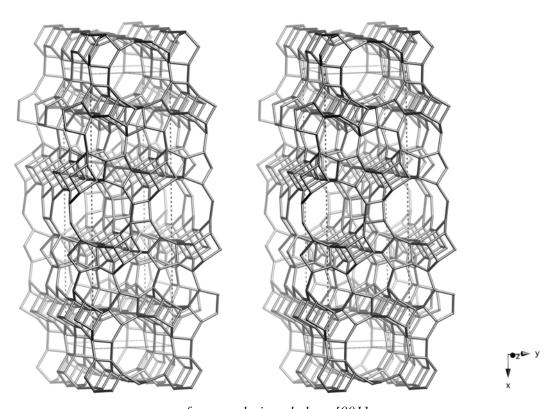
IWW Pbam

Framework Type Data



framework viewed along [001]

orthorhombic, *Pbam*, a = 41.7Å, b = 12.7Å, c = 12.7Å Idealized cell data:

Coordination sequences and vertex symbols:

see Appendix A for a list of the coordination sequences and vertex symbols for the 16 T-atoms

Secondary building units: 1-5-1

Composite building units:

brelau melstf



IWW

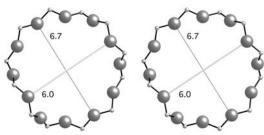
Type Material Data

Crystal chemical data: $[Ge_{22.2}Si_{89.8}O_{224}]$ -IWW

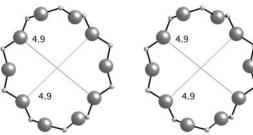
orthorhombic, *Pbam*, $a = 42.1326\text{\AA}$, $b = 12.9885\text{\AA}$, $c = 12.6814\text{\AA}^{(1)}$

Framework density: 16.1 T/1000Å³

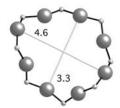
Channels: [001] **12** 6.0 x 6.7* $\leftrightarrow \bot$ [001] **10** 4.9 x 4.9** \leftrightarrow [001] **8** 3.3 x 4.6*



12-ring viewed along [001]



10-ring viewed normal to [001]



3.3

8-ring viewed allong [001]

References:

(1) Corma, A., Rey, F., Valencia, S., Jorda, J.L. and Rius, J. Nature Materials, 2, 493-497 (2003)