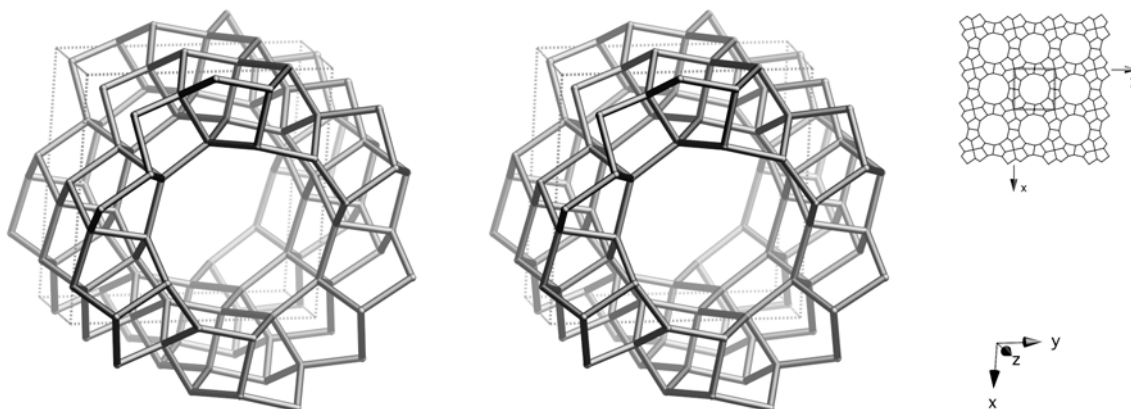


Framework Type Data



framework viewed along [001] (upper right: projection down [001])

Idealized cell data: tetragonal, $P\bar{4}$, $a = 13.0\text{\AA}$, $c = 4.9\text{\AA}$

Coordination sequences and vertex symbols:

$T_1(4,1)$	4	12	24	39	61	93	133	179	209	246	$5\cdot5\cdot5\cdot6_2\cdot5\cdot7$
$T_2(4,1)$	4	12	26	41	65	94	130	169	218	269	$5\cdot5\cdot5\cdot6\cdot7\cdot12_6$
$T_3(4,1)$	4	12	24	41	65	95	128	169	218	270	$5\cdot6\cdot5\cdot6\cdot5\cdot6_2$
$T_4(4,1)$	4	12	23	43	68	94	125	172	226	269	$5\cdot6\cdot5\cdot6\cdot5_2\cdot6$
$T_5(1,4)$	4	12	28	38	60	98	152	182	200	246	$5\cdot5\cdot5\cdot5\cdot8_2\cdot8_2$

Secondary building units: see *Compendium*

Composite building units:

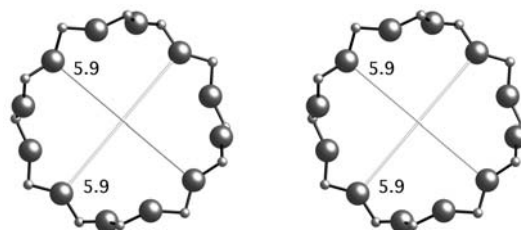
cas

**Materials with this framework type:**

*VPI-8⁽¹⁾

Type Material Data

Crystal chemical data:	[Si ₁₇ O ₃₄]-VET tetragonal, $P\bar{4}$, $a = 13.045 \text{ \AA}$, $c = 5.034 \text{ \AA}$ ⁽¹⁾
Framework density:	19.8 T/1000 \AA^3
Channels:	[001] 12 5.9 x 5.9*



12-ring viewed along [001]

References:

- (1) Freyhardt, C.C., Lobo, R.F., Khodabandeh, S., Lewis, J.E., Tsapatsis, M., Yoshikawa, M., Cambor, M.A., Pan, M., Helmkamp, M.M., Zones, S.I. and Davis, M.E. *J. Am. Chem. Soc.*, **118**, 7299-7310 (1996)