

LiAl₂[NO₃][OH]₆*hP*32(176) *P*6₃/*m* – ih²fba**[LiAl₂(OH)₆]NO₃ [1]**

Structural features: Slabs of edge-linked Li(OH)₆ and Al(OH)₆ octahedra alternate with layers containing NO₃ trigonal units (perpendicular to [001], partial rotational disorder).

Besserguenev A.V. et al. (1997) [1]

Al₂H₆LiNO₉

a = 0.51092, *c* = 1.43738 nm, *c/a* = 2.813, *V* = 0.3249 nm³, *Z* = 2

site	Wyck.	sym.	<i>x</i>	<i>y</i>	<i>z</i>	occ.	atomic environment
(OH)1	12 <i>i</i>	1	0.004	0.3639	0.0717		non-coplanar triangle Al ₂ Li
O2	6 <i>h</i>	<i>m</i> ..	0.063	0.266	¹ / ₄	0.153	non-coplanar triangle NO ₂
O3	6 <i>h</i>	<i>m</i> ..	0.266	0.203	¹ / ₄	0.847	non-coplanar triangle NO ₂
Al4	4 <i>f</i>	3..	¹ / ₃	² / ₃	0.0		octahedron (OH) ₆
Li5	2 <i>b</i>	-3..	0	0	0		octahedron (OH) ₆
N6	2 <i>a</i>	-6..	0	0	¹ / ₄		coplanar hexagon O ₆

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

Experimental: powder, diffractometer, X-rays, R_p = 0.054

Remarks: Hydrogen atoms are not taken into consideration for Pearson symbol, Wyckoff sequence and atomic environments.

References: [1] Besserguenev A.V., Fogg A.M., Francis R.J., Price S.J., O'Hare D., Isupov V.P., Tolochko B.P. (1997), Chem. Mater. 9, 241-247.