

# CdAl<sub>2</sub>S<sub>4</sub> (*E*3) Structure: A2BC4\_tI14\_82\_bc\_a\_g-001

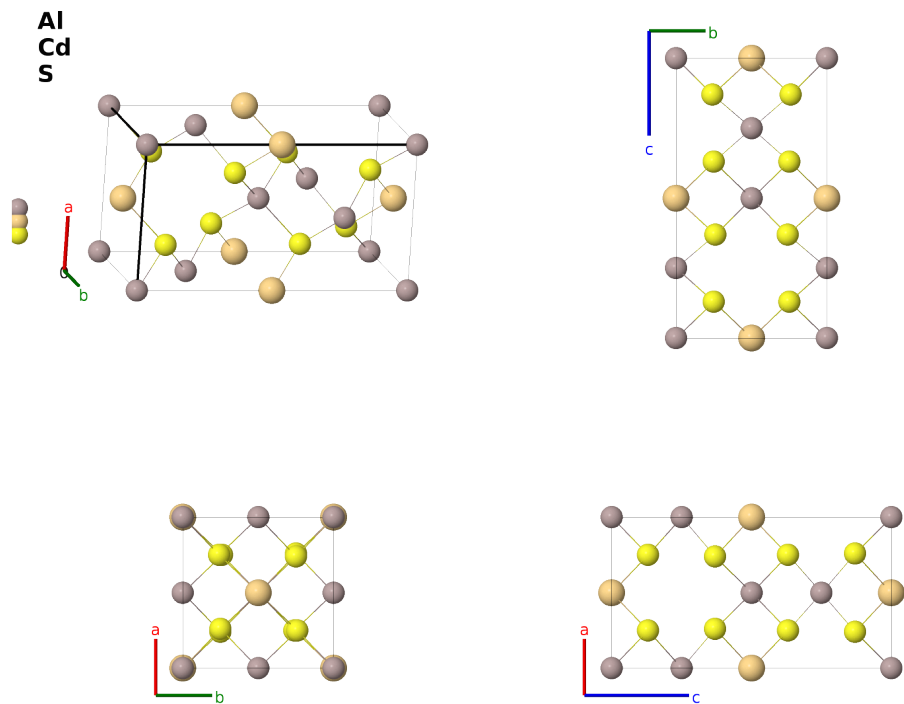
This structure previously used the label(s) **A2BC4\_tI14\_82\_bc\_a\_g**. Calls to these addresses will be redirected here.

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. Osos, 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/S1AD>

[https://aflow.org/prototype-encyclopedia/A2BC4\\_tI14\\_82\\_bc\\_a\\_g-001](https://aflow.org/prototype-encyclopedia/A2BC4_tI14_82_bc_a_g-001)



|                                    |                                                                                                                                  |
|------------------------------------|----------------------------------------------------------------------------------------------------------------------------------|
| Prototype                          | Al <sub>2</sub> CdS <sub>4</sub>                                                                                                 |
| AFLOW prototype label              | A2BC4_tI14_82_bc_a_g-001                                                                                                         |
| <i>Strukturbericht</i> designation | <i>E</i> 3                                                                                                                       |
| ICSD                               | 25634                                                                                                                            |
| CCDC                               | 1601284                                                                                                                          |
| Pearson symbol                     | tI14                                                                                                                             |
| Space group number                 | 82                                                                                                                               |
| Space group symbol                 | $I\bar{4}$                                                                                                                       |
| AFLOW prototype command            | <code>aflow --proto=A2BC4_tI14_82_bc_a_g-001</code><br><code>--params=a, c/a, x<sub>4</sub>, y<sub>4</sub>, z<sub>4</sub></code> |

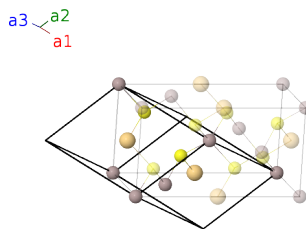
### Other compounds with this structure

CdAl<sub>2</sub>Se<sub>4</sub>, CdAl<sub>2</sub>Te<sub>4</sub>, CdGa<sub>2</sub>S<sub>4</sub>, CdGa<sub>2</sub>Se<sub>4</sub>, CdGa<sub>2</sub>Te<sub>4</sub>, CoGa<sub>2</sub>S<sub>4</sub>, FeGa<sub>2</sub>S<sub>4</sub>, HfGa<sub>2</sub>Se<sub>4</sub>,  $\beta$ -HgAg<sub>2</sub>I<sub>4</sub>, HgAl<sub>2</sub>S<sub>4</sub>, HgAl<sub>2</sub>Se<sub>4</sub>, HgAl<sub>2</sub>Te<sub>4</sub>, HgGa<sub>2</sub>S<sub>4</sub>, HgGa<sub>2</sub>Te<sub>4</sub>, HgIn<sub>2</sub>Se<sub>4</sub>, HgIn<sub>2</sub>Te<sub>4</sub>, ZnGa<sub>2</sub>S<sub>4</sub>, ZnGa<sub>2</sub>Se<sub>4</sub>, ZnGa<sub>2</sub>Te<sub>4</sub>, ZnIn<sub>2</sub>Se<sub>4</sub>, ZnIn<sub>2</sub>Te<sub>4</sub>

- When  $c = 2a$  and  $x = y = 1/4$ , and  $z = 1/8$  the atoms are on the sites of the diamond (A4) structure, but of course there are defects. Removing the Al-I (2b) atom transforms this to the BPO<sub>4</sub> (*H07*) structure.
- (Pearson, 1967) gives this the designation *E3*, without any subscript. It should be defined as *E3<sub>a</sub>*, but as he did not put any other structures into this subcategory, we will leave it as is.

