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# Quantum simulation on near-term hardware

Is there any hope?

Mohan Sarovar

Sandia National Laboratories, Livermore

UC Berkeley AMO seminar, November 2017



U.S. DEPARTMENT OF  
**ENERGY**



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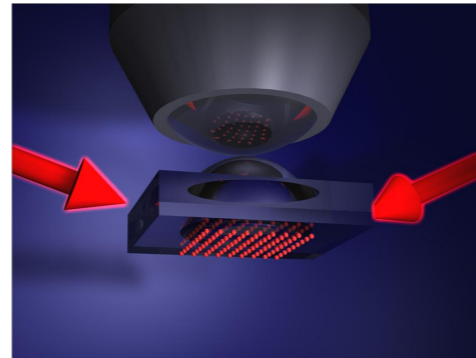
# Analog quantum simulation

- Emulate one quantum system by engineering another to mimic it
- Feynman's original idea for *quantum supremacy*  
Feynman, Int. J. Theor. Phys. **21**, 467 (1982)
- Already an experimental reality:



## Probing the relaxation towards equilibrium in an isolated strongly correlated one-dimensional Bose gas

S. Trotzky<sup>1,2,3\*</sup>, Y.-A. Chen<sup>1,2,3</sup>, A. Fleisch<sup>4\*</sup>, I. P. McCulloch<sup>5</sup>, U. Schollwöck<sup>1,6</sup>, J. Eisert<sup>6,7,8</sup> and I. Bloch<sup>1,2,3</sup>



Greiner Lab, Harvard

## ARTICLE

doi:10.1038/nature09994

## Quantum simulation of antiferromagnetic spin chains in an optical lattice

Jonathan Simon<sup>1</sup>, Waseem S. Bakr<sup>1</sup>, Ruichao Ma<sup>1</sup>, M. Eric Tai<sup>1</sup>, Philipp M. Preiss<sup>1</sup> & Markus Greiner<sup>1</sup>

nature

Vol 465 | 3 June 2010 | doi:10.1038/nature09071

## LETTERS

## Quantum simulation of frustrated Ising spins with trapped ions

K. Kim<sup>1</sup>, M.-S. Chang<sup>1</sup>, S. Korenblit<sup>1</sup>, R. Islam<sup>1</sup>, E. E. Edwards<sup>1</sup>, J. K. Freericks<sup>2</sup>, G.-D. Lin<sup>3</sup>, L.-M. Duan<sup>3</sup> & C. Monroe<sup>1</sup>

## New Journal of Physics

The open-access journal for physics

## Quantum simulation of the transverse Ising model with trapped ions

K Kim<sup>1,5,6</sup>, S Korenblit<sup>1</sup>, R Islam<sup>1</sup>, E E Edwards<sup>1</sup>, M-S Chang<sup>1,7</sup>, C Noh<sup>2,8</sup>, H Carmichael<sup>2</sup>, G-D Lin<sup>3,9</sup>, L-M Duan<sup>3</sup>, C C Joseph Wang<sup>4</sup>, J K Freericks<sup>4</sup> and C Monroe<sup>1</sup>

# Analog quantum simulation

Complex control requirements for any non-trivial quantum simulation experiment

Analog quantum simulation has no natural notion of error correction

How does one assess when an analog quantum simulation result is robust to noise/imperfections?

Rep. Prog. Phys. **75** (2012) 082401 (18pp)

[doi:10.1088/0034-4885/75/8/082401](https://doi.org/10.1088/0034-4885/75/8/082401)

## Can one trust quantum simulators?

Philipp Hauke<sup>1</sup>, Fernando M Cucchietti<sup>1,2</sup>, Luca Tagliacozzo<sup>1</sup>,  
Ivan Deutsch<sup>3,4</sup> and Maciej Lewenstein<sup>1,5</sup>

# But is it so bad?

## Physicist's intuition/consolation

1. For many questions we “care about” (coarse-grained observables) microscopic imperfections in the model (Hamiltonian) should not matter. After all, *no real material corresponds perfectly to an ideal model.*
2. Real materials are at finite temperature, so if we introduce an equilibration/thermalization mechanism then the emergent properties may be stable.

### Aims:

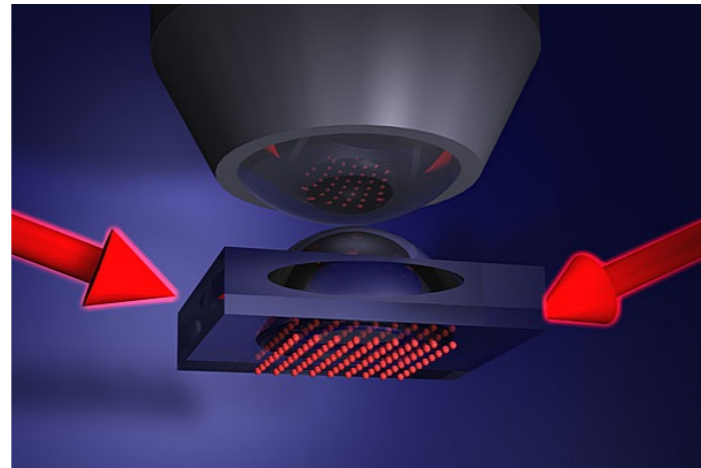
1. To formalize (1)
2. To enable (2)

# Outline

1. Parameter space compression and simulation sensitivity
2. Engineering thermalization in quantum many-body models

# Quantifying parameter sensitivity

A quantum simulator produces parameterized probability distributions

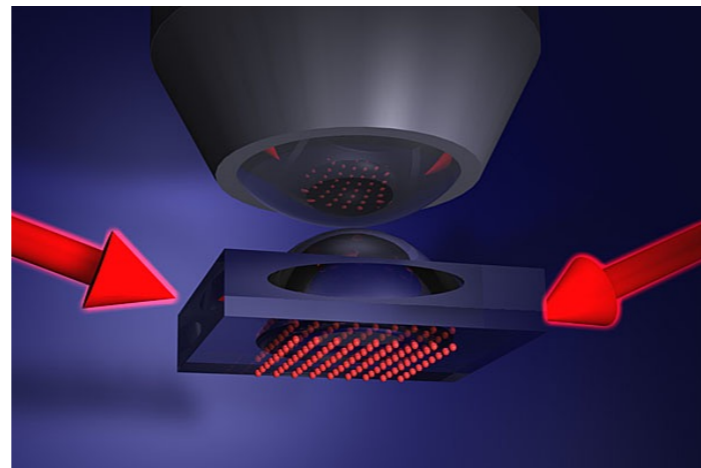


$$\{p_m\}_m$$

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# Quantifying parameter sensitivity

A quantum simulator produces parameterized probability distributions



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$\{P_m\}_m$

Ideal many-body Hamiltonian dependent on  $K$  parameters

$$H(\lambda) = \sum_{k=1}^K \lambda_k H_k \quad \lambda = (\lambda_1, \lambda_2, \dots, \lambda_K)$$

Observable of interest

$$O = \sum_m \theta_m P_m$$

Produces a parameterized probability distribution

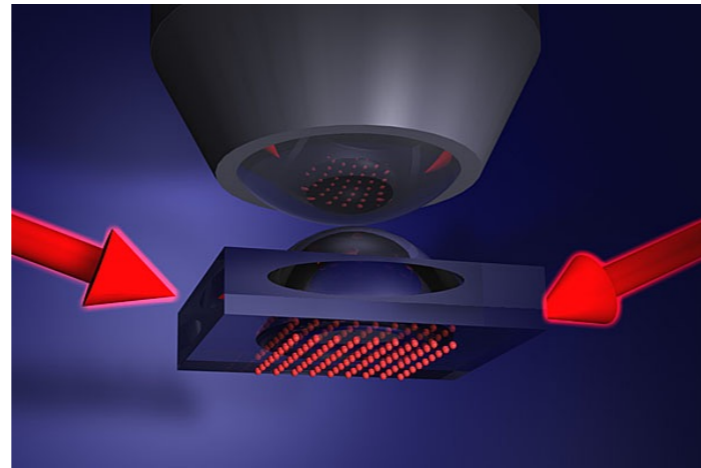
$$p_m(\lambda) = \text{tr}(P_m \rho_{\text{th}}(T))$$

← Or ground state

$$\rho_{\text{th}}(T) = \frac{e^{\beta H(\lambda)}}{\mathcal{Z}}$$
$$\beta = \frac{1}{k_B T}$$

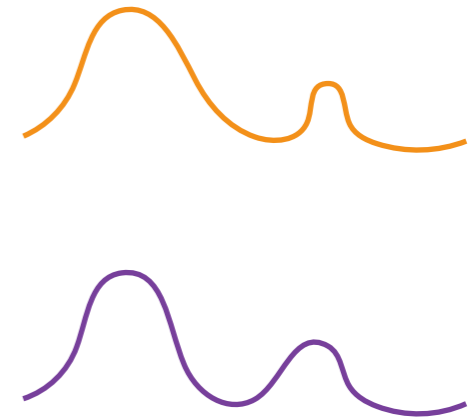
# Quantifying parameter sensitivity

A quantum simulator produces parameterized probability distributions



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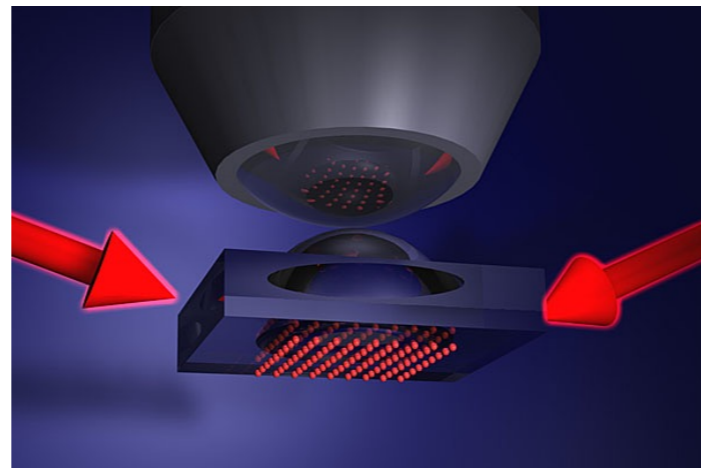
$\lambda^0$   
 $\lambda$



**Robustness: how different are the output distributions under perturbations of the Hamiltonian parameters?**

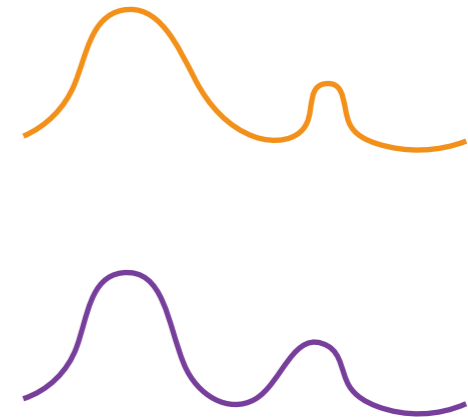
# Quantifying parameter sensitivity

A quantum simulator produces parameterized probability distributions



Greiner Lab, Harvard

$\lambda^0$   
 $\lambda$



**Robustness: how different are the output distributions under perturbations of the Hamiltonian parameters?**

Kullback-Leibler divergence: a measure of difference between probability distributions

$$\begin{aligned} D_{\text{KL}}(p(\lambda) || p(\lambda^0)) &= \sum_m p_m(\lambda) \log \frac{p_m(\lambda)}{p_m(\lambda^0)} \\ &= \frac{1}{2} \Delta\lambda^\top F(\lambda^0) \Delta\lambda + \mathcal{O}(\|\Delta\lambda\|^3) \quad (\text{small deviations}) \end{aligned}$$

Fisher information matrix (FIM)

$$\Delta\lambda = \lambda - \lambda^0$$

# Quantifying parameter sensitivity

Fisher Information matrix (FIM)

$$F_{ij}(\lambda^0) = \sum_{m=1}^M \frac{1}{p_m(\lambda)} \frac{\partial p_m(\lambda)}{\partial \lambda_i} \frac{\partial p_m(\lambda)}{\partial \lambda_j} \Big|_{\lambda=\lambda^0}$$

Spectral analysis of the FIM

$$F = \sum_{k=1}^K \zeta_k v_k v_k^\dagger$$

Eigendecomposition of FIM

$$D_{\text{KL}}(p(\lambda) || p(\lambda^0)) \approx \frac{1}{2} \Delta \lambda^\top F(\lambda^0) \Delta \lambda = \sum_{k=1}^K \frac{\zeta_k}{2} \|v_k^\dagger \Delta \lambda\|^2$$

