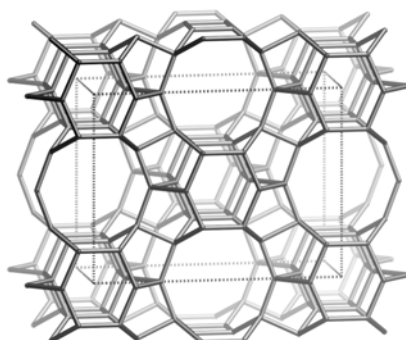
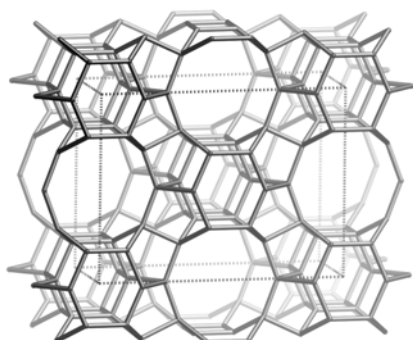


Framework Type Data



framework viewed along [001]

Idealized cell data: orthorhombic, *Cmmm*, $a = 18.9\text{\AA}$, $b = 14.4\text{\AA}$, $c = 7.5\text{\AA}$

Coordination sequences and vertex symbols:

$T_1 (16,1)$	4	10	21	38	61	88	116	149	190	237	289	340	$4\cdot5\cdot4\cdot6\cdot6\cdot8_2$
$T_2 (8,..m)$	4	12	23	38	59	82	119	155	191	231	283	338	$5\cdot8\cdot5\cdot8\cdot5_2\cdot6$
$T_3 (8,..m)$	4	10	19	34	60	89	118	144	179	238	296	342	$4\cdot5\cdot4\cdot5\cdot6\cdot8_2$
$T_4 (4,2mm)$	4	12	24	38	52	88	118	150	192	232	278	340	$5\cdot5\cdot5\cdot5\cdot10_2\cdot10_4$

Secondary building units: 6

Composite building units:

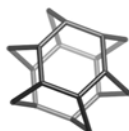
mtt



d6r



mso

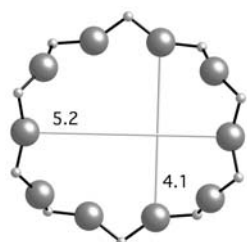


Materials with this framework type:

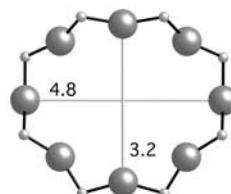
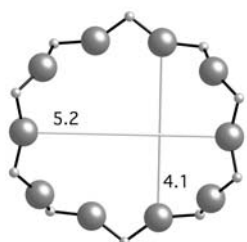
*SUZ-4^(1,2)

Type Material Data

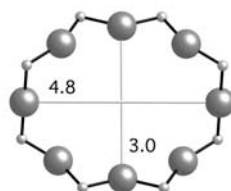
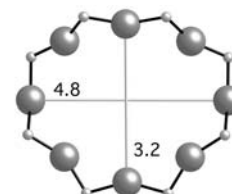
Crystal chemical data:	$\text{IK}_4\text{I} [\text{Al}_4\text{Si}_{32}\text{O}_{72}]$ -SZR orthorhombic, $Cmmm$, $a = 18.8064\text{\AA}$, $b = 14.2298\text{\AA}$, $c = 7.4548\text{\AA}$ ⁽²⁾
Framework density:	18 T/1000 \AA^3
Channels:	{ [001] 10 4.1 x 5.2 \leftrightarrow [010] 8 3.2 x 4.8 \leftrightarrow [110] 8 3.0 x 4.8 }***



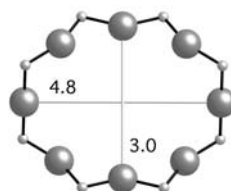
10-ring viewed along [001]



8-ring viewed along [010]



8-ring viewed along [110]



References:

- (1) Lawton, S.L., Bennett, J.M., Schlenker, J.L. and Rubin, M.K. *Chem. Commun.*, 894-896 (1993)
- (2) Strohmaier, K.G., Afeworki, M. and Dorset, D.L. *Z. Kristallogr.*, **221**, 689-698 (2006)