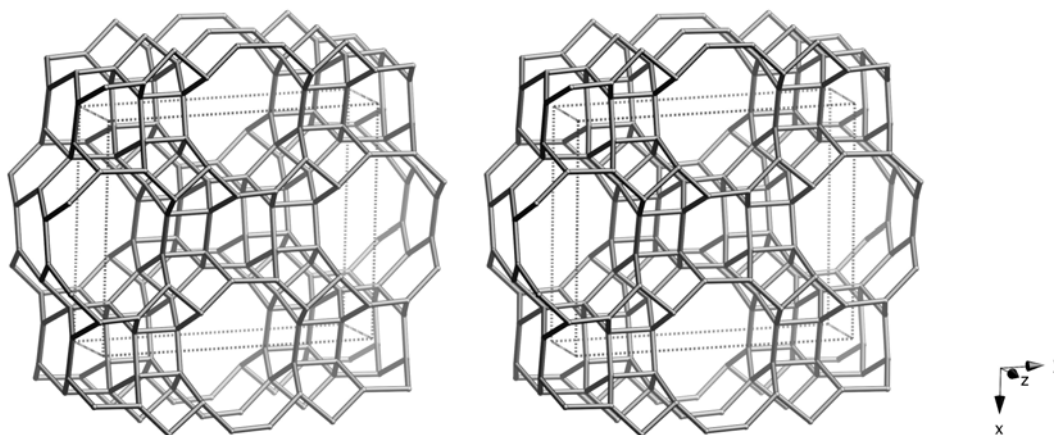


Framework Type Data



framework viewed along [001]

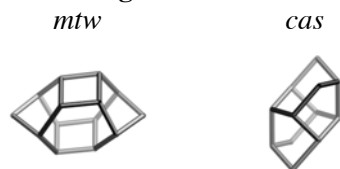
Idealized cell data: orthorhombic, *Cmmm*, $a = 16.9\text{\AA}$, $b = 20.4\text{\AA}$, $c = 5.3\text{\AA}$

Coordination sequences and vertex symbols:

$T_1(8,m)$	4	12	21	39	63	85	117	154	192	242	$5\cdot6\cdot5\cdot6\cdot5_2\cdot6$
$T_2(8,m)$	4	12	25	38	57	86	119	158	194	233	$5\cdot6\cdot5\cdot6\cdot6\cdot12_6$
$T_3(8,m)$	4	11	22	38	58	86	121	156	191	229	$4\cdot6_2\cdot5\cdot6\cdot5\cdot6$
$T_4(8,m)$	4	11	22	38	63	91	115	147	195	244	$4\cdot6_2\cdot5\cdot6\cdot5\cdot6$

Secondary building units: 5-3

Composite building units:

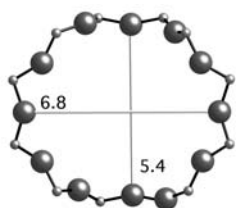


Materials with this framework type:

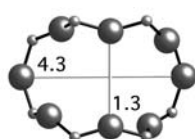
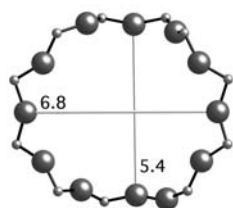
*GUS-1⁽¹⁾

Type Material Data

Crystal chemical data:	[Si ₃₂ O ₆₄]-GON orthorhombic, C222, $a = 16.421\text{\AA}$, $b = 20.054\text{\AA}$, $c = 5.046\text{\AA}$ ⁽¹⁾
Framework density:	19.3 T/1000Å ³
Channels:	[001] 12 5.4 x 6.8*



12-ring viewed along [001]



8-ring viewed along [001]

**References:**

- (1) Plévert, J., Kubota, Y., Honda, T., Okubo, T. and Sugi, Y. *Chem. Commun.*, 2363-2364 (2000)