

Variational counterdiabatic driving of the ferromagnetic p -spin model

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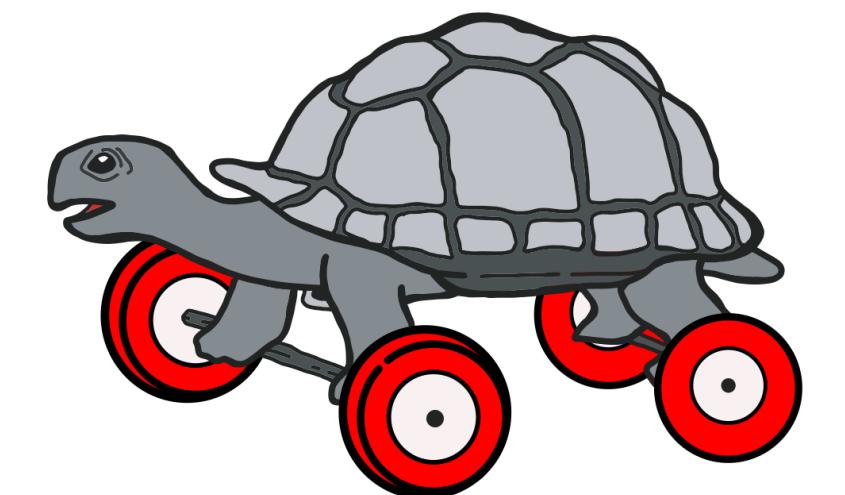
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Abstract

In **adiabatic quantum computation** (AQC), or quantum annealing (QA), the goal is to find the ground state of a many-body Hamiltonian H_p , starting from the simple ground state of a transverse field potential H_{TF} . A **fast** evolution with the time-dependent Hamiltonian $H_0(s) = (1-s)H_{\text{TF}} + sH_p$ renders the system robust against decoherence and thermal noise but can cause **diabatic transition** towards excited states. Adding a **counterdiabatic** (CD) potential H_{CD} to the original Hamiltonian can suppress Landau-Zener excitations and drive the system towards the correct target state, improving the efficiency of QA. The **variational approach** to CD driving can help us build approximate CD operators that are more viable experimentally than the exact CD potential.



Exact counterdiabatic driving

$$H_{\text{CD}}(t) = \dot{s}(t) \sum_{n \neq m} \frac{\langle E_n(s) | \partial_s H_0(s) | E_m(s) \rangle}{E_n(s) - E_m(s)} |E_n(s)\rangle \langle E_m(s)|$$

