



Ion Discrimination by Nanoscale Design



water tower: www.pbase.com/mescaleroman/images/57217408

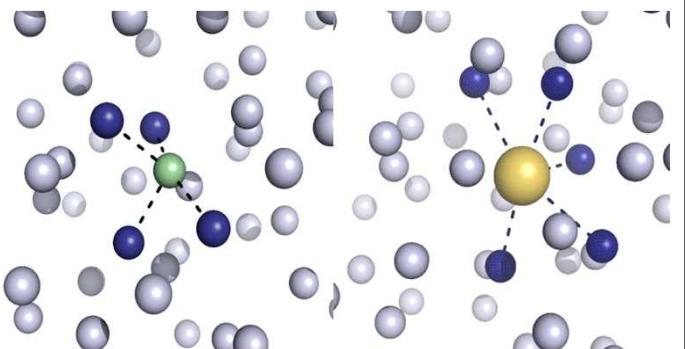
Susan Rempe

Sandia National Labs, Albuquerque, NM



Sandia National Laboratories

Natural protein channels: Not just simple holes!



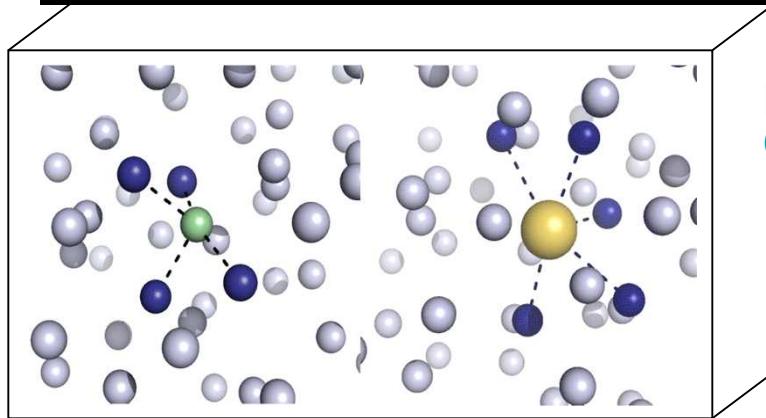
Ions in water
(Varma & Rempe, 2007)

- Ions are 'happy' **coordinated** with water **ligands** in liquid water

| Bare ion | Ion-water bond |
|----------|----------------------------|
| .95 Å | Na^+ 2.4 Å |
| 1.4 Å | H_2O 2.8 Å |
| 1.3 Å | K^+ 2.8 Å |

- K^+/Na^+ exquisite discrimination:
 - same charge
 - same size at kT
 - larger ion transported, fast!

Natural protein channels: Not just simple holes!

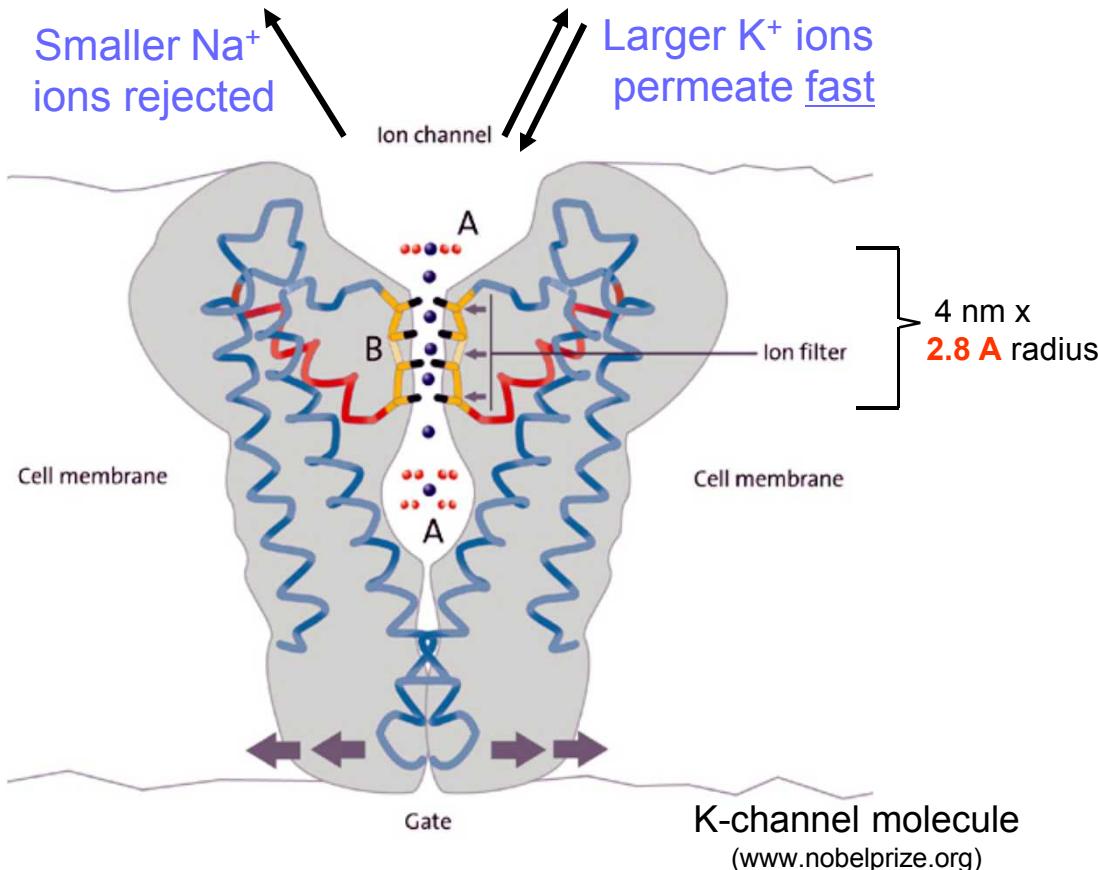


Ions in water
(Varma & Rempe, 2007)

- Ions are 'happy' **coordinated** with water **ligands** in liquid water

Unresolved Questions:

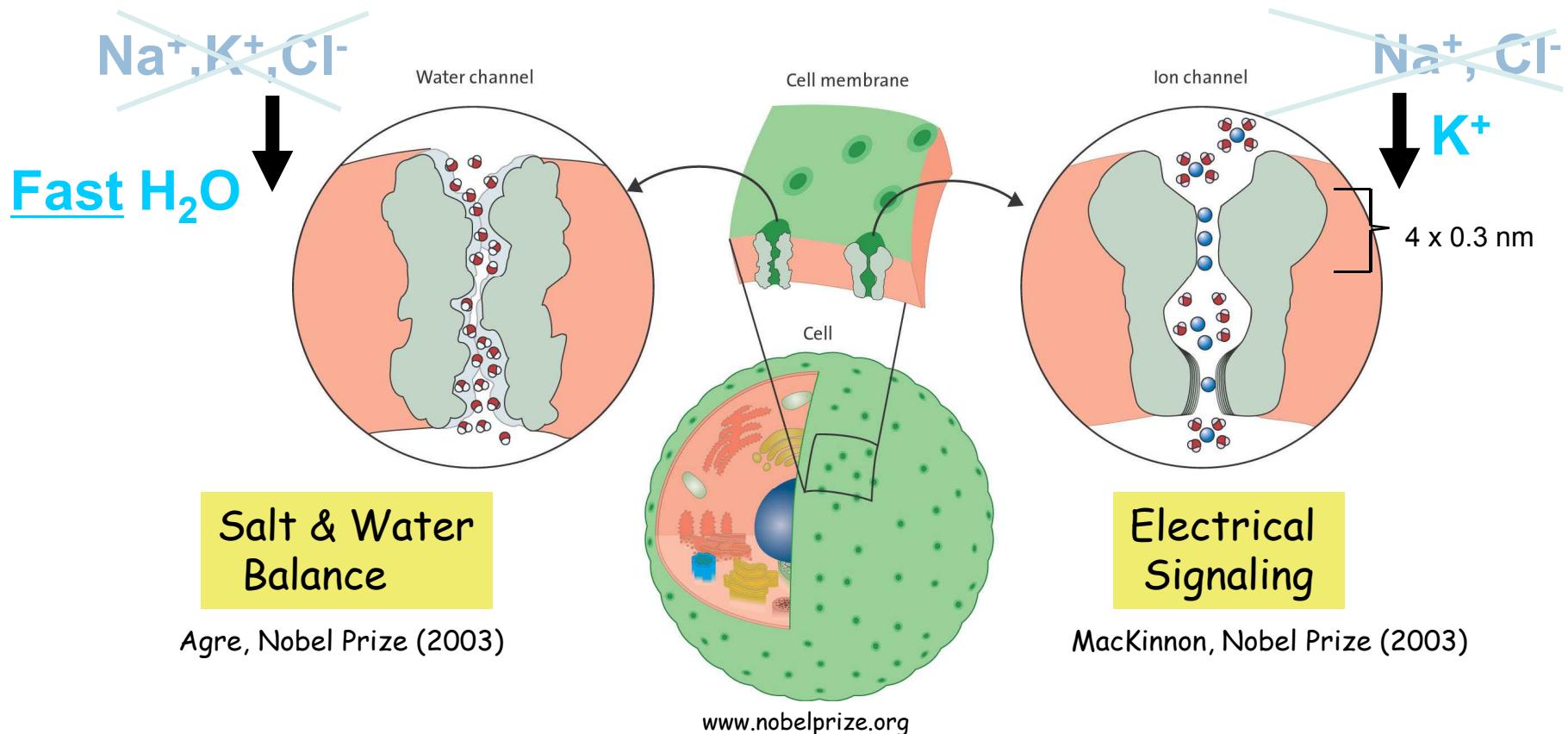
- How do channels work?



| Bare ion | Ion-water bond |
|----------------------|----------------|
| Na^+ | 2.4 Å |
| H_2O | 2.8 Å |
| K^+ | 2.8 Å |

- K^+/Na^+ exquisite discrimination:
 - same charge
 - same size at kT
 - larger ion transported, fast!

