

Speedup of the Quantum Adiabatic Algorithm by Topological Cancellation

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Quantum Adiabatic Algorithm (QAA)

- Unique among possible quantum algorithms because it is “general purpose” – most optimization problems can be formulated to use it (also QAA is technically universal)
- Devices meant to implement it have been constructed (D-Wave), so it may be the first algorithm to run on an actual quantum computer
- Still not clear whether there are any problems for which the QAA is more efficient than various classical optimization algorithms

Outline

- A Very Short Introduction to the Quantum Adiabatic Algorithm (QAA)
- Topology of the QAA
- Topological Cancellation
- Improved QAA: Single-Spin Approximation
- Improved QAA: Cluster Approximation
- Conclusions

Basics of the QAA

A quantum computer evolves according to the time-dependent Hamiltonian

$$H_{qaa} = \left(1 - \frac{t}{t_D}\right) H_i + \frac{t}{t_D} H_f$$

Here H_i is any sufficiently simple Hamiltonian, e.g., : $H_i = h_0 \sum_{n=1}^L \sigma_x^{(n)}$

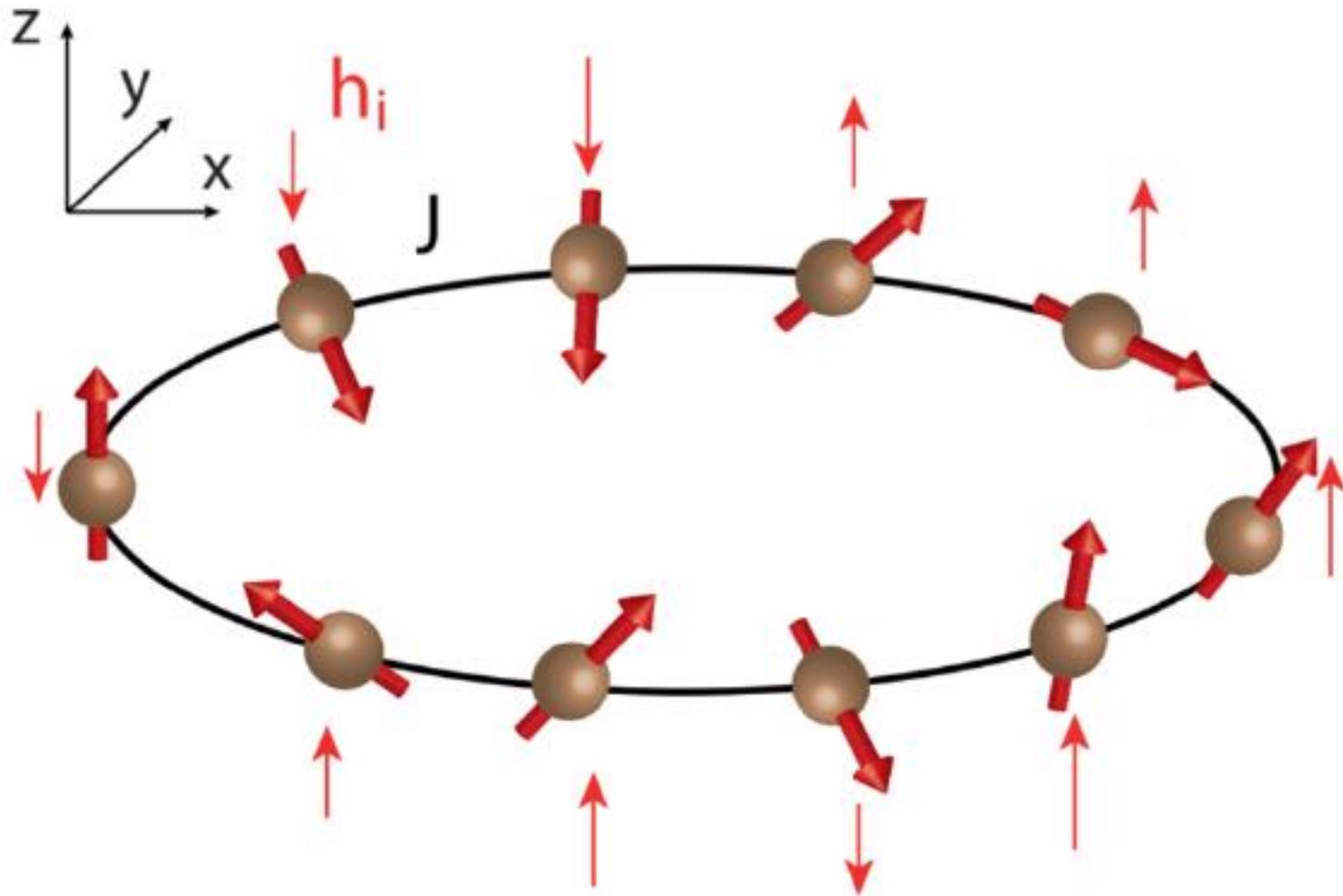
ground state $|\Psi_i\rangle = 2^{-L/2} (|0\rangle - |1\rangle)_1 (|0\rangle - |1\rangle)_2 \cdots (|0\rangle - |1\rangle)_L$

and H_f is a complicated Hamiltonian whose ground state encodes the solution to the optimization problem, e.g.,

$$H_f = \sum_{n=1}^L h_n \sigma_z^{(n)} + J \sum_{n=1}^L \sigma_z^{(n)} \sigma_z^{(n+1)}, h_n \in [-1, 1]$$

Specification of $\{h_n\}$ determines an instance (or a realization) of the optimization problem.

Random-field Ising Model with t-dependent transverse field



Variable Number L of spins on the Ring

Adiabaticity

- If t_D is large enough, then the adiabatic theorem tells that $|\Psi_i\rangle$ will evolve into $|\Psi_f\rangle$ at $t = t_D$, and that $|\Psi_f\rangle$ is the ground state of H_f , thus solving the optimization problem.

