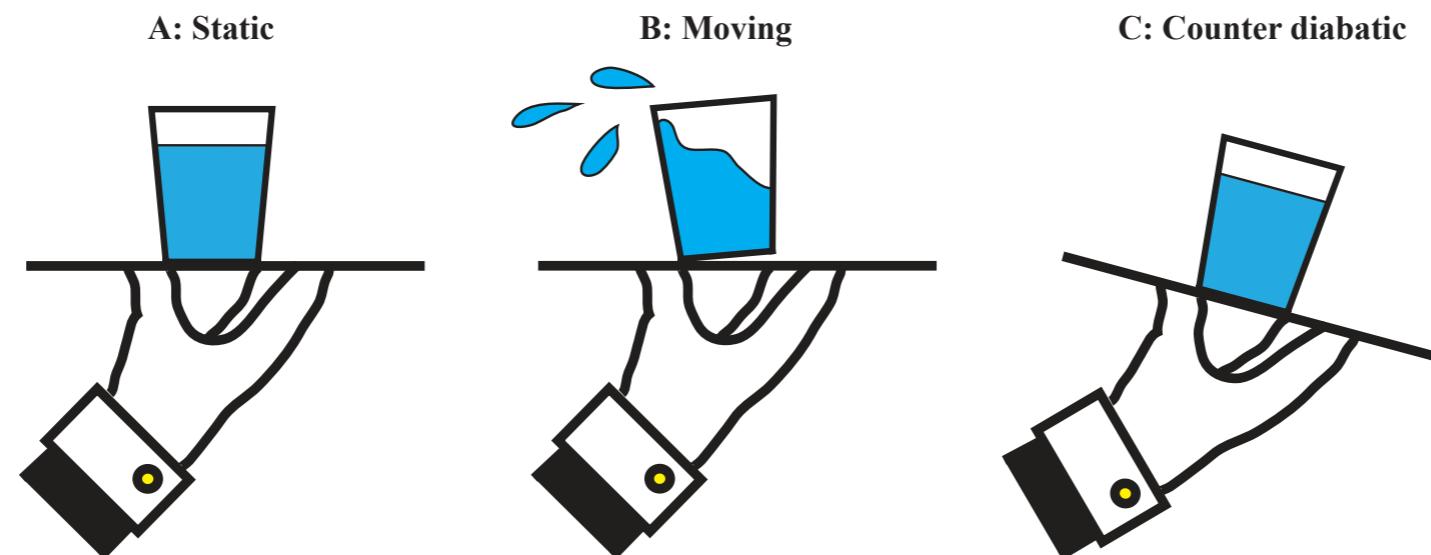


Counter-diabatic driving in quantum many body systems

Dries Sels & Anatoli Polkovnikov
Harvard-BU

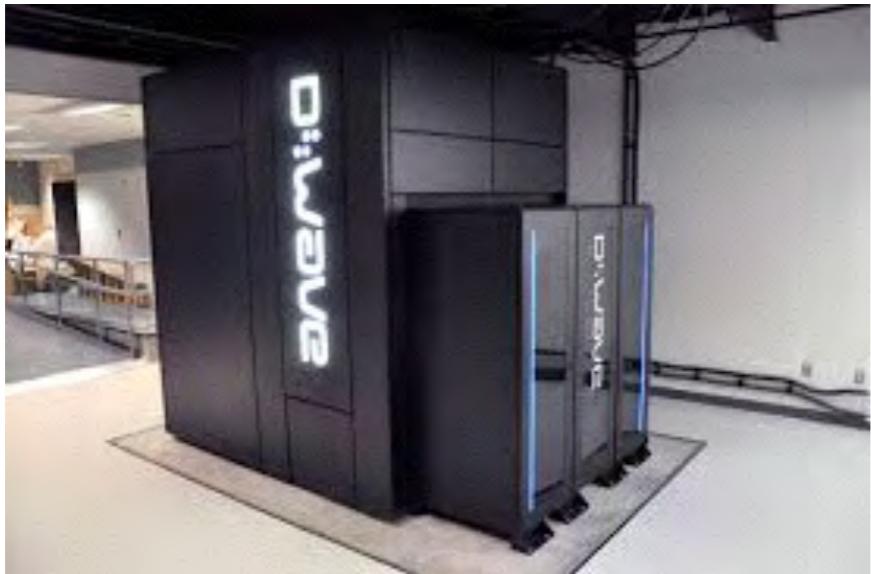


Outline

- Non-adiabatic response and quantum geometry
- Counter-diabatic driving and quantum speed limit
- Variational adiabatic gauge fields for many body systems

