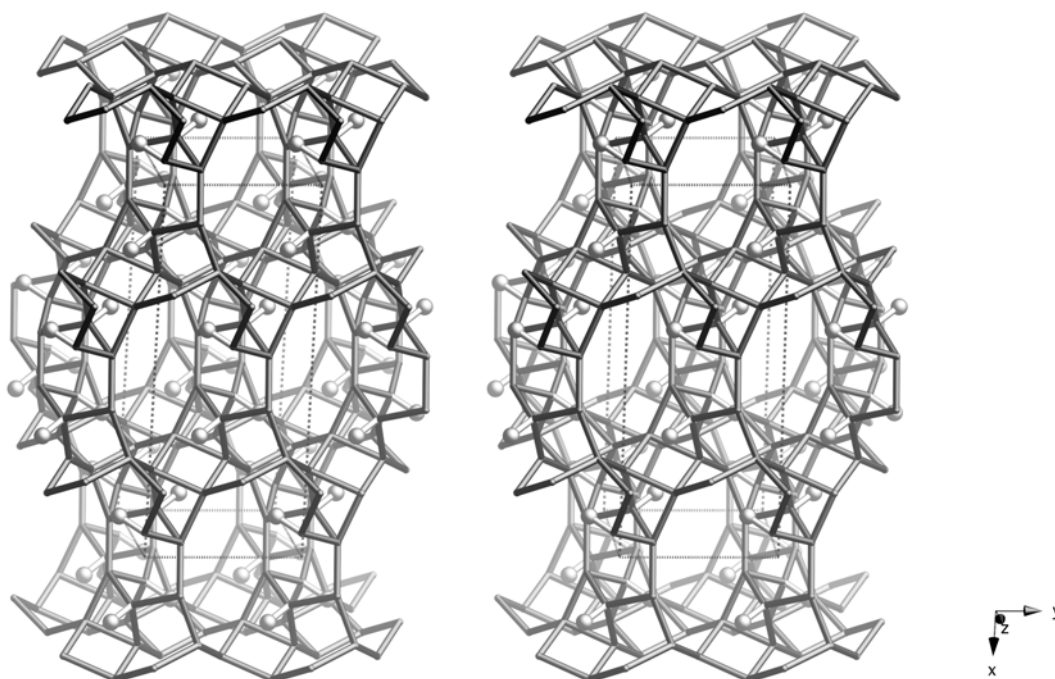


Framework Type Data



framework viewed along [001]

Idealized cell data: monoclinic, $C2/c$, $a = 20.9\text{\AA}$, $b = 9.2\text{\AA}$, $c = 8.6\text{\AA}$, $\beta = 89.7^\circ$

Coordination sequences and vertex symbols:

$T_1(8,1)$	4	9	18	31	48	71	99	132	162	197	$4\cdot6\cdot6\cdot6\cdot10_2$
$T_2(8,1)$	4	9	19	33	51	77	96	126	162	203	$4\cdot6\cdot4\cdot8_2\cdot6\cdot6_2$
$T_3(8,1)$	4	10	19	33	52	72	102	126	161	204	$4\cdot6_2\cdot4\cdot8_2\cdot8_2\cdot10$
$T_4(8,1)$	3	8	16	29	49	68	94	123	162	203	$4\cdot6\cdot6_2$

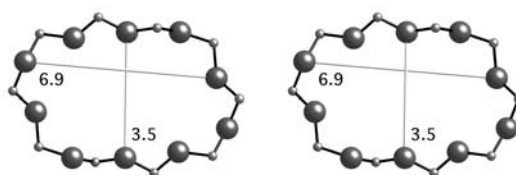
Secondary building units: 4

Materials with this framework type:

*Partheite⁽¹⁾

Type Material Data

Crystal chemical data:	$[\text{Ca}_8(\text{H}_2\text{O})_{16}] [\text{Al}_{16}\text{Si}_{16}\text{O}_{60}(\text{OH})_8]$ - PAR monoclinic, $C2/c$ $a = 21.555\text{\AA}$, $b = 8.761\text{\AA}$, $c = 9.304\text{\AA}$, $\beta = 91.55^\circ$ ⁽¹⁾
Stability:	Stable at 150°C, transforms at 400°C ⁽¹⁾
Framework density:	18.2 T/1000Å ³
Channels:	[001] 10 3.5 x 6.9*



10-ring viewed along [001]

References:

- (1) Engel, N. and Yvon, K. *Z. Kristallogr.*, **169**, 165-175 (1984)