

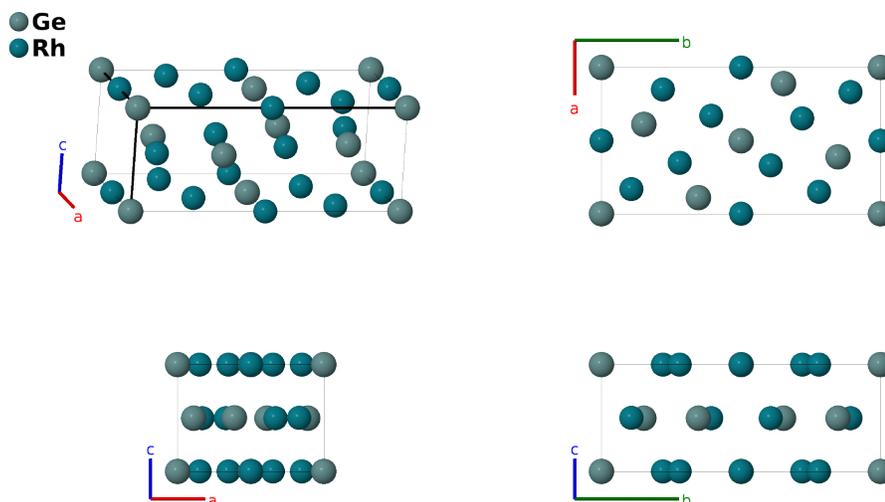
Rh₅Ge₃ Structure: A3B5_oP16_55_ah_cgh-001

This structure originally had the label A3B5_oP16_55_ch_agh. Calls to that address will be redirected here.

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

https://aflow.org/p/A3B5_oP16_55_ah_cgh-001



Prototype	Ge ₃ Rh ₅
AFLOW prototype label	A3B5_oP16_55_ah_cgh-001
ICSD	53891
Pearson symbol	oP16
Space group number	55
Space group symbol	<i>Pbam</i>
AFLOW prototype command	aflow --proto=A3B5_oP16_55_ah_cgh-001 --params= <i>a, b/a, c/a, x₃, y₃, x₄, y₄, x₅, y₅</i>

