

$\text{Cu}_{25.47}\text{Ce}_7$	$hP78$	$(175) P6/m - \bar{1}^3k^{2,4}he$
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CeCu_{3.6} [1]

Structural features: Variant of GdAg_{3.6} with disordered atom arrangement at $z = 0$.

Allibert C. et al. (1984) [1]

$\text{Ce}_7\text{Cu}_{25.47}$

$a = 1.1858$, $c = 0.9107$ nm, $c/a = 0.768$, $V = 1.1090$ nm³, $Z = 2$

site	Wyck.	sym.	x	y	z	occ.	atomic environment
Cu1	12l	1	0.1079	0.4285	0.3064		
Cu2	12l	1	0.2693	0.0734	0.2509		
Cu3	12l	1	0.4939	0.1117	0.1693		
Cu4	6k	$m..$	0.0593	0.2376	$\frac{1}{2}$		icosahedron Cu ₈ Ce ₄
Ce5	6k	$m..$	0.4712	0.1412	$\frac{1}{2}$		14-vertex polyhedron Cu ₁₄
Cu6	6j	$m..$	0.0378	0.1405	0	0.49	
Ce7	6j	$m..$	0.1056	0.3585	0	0.5	
Ce8	6j	$m..$	0.1309	0.4193	0	0.5	
Cu9	6j	$m..$	0.3000	0.6202	0	0.333	
Cu10	4h	3..	$\frac{1}{3}$	$\frac{2}{3}$	0.2618		
Ce11	2e	6..	0	0	0.2974		

Experimental: single crystal, diffractometer, X-rays, $R = 0.074$, $T = 298$ K

Remarks: Short interatomic distances for partly occupied site(s).

References: [1] Allibert C., Wong Ng W., Nyburg S.C. (1984), Acta Crystallogr. C 40, 211-214.