

# Non-Abelian lattice gauge theory on quantum computers

1. Motivation
2. Gauge fields on gated-based quantum computers
3. Gauge fields on a quantum annealer
4. Including quarks
5. Outlook



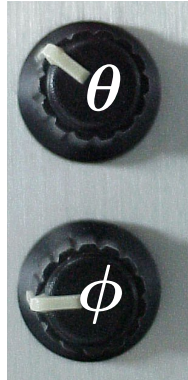
## From bits to qubits

Classical computers use bits.  
One bit is either  $|0\rangle$  or  $|1\rangle$ .



$$= |0\rangle$$

Quantum computers use qubits.  
One qubit is a superposition of  $|0\rangle$  and  $|1\rangle$ .



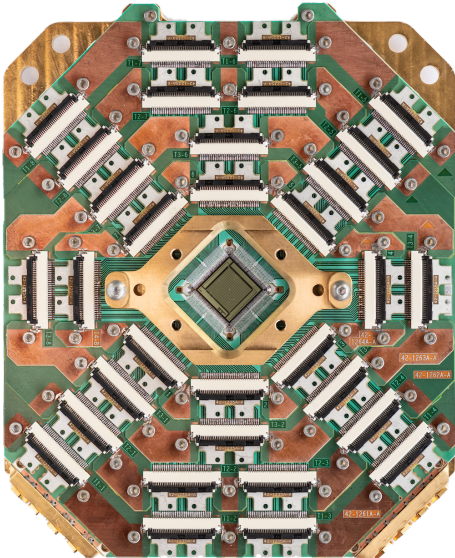
$$= \cos\left(\frac{\theta}{2}\right) |0\rangle + e^{i\phi} \sin\left(\frac{\theta}{2}\right) |1\rangle$$

Multiple bits act independently.

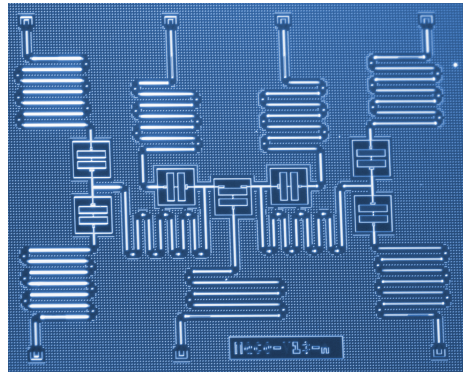
Multiple qubits can be entangled, so measuring one affects the others.

NOTE: Qubits can be in a *superposition of all* classically allowed states.

# I will show results from qubits at IBM and D-Wave

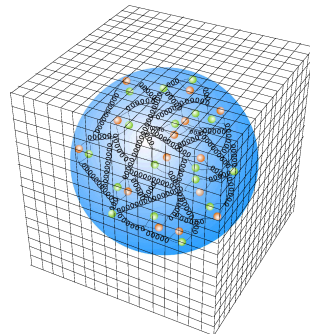
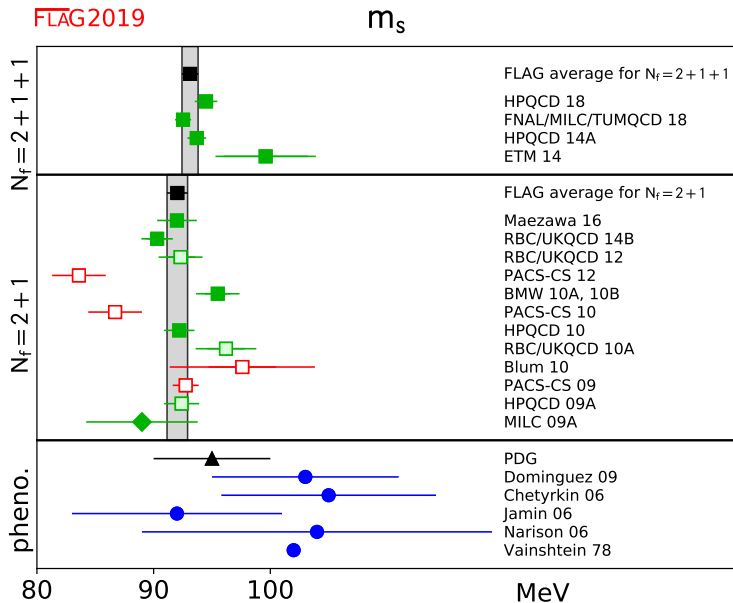


D-Wave, 5760 qubits, no gates



IBM, 7 qubits, universal gate set

# Lattice gauge theory is very successful without qubits



## What qubits might do for lattice gauge theory

Quantum computers offer an efficient Hamiltonian-based approach that might...

...allow us to avoid Euclidean time, thus moving from statics to dynamics.