Quantum simulation of this model is trivially robust if all  $\zeta_k \approx 0$

# Quantifying parameter sensitivity

Spectral analysis of the FIM

$$F = \sum_{k=1}^K \zeta_k v_k v_k^\dagger$$

Eigendecomposition of FIM

## Key observation:

Eigenvalues of FIM prescribe an ordering of parameter influence, *i.e.*  
if  $\zeta_k$  is large, then the parameter

$$\sum_j \lambda_j v_k^j$$

has a large influence on observations

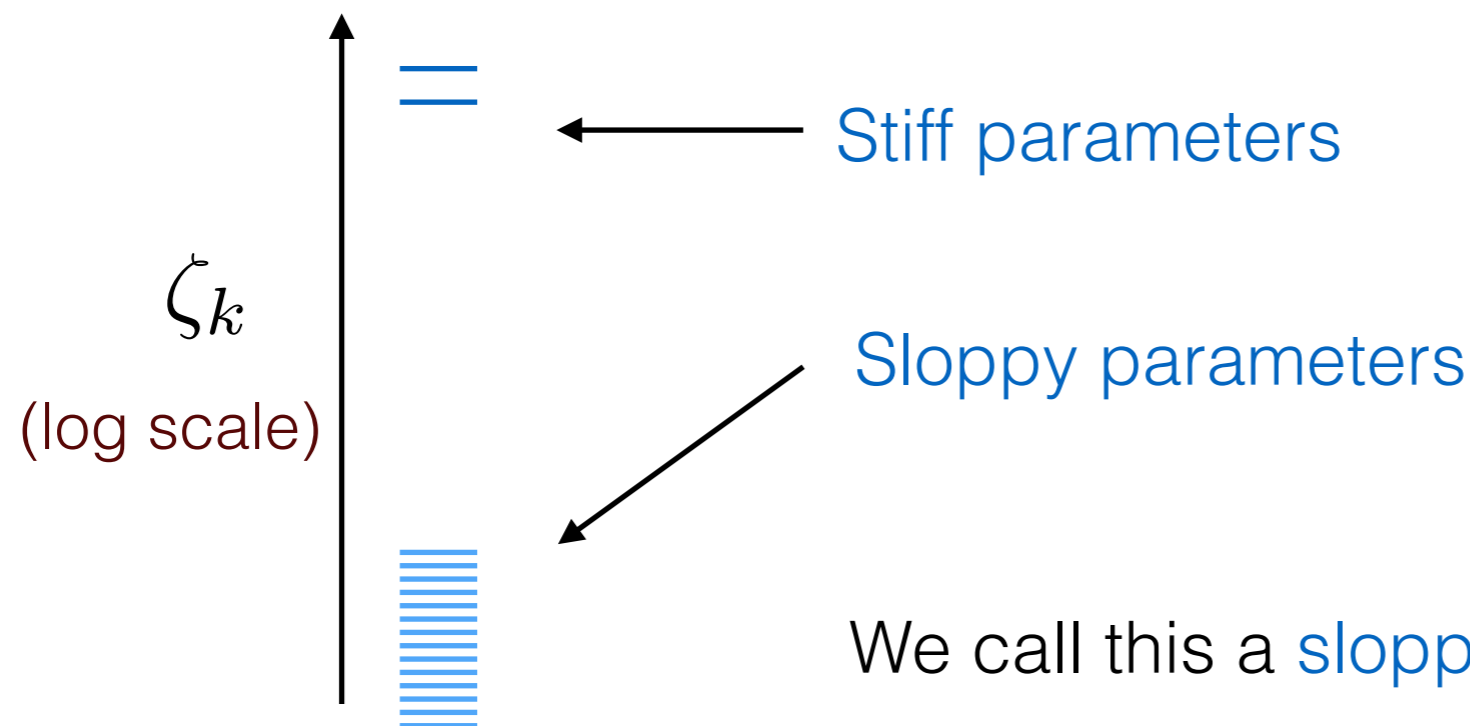
So a *non-trivial* way for a quantum simulation to be robust is that only a few eigenvalues matter, and they correspond to parameters that can be controlled well.

# Quantifying parameter sensitivity

Spectral analysis of the FIM

$$F = \sum_{k=1}^K \zeta_k v_k v_k^\dagger$$

Eigendecomposition of FIM



We call this a sloppy quantum simulation model

c.f.



**Parameter Space Compression Underlies Emergent Theories and Predictive Models**

Benjamin B. Machta *et al.*

*Science* **342**, 604 (2013);

DOI: 10.1126/science.1238723

# Quantifying parameter sensitivity

Spectral analysis of the FIM

$$F = \sum_{k=1}^K \zeta_k v_k v_k^\dagger$$

Eigendecomposition of FIM

Are quantum many-body models sloppy?  
If so, why?

We call this a **sloppy quantum simulation model**

c.f.



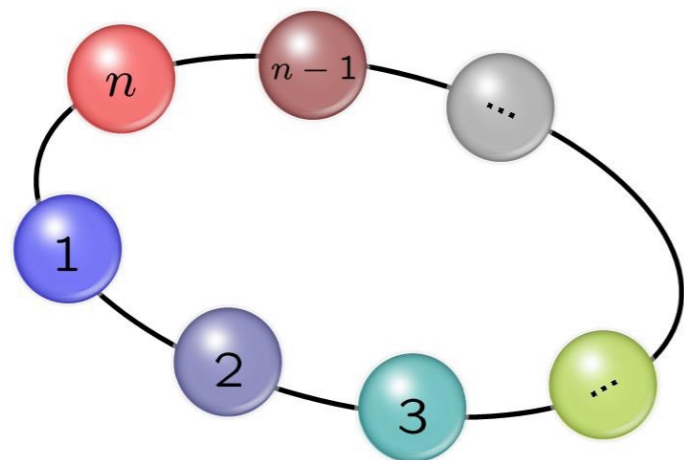
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# Transverse-field Ising model



$$H_{1D \text{ Ising}} = \sum_{i=1}^n B_i \sigma_x^i + \sum_{i=1}^{n-1} J_i \sigma_z^i \sigma_z^{i+1} + J_n \sigma_z^n \sigma_z^1$$

Uniform antiferromagnetic model parameter regime, so we want:

$$B_i = B^0 \quad \forall i$$

$$J_i = J^0 > 0 \quad \forall i$$

Quantum critical point at

$$B^0 = \frac{J^0}{2}$$

Observables of interest

$$S_z = \sum_i \langle \sigma_z^i \rangle$$

Net magnetization

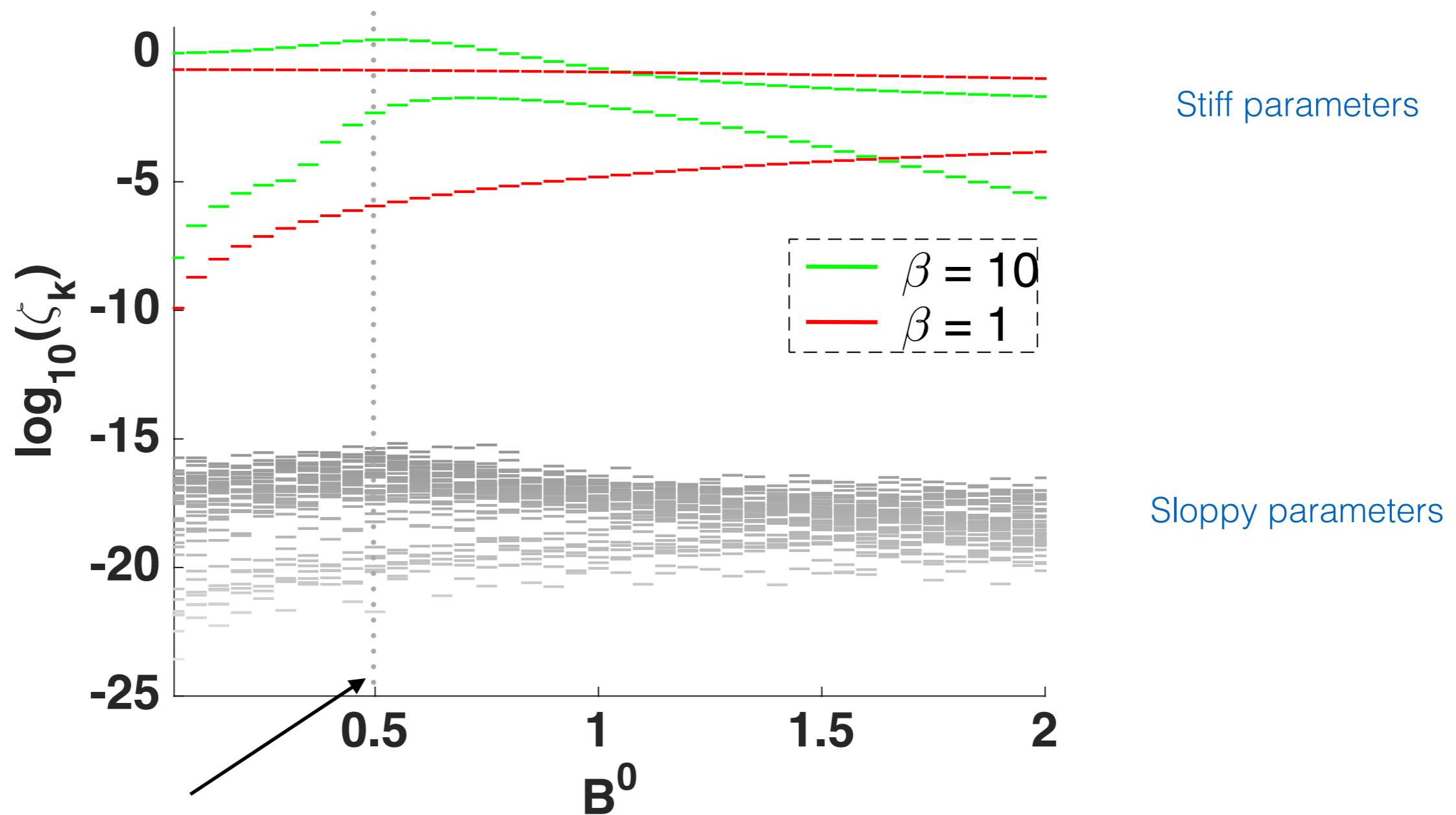
$$C_z(i, j) = \langle \sigma_z^i \sigma_z^j \rangle$$