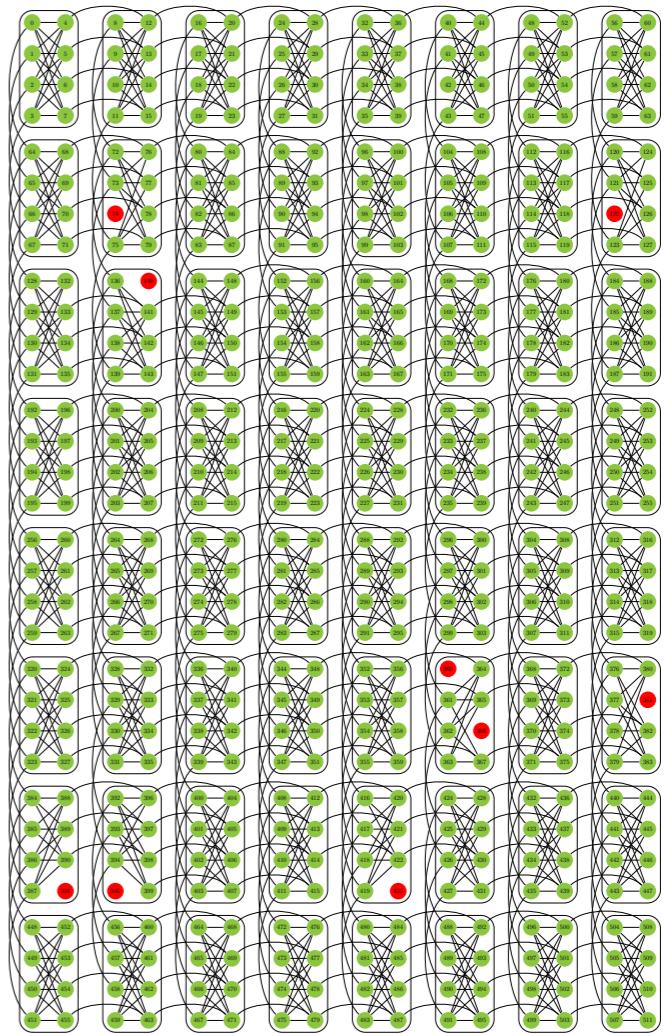
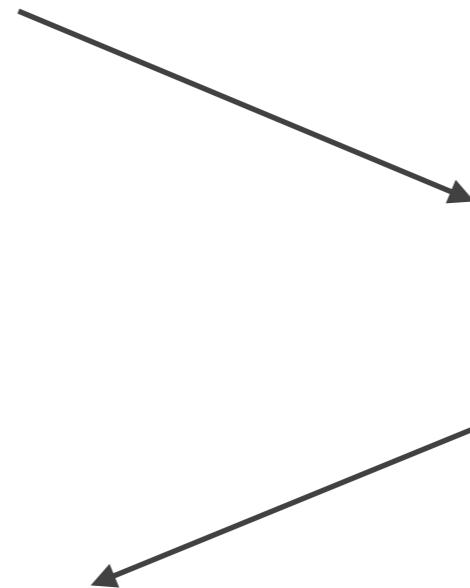
Motivation

- How to drive your system without exciting it?



Example : D-Wave machine

$$\sum_{ij} J_{ij} s_i^z s_j^z + \sum_i h_i(\tau) s_i^x$$



Non-adiabatic forces

- Consider a particle in an optical tweezer:

$$H = \frac{p^2}{2m} + V(x - \lambda(t))$$

- Let's move with the atom: $|\psi\rangle = U(\lambda) |\phi\rangle$

- Dynamics: $i\partial_t |\psi\rangle = H |\psi\rangle$

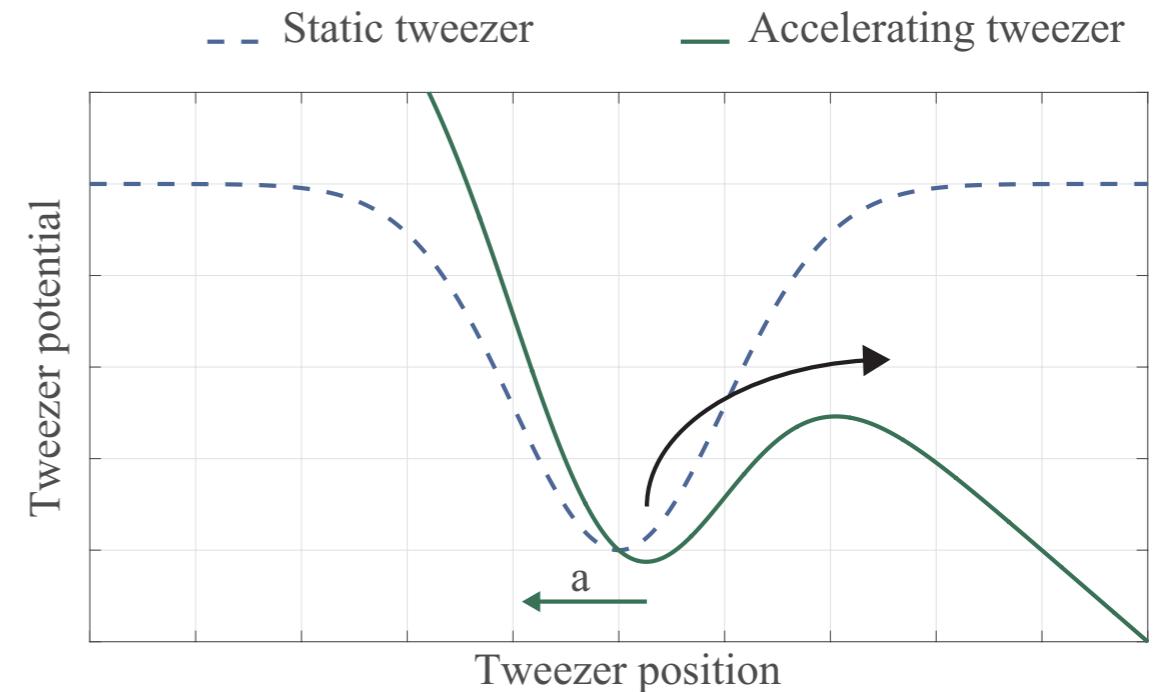
$\rightarrow \tilde{H} = U_\lambda^\dagger H U_\lambda - i U_\lambda^\dagger \partial_t U_\lambda$

- Hence:

$$\tilde{H} = \frac{p^2}{2m} + V(x) - \dot{\lambda} p$$

- Shift the momentum:

$$\tilde{H}' = \frac{p^2}{2m} + V(x) + m\ddot{\lambda}x$$



General

Time-dependent Hamiltonian:

$$H(\lambda_\mu(t))$$

Adiabatic gauge potential:

$$A_\mu = iU^\dagger \partial_{\lambda_\mu} U$$

Moving frame Hamiltonian:

$$\tilde{H} = U^\dagger H U - \dot{\lambda}_\mu A_\mu$$

Adiabatic gauge potential

- Berry connection is the expectation value of the gauge potential

$$\langle 0 | A_\mu | 0 \rangle$$

- Berry curvature:

$$F_{\mu\nu} = -i \langle 0 | [A_\mu, A_\nu] | 0 \rangle$$

- Metric tensor

$$g_{\mu\nu} = \frac{1}{2} \langle 0 | A_\mu A_\nu + A_\nu A_\mu | 0 \rangle$$

- Defines a metric on the ground state manifold. Can be used to detect and classify phase transitions.
- Universal near second order QPT: $g \sim |\lambda - \lambda_c|^{d\nu-2}$