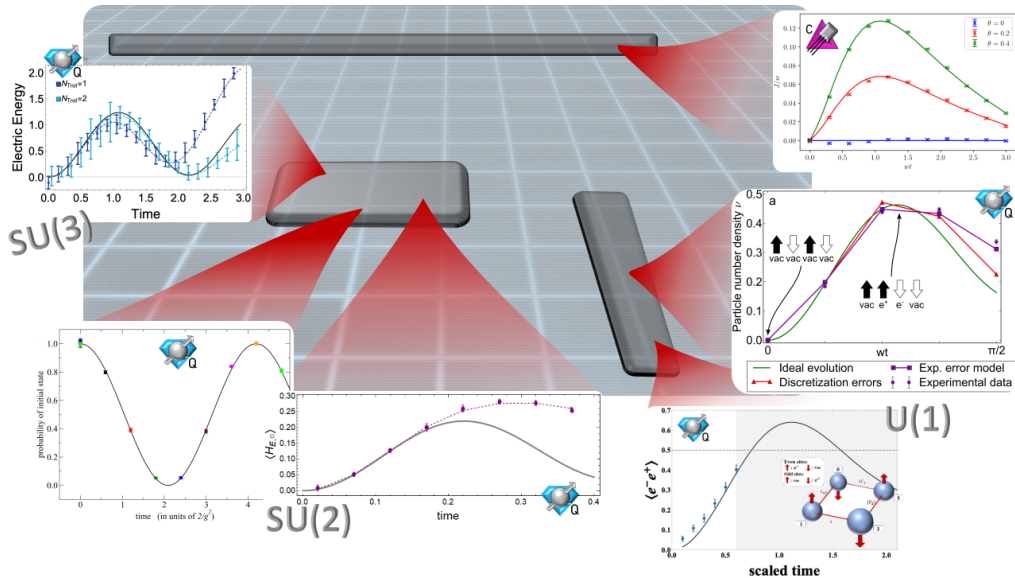
...allow us to include a chemical potential, thus reaching nuclear densities.

*Lattice QCD at non-zero density would be valuable for heavy-ion collisions, the early Universe and neutron-star structure. In practice, simulations at finite  $\mu$  suffer from a “sign problem” and are at a rudimentary stage.*

— paraphrased from Particle Data Group, Review of Lattice QCD

# Time evolution in gauge theories using qubits

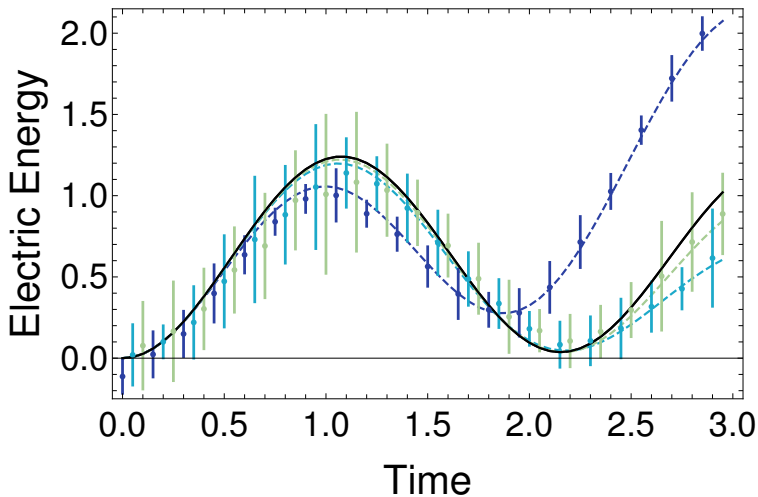
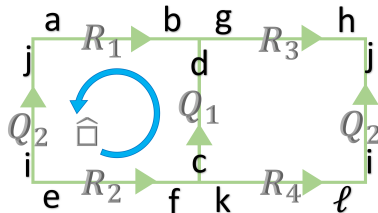
Figure 2 from Klco, Roggero and Savage, arXiv:2107.04769



# SU(3) pure gauge theory on gate-based hardware

Ciavarella, Klco, Savage, Phys.Rev.D103(2021)094501

Two plaquettes with a periodic boundary:



# How the Hamiltonian was constructed

Ciavarella, Klco, Savage, Phys.Rev.D103(2021)094501

The Hamiltonian is

$$\hat{H} = \frac{g^2}{2a^{d-2}} \sum_{b, \text{links}} \left| \hat{\mathbf{E}}^{(b)} \right|^2 + \frac{1}{2a^{4-d}g^2} \sum_{\text{plaquettes}} \left[ 6 - \hat{\square}(\mathbf{x}) - \hat{\square}^\dagger(\mathbf{x}) \right]$$

The chromoelectric term comes from

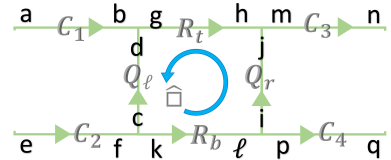
$$\sum_b \left| \hat{\mathbf{E}}^{(b)} \right|^2 |p, q\rangle = \frac{p^2 + q^2 + pq + 3p + 3q}{3} |p, q\rangle$$

