

HoCoGa₅ Structure:

AB5C_tP7_123_b_ci_a-001

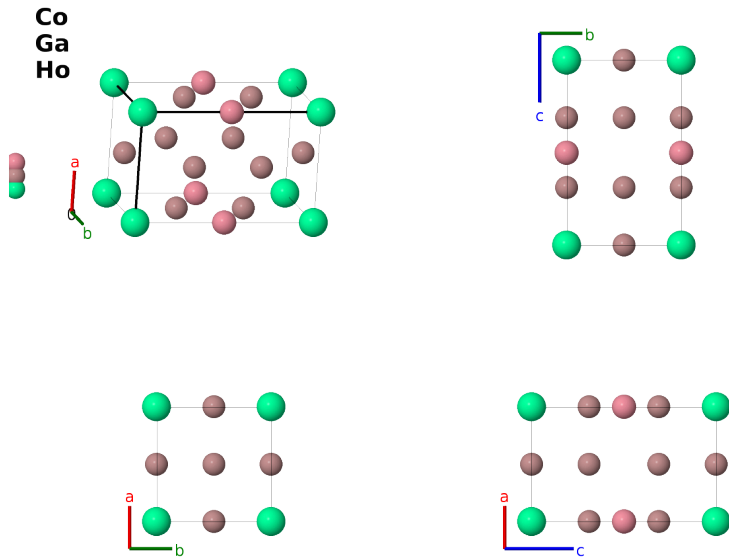
This structure previously used the label(s) **AB5C_tP7_123_b_ci_a**. Calls to these addresses will be redirected here.

Cite this page as:

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- 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/SBCX>

https://aflow.org/prototype-encyclopedia/AB5C_tP7_123_b_ci_a-001



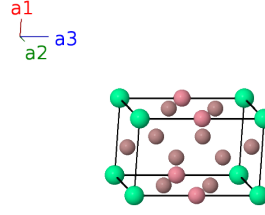
Prototype	CoGa ₅ Ho
AFLOW prototype label	AB5C_tP7_123_b_ci_a-001
ICSD	42427
CCDC	1612093
Pearson symbol	tP7
Space group number	123
Space group symbol	<i>P</i> 4/ <i>mmm</i>
AFLOW prototype command	<code>aflow --proto=AB5C_tP7_123_b_ci_a-001</code> <code>--params=a,c/a,z₄</code>

Other compounds with this structure

CeRhIn₅, CeCoIn₅, DyCoGa₅, ErCoGa₅, GdCoGa₅, LuCoGa₅, TbCoGa₅, TmCoGa₅, UCoGa₅, YCoGa₅, CoIrIn₅, CoRhIn₅, LaCoIn₅, LaIrIn₅, LaRhIn₅, PrCoIn₅, PrIrIn₅, PrRhIn₅

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
$\mathbf{B}_1$	$=$	0	$=$	0	Ho I
$\mathbf{B}_2$	$=$	$\frac{1}{2} \mathbf{a}_3$	$=$	$\frac{1}{2} c \hat{\mathbf{z}}$	Co I
$\mathbf{B}_3$	$=$	$\frac{1}{2} \mathbf{a}_1 + \frac{1}{2} \mathbf{a}_2$	$=$	$\frac{1}{2} a \hat{\mathbf{x}} + \frac{1}{2} a \hat{\mathbf{y}}$	Ga I
$\mathbf{B}_4$	$=$	$\frac{1}{2} \mathbf{a}_2 + z_4 \mathbf{a}_3$	$=$	$\frac{1}{2} a \hat{\mathbf{y}} + cz_4 \hat{\mathbf{z}}$	Ga II
$\mathbf{B}_5$	$=$	$\frac{1}{2} \mathbf{a}_1 + z_4 \mathbf{a}_3$	$=$	$\frac{1}{2} a \hat{\mathbf{x}} + cz_4 \hat{\mathbf{z}}$	Ga II
$\mathbf{B}_6$	$=$	$\frac{1}{2} \mathbf{a}_2 - z_4 \mathbf{a}_3$	$=$	$\frac{1}{2} a \hat{\mathbf{y}} - cz_4 \hat{\mathbf{z}}$	Ga II
$\mathbf{B}_7$	$=$	$\frac{1}{2} \mathbf{a}_1 - z_4 \mathbf{a}_3$	$=$	$\frac{1}{2} a \hat{\mathbf{x}} - cz_4 \hat{\mathbf{z}}$	Ga II

References

- [1] Y. Grin, Y. Yarmolyuk, and E. I. Gladyshevskii, *Kristallicheskie struktury soedinenij $R_2\text{CoGa}_8$ ($R=\text{Sm, Gd, Tb, Dy, Ho, Er, Tm, Lu, Y}$) and $R\text{CoGa}_5$ ($R=\text{Gd, Tb, Dy, Ho, Er, Tm, Lu, Y}$)*, Kristallografiya **24**, 242–246 (1979).

Found in

- [1] P. Villars, *HoCoGa5 Crystal Structure* (2016). PAULING FILE in: Inorganic Solid Phases, SpringerMaterials (online database).

First cited in

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