

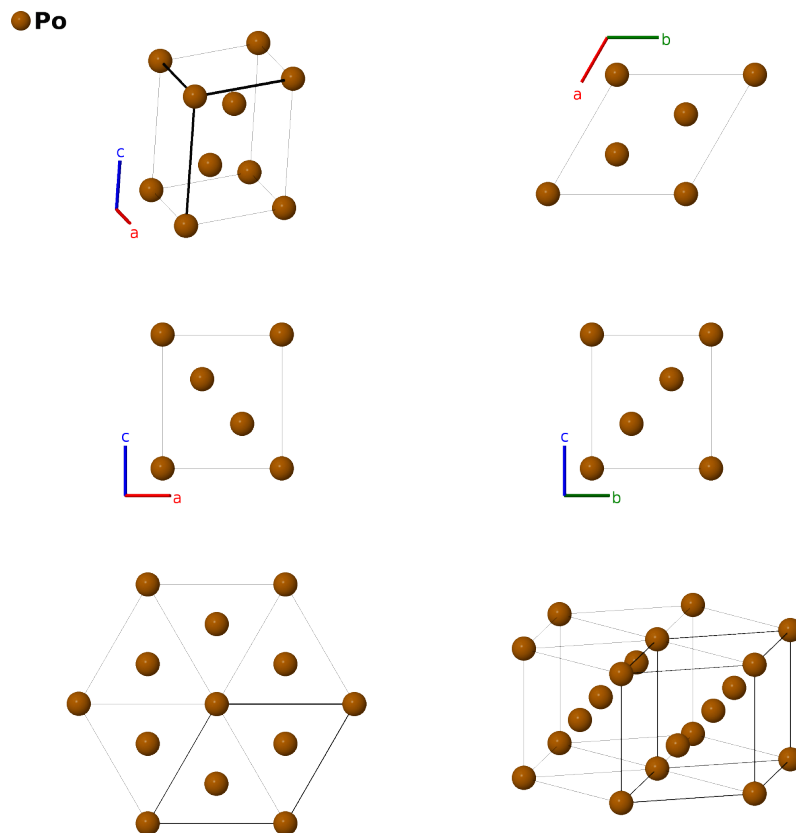
# $\beta$ -Po ( $A_i$ ) Structure: A\_hR1\_166\_a-001

This structure originally had the label A\_hR1\_166\_a.beta-Po. Calls to that address will be redirected here.

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

[https://aflow.org/p/A\\_hR1\\_166\\_a-001](https://aflow.org/p/A_hR1_166_a-001)



Prototype	Po
AFLOW prototype label	A_hR1_166_a-001
<i>Strukturbericht</i> designation	$A_i$
ICSD	649161
Pearson symbol	hR1
Space group number	166
Space group symbol	$R\bar{3}m$
AFLOW prototype command	<code>aflow --proto=A_hR1_166_a-001 --params=a, c/a</code>

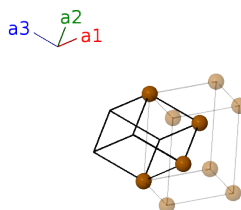
- Originally, Po was assigned Strukturbericht designation A19, which is now considered to be incorrect. (Donohue, 1982, 390)
- This rhombohedral structure becomes cubic at various values of  $c/a$  (or  $\alpha$ )

$c/a$	$\alpha$	Cubic Lattice
$\sqrt{7}$	$60^\circ$	Face-Centered Cubic
$\sqrt{\frac{3}{2}}$	$90^\circ$	Simple Cubic
$\sqrt{\frac{3}{8}}$	$109.47^\circ$	Body-Centered Cubic

- Note that  $\beta$ -Po (A\_hR1\_166\_a) and  $\alpha$ -Hg (A\_hR1\_166\_a) have the same AFLOW prototype label. They are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files. Experimentally,  $\beta$ -Po ( $A_i$ ) has  $c/a$  near 1, or  $\alpha > 90^\circ$ , while  $\alpha$ -Hg (A10) has  $c/a$  near 2, or  $\alpha < 90^\circ$ .
- Hexagonal settings of rhombohedral structures 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
$\mathbf{B}_1$	=	0	=	0	(1a) Po I

### References

- [1] W. H. Beamer and C. L. Maxwell, *Physical Properties of Polonium. II. X-Ray Studies and Crystal Structure*, J. Chem. Phys. **17**, 1293–1298 (1949), doi:10.1063/1.1747155.

### Found in

- [1] J. Donohue, *The Structures of the Elements* (Robert E. Krieger Publishing Company, New York, 1974).