First-Order Phase Transition Of The Schwinger Model With A Quantum Computer

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Why quantum computing **Some weaknesses of alternatives**

Exact diagonalization

N degrees of freedom of dimension d, $dim(\mathcal{H}) = d^N$

Monte Carlo

Sign problem with topological terms or chemical potential,

N_{Sam}

Tensor networks

ples
$$\sim e^{N/7}$$

Exponential increase in bond dimension with time,

$$D \sim e^t$$



Why topological terms Strong CP problem

- Why does QCD preserve CP symmetry (Charge conjugation Parity)?
- Experimentally CP within QCD interactions is conserved
- Theoretically we can have a CP violating topological term

$$\mathscr{L} = -\frac{1}{4}F_{\mu\nu}F^{\mu\nu} + \bar{\psi}\left(i\gamma^{\mu}D_{\mu} - m\right)\psi + \theta\frac{g^2}{32\pi^2}F_{\mu\nu}\tilde{F}^{\mu\nu}$$

Gives a neutron electric dipole moment

$$d_N = \left(5.2 \times 10^{-16} e \cdot cm\right)\theta$$

- Experimentally $d_N < 10^{-26} e \cdot cm \implies \theta < 10^{-10}$

• Range $\theta \in [0, 2\pi]$ raises the strong CP problem of fine-tuning to $\theta < 10^{-10}$

Some related work in the field from 2024

- Scalable Circuits for Preparing Ground States on Digital Quantum Computers: The Schwinger Model Vacuum on 100 Qubits, Farrell et al, PRX Quantum 5, 020315
- Concurrent VQE for Simulating Excited States of the Schwinger Model, Guo et al, arXiv:2407.15629
- Digital Quantum Simulation for Spectroscopy of Schwinger Model, Ghim et al, arXiv:2404.14788

Schwinger model theory **Staggered and Wilson fermions**

$$\Delta H_W = -r \frac{a^2}{2} \sum_n \overline{\psi}_n \left(\partial_1\right)^2 \psi_n$$

$$H_W = \sum_n \left(-\frac{1}{2} \overline{\psi}_n \left(r + i\gamma^1 \right) U_n \psi_{n+1} + \frac{1}{2} \overline{\psi}_n \left(-r + i\gamma^1 \right) U_n^{\dagger} \psi_{n-1} + (am_{lat} + r) \overline{\psi}_n \psi_n + \frac{ag^2}{2} \left(l_0 + L_n \psi_n \right) \right) \left(-\frac{1}{2} \overline{\psi}_n \psi_n + \frac{ag^2}{2} \left(l_0 + L_n \psi_n \right) \right) \left(-\frac{1}{2} \overline{\psi}_n \psi_n + \frac{ag^2}{2} \left(l_0 + L_n \psi_n \right) \right) \left(-\frac{1}{2} \overline{\psi}_n \psi_n + \frac{ag^2}{2} \left(l_0 + L_n \psi_n \right) \right) \left(-\frac{1}{2} \overline{\psi}_n \psi_n + \frac{ag^2}{2} \left(l_0 + L_n \psi_n \right) \right) \left(-\frac{1}{2} \overline{\psi}_n \psi_n + \frac{ag^2}{2} \left(l_0 + L_n \psi_n \right) \right) \left(-\frac{1}{2} \overline{\psi}_n \psi_n + \frac{ag^2}{2} \left(l_0 + L_n \psi_n \right) \right) \left(-\frac{1}{2} \overline{\psi}_n \psi_n + \frac{ag^2}{2} \left(l_0 + L_n \psi_n \right) \right) \left(-\frac{1}{2} \overline{\psi}_n \psi_n + \frac{ag^2}{2} \left(l_0 + L_n \psi_n \right) \right) \left(-\frac{1}{2} \overline{\psi}_n \psi_n + \frac{ag^2}{2} \left(l_0 + L_n \psi_n \right) \right) \left(-\frac{1}{2} \overline{\psi}_n \psi_n + \frac{ag^2}{2} \left(l_0 + L_n \psi_n \right) \right) \left(-\frac{1}{2} \overline{\psi}_n \psi_n + \frac{ag^2}{2} \left(l_0 + L_n \psi_n \right) \right) \left(-\frac{1}{2} \overline{\psi}_n \psi_n + \frac{ag^2}{2} \left(l_0 + L_n \psi_n \right) \right) \left(-\frac{1}{2} \overline{\psi}_n \psi_n + \frac{ag^2}{2} \left(l_0 + L_n \psi_n \right) \right) \left(-\frac{1}{2} \overline{\psi}_n \psi_n + \frac{ag^2}{2} \left(l_0 + L_n \psi_n \right) \right)$$