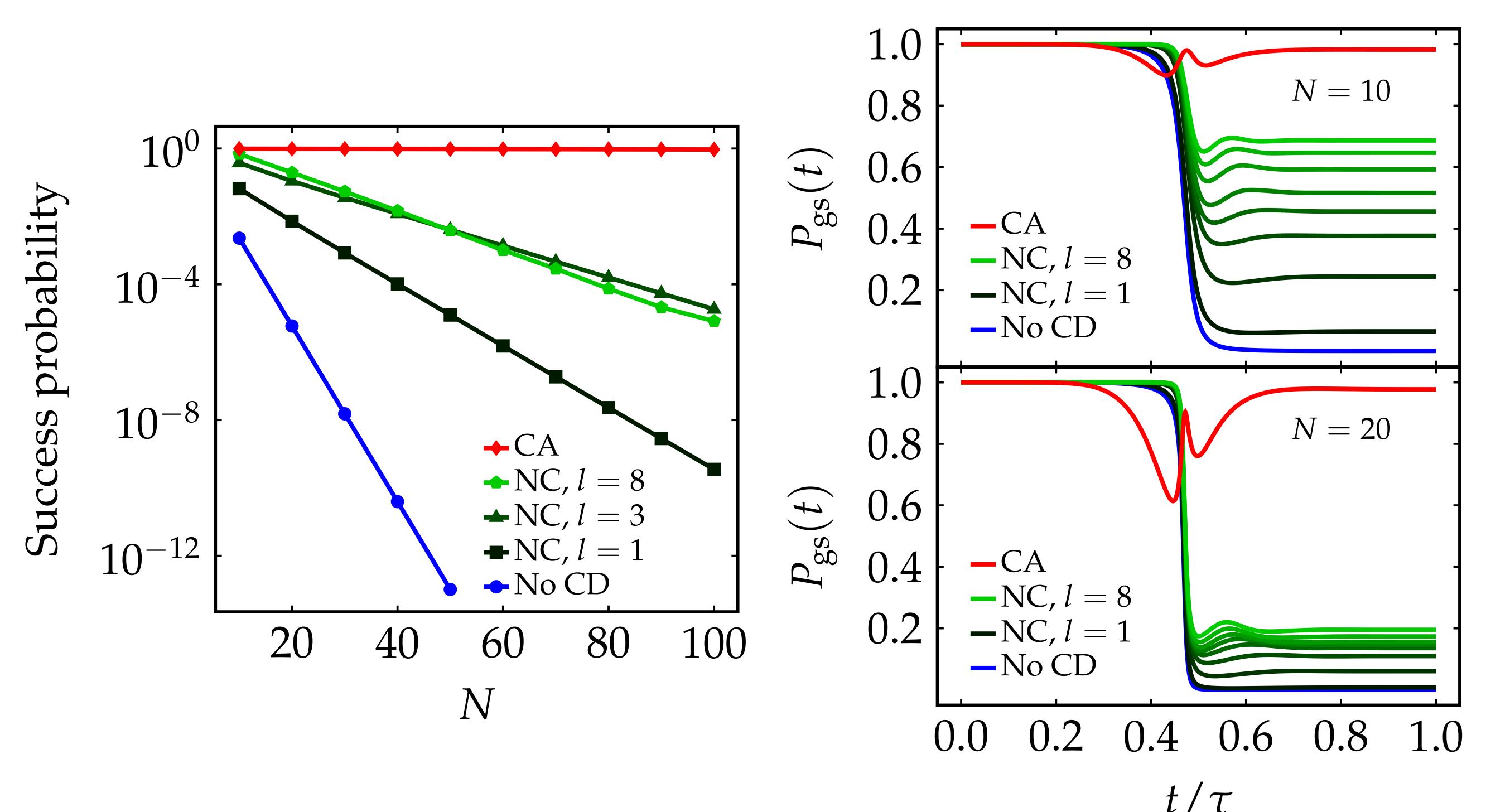
- Nonlocal
- Needs the exact spectrum in advance
- Poorly defined around Quantum Phase Transitions
- **Useless** for practical purposes!



M. V. Berry, J. Phys. A: Math. Theor. 42 (2009) 365303

Results for $p = 3$: the cyclic ansatz (CA)

$$\mathcal{L}_1 = S_y \quad \mathcal{L}_2 = S_y^3 \quad \mathcal{L}_3 = S_x S_y S_z + S_z S_y S_x$$



Variational approach

$$H_{\text{CD}}(t) = \dot{s}(t) A_s(t)$$

$$[\mathrm{i}\partial_s H_0 - [A_s, H_0], H_0] = 0$$

$$\mathcal{S}(A_s) = \text{tr}[G_s^2(A_s)]$$

$$G_s(A_s) = \partial_s H_0 + \mathrm{i}[A_s, H_0]$$

$$\frac{\delta \mathcal{S}(A_s)}{\delta A_s} = 0$$

Local ansatz

$$A_s(t) = \sum_{k=1}^l \alpha_k(t) \mathcal{L}_k$$

$$\mathcal{L}_k \in \{ \{ \sigma_i^\alpha \}, \{ \sigma_i^\alpha \sigma_j^\beta \}, \dots \}$$

$$\alpha, \beta \in \{x, y, z\} \quad i, j \in \{1, \dots, N\}$$

D. Sels, A. Polkovnikov, PNAS May 16, 2017 114 (20)

Nested commutator (NC)

$$A_s^{(l)}(t) = \mathrm{i} \sum_{k=1}^l \alpha_k(t) \mathcal{O}_{2k-1}(t)$$

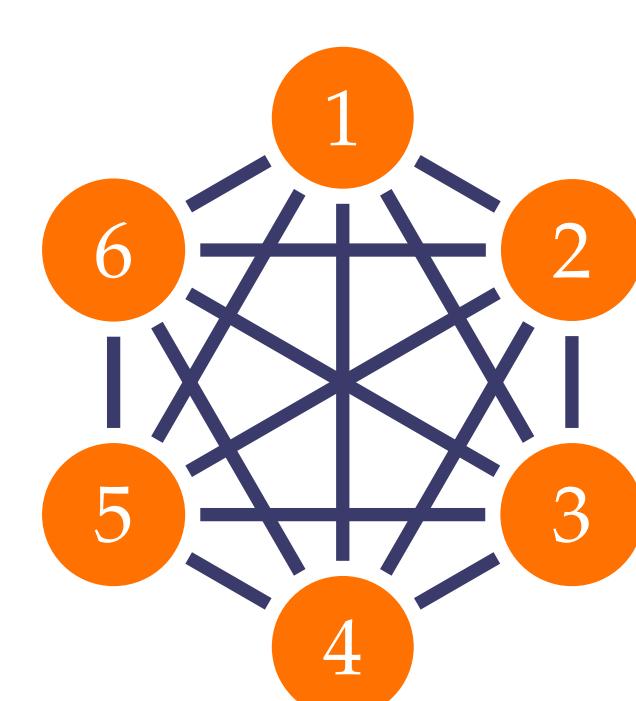
$$\mathcal{O}_k = \begin{cases} \partial_s H_0 & k=0 \\ [H_0, \mathcal{O}_{k-1}] & k>0 \end{cases}$$