The chromomagnetic term comes from

$$\left\langle \begin{pmatrix} C_1, R'_t, C_3 \\ Q'_\ell, Q'_r \\ C_2, R'_b, C_4 \end{pmatrix} \middle| \hat{\square} \middle| \begin{pmatrix} C_1, R_t, C_3 \\ Q_\ell, Q_r \\ C_2, R_b, C_4 \end{pmatrix} \right\rangle =$$

$$\sqrt{\frac{\dim(\mathbf{R}_t) \dim(\mathbf{R}_b)}{\dim(\mathbf{R}'_t) \dim(\mathbf{R}'_b) \dim(\mathbf{Q}_\ell) \dim(\mathbf{Q}_r) \dim(\mathbf{Q}'_\ell)^3 \dim(\mathbf{Q}'_r)^3}}$$

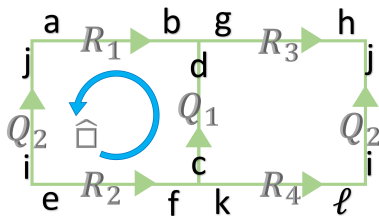
$$\begin{Bmatrix} \bar{\mathbf{R}}_t & C_1 & \bar{\mathbf{Q}}_\ell \\ 3 & 1 & 3 \\ \bar{\mathbf{R}}'_t & C_1 & \bar{\mathbf{Q}}'_\ell \end{Bmatrix} \bullet \square \begin{Bmatrix} \mathbf{R}_t & \bar{C}_3 & \bar{\mathbf{Q}}_r \\ 3 & 1 & 3 \\ \mathbf{R}'_t & \bar{C}_3 & \bar{\mathbf{Q}}'_r \end{Bmatrix} \square \bullet \begin{Bmatrix} \bar{\mathbf{R}}_b & C_2 & \mathbf{Q}_\ell \\ 3 & 1 & 3 \\ \bar{\mathbf{R}}'_b & C_2 & \mathbf{Q}'_\ell \end{Bmatrix} \bullet \square \begin{Bmatrix} \mathbf{R}_b & \bar{C}_4 & \mathbf{Q}_r \\ 3 & 1 & 3 \\ \mathbf{R}'_b & \bar{C}_4 & \mathbf{Q}'_r \end{Bmatrix} \square \bullet$$





# How the gauge links were truncated

Ciavarella, Klco, Savage, Phys.Rev.D103(2021)094501



Each of the 6 links is an irrep of SU(3): 1, 3,  $\bar{3}$ , 6,  $\bar{6}$ , 8, ...

Truncating to only  $\{1, 3, \bar{3}\}$  gives  $3^6 = \boxed{729}$  basis states for the lattice.

Enforcing Gauss's law at every vertex leaves only  $\boxed{27}$  of those basis states.

$\boxed{9}$  of 27 are **global** singlet states. (Apply  $\hat{\square}, \hat{\square}^\dagger$  to the strong-coupling vacuum.)

Spatial translation and parity block diagonalize the Hamiltonian:  $9 \rightarrow \boxed{4} + 2 + 2 + 1$ .

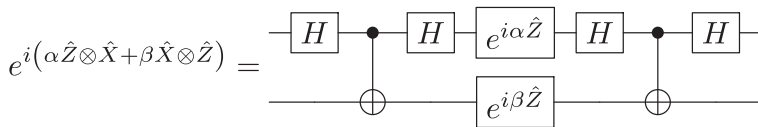
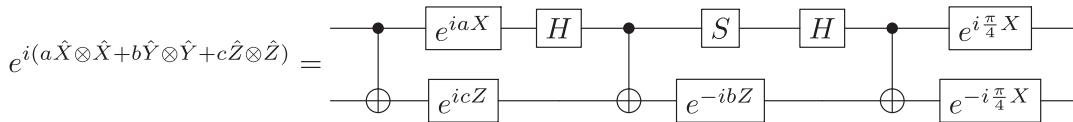
$$\hat{H}^{(1\bar{3}\bar{3};++)} = \frac{g^2}{2} \begin{pmatrix} 0 & 0 & 0 & 0 \\ 0 & \frac{16}{3} & 0 & 0 \\ 0 & 0 & \frac{16}{3} & 0 \\ 0 & 0 & 0 & 8 \end{pmatrix} + \frac{1}{2g^2} \begin{pmatrix} 6 & -2 & 0 & 0 \\ -2 & 5 & -\frac{\sqrt{2}}{9} & -\frac{\sqrt{2}}{3} \\ 0 & -\frac{\sqrt{2}}{9} & 6 & -\frac{2}{3} \\ 0 & -\frac{\sqrt{2}}{3} & -\frac{2}{3} & 6 \end{pmatrix}$$

Note the various Clebsch-Gordan combinations.

# How the circuit was implemented

Ciavarella, Klco, Savage, Phys.Rev.D103(2021)094501

First-order Trotter is used. The circuit has single qubit terms and these:



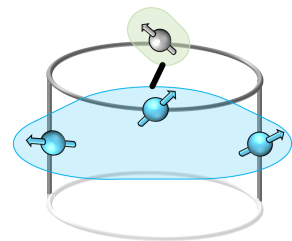
The IBM Athens chip has 5 qubits in total:

2 hold the state of the lattice.