Schwinger model theory Hamiltonians with spin operators



$$\Delta H = \lambda \left(\sum_{\substack{n=0}}^{N-1} \right)$$

$$\begin{split} H_{S} &= \frac{x}{2} \sum_{n=0}^{N-2} \left(X_{n} X_{n+1} + Y_{n} Y_{n+1} \right) \\ &+ \frac{1}{2} \sum_{n=0}^{N-2} \sum_{k=n+1}^{N-1} \left(N - k - 1 + \lambda \right) Z_{n} Z_{k} \\ &+ \sum_{n=0}^{N-2} \left(\frac{N}{4} - \frac{1}{2} \left[\frac{n}{2} \right] + l_{0} (N - n - 1) \\ &+ \frac{m_{lat}}{g} \sqrt{x} \sum_{n=0}^{N-1} (-1)^{n} Z_{n} \\ &+ l_{0}^{2} (N - 1) + \frac{1}{2} l_{0} N + \frac{1}{8} N^{2} + \frac{\lambda}{4} N \,. \end{split}$$





Schwinger model theory **Observables**

Electric field density (EFD)

 $L_W = l_0 + L_{[N]}$

 $P_S = \frac{N}{2} + \frac{1}{2}$

 $P_W = N + \frac{1}{2} \sum_{n=0}^{N-1} \left(X_{2n} X_{2n+1} + Y_{2n} Y_{2n+1} \right)$

Particle number (PN)

$$L_{S} = l_{0} + \frac{1}{2} \left(L_{N/2-2} + L_{N/2-1} \right) = l_{0} + \frac{1}{4} \left(1 + Z_{N/2-1} \right) + \frac{1}{2} \sum_{k=0}^{N/2-1} L_{W} = l_{0} + L_{[N/2]-1} = l_{0} + \sum_{k=0}^{[N/2]-1} Q_{k} = l_{0} + \frac{1}{2} \sum_{k=0}^{[N/2]-1} (Z_{2k} + Z_{2k})$$

$$\sum_{n=0}^{N-1} (-1)^n Z_n$$



Schwinger model theory Phase diagram





Methods for quantum computing Goal outline

- Find ground states with noiseless variational quantum eigensolver (VQE) around 1st order phase transition (PT) line testing different ansatz and gates
- 6, 8, 10, 12 qubits for both fermion discretizations



- Prepare these states on IBM's quantum devices
- Measure the EFD and PN to demonstrate the PT

 $R(\theta_1, \theta_2, \theta_3)$ SO(4) $R(\theta_4, \theta_5, \theta_6)$ R_Y^{\dagger} (a) $R_Y(-\theta/2)$ R_Z R_{XX+YY} $R_Y(-\theta/2)$ R_Z^{\dagger} (b) $|q_1\rangle$ - $|q_1\rangle$. $|q_2\rangle$ — $|q_2\rangle$ – $|q_{N-3}\rangle$ $|q_{N-3}\rangle$ $|q_{N-2}\rangle$ — $|q_{N-2}\rangle$ $|q_{N-1}\rangle$ $|q_{N-1}|$

(d)





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Results for quantum computing VQE performance

