

Dissipative Preparation of Vacuum and Meson States in a 1+1D $\mathbb{Z}_2$ Lattice Gauge Theory

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Quantum Simulation

- Real-time dynamics of lattice gauge theories are largely inaccessible to conventional Euclidean lattice calculations
- Digital and analog quantum simulators offer a direct route to real-time dynamics, but require **state preparation**
- Preparing the vacuum and low-lying states with high fidelity is a prerequisite for any subsequent scattering or real-time experiment
- Goal of this talk: steps towards **dissipative preparation** of such states on quantum computers

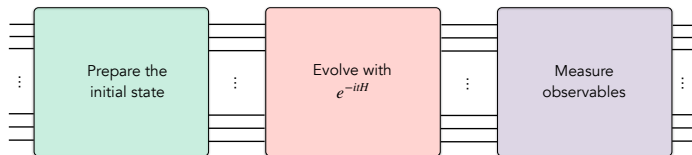


Figure from Davoudi, hep-lat:2507.15840

State Preparation Methods

- Adiabatic State Preparation
 - ▶ Performance guarantees
 - ▶ hindered by phase transitions, long ramp times are difficult for (current) hardware
- VQE Ansätze
 - ▶ Shallow circuits, classical optimization
 - ▶ No performance guarantees
- Spectral Filtering Algorithms
 - ▶ Performance guarantees
 - ▶ initial state overlap influences cost

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- Dissipative State Preparation
 - ▶ Potential for performance guarantees
 - ▶ Can start from **arbitrary initial states**

Dissipative Dynamics

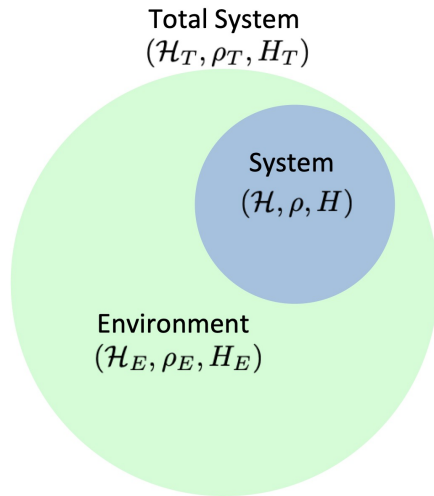


Figure from AIP Advances 1 February 2020; 10 (2): 025106

Dissipative Dynamics

The Lindbladian $\mathcal{L}$ provides the evolution of a system governed by Hamiltonian H in the presence of dissipation:

$$\dot{\rho} = \mathcal{L}[\rho] = -i[H, \rho] + \sum_{\alpha} \left(K_{\alpha} \rho K_{\alpha}^{\dagger} - \frac{1}{2} \{ K_{\alpha}^{\dagger} K_{\alpha}, \rho \} \right)$$

- $\{K_{\alpha}\}$: Jump operators, not necessarily unitary
- **Steady state condition:** $|\psi\rangle$ is a steady state of the dynamics if

$$H |\psi\rangle = E |\psi\rangle, \quad K_{\alpha} |\psi\rangle = 0 \quad \forall \alpha$$

- Lindblad dynamics can be simulated on quantum computers

¹In particular, we follow [Zhiyan Ding, Chi-Fang Chen, Lin Lin, PRR, arXiv:2308.15676]

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- Lindblad dynamics can be simulated on quantum computers
- **Engineer jump operators $\{K_{\alpha}\}$ such that the ground/meson state is the steady state of the dissipative dynamics¹**

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Engineering Jump Operators

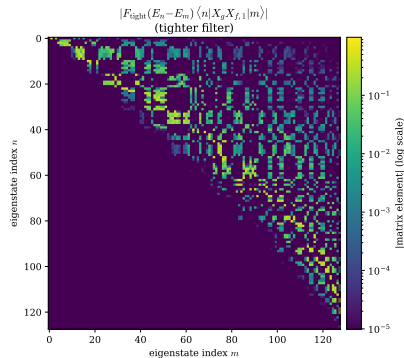
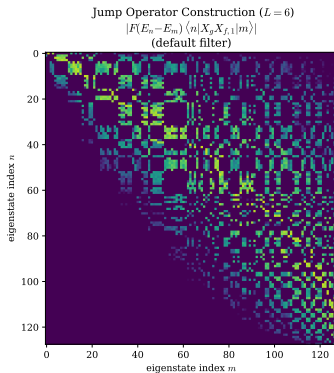
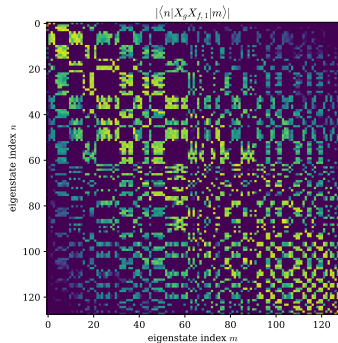
- Choose a **coupling operator** A_α acting on the system
- Apply an **energy filter** f in the eigenbasis of H :

$$K_\alpha = \sum_{i,j} \hat{f}(E_i - E_j) |i\rangle \langle i| A_\alpha |j\rangle \langle j|$$

