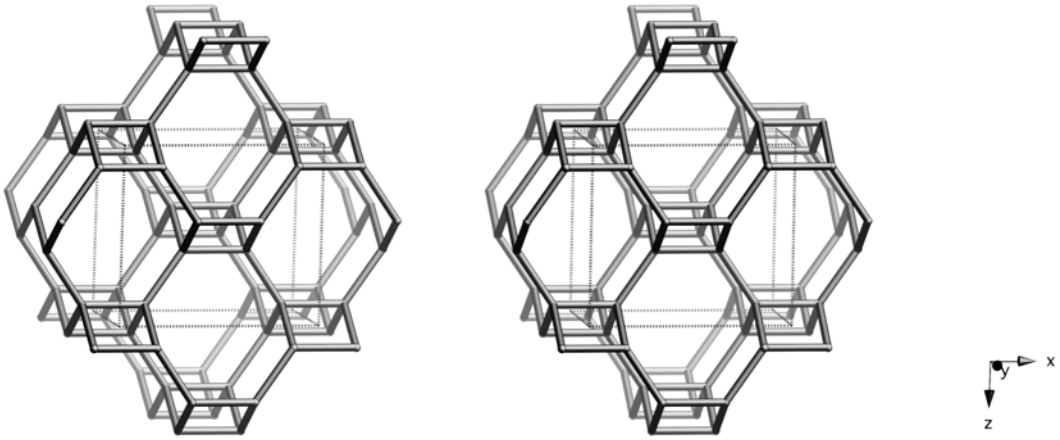
ABW

Imma

Framework Type Data



framework viewed along [010]

Idealized cell data: orthorhombic, *Imma*, $a = 9.9$Å, $b = 5.3$Å, $c = 8.8$Å

Coordination sequences and vertex symbols:

T_1 (8,*m*)	4	10	21	36	54	78	106	136	173	214	$4{\cdot}6{\cdot}4{\cdot}6{\cdot}6{\cdot}8_2$

Secondary building units: 8 or 4

Composite building units:

dzc

double zigzag chain

abw

Materials with this framework type:

*Li-A (Barrer and White)[1-3]
[Be-As-O]-**ABW**[4,5]
[Be-P-O]-**ABW**[4,6,7]
[Ga-Si-O]-**ABW**[8]
[Zn-As-O]-**ABW**[4,9]
[Zn-P-O]-**ABW**[4]
$|(NH_4)$-|[Co-P-O]-**ABW**[10]
$|(NH_4)$-|[Zn-As-O]-**ABW**[11]
$|(NH_4)$-|[Zn-P-O]-**ABW**[12]
|Cs-|[Mg-P-O]-**ABW**[13]
|Cs-|[Al-Si-O]-**ABW**[14,15]
|Cs-|[Al-Ti-O]-**ABW**[16]
|Li-|[Zn-As-O]-**ABW**[17]
|Li-|[Al-Si-O]-**ABW**[18]
|Li-|[Zn-P-O]-**ABW**[19]
|Li-|[Al-Ge-O]-**ABW**[20]
|Na-|[Zn-P-O]-**ABW**[21]
|Na-|[Co-P-O]-**ABW**[22]
|Rb-|[Cu-P-O]-**ABW**[23]
|Rb-|[Ni-P-O]-**ABW**[24]
|Rb-|[Co-P-O]-**ABW**[13]
|Rb-|[Al-Si-O]-**ABW**[14,15]
|Tl-|[Al-Si-O]-**ABW**[25]
UCSB-3 [26]

Type Material: Li-A (Barrer and White) **ABW**

Type Material Data

Crystal chemical data: |Li_4 $(H_2O)_4$| [$Al_4Si_4O_{16}$]-**ABW**
orthorhombic, *Pna*2_1, a = 10.31Å, b = 8.18Å, c = 5.00Å [(2)]
(Relationship to unit cell of Framework Type: $a' = a$, $b' = c$, $c' = b$)

Framework density: 19 T/1000Å^3

Channels: [001] **8** 3.4 x 3.8*

8-ring viewed along [001]

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