Non-adiabatic response

- Let's look at the corrections to the ground state energy

General

$$\tilde{H} = U^\dagger H U - \dot{\lambda}_\mu A_\mu$$

First order:

$$\Delta E_1 = \dot{\lambda}_\mu \langle 0 | A_\mu | 0 \rangle$$

Second order:

$$\Delta E_2 = \sum_{n=0} \frac{|\langle 0 | \dot{\lambda}_\mu A_\mu | n \rangle|^2}{E_0 - E_n}$$

Translation

$$\tilde{H} = U^\dagger H U - \dot{\lambda} p$$

First order:

$$\Delta E_1 = \dot{\lambda} \langle 0 | p | 0 \rangle = 0$$

Second order:

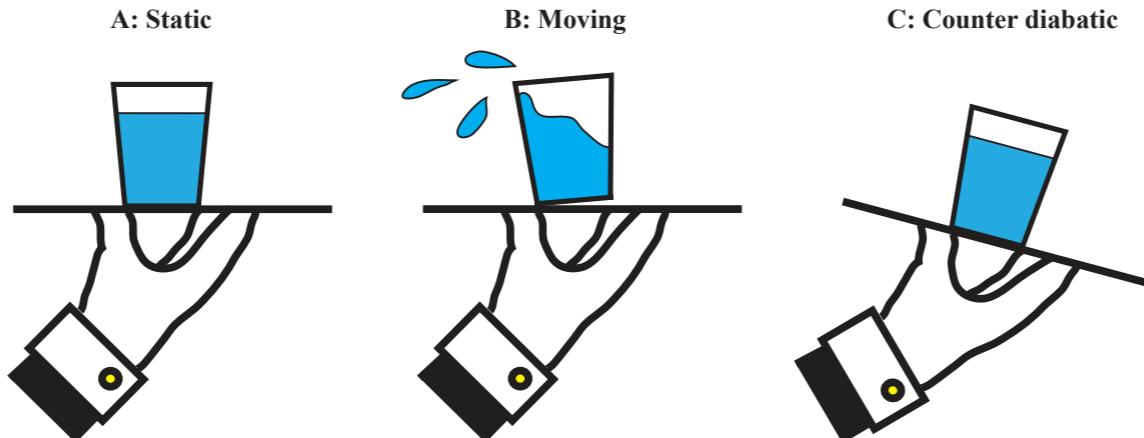
$$\Delta E_2 = \dot{\lambda}^2 \sum_{n \neq 0} \frac{|\langle 0 | p | n \rangle|^2}{E_0 - E_n}$$

Effective mass of the classical parameter: $\Delta E_2 = -\frac{m \dot{\lambda}^2}{2}$

Counter-diabatic drive

- Can we keep the system adiabatic even if we change our parameters fast?
 - Just cancel out the gauge potential!

$$H_{CD} = H(\lambda_t) + \dot{\lambda}_\mu A_\mu$$



- Rotate back into H if possible

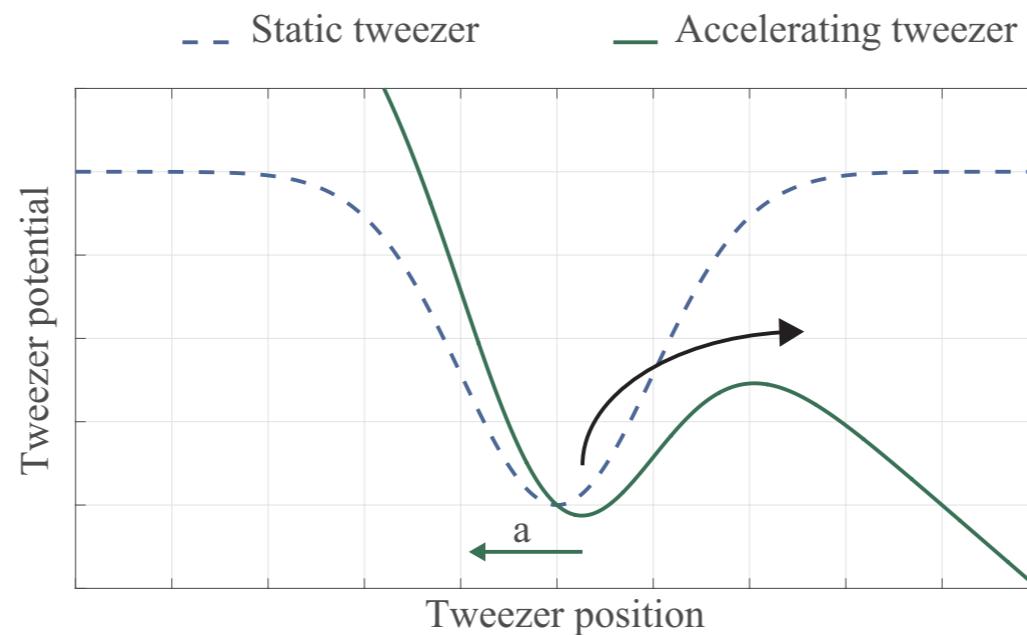
$$H_{FF} = R^\dagger H_{CD}(\lambda_t) R - i R^\dagger \partial_t R$$

