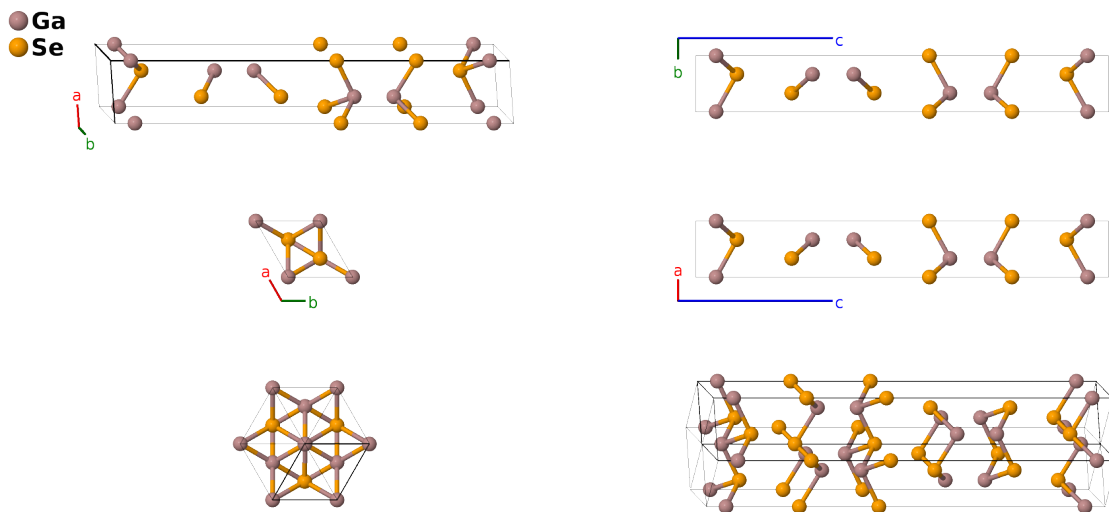


γ -GaSe Structure: AB_hR4_160_2a_2a-001

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<https://aflow.org/p/M3V5>

https://aflow.org/p/AB_hR4_160_2a_2a-001



Prototype	GaSe
AFLOW prototype label	AB_hR4_160_2a_2a-001
ICSD	73388
Pearson symbol	hR4
Space group number	160
Space group symbol	$R\bar{3}m$
AFLOW prototype command	<code>aflow --proto=AB_hR4_160_2a_2a-001 --params=a, c/a, x1, x2, x3, x4</code>

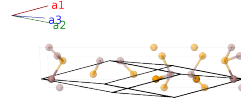
Other compounds with this structure

InSe

- GaSe takes on a variety of structures depending on the stacking of the Ga_2Se_2 layers:
 - β -GaSe is in space group $P6_3/mmc$ #194.
 - γ -GaSe (this structure) is in space group $R\bar{3}m$ #160.
 - δ -GaSe is in space group $P6_3mc$ #186.
 - ϵ -GaSe is in space group $P\bar{6}m2$ #187.
- There is no ICSD entry for (Kuhn, 1975), so we use the earlier work of (Schubert, 1955).
- Hexagonal settings of this structure can be obtained with the option `--hex`.

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) Ga I
\mathbf{B}_2	$=$	$x_2 \mathbf{a}_1 + x_2 \mathbf{a}_2 + x_2 \mathbf{a}_3$	$=$	$cx_2 \hat{\mathbf{z}}$	(1a) Ga II
\mathbf{B}_3	$=$	$x_3 \mathbf{a}_1 + x_3 \mathbf{a}_2 + x_3 \mathbf{a}_3$	$=$	$cx_3 \hat{\mathbf{z}}$	(1a) Se I
\mathbf{B}_4	$=$	$x_4 \mathbf{a}_1 + x_4 \mathbf{a}_2 + x_4 \mathbf{a}_3$	$=$	$cx_4 \hat{\mathbf{z}}$	(1a) Se II

References

- [1] A. Kuhn, A. Chevy, and R. Chevalier, *Crystal Structure and Interatomic Distances in GaSe*, phys. stat. sol. (a) **31**, 469–475 (1975), doi:10.1002/pssa.2210310216.
- [2] K. Schubert, E. Dörre, and M. Kluge, *Zur Kristallchemie der B-Metalle. III. Kristallstruktur von GaSe und InTe*, Z. Metallkd. **46**, 216–224 (1955), doi:10.1515/ijmr-1955-460312.