Natural channel proteins: Small dimensions, large biological impact



Channels are fundamental to life:
severe Health consequences if disrupted.

Nobel Prizes underscore Health significance of Channels....



| Time and Nobel Prize Field | Laureates | Specific Research |
|-------------------------------|--|--|
| 1963 Nobel Prize in Medicine | <u>Alan L. Hodgkin, Andrew F. Huxley</u> | <u>Ion mechanisms of nerve impulse</u> |
| 1985 Nobel Prize in Medicine | Michael S. Brown, Joseph L. Goldstein | Regulation of cholesterol metabolism |
| 1988 Nobel Prize in Chemistry | Johann Deisenhofer, Robert Huber, Hartmut Michel | Structure of photosynthetic reaction center |
| 1991 Nobel Prize in Medicine | <u>Erwin Neher, Bert Sakmann</u> | <u>Function of single ion channels</u> |
| 1997 Nobel Prize in Chemistry | Jens C. Skou | Membrane-bound turnover of ATP |
| 1999 Nobel Prize in Medicine | Günter Blobel | Principles of protein compartmentalization |
| 2003 Nobel Prize in Chemistry | Roderick MacKinnon, Peter Agre | Molecular structural analysis of membrane channels, existence of water channels |
| 2004 Nobel Prize in Medicine | Richard Axel, Linda B. Buck | Odorant receptors |

Global Problem: Desalination

Clean water: a global precious commodity

- water is recyclable
- but RO is expensive, produces sterile water

Salinity Levels:

Seawater: ~35 g/l (0.6 M)

Brackish: ~1-5 g/l (0.08 M)

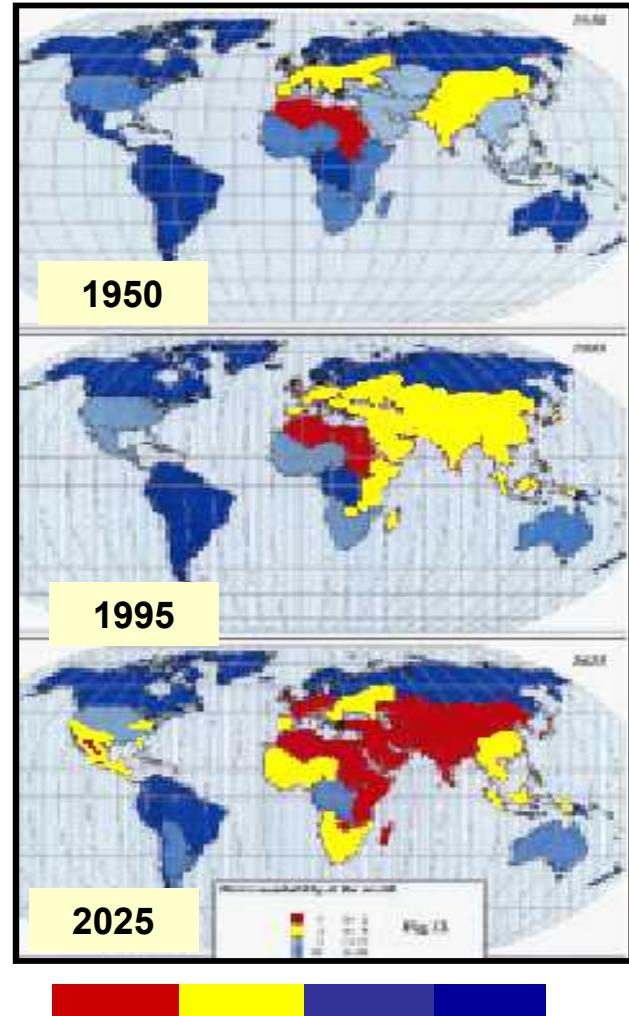
Potable: <0.5 g/l (0.008 M)

Efficient membranes: critical challenge

- fast (barrierless) water transport
- select ion exclusion (mineral water)

“Water promises to be to the 21st century what oil was to the 20th century: the precious commodity that determines the wealth of nations.”

Fortune Magazine, May 15, 2000

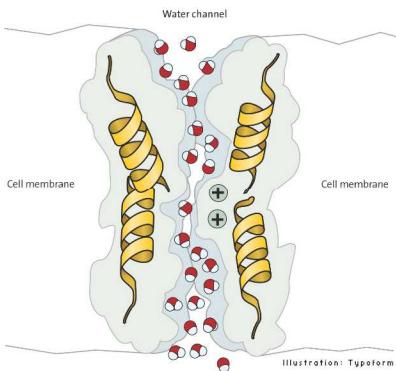


shortage

Efficient Membranes:

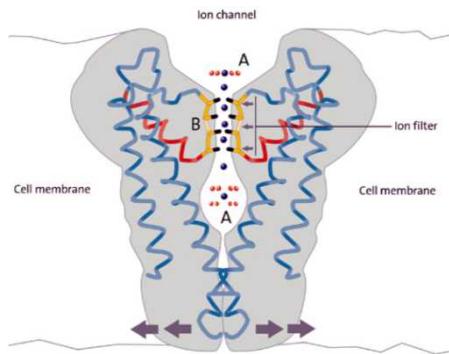
Understand, design, engineer nano-channels for desalination

Water channels



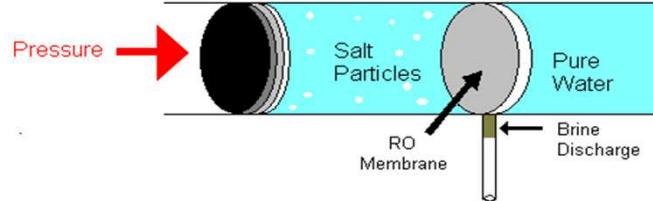
- transport H_2O fast

Ion channels

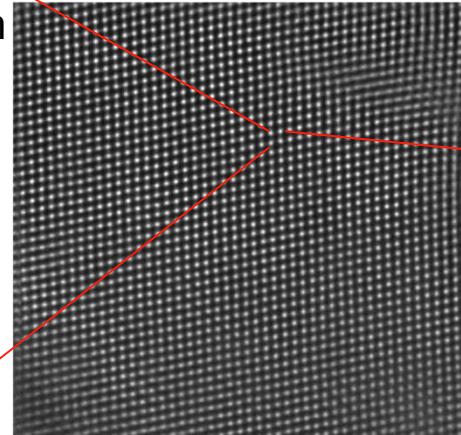


- transport select minerals

- Bio-inspired design



- Theory & modeling



Brinker Lab

- Inorganic membrane synthesis

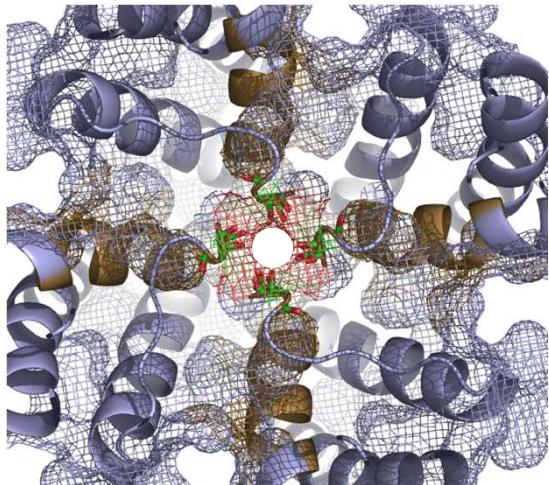


Solution: Harness molecular biomechanisms.
Gain 10x in water flux + minerals.

What parameters do we give our engineers?

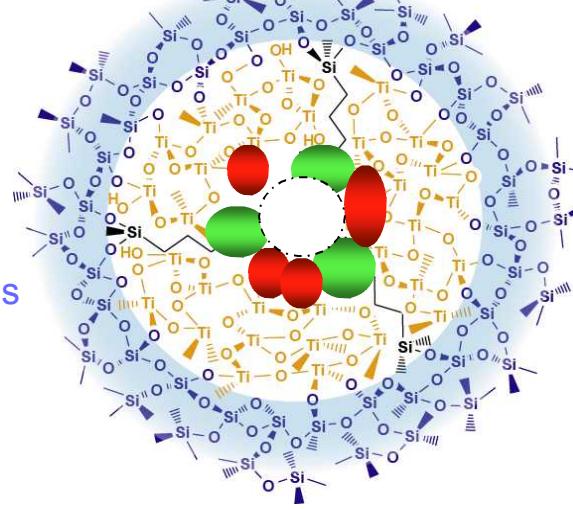
4 x 0.3 nm
Active site

Natural Channels



Bio mechanisms
Engineering solutions

Inorganic Channels



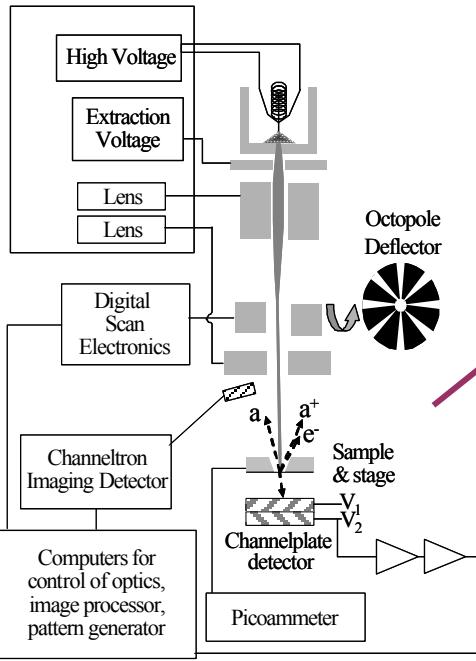
Critical Channel Design Issues:

What's significant about (1) Mouth?
(2) Narrow filter?
(3) Chemistry, architecture?

