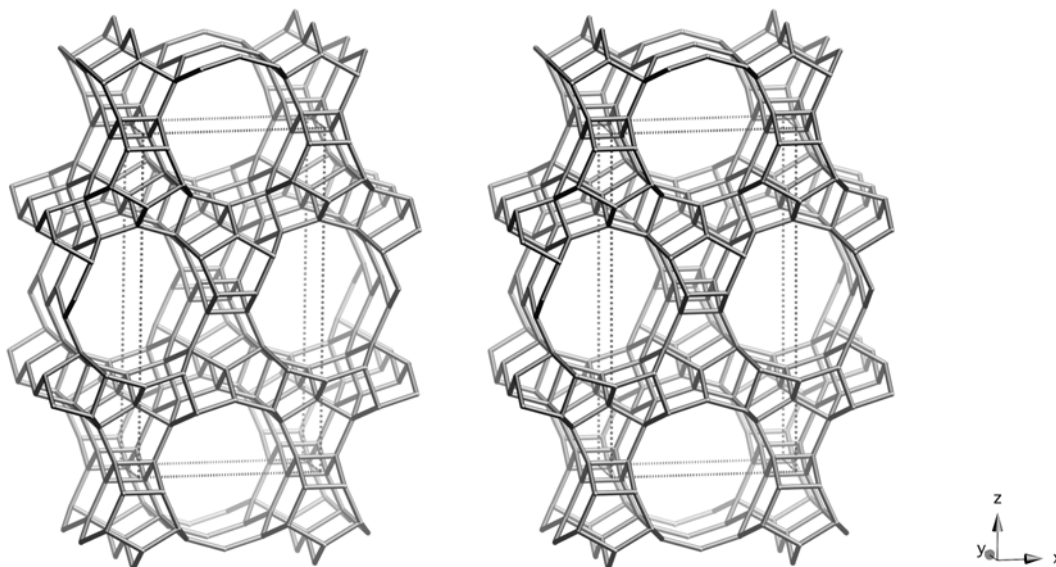


## Framework Type Data



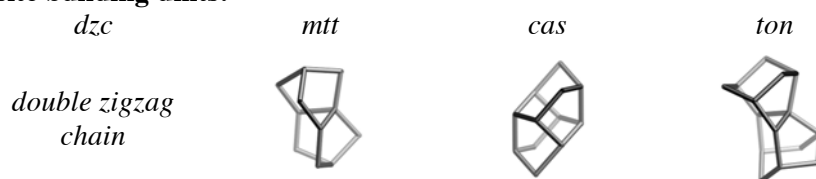
framework viewed along [010]

**Idealized cell data:** orthorhombic, *Imma*,  $a = 14.0\text{\AA}$ ,  $b = 5.3\text{\AA}$ ,  $c = 26.0\text{\AA}$

**Coordination sequences and vertex symbols:**

$T_1(8,m)$	4	10	21	36	56	84	114	143	182	231	4-6-4-6-5-6
$T_2(8,m)$	4	12	23	37	55	83	114	153	195	222	5-6-5-6-5 <sub>2</sub> -6
$T_3(8,m)$	4	12	22	37	57	84	114	156	184	222	5-6-5-6-5-6 <sub>2</sub>
$T_4(4,mm2)$	4	12	24	36	54	79	118	153	190	234	5-5-5-5-14 <sub>14</sub> -*
$T_5(4,mm2)$	4	12	20	34	56	81	116	151	186	220	5-5-5-5-5-6 <sub>2</sub>

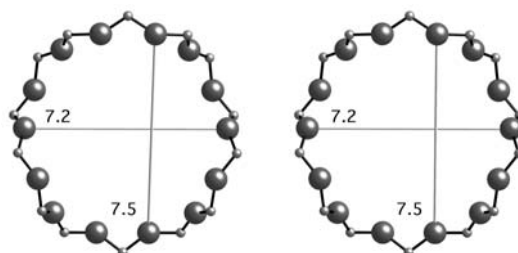
**Secondary building units:** see *Compendium*

**Composite building units:****Materials with this framework type:**

\*CIT-5<sup>(1,2)</sup>

## Type Material Data

<b>Crystal chemical data:</b>	[Si <sub>32</sub> O <sub>64</sub> ]-CFI orthorhombic, $Pmn2_1$ , $a = 13.674\text{\AA}$ , $b = 5.022\text{\AA}$ , $c = 25.488\text{\AA}$ <sup>(2)</sup>
<b>Framework density:</b>	18.3 T/1000 $\text{\AA}^3$
<b>Channels:</b>	[010] <b>14</b> 7.2 x 7.5*



*14-ring viewed along [010]*

**References:**

- (1) Wagner, P., Yoshikawa, M., Lovallo, M., Tsuji, K., Taspatsis, M. and Davis, M.E. *Chem. Commun.*, 2179-2180 (1997)
- (2) Yoshikawa, M., Wagner, P., Lovallo, M., Tsuji, K., Takewaki, T., Chen, C.Y., Beck, L.W., Jones, C., Taspatsis, M., Zones, S.I. and Davis, M.E. *J. Phys. Chem. B*, **102**, 7139-7147 (1998)