Molecular sieves = highly organized matrices of tunable pore shape, size, and polarity for separation, recognition, and organization of molecules with precision of about 1 Å.

detergent builders adsorbents size-shape selective catalysts supramolecular chemistry nanotechnology

**Chemical composition** 

Silica SiO<sub>2</sub>

Aluminosilicates  $M_x^{IA} l_x Si_{2-x} O_4 . nH_2 O$ 

Aluminophosphates AlPO<sub>4</sub> (isoelectronic with Si<sub>2</sub>O<sub>4</sub>)

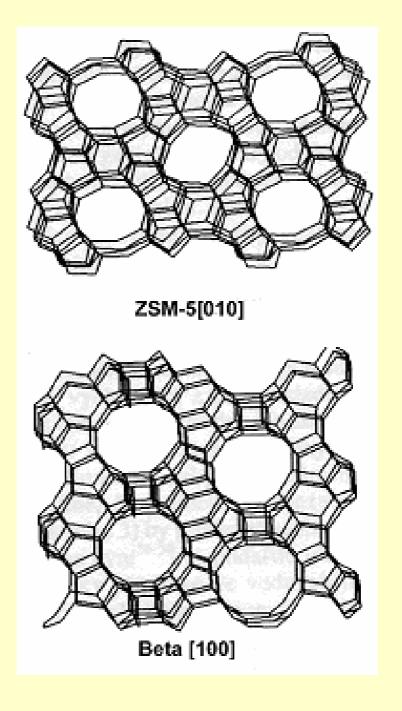
Metallophosphates MPO<sub>4</sub>

Silicoaluminophosphates M<sub>x</sub><sup>I</sup>Si<sub>x</sub>AlP<sub>1-x</sub>O<sub>4</sub>



Pores

Channels



- >40 naturally occurring zeolites
- >139 structure types many hundreds of zeolite compounds

Nomenclature www.iza-structure.org/databases

Structure types - three capital letter codes (Most well known zeolite archetypes: SOD, LTA, FAU, MOR, MFI)

**Four-connected frameworks** 

Interrupted frameworks (denoted by a hyphen: –CLO, cloverite) Structure types do not depend on: chemical composition, element distribution, cell dimensions, symmetry

Several zeolite compounds can belong to the same structure type: FAU – faujasite, Linde X, Y, Beryllophosphate-X, SAPO-37, Zincophosphate-X

Names of zeolite materials:

trivial names - Alpha, Beta, Rho

chemical names – Gallogermanate-A

mineral names – Chabazite, Mordenite, Stilbite, Sodalite

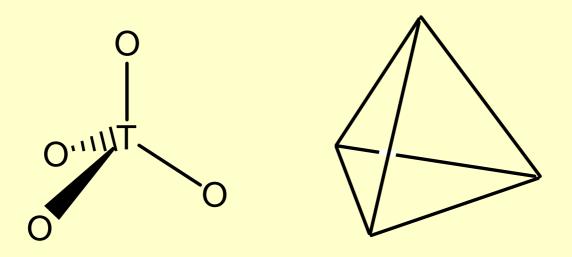
codes – AlPO4-5, 8, 11, ..., 54, ZSM-4, 18, 57, ...

brand names – Linde A, D, F, L, N, Q, R, T, W, X, Y

university names
VPI-5 (Virginia Polytechnical Institute)
ULM (University Le Mans)

**Primary building units:** 

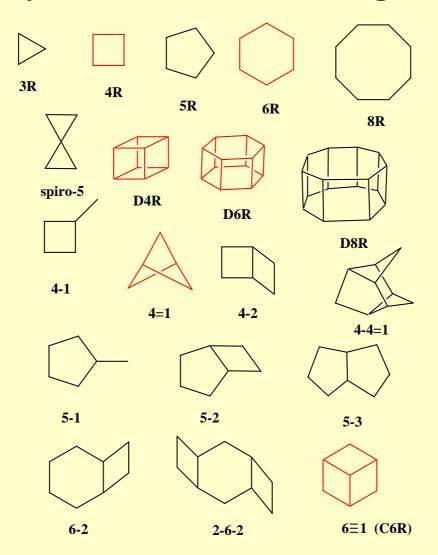
Al(III)O<sub>4</sub>, P(V)O<sub>4</sub> and Si(IV)O<sub>4</sub> tetrahedra



