

ETH

Eidgenössische Technische Hochschule Zürich
Swiss Federal Institute of Technology Zurich

CSCS
Swiss National Supercomputing Centre



The DCA++ Story

How new algorithms, new computers, and innovative software design allow us to solve real simulation problems of high temperature superconductivity

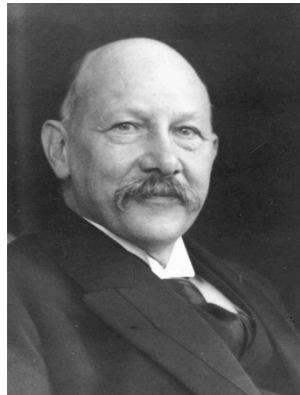
Thomas C. Schulthess

Advance Scientific Computing Advisory Committee
Meeting, Washington DC, March 3-4, 2009

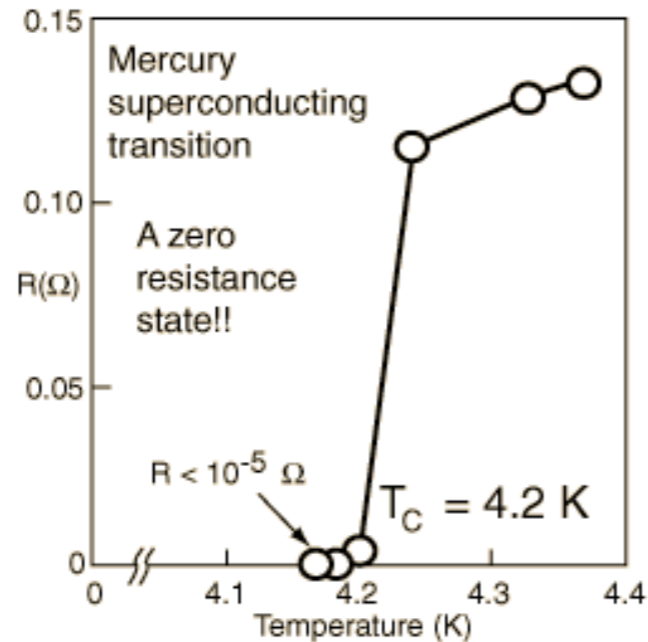


Superconductivity: a state of matter with zero electrical resistivity

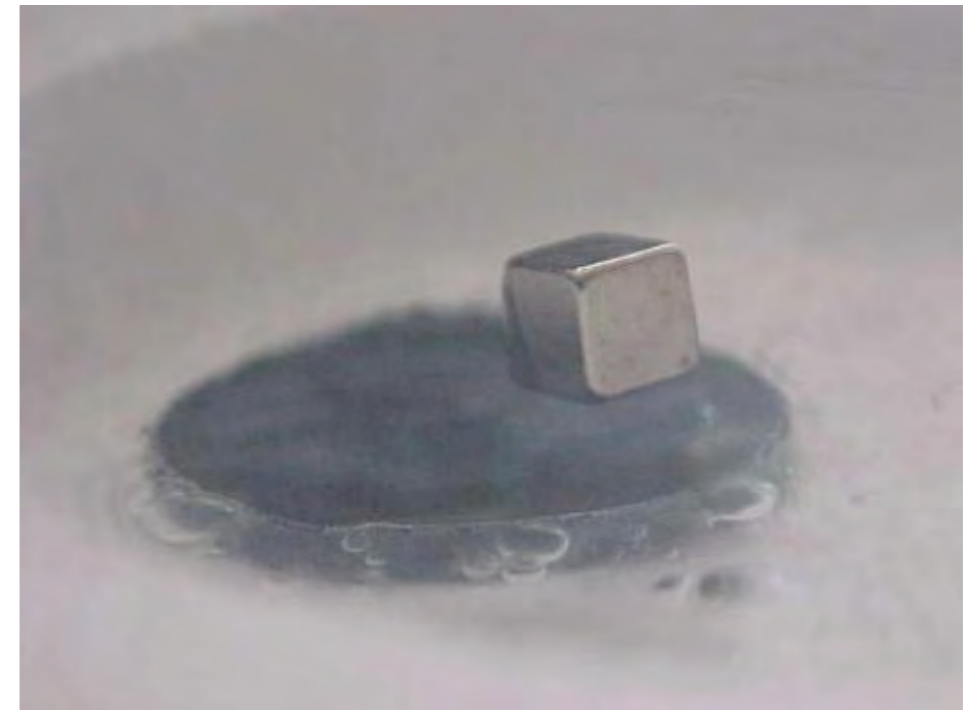
Discovery 1911



Heike Kamerlingh Onnes (1853-1926)



Superconductor repels magnetic field
Meissner and Ochsenfeld, **Berlin 1933**



Microscopic Theory for Superconductivity 1957

PHYSICAL REVIEW

VOLUME 108, NUMBER 5

DECEMBER 1, 1957

Theory of Superconductivity*

J. BARDEEN, L. N. COOPER,† AND J. R. SCHRIEFFER‡
Department of Physics, University of Illinois, Urbana, Illinois
(Received July 8, 1957)

A theory of superconductivity is presented, based on the fact that the interaction between electrons resulting from virtual exchange of phonons is attractive when the energy difference between the electrons states involved is less than the phonon energy, $\hbar\omega$. It is favorable to form a superconducting phase when this attractive interaction dominates the repulsive screened Coulomb interaction. The normal phase is described by the Bloch individual-particle model. The ground state of a superconductor, formed from a linear combination of normal state configurations in which electrons are virtually excited in pairs of opposite spin and momentum, is lower in energy than the normal state by amount proportional to an average $(\hbar\omega)^2$, consistent with the isotope effect. A mutually orthogonal set of excited states in

one-to-one correspondence with those of the normal phase is obtained by specifying occupation of certain Bloch states and by using the rest to form a linear combination of virtual pair configurations. The theory yields a second-order phase transition and a Meissner effect in the form suggested by Pippard. Calculated values of specific heats and penetration depths and their temperature variation are in good agreement with experiment. There is an energy gap for individual-particle excitations which decreases from about $3.5kT_c$ at $T=0^\circ\text{K}$ to zero at T_c . Tables of matrix elements of single-particle operators between the excited-state superconducting wave functions, useful for perturbation expansions and calculations of transition probabilities, are given.



Scanned at the American Institute of Physics



Scanned at the American Institute of Physics

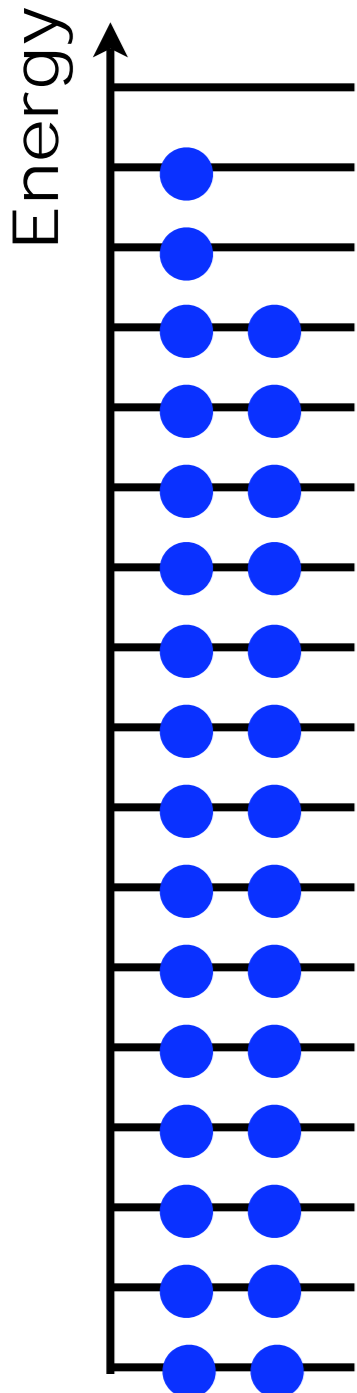


Scanned at the American Institute of Physics

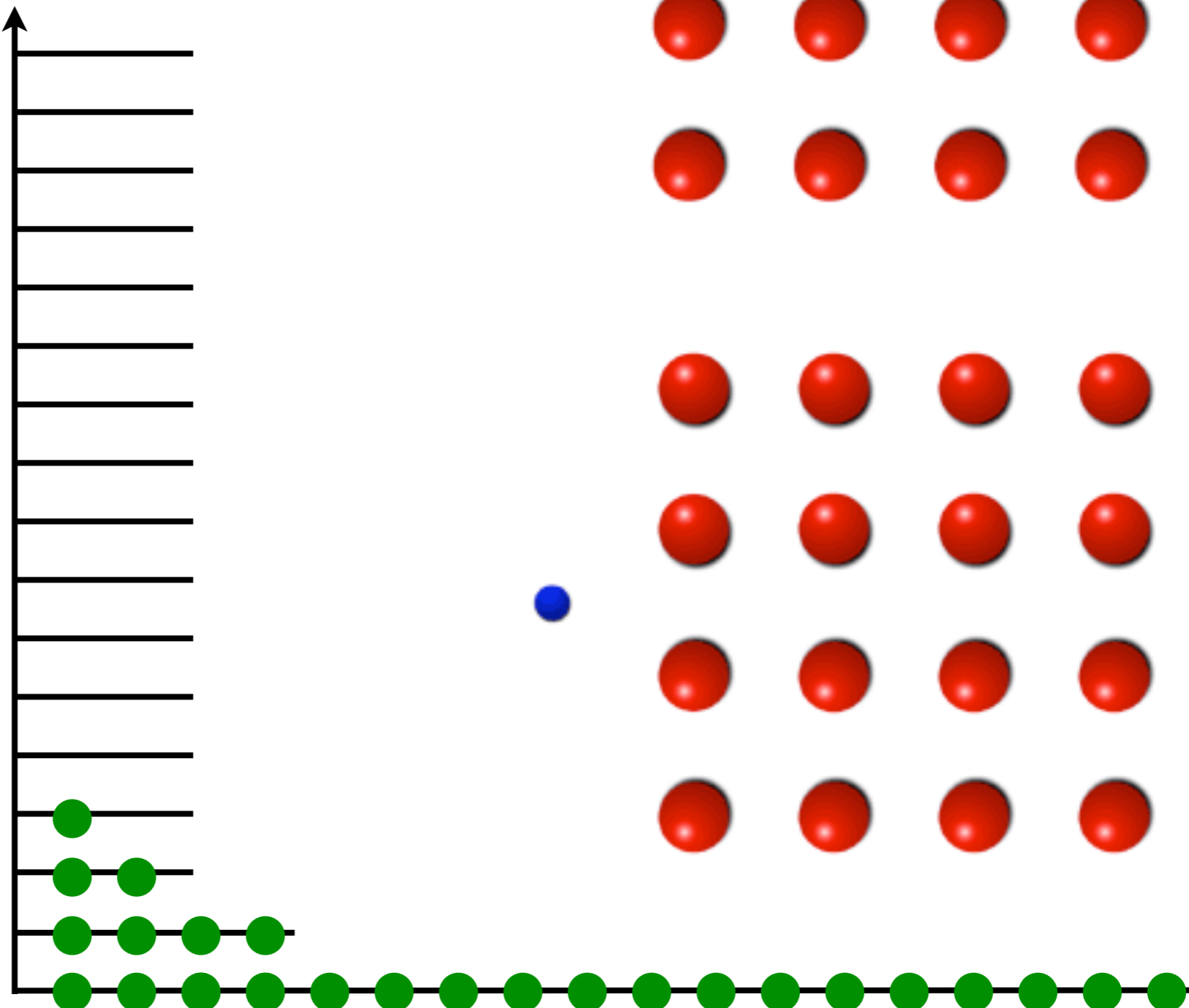
BCS Theory generally accepted in the early 1970s

Fermions, Bosons, and Cooper Pairs

Fermions (electron)