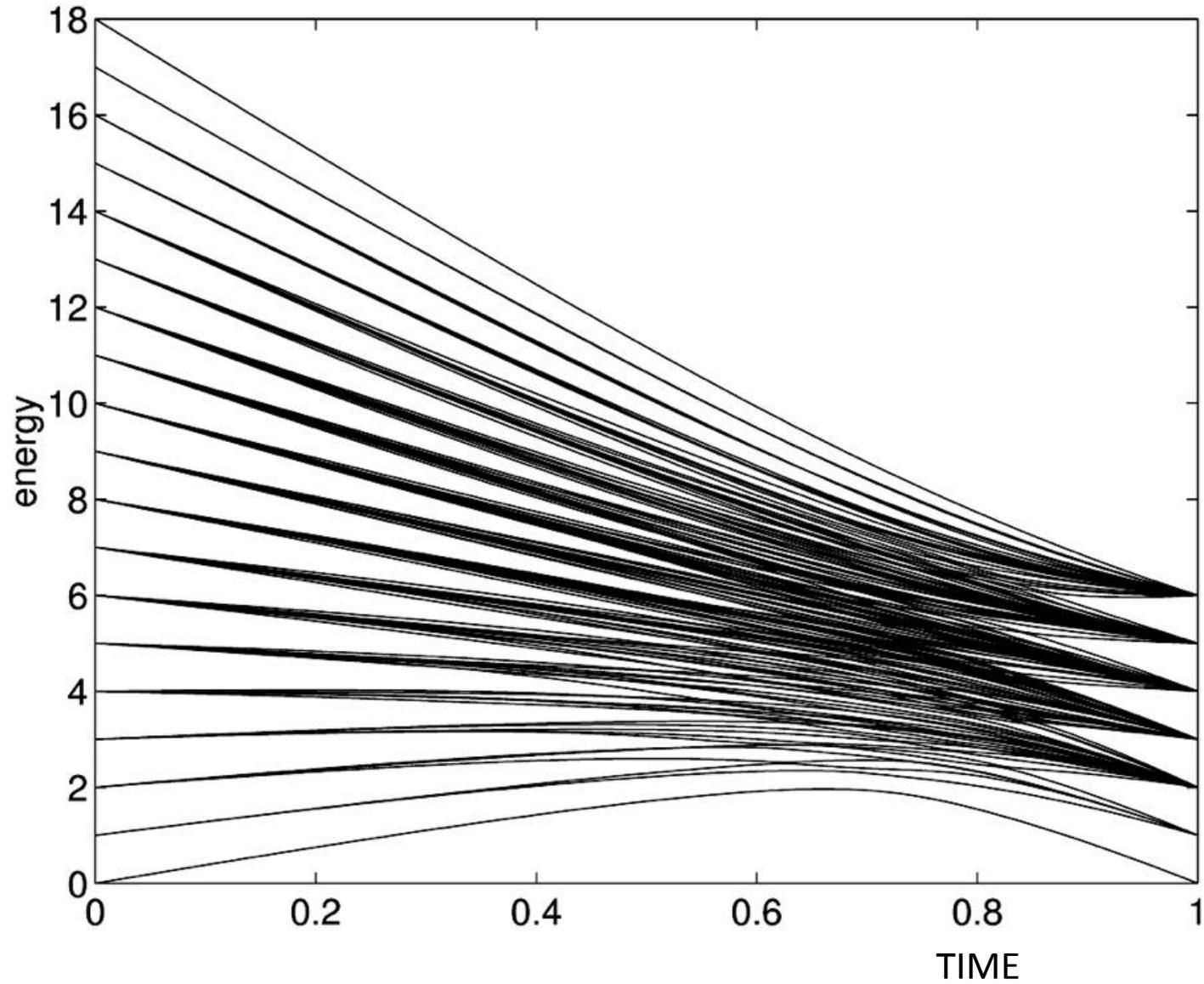
How large is large enough for t_D ? Simple arguments (see, e.g., Messiah, 1961) say that if we define the minimum gap as:

$$\Delta_{\min} = \min_{0 \leq t \leq t_D} [E_1(t) - E_0(t)]$$

Then $t_D \gg C \left\| \left(\frac{dH}{dt} \right)^2 \right\| / \Delta_{\min}^2$

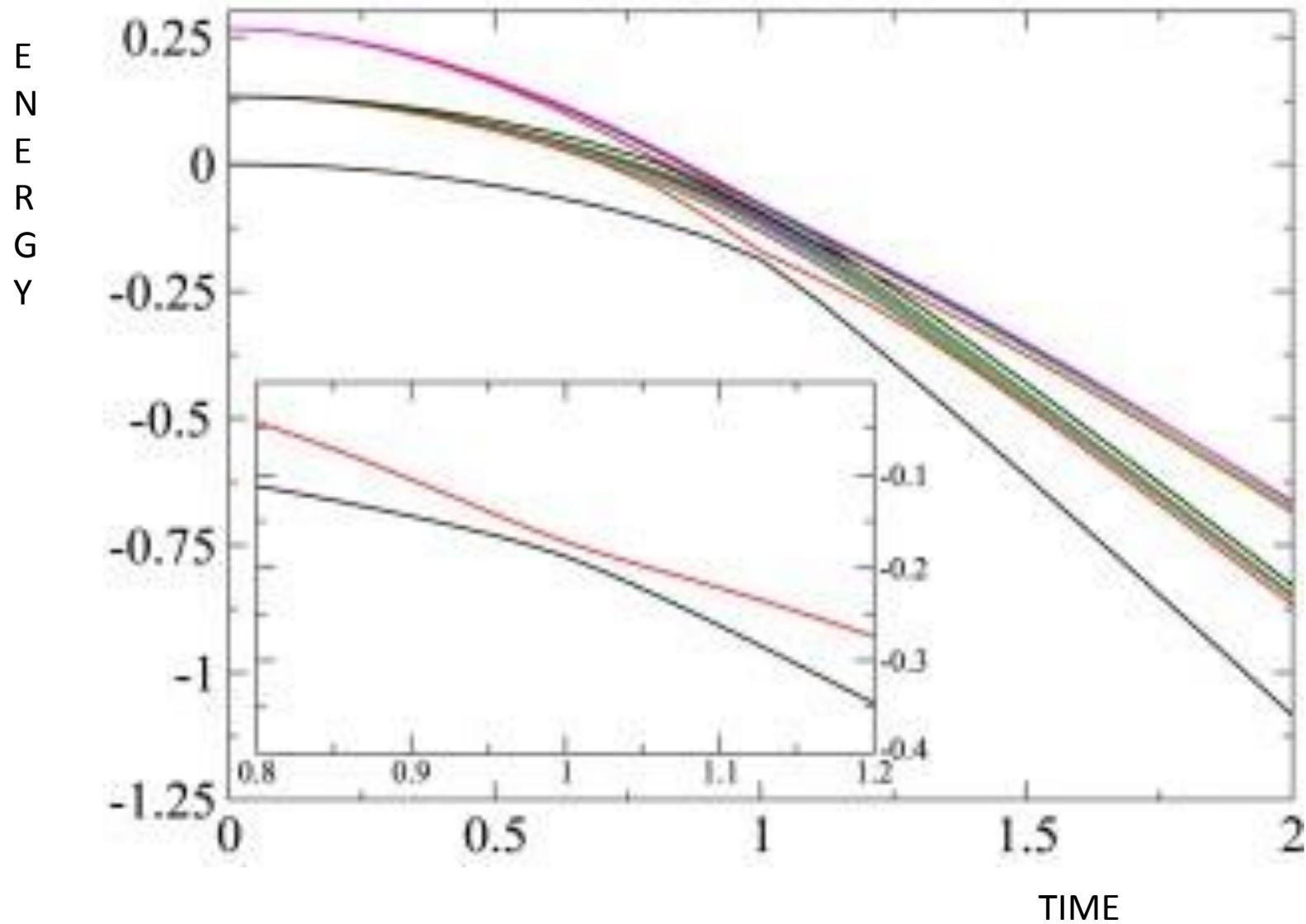
If $\Delta_{\min} = O(L^{-n})$ the algorithm is efficient but if $\Delta_{\min} = O(\exp(-L))$ then the algorithm is inefficient