- Brick ansatz performed better than ladder ansatz
- SO(4) gate ($\lambda \neq 0$) performed better than $R_{XX+YY}(\lambda = 0)$
- Achieved fidelities over 0.99
- 1 layer for $\frac{m_{\text{lat}}}{m_{\text{lat}}} = 10$ and 2 layers
- Number of layers consistent for all system sizes between 6-12 qubits

for
$$\frac{m_{\text{lat}}}{g} = 0$$







Results for quantum computing(a): Wilson $\frac{\tau}{g}$
 $N = 6, m_{lat}/g = 0, l_0 = 0.55$ (a): Wilson $\frac{\tau}{g}$
 $N = 6, m_{lat}/g = 0, l_0 = 0.55$

- Dynamical decoupling
- Readout error mitigation
- Pauli twirling
- Zero noise extrapolation (ZNE)

Noise factors: 1, 3, 5

 $\langle \psi | O | \psi \rangle = \langle 0 | U^{\dagger} O U | 0 \rangle$

 $|\psi\rangle = U|0\rangle = UU^{\dagger}U|0\rangle = UU^{\dagger}UU^{\dagger}U|0\rangle$

Probability



Methods for tensor networks **Goal outline**

Compute ground states as matrix product \bullet states with DMRG and compare to continuum mass perturbation theory for EFD observable

$$\frac{\mathscr{F}}{g} = \frac{e^{\gamma}}{\sqrt{\pi}} \left(\frac{m}{g} \right) \sin\left(2\pi l_0\right) - 8.9139 \frac{e^{2\gamma}}{4\pi} \left(\frac{m}{g}\right)^2$$

• Fixed volume $N/\sqrt{x} = 30$

- Fixed $m_r/g = 0.01$ $\frac{m_r}{g} = \frac{m_{lat}}{g} + MS(N, x, l_0)$ $MS_t = -\frac{m_{lat}}{g}$
- Therefore need the mass shift for each $N \in [70, 100]$ in steps of 10



(a-c) Wilson fermions: $l_0 = 0.1, 0.4, 0.6$

(d-f) Staggered fermions same x as Wilson EFD-MS: with $l_0 = 0.1, 0.4, 0.6$



Results for tensor networks Continuum extrapolation



Summary and outlook

- Analyzed different ansatz and gate performance with VQE
- Remnants of 1st order phase transition with QC
- Variety of error mitigation techniques proved helpful
- Continuum extrapolation comparison of staggered and Wilson fermions
- Explore capabilities of quantum devices in continuum extrapolations
- Looking into topological terms in 2+1 QED
- Investigate the Schwinger model within open quantum systems





Appendices

Variational quantum eigensolver How it works





Update parameters $\vec{\theta}$ to minimize cost function

Variational quantum eigensolver **Details of how algorithm was used**

- L-BFGS-B optimizer
- of the phase transition
- 100k iterations where all gates have their own free parameter
- ones
- high fidelity

• Initial states either generic or representing what is expected from left and right

• Warm up stage with 2k iterations while each layer has 1 free parameter then

• When adding a layer, the existing layers' initial parameters are the optimized

Usually only a few thousand iterations were sufficient for convergence with

Error mitigation techniques Dynamical decoupling

- decoherence occurs (inability of qubits to exhibit quantum behaviour)
- suppress this decoherence, e.g. X gate twice



During idle periods of qubits, the qubits interact with their environment and

Applying externally controlled gates that amount to the identity operation can

Error mitigation techniques Zero noise extrapolation

- 1. Amplify circuit noise for several noise factors
- 2. Run every noise amplified circuit
- 3. Extrapolate back to the zero noise limit



Matrix product states

- $|\psi\rangle = \sum A_{1,\alpha_1}^{\sigma_1} A_{\alpha_1,\alpha_2}^{\sigma_2} \dots A_{\alpha_{N-1},1}^{\sigma_N} |\sigma_1\rangle \otimes |\sigma_2\rangle \dots \otimes |\sigma_N\rangle$ $\sigma_1, \sigma_2, \dots, \sigma_N$
- Physical indices: $\sigma_i \in [1,d]$
- Bond indices: $\alpha_i \in [1,D]$

The MPS ansatz decomposes a quantum state into a set of rank-3 tensors:





Continuum extrapolation Mass shift measurement



Phys. Rev. D **108**, 014516 and references therein

Abstracts for literature in 2024

Scalable Circuits for Preparing Ground States on Digital Quantum Computers: The Schwinger Model Vacuum on 100 Qubits

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good agreement with classical matrix product state simulations.