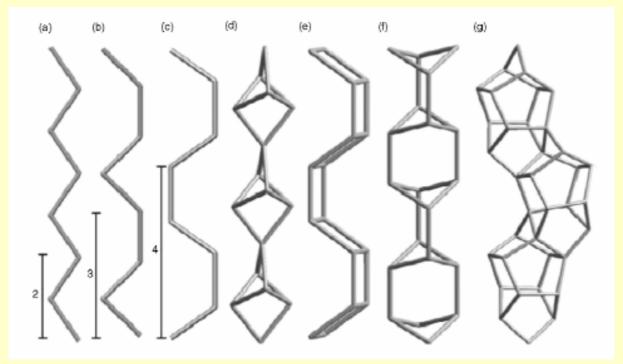
#### **Isoelectronic relationship**

 $(SiO_2)_2$   $[AlSiO_4]^ AlPO_4$ 

#### Secondary (Structural) Building Units (SBU)

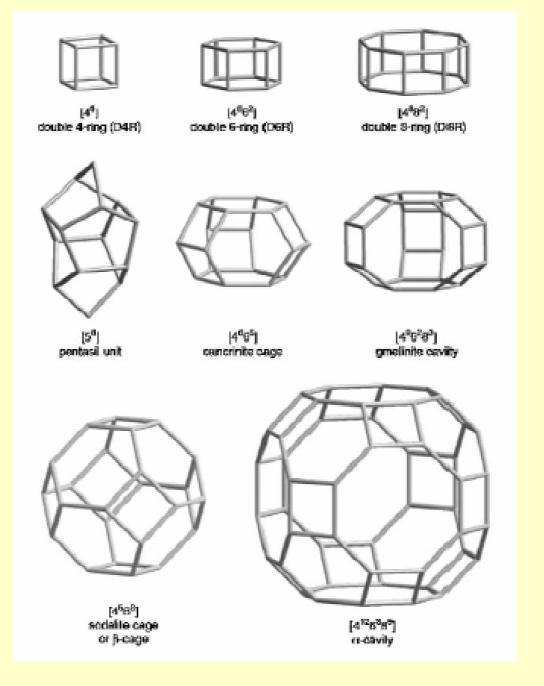


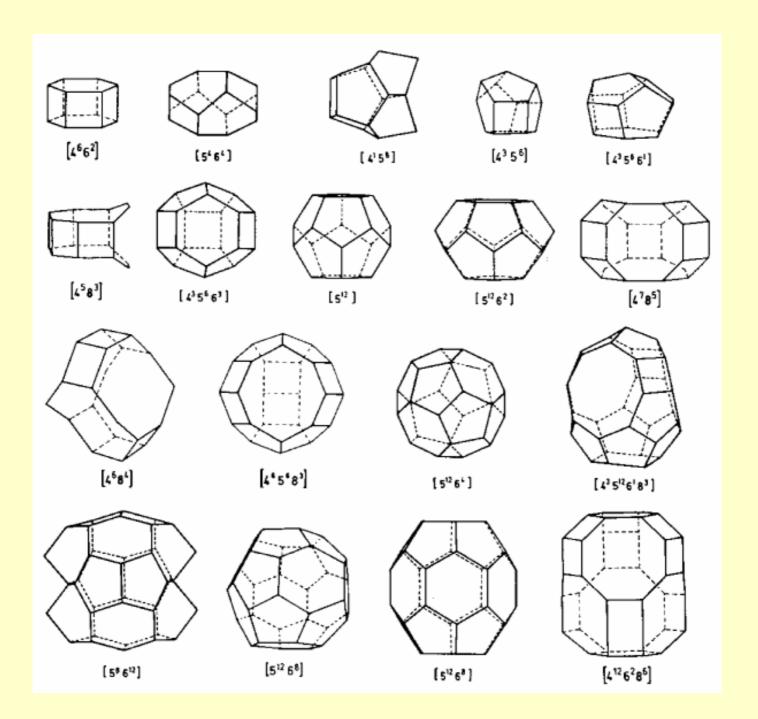
#### Chain composite building units



- (a) zig-zag unbranched single chain, periodicity of two
- (b) sawtooth unbranched single chain, periodicity of three
- (c) crankshaft unbranched single chain, periodicity of four
- (d) natrolite branched single chain
- (e) double crankshaft chain, an unbranched double chain
- (f) narsarsukite chain, a branched double chain
- (g) a pentasil chain

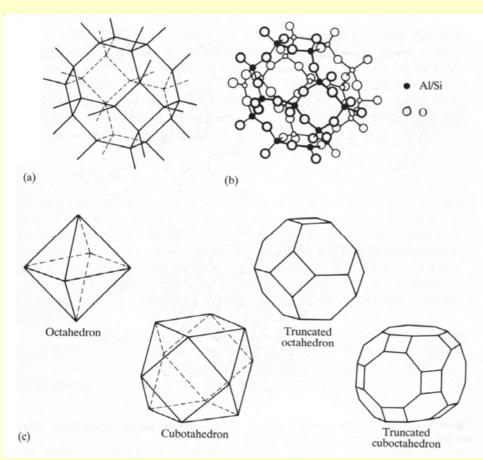
# Polyhedral composite building units





#### **Sodalite Unit**

#### Truncated octahedron





#### **Sodalite Unit**

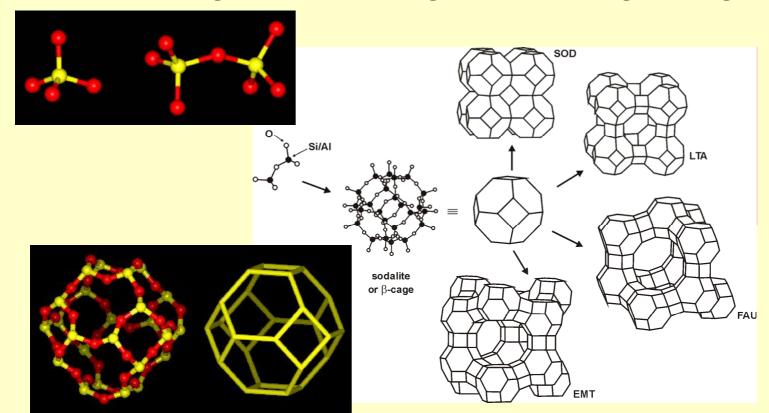
Packing of the sodalite units:

