

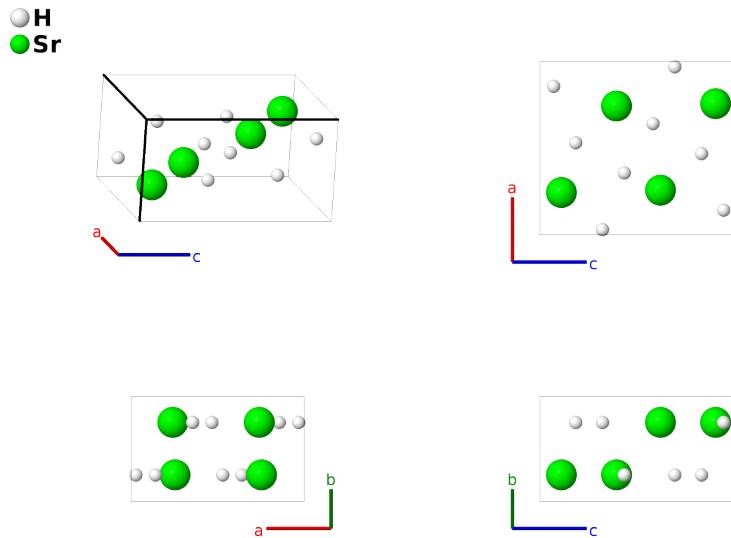
SrH₂ (*C*29) Structure: A2B_oP12_62_2c_c-004

This structure originally had the label A2B_oP12_62_2c_c.SrH2. Calls to that address will be redirected here.

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<https://aflow.org/p/4GNV>

https://aflow.org/p/A2B_oP12_62_2c_c-004



Prototype	H ₂ Sr
AFLOW prototype label	A2B_oP12_62_2c_c-004
Strukturbericht designation	<i>C</i> 29
ICSD	69077
Pearson symbol	oP12
Space group number	62
Space group symbol	<i>Pnma</i>
AFLOW prototype command	<code>aflow --proto=A2B_oP12_62_2c_c-004 --params=<i>a</i>,<i>b/a</i>,<i>c/a</i>,<i>x</i>₁,<i>z</i>₁,<i>x</i>₂,<i>z</i>₂,<i>x</i>₃,<i>z</i>₃</code>

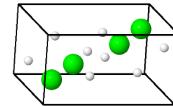
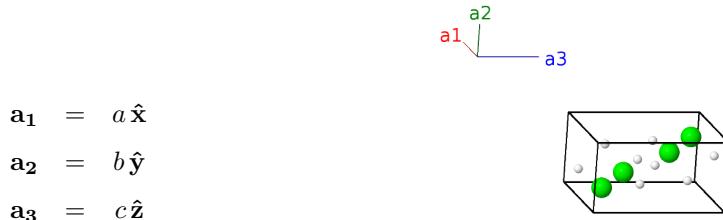
Other compounds with this structure

CaH₂, EuD₂, YbH₂

- We use this SrH₂ as the prototype for the hydrides with this structure.
- The ternary form of this structure is MnCuP. See that structure for similar ternary compounds.

- PbCl_2 (*C23*), HgCl_2 (*C25*), SrH_2 (*C29*), Co_2Si (*C37*), and SrBr_2 (*C53*) all share the same AFLOW label, `A2B_oP12_62_2c_c`. The structures are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.
- We have updated the reported structure from (Roop, 2013) to use the actual structure from (Brese, 1990). The sample studied used deuterium.