P. W. Claeys et al., Phys. Rev. Lett. 123, 090602 (2019)

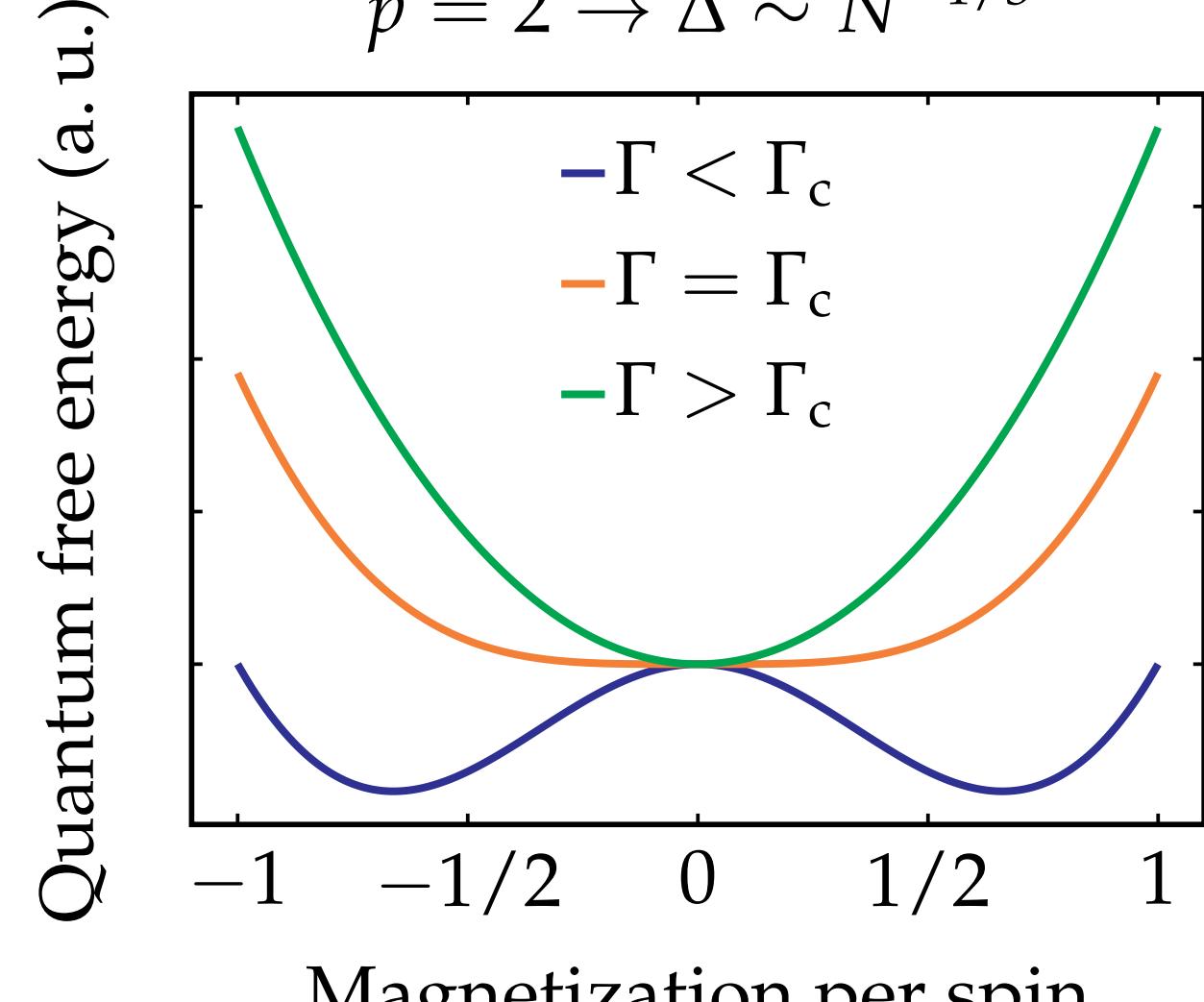
Application: ferromagnetic p -spin model

