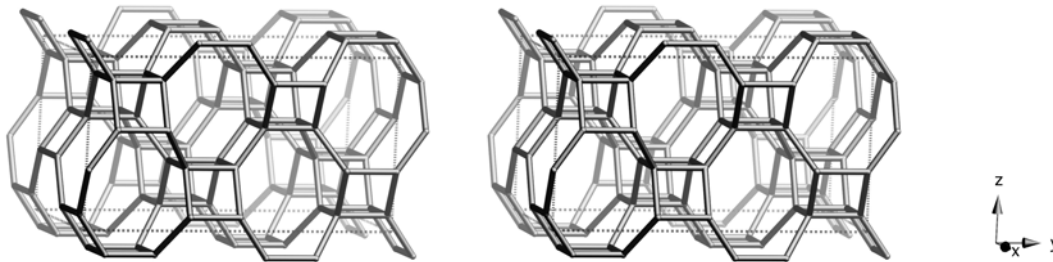


## Framework Type Data



*framework viewed along [100]*

**Idealized cell data:** orthorhombic, *Cmce*,  $a = 9.0\text{\AA}$ ,  $b = 19.4\text{\AA}$ ,  $c = 10.4\text{\AA}$

**Coordination sequences and vertex symbols:**

$T_1(16,1)$	4	9	19	35	53	75	102	132	168	208	$4\cdot4\cdot4\cdot8_2\cdot8\cdot8_2$
$T_2(16,1)$	4	10	20	35	54	76	104	136	171	211	$4\cdot6\cdot4\cdot6\cdot6\cdot8_2$

**Secondary building units:** 8 or 4

**Composite building units:**

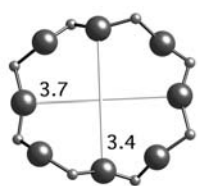
*dcc*  
*double*  
*crankshaft chain*

**Materials with this framework type:**

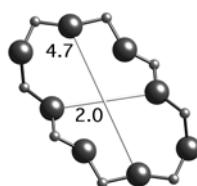
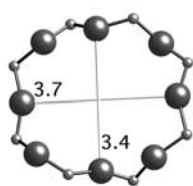
\*AIPO-C<sup>(1,2)</sup>  
 AIPO-H3<sup>(3)</sup>  
 CoAPO-H3<sup>(4)</sup>

## Type Material Data

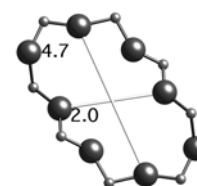
<b>Crystal chemical data:</b>	[Al <sub>16</sub> P <sub>16</sub> O <sub>64</sub> ]-APC orthorhombic, <i>Pbca</i> , $a = 19.821\text{Å}$ , $b = 10.028\text{Å}$ , $c = 8.936\text{Å}$ <sup>(2)</sup> (Relationship to unit cell of Framework Type: $a' = b$ , $b' = c$ , $c' = a$ )
<b>Stability:</b>	Transforms to AIPO-D at ca 250°C <sup>(2)</sup>
<b>Framework density:</b>	18 T/1000Å <sup>3</sup>
<b>Channels:</b>	[001] 8 3.4 x 3.7* ↔ [100] 8 2.0 x 4.7*



8-ring viewed along [001]



distorted 8-ring viewed along [100]

**References:**

- (1) Bennett, J.M., Dytrych, W.J., Pluth, J.J., Richardson Jr., J.W. and Smith, J.V. *Zeolites*, **6**, 349-359 (1986)
- (2) Keller, E.B., Meier, W.M. and Kirchner, R.M. *Solid State Ionics*, **43**, 93-102 (1990)
- (3) Pluth, J.J. and Smith, J.V. *Acta Crystallogr.*, **C42**, 1118-1120 (1986)
- (4) Canesson, L., Arcon, I., Caldarelli, S. and Tuel, A. *Microporous Mesoporous Mat.*, **26**, 117-131 (1998)