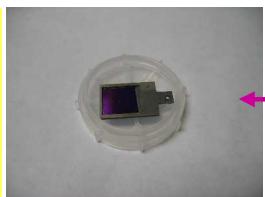
Subtle, challenging questions demand molecular precision:

→ **Molecular Modeling** + Molecular Synthesis

Platforms for Experiments & Modeling

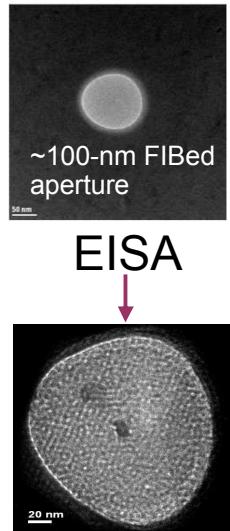


FIB experimental set-up



Special holder device

- TEM observation
- Patch Clamp
- PA-ALD
- Iterate



PA-ALD

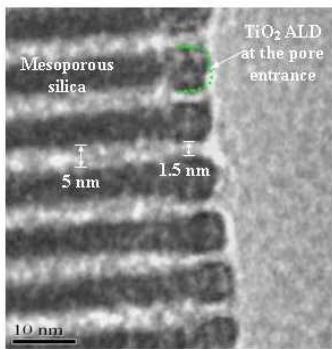
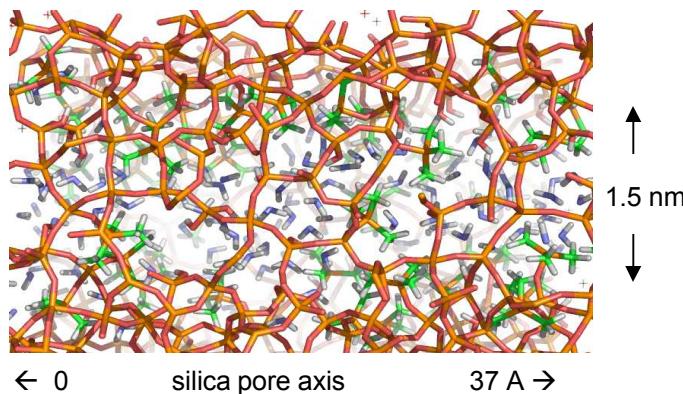


Fig. 1a

- Experimental platform allows successive modification (pore size, and surface chemistry), imaging and transport measurements on identical sample.

- Multiscale modeling essential to understand combined effects of pore size, structure, chemistry and charge.



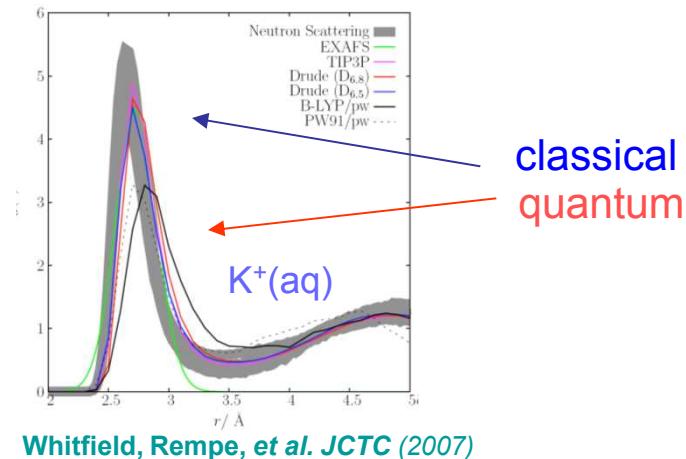
- Classical (large, long times)
- Quantum (accuracy)
 - Thermodynamics (work)
 - Dynamics (transport)

Modeling Approach: Molecular + quantum accuracy

- Quantum (ab initio) interactions:
 - expensive; describe complex interactions (ex. ion-ligand chemical bonds)
- Classical molecular interactions:
 - inexpensive; simplified, parameterized

- Example 1

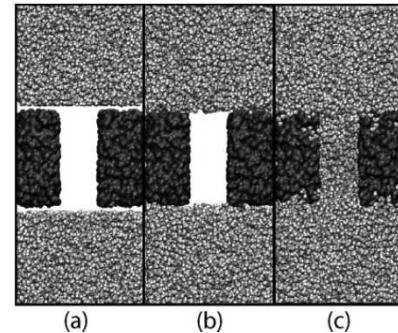
More ligands bind classical ions??



Whitfield, Rempe, et al. *JCTC* (2007)

- Example 3

<10% change in classical parameters:
Water fills/empties?

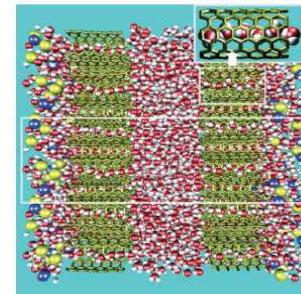


Leung, Rempe, Lorenz *PRL* (2006)
channel construction

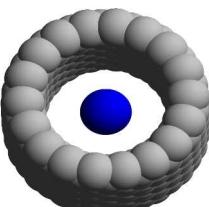
Cruz-Chu et al *JPCB* (2006)

- Example 2

Classical nanotubes
Exclude ions??



Hummer & co *PNAS* (2003)

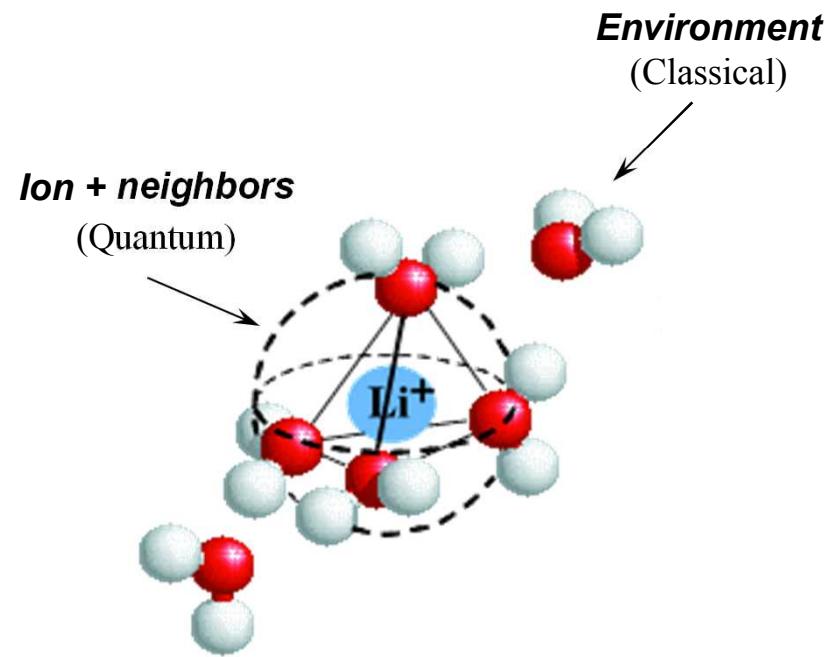
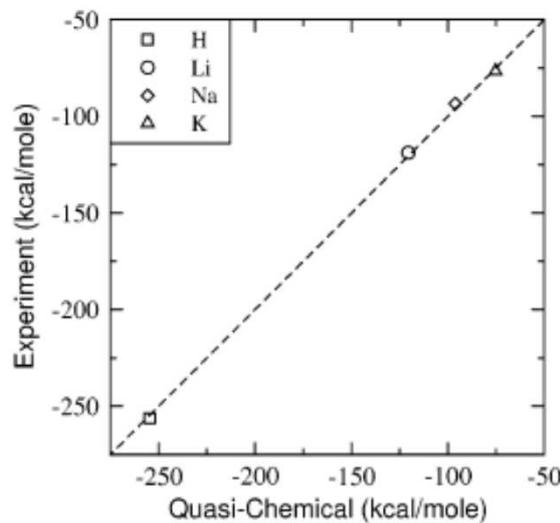


Na⁺ : -210 kJ/mol
Leung, Marsman *JCP* (2007)

Predictions:

Ion transfer thermodynamics using liquid state theory

- Calculate work to transfer ions (water \rightarrow channels), efficiently & accurately
 - 'quasi-chemical' theory¹
- Theory well-tested for ion hydration:²



