

BaS₃ (*D*0₁₇) Structure: AB3_tP8_113_a_ce-001

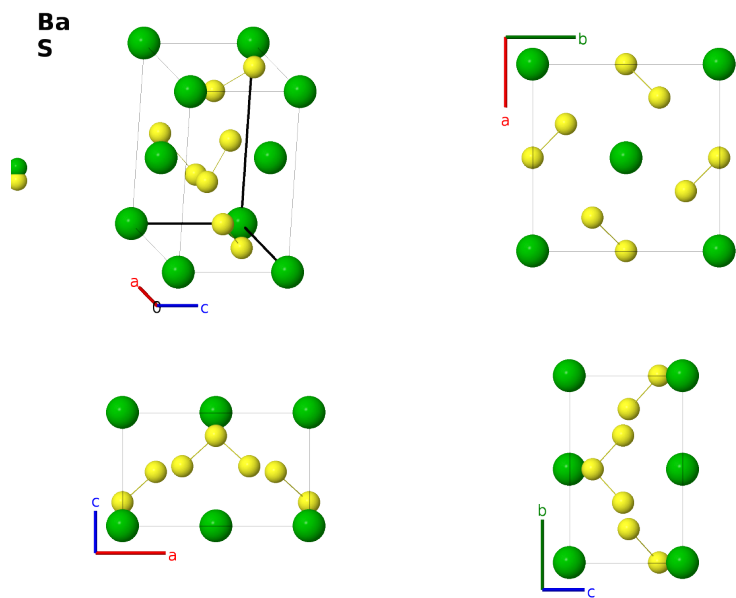
This structure previously used the label(s) **AB3_tP8_113_a_ce**. Calls to these addresses will be redirected here.

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. Oses, 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/0A77>

https://aflow.org/prototype-encyclopedia/AB3_tP8_113_a_ce-001



Prototype	BaS ₃
AFLOW prototype label	AB3_tP8_113_a_ce-001
<i>Strukturbericht</i> designation	<i>D</i> 0 ₁₇
ICSD	70059
CCDC	1631883
Pearson symbol	tP8
Space group number	113
Space group symbol	<i>P</i> $\bar{4}$ ₂ <i>m</i>
AFLOW prototype command	<code>aflow --proto=AB3_tP8_113_a_ce-001</code> <code>--params=<i>a</i>, <i>c/a</i>, <i>z</i>₂, <i>x</i>₃, <i>z</i>₃</code>

Other compounds with this structure

AgDyTe₂, AgHgTe₂, AgErTe₂, AgTe₂Tm, AgGdTe₂, AgTe₂Y, BaSe₃, BaTe₃

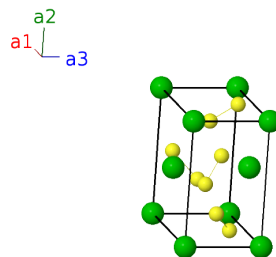
- (Gottfried, 1938) originally gave the $D0_{17}$ label to the $P2_12_12$ #94 BaS₃ structure, however we follow (Partheé, 1993), who uses the current structure as the $D0_{17}$ prototype.

Simple Tetragonal primitive vectors

$$\mathbf{a}_1 = a \hat{\mathbf{x}}$$

$$\mathbf{a}_2 = a \hat{\mathbf{y}}$$

$$\mathbf{a}_3 = c \hat{\mathbf{z}}$$



Basis vectors

	Lattice coordinates		Cartesian coordinates	Wyckoff position	Atom type
$\mathbf{B}_1$	$=$	0	$=$	0	(2a) Ba I
$\mathbf{B}_2$	$=$	$\frac{1}{2} \mathbf{a}_1 + \frac{1}{2} \mathbf{a}_2$	$=$	$\frac{1}{2} a \hat{\mathbf{x}} + \frac{1}{2} a \hat{\mathbf{y}}$	(2a) Ba I
$\mathbf{B}_3$	$=$	$\frac{1}{2} \mathbf{a}_2 + z_3 \mathbf{a}_3$	$=$	$\frac{1}{2} a \hat{\mathbf{y}} + cz_3 \hat{\mathbf{z}}$	(2c) S I
$\mathbf{B}_4$	$=$	$\frac{1}{2} \mathbf{a}_1 - z_3 \mathbf{a}_3$	$=$	$\frac{1}{2} a \hat{\mathbf{x}} - cz_3 \hat{\mathbf{z}}$	(2c) S I
$\mathbf{B}_5$	$=$	$x_3 \mathbf{a}_1 + (x_3 + \frac{1}{2}) \mathbf{a}_2 + z_3 \mathbf{a}_3$	$=$	$ax_3 \hat{\mathbf{x}} + a(x_3 + \frac{1}{2}) \hat{\mathbf{y}} + cz_3 \hat{\mathbf{z}}$	(4e) S II
$\mathbf{B}_6$	$=$	$-x_3 \mathbf{a}_1 - (x_3 - \frac{1}{2}) \mathbf{a}_2 + z_3 \mathbf{a}_3$	$=$	$-ax_3 \hat{\mathbf{x}} - a(x_3 - \frac{1}{2}) \hat{\mathbf{y}} + cz_3 \hat{\mathbf{z}}$	(4e) S II
$\mathbf{B}_7$	$=$	$(x_3 + \frac{1}{2}) \mathbf{a}_1 - x_3 \mathbf{a}_2 - z_3 \mathbf{a}_3$	$=$	$a(x_3 + \frac{1}{2}) \hat{\mathbf{x}} - ax_3 \hat{\mathbf{y}} - cz_3 \hat{\mathbf{z}}$	(4e) S II
$\mathbf{B}_8$	$=$	$-(x_3 - \frac{1}{2}) \mathbf{a}_1 + x_3 \mathbf{a}_2 - z_3 \mathbf{a}_3$	$=$	$-a(x_3 - \frac{1}{2}) \hat{\mathbf{x}} + ax_3 \hat{\mathbf{y}} - cz_3 \hat{\mathbf{z}}$	(4e) S II

References

- [1] S. Yamaoka, J. T. Lemley, J. M. Jenks, and H. Steinfink, *Structural chemistry of the polysulfides dibarium trisulfide and monobarium trisulfide*, Inorg. Chem. **14**, 129–131 (1975), doi:10.1021/ic50143a027.
- [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] C. Gottfried, ed., *Strukturbericht Band IV 1936* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1938).

Found in

- [1] P. Villars and L. Calvert, *Pearson's Handbook of Crystallographic Data for Intermetallic Phases* (ASM International, Materials Park, OH, 1991), 2nd edn.

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