GOOD!



AM Childs talk

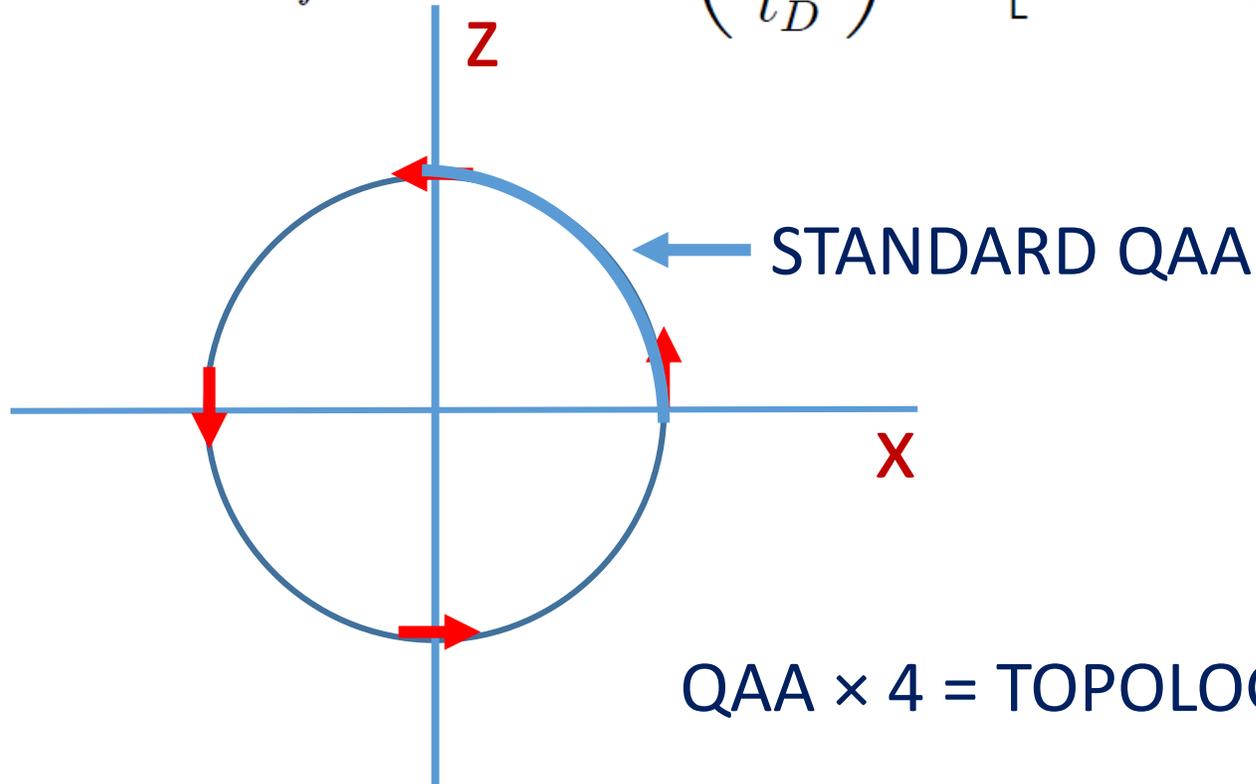
BAD!