- Apply to Bio/Inorganic channels for Nanoscale Design Parameters^{3,4,5}

• ¹Pratt & Rempe (1999); Sabo, Rempe, *et al* *JPCB* (2008)

• ²JACS (2004), *PCCP* (2004), *FPE*(2001), *JACS* (2000)

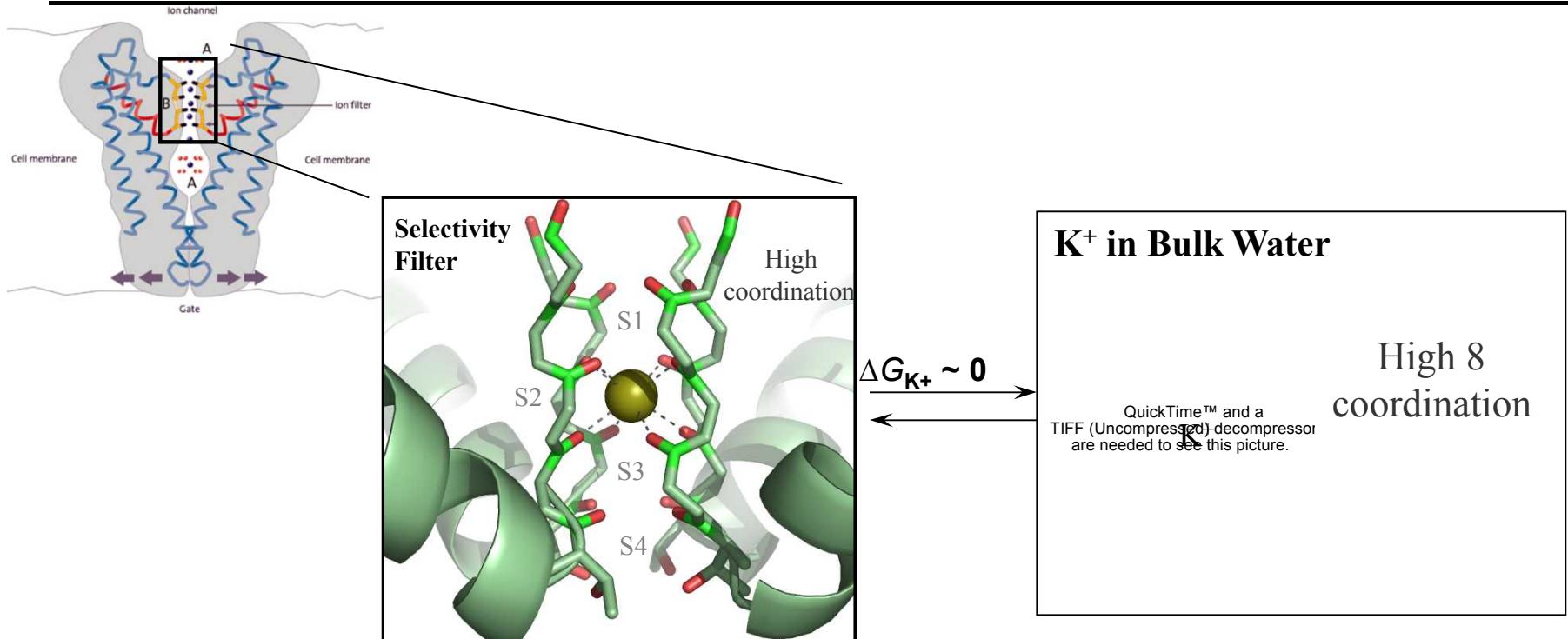
• ³Varma & Rempe *Biophys J* (2007)

• ⁴Varma, Sabo, Rempe *J Molec Bio* (2008)

• ⁵Leung, Rempe, Lorenz *PRL* (2006)

K⁺/Na⁺ Ion Discrimination Problem:

How do K-selective channels work? Prevailing Views

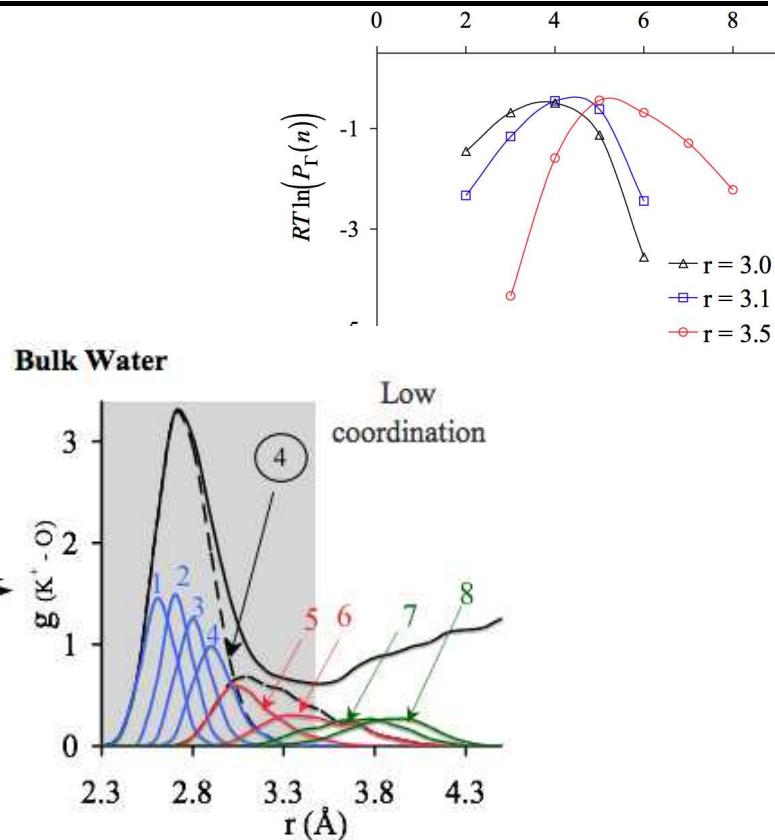
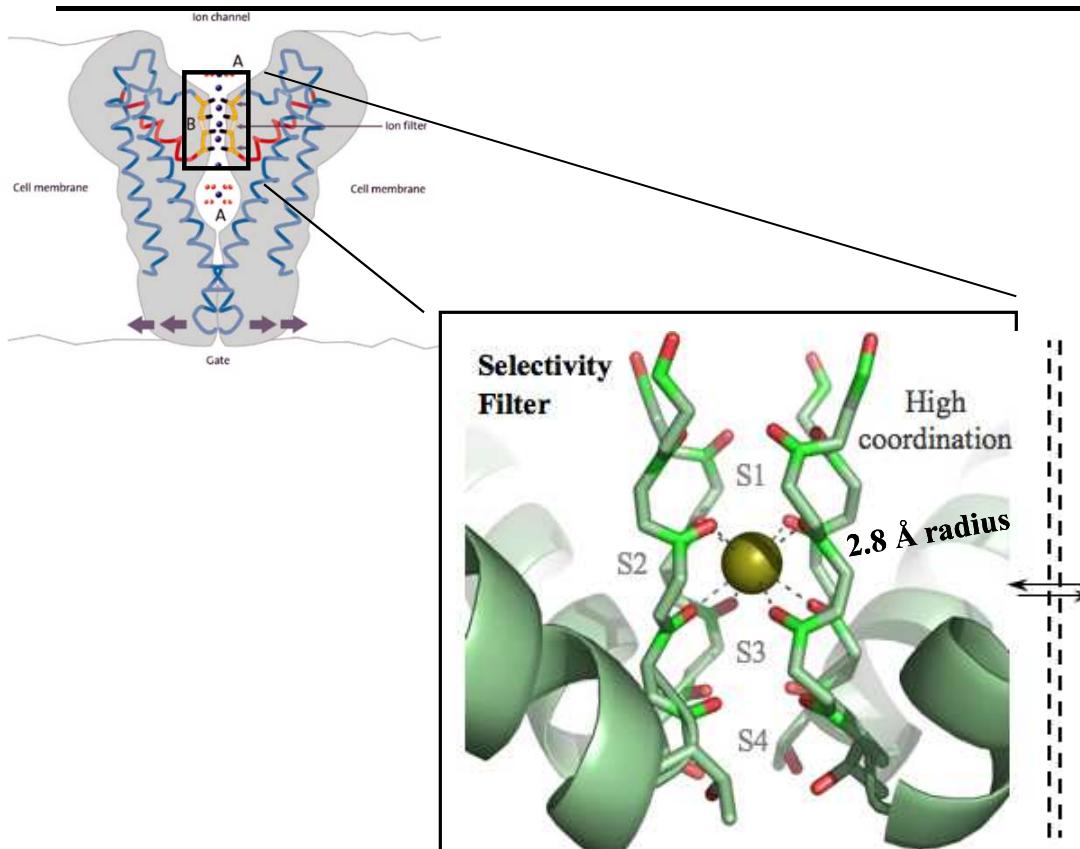


1. Mimic K⁺ ion hydration structures for fast K⁺ transfer^{1,2,3}
2. Specific cavity size fits K⁺^{1,2} vs Liquid-like flexibility³

¹Bezanilla & Armstrong (1972), ²Zhou *et al.* (2001), ³Noskov *et al* (2004)

K^+/Na^+ Ion Discrimination Problem:

How do K-selective channels work? New View #1



1. **Not Mimic** of K⁺ ion hydration structures for fast K⁺ transfer^{1,2}
==> Work avoided by special channel environment (no H-bonders)³

