

Quantum simulation without ancillae

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1. Motivation: what's the matter with ancillae?
2. Quantum simulation without ancillae
 - 2.1. Simulating time evolutions
 - 2.2. Simulating ground-state subspaces
3. Conclusions and open questions



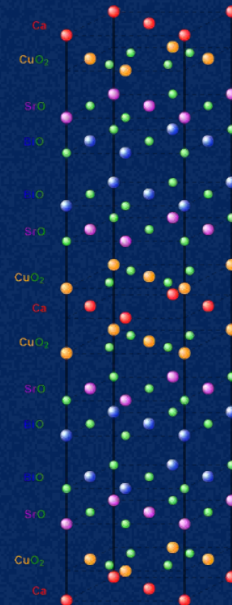
Motivation



Many-body systems described by k -local Hamiltonians:

$$H = \sum_i h_i \in \mathcal{B}((\mathbb{C}^d)^{\otimes n})$$

each term h_i acts nontrivially at most on k of these hilbert spaces



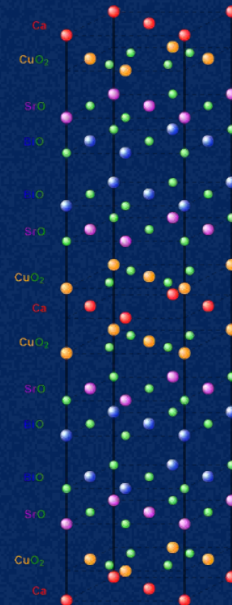


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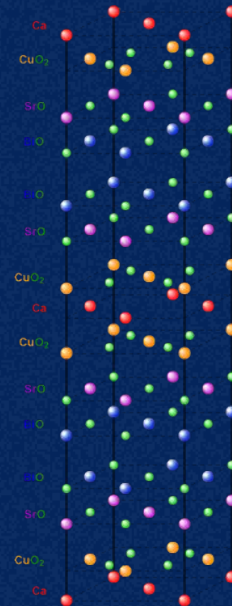


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The larger the k , the more difficult it is to realize the Hamiltonian experimentally.





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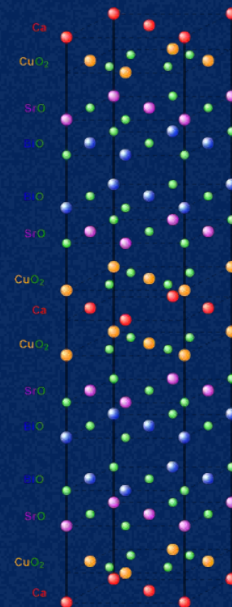
$$H = \sum_i h_i \in \mathcal{B}((\mathbb{C}^d)^{\otimes n})$$



The larger the k , the more difficult it is to realize the Hamiltonian experimentally.



So if you want to measure the properties of such a complex system in the lab, you first need to find a Hamiltonian with a smaller k that describes the same physics, i.e., you need to find the Hamiltonian of your quantum simulator.



Universal quantum Hamiltonians

Toby S. Cubitt , Ashley Montanaro , and Stephen Piddock [Authors Info & Affiliations](#)

Edited by Peter W. Shor, Massachusetts Institute of Technology, Cambridge, MA, and approved August 3, 2018 (received for review March 23, 2018)

August 30, 2018 | 115 (38) 9497-9502 | <https://doi.org/10.1073/pnas.1804949115>

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- H_{QS} reproduces the whole physics of H_T .
- $\dim(H_{QS}) > \dim(H_T)$

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Are there k' -local H_{QS} that reproduce specific physical aspects of a target system such that $k' < k$ and $\dim(H_{QS}) = d'^{n'}$, $\dim(H_T) = d^n$, with $d' = d$ and $n = n'$?

= Can we do quantum simulation without ancillae?

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Time evolution

Ground-state subspace

= Can we do quantum simulation without ancillae?

The background of the slide features a 3D visualization of a wavy, blue surface, possibly representing a potential energy landscape or a simulated time evolution. The surface is covered with numerous small, brown, spherical particles. The overall color scheme is a deep blue with a subtle gradient.

Simulating time evolutions

Simplifying the simulation of local Hamiltonian dynamics

Ayaka Usui, Anna Sanpera, and María García Díaz

Phys. Rev. Research **6**, 023243 – Published 4 June 2024



Ayaka Usui,
UAB



Anna Sanpera,
UAB, ICREA

Definition 1. A Hamiltonian H_{QS} ϵ -simulates a target Hamiltonian H_T at state $|\psi\rangle$ and time t if

$$|\langle\psi|e^{itH_{QS}}e^{-itH_T}|\psi\rangle| \geq 1 - \epsilon, \quad (1)$$

with $\epsilon \in (0, 1]$.

a) Exact simulation ($\epsilon = 0$)

$$[H_T, H_{QS}] = 0$$

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If $h := H_{QS} - H_T$ is degenerate in levels
 $L = \{|\ell_1\rangle, |\ell_2\rangle, \dots, |\ell_m\rangle\}$ and $|\psi\rangle \in \text{span}(L)$,
nontrivial simulation is feasible at all times t .