Topological Picture of the QAA

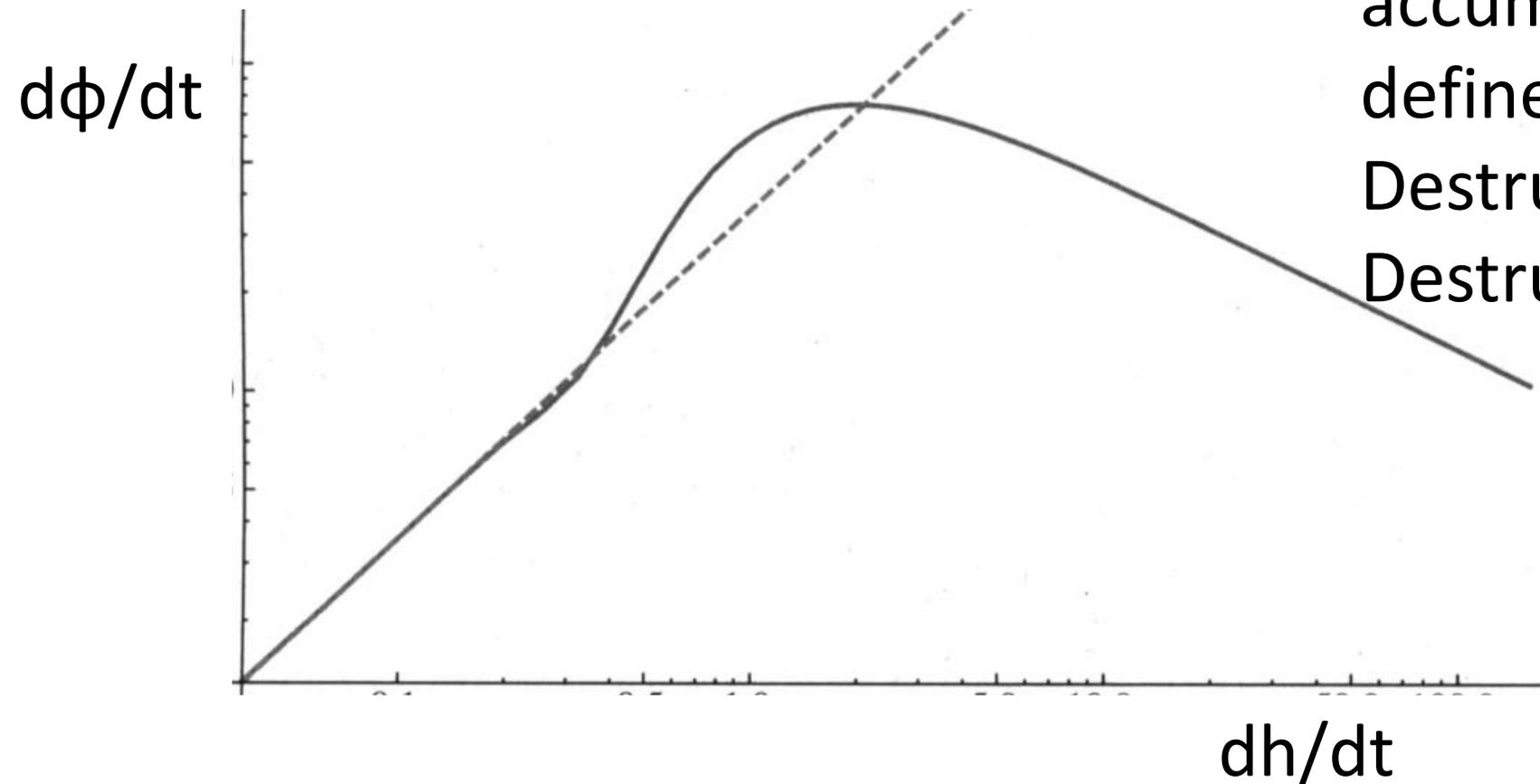
- Consider an *expanded* time evolution: $H_{topo}(t) = \cos\left(\frac{2\pi t}{t_D}\right) H_i + \sin\left(\frac{2\pi t}{t_D}\right) H_j$
- And an effective mean field on the n th spin:

$$h_{mf}^{(n)}(t) = h_0 \cos\left(\frac{2\pi t}{t_D}\right) \hat{x} + \left[h_n + J \left(\sigma_z^{(n+1)}(t) + \sigma_z^{(n-1)}(t) \right) \hat{z} \right] \sin\left(\frac{2\pi t}{t_D}\right)$$



Perfect adiabatic evolution gives $M_y=0$ and Berry phase = π .
Deviations from π indicate breakdown of the algorithm.

Breakdown of Adiabatic Berry Phase Accumulation Spin $\frac{1}{2}$ Particle in a t-dependent B-field.



Sharp Breakdown of adiabatic accumulation at a fairly well-defined driving speed.
Destruction of Topology =
Destruction of QAA!

Gritsev and
Polkovnikov,
PNAS **109**, 6457
(2011)

Topological Cancellation

Define a time-dependent Hamiltonian $H_0(t)$ with instantaneous eigenstates $|n(t)\rangle$:

$$H_0(t) |n(t)\rangle = E_n(t) |n(t)\rangle$$

Then there will be transitions between the $|n(t)\rangle$. But if we define $H = H_0(t) + H_1(t)$, with

$$H_1(t) = i \sum_{m \neq n} \frac{|m\rangle \langle m | \partial_t H_0 | n \rangle \langle n|}{(E_n - E_m)}$$

BERRY CONNECTION TERM
ALSO "COUNTER-DIABATIC"
TERM !

Then if the initial state is $|n(t)\rangle$, the final state is also $|n(t)\rangle$.

For a single spin these equations reduce to

$$H_0(t) = -\vec{B}(t) \cdot \vec{\sigma} / 2 \quad \text{and} \quad H_1(t) = \frac{\vec{B} \times \partial_t \vec{B}}{2B^2} \cdot \vec{\sigma}$$

Application to the QAA

We propose to add the topological term to the Hamiltonian that will suppress the unwanted transitions:

$$H_{total}(t) = f_i\left(\frac{t}{t_D}\right) H_i + f_f\left(\frac{t}{t_D}\right) H_f + H_s(t)$$
$$f_i(0) = f_f(1) = 1; f_i(1) = f_f(0) = H_s(0) = H_s(1) = 0$$

H_s must vanish at both endpoints: at $t=0$ because it is not easy to diagonalize, at $t=t_D$ because it must not change the optimization problem.

H_s involves the instantaneous eigenstates, which we do not know how to calculate efficiently: must approximate!

Single-particle Approximation

Now we choose $f_i(t) = \cos^2\left(\frac{\pi t}{2t_D}\right)$; $f_f(t) = \sin^2\left(\frac{\pi t}{2t_D}\right)$

$$\text{and } H_s(t) = - \sum_n \frac{h_0 h_n \sin\left(\frac{\pi t}{2t_D}\right) \cos\left(\frac{\pi t}{2t_D}\right)}{2t_D \left[h_0^2 \sin^4\left(\frac{\pi t}{2t_D}\right) + h_n^2 \cos^4\left(\frac{\pi t}{2t_D}\right) \right]} \sigma_y^{(n)}$$

as our approximation to H_s .

It's necessary that the turn on and turn off are quadratic
– otherwise it is impossible to get $H_s(t=0) = H_s(t=t_D) = 0$

Non-interacting Spins: $J=0$

$$P_0 = |\langle 0 | \psi(t = t_D) \rangle|^2$$

is the success probability.

