

α -Cristobalite (SiO_2 , low, $C30$) Structure:

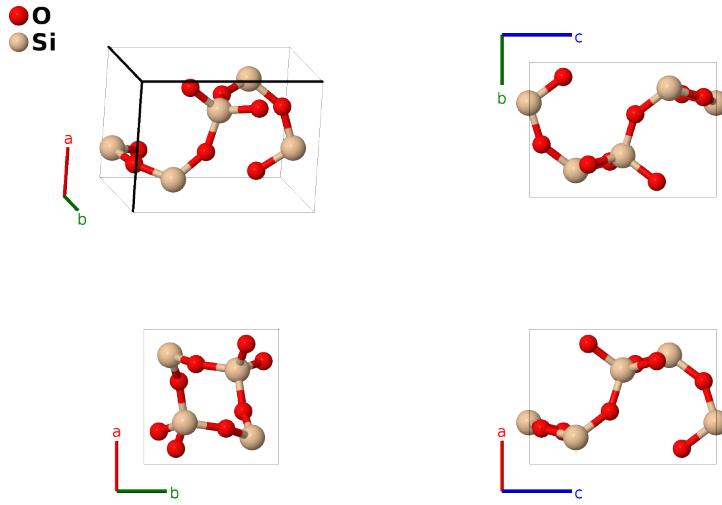
A2B_tP12_92_b_a-001

This structure originally had the label A2B_tP12_92_b_a. Calls to that address will be redirected here.

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<https://aflow.org/p/VHRN>

https://aflow.org/p/A2B_tP12_92_b_a-001



Prototype	O_2Si
AFLOW prototype label	A2B_tP12_92_b_a-001
Strukturbericht designation	$C30$
Mineral name	low (α) cristobalite
ICSD	47221
Pearson symbol	tP12
Space group number	92
Space group symbol	$P4_12_12$
AFLOW prototype command	<code>aflow --proto=A2B_tP12_92_b_a-001 --params=a, c/a, x1, x2, y2, z2</code>

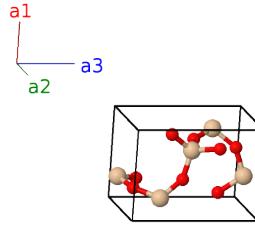
Other compounds with this structure

BeF_2

- If the silicon atoms are replaced by aluminium and phosphorous atoms the structure is very similar to the “low cristobalite” form of AlPO_4 .
- Paralellurite and α -cristobalite ($C30$) have the same AFLOW prototype label, A2B_tP12_92_b_a. They are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Simple Tetragonal primitive vectors

$$\begin{aligned}\mathbf{a}_1 &= a \hat{\mathbf{x}} \\ \mathbf{a}_2 &= a \hat{\mathbf{y}} \\ \mathbf{a}_3 &= c \hat{\mathbf{z}}\end{aligned}$$