**SOD** – bcc, sharing of 4-rings

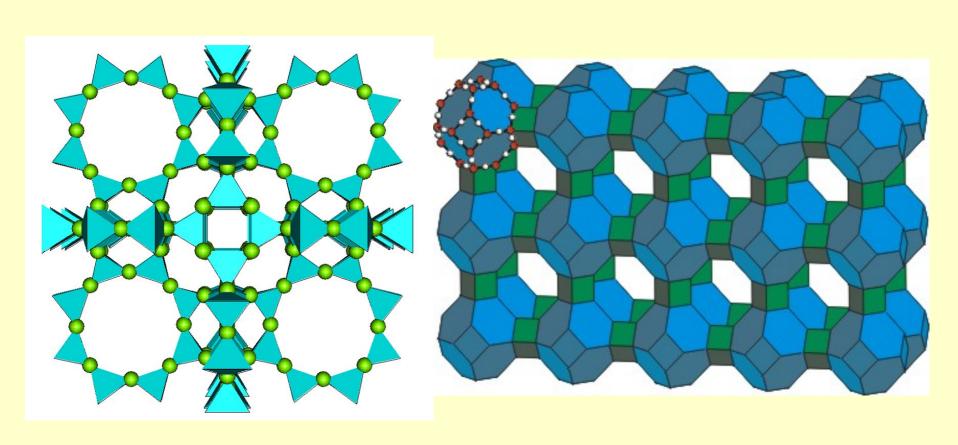
LTA – sc, 4-rings connected through O bridges

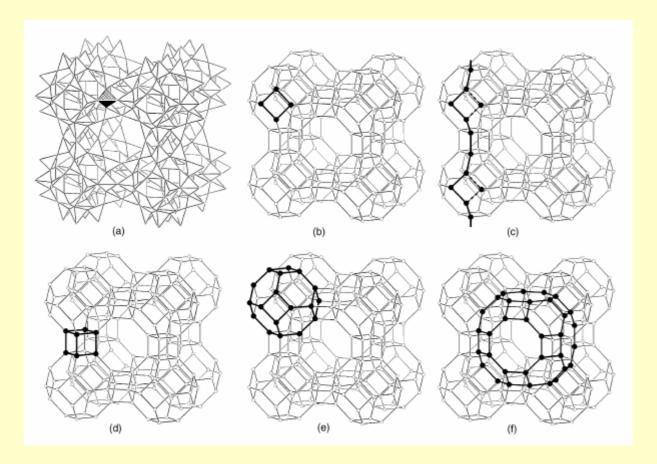
FAU (faujasite) – cubic diamond, 6-rings connected through O bridges

EMT – hexagonal diamond, 6-rings connected through O bridges



# **Zeolite LTA**

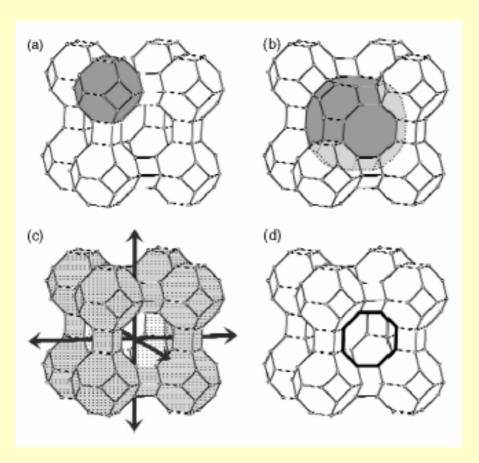




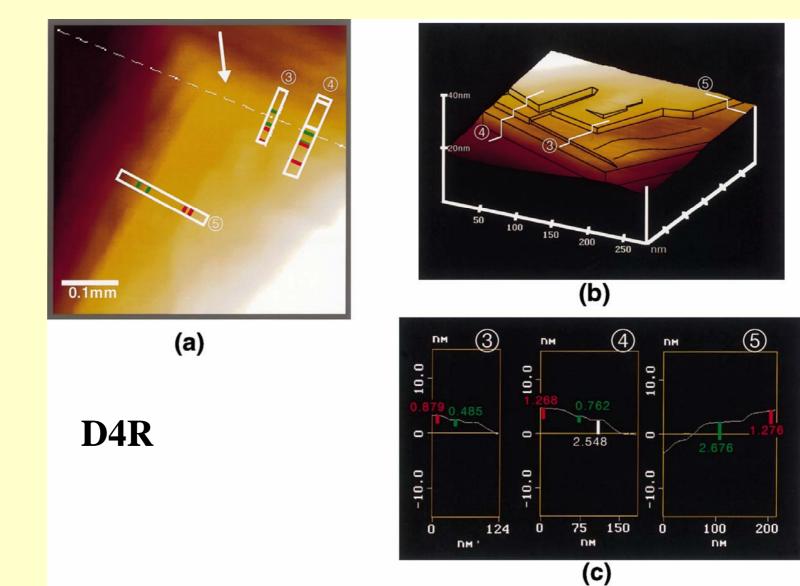
#### **Zeolite A**

- (a) [TO<sub>4</sub>] tetrahedra as BBU
- (b) four-membered single rings
- (c) lB fuenfer chains
- (d) cubes [4<sup>6</sup>]
- (e) truncated octahedra [ $4^66^8$ ] (sodalite- or  $\beta$ -cages)
- (f) truncated cubeoctahedra  $[4^{12}6^88^6]$  ( $\alpha$ -cavities)