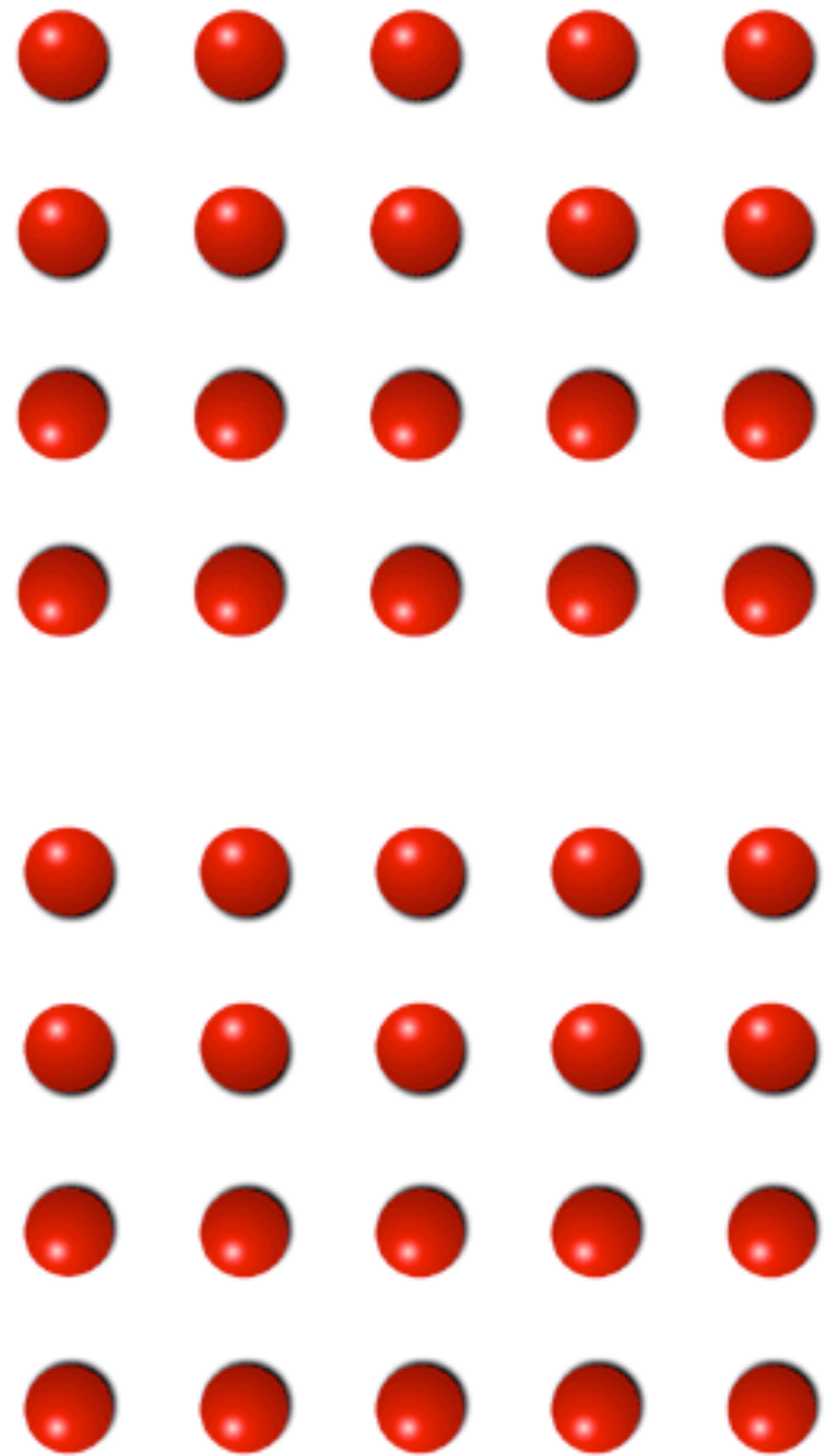


Bosons-like



•

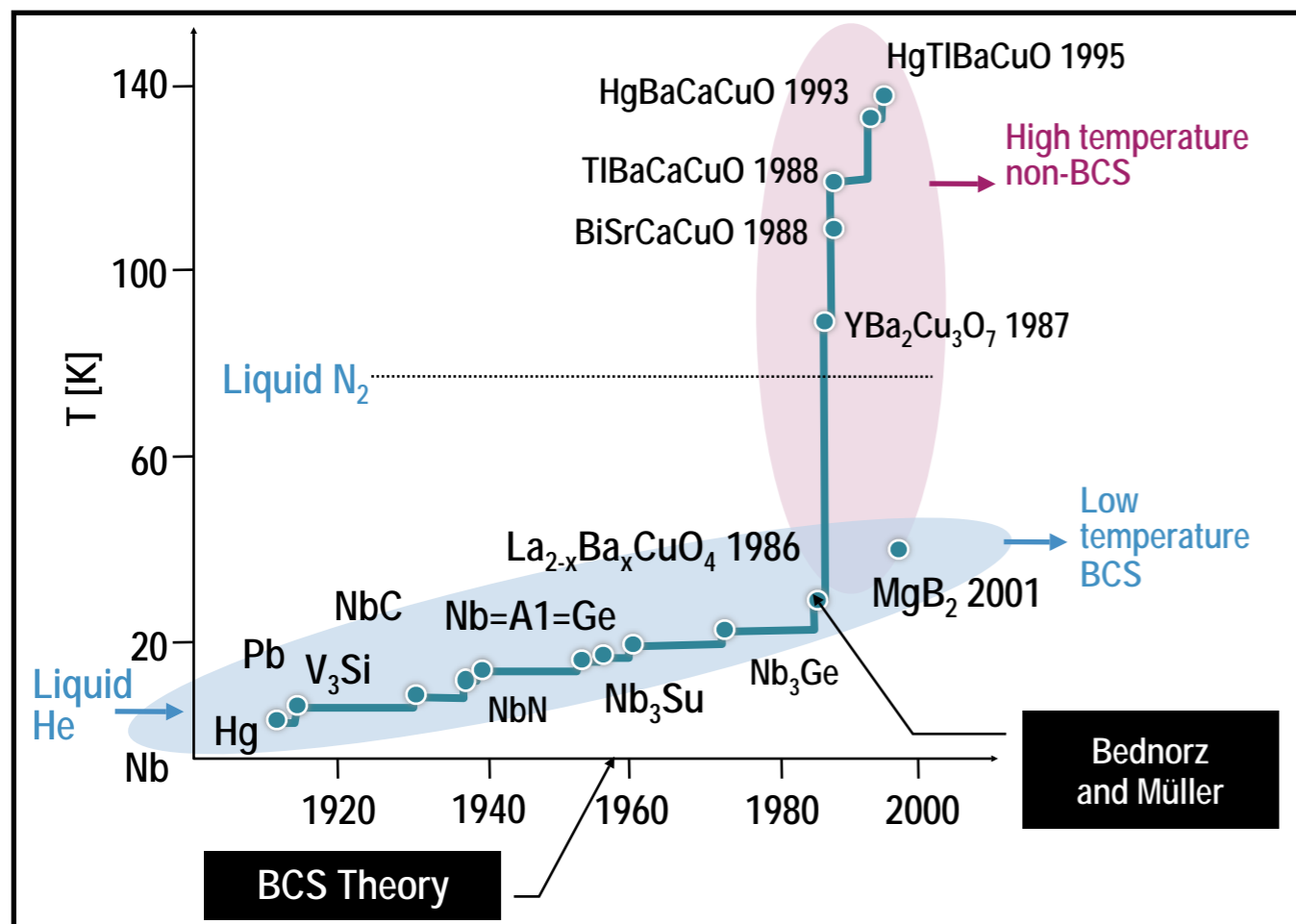
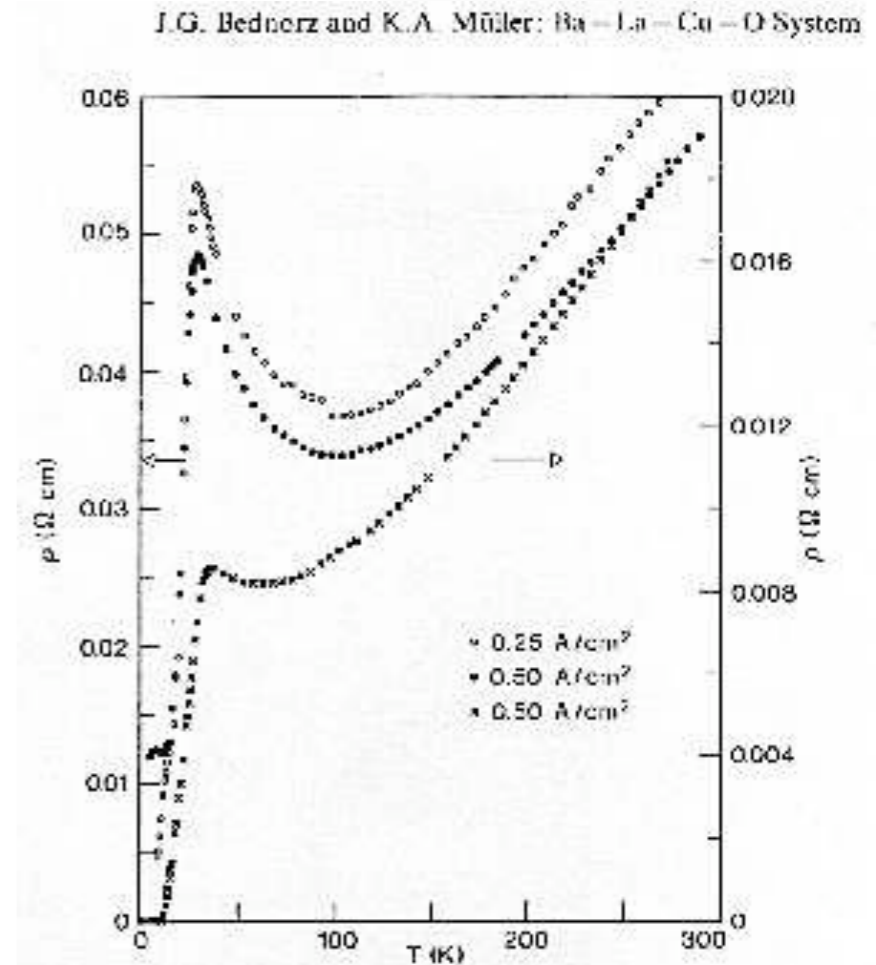
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Superconductivity in the cuprates



- Discovered in 1986 by Bednorz and Müller
- Totally different materials
 - In the normal state conventional superconductors are metals cuprates are insulators or poor conductors



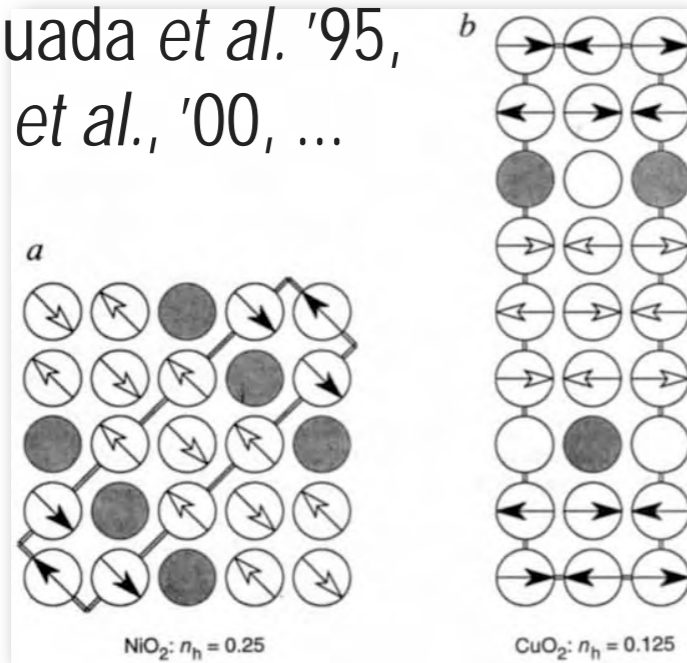
Twenty years later

- ▶ No predictive power for T_c in known materials
- ▶ No predictive power for design of new SC materials
- ▶ No explanation for other unusual properties of cuprates (pseudogap, transport, ...)
- ▶ Only partial consensus on which materials aspects are essential for high- T_c superconductivity
- ▶ No controlled solution for proposed models

The role of inhomogeneities

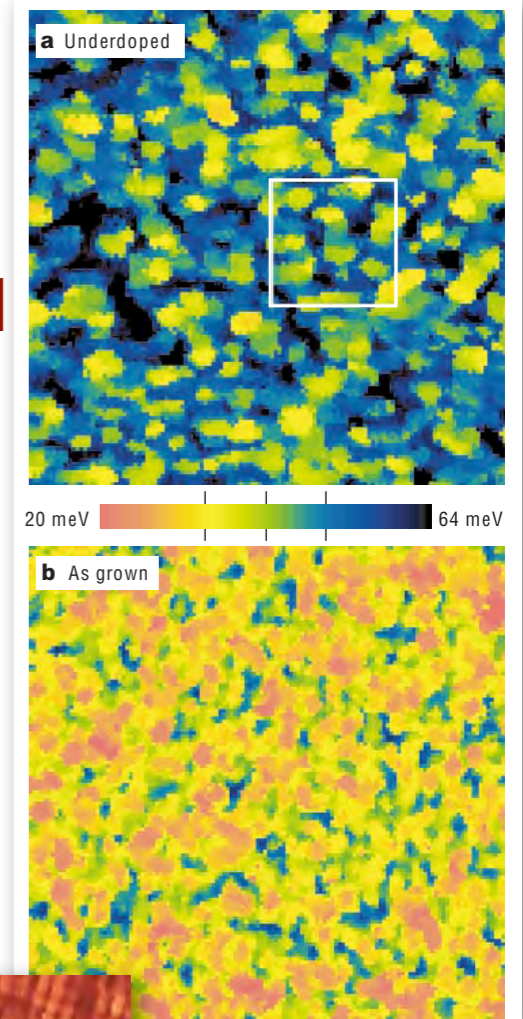
Stripes in neutron scattering:

Tranquada *et al.* '95,
Mook *et al.*, '00, ...



Random SC gap modulations in STM (BSCCO):

Lang *et al.* '02

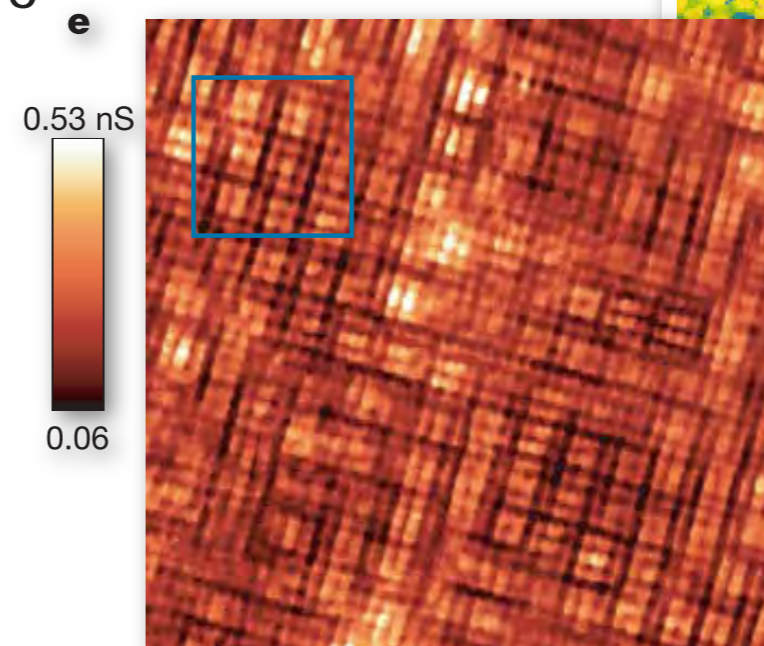
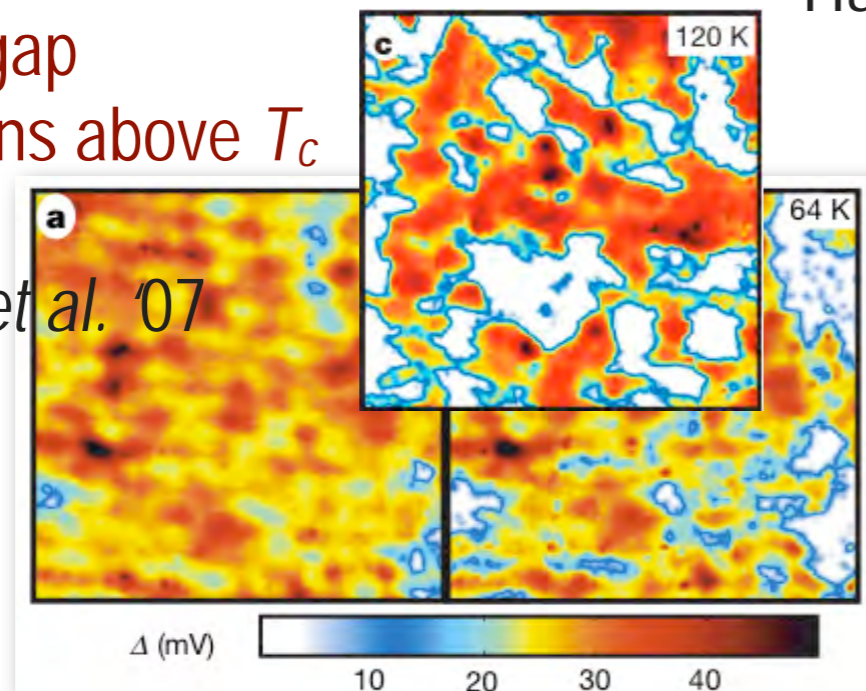


Charge ordered "checkerboard" state (Na doped cuprates):

Hanaguri *et al.* '04

Random gap modulations above T_c (BSCCO):

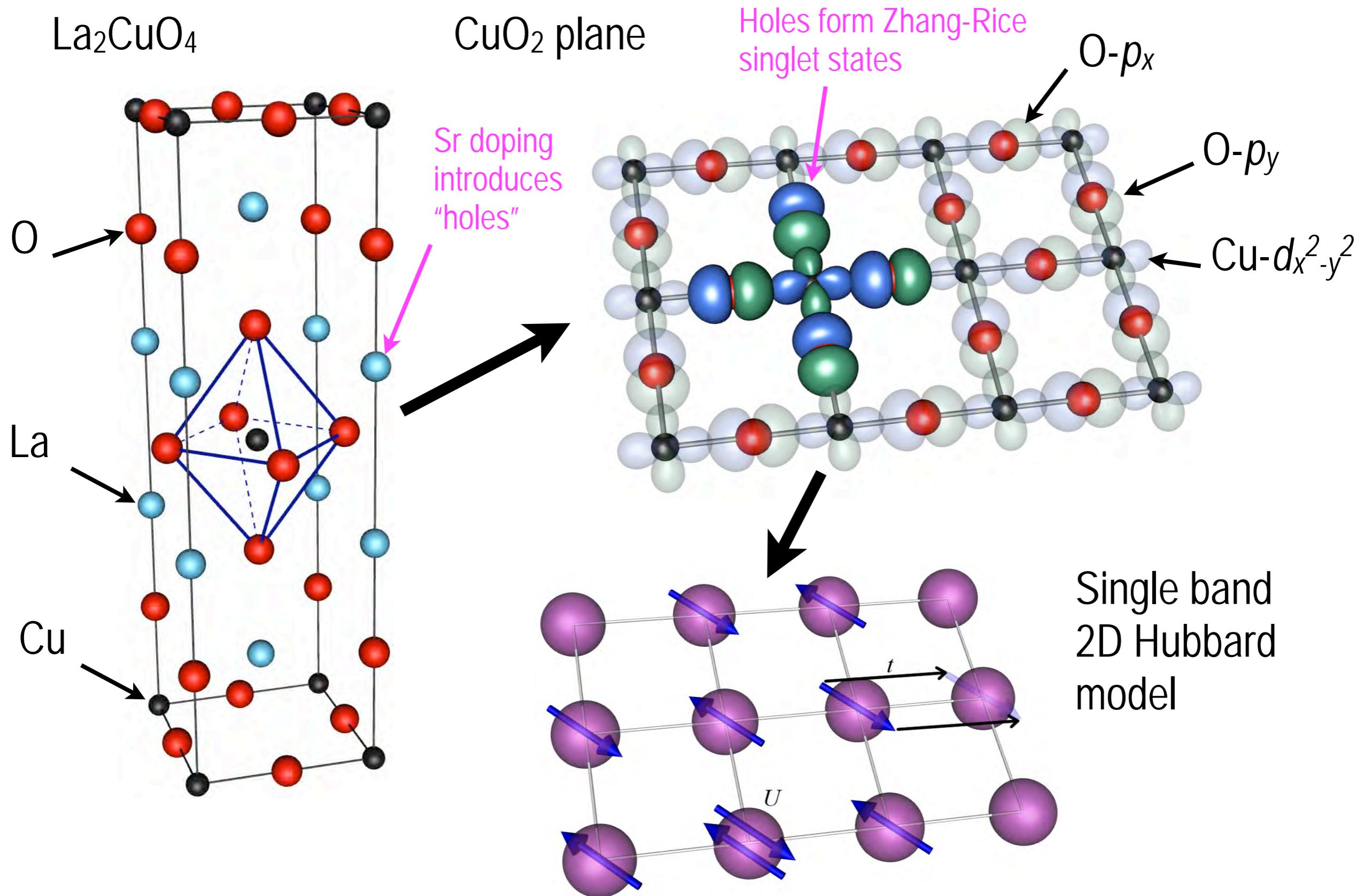
Gomes *et al.* '07



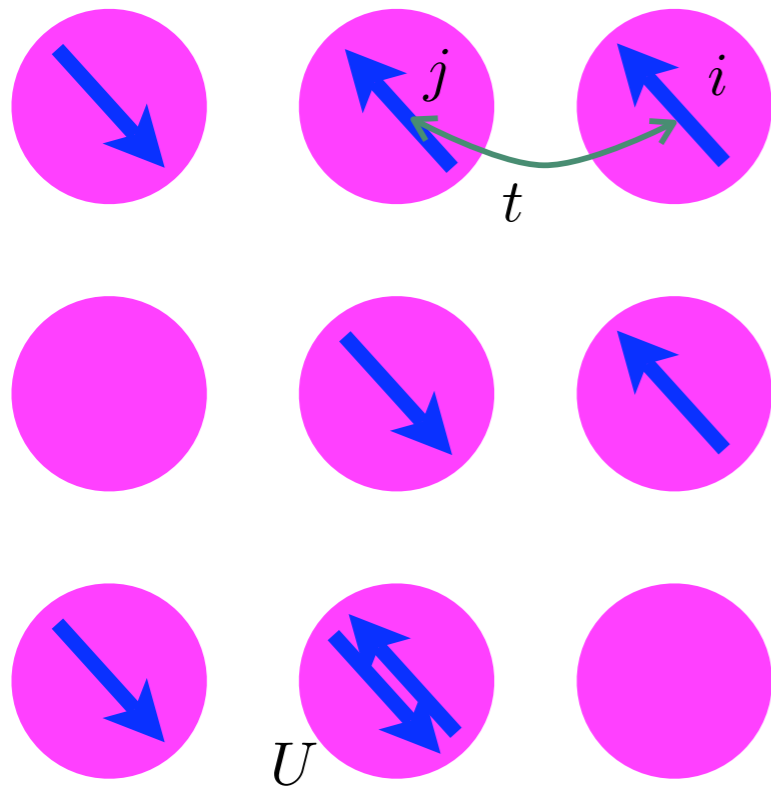
Outline

- Brief introduction into superconductivity and the cuprates
- Background: The two dimensional Hubbard model and the DCA/QMC method
- Simulational studies with the DCA/QMC method
- Algorithmic improvements and a method to study effects of disorder and nanoscale inhomogeneities
 - Accelerating Hirsch-Fye QMC with delayed updates
 - Mixed precision and multithreaded implementations (GPU in particular)
 - Disorder averaging and a first study of how disorder affects the superconducting transition temperature
- DCA++, concurrency, scaling, and performance
 - Results for Cray XT4 and first results for a PF/s scale system
- Summary and conclusions