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Simulating the Same Physics with Two Distinct Hamiltonians

Karol Gietka, Ayaka Usui, Jianqiao Deng, and Thomas Busch
Phys. Rev. Lett. **126**, 160402 – Published 22 April 2021

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$$k = 3, k' = 2$$

$$H_T = \sum_{j=1}^4 (J_3 \sigma_z^j \sigma_z^{j+1} \sigma_z^{j+2} + h_x \sigma_x^j)$$

$$H_{QS} = \sum_{j=1}^4 (J_x \sigma_x^j \sigma_x^{j+1} + J_y \sigma_y^j \sigma_y^{j+1} + J_z \sigma_z^j \sigma_z^{j+1})$$

$$(J_x = J_y = J_z)$$

Heisenberg model

b) Approximate simulation ($\epsilon \neq 0$)

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Theorem: Every H_{QS} ϵ^* -simulates any H_T
at any $|\psi\rangle, t$ with

$$\epsilon^* = \min \left[1, \frac{t \Delta_h}{2} \right]$$

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spectral diameter of h
(expressible by an SDP)



Simulating ground-state subspaces

(soon on arXiv!)



Simulating ground-state subspaces

Given H_T :

- k -local
- $\dim(H_T) = d^n$
- ground-state subspace $\mathcal{S} = \text{span}\{|\psi_i\rangle\}_{i=1}^M$



Simulating ground-state subspaces

Given H_T :

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We want H_{QS} :

- k' -local, $k' < k$
- $\dim(H_{QS}) = d'^{n'}$, $d'=d$, $n'=n$
- ground-state subspace $\mathcal{S}' = \mathcal{S}$



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- ground-state subspace $\mathcal{S}' = \mathcal{S}$

Does it exist?

If so, how is it built?



Simulating ground-state subspaces

a) Semidefinite program:

$$\begin{array}{ll} \text{maximize} & \delta \\ \delta \in \mathbb{R}, \alpha \in \mathbb{R}^{L+1} & \\ \text{subject to} & PH_{QS}P = 0 \\ & PH_{QS}Q = 0 \\ & QH_{QS}Q \succeq \delta Q \\ & H_{QS} = \alpha_0 1_{d^n} + \sum_{j=1}^L \alpha_j \Lambda_j^{(k')} \\ & a 1_{d^n} \preceq H_{QS} \preceq b 1_{d^n} \end{array}$$



Simulating ground-state subspaces

a) Semidefinite program:

maximize δ
 $\delta \in \mathbb{R}, \alpha \in \mathbb{R}^{L+1}$
subject to

$$PH_{QS}P = 0$$
$$PH_{QS}Q = 0$$
$$QH_{QS}Q \succeq \delta Q$$
$$Q = 1_{d^n} - P$$
$$P = \sum_{i=1}^M |\psi_i\rangle\langle\psi_i|$$
$$H_{QS} = \alpha_0 1_{d^n} + \sum_{j=1}^L \alpha_j \Lambda_j^{(k')}$$
$$a 1_{d^n} \preceq H_{QS} \preceq b 1_{d^n}$$

generators of ℓ -body interactions,
 $\ell = 1, 2, \dots, k'$

allowed energy range



Simulating ground-state subspaces

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$\rightarrow \delta^* = 0 : \nexists k'$ -local parent to \mathcal{S}

$\rightarrow \delta^* \neq 0 : H_{QS}^* \text{ } k'$ -local parent to \mathcal{S} with largest spectral gap



Simulating ground-state subspaces

b) Simple algebraic method (built upon

Method to identify parent Hamiltonians for trial states)

[Martin Greiter](#), [Vera Schnells](#), and [Ronny Thomale](#)

Physical Review B



Simulating ground-state subspaces

b) Simple algebraic method:

1) Find α such that

$$\left(\sum_{j=1}^L \alpha_j \Lambda_j^{(k')} \right) |\psi_i\rangle = -\alpha_0 |\psi_i\rangle, \quad \forall i = 1, \dots, M.$$



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Equivalent to finding α such that

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Equivalent to finding α such that