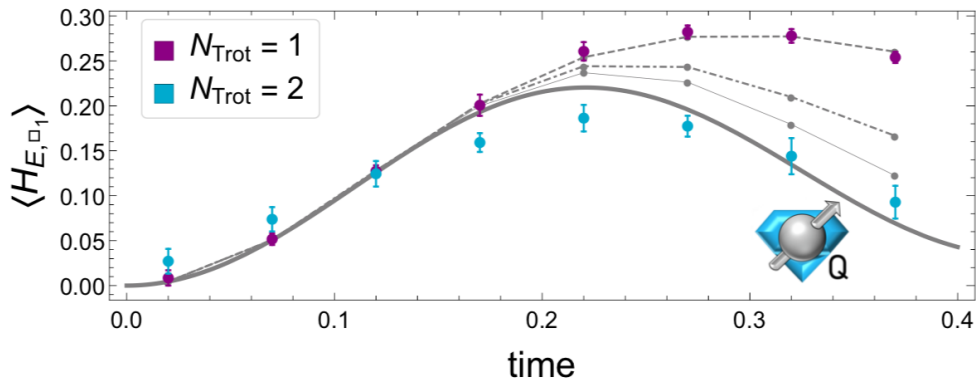
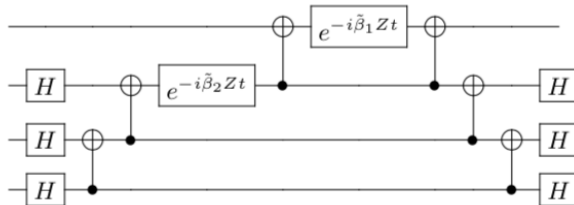
3 were used for post-selection error mitigation.

# SU(2) pure gauge theory on gate-based hardware

Klco, Savage, Stryker, Phys.Rev.D101(2020)074512



$$e^{-i\hat{\square}_2^{(1/2)}t} =$$

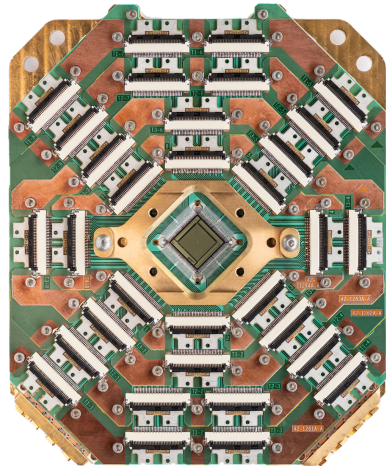


# What a D-Wave quantum annealer calculates

The hardware moves quasi-adiabatically to the ground state of

$$H(q) = \sum_{i=1}^N h_i q_i + \sum_{i=1}^N \sum_{j=i+1}^N J_{ij} q_i q_j$$

The user chooses any real  $h_i$  and  $J_{ij}$ .  
Each  $q_i$  is either 0 or 1.



D-Wave Advantage, 5760 qubits  
(usable as 180 all-to-all qubits)

A Rahman, Lewis, Mendicelli, Powell, Phys.Rev.D104(2021)034501

- Can this be used for a non-Abelian gauge theory? **Yes, some aspects.**
- Can it handle the various Clebsch-Gordan combinations? **Yes!**
- Will the number of qubits scale efficiently to large lattices?  
**No, not with our method on today's hardware.**

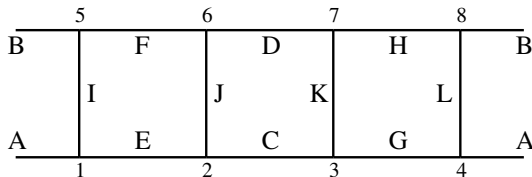
## Constructing the SU(2) Hamiltonian

$$\hat{H} = \frac{g^2}{2} \left( \sum_{i=\text{links}} \hat{E}_i^2 - 2x \sum_{i=\text{plaquettes}} \hat{\square}_i \right)$$

$$x \equiv \frac{2}{g^4}$$

$$\langle \psi | \hat{\mathbf{E}}_i^2 | \psi \rangle = j_i(j_i + 1)$$

$$\begin{aligned} \langle \psi_{\text{final}} | \hat{\square}_1 | \psi_{\text{initial}} \rangle &= (-1)^{j_A + J_E + j_I} \sqrt{(2j_I + 1)(2J_E + 1)} \begin{Bmatrix} j_A & j_E & j_I \\ \frac{1}{2} & J_I & J_E \end{Bmatrix} \\ &\quad (-1)^{j_C + J_E + j_J} \sqrt{(2j_E + 1)(2J_J + 1)} \begin{Bmatrix} j_C & j_E & j_J \\ \frac{1}{2} & J_J & J_E \end{Bmatrix} \\ &\quad (-1)^{j_D + J_F + j_J} \sqrt{(2j_J + 1)(2J_F + 1)} \begin{Bmatrix} j_D & j_F & j_J \\ \frac{1}{2} & J_J & J_F \end{Bmatrix} \\ &\quad (-1)^{j_B + J_F + j_I} \sqrt{(2j_F + 1)(2J_I + 1)} \begin{Bmatrix} j_B & j_F & j_I \\ \frac{1}{2} & J_I & J_F \end{Bmatrix} \end{aligned}$$



We also apply vertical reflection, horizontal reflection, and translation symmetries.