Let's play a game



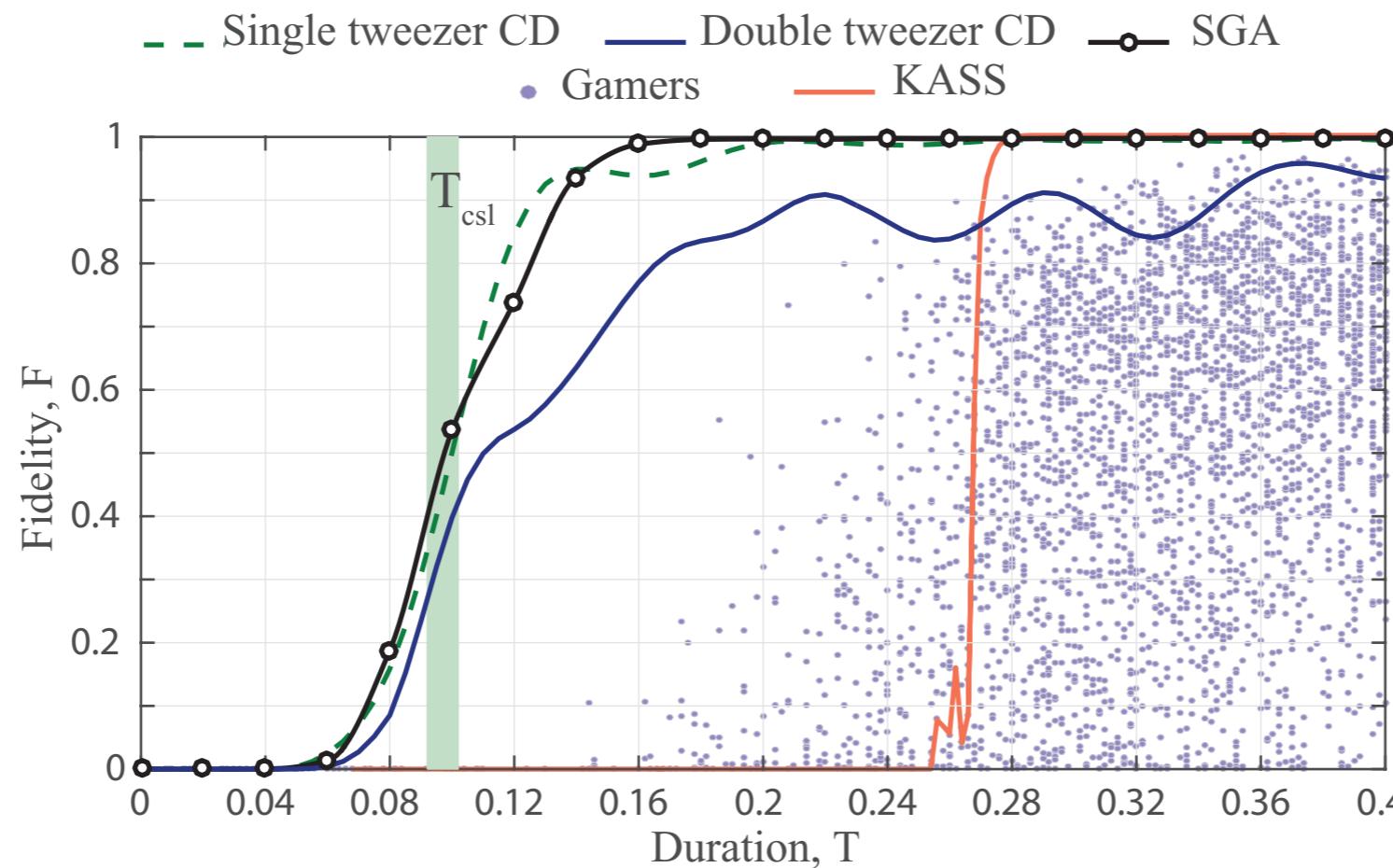
Effective translation

- System is described by: $H = \frac{p^2}{2m} + V_0(x) + V_1(x - \lambda)$
- Let's assume: $A_\lambda = v(\lambda)(xp + px)$
- Transform to: $H_{FF} = H - \partial_t(v(\lambda)\dot{\lambda})x \approx \frac{p^2}{2m} + V_0(x) + V_1(x - \lambda_{FF}(t))$



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Many body systems

- Whole other ballgame

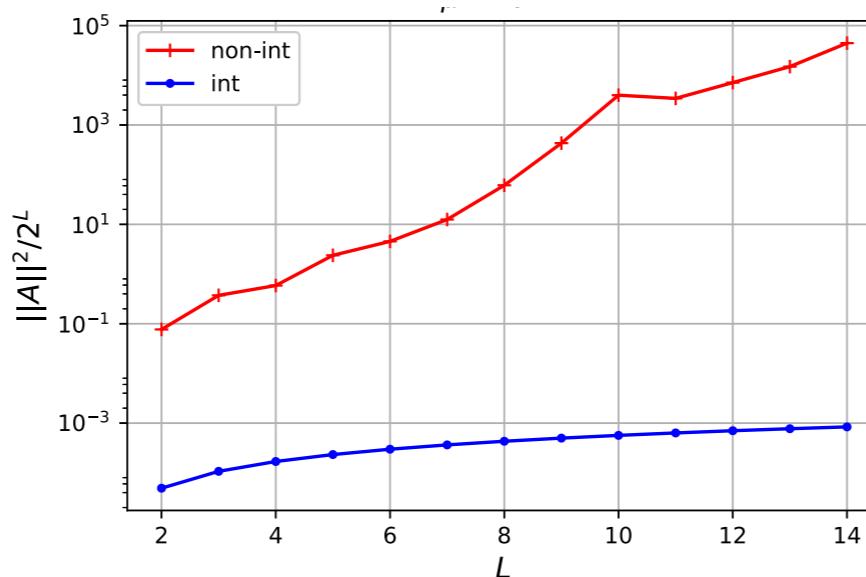
$$\langle n | A_\lambda | m \rangle = \langle n | U^\dagger i\partial_\lambda U | m \rangle = \langle n(\lambda) | i\partial_\lambda | m(\lambda) \rangle$$

- Hence

$$\langle n | A_\lambda | m \rangle = -i \frac{\langle n | \partial_\lambda H | m \rangle}{E_n - E_m}$$

- If the system is ergodic we are in trouble:

$$|\langle n | A_\lambda | m \rangle| = \frac{|\langle n | \partial_\lambda H | m \rangle|}{|E_n - E_m|} \sim \frac{e^{-S/2}}{e^{-S}} = e^{S/2}$$



What now?

- We should be less stringent. Eigenstates of the same energy density are locally indistinguishable anyway.