$$H_{\text{TF}} = -\Gamma \sum_{i=1}^N \sigma_i^x$$

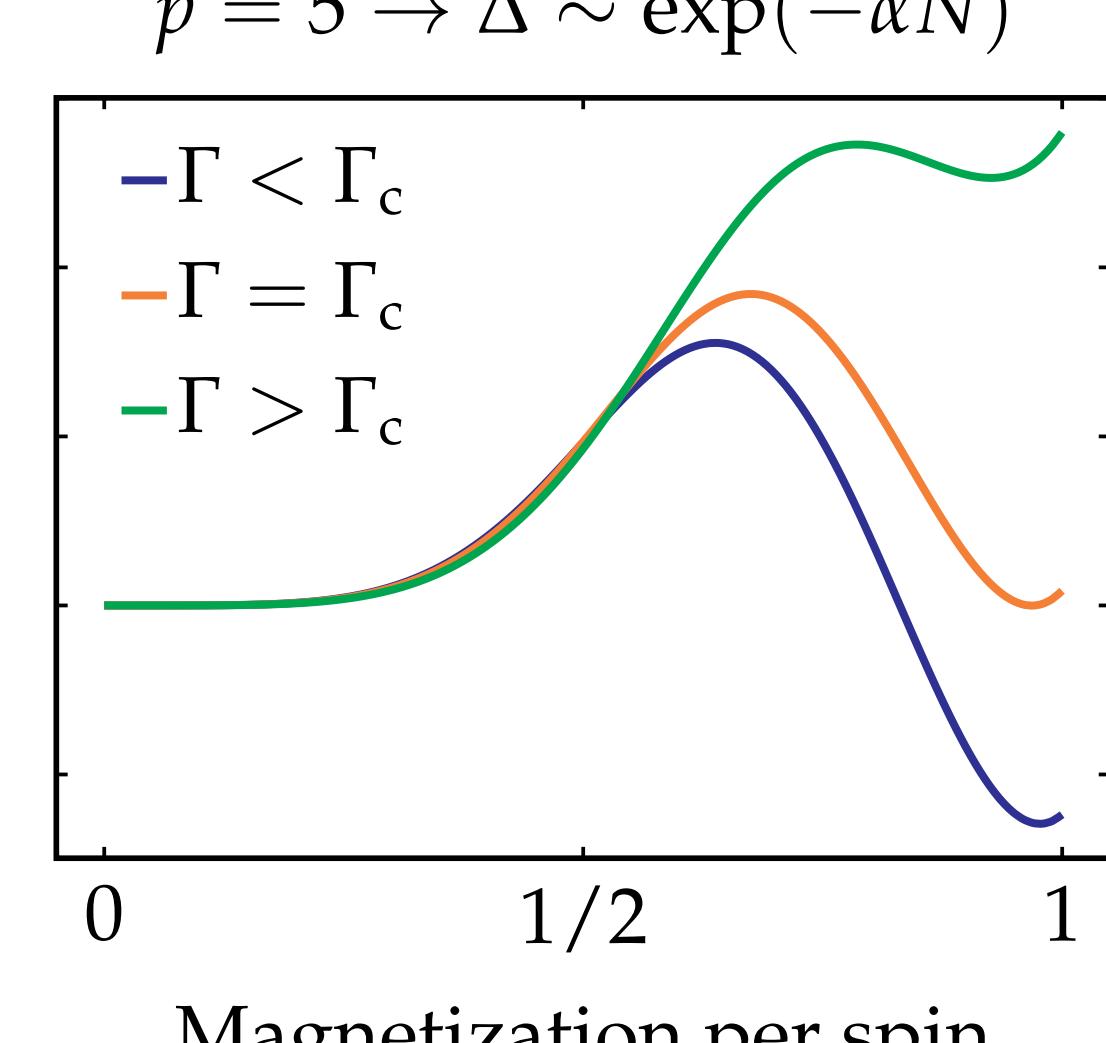
$$H_p = -JN \left(\frac{1}{N} \sum_{i=1}^N \sigma_i^z \right)^p \quad p \geq 2$$