Where $\hat{f}(E_i - E_j) = 0$ if $E_i \geq E_j$

- Consequence: K_α can only *decrement* energy
- In practice, implemented via a smooth approximation to a step function

Engineering Jump Operators



Engineering Jump Operators

We can construct K_α without the spectrum of H , via the inverse Fourier transform of $\hat{f}$, f :

$$\begin{aligned} K_\alpha &= \int_{-\infty}^{\infty} ds f(s) e^{iHs} A_\alpha e^{-iHs} \\ &\approx \sum_{l=-M}^M f(s_l) e^{iHs_l} A_\alpha e^{-iHs_l} w_l, \quad s_l = lt_s, \quad w_{\pm M} = \frac{t_s}{2}, w_l = t_s \end{aligned}$$

We truncate the bounds of the integral, and then approximate it as a discrete sum $\rightarrow$ allows for representation of K_α as a quantum circuit.

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Algorithm couples system register to ancilla qubit $\rightarrow$ measure and reset ancilla qubit $\rightarrow$ non-unitary dynamics on system qubits

In particular, we follow [Zhiyan Ding, Chi-Fang Chen, Lin Lin, PRR, arXiv:2308.15676]

Algorithm Overview

- 1 **Choose** a set of $\{A_\alpha\}$ coupling operators
- 2 Set filter parameters (based on $\Delta, ||H||$)
- 3 Determine the approximation parameters for the integral (bounds, number of samples)
- 4 Apply quantum circuit that approximates Lindbladian evolution

The choice of coupling operators is fundamental to the performance of the algorithm, not all choices of $\{A_\alpha\}$ will converge to the target state.

In this talk, we will consider the exact Lindblad dynamics, rather than the circuit dynamics.

In particular, we follow [Zhiyan Ding, Chi-Fang Chen, Lin Lin, PRR, arXiv:2308.15676]

$\mathbb{Z}_2$ LGT Hamiltonian

Staggered fermions ξ_n on matter sites, $\mathbb{Z}_2$ gauge qubits $X_{g,n}, Z_{g,n}$ on links, with periodic boundary conditions:

$$H = \underbrace{\frac{1}{2a} \sum_n (\xi_n^\dagger Z_{g,n} \xi_{n+1} + \text{h.c.})}_{H_{\text{kin}}} + \underbrace{m \sum_n (-1)^n \xi_n^\dagger \xi_n}_{H_{\text{mass}}} + \underbrace{\varepsilon \sum_n X_{g,n}}_{H_{\text{el}}}$$

- We work in the subspace of physical (Gauss Law preserving), total charge 0 states.
- Charge conjugation C commutes with H ; eigenstates carry momentum k and $c = \pm 1$:
 $C |k, c\rangle = c e^{-ik} |k, c\rangle$
- **Vacuum**: unique ground state, $c = +1$
- **Vector meson**: unique lowest energy state $|k, c = -1\rangle$

Targeting Excited States

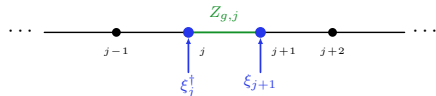
- To target excited states, we target ground states in specific sectors of the Hamiltonian
- If $[A_\alpha, S] = 0$, then $[K_\alpha, S] = 0$, and therefore we can choose $\{A_\alpha\}$ such that symmetries are not broken $\rightarrow$ dissipative dynamics will not mix states across sectors
- Starting state with definite quantum numbers $\rightarrow \{K_\alpha\}$ that respect symmetries $\rightarrow$ Ground state of sector

To obtain the vector meson state, we restrict the space of allowed coupling operators to translationally invariant, charge conjugation invariant operators.