From cuprate materials to the Hubbard model

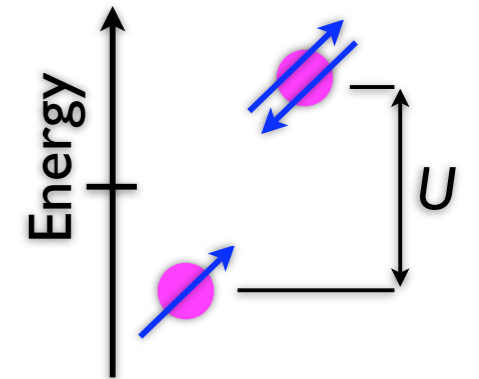


2D Hubbard model and its physics

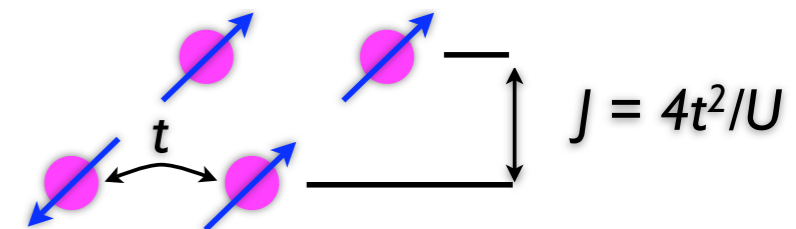


Half filling: number of carriers = number of sites

Formation of a **magnetic moment** when U is large enough

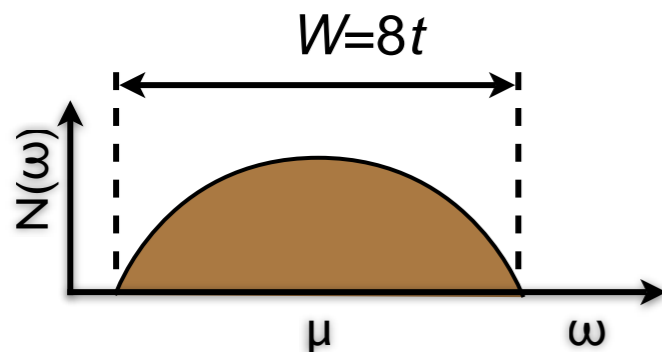


Antiferromagnetic alignment of neighboring moments



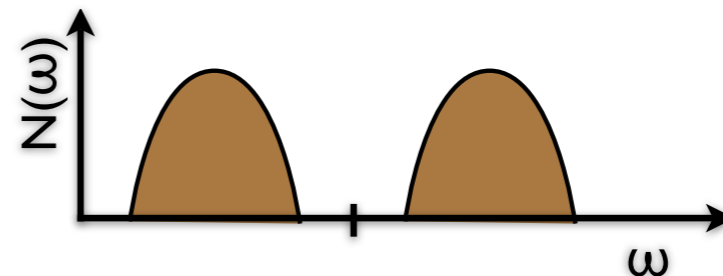
1. When $t \gg U$:

Model describes a metal with band width $W=8t$

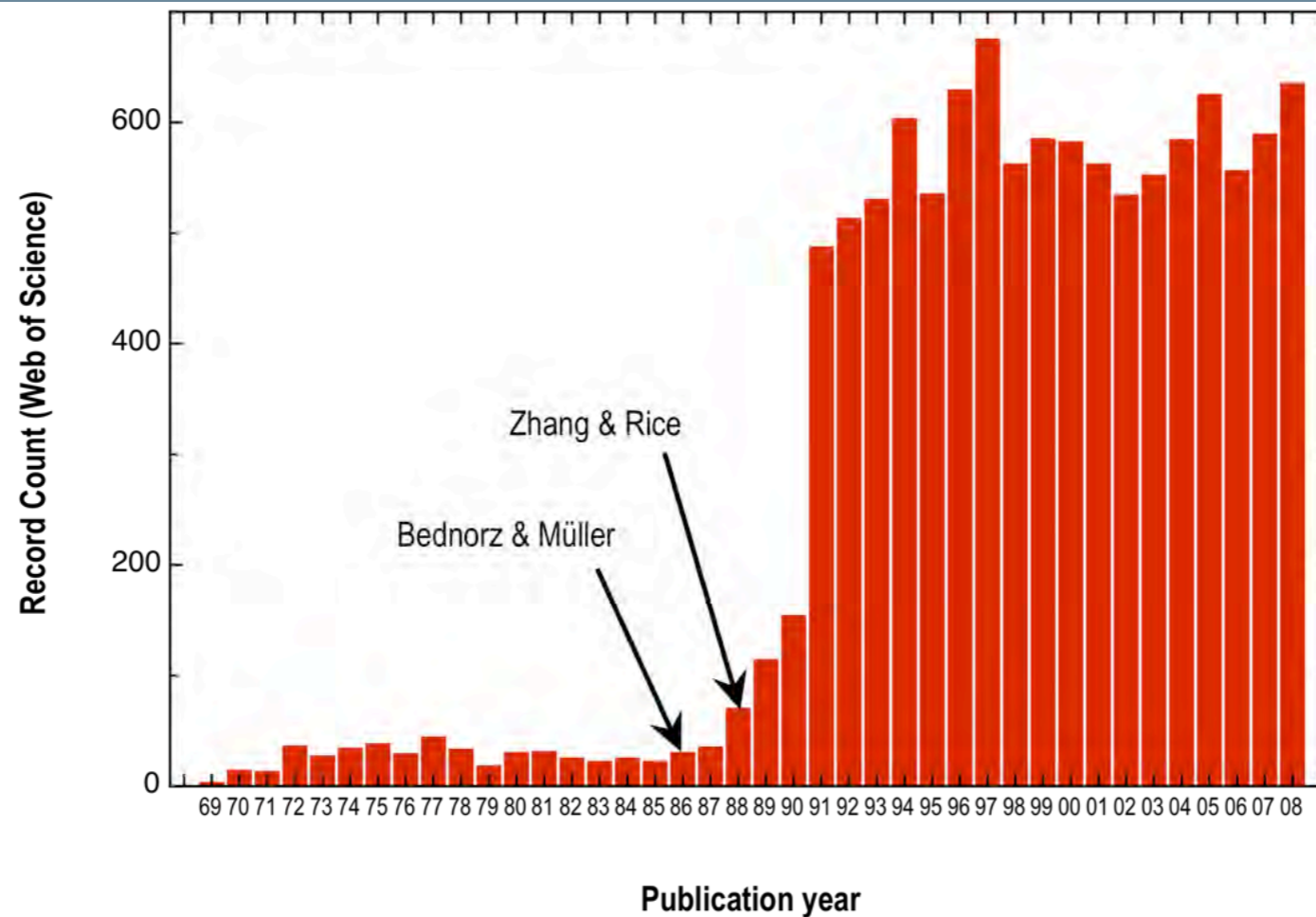


2. When $U \gg 8t$ at half filling (not doped)

Model describes a "Mott Insulator" with antiferromagnetic ground state (as seen experimentally seen in undoped cuprates)



Hubbard model for the cuprates



3. Parameter range relevant for superconducting cuprates

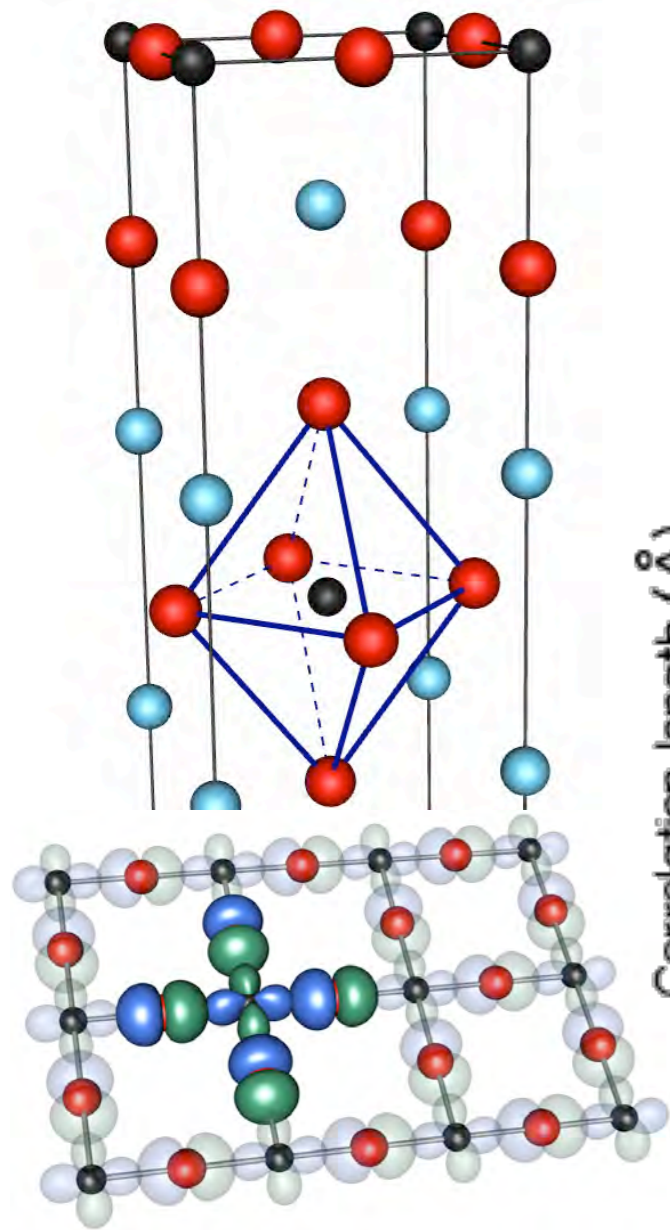
$$U \approx 8t$$

Finite doping levels (0.05 – 0.25)

Typical values: $U \sim 10\text{eV}$; $t \sim 0.9\text{eV}$; $J \sim 0.2\text{eV}$; (0.1eV ~ 10^3 Kelvin)

No simple solution!

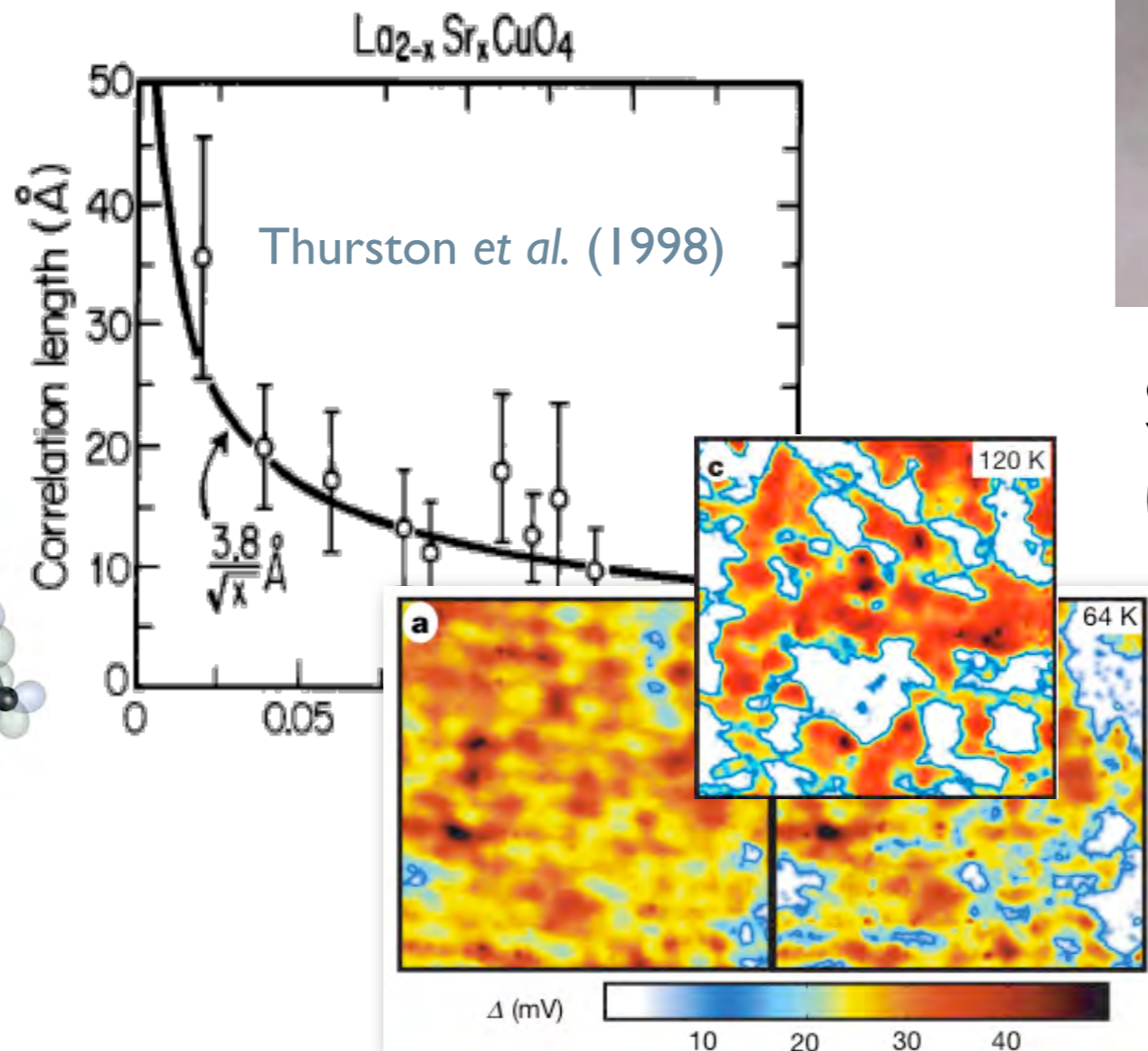
The challenge: a (quantum) multi-scale problem



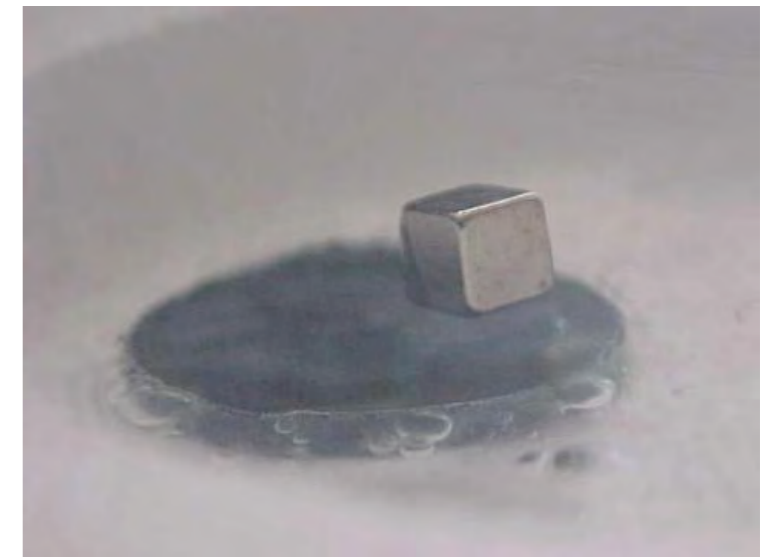
On-site Coulomb repulsion ($\sim A$)

complexity $\sim 4^N$

Antiferromagnetic correlations / nano-scale gap fluctuations



Gomes et al. (2007)