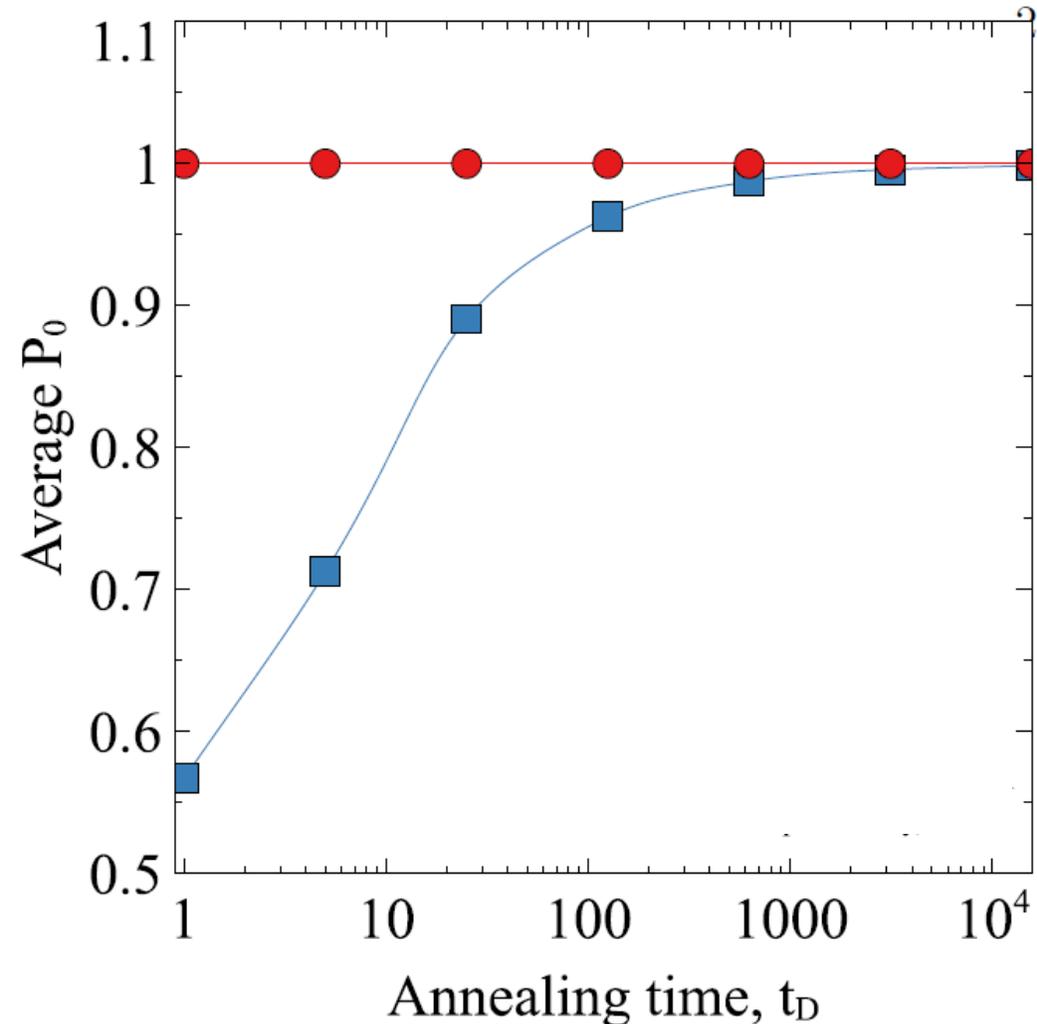
Random-field Ising model with
 L spins on a ring.

Red points have steering.

Blue points have no steering.

10^4 realizations of the disorder.

$L=1$, $h_0 = 10$



Weakly Interacting Spins: (J=0.1)

$$P_n = |\langle n | \psi(t = t_D) \rangle|^2$$

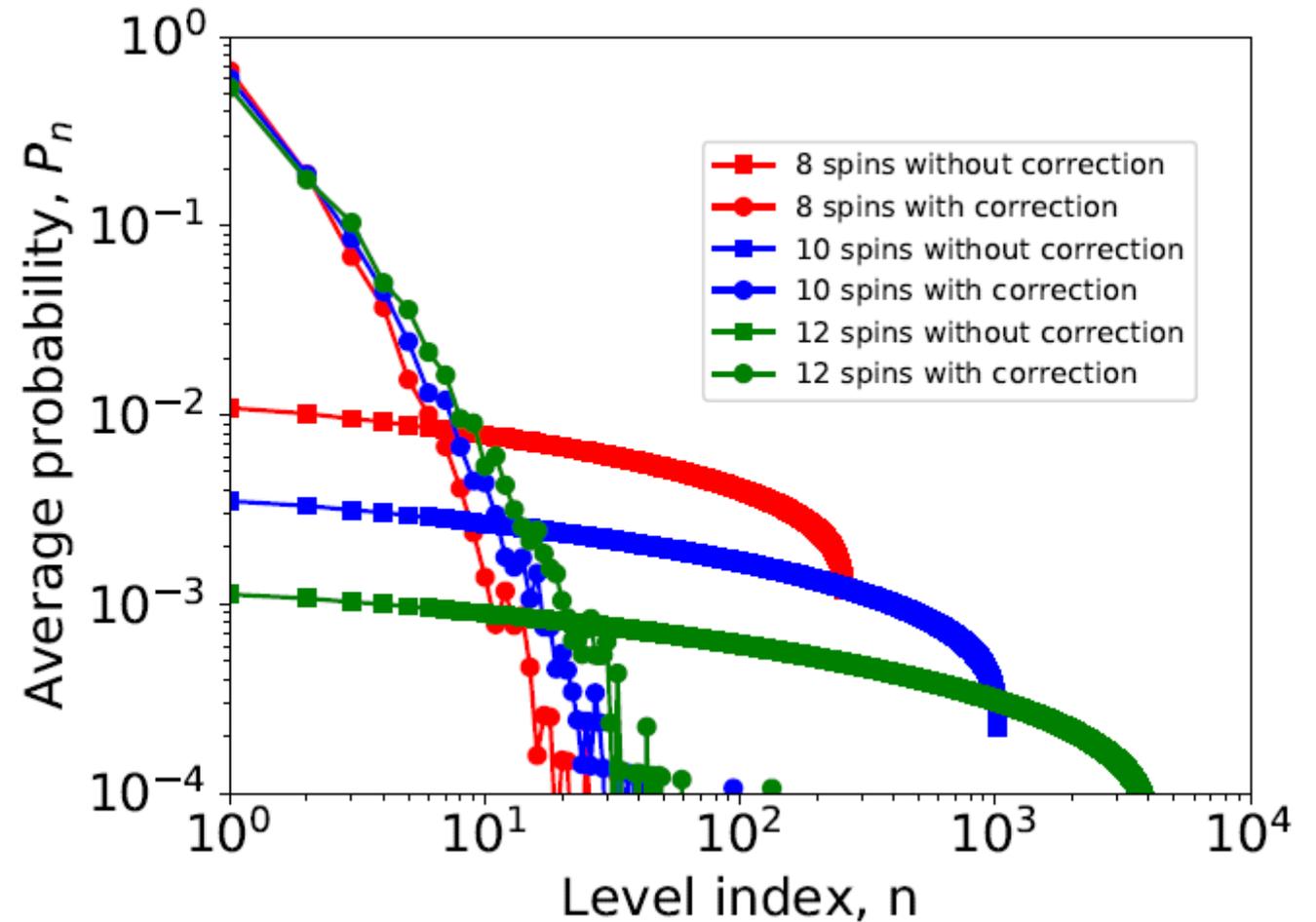
is the probability of ending in the state $|n\rangle$.