### Body-centered Tetragonal primitive vectors

$$\begin{aligned} \mathbf{a}_1 &= -\frac{1}{2}a \hat{\mathbf{x}} + \frac{1}{2}a \hat{\mathbf{y}} + \frac{1}{2}c \hat{\mathbf{z}} \\ \mathbf{a}_2 &= \frac{1}{2}a \hat{\mathbf{x}} - \frac{1}{2}a \hat{\mathbf{y}} + \frac{1}{2}c \hat{\mathbf{z}} \\ \mathbf{a}_3 &= \frac{1}{2}a \hat{\mathbf{x}} + \frac{1}{2}a \hat{\mathbf{y}} - \frac{1}{2}c \hat{\mathbf{z}} \end{aligned}$$



### Basis vectors

|                | Lattice coordinates                                                                 |     | Cartesian coordinates                                                    | Wyckoff position | Atom type |
|----------------|-------------------------------------------------------------------------------------|-----|--------------------------------------------------------------------------|------------------|-----------|
| $\mathbf{B}_1$ | $= 0$                                                                               | $=$ | $0$                                                                      | (2a)             | Al I      |
| $\mathbf{B}_2$ | $= \frac{1}{2} \mathbf{a}_1 + \frac{1}{2} \mathbf{a}_2$                             | $=$ | $\frac{1}{2}c \hat{\mathbf{z}}$                                          | (2b)             | Cd I      |
| $\mathbf{B}_3$ | $= \frac{3}{4} \mathbf{a}_1 + \frac{1}{4} \mathbf{a}_2 + \frac{1}{2} \mathbf{a}_3$  | $=$ | $\frac{1}{2}a \hat{\mathbf{y}} + \frac{1}{4}c \hat{\mathbf{z}}$          | (2c)             | Al II     |
| $\mathbf{B}_4$ | $= (y_4 + z_4) \mathbf{a}_1 + (x_4 + z_4) \mathbf{a}_2 + (x_4 + y_4) \mathbf{a}_3$  | $=$ | $ax_4 \hat{\mathbf{x}} + ay_4 \hat{\mathbf{y}} + cz_4 \hat{\mathbf{z}}$  | (8g)             | S I       |
| $\mathbf{B}_5$ | $= -(y_4 - z_4) \mathbf{a}_1 - (x_4 - z_4) \mathbf{a}_2 - (x_4 + y_4) \mathbf{a}_3$ | $=$ | $-ax_4 \hat{\mathbf{x}} - ay_4 \hat{\mathbf{y}} + cz_4 \hat{\mathbf{z}}$ | (8g)             | S I       |
| $\mathbf{B}_6$ | $= -(x_4 + z_4) \mathbf{a}_1 + (y_4 - z_4) \mathbf{a}_2 - (x_4 - y_4) \mathbf{a}_3$ | $=$ | $ay_4 \hat{\mathbf{x}} - ax_4 \hat{\mathbf{y}} - cz_4 \hat{\mathbf{z}}$  | (8g)             | S I       |
| $\mathbf{B}_7$ | $= (x_4 - z_4) \mathbf{a}_1 - (y_4 + z_4) \mathbf{a}_2 + (x_4 - y_4) \mathbf{a}_3$  | $=$ | $-ay_4 \hat{\mathbf{x}} + ax_4 \hat{\mathbf{y}} - cz_4 \hat{\mathbf{z}}$ | (8g)             | S I       |

### References

- [1] H. Hahn, G. Frank, W. Klingler, A.-D. Störger, and G. Störger, *Über ternäre Chalogenide des Aluminiums, Galliums und Indiums mit Zink, Cadmium und Quecksilber*, Z. Anorganische und Allgemeine Chemie **279**, 241–270 (1955), doi:10.1002/zaac.19552790502.
- [2] E. Parthé, L. Gelato, B. Chabot, M. Penso, K. Cenzula, and R. Gladyshevskii, *Standardized Data and Crystal Chemical Characterization of Inorganic Structure Types*, Gmelin Handbook of Inorganic and Organometallic Chemistry, vol. 2 (Springer-Verlag, Berlin, Heidelberg, 1993), 8 edn., doi:10.1007/978-3-662-02909-1\_3.
- [3] P. Villars and L. Calvert, *Pearson's Handbook of Crystallographic Data for Intermetallic Phases* (ASM International, Materials Park, OH, 1991), 2nd edn.
- [4] C. Hermann, O. Lohrmann, and H. Philipp, eds., *Strukturbericht Band II 1928-1932* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).

- [5] W. B. Pearson, *A Handbook of Lattice Spacings and Structures of Metals and Alloys, Volume 2, International Series of Monographs on Metal Physics and Physical Metallurgy*, vol. 8 (Pergamon Press, Oxford, London, Edinburgh, New York, Toronto, Sydney, Paris, Braunschweig, 1967).

**Found in**

- [1] P. Villars, *CdAl<sub>2</sub>S<sub>4</sub> Crystal Structure* (2016). PAULING FILE in: Inorganic Solid Phases, SpringerMaterials (online database), Springer, Heidelberg (ed.) SpringerMaterials.

**First cited in**

M. J. Mehl, D. Hicks, C. Toher, O. Levy, R. M. Hanson, G. L. W. Hart, and S. Curtarolo, *The AFLOW Library of Crystallographic Prototypes: Part 1*, Comput. Mater. Sci. **136**, S1-828 (2017). doi: 10.1016/j.commatsci.2017.01.017