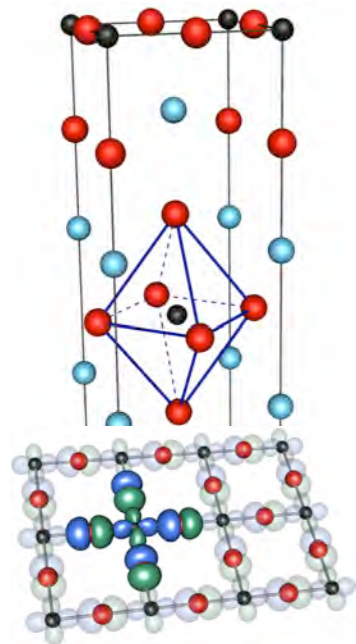


Superconductivity (macroscopic)

$N \sim 10^{23}$

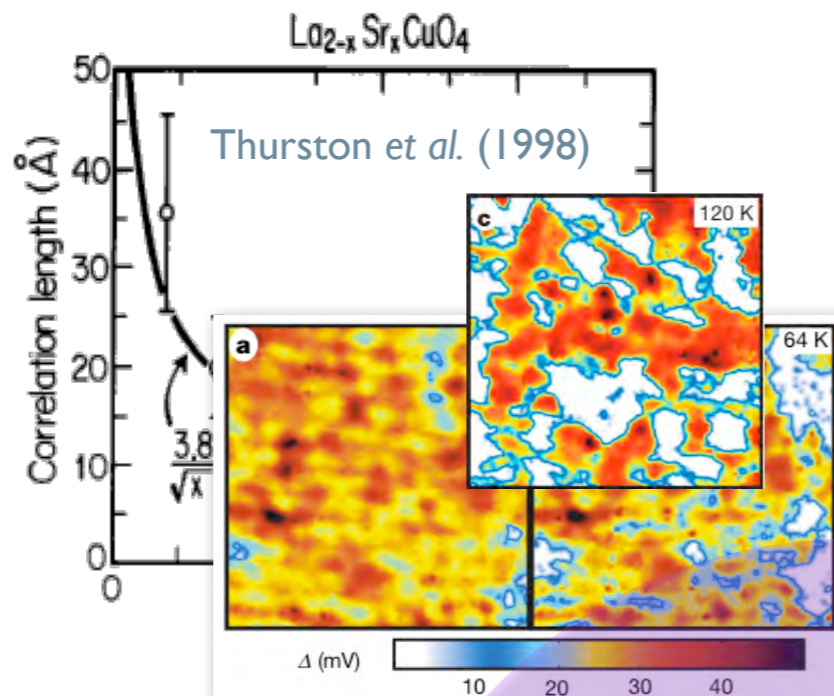
Quantum cluster theories

Maier *et al.*, Rev. Mod. Phys. '05

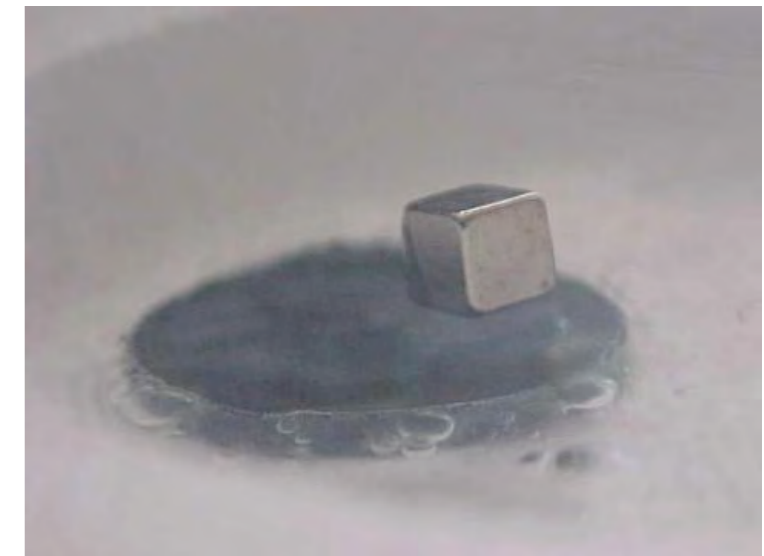


On-site Coulomb repulsion ($\sim A$)

Antiferromagnetic correlations / nano-scale gap fluctuations



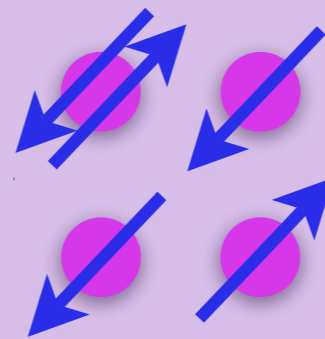
Gomes *et al.* (2007)



Superconductivity (macroscopic)

Explicitly treat correlations within a localized cluster

Treat macroscopic scales within mean-field



Coherently embed cluster into effective medium

Green's functions in quantum many-body theory

Noninteracting Hamiltonian &
Green's function

$$H_0 = \left[-\frac{1}{2} \nabla^2 + V(\vec{r}) \right]$$

$$\left[i \frac{\partial}{\partial t} - H_0 \right] G_0(\vec{r}, t, \vec{r}', t') = \delta(\vec{r} - \vec{r}') \delta(t - t')$$

Fourier transform & analytic continuation: $z^\pm = \omega \pm i\epsilon$ $G_0^\pm(\vec{r}, z) = [z^\pm - H_0]^{-1}$

Hubbard Hamiltonian

$$H = -t \sum_{\langle ij \rangle, \sigma} c_{i\sigma}^\dagger c_{j\sigma} + U \sum_i n_{i\uparrow} n_{i\downarrow} \quad n_{i\sigma} = c_{i\sigma}^\dagger c_{i\sigma}$$

Hide symmetry in algebraic properties of field operators

$$c_{i\sigma} c_{j\sigma'} + c_{j\sigma'} c_{i\sigma} = 0$$
$$c_{i\sigma} c_{j\sigma'}^\dagger + c_{j\sigma'}^\dagger c_{i\sigma} = \delta_{ij} \delta_{\sigma\sigma'}$$

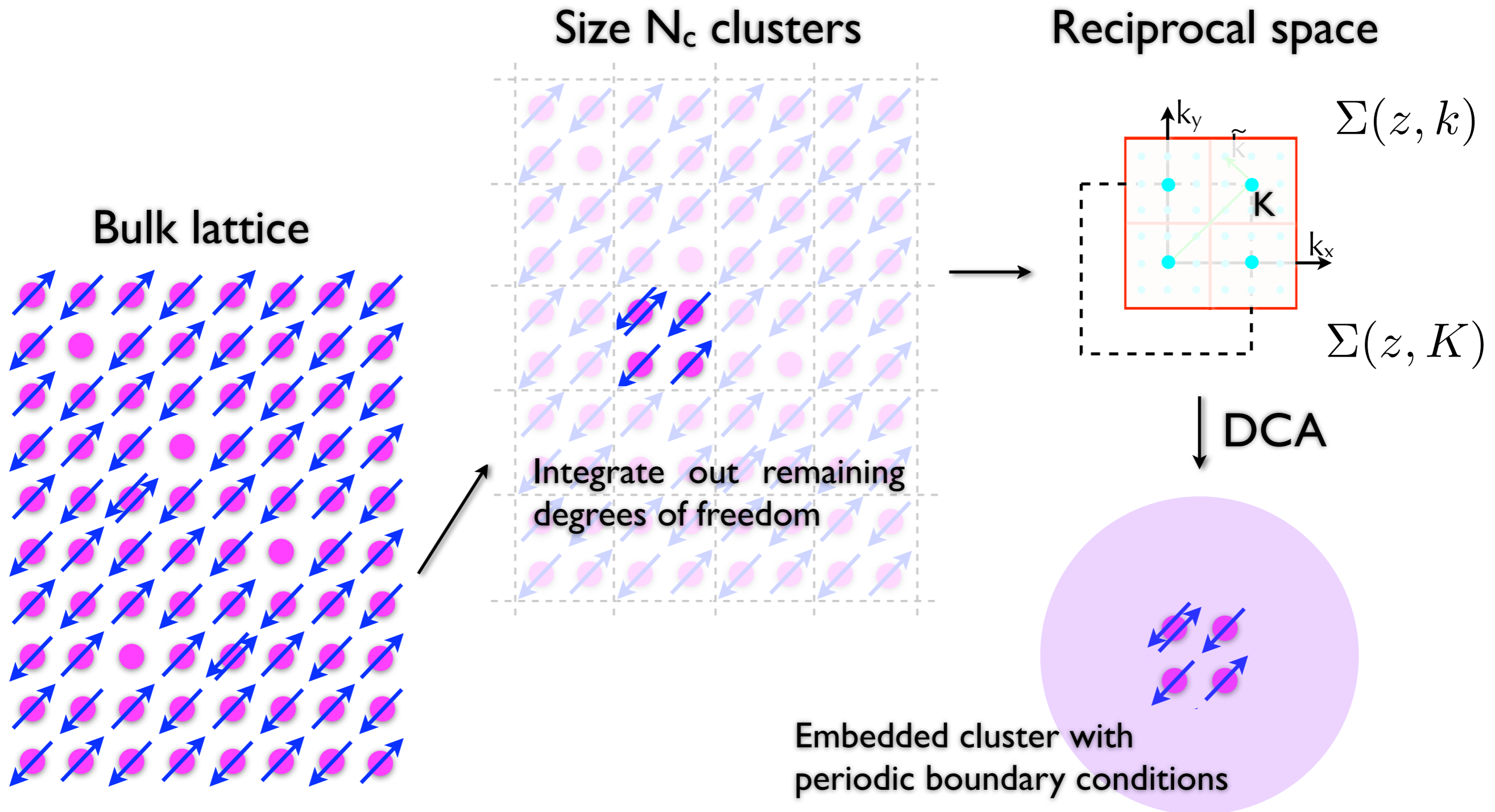
Green's function $G_\sigma(r_i, \tau; r_j, \tau') = - \left\langle \mathcal{T} c_{i\sigma}(\tau) c_{j\sigma}^\dagger(\tau') \right\rangle$

Spectral representation

$$G_0(k, z) = [z - \epsilon_0(k)]^{-1}$$

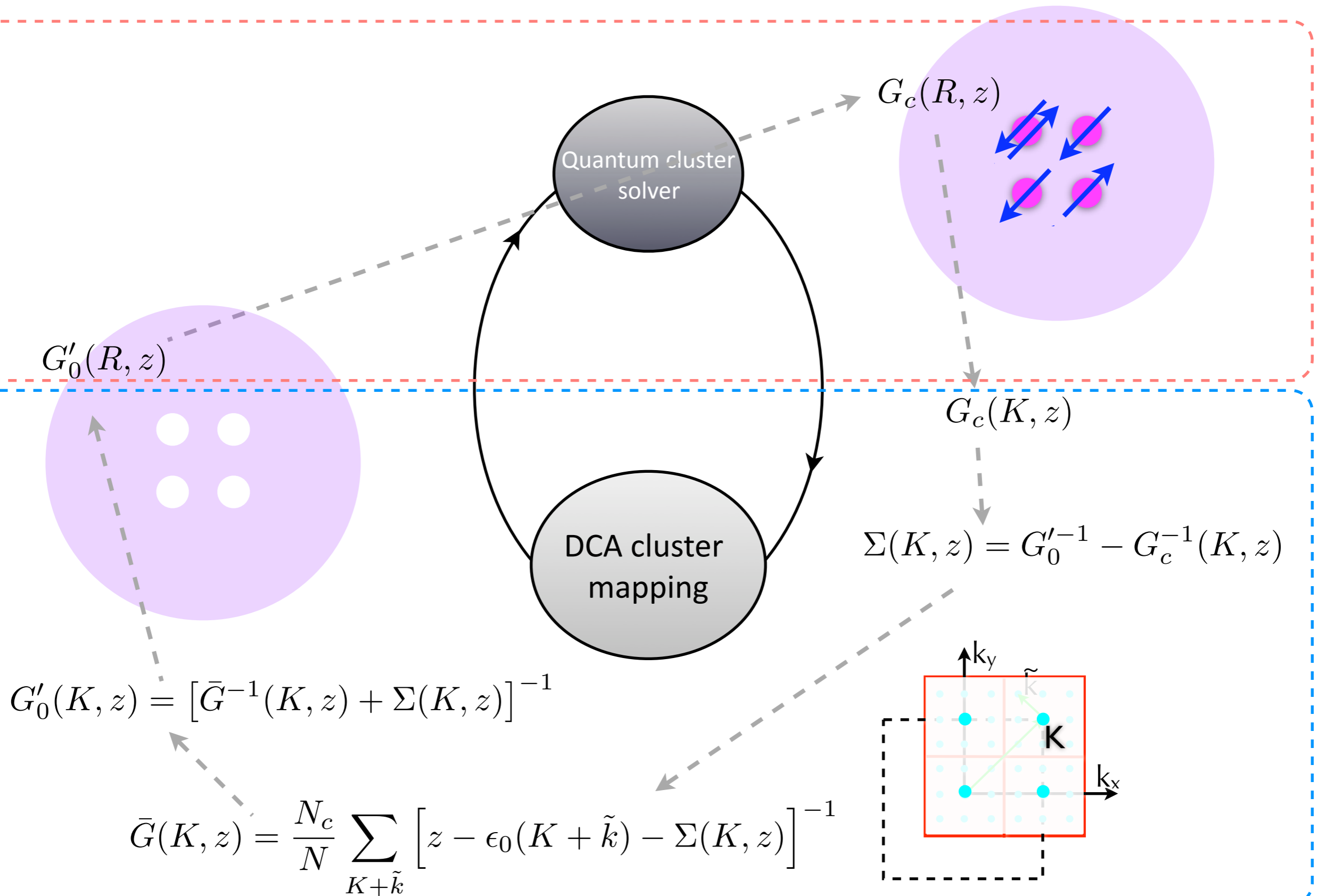
$$G(k, z) = [z - \epsilon_0(k) - \Sigma(k, z)]^{-1}$$

Sketch of the Dynamical Cluster Approximation



Solve many-body problem with quantum Monte Carlo on cluster
➤ Essential assumption: Correlations are short ranged

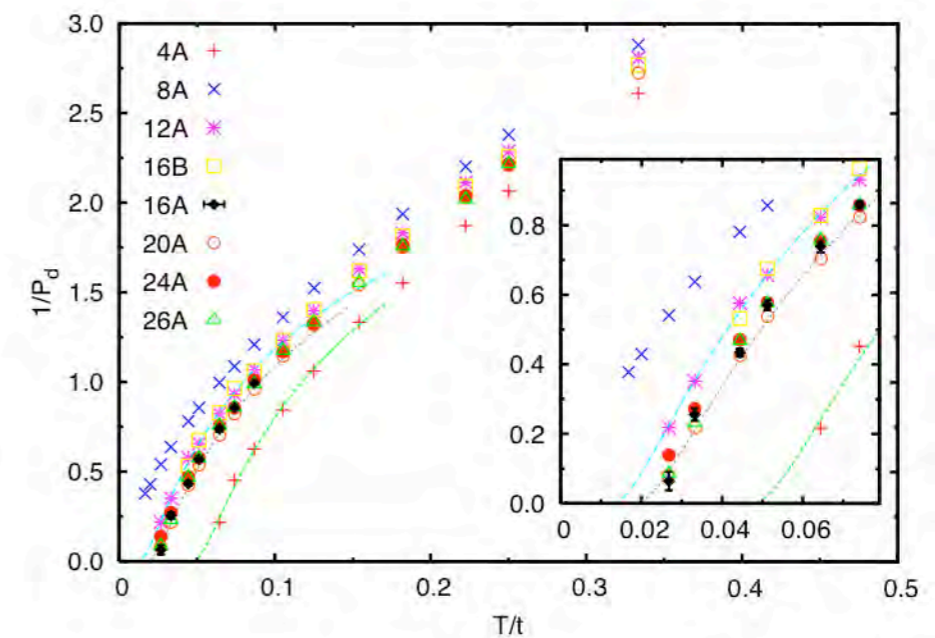
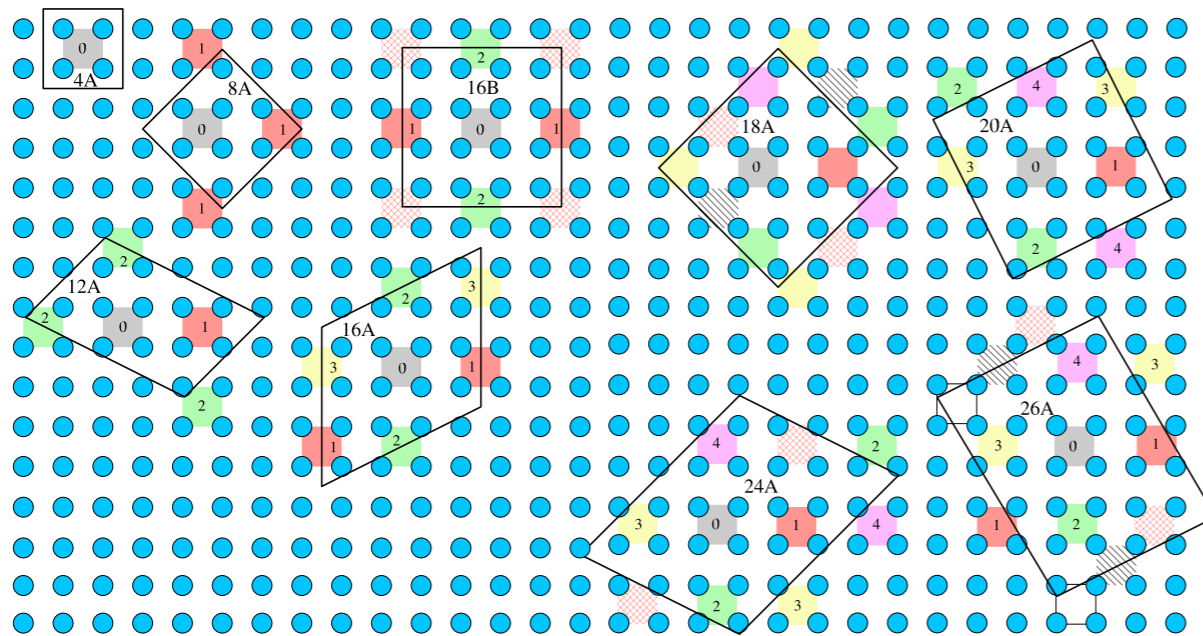
DCA method: self-consistently determine the “effective” medium