$$p=2 \rightarrow \Delta \sim N^{-1/3}$$



$$p=5 \rightarrow \Delta \sim \exp(-\alpha N)$$

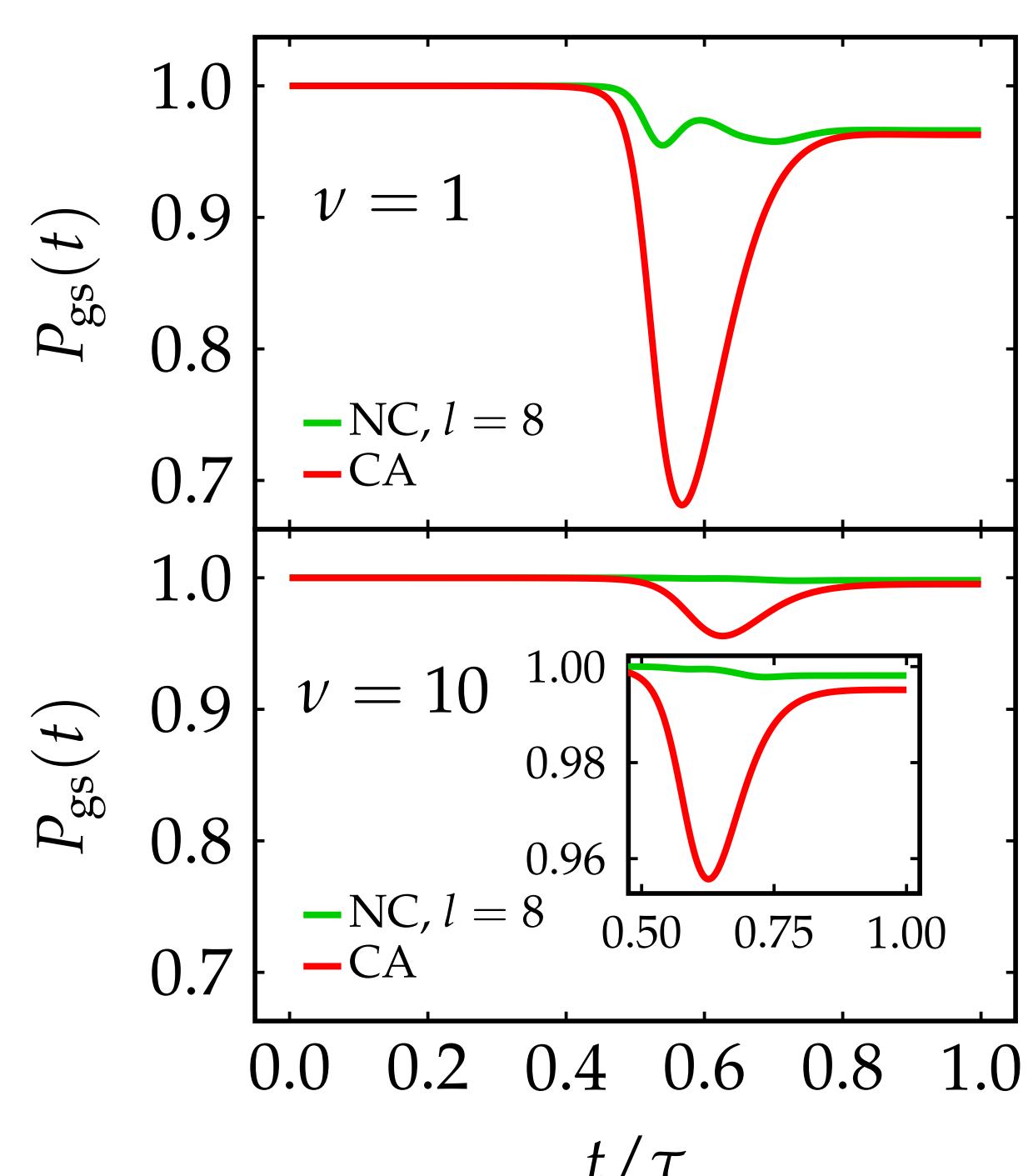


Finite-range and random models

Finite-range p -spin model

$$H_p = -\frac{1}{N^{p-1}} \sum_{i_1, \dots, i_p} \frac{J}{d_{i_1, \dots, i_p}^v} \sigma_{i_1}^z \dots \sigma_{i_p}^z$$

$$v \begin{cases} = 0 & \text{infinite-range} \\ > 0 & \text{finite-range} \end{cases}$$