Correlation functions

# Transverse-field Ising model

Net magnetization observable

$$n = 10, J^0 = 1$$

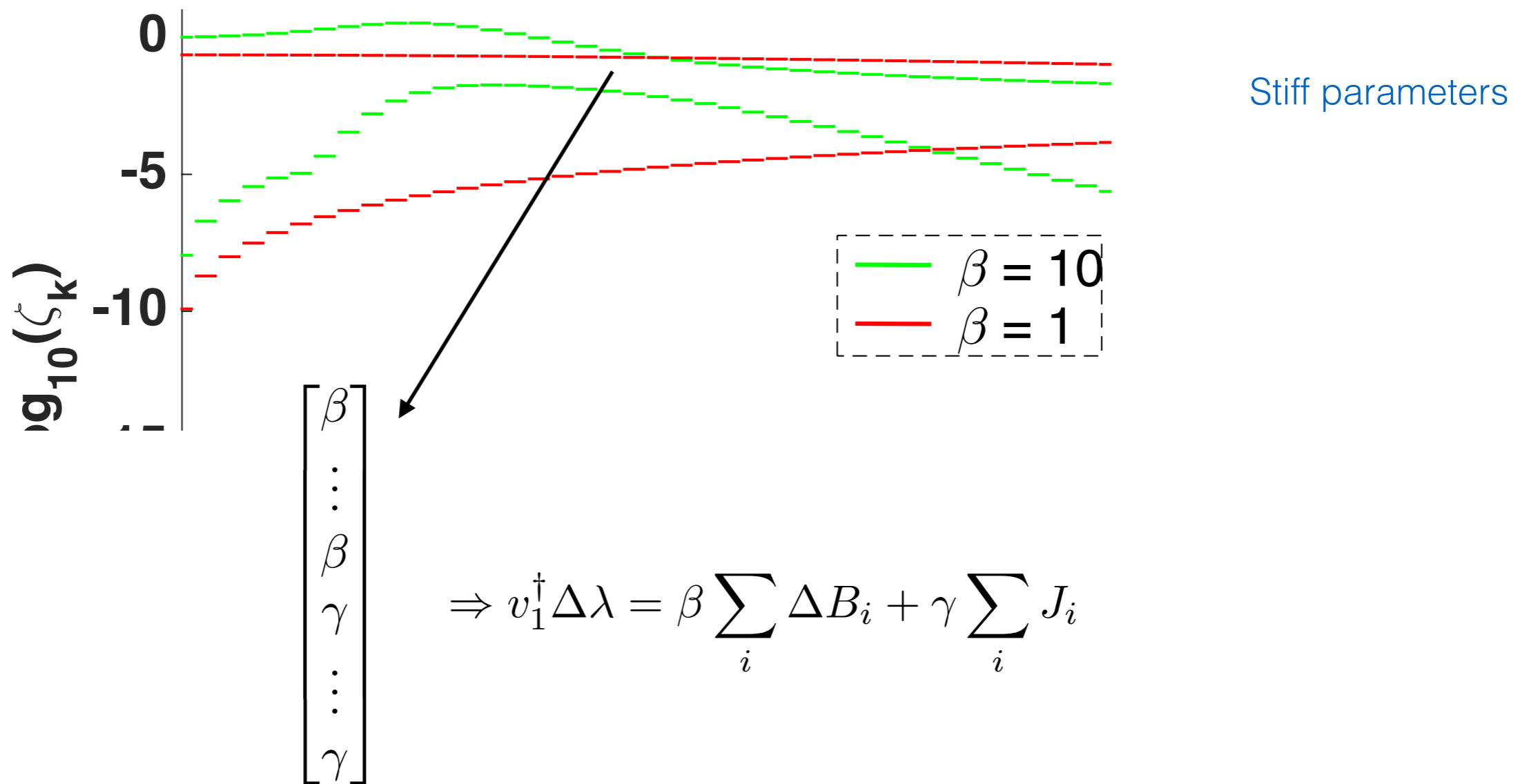


Quantum critical point

# Transverse-field Ising model

Net magnetization observable

What is the dominant stiff parameter?  $n = 10, J^0 = 1$



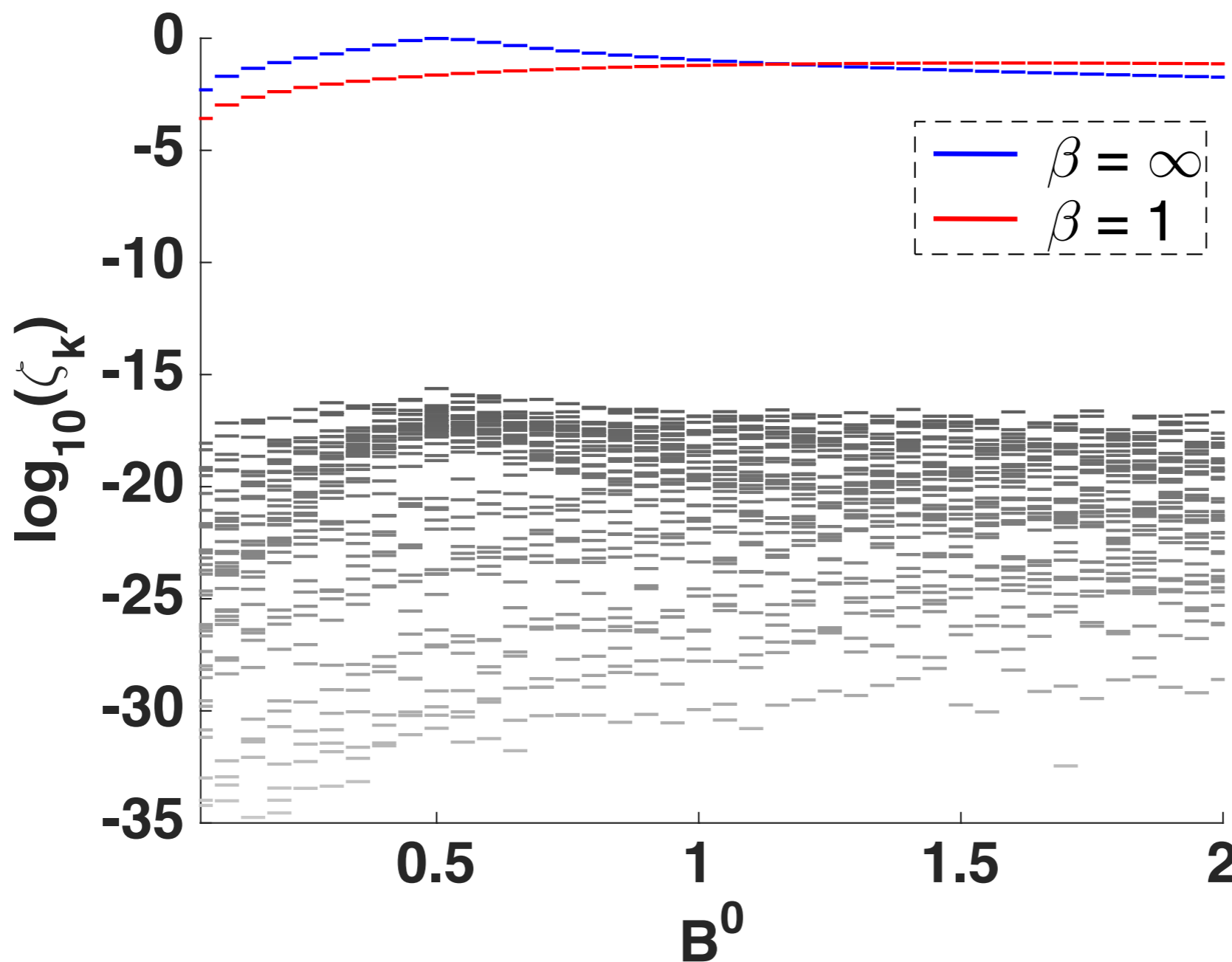
Fluctuations in  $B_i$  and  $J_i$  do not matter, as long as these (spatially) average to zero

# Transverse-field Ising model

Correlation function observable

$$C_z(2, 6)$$

$$n = 10, J^0 = 1$$



Stiff parameter

Sloppy parameters

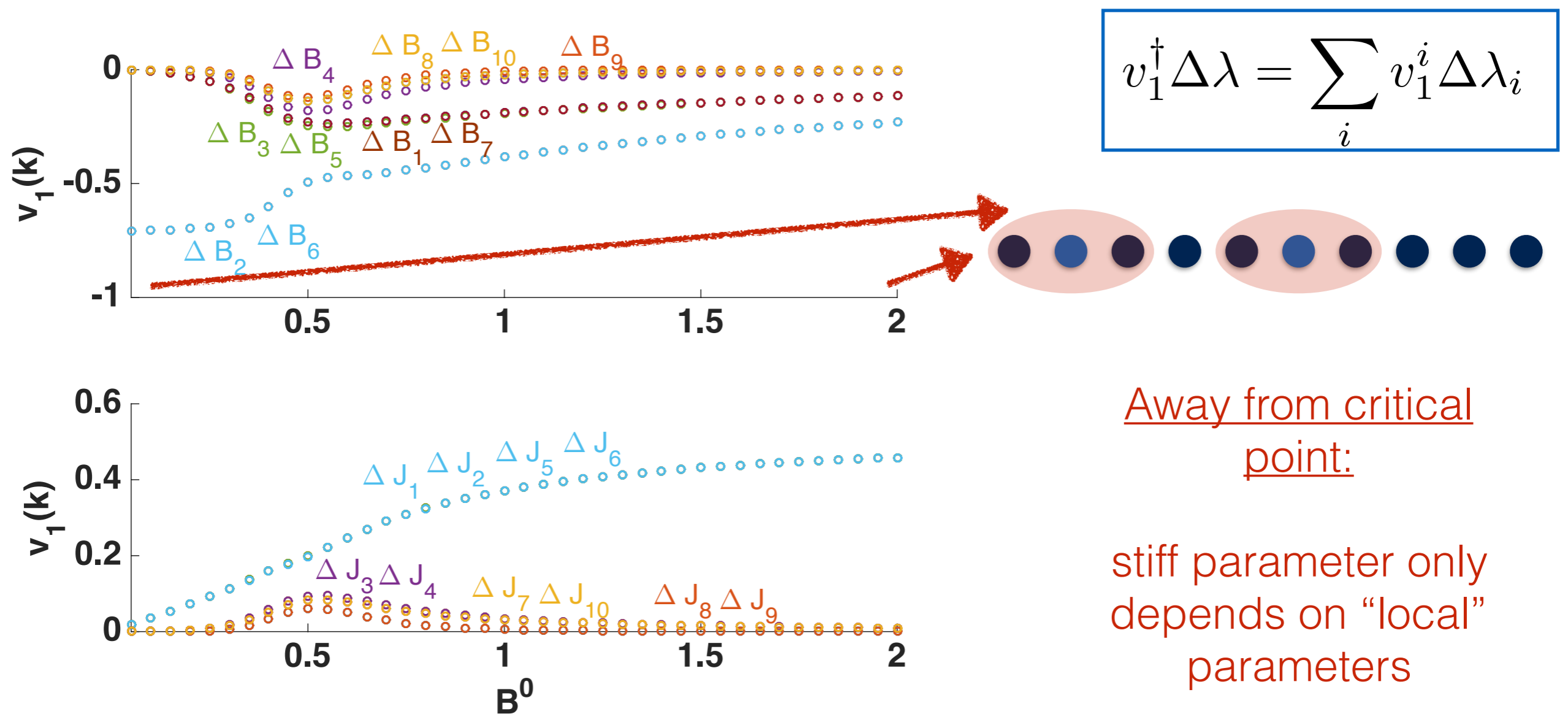
# Transverse-field Ising model

Correlation function observable

$$C_z(2, 6)$$

$$n = 10, J^0 = 1$$

Composition of stiff parameter



Away from critical point:

stiff parameter only depends on “local” parameters

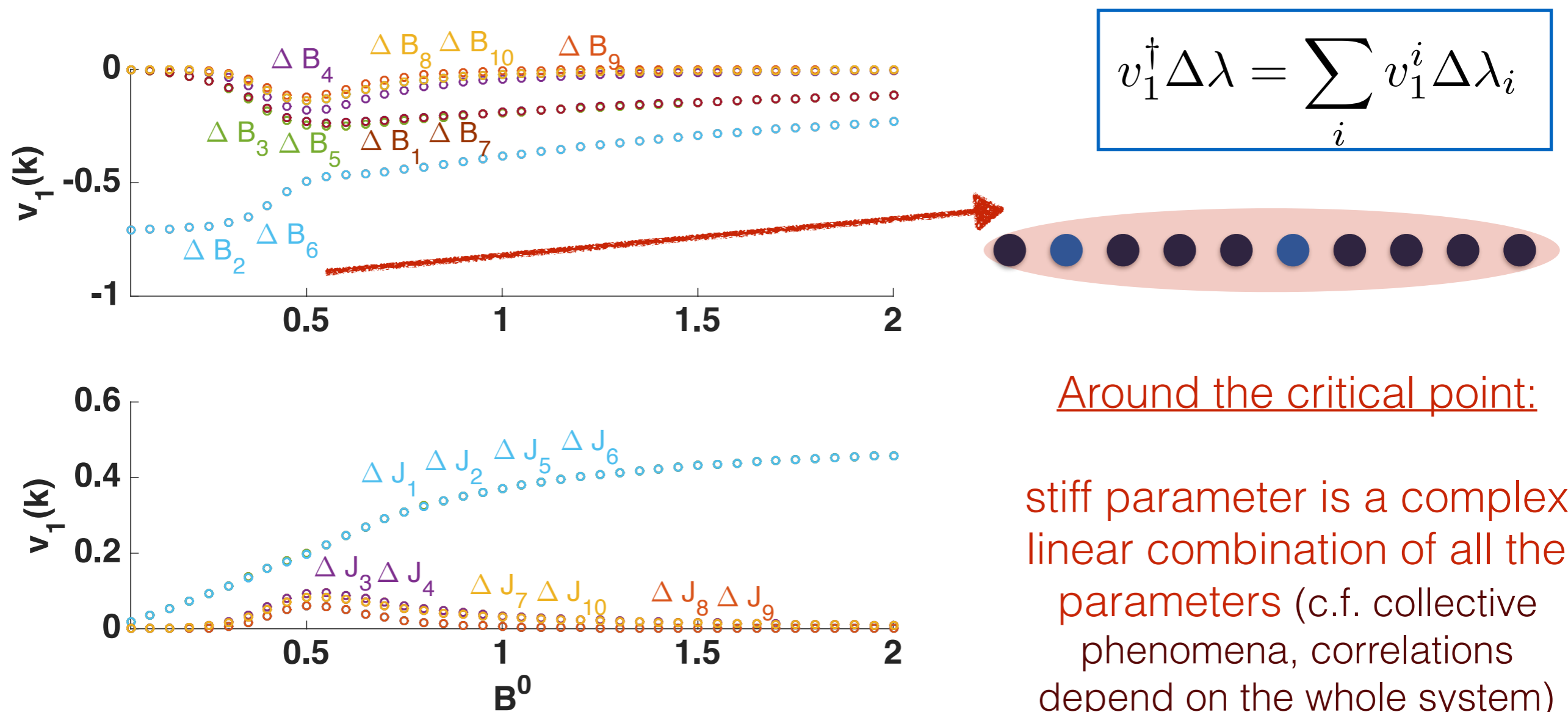
# Transverse-field Ising model

Correlation function observable

$$C_z(2, 6)$$

$$n = 10, J^0 = 1$$

Composition of stiff parameter



# Transverse-field Ising model

## Lessons learnt:

- **The observable matters.** Collective, global phenomena can be very robust to local fluctuations.
- **Quantum phase transitions** can lead to complex dependence of stiff parameters on microscopic parameters.
- Generally: sloppiness increases with **temperature**.
- **Specifically for Ising model:** detect phase transition by measuring magnetic susceptibility as opposed to measuring a correlation function.

# Other models

We have studied many other quantum-many body models of interest:

1. 1D and 2D transverse-field Ising models with varying boundary conditions
2. Several flavors of the Heisenberg model
3. J1-J2-anti-ferromagnetic Heisenberg model (non-nearest neighbor interactions)
4. Fermi-Hubbard model
5. Random/disordered transverse-field Ising model (not sloppy)

See:

Reliability of analog quantum simulation,  
M. S., J. Zhang, L. Zeng. EPJ Quantum Information 4,1 (2017)  
[arXiv:1603.09283](https://arxiv.org/abs/1603.09283)

# Why so sloppy?

Fisher Information matrix (FIM)

$$F_{ij}(\lambda^0) = \sum_{m=1}^M \frac{1}{p_m(\lambda)} \frac{\partial p_m(\lambda)}{\partial \lambda_i} \frac{\partial p_m(\lambda)}{\partial \lambda_j} \Big|_{\lambda=\lambda^0}$$

Use rank as a proxy for sloppiness

low rank => sloppy


# Why so sloppy?

Fisher Information matrix (FIM)

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Write as:

$$F = V \Lambda^{-1} V^T$$


$$V = \begin{pmatrix} \frac{\partial p_1(\lambda)}{\partial \lambda_1} & \frac{\partial p_2(\lambda)}{\partial \lambda_1} & \dots & \frac{\partial p_M(\lambda)}{\partial \lambda_1} \\ \vdots & \vdots & \ddots & \vdots \\ \frac{\partial p_1(\lambda)}{\partial \lambda_K} & \dots & \dots & \frac{\partial p_M(\lambda)}{\partial \lambda_K} \end{pmatrix} \quad \Lambda = \begin{pmatrix} p_1(\lambda) & & & \\ & \ddots & & \\ & & \ddots & \\ & & & p_M(\lambda) \end{pmatrix}$$

$K \times M$   $M \times M$

$\Lambda$  is full-rank, therefore  $\text{rank}(F) \leq \text{rank}(V)$

What dictates the rank of  $V$ ?