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$$C_{\ell j}^{(i)} := \langle \psi_i | \Lambda_{\ell}^{(k')} \Lambda_j^{(k')} | \psi_i \rangle$$

$$\Lambda_0^{(k')} := \mathbf{1}_{d^n}$$



Simulating ground-state subspaces

b) Simple algebraic method:

2) Pick some $\alpha \in \mathcal{K}$ and build H_{QS} :

$$H_{QS} = \sum_{j=1}^L \alpha_j \Lambda_j^{(k')}$$



Simulating ground-state subspaces

b) Simple algebraic method:

2) Pick some $\alpha \in \mathcal{K}$ and build H_{QS} :

$$H_{QS} = \sum_{j=1}^L \alpha_j \Lambda_j^{(k')}$$

3) Diagonalize H_{QS} . It will be a suitable parent Hamiltonian if

$$-\alpha_0 \text{ is the ground energy and } \mathcal{S}' = \mathcal{S}$$



Simulating ground-state subspaces

b) Simple algebraic method:

-Cons: Diagonalization needed!



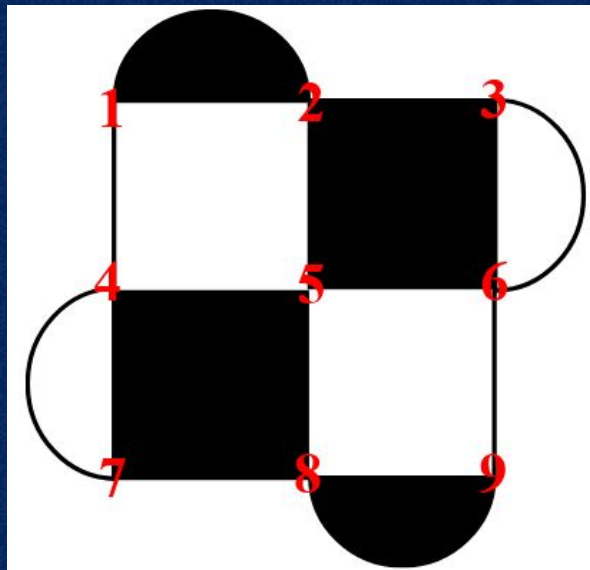
Simulating ground-state subspaces

b) Simple algebraic method:

-**Cons:** Diagonalization needed!

-**Pros:** Allows to check whether the target subspace appears as an excited eigenspace of some H_{QS} .

Example: the Rotated Surface Code Hamiltonian



$$\begin{aligned}
 H_{RSC}^{(3)} = & -X_1 X_2 X_4 X_5 - X_5 X_6 X_8 X_9 \\
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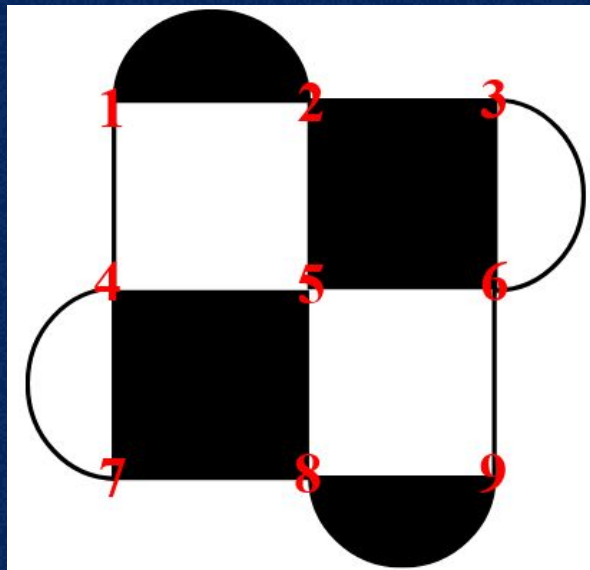
$$H_{RSC}^{(L)} = - \sum_{f=1}^{F_w} A_f - \sum_{f=1}^{F_b} B_f$$

where

$$\begin{aligned}
 A_f &= \prod_{j \in \partial f} X_j \\
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- Well-known 4-local Hamiltonian.