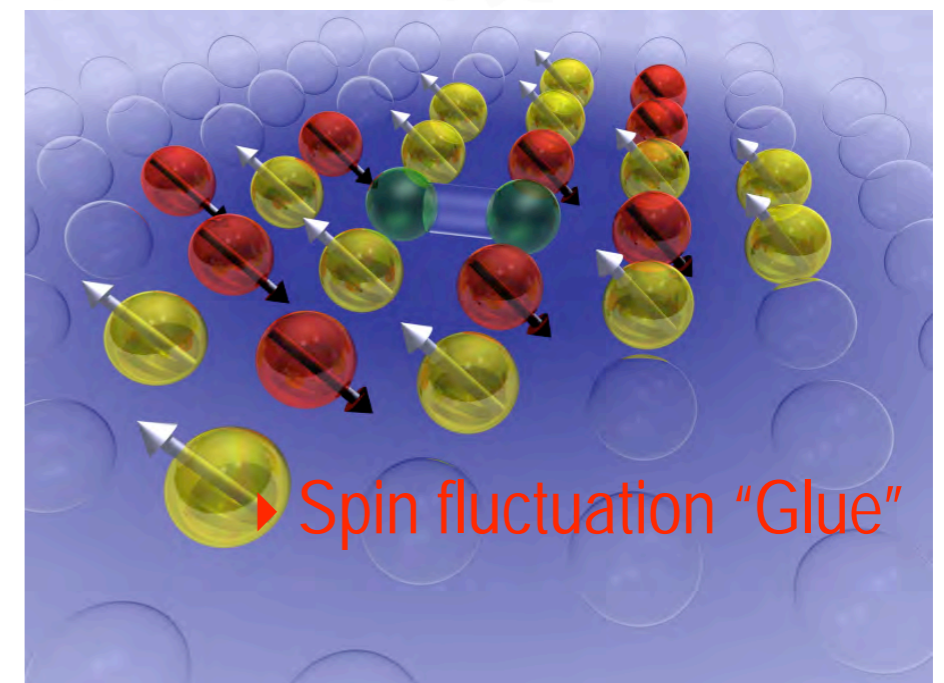
Systematic solution and analysis of the pairing mechanism in the 2D Hubbard Model



- First systematic solution demonstrates existence of a superconducting transition in 2D Hubbard model Maier, et al., Phys. Rev. Lett. 95, 237001 (2005)

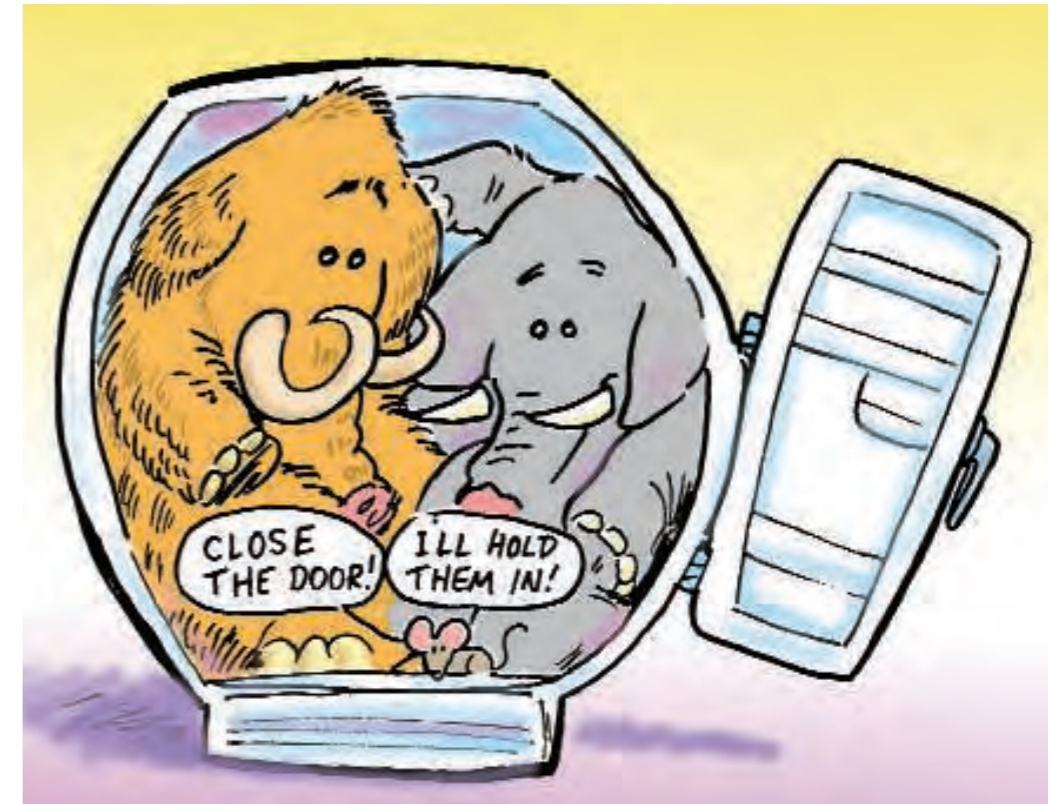


- Study the mechanism responsible for pairing in the model
 - Analyze the particle-particle vertex
 - Pairing is mediated by spin fluctuations Maier, et al., Phys. Rev. Lett. 96 47005 (2006)

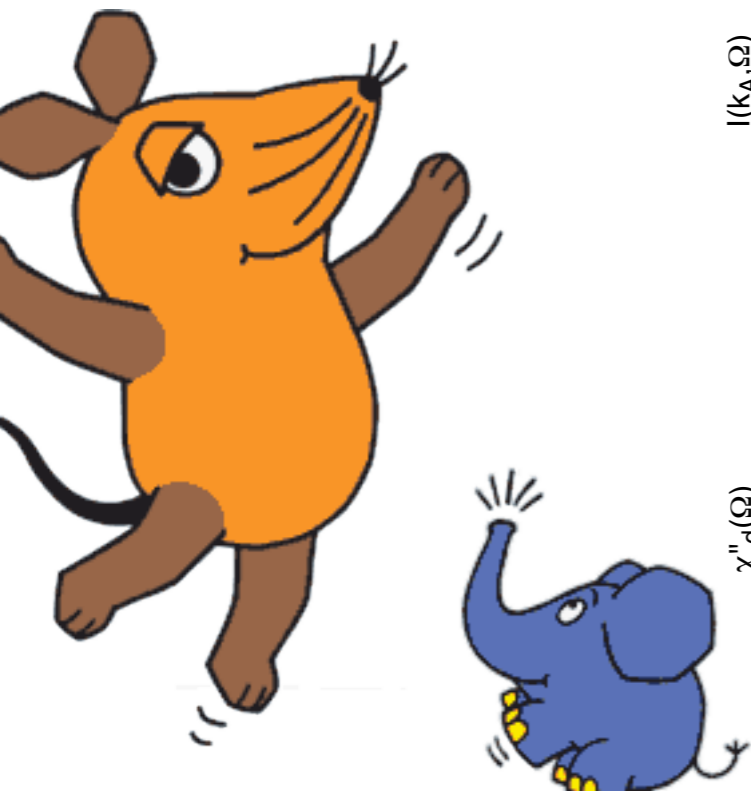
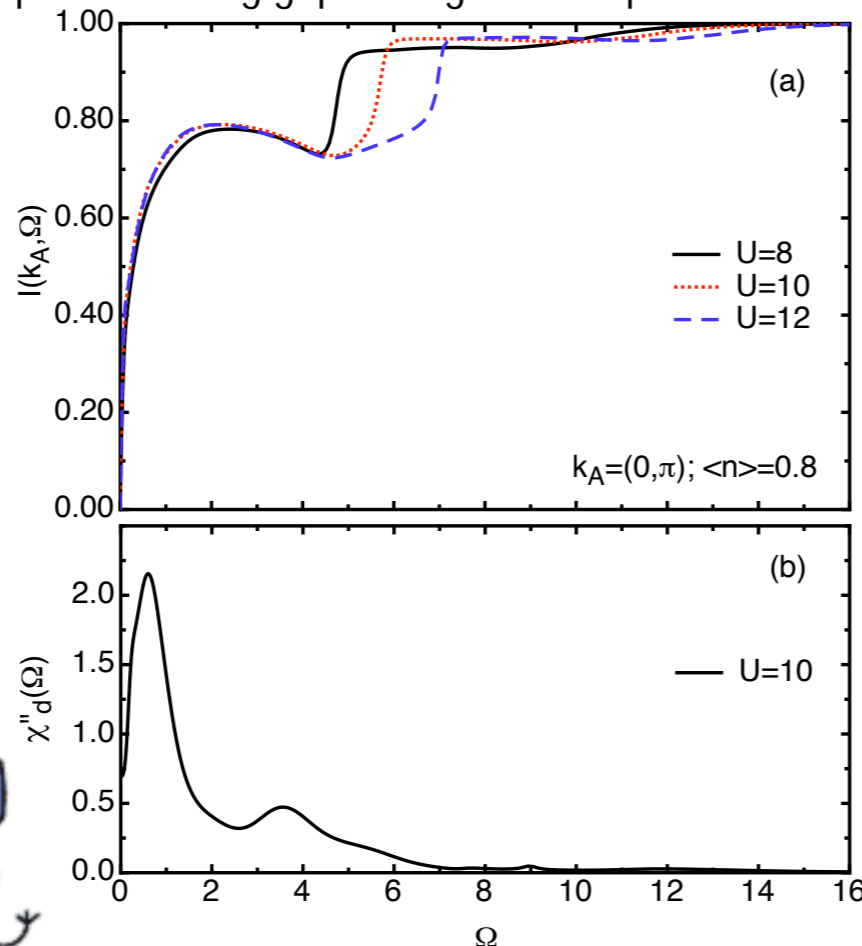


Moving toward a resolution of debate over pairing mechanism in the 2D Hubbard model

- “We have a mammoth (U) and an elephant (J) in our refrigerator - do we care much if there is also a mouse?”
 - P.W. Anderson, Science **316**, 1705 (2007)
 - see also www.sciencemag.org/cgi/eletters/316/5832/1705 “Scalapino is not a glue sniffer”
- Relative importance of resonant valence bond and spin-fluctuation mechanisms
 - Maier et al., Phys. Rev. Lett. **100** 237001 (2008)



Fraction of superconducting gap arising from frequencies $\leq \Omega$



Both retarded spin-fluctuations and non-retarded exchange interaction J contribute to the pairing interaction

Dominant contribution comes from spin-fluctuations!

Hirsch-Fye Quantum Monte Carlo (HF-QMC) for the quantum cluster solver

Hirsch & Fye, Phys. Rev. Lett. 56, 2521 (1998)

Partition function & Metropolis Monte Carlo $Z = \int e^{-E[\mathbf{x}]/k_B T} d\mathbf{x}$

Acceptance criterion for M-MC move: $\min\{1, e^{E[\mathbf{x}_k] - E[\mathbf{x}_{k+1}]}\}$

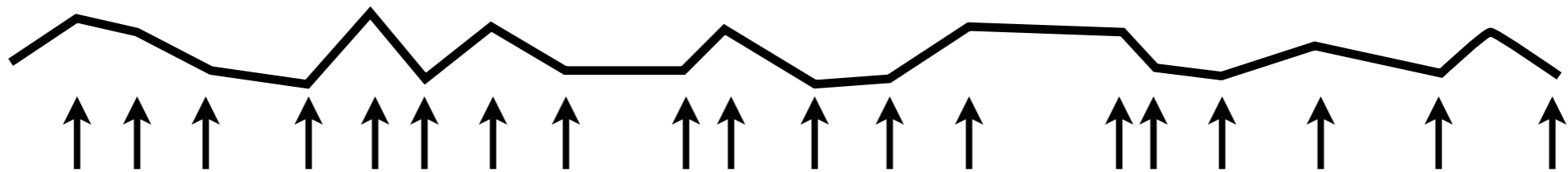
Partition function & HF-QMC: $Z \sim \sum_{s_i, l} \det[\mathbf{G}_c(s_i, l)^{-1}]$

N_c $N_l \approx 10^2$

matrix of dimensions $N_t \times N_t$

$N_t = N_c \times N_l \approx 2000$

Acceptance: $\min\{1, \det[\mathbf{G}_c(\{s_i, l\}_k)] / \det[\mathbf{G}_c(\{s_i, l\}_{k+1})]\}$

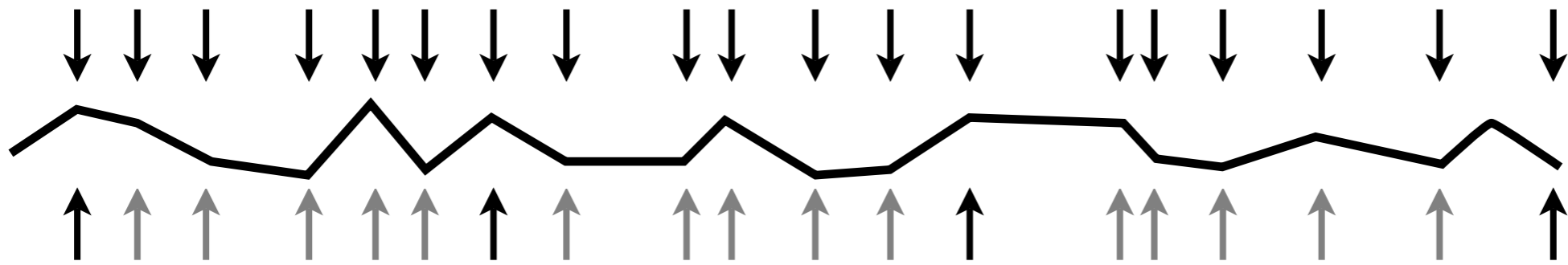


Update of accepted Green's function:

$$\mathbf{G}_c(\{s_i, l\}_{k+1}) = \mathbf{G}_c(\{s_i, l\}_k) + \mathbf{a}_k \times \mathbf{b}_k$$

HF-QMC with Delayed updates (or Ed updates)

$$\mathbf{G}_c(\{s_i, l\}_{k+1}) = \mathbf{G}_c(\{s_i, l\}_k) + \mathbf{a}_k \times \mathbf{b}_k^t$$