# Quantum simulation model symmetries

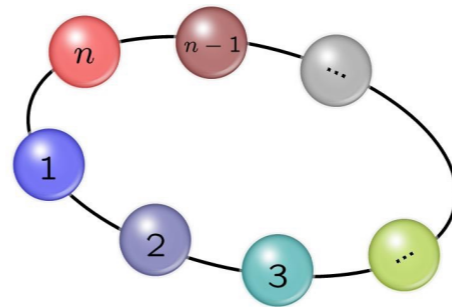
Let  $G$  be the group of symmetries of the quantum simulation model, *i.e.*

$$\begin{aligned} [U_g, H(\lambda)] &= 0 \\ [U_g, O] &= 0 \end{aligned} \quad \forall g$$

$\{U_g\}_g$  a faithful unitary representation of  $G$

*e.g.*

Translational invariance



Recall

$$H(\lambda) = \sum_{k=1}^K \lambda_k H_k$$

$$O = \sum_m \theta_m P_m$$

$$p_m(\lambda) = \text{tr} (P_m \rho_{\text{th}}(T))$$

# Quantum simulation model symmetries

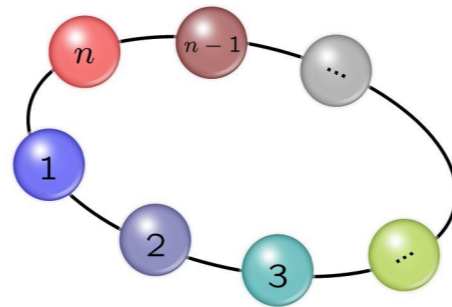
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Suppose

$$U_g H_k U_g^\dagger = H_j$$

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# Quantum simulation model symmetries

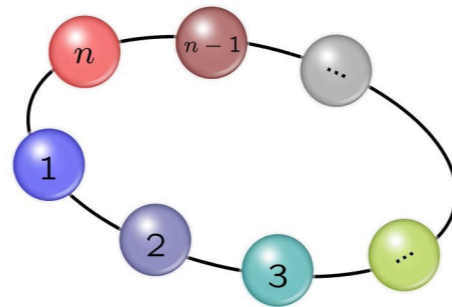
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*e.g.*

Translational invariance



Suppose

$$U_g H_k U_g^\dagger = H_j$$

Then

$$\frac{\partial p_m(\lambda)}{\partial \lambda_k} = \frac{\partial p_m(\lambda)}{\partial \lambda_j} \quad \forall m$$

Repeated rows in  $V$

$$\begin{pmatrix} \frac{\partial p_1(\lambda)}{\partial \lambda_1} & \frac{\partial p_2(\lambda)}{\partial \lambda_1} & \dots & \frac{\partial p_M(\lambda)}{\partial \lambda_1} \\ \vdots & \vdots & \vdots & \vdots \\ \frac{\partial p_1(\lambda)}{\partial \lambda_K} & \dots & \dots & \frac{\partial p_M(\lambda)}{\partial \lambda_K} \end{pmatrix}$$

Model symmetries reduce rank of  $F$

# Quantum simulation model symmetries

Model symmetries reduce rank of  $V$ , and hence of  $F$

Algorithm for computing an upper bound on the rank of  $F$ :

1. For all  $1 \leq k \leq K$ , compute the orbit of  $H_k$  under the symmetry group

$$\text{Orbit: } \{U_g H_k U_g^\dagger \mid g \in G\}$$

2. Number of distinct orbits will be the number of unique rows in  $V$ , and therefore a **bound on the rank of  $V$  and  $F$**

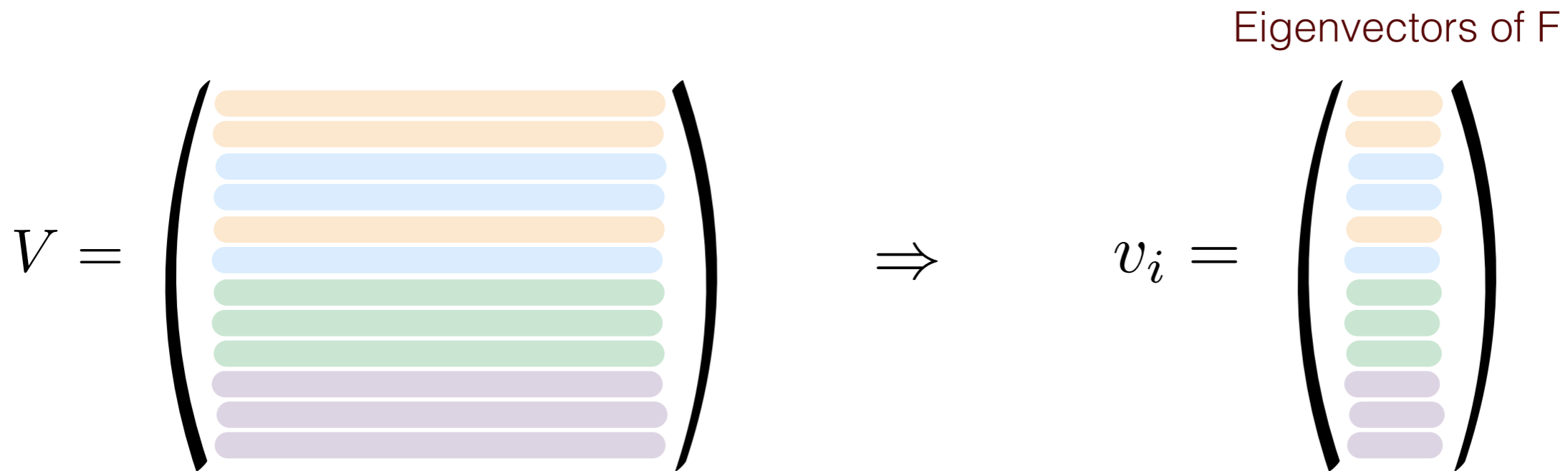
Recall

$$H(\lambda) = \sum_{k=1}^K \lambda_k H_k$$

# Quantum simulation model symmetries

Such a symmetry analysis allows us to extract the form of the stiff/influential parameter as well

$$F = V \Lambda^{-1} V^T$$

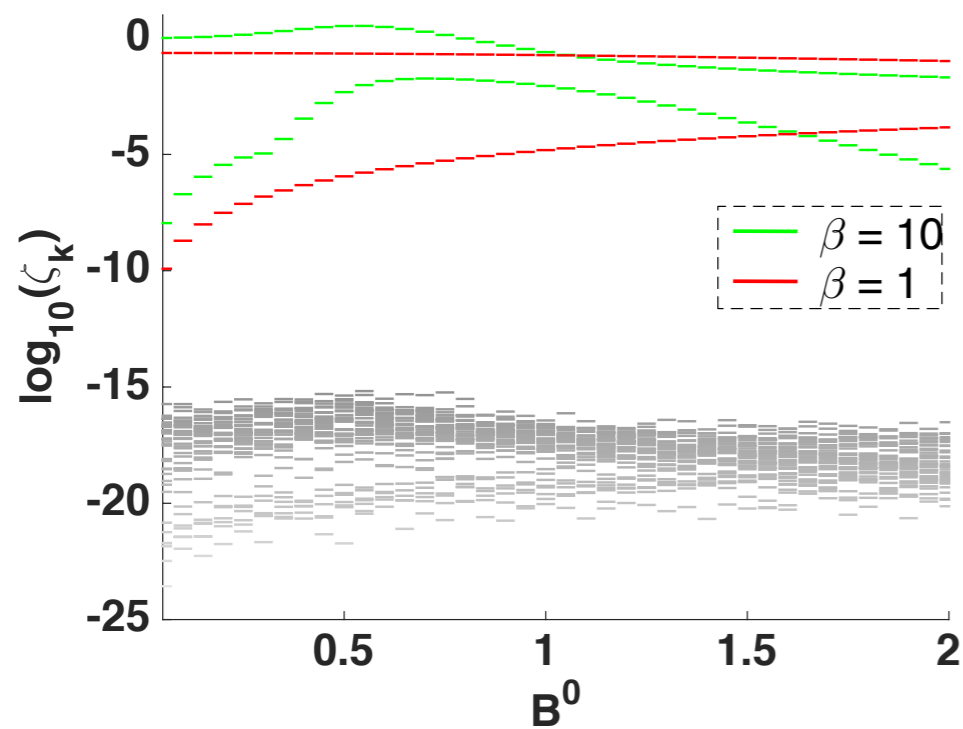


$$v_k^\dagger \Delta \lambda = \sum_s \mu_s^k(\lambda^0, \beta) \sum_{l: H_l \in \text{Orbit } s} \Delta \lambda_l$$

Sum over orbits

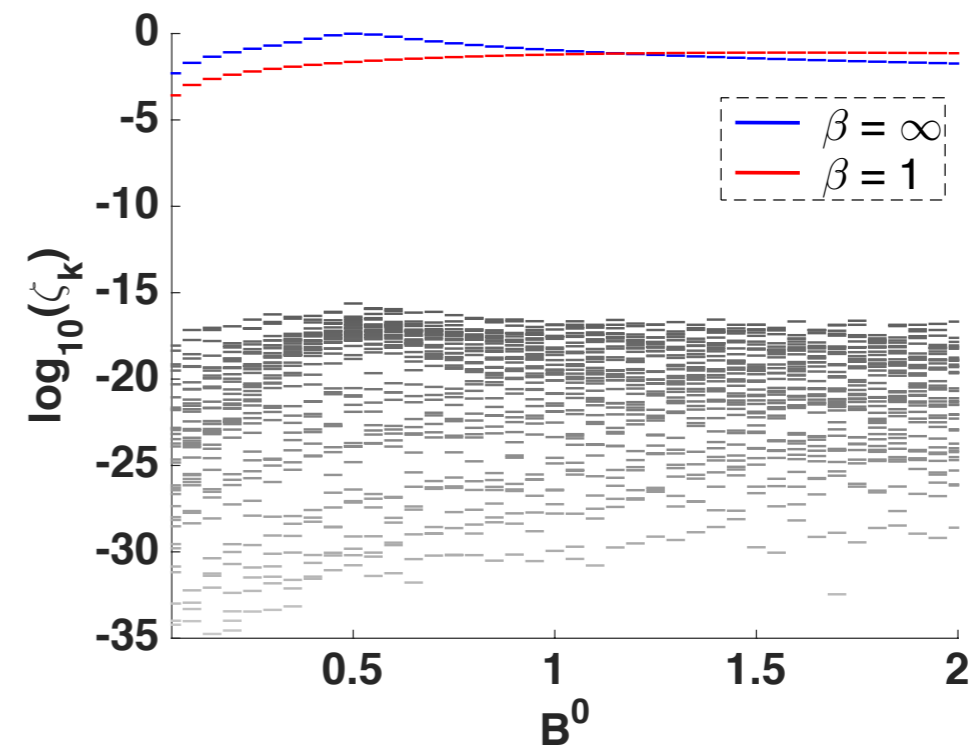
# Transverse-field Ising model *revisited*

Net magnetization observable



Rank(F)  $\leq 2$

Correlation function observable



Rank(F)  $\leq 1$

For any value of the model parameters, temperature, and any number of spins

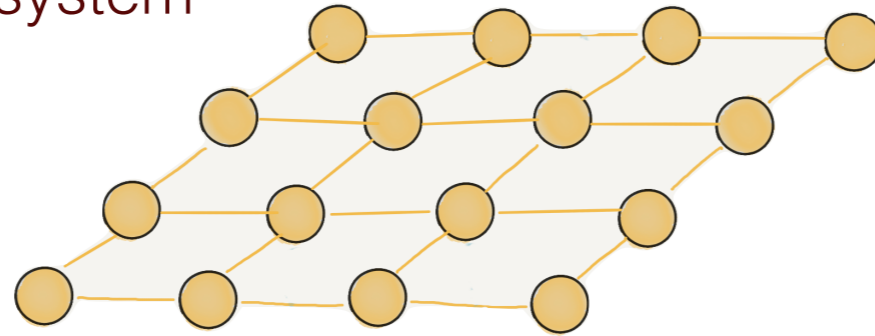
This is a very sloppy model, and hence potentially robustly simulatable.

# Outline

1. Parameter space compression and simulation sensitivity
2. Engineering thermalization in quantum many-body models

# Simulation of many-body systems

Many-body lattice system



Engineered Hamiltonian

$$H$$

How do we prepare a thermal state at temperature  $T$ ?  $\rho(\beta) = \frac{e^{-\beta H}}{\mathcal{Z}}$

The apparatus has a physical temperature, but this is always relevant.

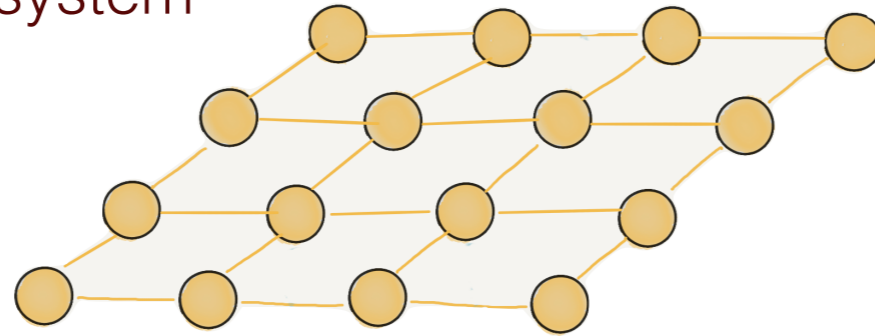
The errors/noise on the logical degrees of freedom may be non-equilibrating.