#### Pores in Zeolite A (LTA)

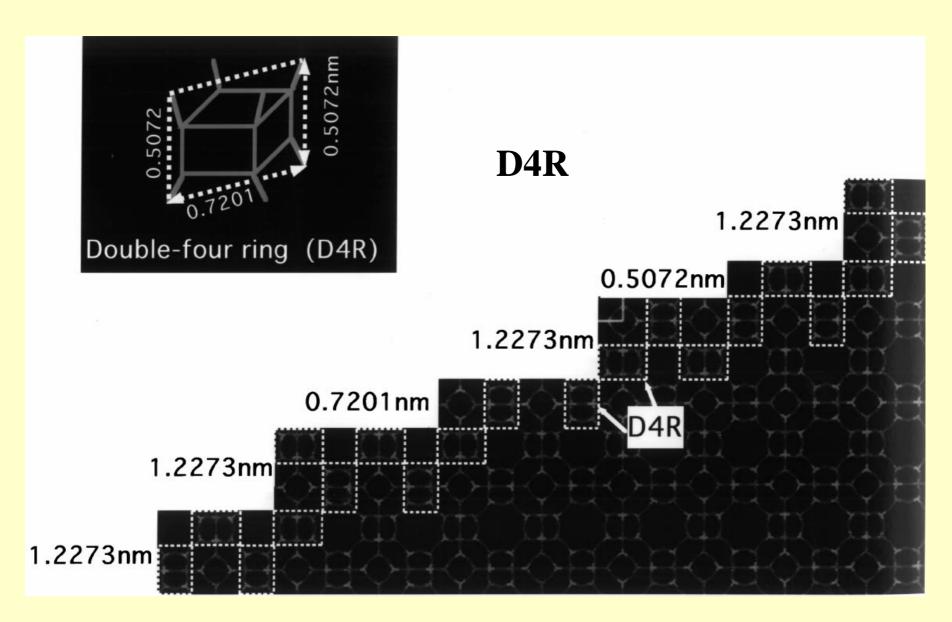


- (a) the sodalite cage  $[4^66^8]$
- (b) the  $\alpha$ -cavity [4<sup>12</sup>6<sup>8</sup>8<sup>6</sup>]
- (c) the 3-dimensional channel system
- (d) the 8-ring defining the 0.41 nm effective channel width



### **AFM growth studies of LTA**

S. Sugiyama et. al. Microporous and Mesoporous Materials 28 (1999) 1–7



#### Zeolite FAU (X and Y) and EMT

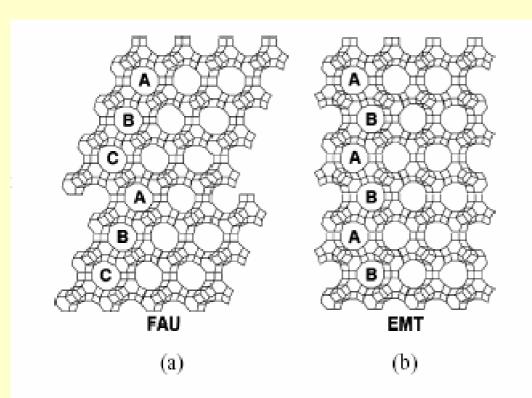


Fig. 1. Structure of zeolite Y: (a) cubic polymorph known as FAU with ABCABC... stacking, (b) hexagonal polymorph known as EMT with ABABAB... stacking.

FAU				
Cubic	ABCABC stacking of layers agent	analagous to zinc blende	15-crown-5 structure directing agent	
	EN	МT		
Hexagonal	ABABAB stacking of layers	analagous to wurtzite	18-crown-6 structure directing agent	

#### **Molecular Sieves**

Zeolite	Cation	Code	Pore diameter
Zeolite A:	Na	4A	0.42 nm
	Ca	5A	0.48 nm
	Na, K	3A	0.38 nm
Zeolite X:	Na	13X	0.8-1.0 nm
	Ca	10X	0.7 nm

Zeolite Y contains more Si

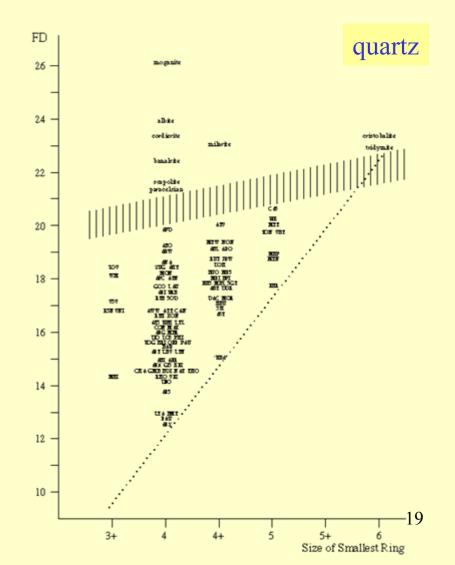
#### **Framework Density**

Framework density (FD)

Defined as the number of tetrahedral atoms (T-atoms) per cubic nanometer  $(1000 \ A^3)$ 

