

$\text{K}_{0.07}\text{Ho}[\text{PO}_4][\text{OH}]_{0.07}$ *hP*24(180) *P*6₂22 – kcdc**HoPO₄:KOH** [1]

Structural features: Distorted HoO₈ square antiprisms are interconnected via common edges and PO₄ tetrahedra to form a 3D-framework; (K,OH) in channels parallel to [001] (partial disorder). Filled-up derivative of hexagonal CePO₄.

Zaki M. et al. (1994) [1]

 $\text{H}_{0.07}\text{HoK}_{0.07}\text{O}_{4.07}\text{P}$ $a = 0.6936$, $c = 0.6418$ nm, $c/a = 0.925$, $V = 0.2674$ nm³, $Z = 3$

site	Wyck.	sym.	<i>x</i>	<i>y</i>	<i>z</i>	occ.	atomic environment
O1	12 <i>k</i>	1	0.311	0.4401	0.20023		single atom P
M2	6 <i>e</i>	2..	0	0	0.05467	0.072	
Ho3	3 <i>d</i>	222	1/2	0	1/2		8-vertex polyhedron O ₈
P4	3 <i>c</i>	222	1/2	0	0		tetrahedron O ₄

 $\text{M2} = 0.5\text{K} + 0.5\text{OH}$

Transformation from published data: origin shift 0 0 1/2

Experimental: powder, diffractometer, X-rays, $R_B = 0.056$

Remarks: Short interatomic distances for partly occupied site(s). Hydrogen atoms are not taken into consideration for Pearson symbol, Wyckoff sequence and atomic environments.

References: [1] Zaki M., Aamili A., El Ghazzi M., Zambon D., Zahir M., Sadel A., Cousseins J.C. (1994), Adv. Mater. Res. (Zuerich) 1/2, 201-208.