*e.g.* Spin lattice, and independent depolarizing channel on each spin

$$\rho(t) \rightarrow \frac{\mathbf{1}}{n}$$

# How to get thermalization?

Many-body lattice system



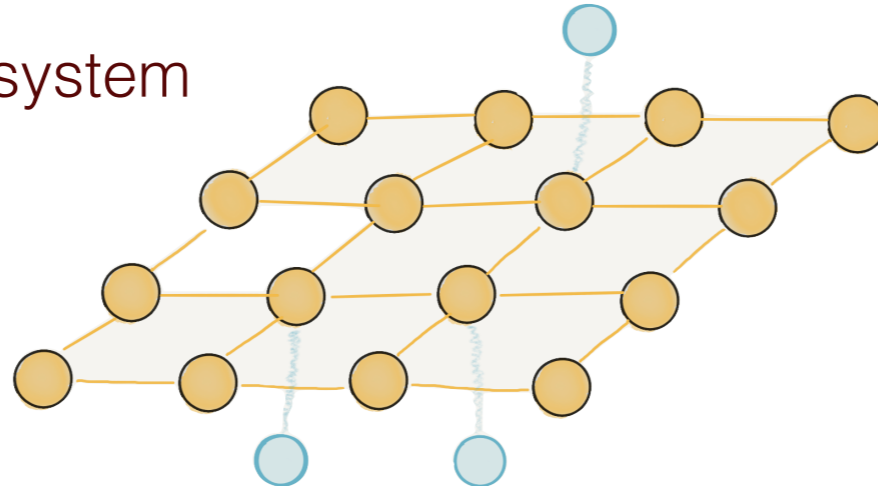
Engineered Hamiltonian

$H$

Need to *engineer* thermalization also

# How to get thermalization?

Many-body lattice system



Engineered Hamiltonian

$$H$$

Engineered dissipative evolution

$$\dot{\rho}(t) = \mathcal{L}\rho(t)$$

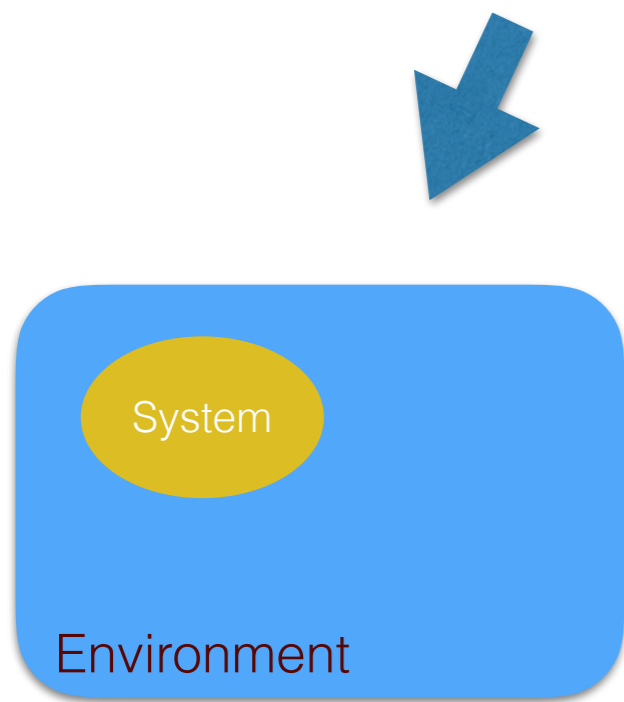
But what kind of evolution will result in a thermal state?

Only general characterization known [e.g. Breuer & Petruccione, The theory of open quantum systems]:

Born-Markov master equation + ergodicity + KMS conditions  $\Rightarrow$  thermal steady state

# How to get thermalization?

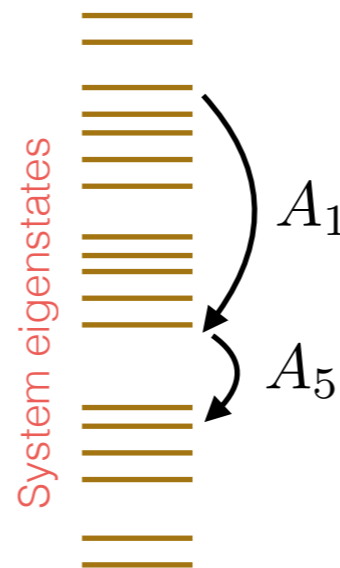
Born-Markov master equation + ergodicity + KMS conditions  $\Rightarrow$  thermal steady state



- Weak coupling to many degrees of freedom
- Fast relaxing environment

$$[A_\alpha, X] = [A_\alpha^\dagger, X] = 0$$

$$\Rightarrow X \propto 1$$



$$\frac{\gamma_{\alpha\beta}(-\omega)}{\gamma_{\beta\alpha}(\omega)} = e^{-\beta\omega}$$

Bath induced fluctuations and dissipation satisfy detailed balance

$$\frac{d\rho}{dt} = -i[H, \rho] + \sum_{\omega, \alpha, \beta} \gamma_{\alpha\beta}(\omega) \left( A_\beta(\omega) \rho A_\alpha^\dagger(\omega) - \frac{1}{2} \{ A_\alpha^\dagger(\omega) A_\beta(\omega) \} \right)$$

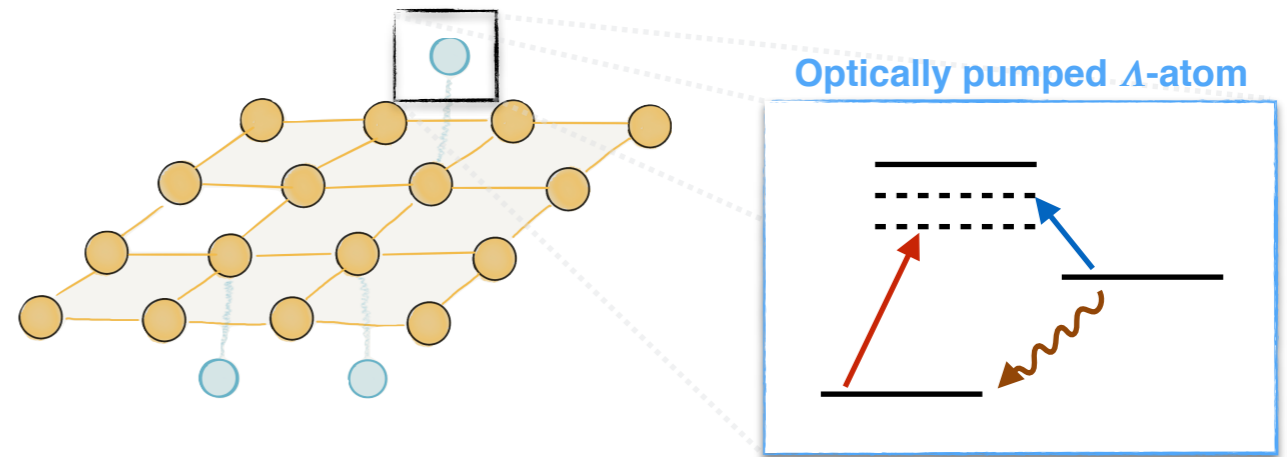
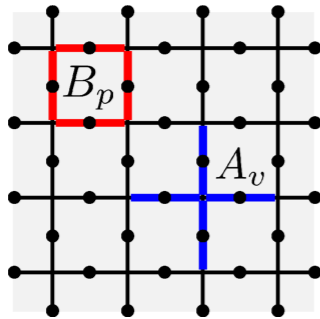
# How to get thermalization?

A constructive approach to thermalization for stabilizer Hamiltonians

K. Young, M.S. *et al.* J. Phys. B, **45**, 154012 (2012)

e.g. toric code

$$H_{TC} = -\lambda_e \sum_v A_v - \lambda_m \sum_p B_p$$



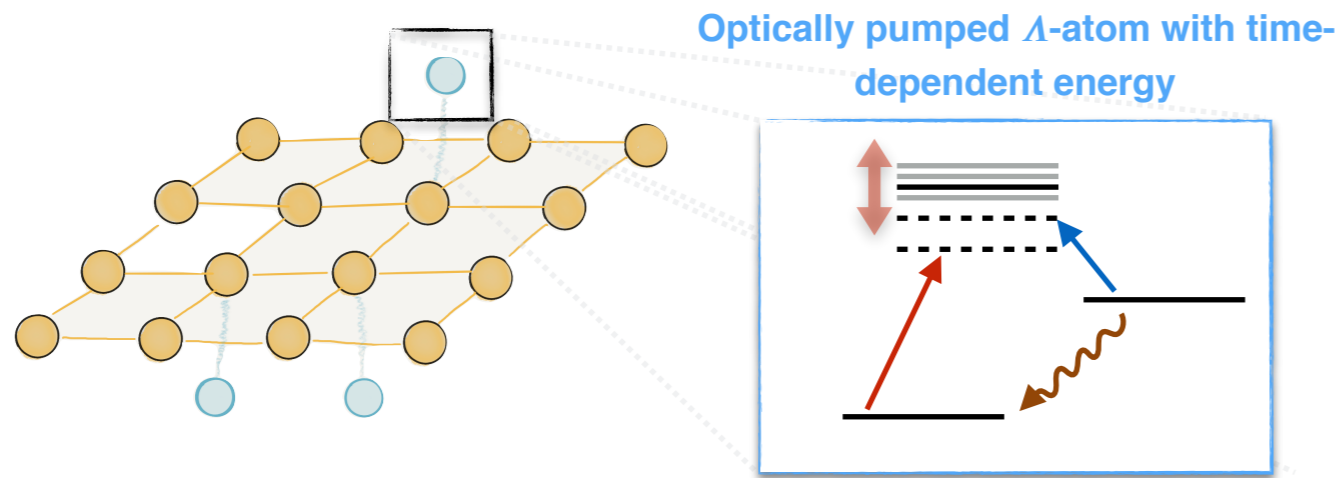
A number of special properties of stabilizer Hamiltonians make this work:

1. Easily predicted spectra  $\Rightarrow$  we can choose the energies of the ancillas to achieve resonant energy exchange
2. Local perturbations create energy excitations  $\Rightarrow$  so local couplings to ancilla move one up and down between eigenstates (ergodicity easy)

# How to get thermalization?

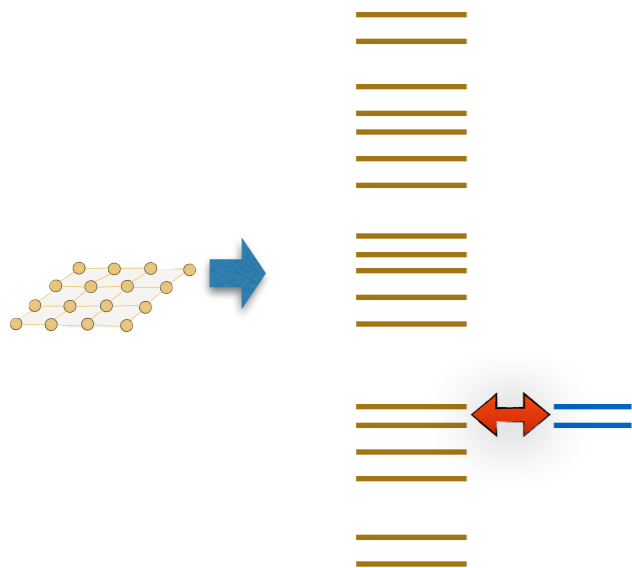
A generalization of this approach

M.S., Jonathan Moussa, C. Daniel Freeman



We assume ergodicity is satisfied by the system-ancilla couplings

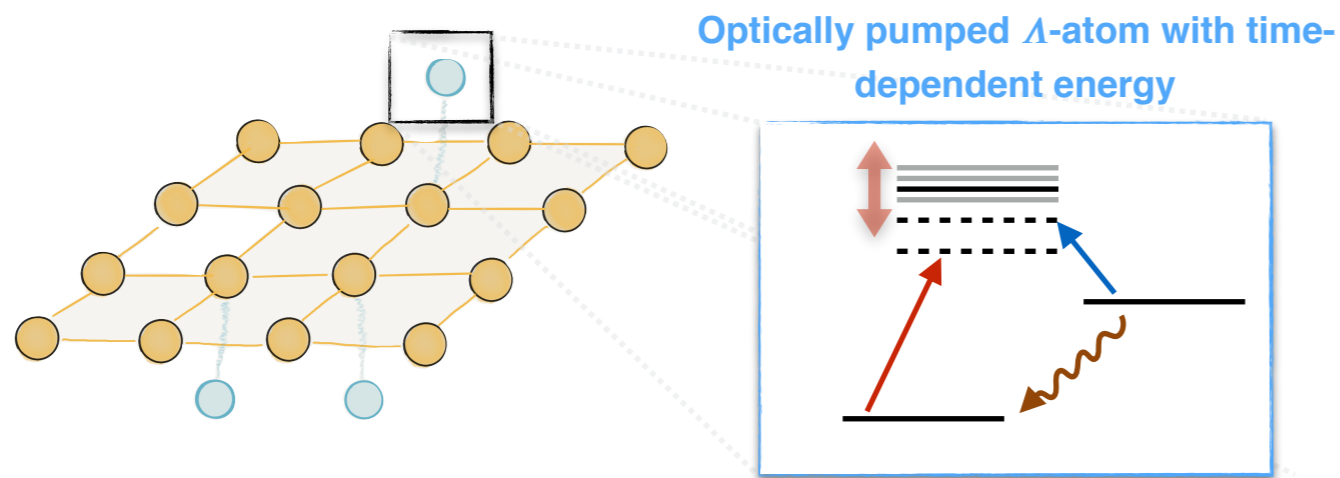
This mimics a macroscopic bath over some longer timescale