FD is related to the void volume of the crystal: as the FD value decreases, the void volume and capacity for adsorption increases

FD < 20 are characteristic of microporous structures the minimum known FD is 12.5 with the void occupying just over half of the crystal volume



#### **Pores**

Various sizes (4 - 13 Å), shapes (circular, elliptical, cloverleaf-like), and connectivity (1-3D)

The size of the rings formed by the  ${\rm TO_4}$  tetrahedra ranges from 4 to 18 of the T-atoms and determines the pore aperture

Extraframework charge-balancing cations Ion-exchangeable, size, charge, positions, distribution, ordering, coordination number

Si-to-Al ratio

Influences cation content, hydro-phobicity/-philicity, acidity

Löwenstein rule:

absence of the Al-O-Al moieties, in aluminosilicates Si/Al > 1

Linde A (LTA) Si/Al = 1

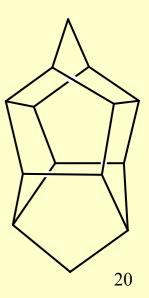
ZK-4 (LTA) Si/Al = 2.5

ZSM-5 Si/Al =  $20 - \infty$ 

Pure  $SiO_2$   $Si/Al = \infty$ 

#### **Pentasils**

ZSM-5



#### **Zeolite Synthesis**

**Synthesis** - an empirical and heuristic process, new phases are often discovered by serendipity

**Aluminosilicates – high pH** 

altering it

extraction (neutral templates)

#### **Zeolite Synthesis**

Structure of the zeolite product depends on:

- Composition
- Concentrations and reactant ratios
- Order of mixing
- Temperature
- Ageing time (hours to weeks)
- Crystallization time (days to weeks, kinetics of the structure-directing process is slow)
- pH
- Stirring/no stirring
- Pressure
- Seeding
- Reactor material (PTFE, glass, steel)
- Templates

Templates: Organic cationic quaternary alkylammonium salts, alkylamines, aminoalcohols, crownethers,

structure-directing, space-filling, charge-balancing

Vary the template - discover new structures!

#### **Templates**

Template or guest compounds

Three levels of the guest action with increasing structure-directing specificity:

- Space-filling the least specific, observed, for example, in the synthesis of AlPO<sub>4</sub>-5, 23 different, structurally unrelated compounds, could be employed, they pack in the channels of the structure thereby increasing its stability.
- Structure-directing a higher degree of specificity, only tetramethylammonium hydroxide is effective in the synthesis of  $AIPO_4$ -20
- -elongated molecules, such as linear diamines, initiate the formation of channels
- -nondirectional-shaped guests leads to the formation of cage-like cavities, the size of these cavities correlates with the size of freely rotating guests
- True templating very rare, it requires even more precise hostguest fit which results in the cessation of the free guest-molecule rotation

A curiosity: aluminophosphate VPI-5 does not require any guest for its formation!

# **Templates**

The ratio  $TO_2/(C + N + O)$  is a measure of space-filling of the framework by the guest molecules, characteristic for a specific guest and structure.

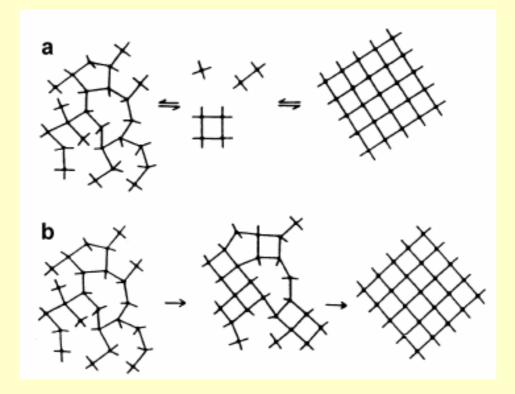
Existence of primary and secondary units in a synthesis mixture

4R, 6R, 8R, D4R, D6R, 5-1, cubooctahedron

#### **Zeolite Synthesis Mechanisms**

(a) gel dissolution and solution mediated crystallization (SBU in

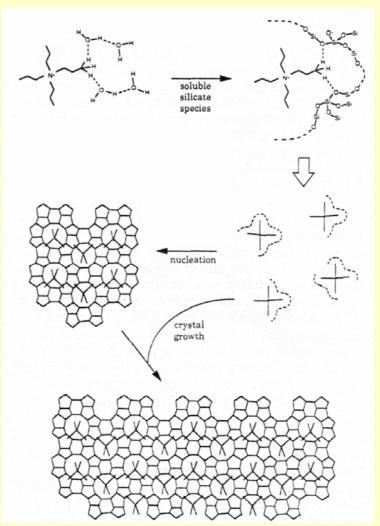
solution)



(b) "in situ" rearrangement of the gel

# **Zeolite Synthesis Mechanisms**

Mechanism of structure-directing action of the TPA template



Wide range of solid state characterization methods for zeolites: diffraction, microscopy, spectroscopy, thermal, adsorption and so forth

**Zeolite post modification for controlling properties of zeolites** 

Tailoring channel, cage, window dimensions:

- **♦** Cation choice (Ca<sup>2+</sup> exchanged for Na<sup>+</sup>)
- **★**Larger Si/Al decreases unit cell :

decreases unit cell parametrs, window size decreases number of cations, free space increases hydrophobicity