Basis vectors

	Lattice coordinates	Cartesian coordinates	Wyckoff position	Atom type
\mathbf{B}_1	$x_1 \mathbf{a}_1 + x_1 \mathbf{a}_2$	$a x_1 \hat{\mathbf{x}} + a x_1 \hat{\mathbf{y}}$	(4a)	Si I
\mathbf{B}_2	$-x_1 \mathbf{a}_1 - x_1 \mathbf{a}_2 + \frac{1}{2} \mathbf{a}_3$	$-a x_1 \hat{\mathbf{x}} - a x_1 \hat{\mathbf{y}} + \frac{1}{2} c \hat{\mathbf{z}}$	(4a)	Si I
\mathbf{B}_3	$-(x_1 - \frac{1}{2}) \mathbf{a}_1 + (x_1 + \frac{1}{2}) \mathbf{a}_2 + \frac{1}{4} \mathbf{a}_3$	$-a(x_1 - \frac{1}{2}) \hat{\mathbf{x}} + a(x_1 + \frac{1}{2}) \hat{\mathbf{y}} + \frac{1}{4} c \hat{\mathbf{z}}$	(4a)	Si I
\mathbf{B}_4	$(x_1 + \frac{1}{2}) \mathbf{a}_1 - (x_1 - \frac{1}{2}) \mathbf{a}_2 + \frac{3}{4} \mathbf{a}_3$	$a(x_1 + \frac{1}{2}) \hat{\mathbf{x}} - a(x_1 - \frac{1}{2}) \hat{\mathbf{y}} + \frac{3}{4} c \hat{\mathbf{z}}$	(4a)	Si I
\mathbf{B}_5	$x_2 \mathbf{a}_1 + y_2 \mathbf{a}_2 + z_2 \mathbf{a}_3$	$a x_2 \hat{\mathbf{x}} + a y_2 \hat{\mathbf{y}} + c z_2 \hat{\mathbf{z}}$	(8b)	O I
\mathbf{B}_6	$-x_2 \mathbf{a}_1 - y_2 \mathbf{a}_2 + (z_2 + \frac{1}{2}) \mathbf{a}_3$	$-a x_2 \hat{\mathbf{x}} - a y_2 \hat{\mathbf{y}} + c(z_2 + \frac{1}{2}) \hat{\mathbf{z}}$	(8b)	O I
\mathbf{B}_7	$-(y_2 - \frac{1}{2}) \mathbf{a}_1 + (x_2 + \frac{1}{2}) \mathbf{a}_2 + (z_2 + \frac{1}{4}) \mathbf{a}_3$	$-a(y_2 - \frac{1}{2}) \hat{\mathbf{x}} + a(x_2 + \frac{1}{2}) \hat{\mathbf{y}} + c(z_2 + \frac{1}{4}) \hat{\mathbf{z}}$	(8b)	O I
\mathbf{B}_8	$(y_2 + \frac{1}{2}) \mathbf{a}_1 - (x_2 - \frac{1}{2}) \mathbf{a}_2 + (z_2 + \frac{3}{4}) \mathbf{a}_3$	$a(y_2 + \frac{1}{2}) \hat{\mathbf{x}} - a(x_2 - \frac{1}{2}) \hat{\mathbf{y}} + c(z_2 + \frac{3}{4}) \hat{\mathbf{z}}$	(8b)	O I
\mathbf{B}_9	$-(x_2 - \frac{1}{2}) \mathbf{a}_1 + (y_2 + \frac{1}{2}) \mathbf{a}_2 - (z_2 - \frac{1}{4}) \mathbf{a}_3$	$-a(x_2 - \frac{1}{2}) \hat{\mathbf{x}} + a(y_2 + \frac{1}{2}) \hat{\mathbf{y}} - c(z_2 - \frac{1}{4}) \hat{\mathbf{z}}$	(8b)	O I
\mathbf{B}_{10}	$(x_2 + \frac{1}{2}) \mathbf{a}_1 - (y_2 - \frac{1}{2}) \mathbf{a}_2 - (z_2 - \frac{3}{4}) \mathbf{a}_3$	$a(x_2 + \frac{1}{2}) \hat{\mathbf{x}} - a(y_2 - \frac{1}{2}) \hat{\mathbf{y}} - c(z_2 - \frac{3}{4}) \hat{\mathbf{z}}$	(8b)	O I
\mathbf{B}_{11}	$y_2 \mathbf{a}_1 + x_2 \mathbf{a}_2 - z_2 \mathbf{a}_3$	$a y_2 \hat{\mathbf{x}} + a x_2 \hat{\mathbf{y}} - c z_2 \hat{\mathbf{z}}$	(8b)	O I
\mathbf{B}_{12}	$-y_2 \mathbf{a}_1 - x_2 \mathbf{a}_2 - (z_2 - \frac{1}{2}) \mathbf{a}_3$	$-a y_2 \hat{\mathbf{x}} - a x_2 \hat{\mathbf{y}} - c(z_2 - \frac{1}{2}) \hat{\mathbf{z}}$	(8b)	O I

References

- [1] J. J. Pluth, J. V. Smith, and J. J. Faber, *Crystal structure of low cristobalite at 10, 293, and 473 K: Variation of framework geometry with temperature*, J. Appl. Phys. **57**, 1045–1049 (1985), doi:10.1063/1.334545.

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.