## The quantum annealer eigensolver (QAE)

Recall the variational method:  $E_0 \leq \frac{\langle \psi | H | \psi \rangle}{\langle \psi | \psi \rangle}$ .

Recall that D-Wave finds the minimum of  $H(q) = \sum_{i=1}^N h_i q_i + \sum_{i=1}^N \sum_{j=i+1}^N J_{ij} q_i q_j$ .

If the  $|\psi\rangle$  vector has only 0 and 1 as entries, then **those are basically the same**.

$q \rightarrow |\psi\rangle$

$h_i \rightarrow$  on-diagonals of  $H$

$J_{ij} \rightarrow$  off-diagonals of  $H$

All Hamiltonian entries can be entered directly and easily.

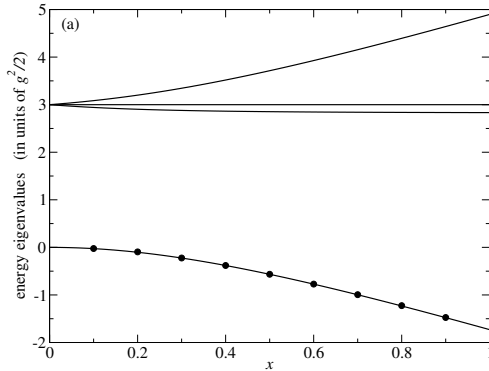
QAE handles a general vector (fixed-point representation) and the normalization. It uses one penalty term (called  $\lambda$ ) to avoid the null vector.

Tepelukhin, Kendrick, Babikov, J. Chem. Theory & Comp 15, 4555 (2019)

**We built an adaptive QAE** to use fewer qubits and solve larger Hamiltonians. Its only parameter is the  $\lambda$  from original QAE.

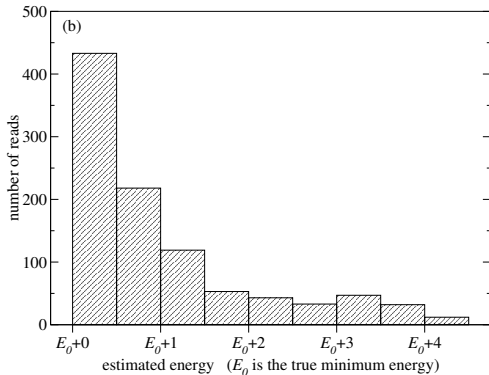
# Ground state eigenvalue for two plaquettes and $j_{\max} = \frac{1}{2}$

A Rahman, Lewis, Mendicelli, Powell, Phys. Rev. D104(2021)034501



Data points are from QAE.

Curves are exact eigenvalues.



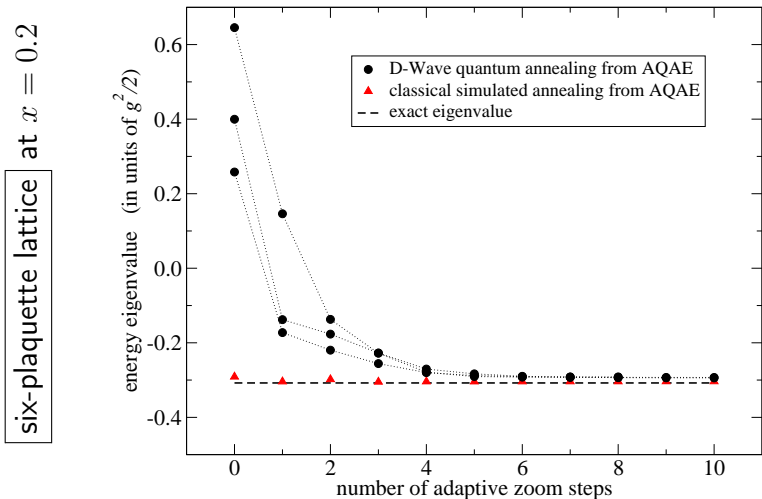
Raw data for  $x = 0.5$  in the graph above.

1000 anneals were used.

Each anneal took 20 microseconds.

# The importance of our adaptive algorithm

A Rahman, Lewis, Mendicelli, Powell, Phys.Rev.D104(2021)034501



The original QAE has no adaptive step, so zoom=0.

