

Hexagonal WO_3 Structure: A3B_hP12_191_gl_f-001

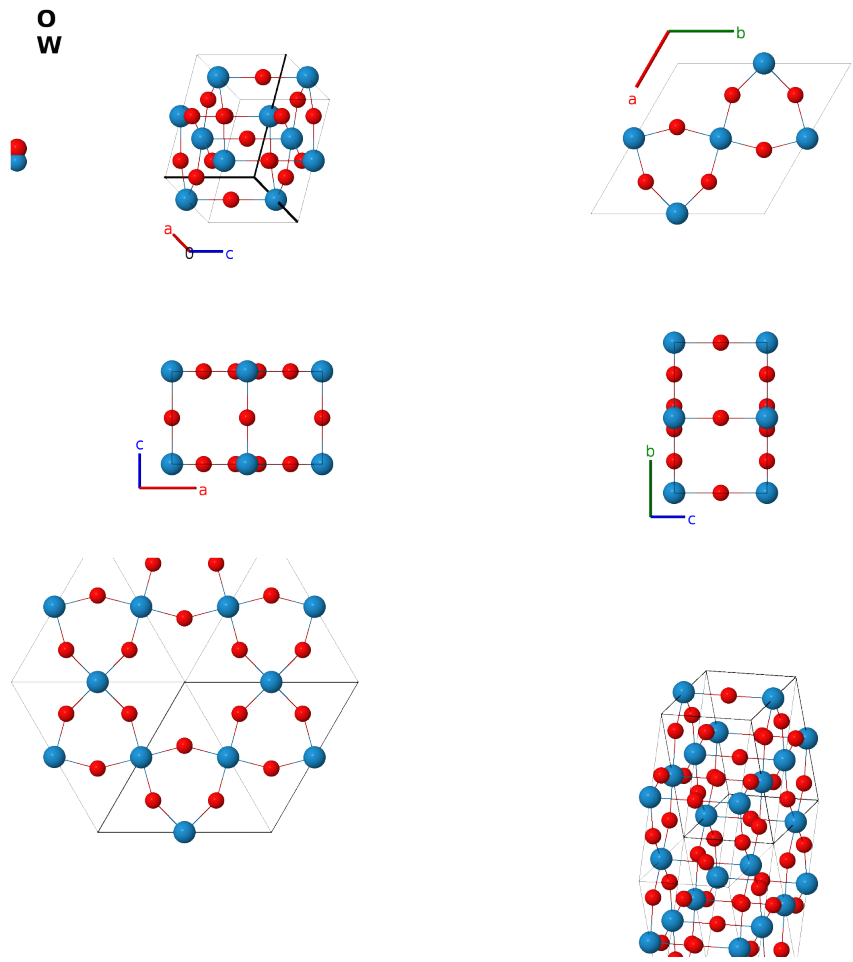
This structure previously used the label(s) **A3B_hP12_191_gl_f**. Calls to these addresses will be redirected here.

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<https://aflow.org/prototype-encyclopedia/SQAX>

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Prototype	O_3W
AFLOW prototype label	A3B_hP12_191_gl_f-001
ICSD	32001
CCDC	1605913
Pearson symbol	hP12
Space group number	191

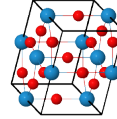
Space group symbol $P6/mmm$

AFLOW prototype command `aflow --proto=A3B_hP12_191_g1_f-001`
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- All stable phases of WO_3 are distortions of the cubic $\alpha\text{-ReO}_3$ ($D0_9$) phase. Based on (Woodward, 1997 and Vogt, 1999), the known stable phases and their approximate temperature ranges are:
 - $\alpha\text{-WO}_3$ (1010-1170 K) (Vogt, 1999)
 - $\beta\text{-WO}_3$ (600-1170 K) (Vogt, 1999)
 - $\gamma\text{-WO}_3$ (290-600 K) (Vogt, 1999)
 - $\delta\text{-WO}_3$ (230-290 K) (Diehl, 1978)
 - $\epsilon\text{-WO}_3$ (below 23 K) (Woodward, 1997)
- Woodward notes that “The transition temperatures display large hysteresis effects and universal agreement is not found in the literature.”
- In addition, several other structures have been proposed and/or found:
 - The original $D0_{10}$ structure (Bräkken, 1931; Hermann, 1937), superseded by $\delta\text{-WO}_3$
 - The original $\beta\text{-WO}_3$ (Salje, 1977)
 - Hexagonal WO_3 , presumably metastable, found by (Gerand, 1979) while dehydrating $\text{WO}_3\cdot\text{H}_2\text{O}$ (this structure)
- (Gerand, 1979) determined the structure of hexagonal WO_3 by X-ray diffraction of powdered samples. They found evidence of super-reflection, indicating that the unit cell should be doubled along the c -axis, but were unable to determine the shift in atomic positions for the double unit cell. We report what they called the “half-cell” structure.