¹Rempe *et al* PCCP (2004), ²Varma & Rempe *Biophys Chem* (2006), ³Varma & Rempe *J Molec Bio* (2008)

Special Channel Environment:

No local competitors for O-ligands stabilizes high ion coordinations

- Coordination preferences controlled by ligand desolvation penalty

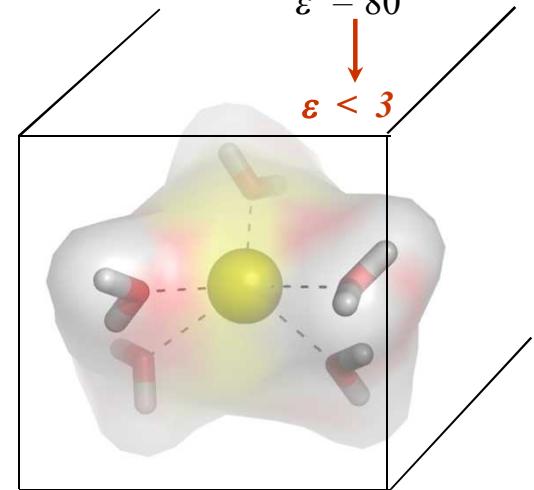


In this “quasi-liquid” environment, preferred ion coordination is higher

$$\text{K}^+ = 8 \quad \text{Na}^+ = 6$$

Surroundings

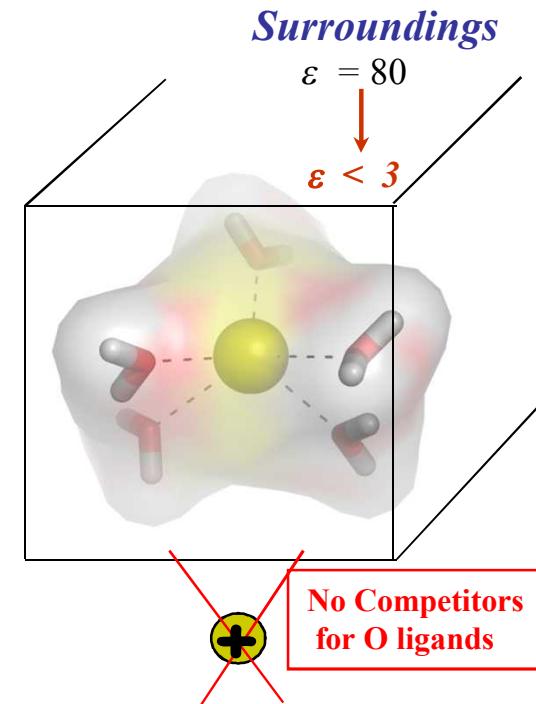
$$\begin{array}{l} \varepsilon = 80 \\ \downarrow \\ \varepsilon < 3 \end{array}$$



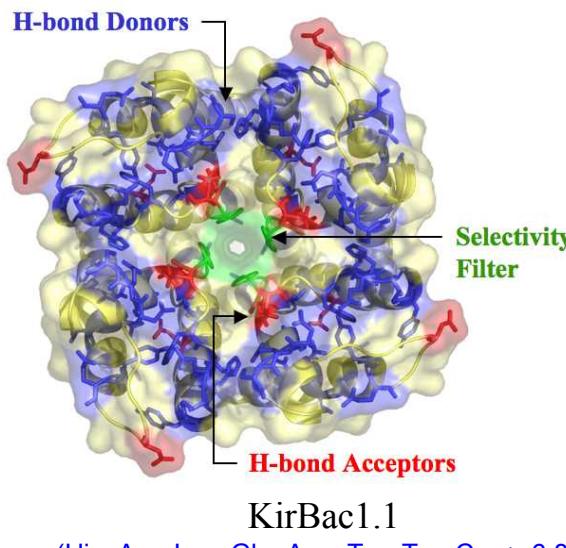
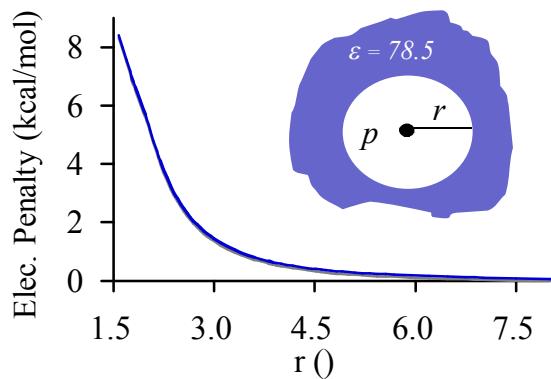
Special Channel Environment:

No local competitors for O-ligands stabilizes high ion coordinations

- Coordination preferences controlled by ligand desolvation penalty



- Ligand desolvation penalty eliminated locally (6 Å)



Highly selective K-channels lack H-bond donors locally

- explains how 8 ligands bind K^+ with no work

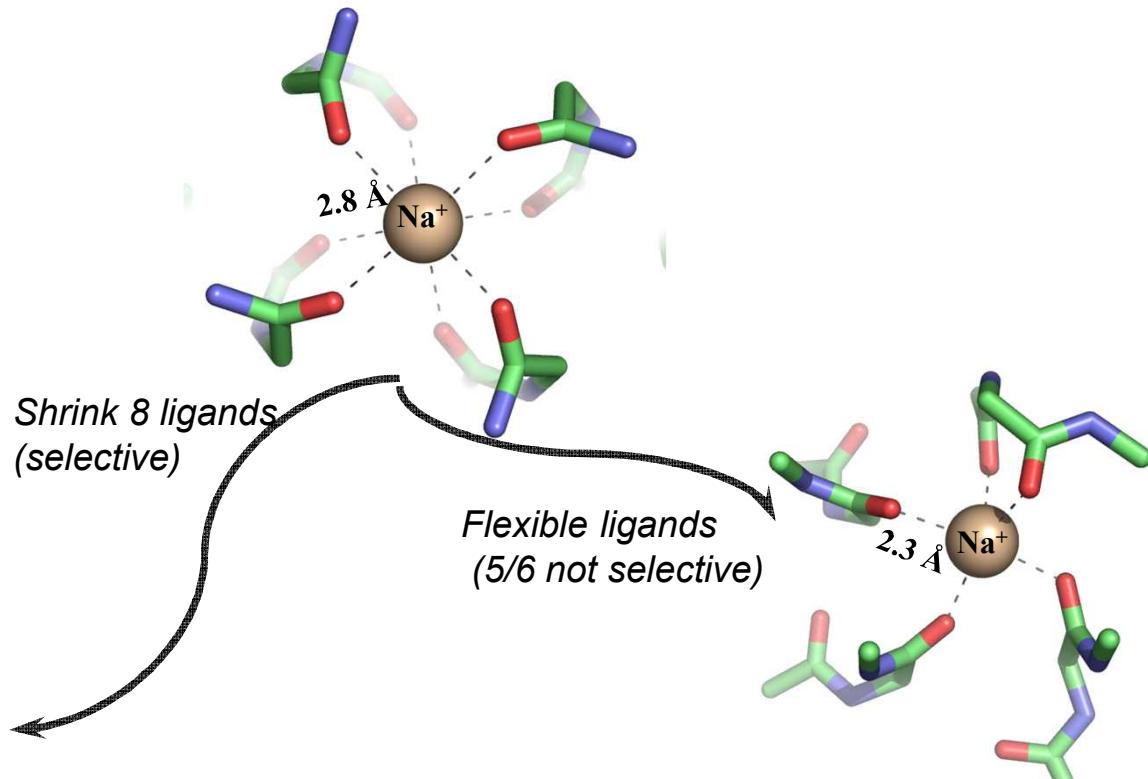
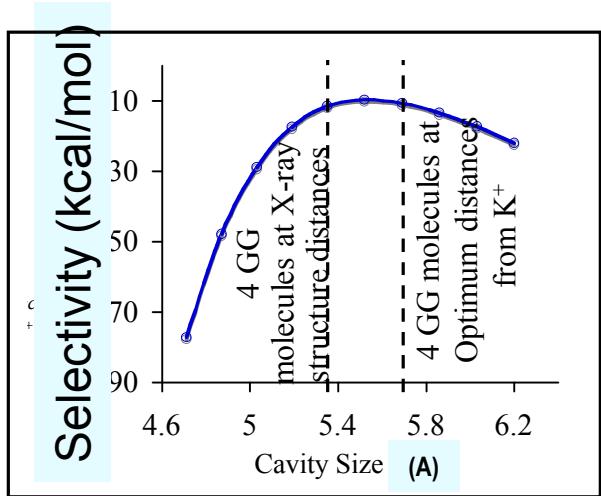
Special Channel Environment:

Structural transitions in ion coordination driven by
changes in competition for ligand binding

QuickTime™ and a
TIFF (Uncompressed) decompressor
are needed to see this picture.