**✦**Reaction temperature, higher T, larger pores

# **Stability Rules**

Löwenstein rule: never Al-O-Al

Dempsey rule: Al-O-Si-O-Si-O-Al

is more stable than

Al-O-Si-O-Al

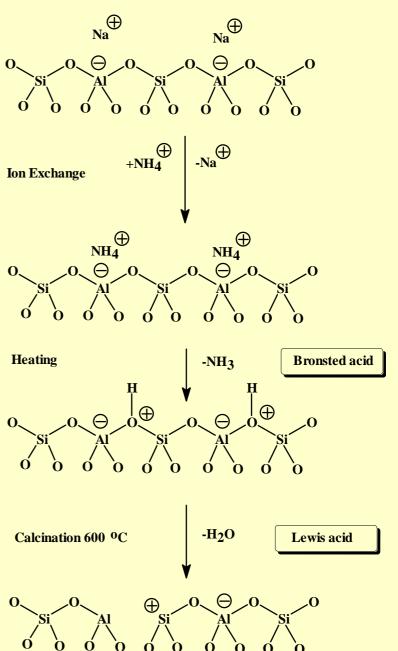
# **Brønsted Acidity**

**Tuning Bronsted acidity:** 

Ion exchange for NH<sub>4</sub><sup>+</sup> Pyrolysis to expel NH<sub>3</sub> Calcination to expel H<sub>2</sub>O

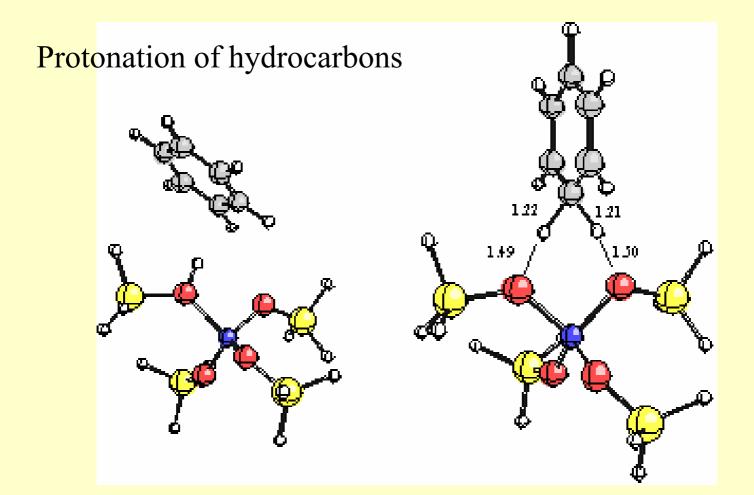
Solid acid for the hydrocarbon cracking

The larger the Si/Al ratio, the more acidic is the zeolite



29

# **Brønsted Acidity**

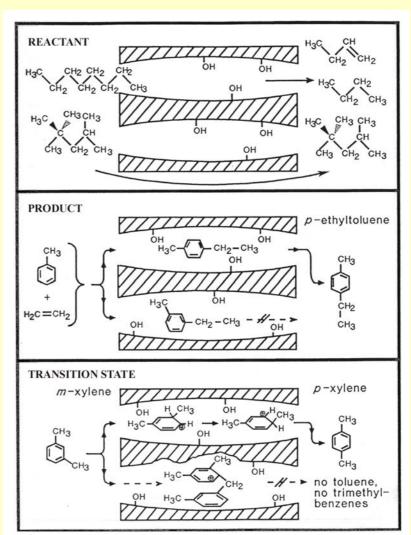


#### **Size-Shape Selectivity**

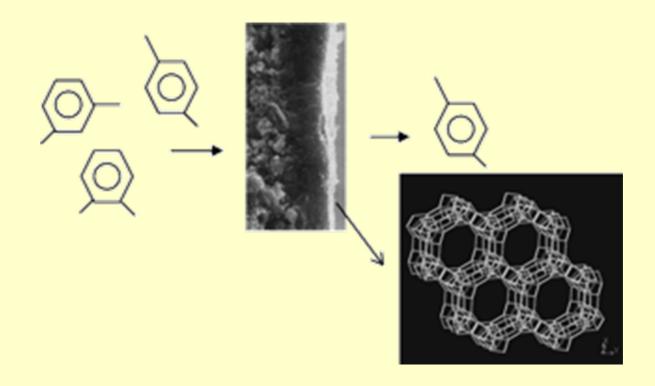
Size-shape selective catalysis, separations, sensing

#### **Selectivity at:**

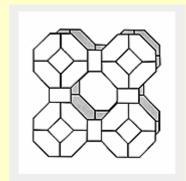
- Reactants
- Products
- •Transition state