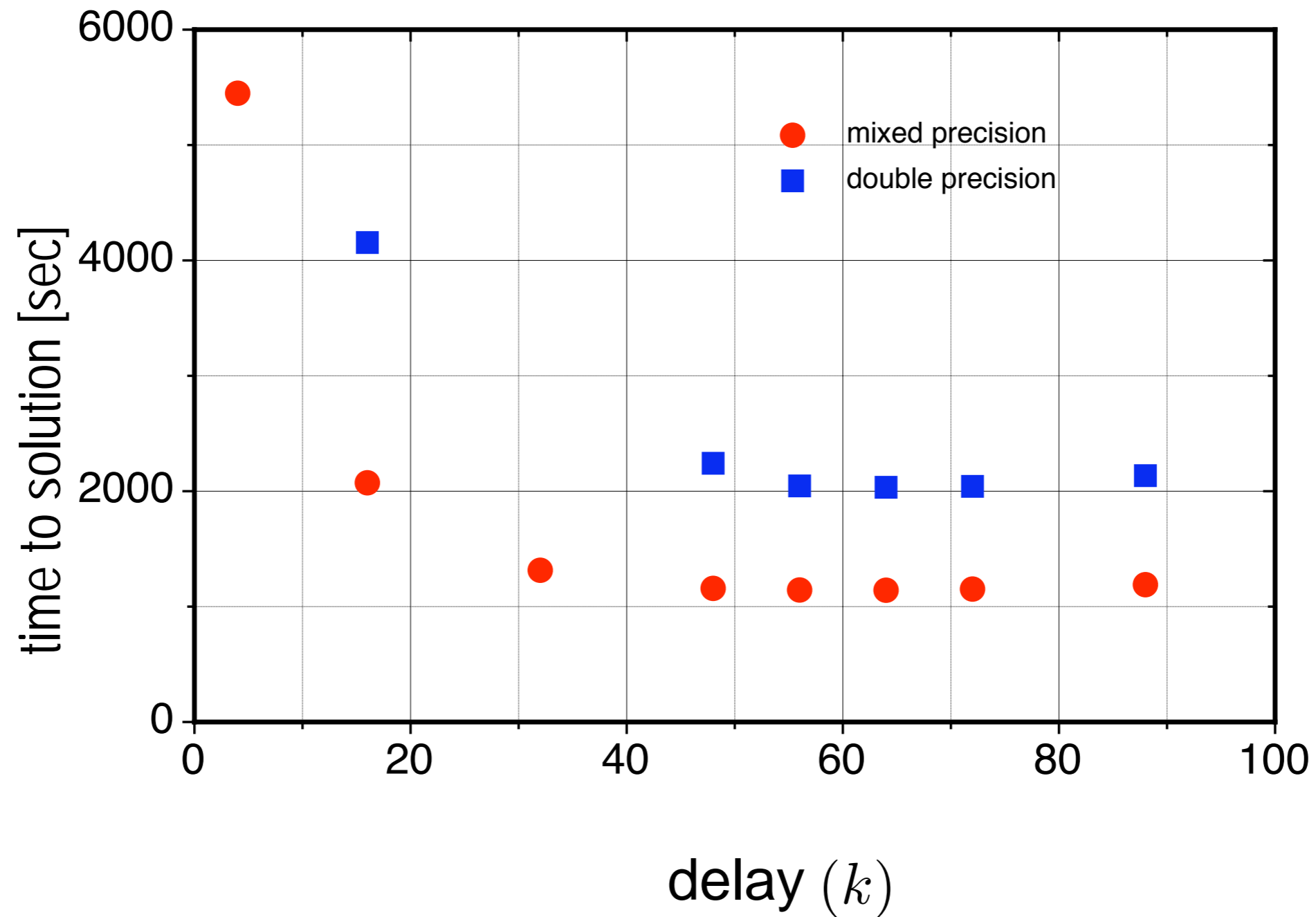
$$\mathbf{G}_c(\{s_i, l\}_{k+1}) = \mathbf{G}_c(\{s_i, l\}_0) + [\mathbf{a}_0 | \mathbf{a}_1 | \dots | \mathbf{a}_k] \times [\mathbf{b}_0 | \mathbf{b}_1 | \dots | \mathbf{b}_k]^t$$

Complexity for k updates remains $\mathcal{O}(kN_t^2)$

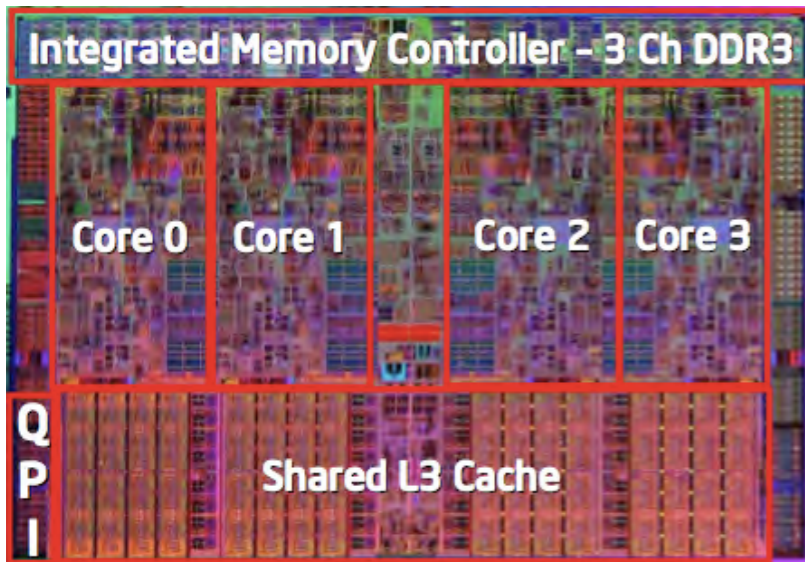
But we can replace k rank-1 updates with one matrix-matrix multiply plus some additional bookkeeping.

Performance improvement with delayed updates

$N_c = 16$ $N_l = 150$ $N_t = 2400$

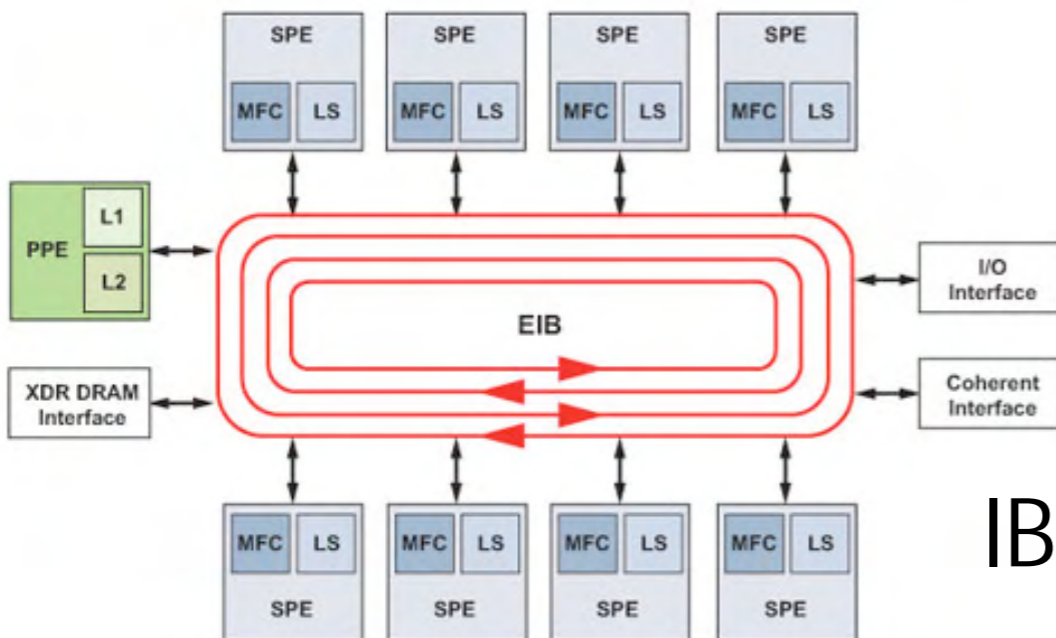
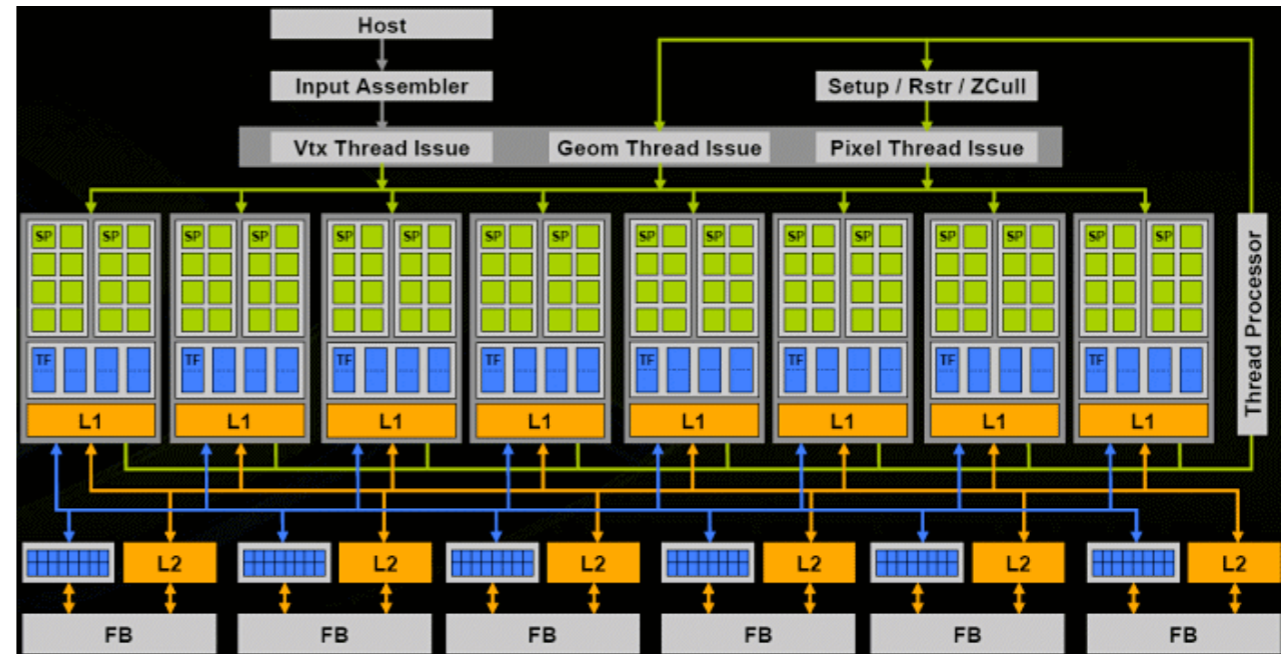


MultiCore/GPU/Cell: threaded programming



Multi-core processors: OpenMP (or just MPI)

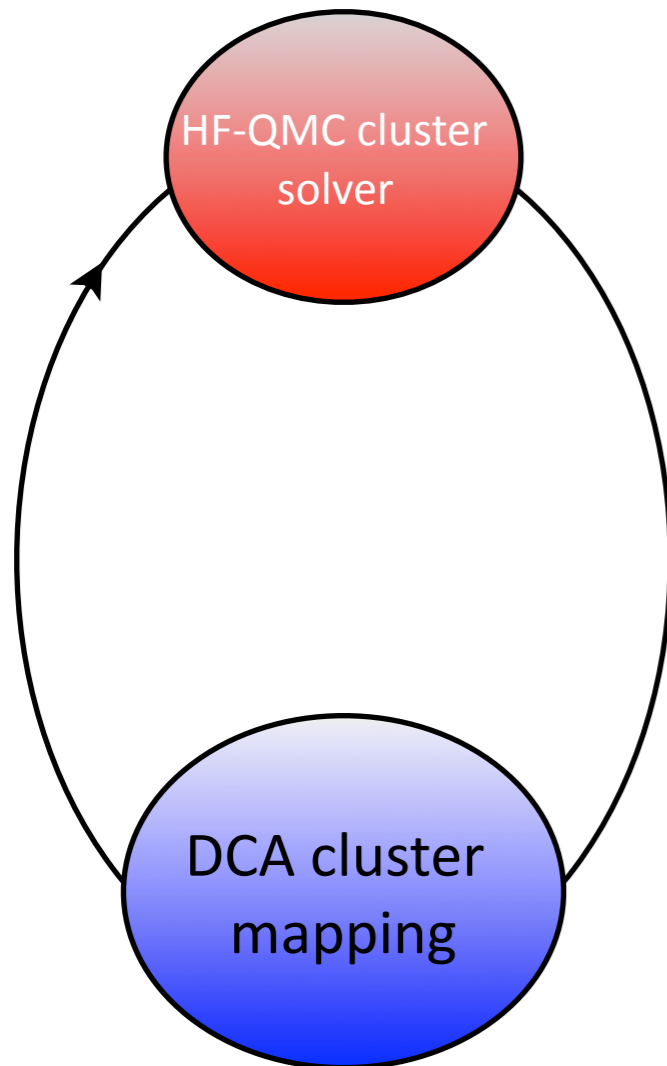
NVIDIA G80 GPU: CUDA, cuBLAS



IBM Cell BE: SIMD, threaded prog.

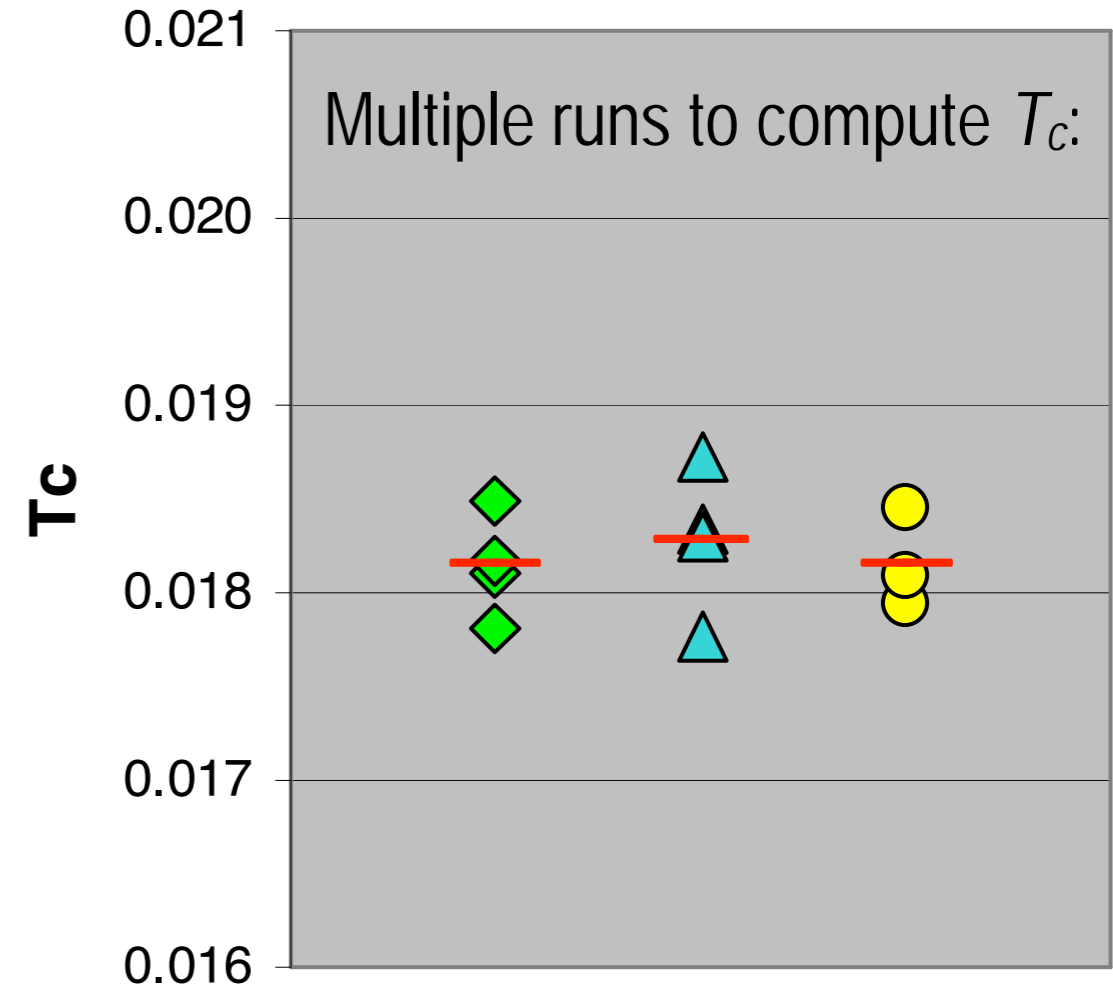
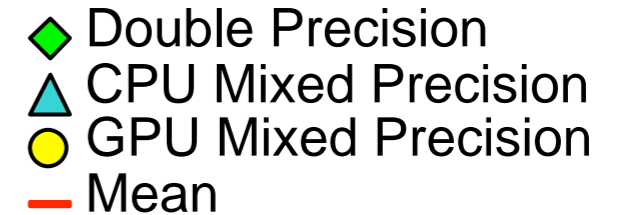
DCA++ with mixed precision

Run HF-QMC in single precision



Results for mixed and double precision runs are identical for same random number sequence!

