

H₃S (130 GPa) Structure: A3B_hR4_160_b_a-001

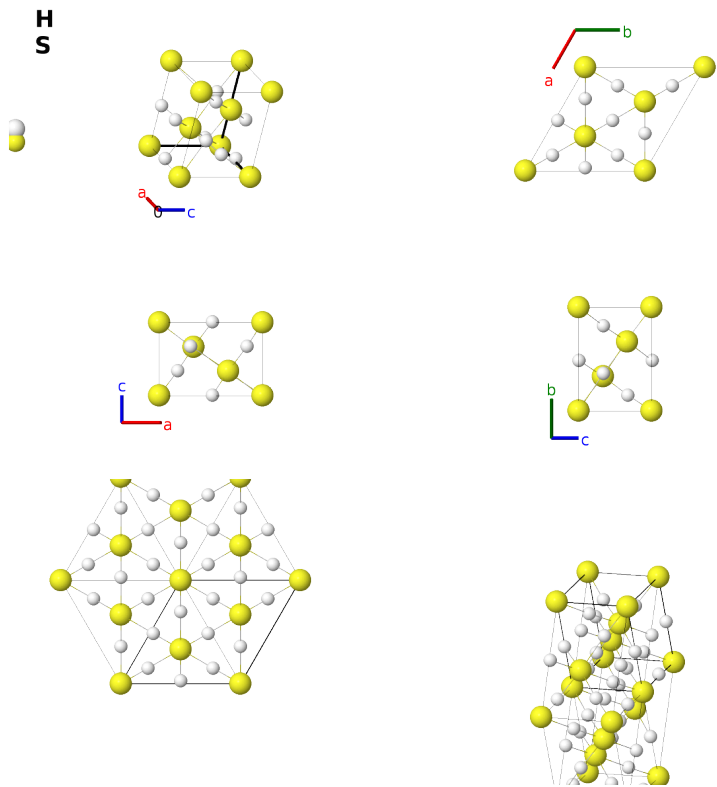
This structure previously used the label(s) **A3B_hR4_160_b_a**. Calls to these addresses will be redirected here.

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<https://aflow.org/prototype-encyclopedia/P2LA>

https://aflow.org/prototype-encyclopedia/A3B_hR4_160_b_a-001

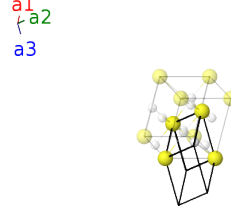


Prototype	H ₃ S
AFLOW prototype label	A3B_hR4_160_b_a-001
ICSD	291501
CCDC	1723598
Pearson symbol	hR4
Space group number	160
Space group symbol	<i>R</i> 3 <i>m</i>
AFLOW prototype command	<code>aflow --proto=A3B_hR4_160_b_a-001</code> <code>--params=<i>a</i>, <i>c/a</i>, <i>x</i>₁, <i>x</i>₂, <i>z</i>₂</code>

- This structure was found by first-principles electronic structure calculations and is predicted to be the stable structure of H_3S for pressures between 90 and 150 GPa. When $c/a \rightarrow \sqrt{8}$, $x_2 \rightarrow 1/2$ and $z_2 \rightarrow 0$ this structure continuously evolves into the cubic 200 GPa H_3S structure.
- The data presented here was computed at 130 GPa.

Rhombohedral primitive vectors

$$\begin{aligned}\mathbf{a}_1 &= \frac{1}{2}a \hat{\mathbf{x}} - \frac{\sqrt{3}}{6}a \hat{\mathbf{y}} + \frac{1}{3}c \hat{\mathbf{z}} \\ \mathbf{a}_2 &= \frac{1}{\sqrt{3}}a \hat{\mathbf{y}} + \frac{1}{3}c \hat{\mathbf{z}} \\ \mathbf{a}_3 &= -\frac{1}{2}a \hat{\mathbf{x}} - \frac{\sqrt{3}}{6}a \hat{\mathbf{y}} + \frac{1}{3}c \hat{\mathbf{z}}\end{aligned}$$



Basis vectors

	Lattice coordinates		Cartesian coordinates	Wyckoff position	Atom type
$\mathbf{B}_1$	$= x_1 \mathbf{a}_1 + x_1 \mathbf{a}_2 + x_1 \mathbf{a}_3$	$=$	$cx_1 \hat{\mathbf{z}}$	(1a)	S I
$\mathbf{B}_2$	$= x_2 \mathbf{a}_1 + x_2 \mathbf{a}_2 + x_2 \mathbf{a}_3$	$=$	$\frac{1}{2}a(x_2 - z_2) \hat{\mathbf{x}} + \frac{\sqrt{3}}{6}a(x_2 - z_2) \hat{\mathbf{y}} + \frac{1}{3}c(2x_2 + z_2) \hat{\mathbf{z}}$	(3b)	H I
$\mathbf{B}_3$	$= z_2 \mathbf{a}_1 + x_2 \mathbf{a}_2 + x_2 \mathbf{a}_3$	$=$	$-\frac{1}{2}a(x_2 - z_2) \hat{\mathbf{x}} + \frac{\sqrt{3}}{6}a(x_2 - z_2) \hat{\mathbf{y}} + \frac{1}{3}c(2x_2 + z_2) \hat{\mathbf{z}}$	(3b)	H I
$\mathbf{B}_4$	$= x_2 \mathbf{a}_1 + z_2 \mathbf{a}_2 + x_2 \mathbf{a}_3$	$=$	$-\frac{1}{\sqrt{3}}a(x_2 - z_2) \hat{\mathbf{y}} + \frac{1}{3}c(2x_2 + z_2) \hat{\mathbf{z}}$	(3b)	H I

References

- [1] D. Duan, Y. Liu, F. T. D. Li, X. Huang, Z. Zhao, H. Yu, B. Liu, W. Tian, and T. Cui, *Pressure-induced metallization of dense $(\text{H}_2\text{S})_2\text{H}_2$ with high- T_c superconductivity* **4**, 6968 (2014), doi:10.1038/srep06968.

First cited in

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