- Here is our trick:

- Observe: $\langle n | A_\lambda | m \rangle = -i \frac{\langle n | \partial_\lambda H | m \rangle}{E_n - E_m}$

$$\rightarrow \partial_\lambda H + i[A, H] = F_{BO}$$

$$\rightarrow [\partial_\lambda H + i[A, H], H] = 0$$

$$S(O) = \text{Tr}(\partial_\lambda H + i[O, H])^2 \rightarrow A = \text{argmin} S(O)$$

Variational method

$$S(O) = \text{Tr}(\partial_\lambda H + i[O, H])^2 \rightarrow A = \operatorname{argmin} S(O)$$

- Now we can just restrict to sensible, local, experimentally relevant operators.
- No local approximation can ever be close to the exact result, so is minimizing S sensible?
 - Yes! $S(O)$ is a measure for the rate of increase of the energy fluctuations:

$$\frac{\partial \delta E_n^2}{\partial t} \propto \langle n | (\partial_\lambda H + i[A, H])^2 | n \rangle$$

Free fermions

- Let's see what happens for free fermions



$$H = - \sum_i (c_{i+1}^\dagger c_i + h.c.) + \sum_i V(i, \lambda) c_i^\dagger c_i$$

- Let's take the most local Ansatz

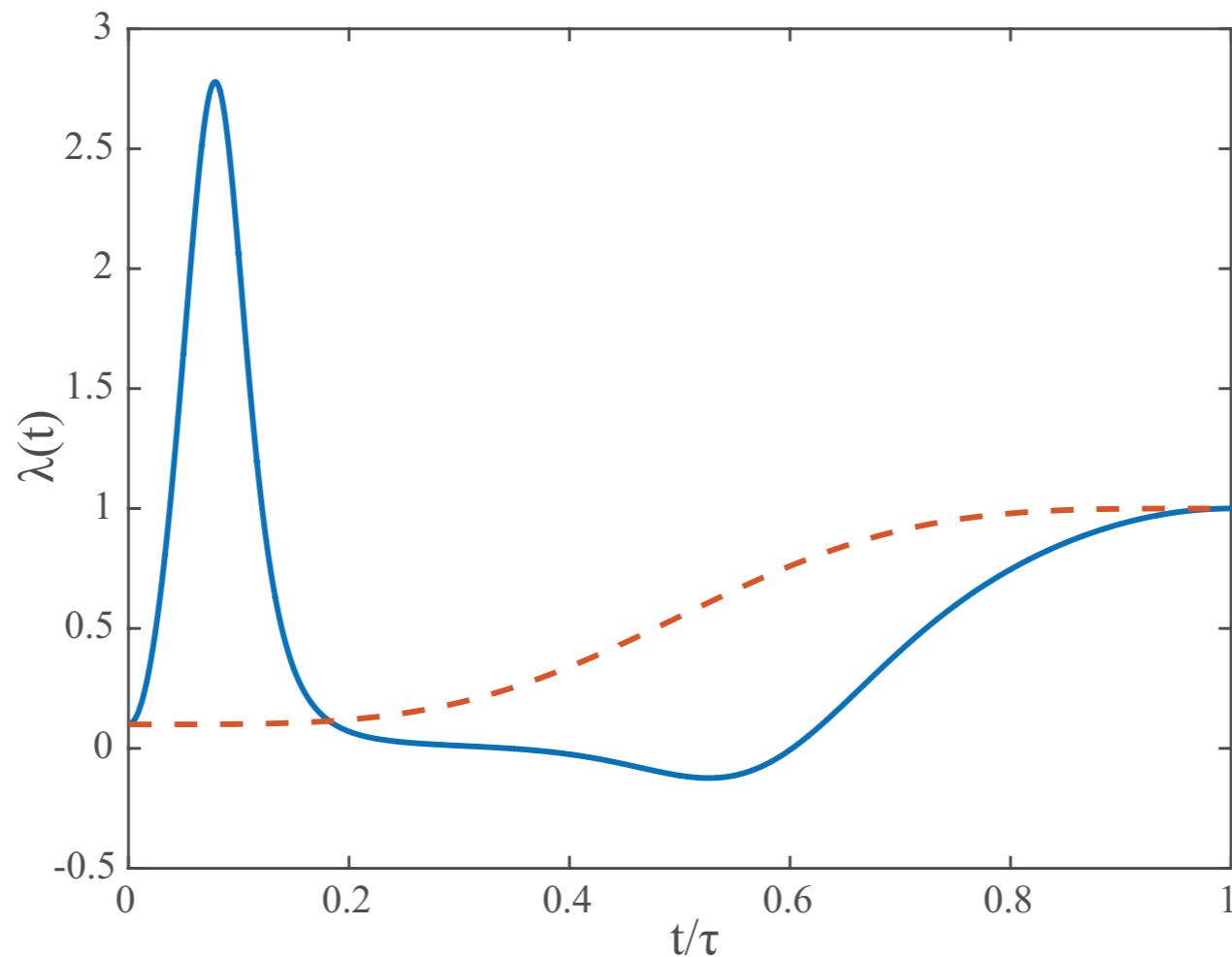
$$A = i \sum_i \alpha_i (c_{i+1}^\dagger c_i - h.c.)$$

- Compute some traces

$$-3\partial_i^2 \alpha + (\partial_i V)^2 \alpha = \partial_i \partial_\lambda V$$