# Separation of xylene isomers by pervaporation thru a MFI membrane







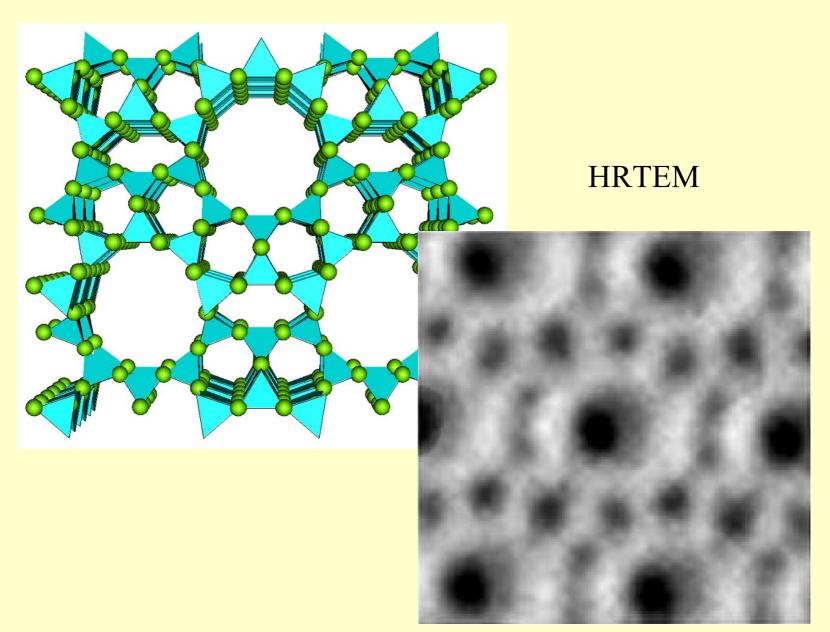
Odor control, adsorbents

Ion exchange capacity, water softening, detergents (25wt% zeolite)

Host-guest inclusion, atoms, ions, molecules, radicals, organometallics, coordination compounds, clusters, polymers (conducting, insulating)

**Nanoreaction chambers** 

Advanced zeolite devices, electronic, optical, magnetic applications, nanoscale materials, size tunable properties, QSEs



Aquaculture

Ammonia filtration in fish hatcheries Biofilter media

Agriculture

Odor control Confined animal environmental control Livestock feed additives

Horticulture Nurseries, Greenhouses

Floriculture

Vegetables/herbs

Foliage

Tree and shrub transplanting

Turf grass soil amendment

Reclamation, revegetation, landscaping

Silviculture (forestry, tree plantations)

Medium for hydroponic growing

Household Products Household odor control Pet odor control

Industrial Products Absorbents for oil and spills Gas separations

Radioactive Waste Site remediation/decontamination

Water Treatment Water filtration Heavy metal removal Swimming pools

Wastewater Treatment Ammonia removal in municipal sludge/wastewater Heavy metal removal Septic leach fields

#### Aluminophosphates

- **+**Isoelectronic relationship of AlPO<sub>4</sub> to (SiO<sub>2</sub>)<sub>2</sub>
- **★**Ionic radius of Si<sup>4+</sup> (0.26 Å) is very close to the average of the ionic radii of Al<sup>3+</sup> (0.39 Å) and P<sup>5+</sup> (0.17 Å)

Many similarities between aluminosilicate and AlPO<sub>4</sub> molecular sieves

Dense AlPO<sub>4</sub> phases are isomorphic with the structural forms of SiO<sub>2</sub>: quartz, tridymite, and cristobalite Aluminosilicate framework charge balanced by extraframework cations

Aluminophosphate frameworks neutral  $(AlO_2^-)(PO_2^+) = AlPO_4$ 

#### Aluminophosphates

Some AlPO<sub>4</sub> structures are analogous to zeolites while other are novel and unique to this class of molecular sieves.

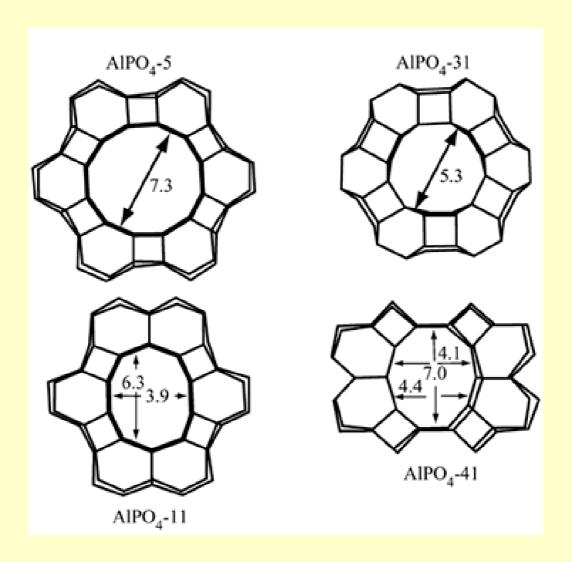
Only even-number rings = the strict alternation of Al and P atoms

Incorporation of elements such as Si, Mg, Fe, Ti, Co, Zn, Mn, Ga, Ge, Be, Li, As, and B into the tetrahedral sites of AlPO $_4$  gives a vast number of element-substituted molecular sieves (MeAPO, MeAPSO, SAPO) important heterogeneous catalysts  $M^{1+}$ ,  $M^{2+}$ , and  $M^{3+}$  incorporate into the Al sites

M<sup>5+</sup> elements incorporate into the P sites

This substitution introduces a negative charge on these frameworks. Si<sup>4+</sup>, Ti<sup>4+</sup>, and Ge<sup>4+</sup> can either replace P and introduce a negative charge or a pair of these atoms can replace an Al/P pair and retain the charge neutrality.