# How to get thermalization?

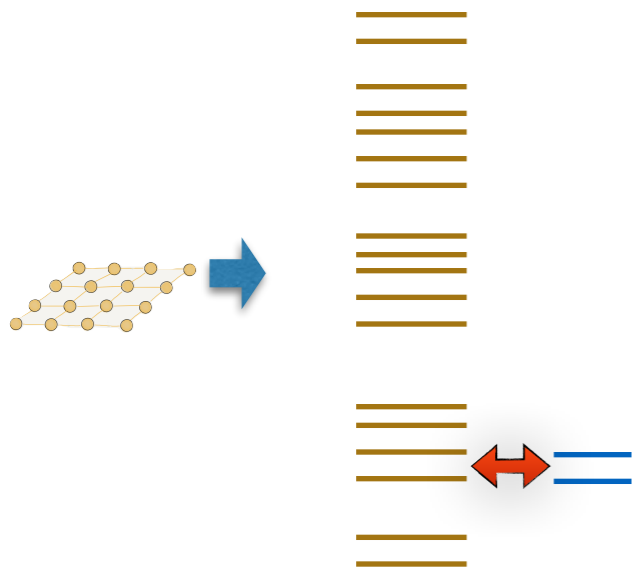
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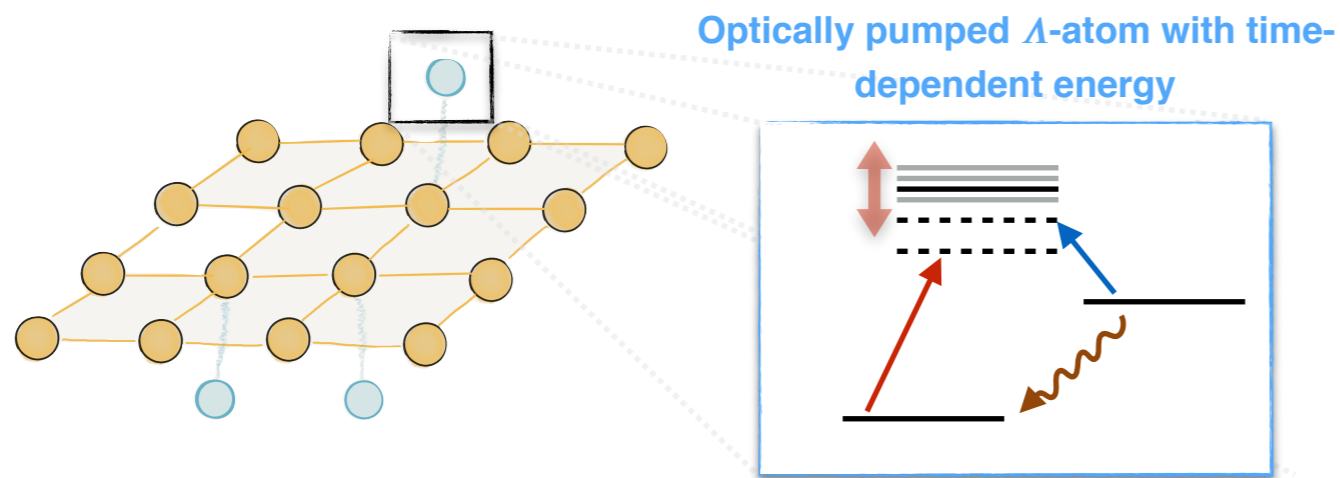
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# How to get thermalization?

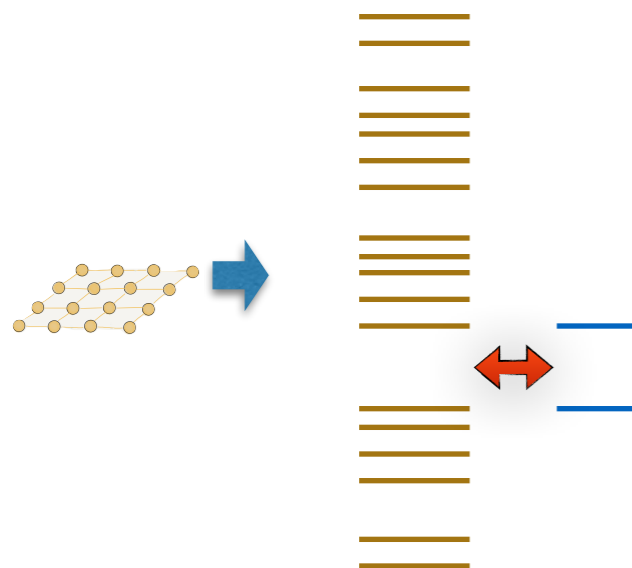
A generalization of this approach

M.S., Jonathan Moussa, C. Daniel Freeman



We assume ergodicity is satisfied by the system-ancilla couplings

This mimics a macroscopic bath over some longer timescale

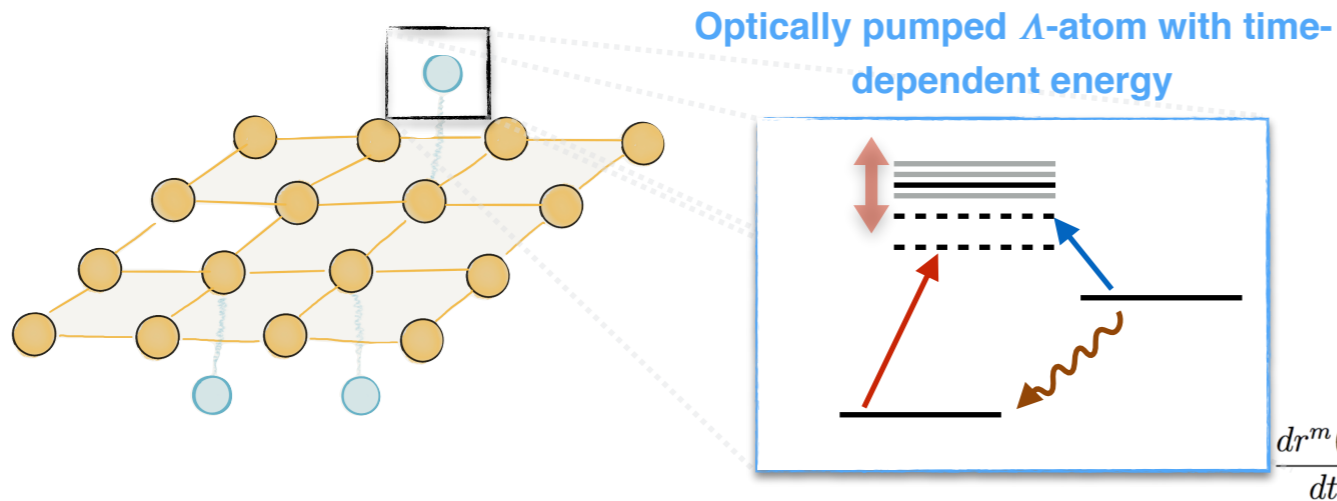


Ancillas will be resonant with different energy transitions in the system at different times, so as long as we maintain Boltzmann populations in the ancilla at all time, we will induce transitions that thermalize the system.

# How to get thermalization?

## Spin lattice example

M.S., Jonathan Moussa, C. Daniel Freeman



$$\frac{dr^m(t)}{dt} = \mathcal{L}_m(t)[r^m(t)] = \gamma_+^m(t)\mathcal{D}[\tau_+^m]r^m(t) + \gamma_-^m(t)\mathcal{D}[\tau_-^m]r^m(t),$$

$$H_T(t) = H_{\text{sys}} - \sum_{m=1}^M \frac{\Omega_m(t)}{2} \tau_z^m + \sum_{m=1}^M g_m (\sigma_x^{k_m} \tau_x^m),$$

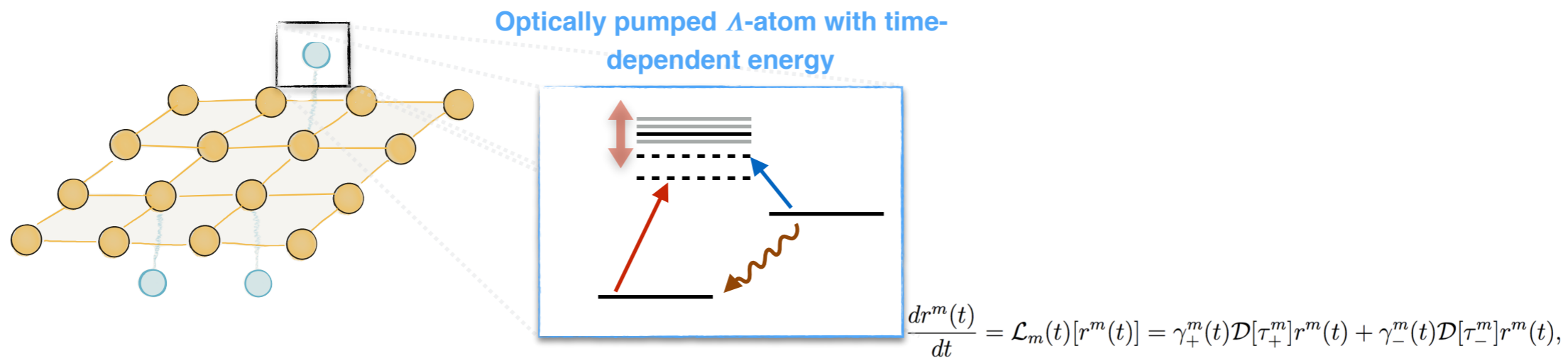
We require the parameter regime

$$\left| \frac{d\Omega_m(t)}{dt} \right| \ll g_m \sim \Gamma^m \ll \|H_{\text{sys}}\|, \quad \forall m, t$$

$$\Gamma^m \equiv \gamma_+^m + \gamma_-^m$$

# How to get thermalization?

## Spin lattice example



$$H_T(t) = H_{\text{sys}} - \sum_{m=1}^M \frac{\Omega_m(t)}{2} \tau_z^m + \sum_{m=1}^M g_m (\sigma_x^{k_m} \tau_x^m),$$

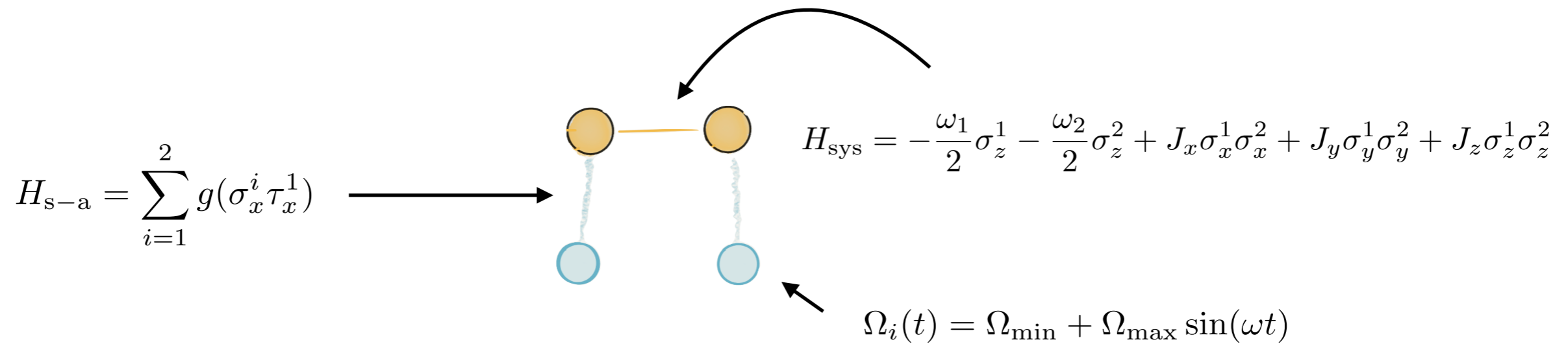
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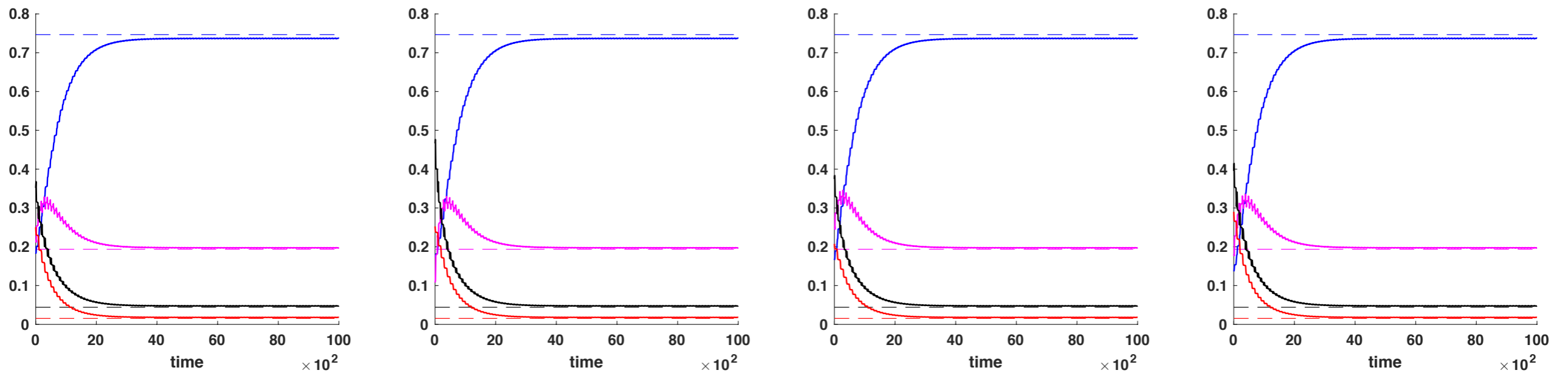
$$\Gamma^m \equiv \gamma_+^m + \gamma_-^m$$

In this regime we can derive a reduced master equation for the system alone that describes its evolution when the ancilla dynamics is averaged over.