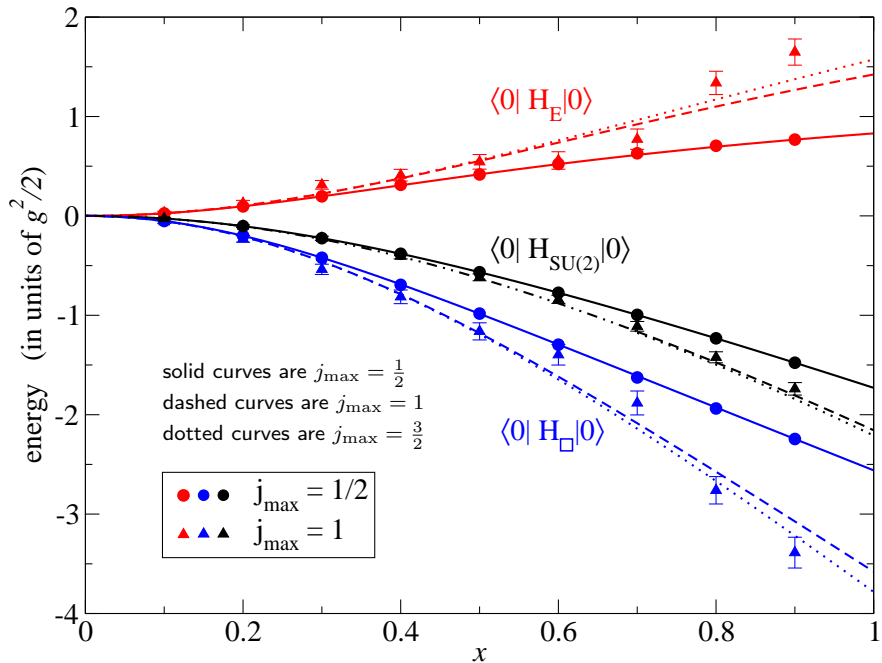
Our AQAE is helpful on a classical simulator.

Our AQAE is necessary for larger Hamiltonians on noisy quantum hardware.



# Assessing the gauge truncation

A Rahman, Lewis, Mendicelli, Powell, Phys. Rev. D104(2021)034501



# Time evolution as a minimization problem

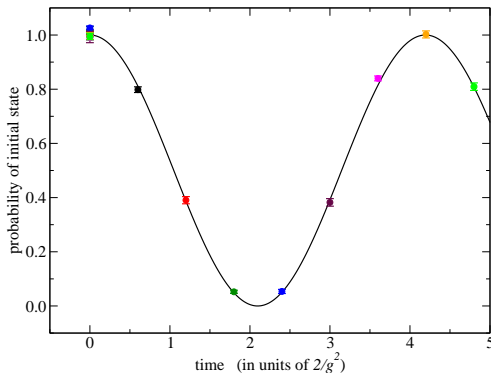
The TEDVP algorithm minimizes this functional:

$$\mathcal{L} = \sum_{t,t'} \langle t' | \langle \Psi_{t'} | \mathcal{C} | \Psi_t \rangle | t \rangle - \lambda \left( \sum_{t,t'} \langle t' | \langle \Psi_{t'} | \Psi_t \rangle | t \rangle - 1 \right)$$
$$\mathcal{C} = C_0 + \frac{1}{2} \sum_t \left( I \otimes |t\rangle \langle t| - e^{-i\epsilon H_t} \otimes |t+\epsilon\rangle \langle t| - e^{i\epsilon H_t} \otimes |t\rangle \langle t+\epsilon| + I \otimes |t+\epsilon\rangle \langle t+\epsilon| \right)$$

McClellan, Parkill, Aspuru-Guzik, Proc. Natl. Acad. Sci. 110, E3901 (2013)

For D-Wave hardware, we

- express  $H_t$  as imaginary so coefficients are real.
- use a combined QAE+TEDVP algorithm.



A Rahman, Lewis, Mendicelli, Powell, Phys. Rev. D104(2021)034501

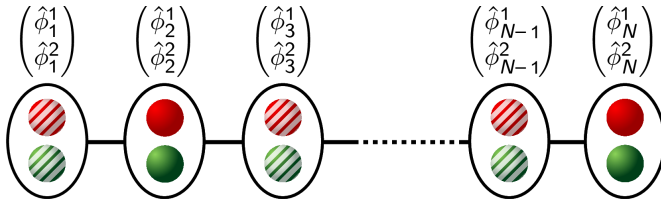
# Including quarks

Atas, Zhang, Lewis, Jahanpour, Haase, Muschik, accepted for publication.  
*SU(2) hadrons on a quantum computer via a variational approach*

Consider a one-dimensional lattice. It will have no colour-magnetic fields.

Put quarks and antiquarks on alternating sites ("staggered fermions").

We need two qubits per lattice site.