K^+/Na^+ Ion Discrimination Problem:

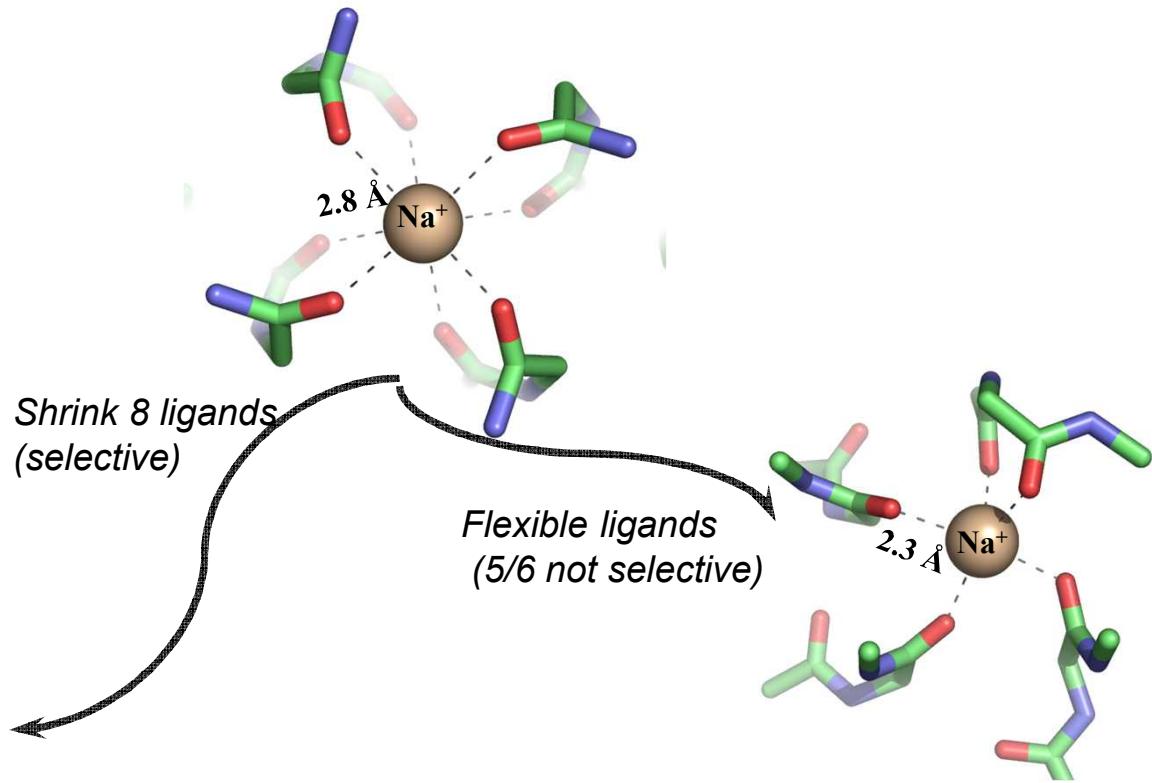
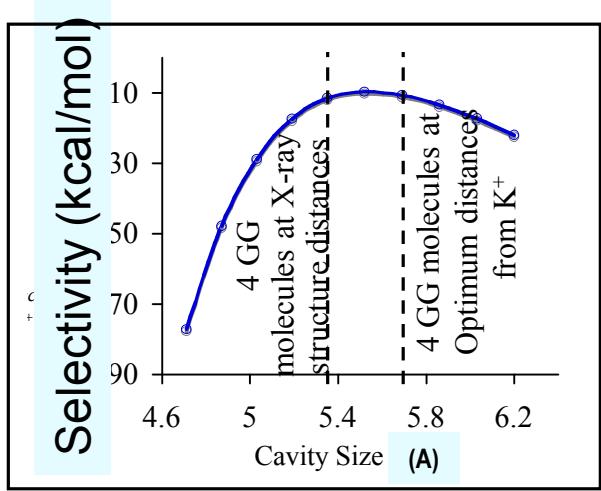
How do K-selective channels work? New View #2



1. Cavity size important
2. Flexibility important

K^+/Na^+ Ion Discrimination Problem:

How do K-selective channels work? New View #2



1. Cavity size important, **but specific size not necessary**
2. Flexibility important, **but can reduce ligands & eliminate selectivity**

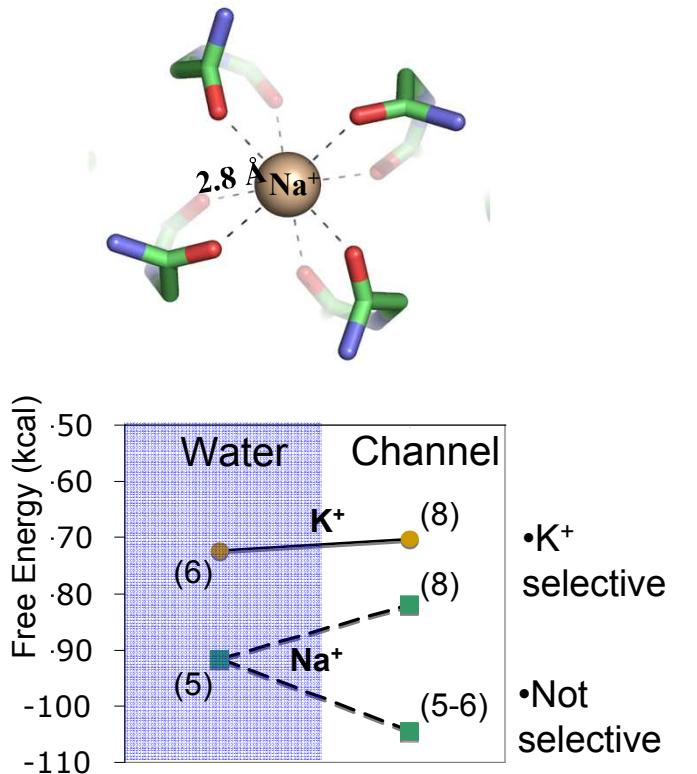
==> **Rigidity¹⁻⁵ important to maintain Over-Coordinated¹⁻⁴ ions (>6 ligands)**

¹Varma & Rempe *Biophys J* (2007); ²Varma, Sabo, Rempe *J Molec Bio* (2008);

³Bostick & Brooks *PNAS* (2007); ⁴Thomas *et al. Biophys J* (2007); ⁵Asthagiri *et al. JCP* (2006)

Fast K⁺/Na⁺ discrimination: Mechanism & Translation Strategy #1

Natural Channels¹



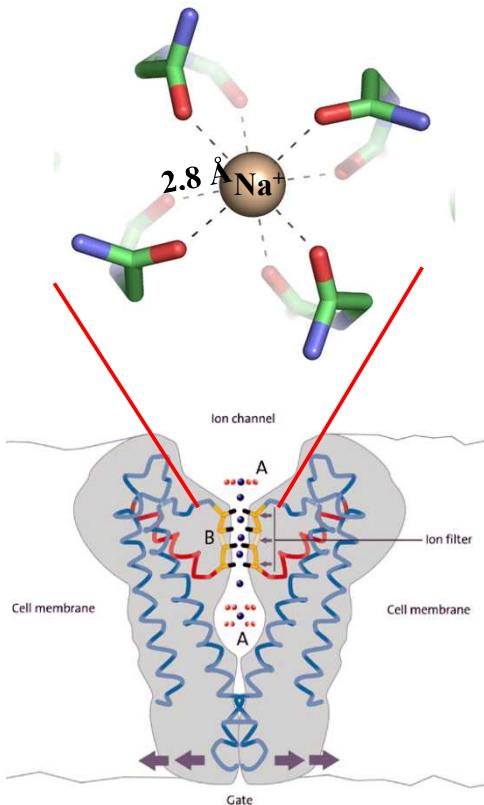
Nanoscale design parameters (natural channel):

“The caress of the surroundings,² the crowding of the ligands”

¹Varma & Rempe *Biophys J* (2007); ²Jordan *Biophys J* (2007); ³Jiang, Brinker, et al. *JACS* (2006)

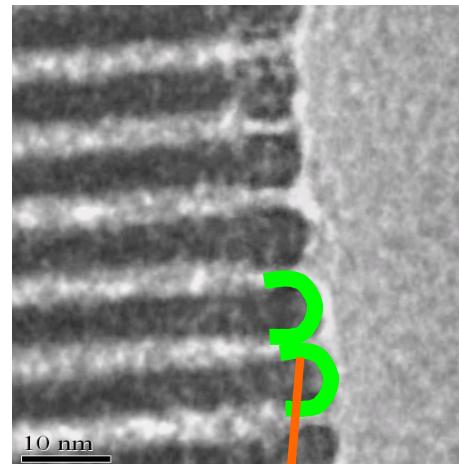
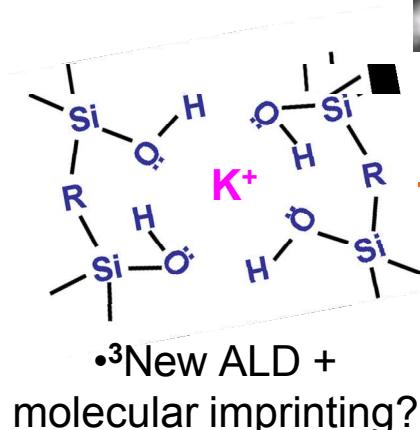
Fast K⁺/Na⁺ discrimination: Mechanism & Translation Strategy #1

Natural Channels¹

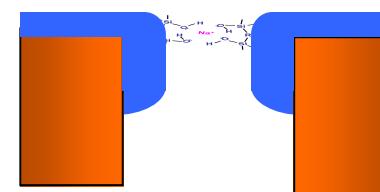


Inorganic Channels

Bio mechanisms
→
Engineered Mimicry



4 nm x 1.5 nm



Nanoscale design parameters (natural channel):

“The caress of the surroundings,² the crowding of the ligands”

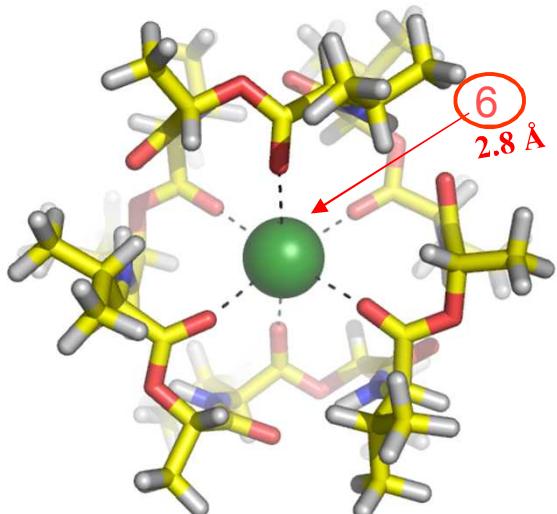
¹Varma & Rempe *Biophys J* (2007); ²Jordan *Biophys J* (2007); ³Jiang, Brinker, et al. *JACS* (2006,2007)

Fast K⁺/Na⁺ discrimination:

Mechanism #2

Natural K-selective Molecule

- 6 C=O ligands, not >6?