Round points are with steering

Square points are without

10^4 realizations of the disorder

$L = 8, 10, 12$; $t_D = 1$ (short!)



Interacting Spins: (all J)

$$P_0 = |\langle 0 | \psi(t = t_D) \rangle|^2$$

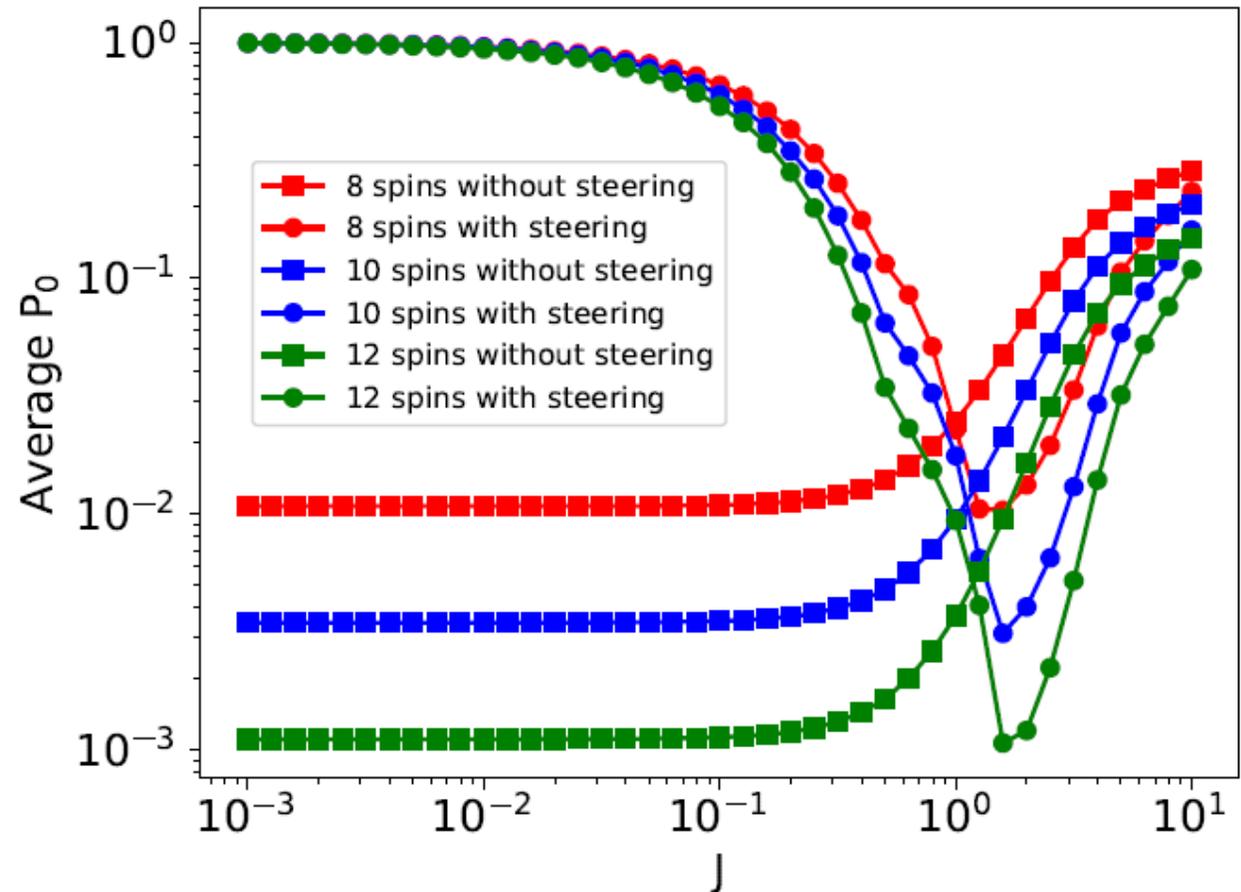
is the probability of ending in the state $|0\rangle$.

Round points are with steering

Square points are without

10^4 realizations of the disorder

$L = 8, 10, 12$; $t_D = 1$ (short!)