Hexagonal primitive vectors

$$\begin{aligned}\mathbf{a}_1 &= \frac{1}{2}a \hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a \hat{\mathbf{y}} \\ \mathbf{a}_2 &= \frac{1}{2}a \hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a \hat{\mathbf{y}} \\ \mathbf{a}_3 &= c \hat{\mathbf{z}}\end{aligned}$$



Basis vectors

	Lattice coordinates		Cartesian coordinates	Wyckoff position	Atom type
$\mathbf{B}_1$	$= \frac{1}{2} \mathbf{a}_1$	$=$	$\frac{1}{4}a \hat{\mathbf{x}} - \frac{\sqrt{3}}{4}a \hat{\mathbf{y}}$	(3f)	W I
$\mathbf{B}_2$	$= \frac{1}{2} \mathbf{a}_2$	$=$	$\frac{1}{4}a \hat{\mathbf{x}} + \frac{\sqrt{3}}{4}a \hat{\mathbf{y}}$	(3f)	W I
$\mathbf{B}_3$	$= \frac{1}{2} \mathbf{a}_1 + \frac{1}{2} \mathbf{a}_2$	$=$	$\frac{1}{2}a \hat{\mathbf{x}}$	(3f)	W I
$\mathbf{B}_4$	$= \frac{1}{2} \mathbf{a}_1 + \frac{1}{2} \mathbf{a}_3$	$=$	$\frac{1}{4}a \hat{\mathbf{x}} - \frac{\sqrt{3}}{4}a \hat{\mathbf{y}} + \frac{1}{2}c \hat{\mathbf{z}}$	(3g)	O I
$\mathbf{B}_5$	$= \frac{1}{2} \mathbf{a}_2 + \frac{1}{2} \mathbf{a}_3$	$=$	$\frac{1}{4}a \hat{\mathbf{x}} + \frac{\sqrt{3}}{4}a \hat{\mathbf{y}} + \frac{1}{2}c \hat{\mathbf{z}}$	(3g)	O I
$\mathbf{B}_6$	$= \frac{1}{2} \mathbf{a}_1 + \frac{1}{2} \mathbf{a}_2 + \frac{1}{2} \mathbf{a}_3$	$=$	$\frac{1}{2}a \hat{\mathbf{x}} + \frac{1}{2}c \hat{\mathbf{z}}$	(3g)	O I
$\mathbf{B}_7$	$= x_3 \mathbf{a}_1 + 2x_3 \mathbf{a}_2$	$=$	$\frac{3}{2}ax_3 \hat{\mathbf{x}} + \frac{\sqrt{3}}{2}ax_3 \hat{\mathbf{y}}$	(6l)	O II
$\mathbf{B}_8$	$= -2x_3 \mathbf{a}_1 - x_3 \mathbf{a}_2$	$=$	$-\frac{3}{2}ax_3 \hat{\mathbf{x}} + \frac{\sqrt{3}}{2}ax_3 \hat{\mathbf{y}}$	(6l)	O II
$\mathbf{B}_9$	$= x_3 \mathbf{a}_1 - x_3 \mathbf{a}_2$	$=$	$-\sqrt{3}ax_3 \hat{\mathbf{y}}$	(6l)	O II

$$\begin{aligned}
\mathbf{B}_{10} &= -x_3 \mathbf{a}_1 - 2x_3 \mathbf{a}_2 &= -\frac{3}{2}ax_3 \hat{\mathbf{x}} - \frac{\sqrt{3}}{2}ax_3 \hat{\mathbf{y}} & (6l) & \text{O II} \\
\mathbf{B}_{11} &= 2x_3 \mathbf{a}_1 + x_3 \mathbf{a}_2 &= \frac{3}{2}ax_3 \hat{\mathbf{x}} - \frac{\sqrt{3}}{2}ax_3 \hat{\mathbf{y}} & (6l) & \text{O II} \\
\mathbf{B}_{12} &= -x_3 \mathbf{a}_1 + x_3 \mathbf{a}_2 &= \sqrt{3}ax_3 \hat{\mathbf{y}} & (6l) & \text{O II}
\end{aligned}$$

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Found in

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