

CdI ₂	<i>hP</i> 36	(186) <i>P</i> 6 ₃ <i>mc</i> – b ¹² a ⁶
------------------	--------------	--

CdI₂ 24H₇ [1]

Structural features: Close-packed I layers in h₂(hc)₅ stacking; Cd occupies all octahedral voids in every second interlayer. Layer structure with infinite slabs of edge-linked CdI₆ octahedra.

Chadha G.K., Trigunayat G.C. (1967) [1]

CdI₂

a = 0.424, *c* = 8.202 nm, *c/a* = 19.344, *V* = 1.2770 nm³, *Z* = 12

site	Wyck.	sym.	<i>x</i>	<i>y</i>	<i>z</i>	occ.	atomic environment
I1	2 <i>b</i>	3 <i>m.</i>	$\frac{1}{3}$	$\frac{2}{3}$	0.04167		non-coplanar triangle Cd ₃
I2	2 <i>b</i>	3 <i>m.</i>	$\frac{1}{3}$	$\frac{2}{3}$	0.125		non-coplanar triangle Cd ₃
Cd3	2 <i>b</i>	3 <i>m.</i>	$\frac{1}{3}$	$\frac{2}{3}$	0.1875		octahedron I ₆
I4	2 <i>b</i>	3 <i>m.</i>	$\frac{1}{3}$	$\frac{2}{3}$	0.29167		non-coplanar triangle Cd ₃
Cd5	2 <i>b</i>	3 <i>m.</i>	$\frac{1}{3}$	$\frac{2}{3}$	0.35417		octahedron I ₆
I6	2 <i>b</i>	3 <i>m.</i>	$\frac{1}{3}$	$\frac{2}{3}$	0.45833		non-coplanar triangle Cd ₃
Cd7	2 <i>b</i>	3 <i>m.</i>	$\frac{1}{3}$	$\frac{2}{3}$	0.52083		octahedron I ₆
Cd8	2 <i>b</i>	3 <i>m.</i>	$\frac{1}{3}$	$\frac{2}{3}$	0.60417		octahedron I ₆
I9	2 <i>b</i>	3 <i>m.</i>	$\frac{1}{3}$	$\frac{2}{3}$	0.70833		non-coplanar triangle Cd ₃
Cd10	2 <i>b</i>	3 <i>m.</i>	$\frac{1}{3}$	$\frac{2}{3}$	0.77083		octahedron I ₆
I11	2 <i>b</i>	3 <i>m.</i>	$\frac{1}{3}$	$\frac{2}{3}$	0.875		non-coplanar triangle Cd ₃
Cd12	2 <i>b</i>	3 <i>m.</i>	$\frac{1}{3}$	$\frac{2}{3}$	0.9375		octahedron I ₆
I13	2 <i>a</i>	3 <i>m.</i>	0	0	0.0		non-coplanar triangle Cd ₃
I14	2 <i>a</i>	3 <i>m.</i>	0	0	0.08333		non-coplanar triangle Cd ₃
I15	2 <i>a</i>	3 <i>m.</i>	0	0	0.16667		non-coplanar triangle Cd ₃
I16	2 <i>a</i>	3 <i>m.</i>	0	0	0.25		non-coplanar triangle Cd ₃
I17	2 <i>a</i>	3 <i>m.</i>	0	0	0.33333		non-coplanar triangle Cd ₃
I18	2 <i>a</i>	3 <i>m.</i>	0	0	0.41667		non-coplanar triangle Cd ₃

Transformation from published data: -*x*, -*y*, -*z*; origin shift 0 0 0.95833

Experimental: single crystal, Weissenberg photographs, X-rays

Remarks: Zhdanov notation (2222211)₂. We derived idealized atom coordinates from the stacking sequence.

References: [1] Chadha G.K., Trigunayat G.C. (1967), Acta Crystallogr. 23, 726-729.