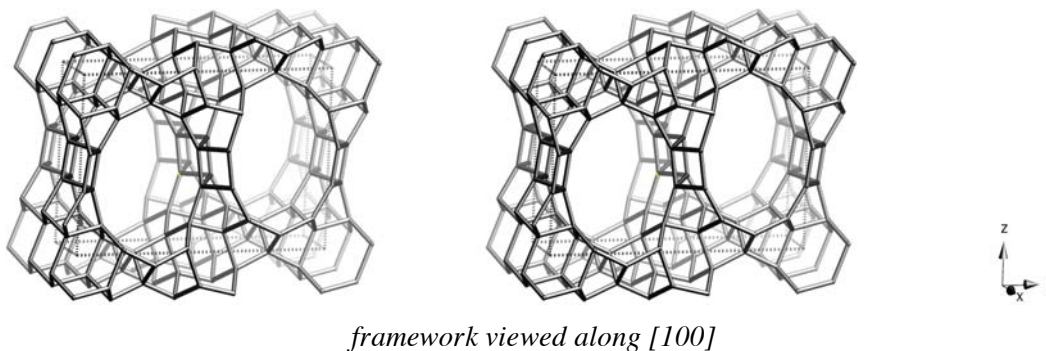


## Framework Type Data

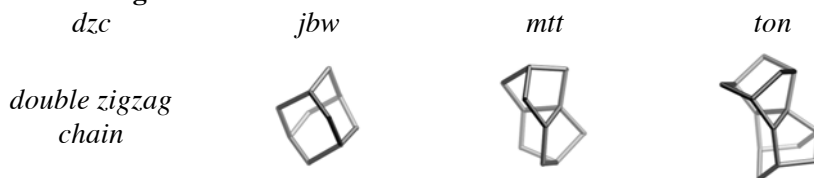


**Idealized cell data:** orthorhombic, *Pmmn* (origin choice 2),  $a = 5.3\text{\AA}$ ,  $b = 22.6\text{\AA}$ ,  $c = 14.0\text{\AA}$

**Coordination sequences and vertex symbols:**

T <sub>1</sub> (4, <i>m</i> ..)	4	12	20	36	59	84	115	152	183	236	286	344	5·5·5·5·6 <sub>2</sub> ·12 <sub>2</sub>
T <sub>2</sub> (4, <i>m</i> ..)	4	10	19	35	57	84	113	147	179	229	289	337	4·5·4·5·6·12 <sub>2</sub>
T <sub>3</sub> (4, <i>m</i> ..)	4	12	24	38	57	79	116	156	193	233	280	330	5 <sub>2</sub> ·6 <sub>2</sub> ·6·6 <sub>2</sub> ·6·6 <sub>2</sub>
T <sub>4</sub> (4, <i>m</i> ..)	4	10	20	35	57	83	115	146	183	227	282	348	4·5·4·5·12 <sub>2</sub> ·*
T <sub>5</sub> (4, <i>m</i> ..)	4	12	22	37	57	87	117	152	185	234	286	345	5 <sub>2</sub> ·6 <sub>2</sub> ·6·6 <sub>2</sub> ·6·6 <sub>2</sub>
T <sub>6</sub> (4, <i>m</i> ..)	4	12	22	37	55	87	120	148	180	234	294	344	5·5·5·5·6 <sub>2</sub> ·*
T <sub>7</sub> (2, <i>mm</i> 2)	4	12	24	44	56	78	110	158	206	234	262	336	5·5·5·5·12 <sub>2</sub> ·*
T <sub>8</sub> (2, <i>mm</i> 2)	4	12	22	38	60	78	114	150	200	232	280	332	5·5·5·5·6 <sub>2</sub> ·12 <sub>2</sub>

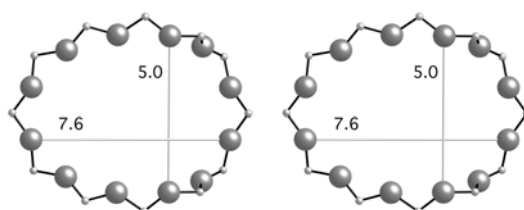
**Secondary building units:** see *Compendium*

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

\*SSZ-60<sup>(1)</sup>

## Type Material Data

<b>Crystal chemical data:</b>	[B <sub>0.6</sub> Si <sub>27.4</sub> O <sub>56</sub> ]-SSY orthorhombic, $Pmn2_1$ , $a = 21.951\text{\AA}$ , $b = 13.698\text{\AA}$ , $c = 5.012\text{\AA}$ <sup>(1)</sup> (Space group changed to standard setting) (Relationship to unit cell of Framework Type: $a' = b$ , $b' = c$ , $c' = a$ )
<b>Framework density:</b>	18.6 T/1000 $\text{\AA}^3$
<b>Channels:</b>	[001] 12 5.0 x 7.6*



12-ring viewed along [001]

**References:**

- (1) Burton, A. and Elomari, S. *Chem. Commun.*, 2618-2619 (2004)