- Exact solution for linear potential

Compensating electric field



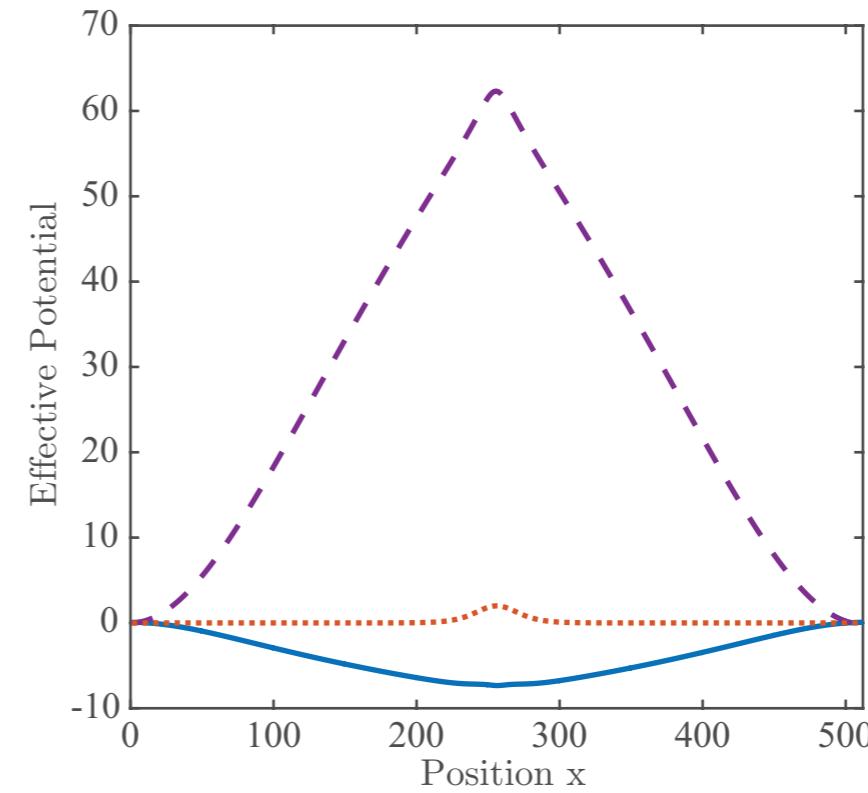
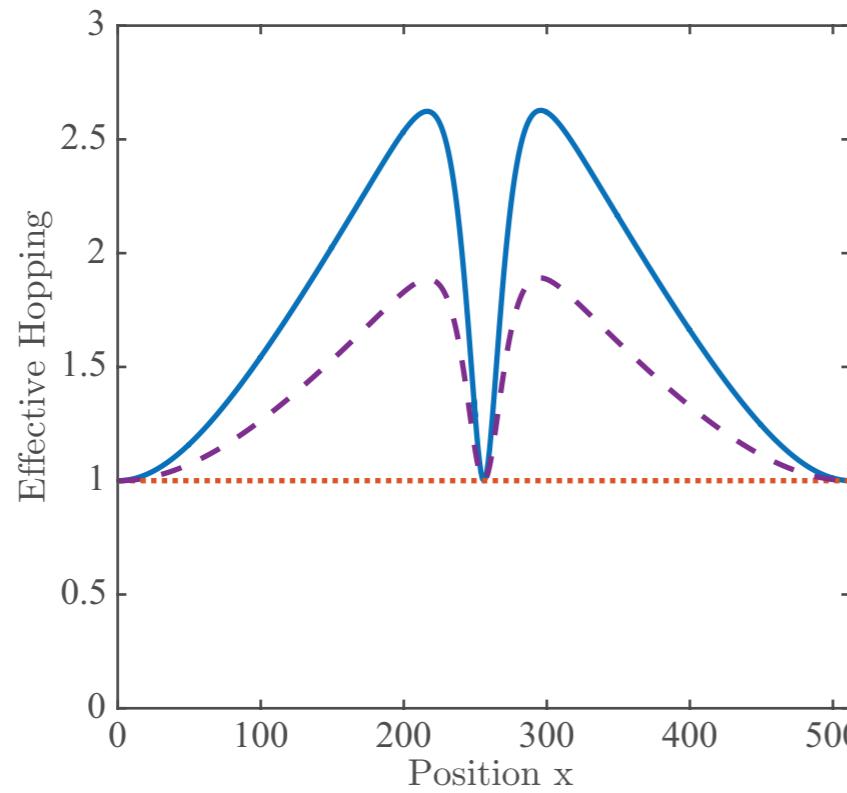
Anderson orthogonality I

- Recall:

$$H_{CD} = H + \dot{\lambda} A_\lambda = - \sum_i ((1 - i\alpha_i) c_i^\dagger c_i + h.c.) + \sum_i V(i, \lambda) c_i^\dagger c_i$$

