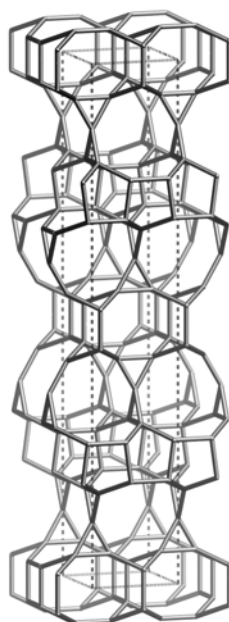
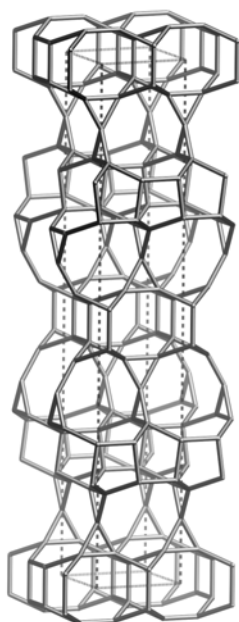


## Framework Type Data



framework viewed along [001]

**Idealized cell data:** orthorhombic, *Cmmm*,  $a = 7.2\text{\AA}$ ,  $b = 41.9\text{\AA}$ ,  $c = 7.2\text{\AA}$

**Coordination sequences and vertex symbols:**

T <sub>1</sub> (8,..m)	4	9	19	39	59	78	117	155	189	235	288	330	3·4·8 <sub>3</sub> ·9 <sub>4</sub> ·8 <sub>5</sub> ·9 <sub>4</sub>
T <sub>2</sub> (8,m..)	4	10	21	37	58	91	117	144	194	241	287	343	4·4·6 <sub>2</sub> ·8·6 <sub>2</sub> ·8
T <sub>3</sub> (8,m..)	4	9	21	42	57	82	119	151	188	239	277	345	3·4·8 <sub>2</sub> ·9 <sub>4</sub> ·8 <sub>2</sub> ·9 <sub>4</sub>
T <sub>4</sub> (8,..m)	4	11	21	40	61	89	116	145	191	239	294	339	4·5 <sub>2</sub> ·5·8·5·8
T <sub>5</sub> (4,m2m)	4	8	20	44	55	80	118	152	204	228	272	354	3·3·9 <sub>4</sub> ·9 <sub>4</sub> ·9 <sub>4</sub> ·9 <sub>4</sub>

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

**Composite building units:**

*lov*



*vsv*

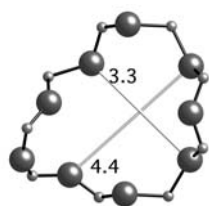
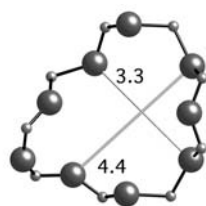
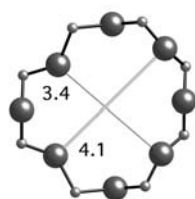
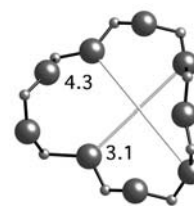
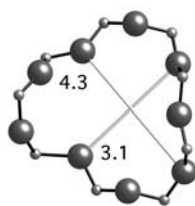
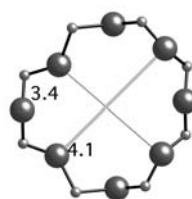


**Materials with this framework type:**

\*RUB-17<sup>(1)</sup>

## Type Material Data

<b>Crystal chemical data:</b>	$\text{K}_4\text{Na}_{12}(\text{H}_2\text{O})_{18}[\text{Zn}_8\text{Si}_{28}\text{O}_{72}]$ -RSN monoclinic, $Cm$ , $a = 7.238\text{\AA}$ , $b = 40.56\text{\AA}$ , $c = 7.308\text{\AA}$ , $\beta = 91.8^\circ$ <sup>(1)</sup>
<b>Stability:</b>	Complete dehydration leads to destruction of the framework <sup>(1)</sup>
<b>Framework density:</b>	$16.8 \text{ T}/1000\text{\AA}^3$
<b>Channels:</b>	$[100] \text{ 9 } 3.3 \times 4.4^* \leftrightarrow [001] \text{ 9 } 3.1 \times 4.3^* \leftrightarrow [010] \text{ 8 } 3.4 \times 4.1^*$

9-ring viewed along  $[100]$ 9-ring viewed along  $[001]$ 8-ring viewed along  $[010]$ **References:**

- (1) Röhrig, C. and Gies, H. *Angew. Chem., Int. Ed.*, **34**, 63-65 (1995)