RIR: -

**Status, subfiles and quality**

Status: Marked as deleted by ICDD  
Subfiles: Inorganic  
Quality: Blank (B)

**Comments**

Deleted by: Deleted by NBS card.  
Color: Red  
Melting point: 1083

**References**

Primary reference: Hanawalt et al., *Anal. Chem.*, **10**, 475, (1938)

**Peak list**

No.	h	k	l	d [Å]	2Theta[deg]	I [%]
1	1	1	1	2,08000	43,473	100,0
2	2	0	0	1,81000	50,375	53,0
3	2	2	0	1,28000	73,997	33,0
4	3	1	1	1,09000	89,934	33,0
5	2	2	2	1,04000	95,578	9,0
6	4	0	0	0,91000	115,662	3,0

**Stick Pattern**