Vacuum Coupling Operators

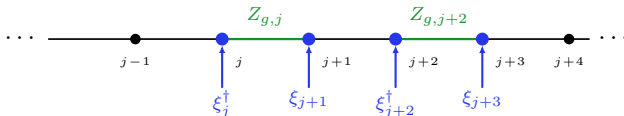
Single hop

$$K_j \sim \xi_j^\dagger Z_{g,j} \xi_{j+1} + \text{h.c.}$$



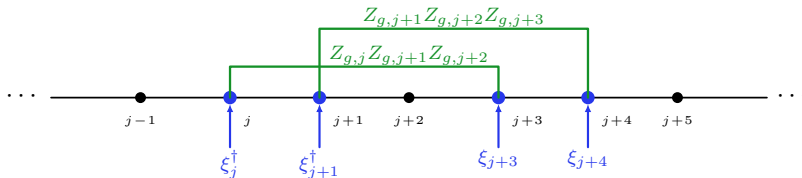
Double hop

$$\text{DH}_{j,2} \sim K_j K_{j+2} + \text{h.c.}$$



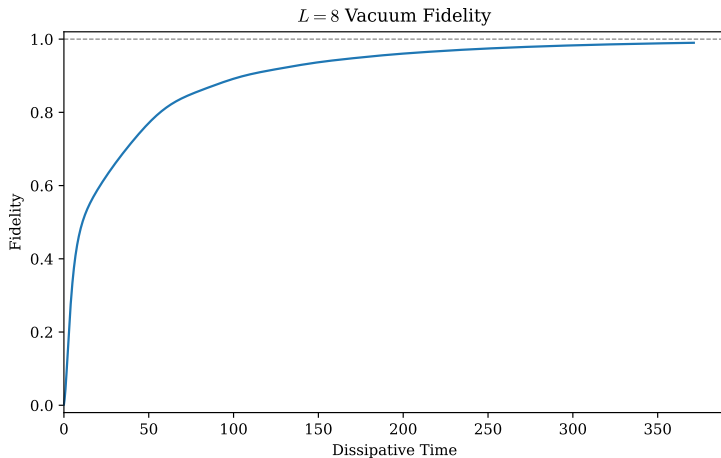
Pair hop

$$\text{PH}_{j,3} \sim \xi_j^\dagger \xi_{j+1}^\dagger \left(\prod_{m=j}^{j+2} Z_{g,m} \right) \left(\prod_{m=j+1}^{j+3} Z_{g,m} \right) \xi_{j+3} \xi_{j+4} + \text{h.c.}$$

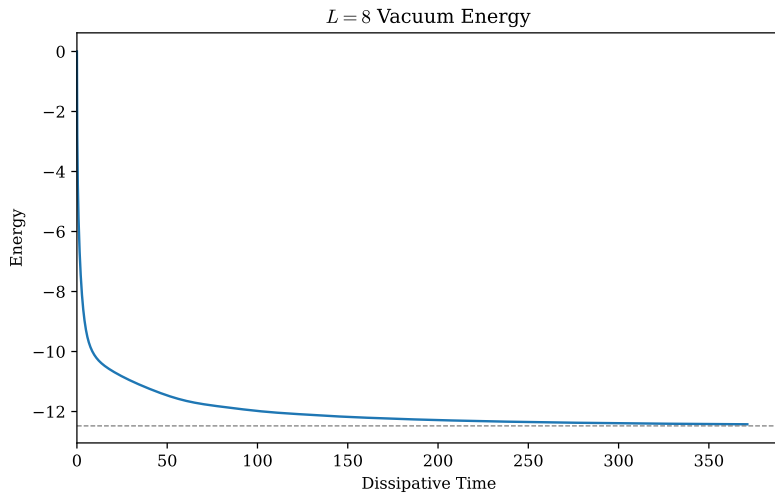


Vacuum State Preparation

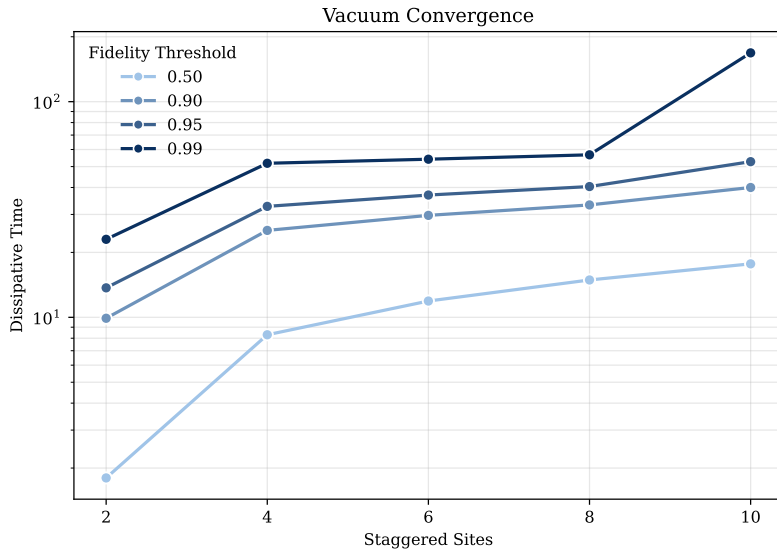
Starting state is the maximally mixed state in the physical Hilbert space