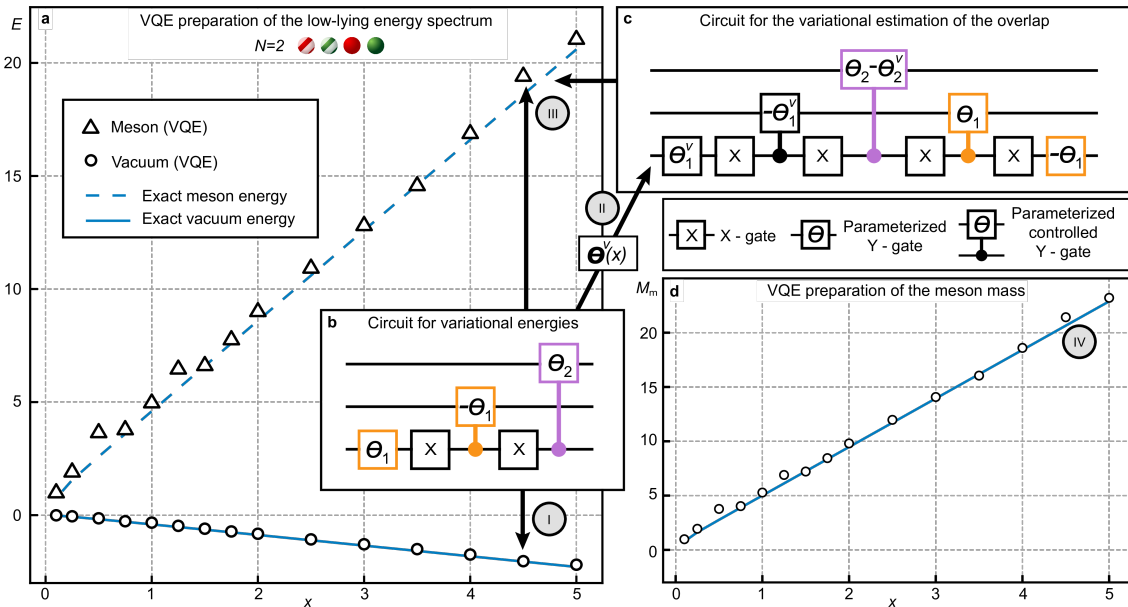
## Absorbing the gauge fields

There are two physics parameters: the gauge coupling and the quark mass.

With open lattice boundaries, gauge field effects are long-range quark interactions.

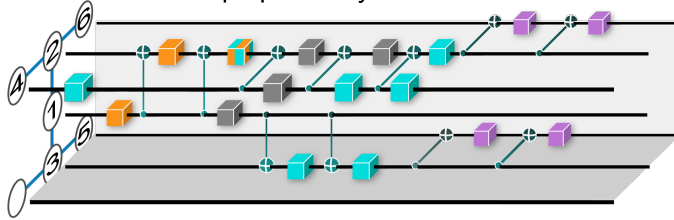
$$\begin{aligned}
 \hat{H} &= x\tilde{m}\hat{H}_m + \hat{H}_{\text{el}} + x\hat{H}_{\text{kin}} \\
 \hat{H}_m &= 2 \sum_{n=1}^N \left( \frac{(-1)^n}{2} (\hat{\sigma}_{2n-1}^z + \hat{\sigma}_{2n}^z) + 1 \right) \\
 \hat{H}_{\text{kin}} &= - \sum_{n=1}^{N-1} (\hat{\sigma}_{2n-1}^+ \hat{\sigma}_{2n}^z \hat{\sigma}_{2n+1}^- + \hat{\sigma}_{2n}^+ \hat{\sigma}_{2n+1}^z \hat{\sigma}_{2n+2}^- + \text{h.c.}) \\
 \hat{H}_{\text{el}} &= \frac{3}{8} \sum_{n=1}^{N-1} (N-n)(1 - \hat{\sigma}_{2n-1}^z \hat{\sigma}_{2n}^z) \\
 &\quad + \frac{1}{8} \sum_{n=1}^{N-2} \sum_{m>n}^{N-1} (N-m) (\hat{\sigma}_{2n-1}^z - \hat{\sigma}_{2n}^z) (\hat{\sigma}_{2m-1}^z - \hat{\sigma}_{2m}^z) \\
 &\quad + \sum_{n=1}^{N-2} \sum_{m>n}^{N-1} (N-m) (\hat{\sigma}_{2n-1}^+ \hat{\sigma}_{2n}^- \hat{\sigma}_{2m}^+ \hat{\sigma}_{2m-1}^- + \text{h.c.})
 \end{aligned}$$