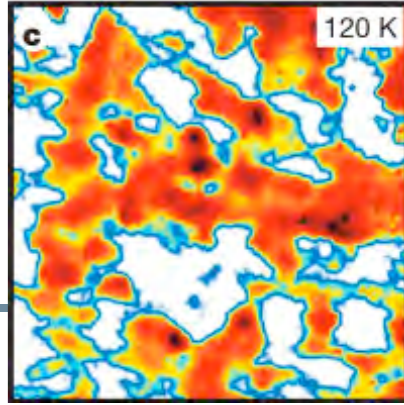
Keep the rest of the code, in particular cluster mapping in double precision



Speedup of HF-QMC updates (2GHz Opteron vs. NVIDIA 8800GTS GPU):

- 9x for offloading BLAS to GPU & transferring all data
- 13x for offloading BLAS to GPU & lazy data transfer
- 19x for full offload HF-updates & full lazy data transfer

Disorder and inhomogeneities



Hubbard Model with random disorder (eg. in U)

$$H^{(\nu)} = -t \sum_{\langle ij \rangle, \sigma} c_{i\sigma}^\dagger c_{j\sigma} + \sum_i U_i^{(\nu)} n_{i\uparrow} n_{i\downarrow}$$

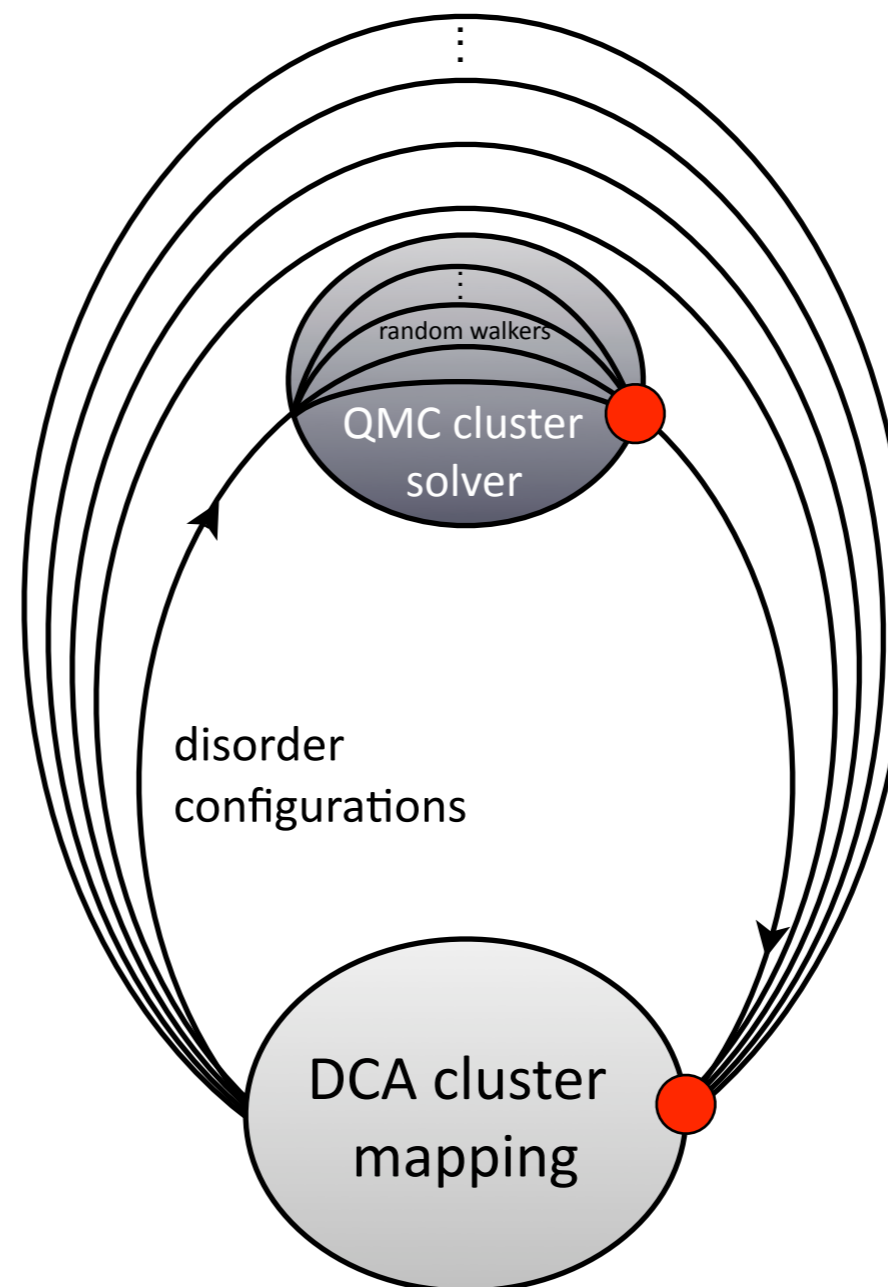
$$U_i^{(\nu)} \in \{U, 0\}; N_c = 16 \rightarrow N_d = 2^{16}$$

... need to disorder-average cluster Green function

$$G_c(X_i - X_j, z) = \frac{1}{N_c} \sum_{\nu=1}^{N_d} G_c^\nu(X_i, X_j, z)$$

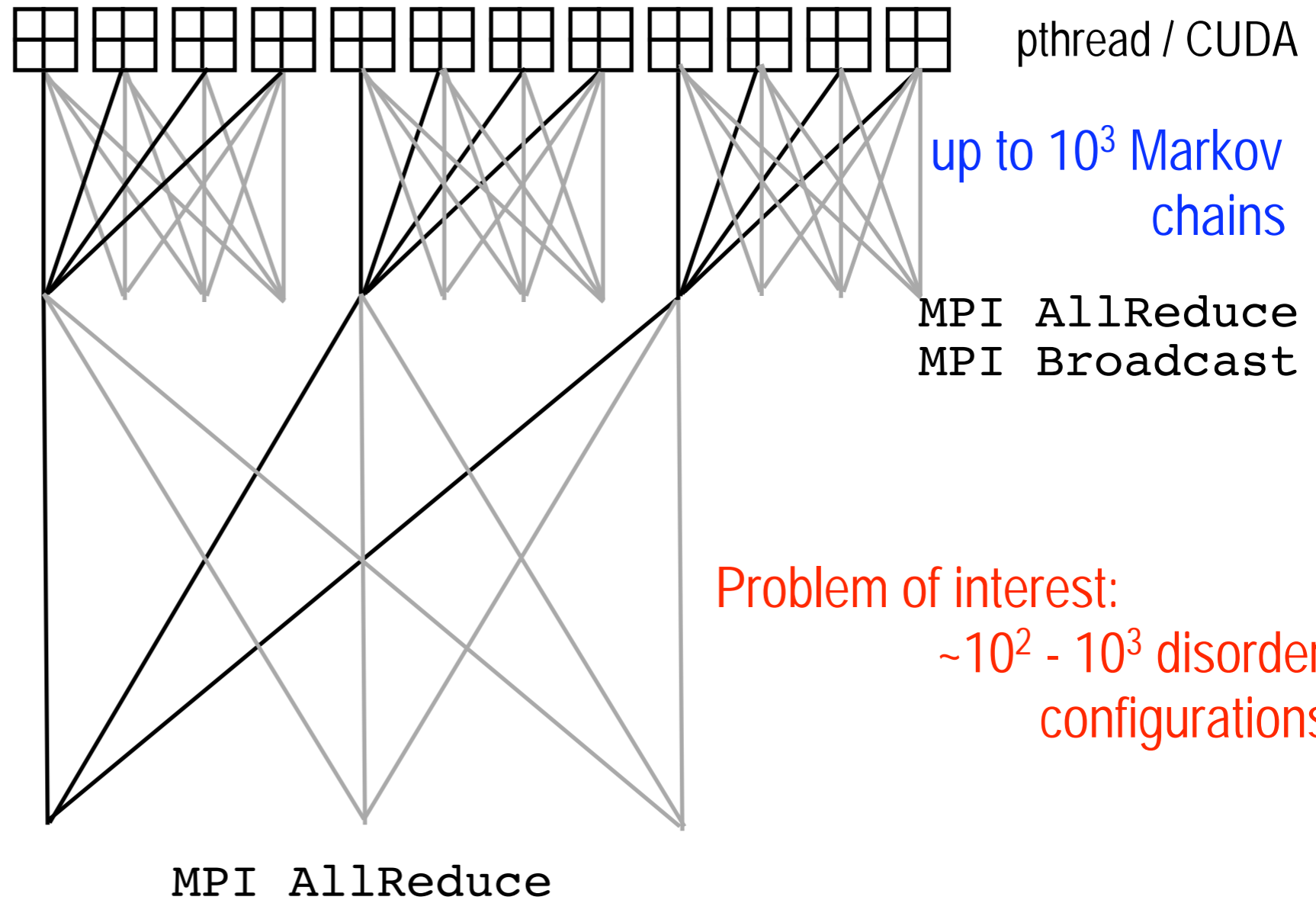
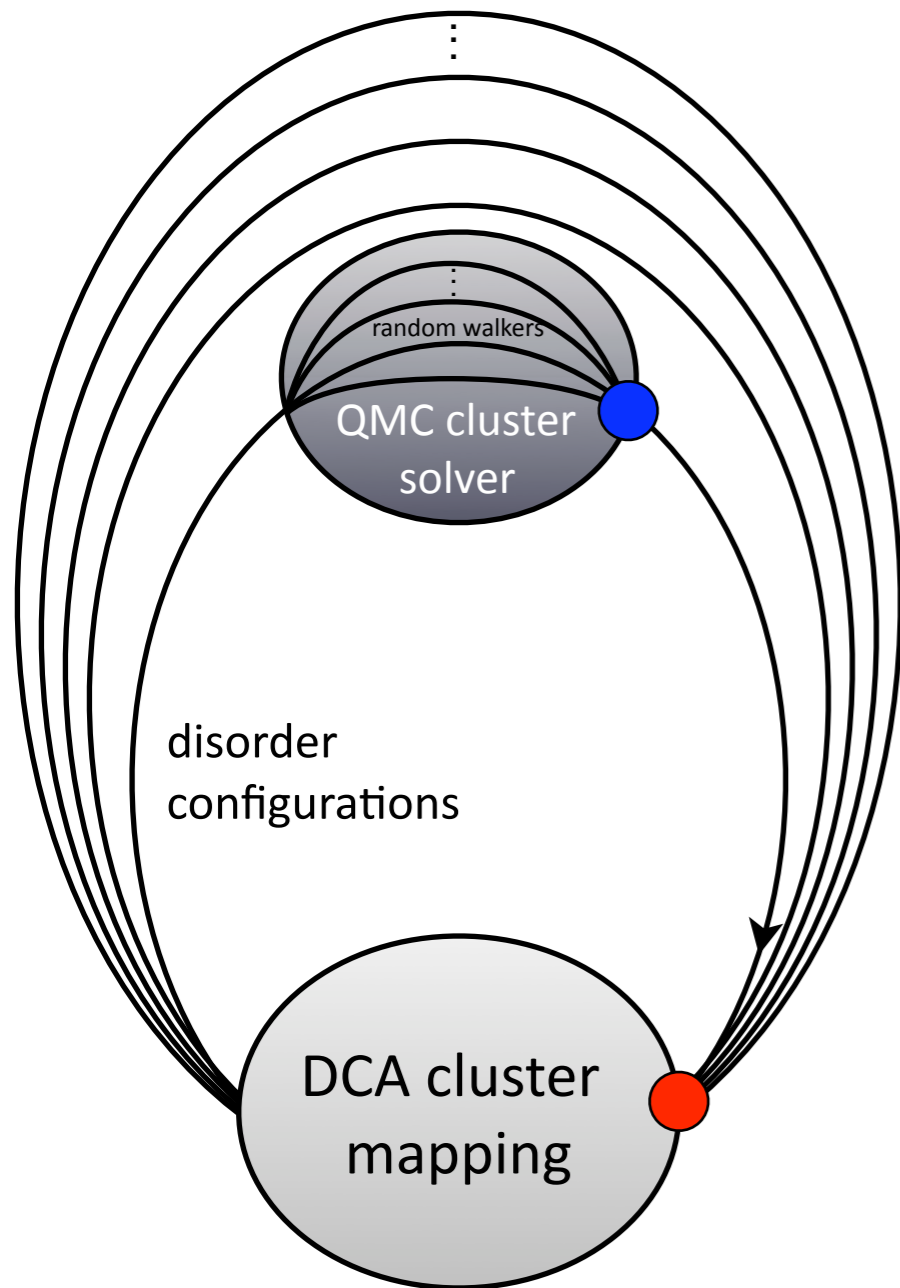
Algorithm 1 DCA/QMC Algorithm with QMC cluster solver (lines 5-10), disorder averaging (lines 4, 11-12), and DCA cluster mapping (line 3, 13)

- 1: Set initial self-energy
 - 2: **repeat**
 - 3: Compute the coarse-grained Green Function
 - 4: **for** Every disorder configuration (in parallel) **do**
 - 5: Perform warm-up steps
 - 6: **for** Every Markov chain (in parallel) **do**
 - 7: Update auxiliary fields
 - 8: Measure Green Function and observables
 - 9: **end for**
 - 10: Accumulate measurements over Markov chains
 - 11: **end for**
 - 12: Accumulate measurements over disorder configurations.
 - 13: Re-compute the self-energy
 - 14: **until** self consistency is reached
-

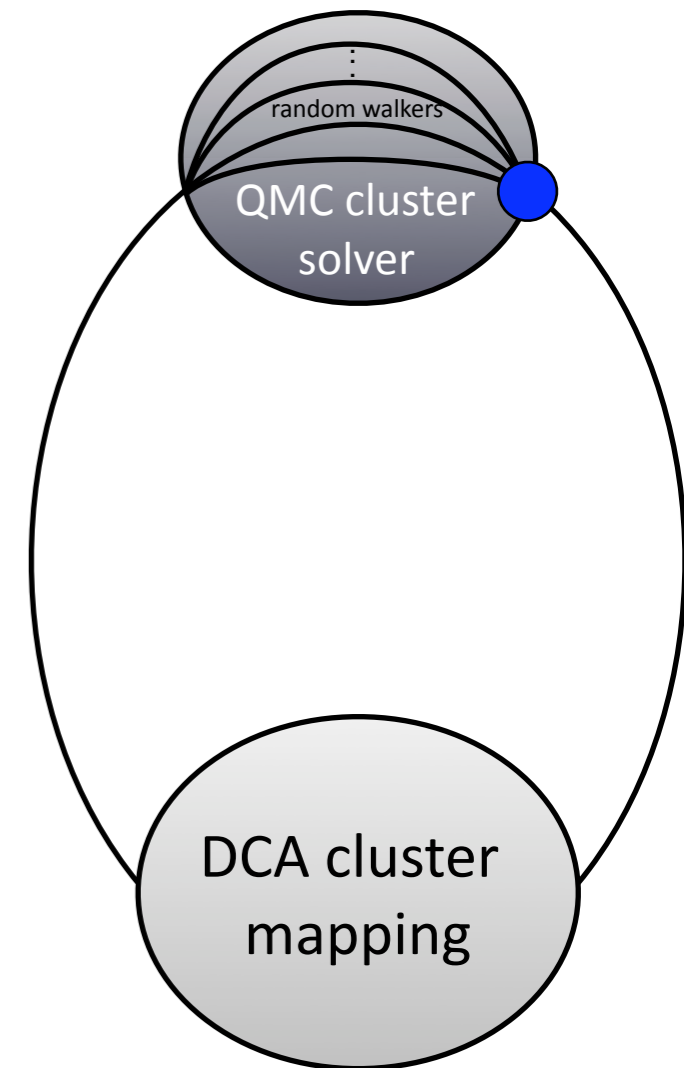
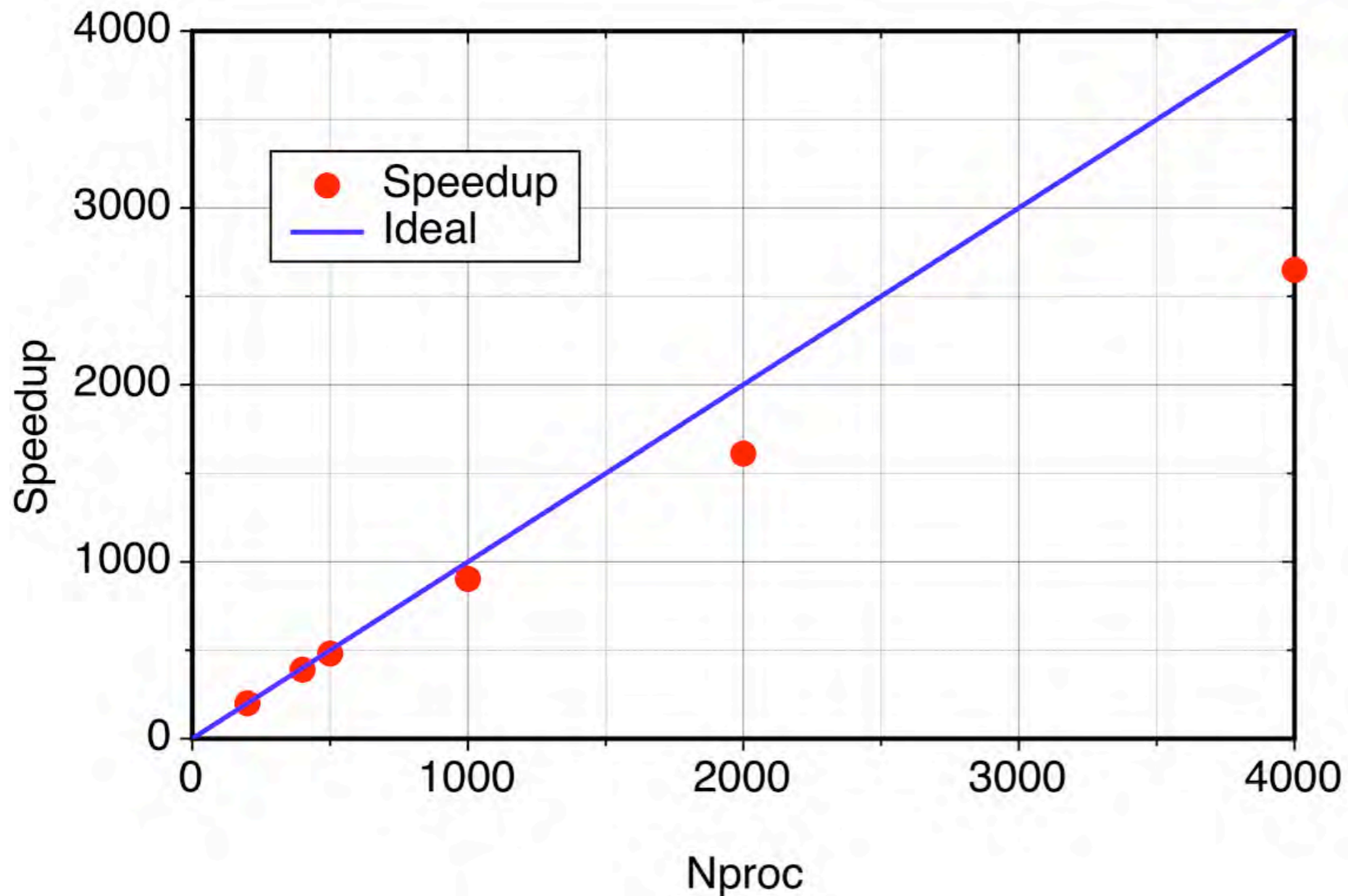
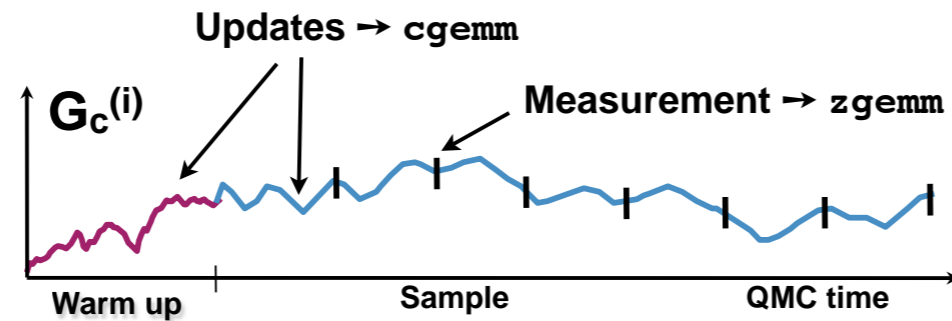


