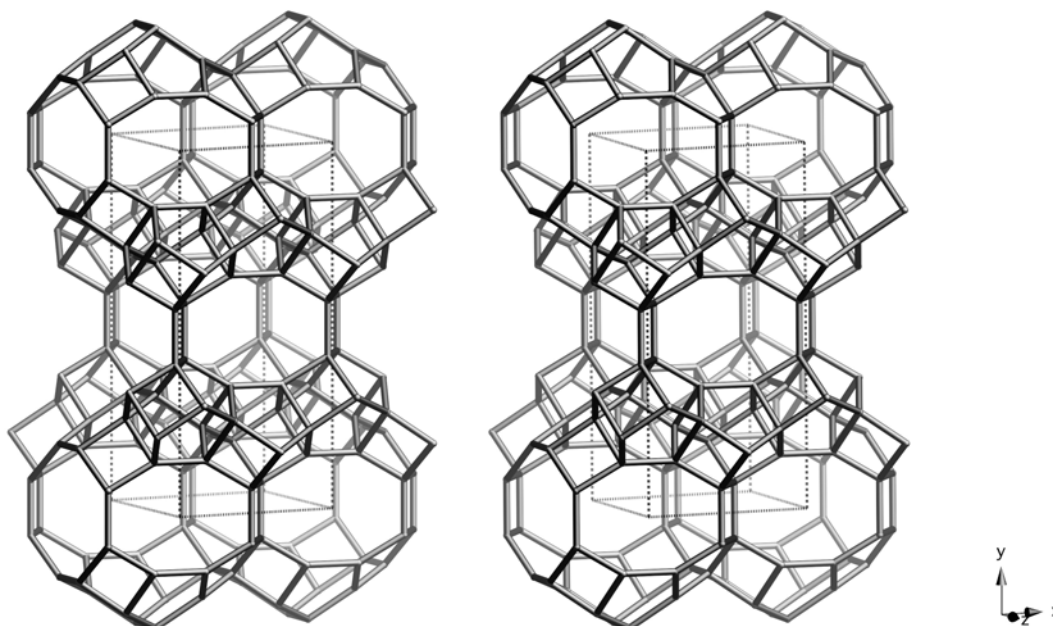


Framework Type Data



framework viewed along [001]

Idealized cell data: monoclinic, $C2/m$, $a = 9.8\text{\AA}$, $b = 20.5\text{\AA}$, $c = 10.0\text{\AA}$, $\beta = 96.9^\circ$

Coordination sequences and vertex symbols:

| | | | | | | | | | | | |
|------------|---|----|----|----|----|----|-----|-----|-----|-----|-------------|
| $T_1(8,1)$ | 4 | 11 | 21 | 34 | 53 | 78 | 108 | 137 | 165 | 207 | 4-6-5-6-5-8 |
| $T_2(8,1)$ | 4 | 10 | 21 | 36 | 54 | 75 | 100 | 136 | 181 | 214 | 4-4-5-8-5-8 |
| $T_3(8,1)$ | 4 | 10 | 19 | 31 | 50 | 82 | 106 | 130 | 168 | 203 | 4-5-4-6-5-5 |
| $T_4(8,1)$ | 4 | 10 | 18 | 31 | 55 | 77 | 103 | 134 | 165 | 214 | 4-5-4-8-5-5 |

Secondary building units: 4

Composite building units:

rth



Materials with this framework type:

*RUB-13⁽¹⁾

SSZ-36 (ITE-RTH structural intermediate)⁽²⁾

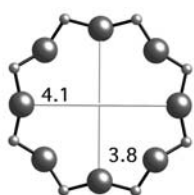
SSZ-50 ([Al-Si-O]-RTH)⁽³⁾

Type Material Data

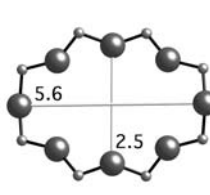
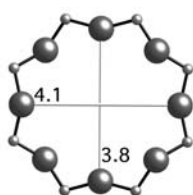
Crystal chemical data: I (C₁₀H₂₁N)₂I [B₂Si₃₀O₆₄]-RTH
 C₁₀H₂₁N = pentamethylpiperidinium
 monoclinic, C2/m
 $a = 9.659\text{\AA}$, $b = 20.461\text{\AA}$, $c = 9.831\text{\AA}$, $\beta = 96.58^\circ$ ⁽¹⁾

Framework density: 16.6 T/1000Å³

Channels: [100] 8 3.8 x 4.1* ↔ [001] 8 2.5 x 5.6*



8-ring viewed along [100]



8-ring viewed along [001]

References:

- (1) Vortmann, S., Marler, B., Gies, H. and Daniels, P. *Microporous Materials*, **4**, 111-121 (1995)
- (2) Wagner, P., Nakagawa, Y., Lee, G.S., Davis, M.E., Elomari, S., Medrud, R.C. and Zones, S.I. *J. Am. Chem. Soc.*, **122**, 263-273 (2000)
- (3) Lee, G.S. and Zones, S.I. *J. Solid State Chem.*, **167**, 289-298 (2002)