

α -ReO₃ ($D0_9$) Structure: A3B_cP4_221_c_b-001

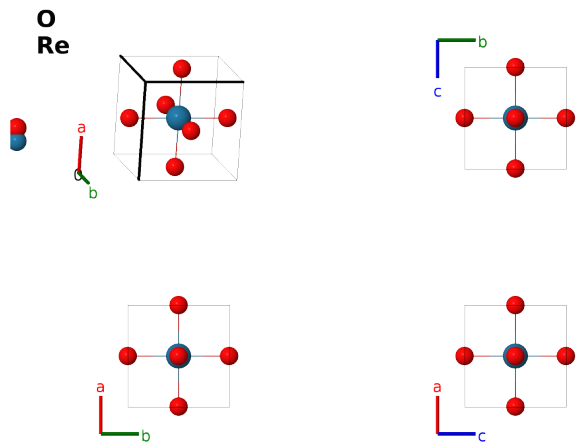
This structure previously used the label(s) **A3B_cP4_221_d_a**. Calls to these addresses will be redirected here.

Cite this page as:

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<https://aflow.org/prototype-encyclopedia/3DRV>

https://aflow.org/prototype-encyclopedia/A3B_cP4_221_c_b-001



Prototype	O ₃ Re
AFLOW prototype label	A3B_cP4_221_c_b-001
<i>Strukturbericht</i> designation	$D0_9$
ICSD	16810
CCDC	1597348
Pearson symbol	cP4
Space group number	221
Space group symbol	$Pm\bar{3}m$
AFLOW prototype command	<code>aflow --proto=A3B_cP4_221_c_b-001 --params=a</code>

Other compounds with this structure

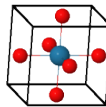
AlF₃, HfF₃, MoF₃, NCu₃, NNa₃, NbF₃, PbI₃, ScF₃, TaF₃, UO₃, WO₃

- Placing an atom on the origin, *aka* the (1a) Wyckoff position, transforms this into the cubic perovskite ($E2_1$) structure.

Simple Cubic primitive vectors

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

$\mathbf{a}_1$
 $\mathbf{a}_2$
 $\mathbf{a}_3$



Basis vectors

	Lattice coordinates		Cartesian coordinates	Wyckoff position	Atom type
$\mathbf{B}_1$	$= \frac{1}{2} \mathbf{a}_1 + \frac{1}{2} \mathbf{a}_2 + \frac{1}{2} \mathbf{a}_3$	$=$	$\frac{1}{2} a \hat{\mathbf{x}} + \frac{1}{2} a \hat{\mathbf{y}} + \frac{1}{2} a \hat{\mathbf{z}}$	(1b)	Re I
$\mathbf{B}_2$	$= \frac{1}{2} \mathbf{a}_2 + \frac{1}{2} \mathbf{a}_3$	$=$	$\frac{1}{2} a \hat{\mathbf{y}} + \frac{1}{2} a \hat{\mathbf{z}}$	(3c)	O I
$\mathbf{B}_3$	$= \frac{1}{2} \mathbf{a}_1 + \frac{1}{2} \mathbf{a}_3$	$=$	$\frac{1}{2} a \hat{\mathbf{x}} + \frac{1}{2} a \hat{\mathbf{z}}$	(3c)	O I
$\mathbf{B}_4$	$= \frac{1}{2} \mathbf{a}_1 + \frac{1}{2} \mathbf{a}_2$	$=$	$\frac{1}{2} a \hat{\mathbf{x}} + \frac{1}{2} a \hat{\mathbf{y}}$	(3c)	O I

References

- [1] K. Meisel, *Rheniumtrioxyd. III. Mitteilung. Über die Kristallstruktur des Rheniumtrioxyds*, Z. Anorganische und Allgemeine Chemie **207**, 121–128 (1932), doi:10.1002/zaac.19322070113.

Found in

- [1] R. T. Downs and M. Hall-Wallace, *The American Mineralogist Crystal Structure Database*, Am. Mineral. **88**, 247–250 (2003).

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