Other compounds with this structure

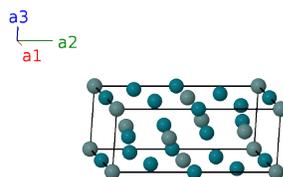
Pd₅Al₃, Pd₅Ga₃, Pd₅In₃, Pt₅Al₃, Rh₅Si₃

Simple Orthorhombic primitive vectors

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

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

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



Basis vectors

	Lattice coordinates		Cartesian coordinates	Wyckoff position	Atom type
\mathbf{B}_1	$=$	0	$=$	0	(2a) Ge I
\mathbf{B}_2	$=$	$\frac{1}{2}\mathbf{a}_1 + \frac{1}{2}\mathbf{a}_2$	$=$	$\frac{1}{2}a\hat{\mathbf{x}} + \frac{1}{2}b\hat{\mathbf{y}}$	(2a) Ge I
\mathbf{B}_3	$=$	$\frac{1}{2}\mathbf{a}_2$	$=$	$\frac{1}{2}b\hat{\mathbf{y}}$	(2c) Rh I
\mathbf{B}_4	$=$	$\frac{1}{2}\mathbf{a}_1$	$=$	$\frac{1}{2}a\hat{\mathbf{x}}$	(2c) Rh I
\mathbf{B}_5	$=$	$x_3\mathbf{a}_1 + y_3\mathbf{a}_2$	$=$	$ax_3\hat{\mathbf{x}} + by_3\hat{\mathbf{y}}$	(4g) Rh II
\mathbf{B}_6	$=$	$-x_3\mathbf{a}_1 - y_3\mathbf{a}_2$	$=$	$-ax_3\hat{\mathbf{x}} - by_3\hat{\mathbf{y}}$	(4g) Rh II
\mathbf{B}_7	$=$	$-(x_3 - \frac{1}{2})\mathbf{a}_1 + (y_3 + \frac{1}{2})\mathbf{a}_2$	$=$	$-a(x_3 - \frac{1}{2})\hat{\mathbf{x}} + b(y_3 + \frac{1}{2})\hat{\mathbf{y}}$	(4g) Rh II
\mathbf{B}_8	$=$	$(x_3 + \frac{1}{2})\mathbf{a}_1 - (y_3 - \frac{1}{2})\mathbf{a}_2$	$=$	$a(x_3 + \frac{1}{2})\hat{\mathbf{x}} - b(y_3 - \frac{1}{2})\hat{\mathbf{y}}$	(4g) Rh II
\mathbf{B}_9	$=$	$x_4\mathbf{a}_1 + y_4\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	$=$	$ax_4\hat{\mathbf{x}} + by_4\hat{\mathbf{y}} + \frac{1}{2}c\hat{\mathbf{z}}$	(4h) Ge II
\mathbf{B}_{10}	$=$	$-x_4\mathbf{a}_1 - y_4\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	$=$	$-ax_4\hat{\mathbf{x}} - by_4\hat{\mathbf{y}} + \frac{1}{2}c\hat{\mathbf{z}}$	(4h) Ge II
\mathbf{B}_{11}	$=$	$-(x_4 - \frac{1}{2})\mathbf{a}_1 + (y_4 + \frac{1}{2})\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	$=$	$-a(x_4 - \frac{1}{2})\hat{\mathbf{x}} + b(y_4 + \frac{1}{2})\hat{\mathbf{y}} + \frac{1}{2}c\hat{\mathbf{z}}$	(4h) Ge II
\mathbf{B}_{12}	$=$	$(x_4 + \frac{1}{2})\mathbf{a}_1 - (y_4 - \frac{1}{2})\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	$=$	$a(x_4 + \frac{1}{2})\hat{\mathbf{x}} - b(y_4 - \frac{1}{2})\hat{\mathbf{y}} + \frac{1}{2}c\hat{\mathbf{z}}$	(4h) Ge II
\mathbf{B}_{13}	$=$	$x_5\mathbf{a}_1 + y_5\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	$=$	$ax_5\hat{\mathbf{x}} + by_5\hat{\mathbf{y}} + \frac{1}{2}c\hat{\mathbf{z}}$	(4h) Rh III
\mathbf{B}_{14}	$=$	$-x_5\mathbf{a}_1 - y_5\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	$=$	$-ax_5\hat{\mathbf{x}} - by_5\hat{\mathbf{y}} + \frac{1}{2}c\hat{\mathbf{z}}$	(4h) Rh III
\mathbf{B}_{15}	$=$	$-(x_5 - \frac{1}{2})\mathbf{a}_1 + (y_5 + \frac{1}{2})\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	$=$	$-a(x_5 - \frac{1}{2})\hat{\mathbf{x}} + b(y_5 + \frac{1}{2})\hat{\mathbf{y}} + \frac{1}{2}c\hat{\mathbf{z}}$	(4h) Rh III
\mathbf{B}_{16}	$=$	$(x_5 + \frac{1}{2})\mathbf{a}_1 - (y_5 - \frac{1}{2})\mathbf{a}_2 + \frac{1}{2}\mathbf{a}_3$	$=$	$a(x_5 + \frac{1}{2})\hat{\mathbf{x}} - b(y_5 - \frac{1}{2})\hat{\mathbf{y}} + \frac{1}{2}c\hat{\mathbf{z}}$	(4h) Rh III

References

- [1] S. Geller, *The rhodium-germanium system. I. The crystal structures of Rh₂Ge, Rh₅Ge₃ and RhGe*, Acta Cryst. **8**, 15–21 (1955), doi:10.1107/S0365110X55000030.

Found in

- [1] P. Villars and K. Cenzual, *Pearson's Crystal Data – Crystal Structure Database for Inorganic Compounds* (2013). ASM International.