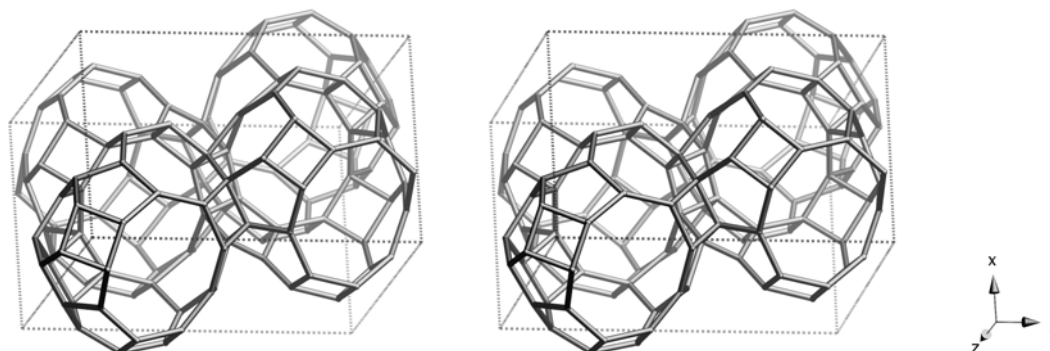


Framework Type Data



framework viewed normal to [100]

Idealized cell data: monoclinic, $P2_1/n$, $a = 13.1\text{\AA}$, $b = 21.9\text{\AA}$, $c = 13.6\text{\AA}$, $\beta = 102.9^\circ$

Coordination sequences and vertex symbols:

T ₁ (4,1)	4	11	22	36	55	84	111	142	179	233	4-5-5-7-6-9
T ₂ (4,1)	4	10	23	37	57	78	108	146	187	225	4-4-5-9-6-7
T ₃ (4,1)	4	11	20	35	54	84	108	142	178	225	4-6-5-6-5-7
T ₄ (4,1)	4	10	19	32	55	81	109	141	174	223	4-5-4-6-5-5
T ₅ (4,1)	4	9	21	38	57	78	104	144	195	220	4-4-4-9-5-7
T ₆ (4,1)	4	10	20	40	57	78	105	143	191	226	4-5-4-7-5-9
T ₇ (4,1)	4	10	20	35	57	77	108	139	184	230	4-6-4-6-5-7
T ₈ (4,1)	4	10	19	35	54	77	108	146	174	215	4-5-4-9-5-6
T ₉ (4,1)	4	11	19	34	53	81	115	140	173	219	4-6-5-5-5-5
T ₁₀ (4,1)	4	10	21	31	51	83	115	134	172	216	4-6-4-6-5-5
T ₁₁ (4,1)	4	12	20	31	50	88	117	137	167	221	5-5 ₂ -5-6-6-6
T ₁₂ (4,1)	4	10	18	35	59	77	107	150	179	221	4-5-4-9-5-5
T ₁₃ (4,1)	4	10	20	38	55	75	104	149	185	218	4-4-5-6-5-9
T ₁₄ (4,1)	4	12	21	33	58	86	112	141	174	226	5-5-5-6-5-7
T ₁₅ (4,1)	4	11	21	33	55	78	111	145	175	217	4-5-5-6-5-9
T ₁₆ (4,1)	4	11	20	32	54	89	115	138	172	221	4-6 ₂ -5-5-5-5

Secondary building units: 5-3

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

*SSZ-23⁽¹⁾

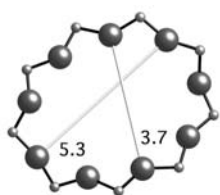
Type Material: SSZ-23

Type Material Data

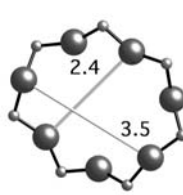
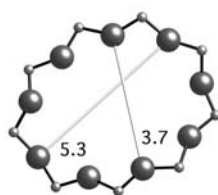
Crystal chemical data: $[(C_{13}H_{24}N)_{4.1} F_{3.3}(OH)_{0.8}] [Si_{64}O_{128}]$ -STT
 $C_{13}H_{24}N = N,N,N$ -trimethyl-1-adamantammonium
 monoclinic, $P2_1/n$
 $a = 12.959\text{\AA}$, $b = 21.792\text{\AA}$, $c = 13.598\text{\AA}$, $\beta = 101.85^\circ$ ⁽¹⁾

Framework density: 17 T/1000 \AA^3

Channels: [101] 9 3.7 x 5.3* \leftrightarrow [001] 7 2.4 x 3.5*



9-ring viewed along [101]



7-ring viewed along [001]

References:

- (1) Cambor, M.A., Diaz-Cabanas, M.-J., Perez-Pariente, J., Teat, S.J., Clegg, W., Shannon, I.J., Lightfoot, P., Wright, P.A. and Morris, R.E. *Angew. Chem., Int. Ed.*, **37**, 2122-2126 (1998)