# Aluminophosphates



# Gallophosphates

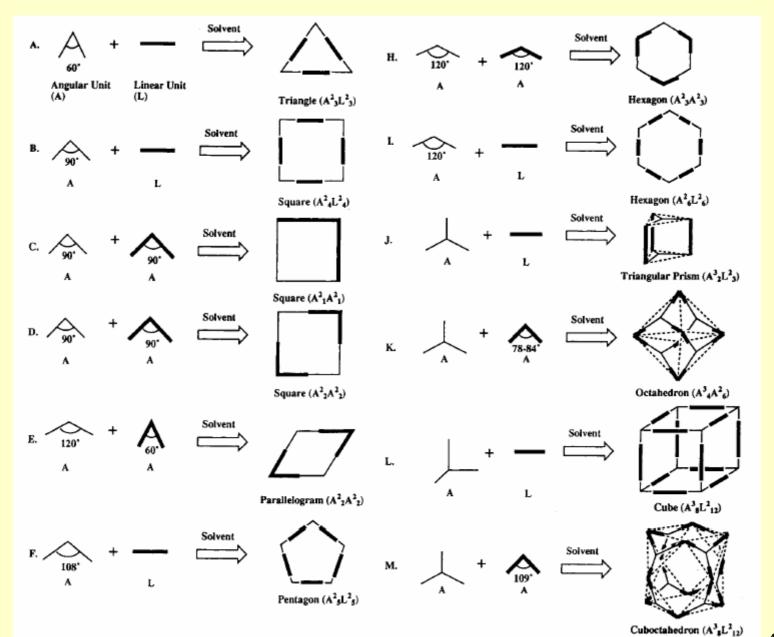
#### **Aluminophosphate Synthesis**

Aluminophosphates prepared by the hydrothermal synthesis Source of Al: pseudoboehmite, Al(O)(OH),  $Al(Oi-Pr)_3$ 

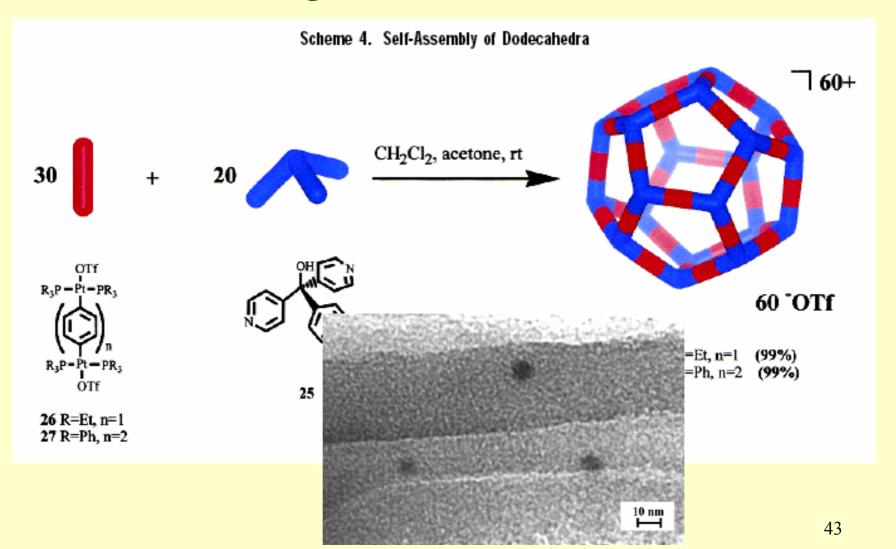
Mixing with aqueous  $H_3PO_4$  in the equimolar ratio – low pH! Forms an  $AlPO_4$  gel, left to age One equivalent of a guest compound = template Crystallization in a reactor Separated by filtration, washed with water Calcination

Other zeolite materials
Oxide and non-oxide frameworks, sulfides, selenides
Coordination frameworks, supramolecular zeolites
The quest for larger and larger pore sizes

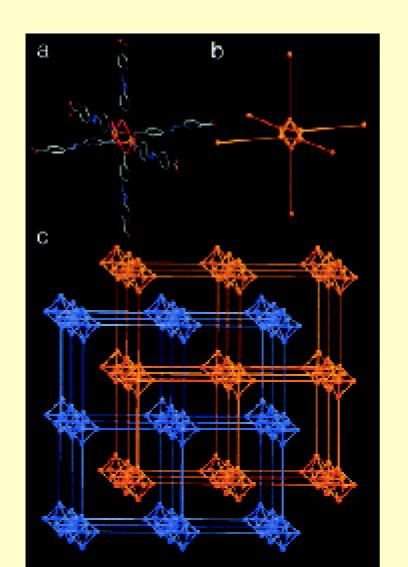
# **Metallo-Organic Framework Structures**

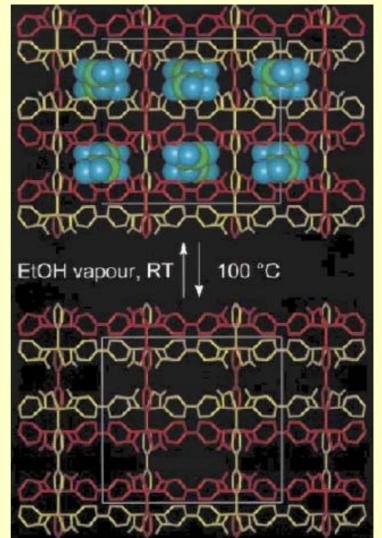


#### **Metallo-Organic Framework Structures**



# **Metallo-Organic Framework Structures**





# **Inorganic and Metallo-Organic Quartz**