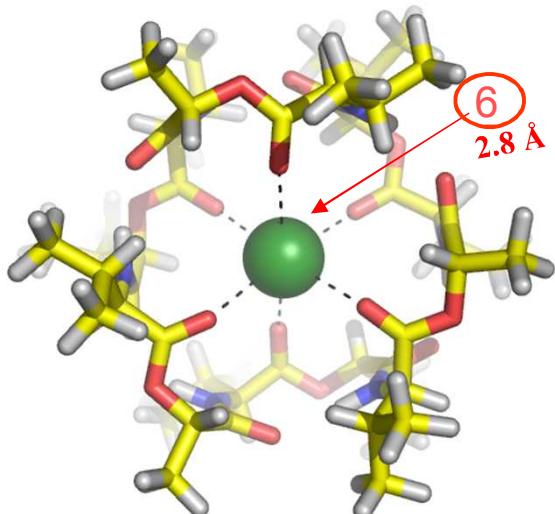


¹Varma, Sabo, Rempe *J Molec Biol* (2008)

Fast K⁺/Na⁺ discrimination: Mechanism #2

Natural K-selective Molecule

- 6 C=O ligands, not >6?



QuickTime™ and a
TIFF (Uncompressed) decompressor
are needed to see this picture.

QuickTime™ and a
TIFF (Uncompressed) decompressor
are needed to see this picture.

$\Delta\Delta g_n(\epsilon)$: change in complexation energy for 2 values of ϵ , which represent dielectric constant lipid membranes. 6-fold coordination is also more stable for higher values of ϵ , such as $\epsilon=80$. RMSD_n reflects change in backbone structure of K⁺-bound valinomycin due to complexation by n water molecules.

Fast K⁺/Na⁺ discrimination: Mechanism #2

Natural K-selective Molecule

- Special C=O chemistry?

QuickTime™ and a
TIFF (Uncompressed) decompressor
are needed to see this picture.

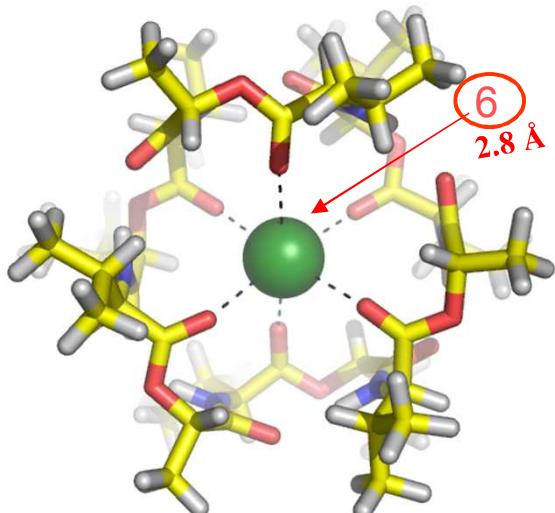
QuickTime™ and a
TIFF (Uncompressed) decompressor
are needed to see this picture.

Fast K⁺/Na⁺ discrimination:

Mechanism #2

Natural K-selective Molecule

- H-bonds maintain cavity size



QuickTime™ and a
TIFF (Uncompressed) decompressor
are needed to see this picture.

¹Varma, Sabo, Rempe *J Molec Biol* (2008)

5.1 Å

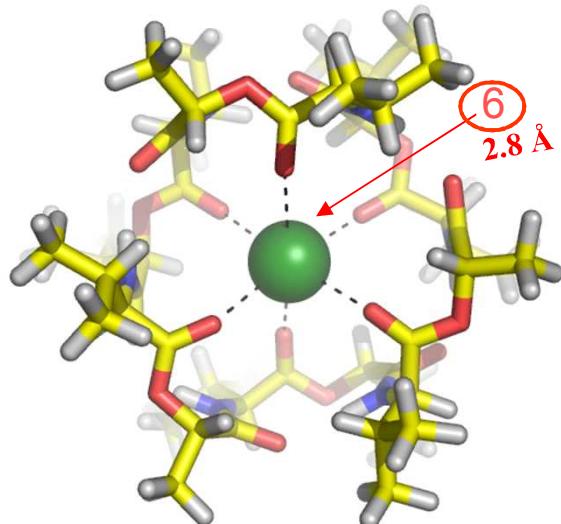
2.3 Å

2.9 Å

H-bonds turned off by replacing proton acceptor atoms =O with =CH₂ groups. QC optimization results in small K⁺ complex changes, but large changes in Na⁺ complex. Absence of H-bonds also increases free energy difference between Na⁺ & K⁺ complexes from 12.3 to 17.8 kcal/mol, thus reducing K⁺/Na⁺ selectivity relative to liquid water.

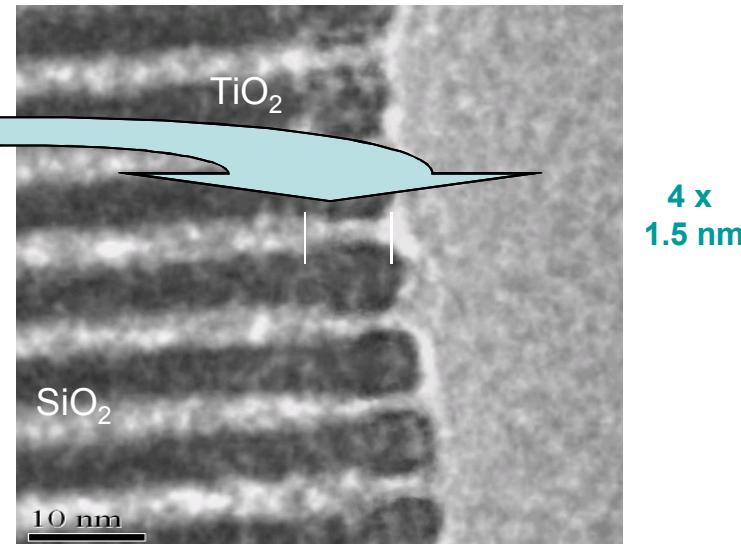
Fast K⁺/Na⁺ discrimination: Mechanism & Translation Strategy #2

Natural K-selective Molecule



- Specific cavity fits K⁺, not Na⁺
- Less selective in water

Inorganic Channels



- Ion-selective aperture?

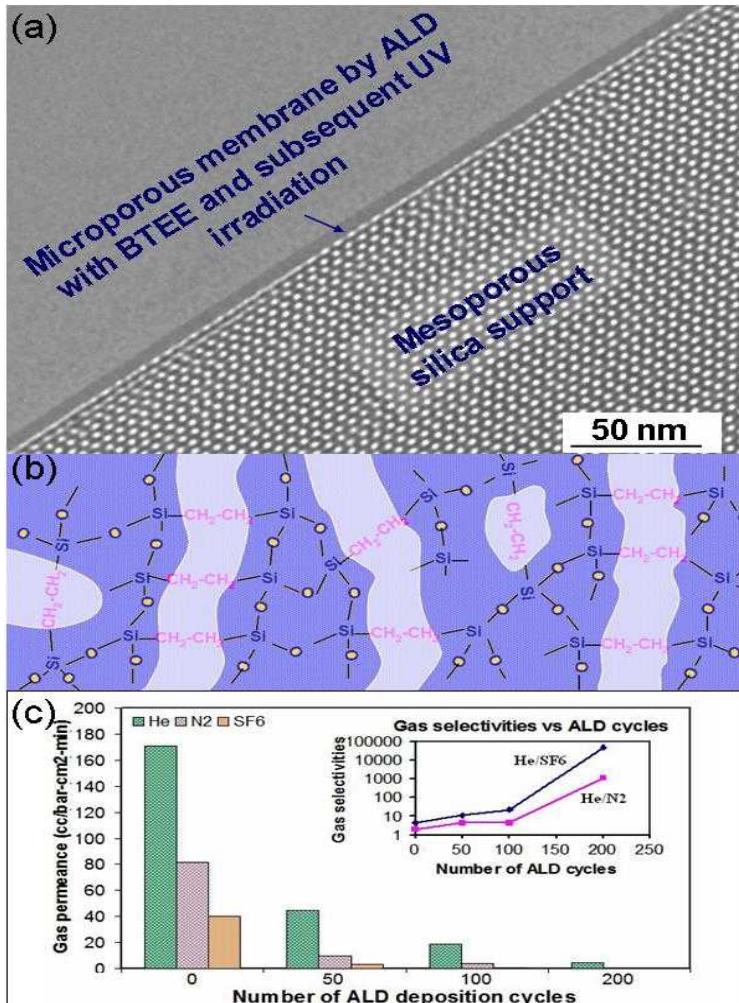
- Outstanding questions:
 - Queue of ion binding sites?
 - Avoid block?
 - Fast water transport + ion selectivity?