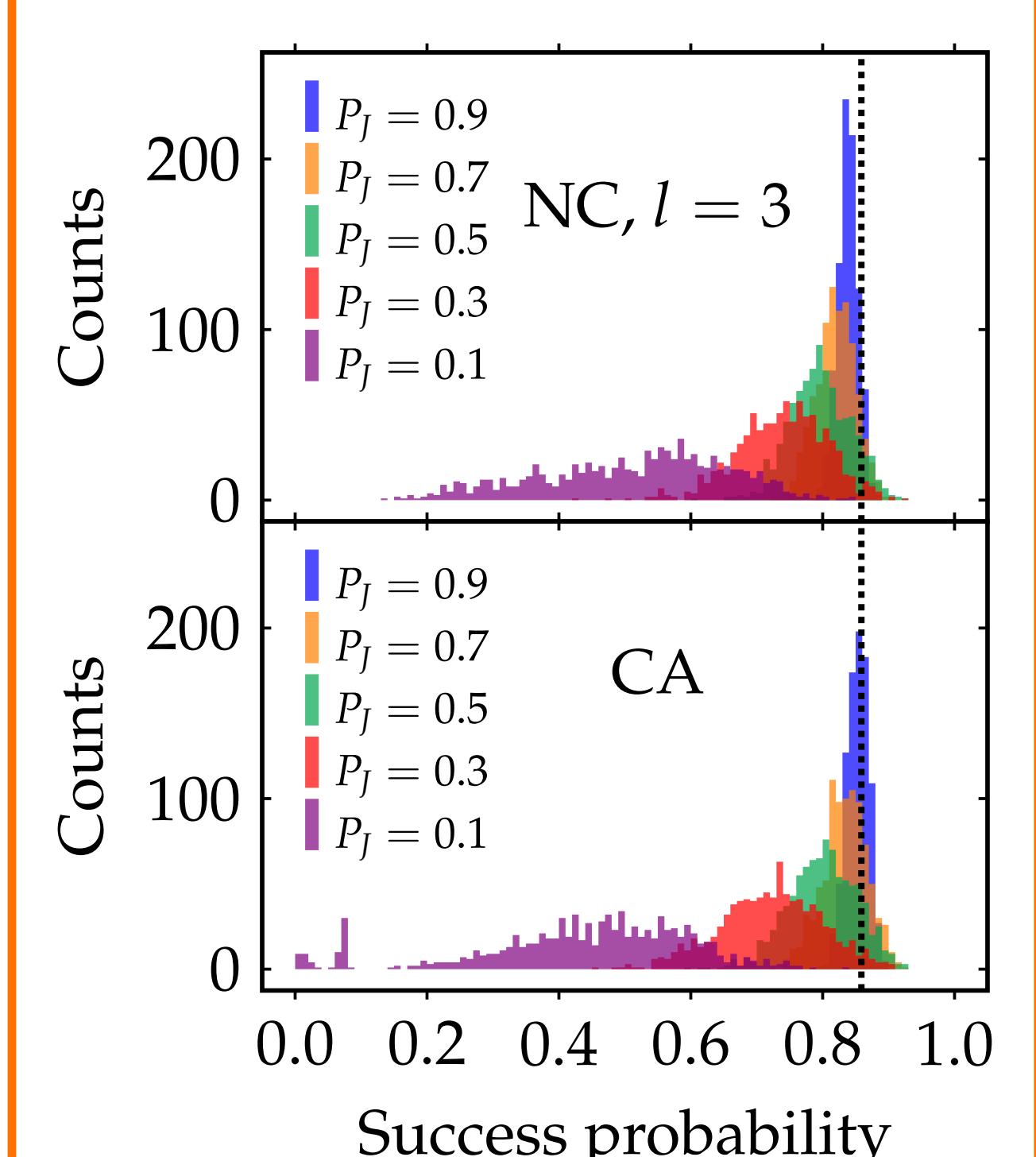


G. Passarelli et al., Phys. Rev. Research 2, 013283 (2020)

Random model

$$J \rightarrow K = \begin{cases} J & \text{probability } P_J \\ 0 & \text{probability } 1 - P_J \end{cases}$$

$P_J = 1$: standard model (dashed line)



Conclusions

- The NC ansatz can improve the annealing performance of the p -spin model
- Its efficiency is bound to decrease in the thermodynamic limit
- For $p = 3$, the **cyclic ansatz** is a viable alternative
- The CA yields optimal fidelity **independently of the system size**
- Analyzing **finite-range** and **random** models can provide some insights into the efficiency of the CA