# A simple example



## Thermalization from four random initial states



# Ongoing work

1. Bounds on thermalization time of this protocol
2. Comparison to discrete-time (gate-based) thermalization protocols (e.g. quantum Metropolis)

# Thanks!

## Collaborators

Reliability of analog  
quantum simulation

Jun Zhang, Lishan Zeng  
Shanghai Jiao Tong University

Engineered  
thermalization

Jonathan Moussa  
Sandia National Labs, Albuquerque

C. Daniel Freeman  
Univ. California, Berkeley



Sandia National Laboratories is a multimission laboratory managed and operated by National Technology and Engineering Solutions of Sandia, LLC., a wholly owned subsidiary of Honeywell International, Inc., for the U.S. Department of Energy's National Nuclear Security Administration under contract DE-NA-0003525.

Backup slides

# Interlude: parameter space compression

Sethna group, Cornell University

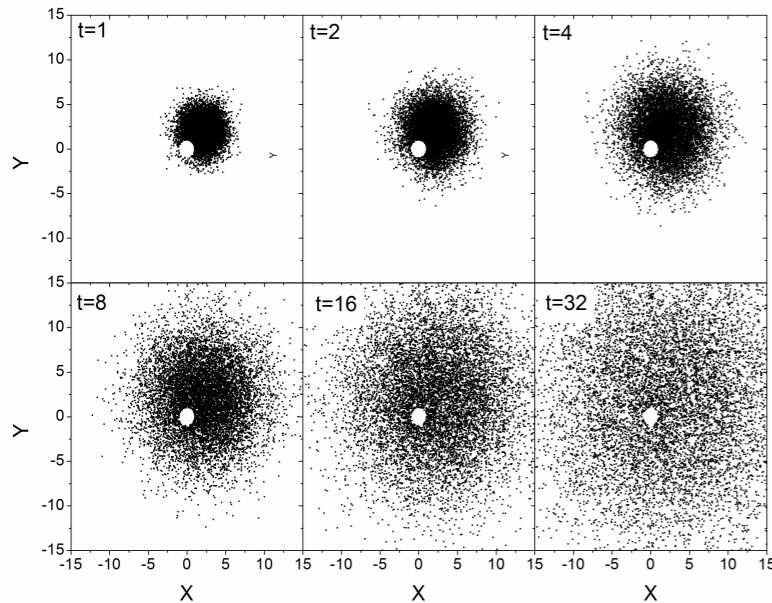
Why the unreasonable effectiveness of simple models in physics?



**Parameter Space Compression Underlies Emergent Theories and Predictive Models**

Benjamin B. Machta *et al.*  
*Science* **342**, 604 (2013);  
DOI: 10.1126/science.1238723

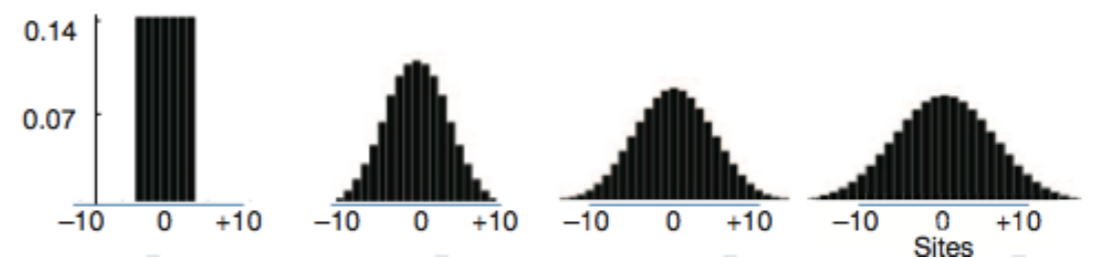
*e.g. particle diffusion*



Plante & Cucinotta

Equation for particle density

$$\frac{\partial}{\partial t}\rho(x, t) = D \frac{\partial^2}{\partial x^2}\rho(x, t) - V \frac{\partial}{\partial x}\rho(x, t) + R\rho(x, t)$$



Only 3 parameters!

# PSC and Fisher Information

## General setup

Model dependent on  $K$  parameters

$$\lambda = (\lambda_1, \lambda_2, \dots, \lambda_K)$$

Interested in some set of observables

$$\vec{x}$$

Parameterized distribution over observables

$$P_\lambda(\vec{x})$$

# PSC and Fisher Information

## General setup

Model dependent on  $K$  parameters

$$\lambda = (\lambda_1, \lambda_2, \dots, \lambda_K)$$

Interested in some set of observables

$$\vec{x}$$

Parameterized distribution over observables  $P_\lambda(\vec{x})$

## Fisher Information matrix (FIM)

$$F_{ij}(\lambda) = \sum_{\vec{x}} \frac{1}{P_\lambda(\vec{x})} \left( \frac{\partial P_\lambda(\vec{x})}{\partial \lambda_i} \right) \left( \frac{\partial P_\lambda(\vec{x})}{\partial \lambda_j} \right)$$

$$F = \sum_{k=1}^K \zeta_k v_k v_k^\dagger$$

### Key observation:

Eigenvalues of FIM prescribe an ordering of parameter influence, *i.e.*  
if  $\zeta_k$  is large, then the parameter

$$\sum_j \lambda_j v_k^j$$

has a large influence on observations

# PSC and Fisher Information

## General setup

Model dependent on  $K$  parameters

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### Key observation:

*Another way to think about this:*

Sensitivity analysis — FIM is the expected value of a Hessian:

$$F_{ij} = - \left\langle \frac{\partial^2 \log(p_m(\lambda))}{\partial \lambda_i \partial \lambda_j} \right\rangle$$

# Parameter space compression

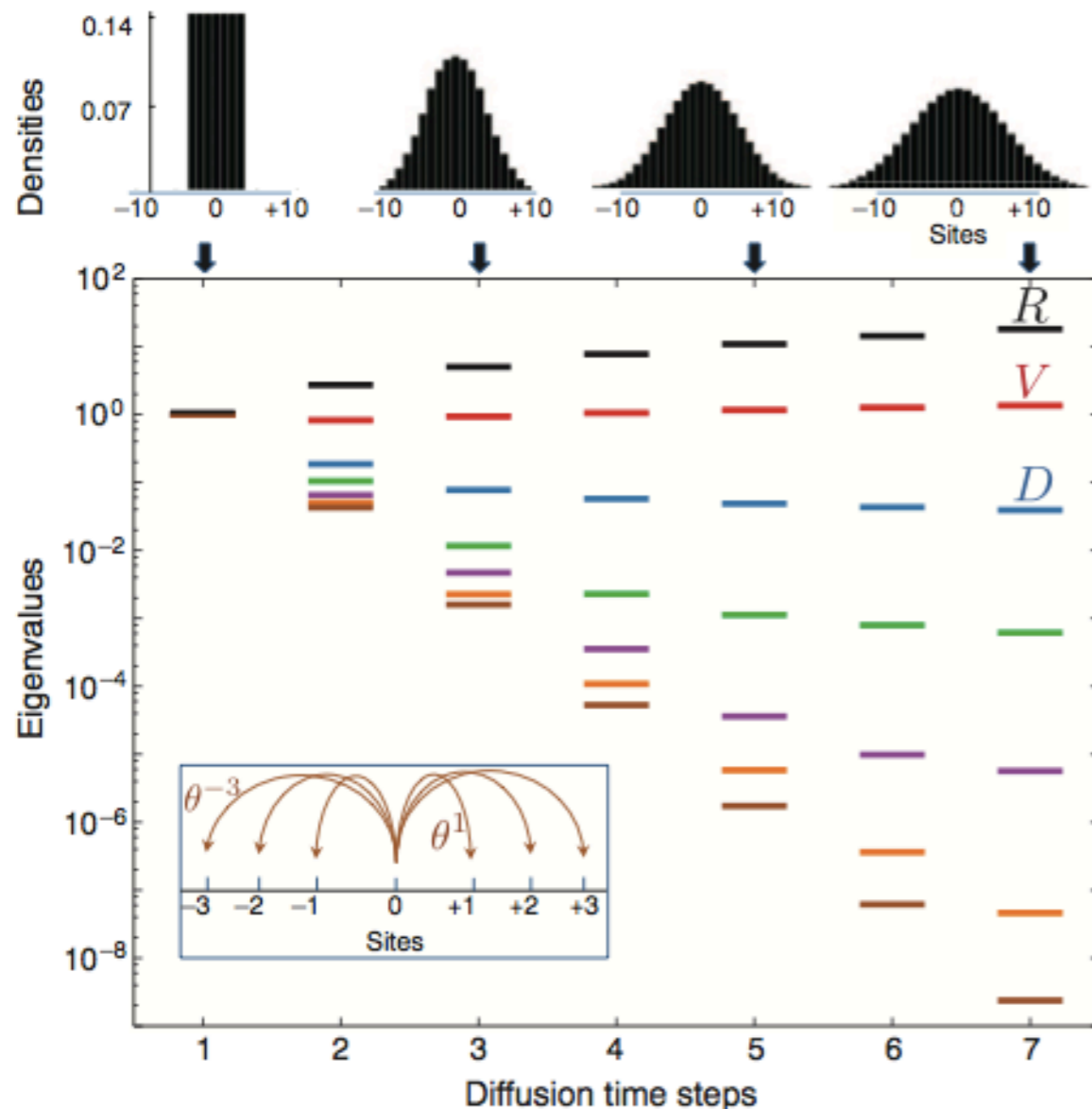
*e.g.* particle diffusion

Machta et al. Science, 342, 604 (2013)

Equation for particle density

$$\frac{\partial}{\partial t}\rho(x, t) = D\frac{\partial^2}{\partial x^2}\rho(x, t) - V\frac{\partial}{\partial x}\rho(x, t) + R\rho(x, t)$$

Discrete time/space diffusion in 1D



Observable here is the full particle density.

If we ask what the particle density is at every time step, then eigenvalues are clustered.

But if we only ask every few time steps, then eigenvalues separate, and a few important parameters emerge, and these are  $R$ ,  $V$ ,  $D$ .

# Parameter space compression

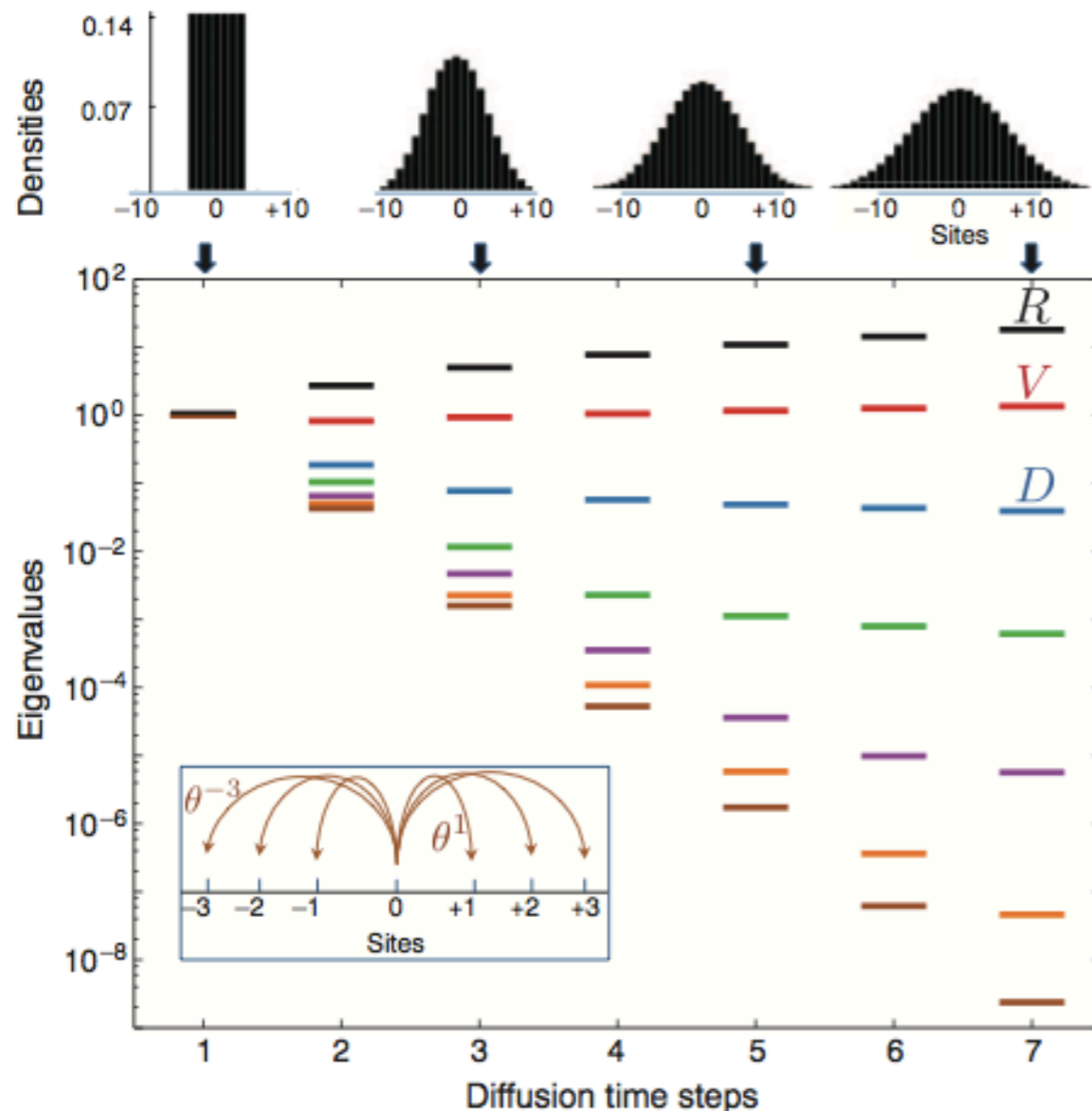
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## Some terminology