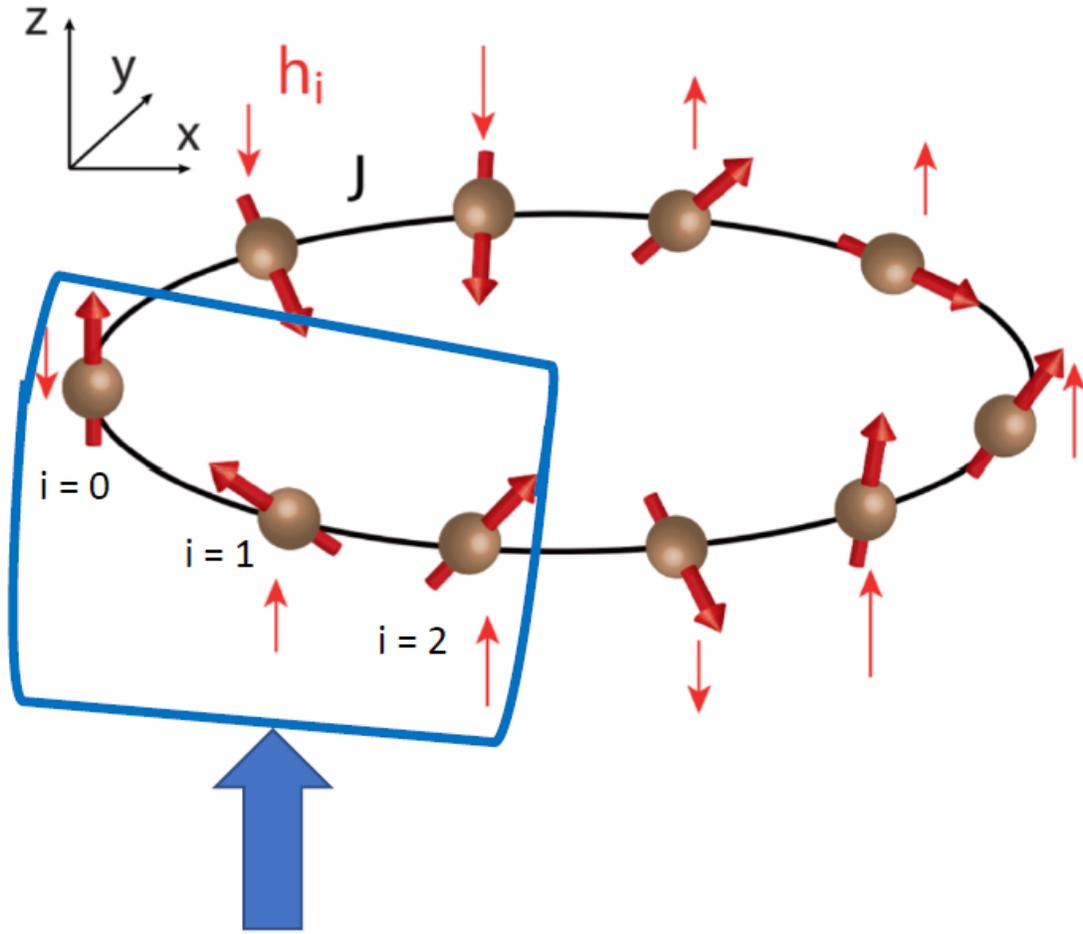
Single-spin Approximation: Remarks

- The algorithm is straightforward to implement.
- The calculation of the steering Hamiltonian is efficient.
- The speedups are very considerable for small J , say $J < 0.5$.
- At intermediate coupling $1 < J < 10$, single-spin steering is actually detrimental (but see below).
- At very large J , single-spin steering makes no difference.
- Can we improve?

Cluster Approximation

- We search for the spin with smallest h_n .
- The idea is that this spin is determined by the interaction and neighboring spins: it “feels” the interaction most strongly.
- This spin is chosen as the center of a cluster for which H_s is computed numerically using the exact many-spin Berry formula.
- The other spins are treated as before.
- This is the first step in a cluster expansion for the method.

Cluster approach



- For a spin with weak z -field, dynamics is governed by neighbors.
- Number spins so that h_1 is the smallest z -field.
 - Apply exact Berry steering to spins numbered as 0, 1 and 2.
- To the rest of the spins, apply 1-spin steering.

Interacting Spins: Cluster Approximation

Percentage of realizations with

$$P_0 = |\langle 0 | \psi(t = t_D) \rangle|^2 > 0.99 .$$

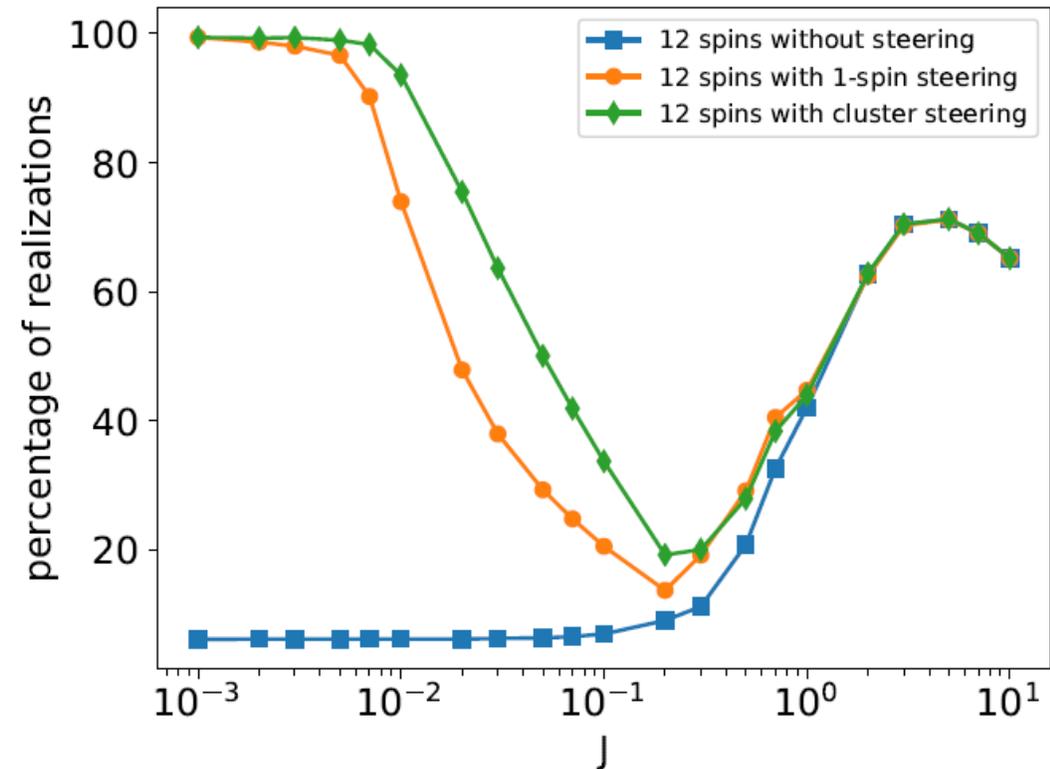
Diamond points are with cluster steering

