Hydrophilite ($CaCl_2$, $C35$) Structure: AB2_oP6_58_a_g-001

This structure previously used the label(s) `AB2_oP6_58_a_g.CaCl2`. Calls to these addresses will be redirected here.

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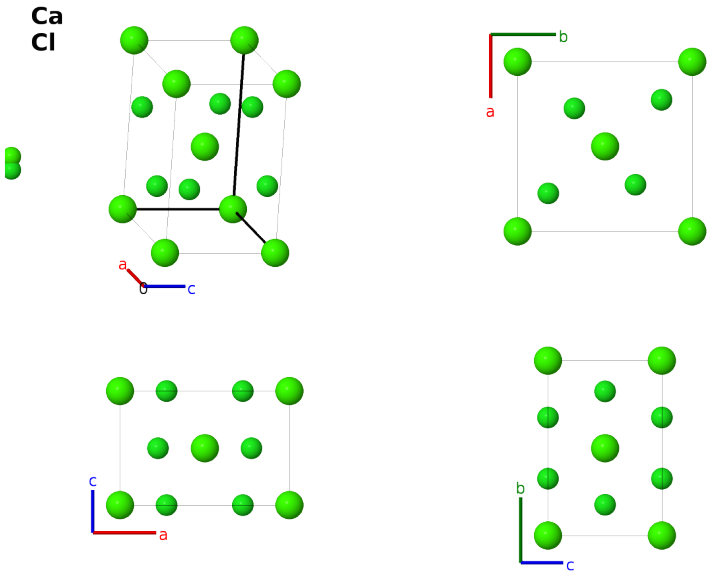
https://aflow.org/prototype-encyclopedia/YTPF

https://aflow.org/prototype-encyclopedia/AB2_oP6_58_a_g-001



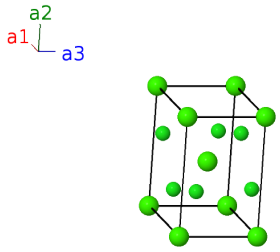
Prototype	$CaCl_2$
AFLOW prototype label	AB2_oP6_58_a_g-001
***Strukturbericht* designation**	$C35$
Mineral name	hydrophilite
ICSD	26686
CCDC	1602076
Pearson symbol	oP6
Space group number	58
Space group symbol	$Pnnm$
AFLOW prototype command	`aflow --proto=AB2_oP6_58_a_g-001 --params=`$a, b/a, c/a, x_2, y_2$

Other compounds with this structure
BeF_2, $CaBr_2$, $CaCl_2$, Co_2C, $CrCl_2$, Co_2N, CrO_2, MgF_2, FeO_2, GeO_2, MgF_2, MnF_2, MnO_2, NiF_2, PbO_2, $PtGe_2$, PtO_2, RuO_2, SiO_2, SnO_2, VO_2, $YbBr_2$, ZrO_2

- Hydrophilite ($CaCl_2$, $C35$), η–Fe_2C, and marcasite (FeS_2, $C18$) have the same AFLOW prototype label, AB2_oP6_58_a_g. They are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Simple Orthorhombic primitive vectors

$$\mathbf{a_1} = a\,\hat{\mathbf{x}}$$
$$\mathbf{a_2} = b\,\hat{\mathbf{y}}$$
$$\mathbf{a_3} = c\,\hat{\mathbf{z}}$$



Basis vectors

		Lattice coordinates		Cartesian coordinates	Wyckoff position	Atom type
$\mathbf{B_1}$	=	0	=	0	(2a)	Ca I
$\mathbf{B_2}$	=	$\frac{1}{2}\,\mathbf{a}_1 + \frac{1}{2}\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$\frac{1}{2}a\,\hat{\mathbf{x}} + \frac{1}{2}b\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(2a)	Ca I
$\mathbf{B_3}$	=	$x_2\,\mathbf{a}_1 + y_2\,\mathbf{a}_2$	=	$ax_2\,\hat{\mathbf{x}} + by_2\,\hat{\mathbf{y}}$	(4g)	Cl I
$\mathbf{B_4}$	=	$-x_2\,\mathbf{a}_1 - y_2\,\mathbf{a}_2$	=	$-ax_2\,\hat{\mathbf{x}} - by_2\,\hat{\mathbf{y}}$	(4g)	Cl I
$\mathbf{B_5}$	=	$-\left(x_2 - \frac{1}{2}\right)\,\mathbf{a}_1 + \left(y_2 + \frac{1}{2}\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$-a\left(x_2 - \frac{1}{2}\right)\,\hat{\mathbf{x}} + b\left(y_2 + \frac{1}{2}\right)\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4g)	Cl I
$\mathbf{B_6}$	=	$\left(x_2 + \frac{1}{2}\right)\,\mathbf{a}_1 - \left(y_2 - \frac{1}{2}\right)\,\mathbf{a}_2 + \frac{1}{2}\,\mathbf{a}_3$	=	$a\left(x_2 + \frac{1}{2}\right)\,\hat{\mathbf{x}} - b\left(y_2 - \frac{1}{2}\right)\,\hat{\mathbf{y}} + \frac{1}{2}c\,\hat{\mathbf{z}}$	(4g)	Cl I

References

[1] A. van Bever and W. Nieuwenkamp, *Die Kristallstruktur von Calciumchlorid, $CaCl_2$*, Z. Kristallogr. **90**, 374–376 (1935), doi:10.1524/zkri.1935.90.1.374.

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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