¹Varma, Sabo, Rempe *J Molec Biol* (2008)

Fast size discrimination:

Translated (Fabricated) by Atomic Layer Deposition



- **1 biomimetic pores (0.3 nm diameter)** via new ALD & molecular templating
 - high flux + high size-selectivity (He/N₂)
- template-based approach for uniform molecular-sized pores established
- mimic biological K⁺/Na⁺ selectivity?

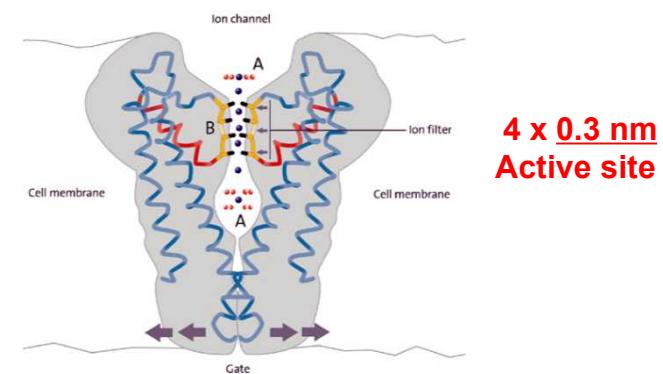
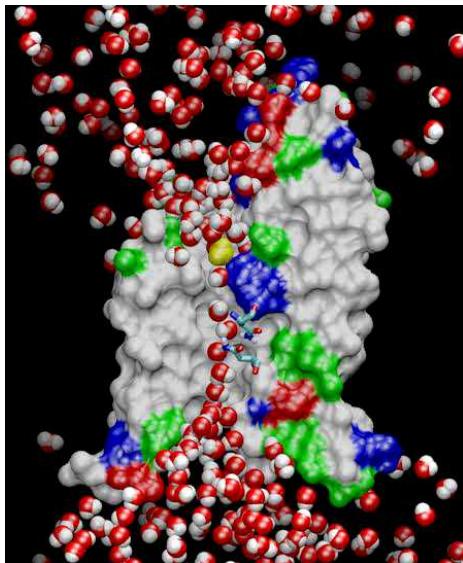


Fig. 1 a) TEM: micropore membrane self-assembled on mesopore support;
b) UV/ozone removes C₂ bridging ligands (pore templates), from dense hybrid film;
c) gas permeance vs ALD cycle; after 200 cycles, N₂ excluded/He transports;
thus pore size ~0.3 nm.

¹Jiang, Brinker, et al JACS, 2007

Summary: Nanoscale Channel Structures for Big Problems

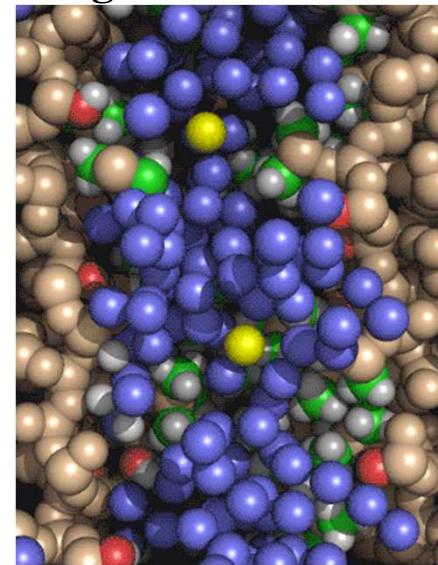
Natural Water Channel



(Tajkhorshid & co)

Bio mechanisms
→
Engineering solutions

Inorganic Water Channel



(Desal Team, Sandia)

- **Channels:** molecular structures & subtle, important functions
- **Solve problems:**
 - **Health** (medicine + biodefense)
 - **Nanomedicine** (smart dialysis)
 - **Water-Energy** (mineral water, efficiently)

Biology ↔ inorganic nanostructures
Quantum Modeling ↔ experiments

Acknowledgements

- **Funding**
 1. DOE: Sandia's Water Desalination Program
 2. DOE: Sandia's LDRD program (BST, ERN)
 3. NIH: National Nanomedicine Center
- **Compute time (~ 300,000 hours cpu)**
 1. Sandia Computing: Thunderbird
 2. National Center of Supercomputing Applications (NCSA), UIUC



- **Collaborative Science, Engineering, Technology Teams**

Desalination Team at Sandia

PI: Susan Rempe

- Jeff Brinker
- Kevin Leung
- Steve Plimpton
- Dubravko Sabo (*postdoc*)
- Seema Singh
- Sameer Varma (*postdoc*)
- Ying-Bing Jiang (*postdoc*)
- Tom Mayer (project manager)

NIH Center for Design of Biomimetic Nanoconductors

(<http://www.nanoconductor.org>)

PI: Eric Jakobsson, UIUC

Senior Scientists

- Narayan Aluru (UIUC)
- Hagan Baylay (U Oxford)
- Jeff Brinker (SNL)
- Millicent Firestone (ANL)
- David LaVan (Yale)
- Kevin Leung (SNL)
- Steve Plimpton (SNL)
- Atul Parikh (UC Davis)
- Umberto Ravaglioli (UIUC)
- Susan Rempe (SNL PI)
- Benoit Roux (U Chicago)
- Marco Saraniti (IIT)
- Larry Scott (IIT)



Ion Discrimination by Nanoscale Design



**Susan Rempe, Sameer Varma, Dubravko Sabo,
Kevin Leung, Ying-bing Jiang, Jeff Brinker**
Sandia National Labs

K^+ transfer from Water into fully flexible K-channel binding sites



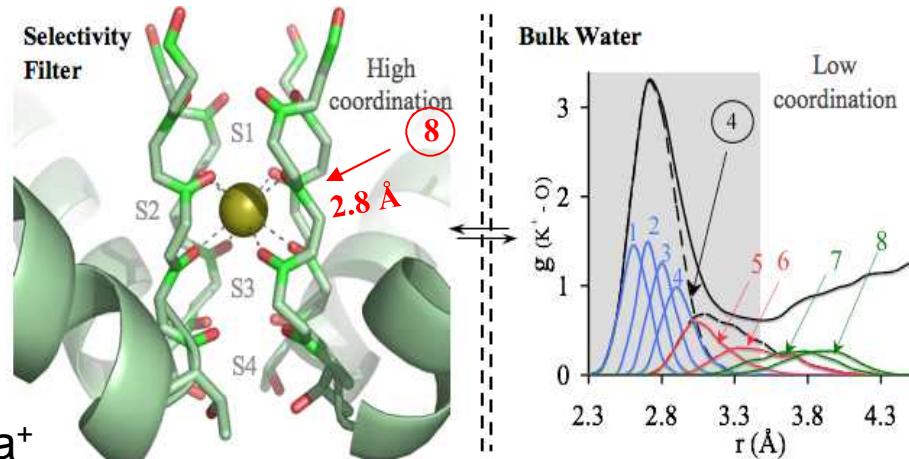
- Build molecular models and predict:
 - Optimized complexes; overlay X-ray structure
 - K⁺ transfer thermodynamics
 - Ion selectivity
- **Achieve a model that represents known data:**
 - reproduces measured ion channel properties
 - reveals new determinants of selectivity: environment & coordination

K^+/Na^+ Ion Discrimination Problem:

How do K-selective channels work?

Traditional Mechanism

- Ion coordination assumed **fixed, mimicked**
- Specific cavity size fits permeant K^+ , not Na^+



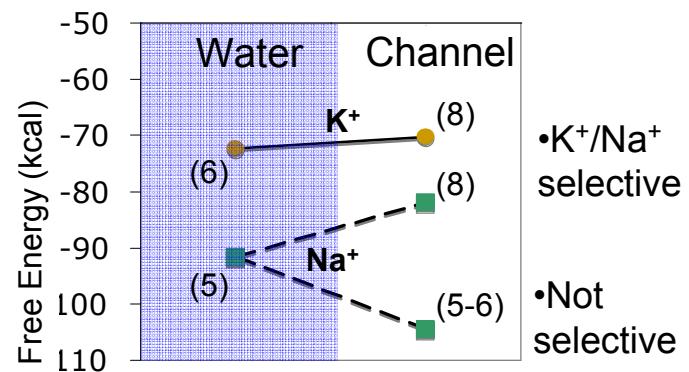
Over-coordination Mechanism

- Ion coordination linked to environment
- Specific (C=O) ligand number 'fits' K^+ , not Na^+

• "The caress of the surroundings, the crowding of the ligands." Jordan (2007), New & Notable *BJ*

Impact

- New explanation of K-channel selectivity
- Engineering parameters
- Health, Water, Nanoengineering



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