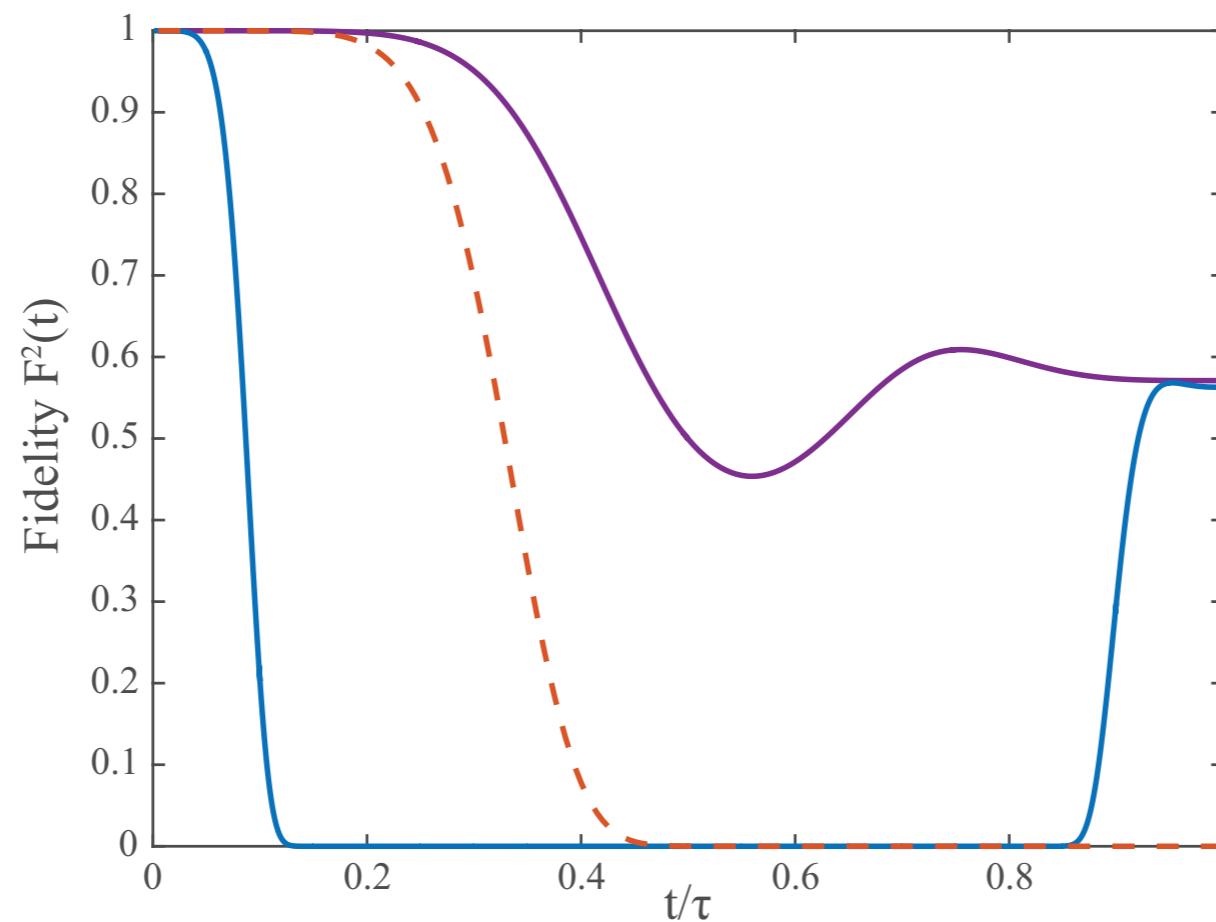
- Gauge transform

$$H_{FF} = - \sum_i (J_{eff}(i) c_i^\dagger c_i + h.c.) + \sum_i V_{eff}(i, \lambda) c_i^\dagger c_i$$

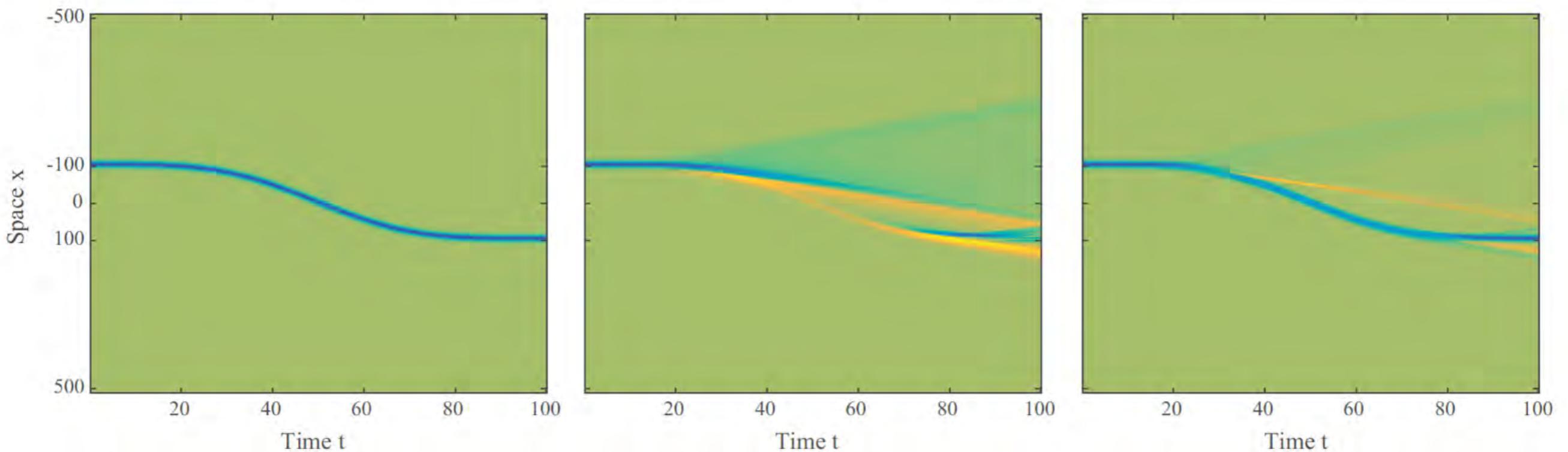


Anderson II

- Chain of 512 sites, half filling, final strength is $2J$ and time is $10/J$.



