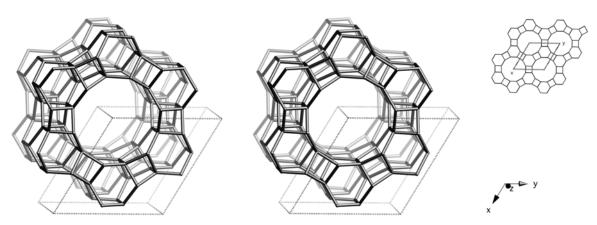
AFI P6/mcc

Framework Type Data



framework viewed along [001] (upper right: projection down [001])

Idealized cell data: hexagonal, P6/mcc, a = 13.8Å, c = 8.6Å

Coordination sequences and vertex symbols:

 $T_1(24,1)$ 4 11 21 35 53 77 105 137 172 212

 $4.6_{2}.6.6_{3}.6_{2}.6_{3}$

Secondary building units: 12 or 6 or 4

Composite building units:

nsc afi

narsarsukite chain



bog



Materials with this framework type:

*AlPO-5⁽¹⁾
[Sn-Al-P-O]-**AFI**⁽²⁾
l(Ni(deta)₂)-l[Al-P-O-F-]-**AFI**⁽³⁾
lTPAF-l[Al-P-O]-**AFI**⁽⁴⁾
CoAPO-5^(5,6)
CrAPO-5⁽⁷⁾
FAPO-5⁽⁸⁾

$$\begin{split} MAPO-5,\,M&=Cd,\,Cu,\,Mo,\,V/Mo,\,Zr^{(9)}\\ MAPO-5,\,M&=Mg,\,Mn^{(10)}\\ SAPO-5 \text{ and numerous compositional variants}^{(11,12)}\\ SSZ-24^{(13)}\\ VAPO-5^{(14)}\\ ZnAPO-5^{(15)} \end{split}$$

Type Material: AIPO-5

Type Material Data

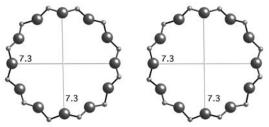
Crystal chemical data: $I(C_{12}H_{28}N) (H_2O)_x (OH)I [Al_{12}P_{12} O_{48}]$ -AFI

 $C_{12}H_{28}N = tetrapropylammonium$

hexagonal, P6cc, a = 13.726Å, c = 8.484Å⁽¹⁾

Framework density: 17.3 T/1000Å³

Channels: [001] **12** 7.3 x 7.3*



12-ring viewed along [001]

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