

Ga₃Pt₅ Structure:

A3B5_oC16_65_ah_bej-001

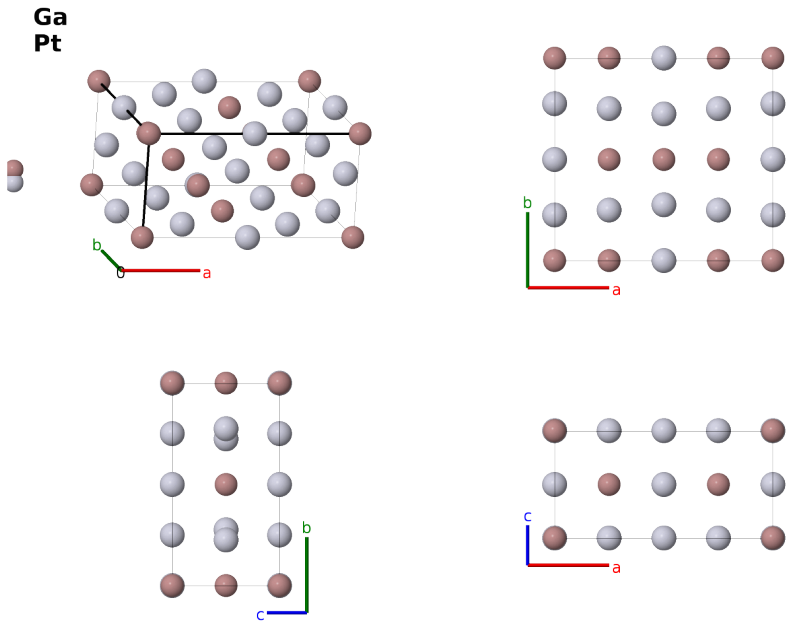
This structure previously used the label(s) **A3B5_oC16_65_ah_bej**. Calls to these addresses will be redirected here.

Cite this page as:

- H. Eckert, S. Divilov, M. J. Mehl, D. Hicks, A. C. Zettl, M. Esters, X. Campilongo, and S. Curtarolo, *The AFLOW Library of Crystallographic Prototypes: Part 4*, Comput. Mater. Sci. **240**, 112988 (2024), doi: 10.1016/j.commatsci.2024.112988.
- D. Hicks, M. J. Mehl, M. Esters, C. Oses, O. Levy, G. L. W. Hart, C. Toher, and S. Curtarolo, *The AFLOW Library of Crystallographic Prototypes: Part 3*, Comput. Mater. Sci. **199**, 110450 (2021), doi: 10.1016/j.commatsci.2021.110450.

<https://aflow.org/prototype-encyclopedia/36DL>

https://aflow.org/prototype-encyclopedia/A3B5_oC16_65_ah_bej-001



Prototype	Ga ₃ Pt ₅
AFLOW prototype label	A3B5_oC16_65_ah_bej-001
ICSD	103927
CCDC	1661858
Pearson symbol	oC16
Space group number	65
Space group symbol	<i>Cmmm</i>
AFLOW prototype command	<code>aflow --proto=A3B5_oC16_65_ah_bej-001</code> <code>--params=a, b/a, c/a, x₄, y₅</code>

Other compounds with this structure

Al₃Ni₅, Ga₃Ni₅, Mn₃Pd₅

Base-centered Orthorhombic primitive vectors



Basis vectors

	Lattice coordinates		Cartesian coordinates	Wyckoff position	Atom type
$\mathbf{B}_1$	=	0	=	0	Ga I
$\mathbf{B}_2$	=	$\frac{1}{2} \mathbf{a}_1 + \frac{1}{2} \mathbf{a}_2$	=	$\frac{1}{2}a \hat{\mathbf{x}}$	Pt I
$\mathbf{B}_3$	=	$\frac{1}{2} \mathbf{a}_2$	=	$\frac{1}{4}a \hat{\mathbf{x}} + \frac{1}{4}b \hat{\mathbf{y}}$	Pt II
$\mathbf{B}_4$	=	$\frac{1}{2} \mathbf{a}_1$	=	$\frac{1}{4}a \hat{\mathbf{x}} - \frac{1}{4}b \hat{\mathbf{y}}$	Pt II
$\mathbf{B}_5$	=	$x_4 \mathbf{a}_1 + x_4 \mathbf{a}_2 + \frac{1}{2} \mathbf{a}_3$	=	$ax_4 \hat{\mathbf{x}} + \frac{1}{2}c \hat{\mathbf{z}}$	Ga II
$\mathbf{B}_6$	=	$-x_4 \mathbf{a}_1 - x_4 \mathbf{a}_2 + \frac{1}{2} \mathbf{a}_3$	=	$-ax_4 \hat{\mathbf{x}} + \frac{1}{2}c \hat{\mathbf{z}}$	Ga II
$\mathbf{B}_7$	=	$-y_5 \mathbf{a}_1 + y_5 \mathbf{a}_2 + \frac{1}{2} \mathbf{a}_3$	=	$by_5 \hat{\mathbf{y}} + \frac{1}{2}c \hat{\mathbf{z}}$	Pt III
$\mathbf{B}_8$	=	$y_5 \mathbf{a}_1 - y_5 \mathbf{a}_2 + \frac{1}{2} \mathbf{a}_3$	=	$-by_5 \hat{\mathbf{y}} + \frac{1}{2}c \hat{\mathbf{z}}$	Pt III

References

- [1] K. Schubert, S. Bhan, W. Burkhardt, R. Gohle, H. G. Meissner, M. Pötzschke, and E. Stolz, *Einige strukturelle Ergebnisse an metallischen Phasen (5)*, *Naturwissenschaften* **47**, 303 (1960), doi:10.1007/BF00600960.

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.

First cited in

- M. J. Mehl, D. Hicks, C. Toher, O. Levy, R. M. Hanson, G. L. W. Hart, and S. Curtarolo, *The AFLOW Library of Crystallographic Prototypes: Part 1*, *Comput. Mater. Sci.* **136**, S1-828 (2017). doi: 10.1016/j.commatsci.2017.01.017