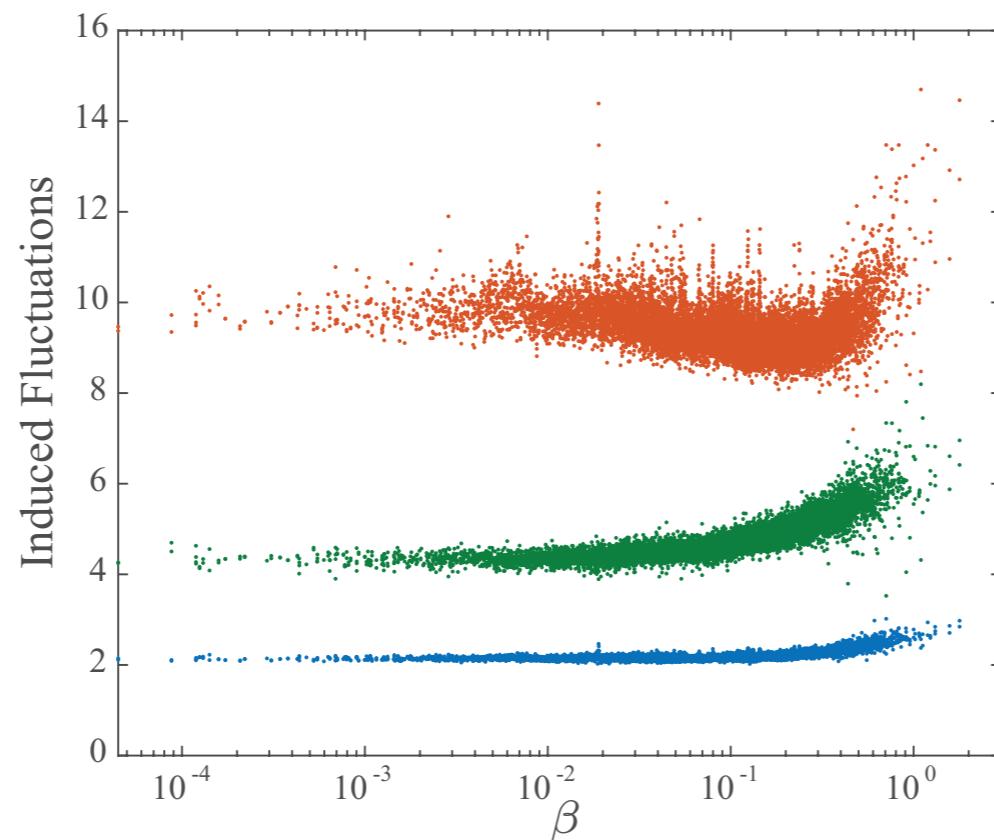
Moving impurity



Flipping spins

- Consider
- $$H = \sum_i (\sigma_{i+1}^z \sigma_i^z + 0.8\sigma_j^z + 0.9\sigma_j^x + h_x(t)\sigma_0^x)$$
- Then we can expand in terms of strings of spins

$$H = \sum_i \alpha_i \sigma_i^y + \sum_{ij} (\beta_{i,j} \sigma_i^x \sigma_j^y + \gamma_{i,j} \sigma_i^z \sigma_j^y)$$



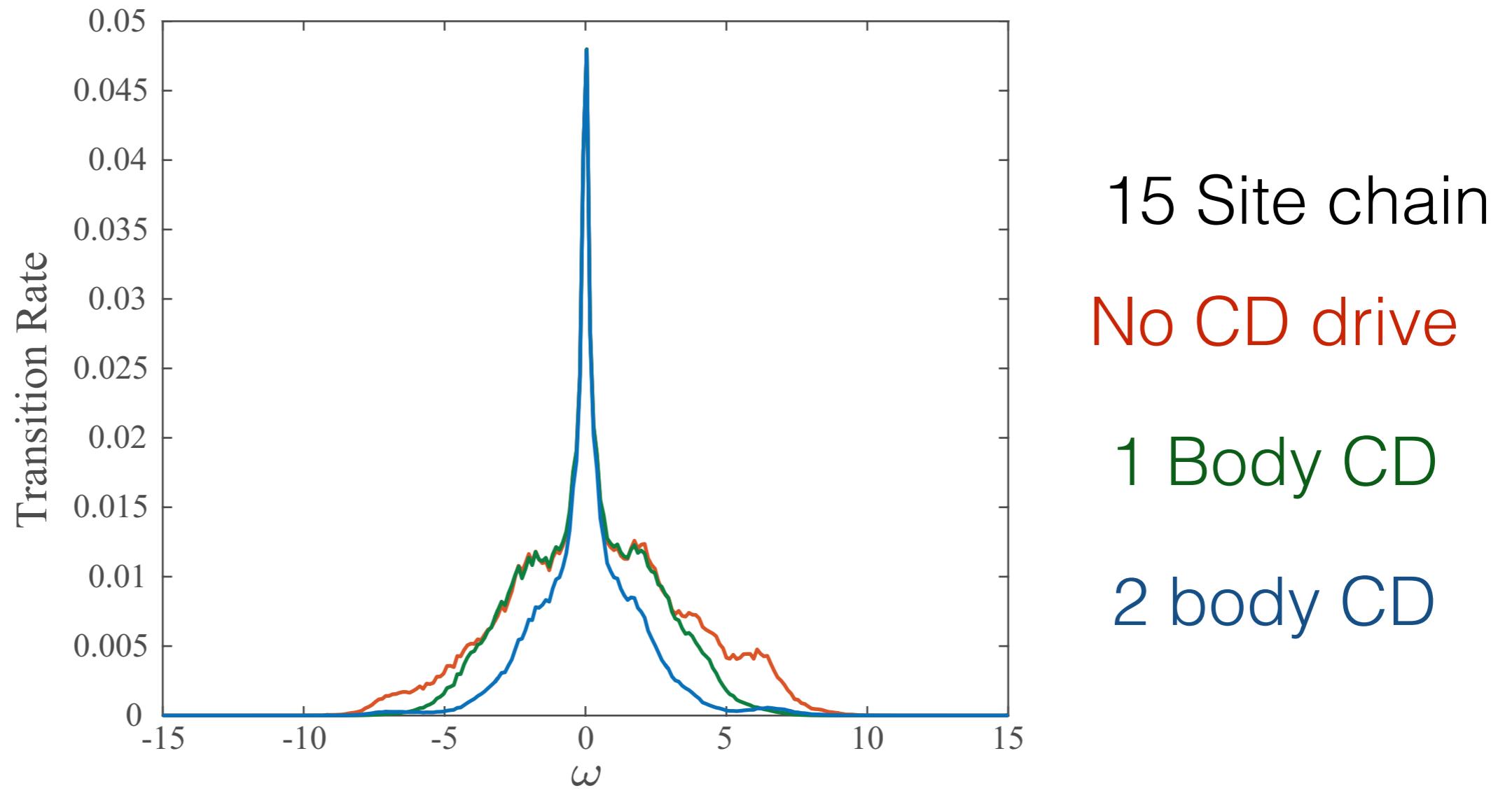
15 Site chain

No CD drive

1 Body CD

2 body CD

Transition rate



Conclusion

- Deep connection between non-adiabatic response and geometry
- Counter-diabatic driving can be used to suppress dissipation in many-body systems
- Simple local expansion for gapped systems
- Physical gauge potential for ergodic system seems only polynomially complex

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M. Kolodrubetz, DS, Pankaj Mehta, AP, Physics Reports 697, 1-88 (2017)