Round points are with single-spin steering

Square points are without steering

10^4 realizations of the disorder

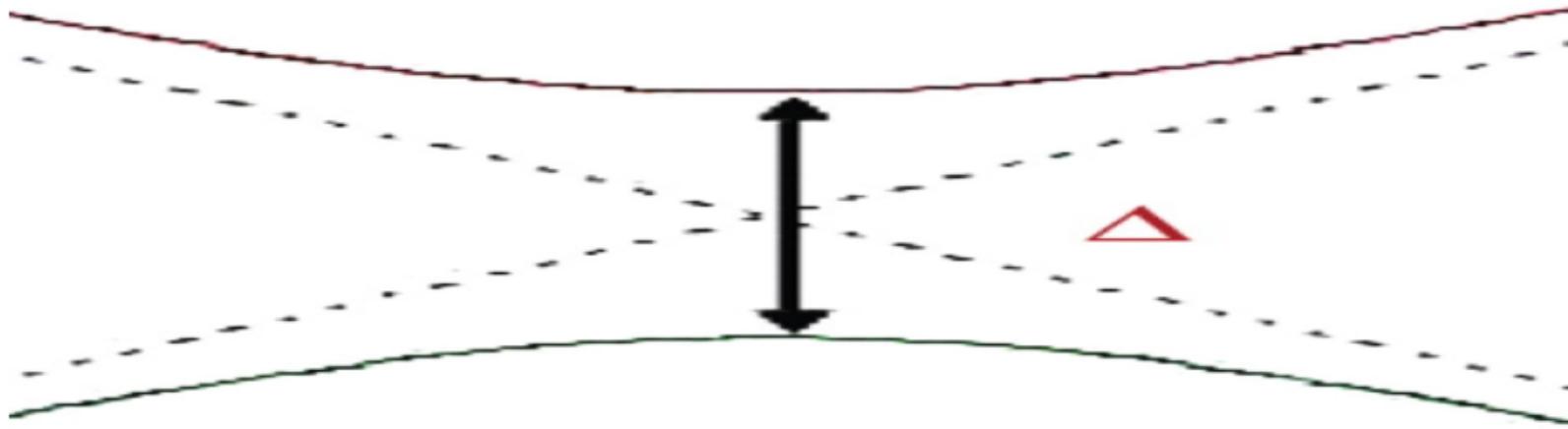
$L = 12$; $t_D = 128$



Conclusions, Discussion

- We add a topological cancellation term to the QAA Hamiltonian that suppresses single-spin transitions.
- The single-spin method shows significant improvements on the standard QAA for random spin systems with weak to moderate interaction strengths.
- The simplest cluster expansion of the cancellation term also gives a speedup in the range of moderate interactions – this suggests that systematic improvement is possible.
- H_i and H_f are stoquastic, but H_s is not.
- We have not investigated the scaling behavior of the method – its local character so far would suggest a constant speedup, but the method in general need not be local.

Landau-Zener-Majorana tunneling

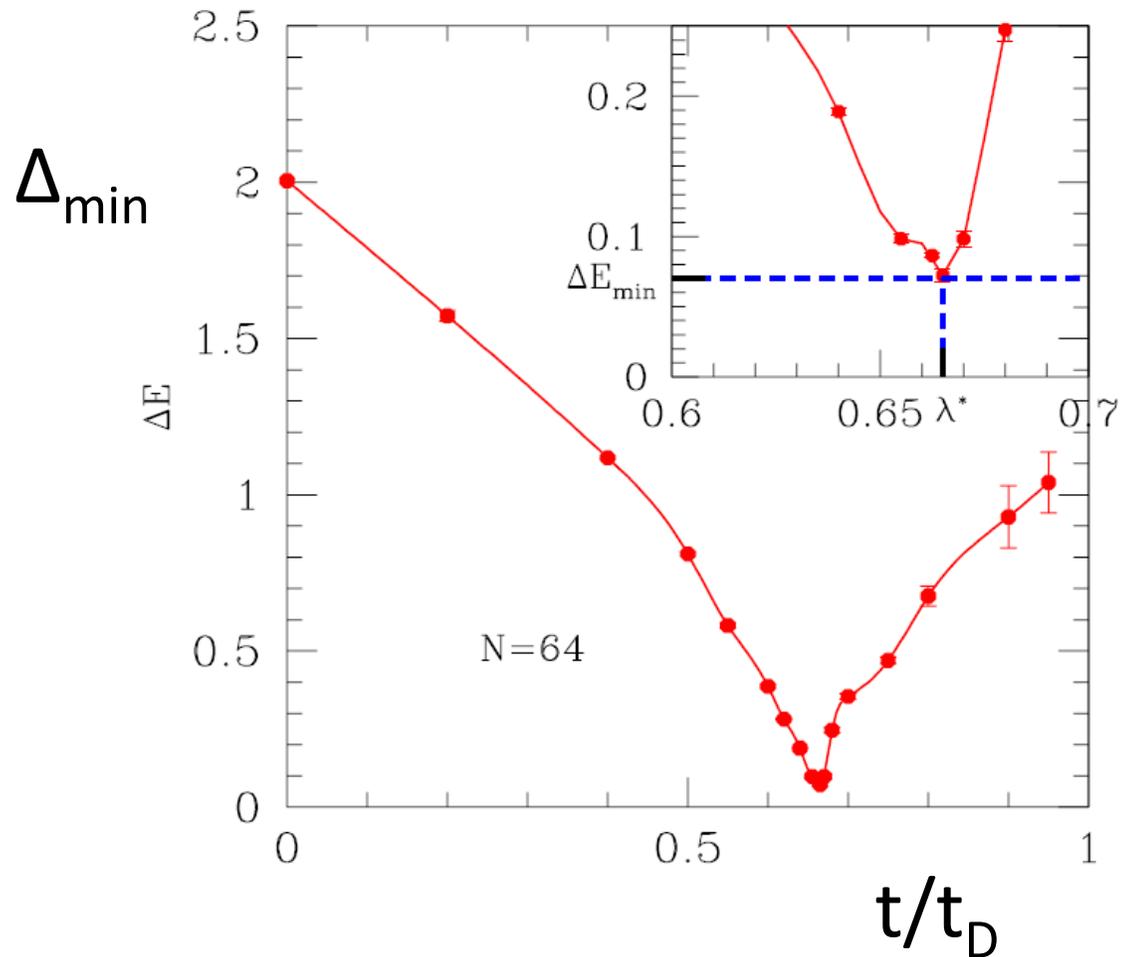


$$H_{LZ} = -\frac{vt}{2}\sigma_z - \frac{\Delta}{2}\sigma_x$$

Δ is minimum gap.

$$P_{transition} = e^{-\frac{\pi\Delta^2}{2v}}$$

LZM Tunneling *is* the Problem



$$t_D \gg \hbar \frac{|H_{10}(t = t_{\min})|}{\Delta_{\min}^2}$$

Seems to be a very accurate criterion in the problems known to date, indicating that individual avoided Level crossings are the main obstacle to the efficiency of the QAA.