● required communication

DCA++ code from a concurrency point of view

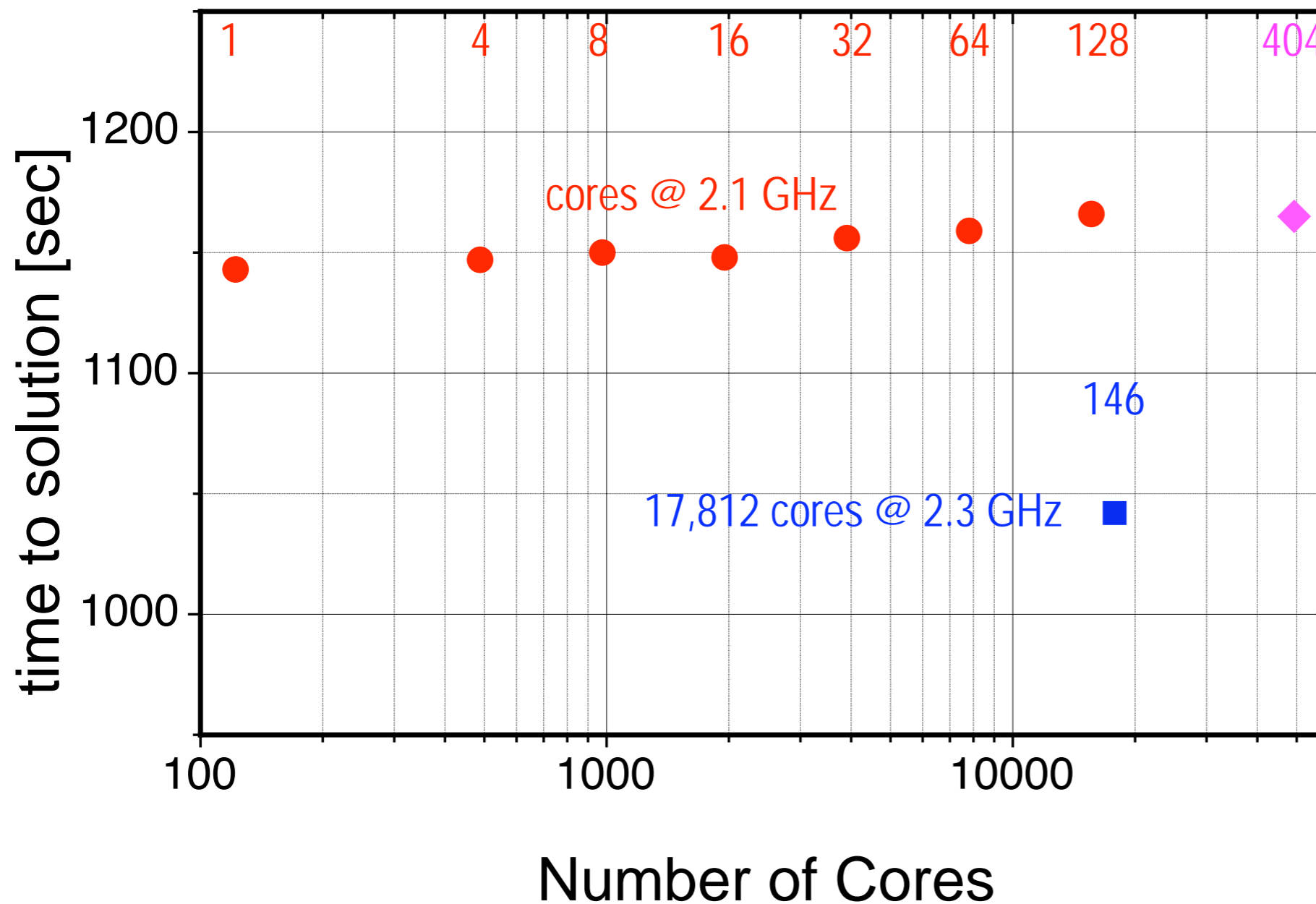
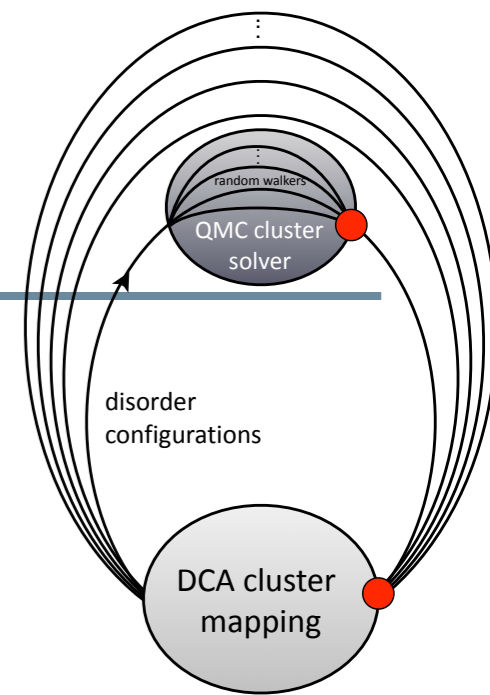


DCA++: strong scaling on HF-QMC



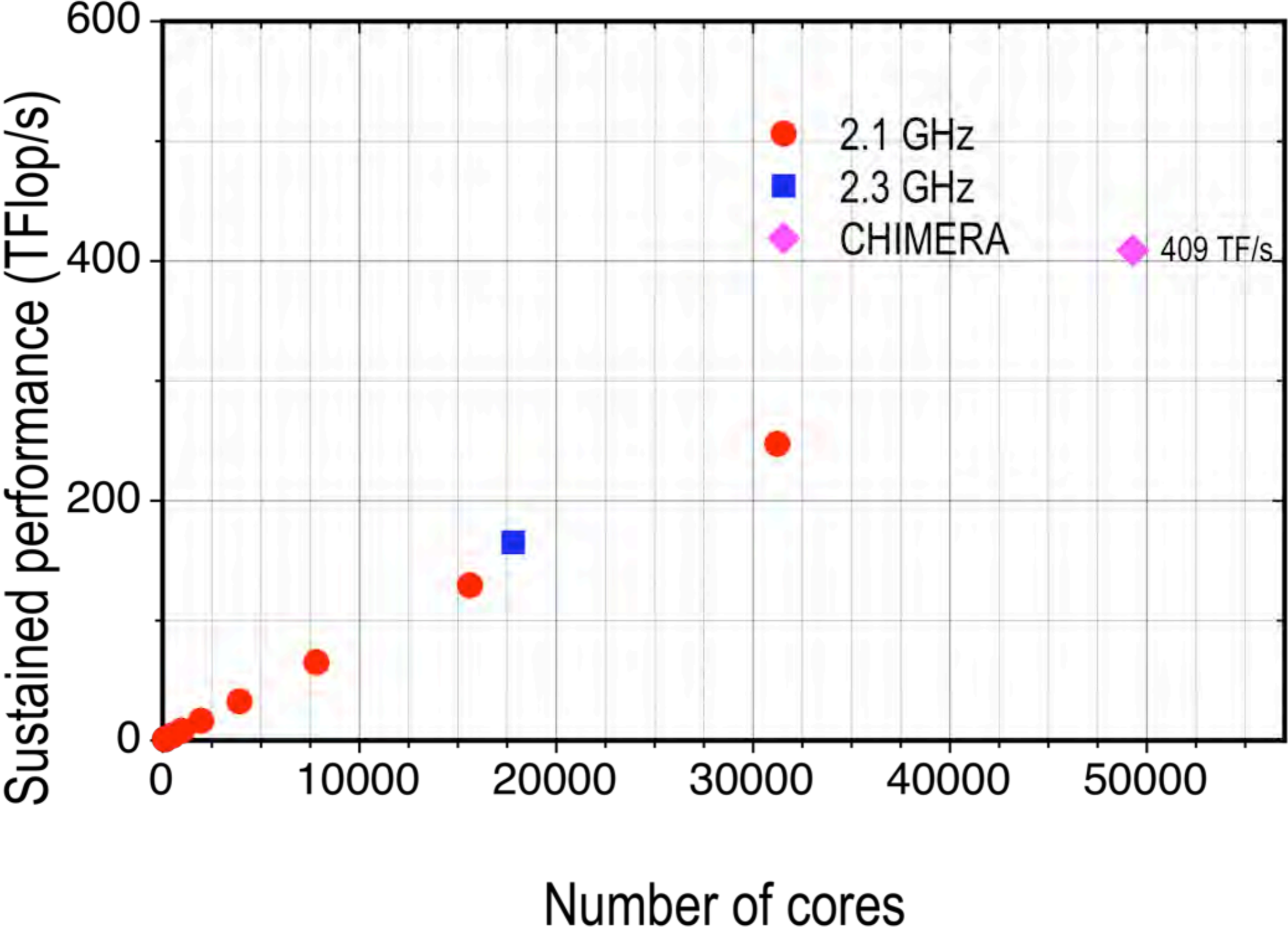
Weak scaling on Cray XT4

- HF-QMC: 122 Markov chains on 122 cores
- Weak scaling over disorder configurations



31,232 cores @ 2.1 GHz +
17,812 cores @ 2.3 GHz =
49,044-core chimera

Sustained performance of DCA++ on Cray XT4



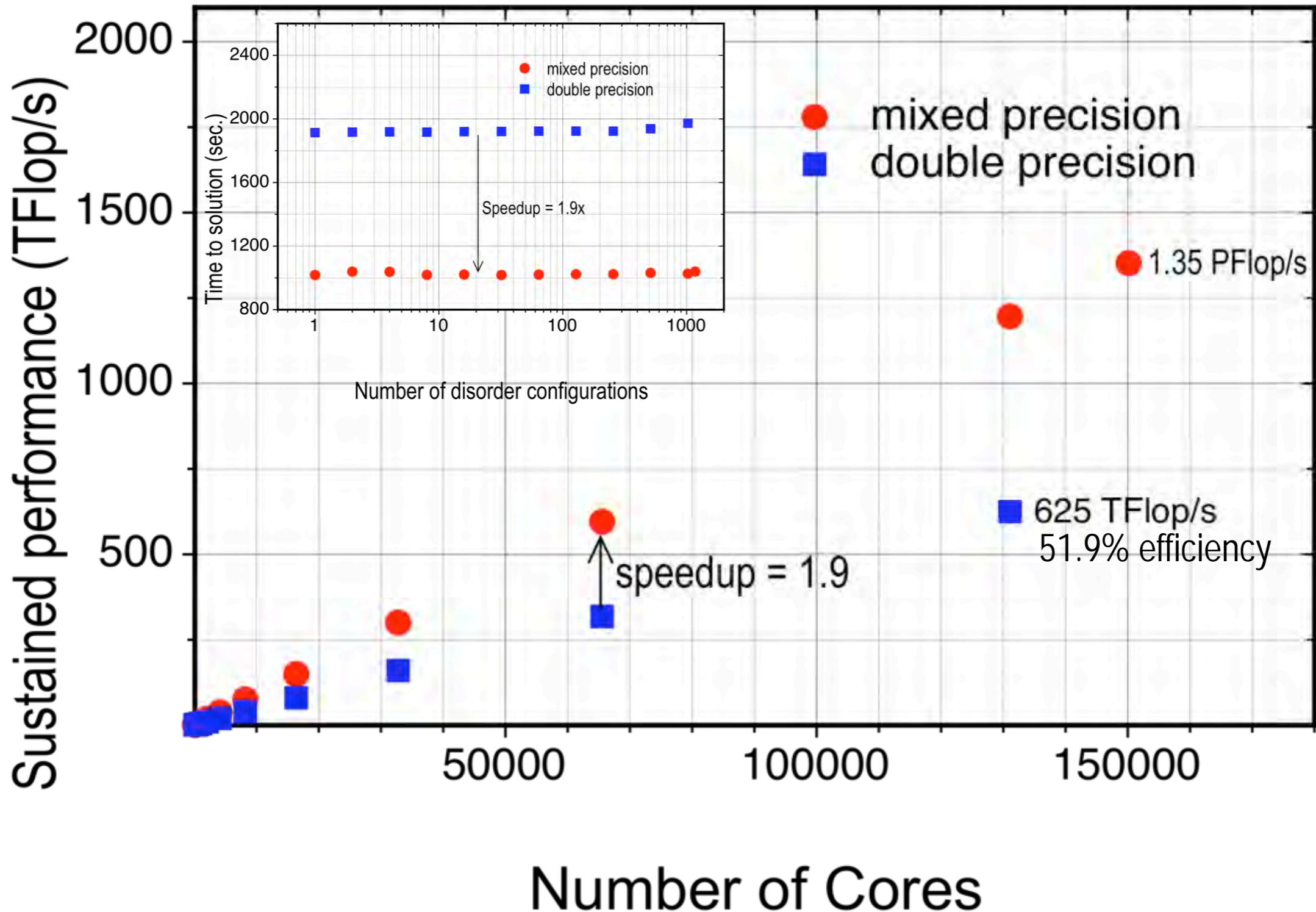
Cray XT5 portion of Jaguar @ NCCS



Peak: 1.382 TF/s
Quad-Core AMD
Freq.: 2.3 GHz
150,176 cores
Memory: 300 TB
For more details, go to
www.nccs.gov

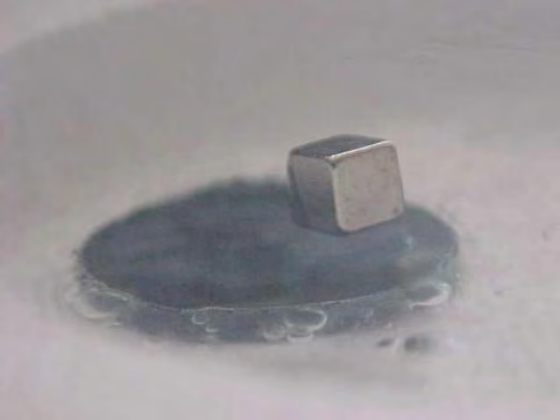
Sustained performance of DCA++ on Cray XT5

Weak scaling with number disorder configurations, each running on 128 Markov chains on 128 cores (16 nodes) - 16 site cluster and 150 time slides



Summary

- Today's methods and computational capabilities allow us to take a deep look into the mechanisms of high- T_c superconductivity
 - Simulations of superconducting transition in model without phonons
 - Dominant contribution to pairing mechanism: "glue" due to spin fluctuations
- DCA++ - optimally mapping DCA/QMC method onto today's hardware architectures
 - Algorithm: Hirsch-Fye QMC with delayed updates (>10x speedup)
 - Accelerator work motivated: mixed precision (almost 2x speedup)
 - Highly scalable implementation to study disorder and nanoscale inhomogeneities
 - Extensible implementation based on C++/STL generic programming model
- Sustained 1.35 PF/s on 150K cores of Cray XT5 portion of NCCS/Jaguar
 - Sustained 625 TF/s on 130K cores in double precision (52% efficiency)
- More than 1000 fold capability enhancement since 2004:
 - NCCS 2004: Cray X1 with 5 TF/s peak, DCA/QMC sustained about 2 TF/s (required high memory bandwidth)
 - NCCS 2008: factor 300 more in peak Flop/s & at least 20x due to algorithms
 - Future: Continuous time QMC - a new class of QMC algorithms



The DCA++ Story:

How new algorithms, new computers, and innovative software design allow us to solve real simulation problems of high high temperature superconductivity

Team, collaborators, computing resources, and funding



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