Simple Orthorhombic primitive vectors



Basis vectors

	Lattice coordinates	Cartesian coordinates	Wyckoff position	Atom type
\mathbf{B}_1	$x_1 \mathbf{a}_1 + \frac{1}{4} \mathbf{a}_2 + z_1 \mathbf{a}_3$	$a x_1 \hat{\mathbf{x}} + \frac{1}{4} b \hat{\mathbf{y}} + c z_1 \hat{\mathbf{z}}$	(4c)	H I
\mathbf{B}_2	$-(x_1 - \frac{1}{2}) \mathbf{a}_1 + \frac{3}{4} \mathbf{a}_2 + (z_1 + \frac{1}{2}) \mathbf{a}_3$	$-a(x_1 - \frac{1}{2}) \hat{\mathbf{x}} + \frac{3}{4} b \hat{\mathbf{y}} + c(z_1 + \frac{1}{2}) \hat{\mathbf{z}}$	(4c)	H I
\mathbf{B}_3	$-x_1 \mathbf{a}_1 + \frac{3}{4} \mathbf{a}_2 - z_1 \mathbf{a}_3$	$-a x_1 \hat{\mathbf{x}} + \frac{3}{4} b \hat{\mathbf{y}} - c z_1 \hat{\mathbf{z}}$	(4c)	H I
\mathbf{B}_4	$(x_1 + \frac{1}{2}) \mathbf{a}_1 + \frac{1}{4} \mathbf{a}_2 - (z_1 - \frac{1}{2}) \mathbf{a}_3$	$a(x_1 + \frac{1}{2}) \hat{\mathbf{x}} + \frac{1}{4} b \hat{\mathbf{y}} - c(z_1 - \frac{1}{2}) \hat{\mathbf{z}}$	(4c)	H I
\mathbf{B}_5	$x_2 \mathbf{a}_1 + \frac{1}{4} \mathbf{a}_2 + z_2 \mathbf{a}_3$	$a x_2 \hat{\mathbf{x}} + \frac{1}{4} b \hat{\mathbf{y}} + c z_2 \hat{\mathbf{z}}$	(4c)	H II
\mathbf{B}_6	$-(x_2 - \frac{1}{2}) \mathbf{a}_1 + \frac{3}{4} \mathbf{a}_2 + (z_2 + \frac{1}{2}) \mathbf{a}_3$	$-a(x_2 - \frac{1}{2}) \hat{\mathbf{x}} + \frac{3}{4} b \hat{\mathbf{y}} + c(z_2 + \frac{1}{2}) \hat{\mathbf{z}}$	(4c)	H II
\mathbf{B}_7	$-x_2 \mathbf{a}_1 + \frac{3}{4} \mathbf{a}_2 - z_2 \mathbf{a}_3$	$-a x_2 \hat{\mathbf{x}} + \frac{3}{4} b \hat{\mathbf{y}} - c z_2 \hat{\mathbf{z}}$	(4c)	H II
\mathbf{B}_8	$(x_2 + \frac{1}{2}) \mathbf{a}_1 + \frac{1}{4} \mathbf{a}_2 - (z_2 - \frac{1}{2}) \mathbf{a}_3$	$a(x_2 + \frac{1}{2}) \hat{\mathbf{x}} + \frac{1}{4} b \hat{\mathbf{y}} - c(z_2 - \frac{1}{2}) \hat{\mathbf{z}}$	(4c)	H II
\mathbf{B}_9	$x_3 \mathbf{a}_1 + \frac{1}{4} \mathbf{a}_2 + z_3 \mathbf{a}_3$	$a x_3 \hat{\mathbf{x}} + \frac{1}{4} b \hat{\mathbf{y}} + c z_3 \hat{\mathbf{z}}$	(4c)	Sr I
\mathbf{B}_{10}	$-(x_3 - \frac{1}{2}) \mathbf{a}_1 + \frac{3}{4} \mathbf{a}_2 + (z_3 + \frac{1}{2}) \mathbf{a}_3$	$-a(x_3 - \frac{1}{2}) \hat{\mathbf{x}} + \frac{3}{4} b \hat{\mathbf{y}} + c(z_3 + \frac{1}{2}) \hat{\mathbf{z}}$	(4c)	Sr I
\mathbf{B}_{11}	$-x_3 \mathbf{a}_1 + \frac{3}{4} \mathbf{a}_2 - z_3 \mathbf{a}_3$	$-a x_3 \hat{\mathbf{x}} + \frac{3}{4} b \hat{\mathbf{y}} - c z_3 \hat{\mathbf{z}}$	(4c)	Sr I
\mathbf{B}_{12}	$(x_3 + \frac{1}{2}) \mathbf{a}_1 + \frac{1}{4} \mathbf{a}_2 - (z_3 - \frac{1}{2}) \mathbf{a}_3$	$a(x_3 + \frac{1}{2}) \hat{\mathbf{x}} + \frac{1}{4} b \hat{\mathbf{y}} - c(z_3 - \frac{1}{2}) \hat{\mathbf{z}}$	(4c)	Sr I

References

- [1] N. E. Brese, M. O'Keeffe, and R. B. von Dreele, *Synthesis and crystal structure of SrD_2 and SrND and bond valence parameters for hydrides*, J. Solid State Chem. **88**, 571–576 (1990), doi:10.1016/0022-4596(90)90255-V.
- [2] R. C. Ropp, *Encyclopedia of the Alkaline Earth Compounds* (Elsevier, Oxford, 2013), chap. 2, pp. 30–31.