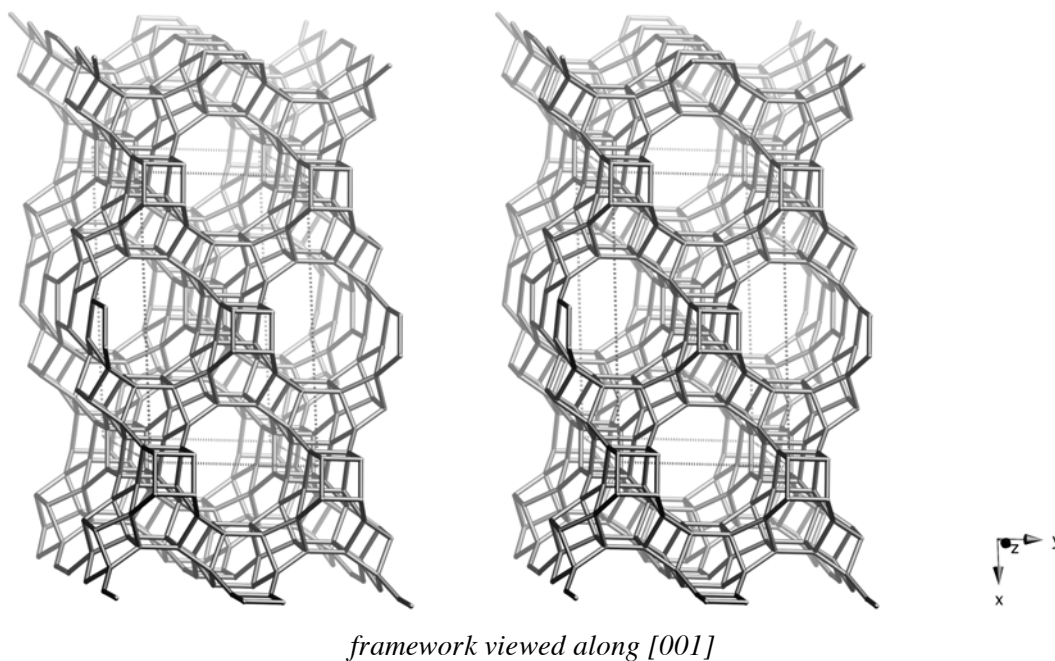


## Framework Type Data

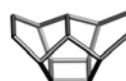


**Idealized cell data:** monoclinic,  $C2/m$ ,  $a = 22.7\text{\AA}$ ,  $b = 13.4\text{\AA}$ ,  $c = 12.6\text{\AA}$ ,  $\beta = 69.5^\circ$

**Coordination sequences and vertex symbols:**

$T_1(8,1)$	4	10	19	32	50	73	101	132	164	199	$4\cdot5\cdot4\cdot10_4\cdot5\cdot6_2$
$T_2(8,1)$	4	11	18	31	52	77	98	126	164	205	$4\cdot5\cdot5\cdot6\cdot5\cdot10_2$
$T_3(8,1)$	4	10	21	32	47	74	105	134	159	196	$4\cdot6\cdot4\cdot10_4\cdot5\cdot6_2$
$T_4(8,1)$	4	10	19	32	51	74	100	130	165	203	$4\cdot4\cdot5\cdot6_2\cdot5\cdot12_7$
$T_5(8,1)$	4	10	19	32	51	74	101	130	164	203	$4\cdot4\cdot5\cdot6_2\cdot5\cdot12_4$
$T_6(8,1)$	4	11	18	28	49	77	103	126	155	201	$4\cdot5_2\cdot5\cdot6_2\cdot6\cdot6_2$
$T_7(8,1)$	4	11	19	32	49	75	105	131	159	196	$4\cdot6_2\cdot5\cdot6\cdot5\cdot10_2$

**Secondary building units:** 1-5-1

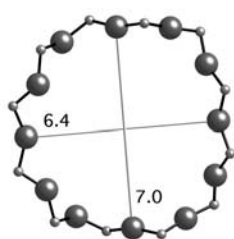
**Composite building units:***bea**bre**lau**mel***Materials with this framework type:**\*CIT-1<sup>(1)</sup>SSZ-26 (related to CON)<sup>(1,2)</sup>SSZ-33 (related to CON)<sup>(1,2)</sup>

## Type Material Data

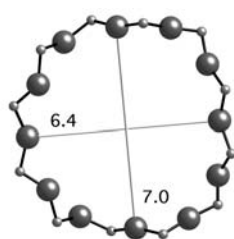
**Crystal chemical data:**  $[\text{H}_2] [\text{B}_2\text{Si}_{54}\text{O}_{112}]$ -CON  
 monoclinic,  $C2/m$   
 $a = 22.624\text{\AA}$ ,  $b = 13.350\text{\AA}$ ,  $c = 12.364\text{\AA}$ ,  $\beta = 68.91^\circ$  <sup>(1)</sup>

**Framework density:**  $16.1 \text{ T}/1000\text{\AA}^3$

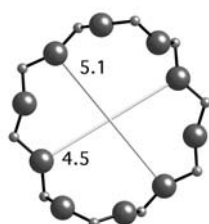
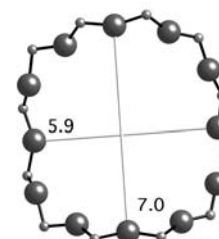
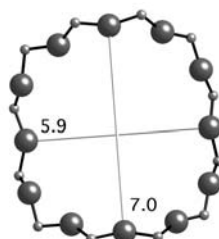
**Channels:**  $[001] \mathbf{12} \ 6.4 \times 7.0^* \leftrightarrow [100] \mathbf{12} \ 7.0 \times 5.9^* \leftrightarrow [010] \mathbf{10} \ 5.1 \times 4.5^*$



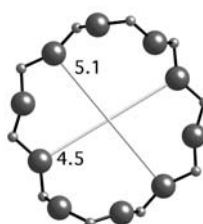
*12-ring viewed along [001]*



*12-ring viewed along [100]*



*10-ring viewed along [010]*

**References:**

- (1) Lobo, R.F. and Davis, M.E. *J. Am. Chem. Soc.*, **117**, 3764-3779 (1995)
- (2) Lobo, R.F., Pan, M., Chan, I., Li, H.X., Medrud, R.C., Zones, S.I., Crozier, P.A. and Davis, M.E. *Science*, **262**, 1543-1546 (1993)