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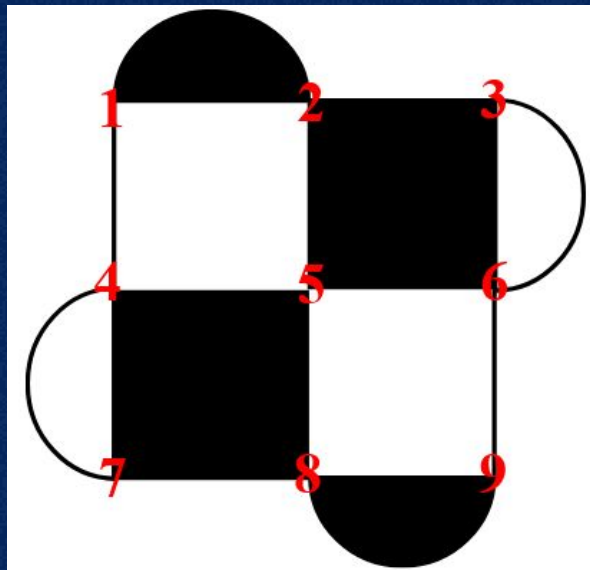
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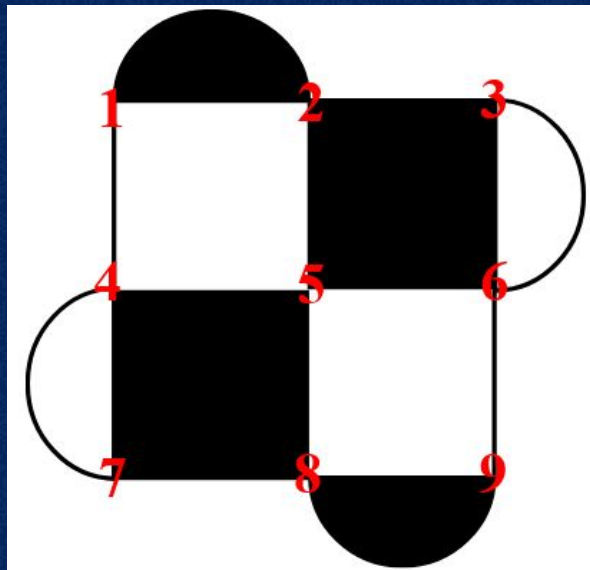
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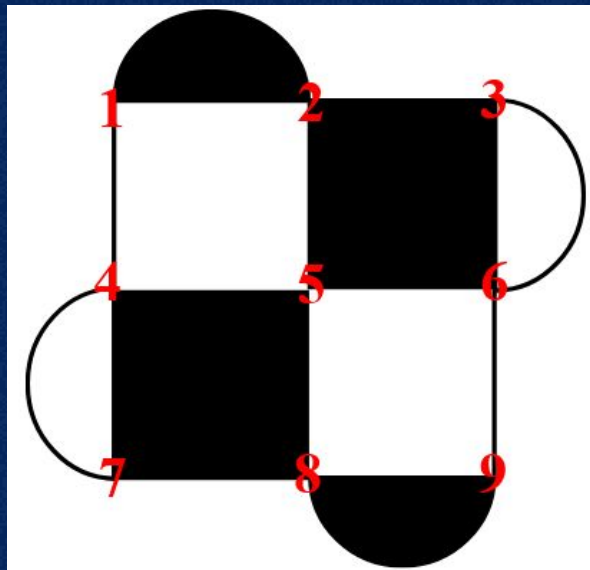
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- **2 degenerate ground states**.
- Ground space exhibits **topological order** (gap, ground states locally indistinguishable).
- **1 protected qubit** can be encoded in ground space.
- Ground states exhibit **long-range entanglement**.

Example: the Rotated Surface Code Hamiltonian

Results:

- **Sanity check:** no 2,3-local n -qubit parent Hamiltonian for the ground space of H_{RSC} (proven analytically in

Graph states as ground states of many-body spin-1 / 2 Hamiltonians

[M. Van den Nest](#)¹, [K. Luttmer](#)¹, [W. Dür](#)^{1,2}, and [H. J. Briegel](#)^{1,2}

Physical Review A

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Physical Review A

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Physical Review A

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- We obtain the 4-local n -qubit parent Hamiltonian with **largest spectral gap** for a fixed energy range.
- For 2x2 and 3x3 lattices, the ground space of H_{RSC} appears as an **excited eigenspace** of 2,3-local n -qubit Hamiltonians, up to errors of order machine precision.



Conclusions and open questions

- Instances can be found of k' -local Hamiltonians leading to the same dynamics as given k -local Hamiltonians of the same dimension, with $k' < k$, for a particular subspace of initial states. Q: given any H_T , is there always a suitable H_{OS} ? If not, what is the relative volume of the manifold of simulatable H_T ? What is the largest dimension that the subspace of initial states can have for each case?



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- Given a target ground space, an SDP can be solved to i) determine whether a k' -local parent Hamiltonian exists, ii) if so, construct the one maximizing the spectral gap. Q: aside from the ground space of H_{RSC} , which other interesting states would benefit from having a more local parent Hamiltonian?

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- The ground space of H_{RSC} may appear as an excited subspace of 2,3-local Hamiltonians, up to errors of order machine precision. Q: is this true for other stabilizer states, e.g. GHZ?

Thanks for your attention



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lecturer positions opening soon!