The vacuum of the lattice Schwinger model is prepared on up to 100 qubits of IBM's Eagle-processor quantum computers. A new algorithm to prepare the ground state of a gapped translationally invariant system on a quantum computer is presented, which we call "scalable circuits ADAPT-VQE" (SC-ADAPT-VQE). This algorithm uses the exponential decay of correlations between distant regions of the ground state, together with ADAPT-VQE, to construct quantum circuits for state preparation that can be scaled to arbitrarily large systems. These scalable circuits can be determined with use of classical computers, avoiding the challenging task of optimizing parameterized circuits on a quantum computer. SC-ADAPT-VQE is applied to the Schwinger model, and is shown to be systematically improvable, with an accuracy that converges exponentially with circuit depth. Both the structure of the circuits and the deviations of prepared wave functions are found to become independent of the number of spatial sites, L. This allows a controlled extrapolation of the circuits, determined with use of small or modest-sized systems, to arbitrarily large L. The circuits for the Schwinger model are determined on lattices up to L = 14 (28 qubits) with the Qiskit classical simulator, and are subsequently scaled up to prepare the L = 50 (100 qubits) vacuum on IBM's 127-superconducting-qubit quantum computers ibm_brisbane and ibm_cusco. After introduction of an improved error-mitigation technique, which we call "operator decoherence renormalization", the chiral condensate and charge-charge correlators obtained from the quantum computers are found to be in

Abstracts for literature in 2024

Concurrent VQE for Simulating Excited States of the Schwinger Model

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This work explores the application of the concurrent variational quantum eigensolver (cVQE) for computing excited states of the Schwinger model. By designing suitable ansatz circuits utilizing universal SO(4) or SO(8) qubit gates, we demonstrate how to efficiently obtain the lowest two, four, and eight eigenstates with one, two, and three ancillary qubits for both vanishing and non-vanishing background electric field cases. Simulating the resulting quantum circuits classically with tensor network techniques, we demonstrate the capability of our approach to compute the two lowest eigenstates of systems with up to $\mathcal{O}(100)$ qubits. Given that our method allows for measuring the low-lying spectrum precisely, we also present a novel technique for estimating the additive mass renormalization of the lattice based on the energy gap. As a proof-of-principle calculation, we prepare the ground and first-excited states with one ancillary and four physical qubits on quantum hardware, demonstrating the practicality of using the cVQE to simulate excited states.

Abstracts for literature in 2024

Digital Quantum Simulation for Spectroscopy of Schwinger Model

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This note discusses a method for computing the energy spectra of quantum field theory utilizing digital quantum simulation. A quantum algorithm, called coherent imaging spectroscopy, quenches the vacuum with a time-oscillating perturbation and then reads off the excited energy levels from the loss in the vacuum-to-vacuum probability following the quench. As a practical demonstration, we apply this algorithm to the (1+1)-dimensional quantum electrodynamics with a topological term known as the Schwinger model, where the conventional Monte Carlo approach is practically inaccessible. In particular, on a classical simulator, we prepare the vacuum of the Schwinger model on a lattice by adiabatic state preparation and then apply various types of quenches to the approximate vacuum through Suzuki-Trotter time evolution. We discuss the dependence of the simulation results on the specific types of quenches and introduce various consistency checks, including the exact diagonalization and the continuum limit extrapolation. The estimation of the computational complexity required to obtain physically reasonable results implies that the method is likely efficient in the coming era of early fault-tolerant quantum computers.

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