Stiff parameters

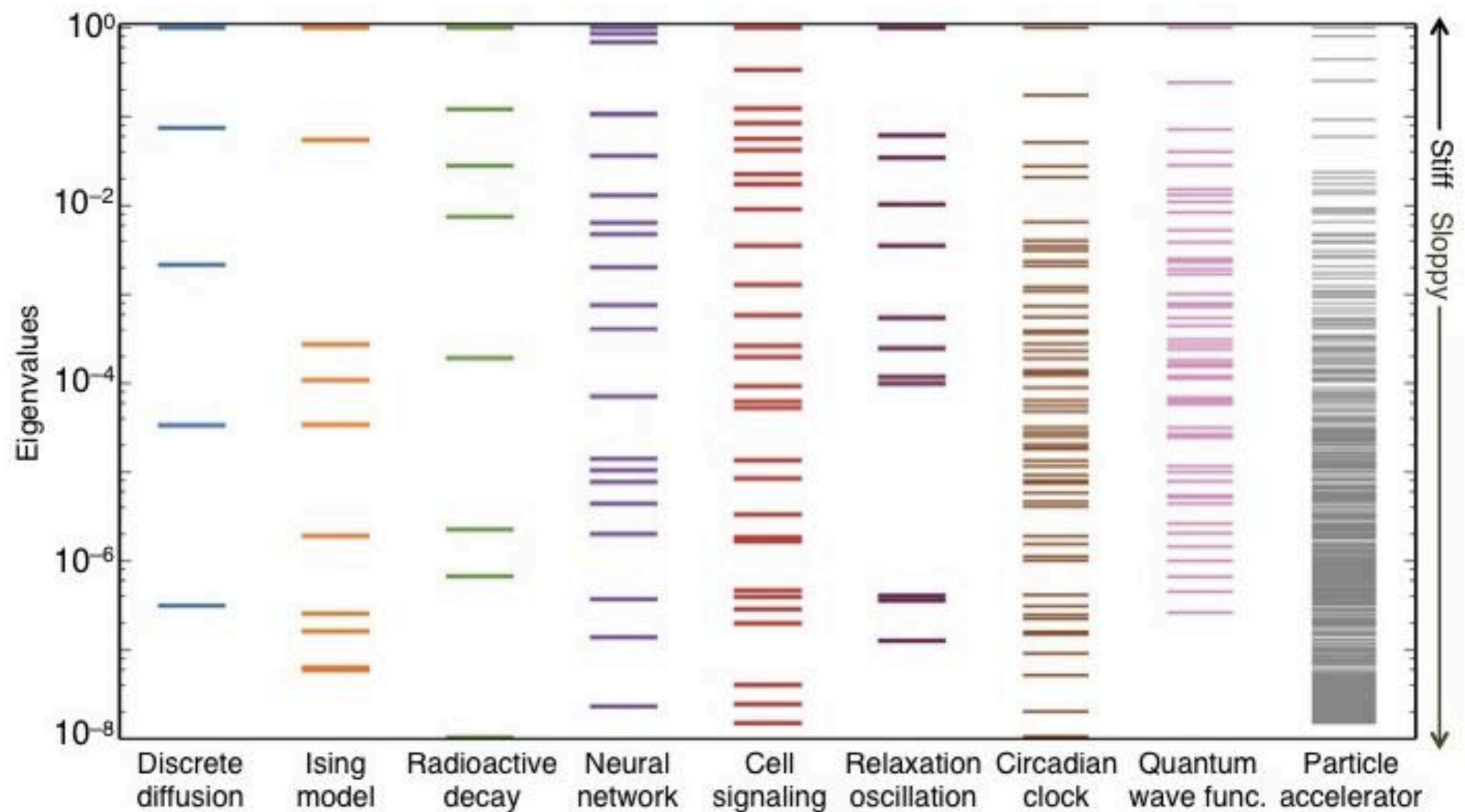
Sloppy parameters

Diffusion is a sloppy model

# Sloppiness abounds!

This phenomenon has been demonstrated for many models in physics and biophysics.

Sethna *et al.*: Models of nature tend to be sloppy, and this is why science is even possible



From the Sethna group website

# Scaling the analysis

The true value of a quantum simulator is to extract properties of models that are not classically tractable.

But the FIM can only be explicitly calculated for small versions of a model.

Strategies for scaling this analysis technique [[arXiv:1603.09283](#)]:

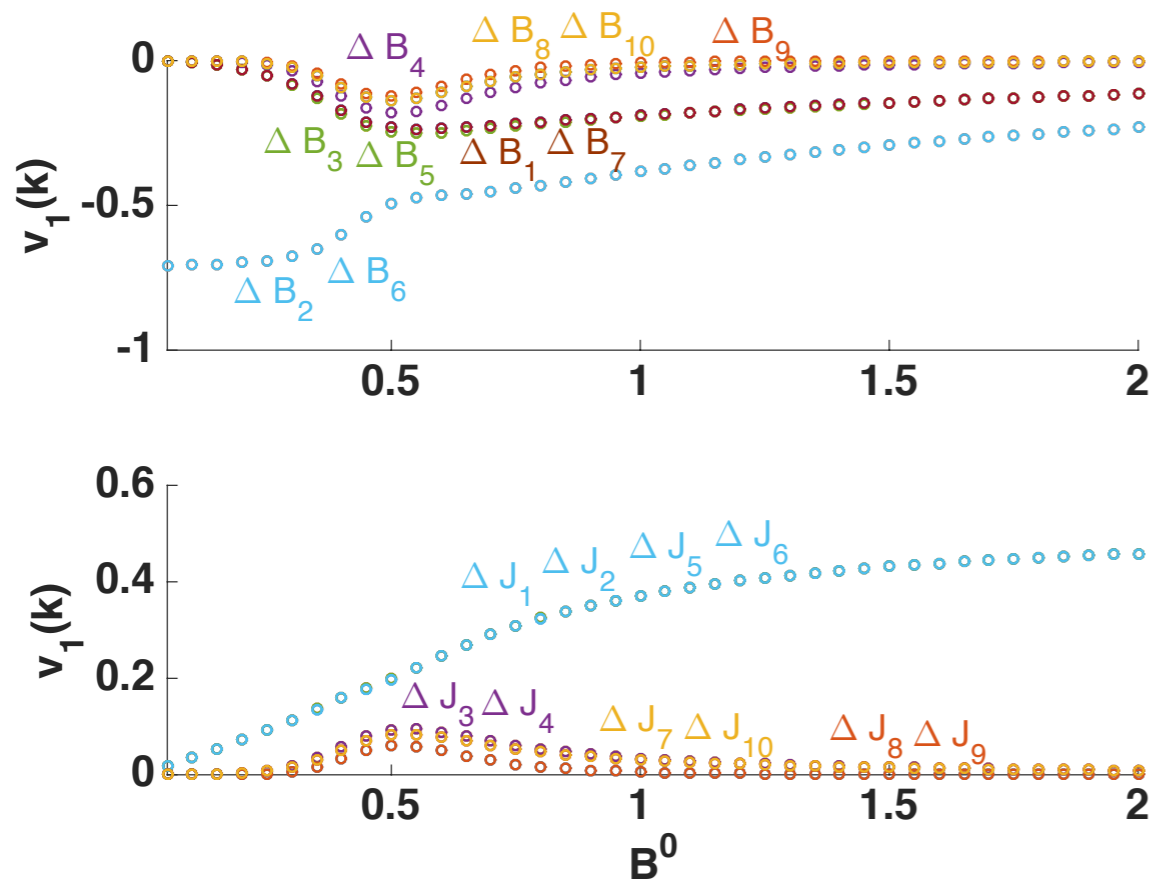
1. Symmetry analysis can be done analytically, independent of system size.  
*e.g.* translational invariance a powerful symmetry that implies sensitivity *only to* collective parameters for any model size.
2. Empirically, sloppiness and form of stiff parameters carries over from small-scale models to large-scale versions.  
*e.g.* TFIM with correlation function

# Transverse field Ising model

Correlation function observable

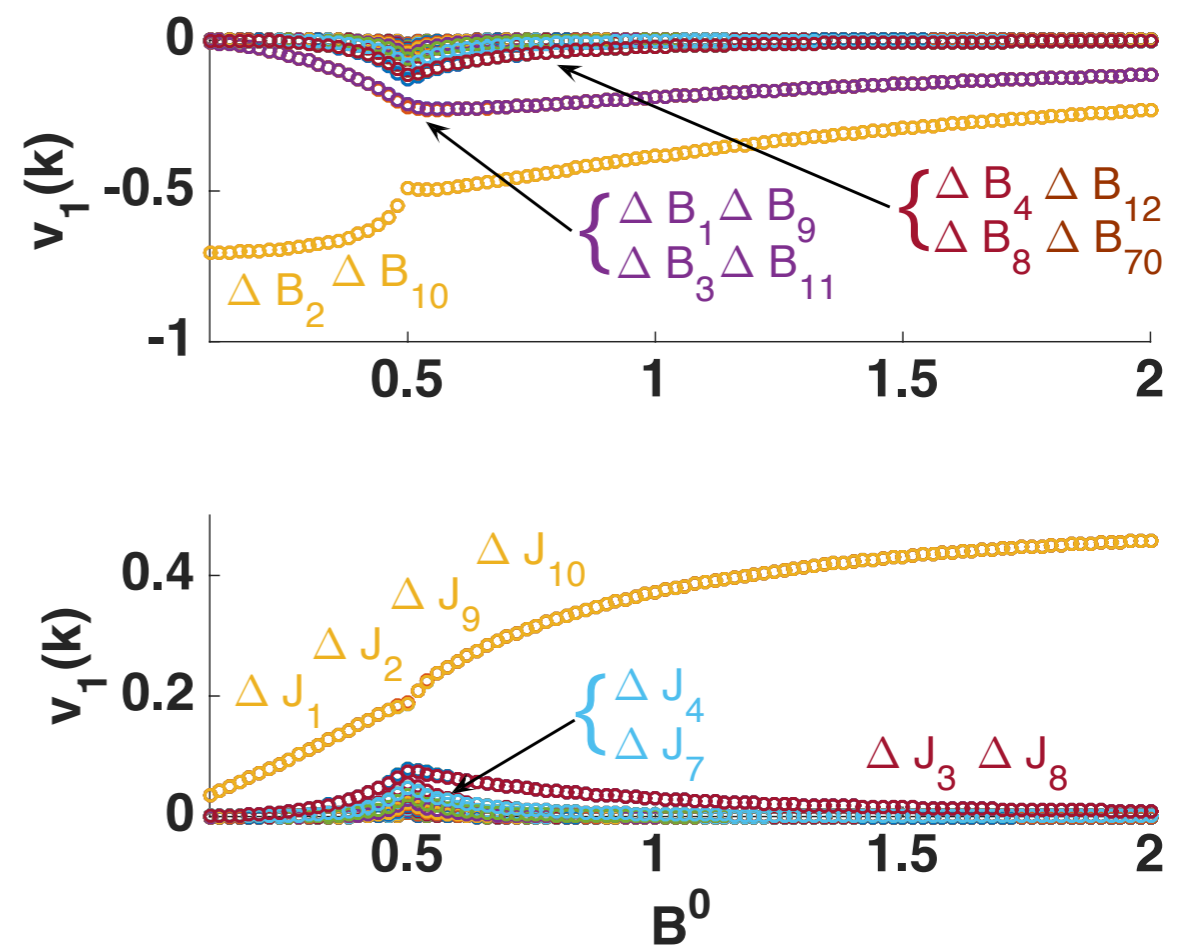
$$n = 10, J^0 = 1$$

$$C_z(2, 6)$$



$$n = 70, J^0 = 1$$

$$C_z(2, 10)$$



Qualitatively the same

# Scaling the analysis

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1. Symmetry analysis can be done analytically, independent of system size.  
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2. Empirically, sloppiness and form of stiff parameters carries over from small-scale models to large-scale versions.  
*e.g.* TFIM with correlation function
3. Conjecture:  
If a small-scale model is sloppy, then its large-scale version will also be sloppy.