# Computing the meson mass

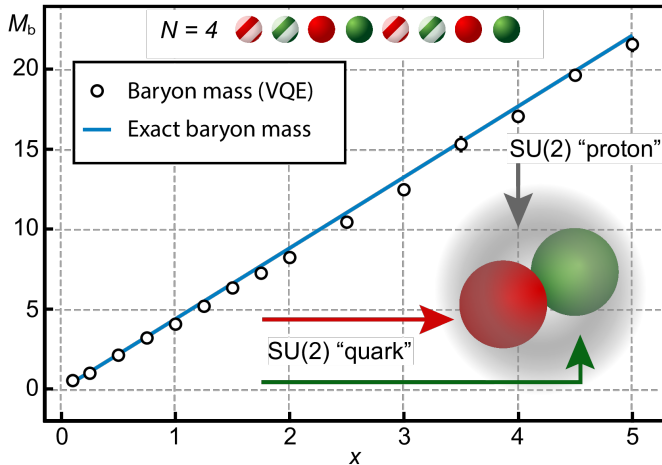


# Computing the baryon mass

a VQE circuit to prepare baryon and vacuum states



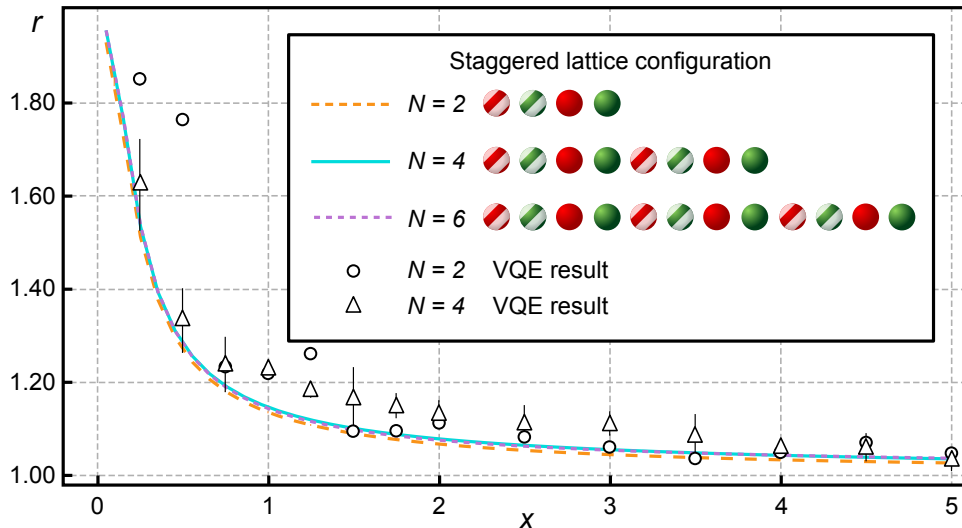
b VQE preparation of the baryon mass



# Computing the meson-to-baryon mass ratio

For continuum SU(2), the meson and baryon are exactly degenerate.

Our staggered lattice calculation is consistent with this continuum limit.



# Comparing several formulations

DAVOUDI, RAYCHOWDHURY, and SHAW

PHYS. REV. D **104**, 074505 (2021)

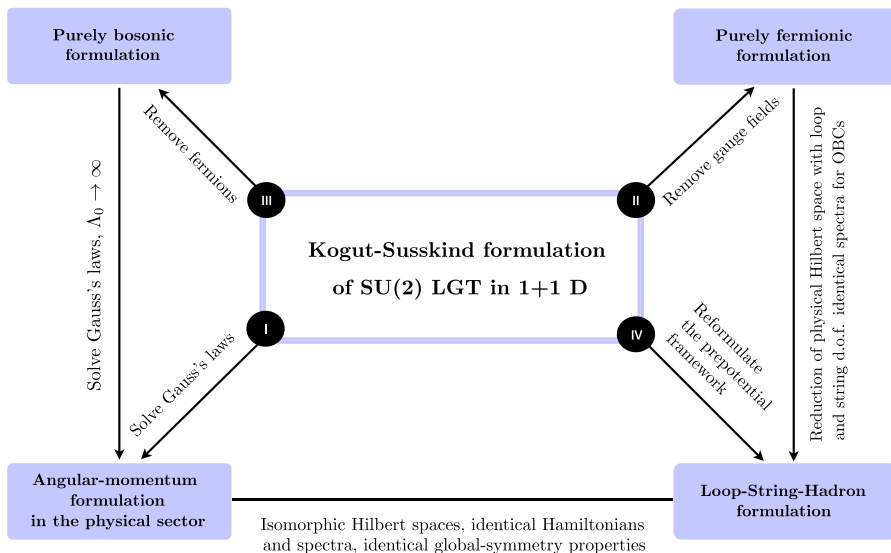
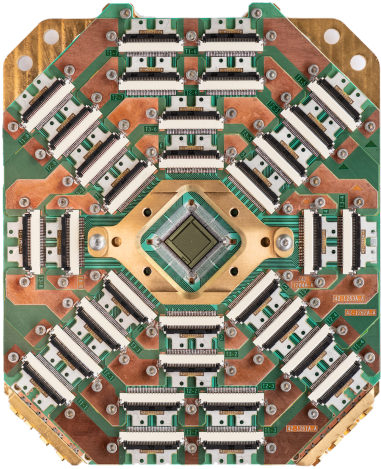


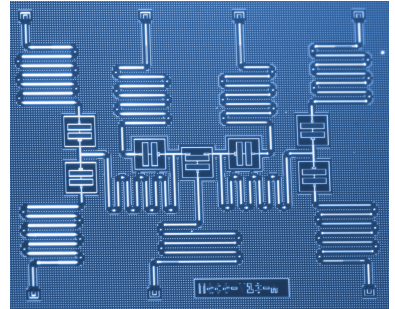
FIG. 1. Various formulations of the KS  $SU(2)$  LGT in 1 + 1D studied in this work and the connection among them.



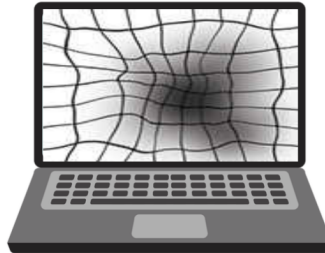
# Many ideas remain to be explored



D-Wave, 5760 qubits, no gates



IBM, 7 qubits, universal gate set



You are here!