Vacuum State Preparation



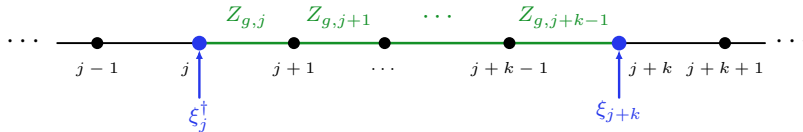
Vacuum State Preparation



Vector Meson Jump Operators

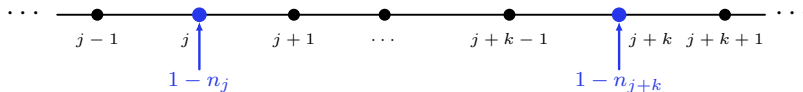
Extended fermion-hopping operators

$$W_k \sim \sum_{j=1}^L \left(\xi_j^\dagger \left(\prod_{m=j}^{j+k-1} Z_{g,m} \right) \xi_{j+k} + \text{h.c.} \right)$$

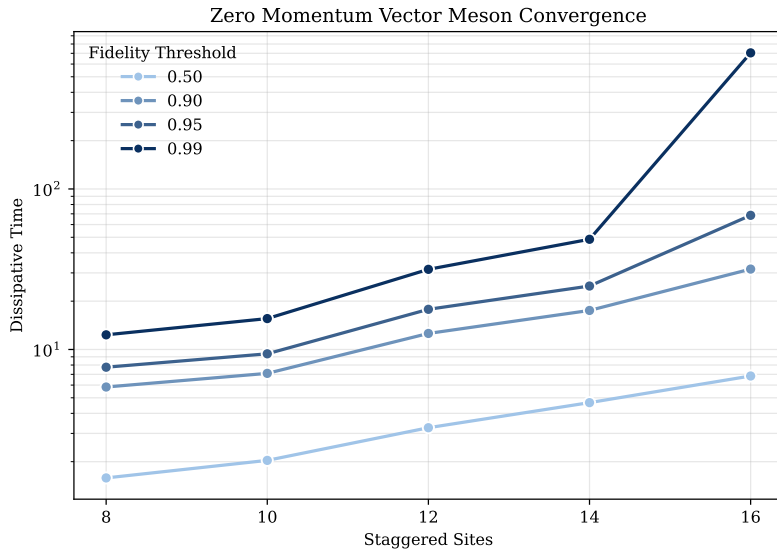


Extended density-density operators

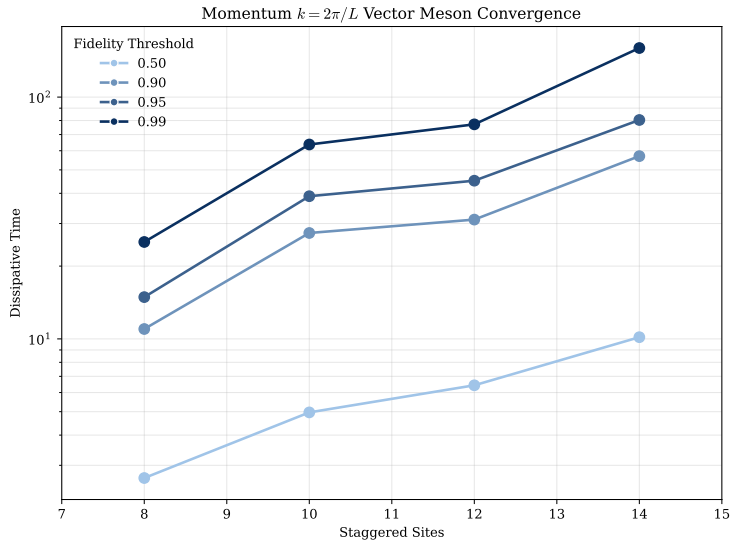
$$D_k \sim \sum_{j=1}^L (1 - n_j)(1 - n_{j+k}), \quad n_j = \xi_j^\dagger \xi_j$$



Vector Meson at $k = 0$



Vector Meson at $k \neq 0$



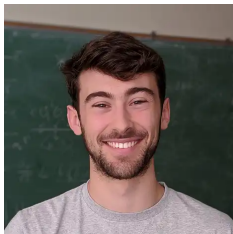
Future Work

- Extend Lindblad dynamics to larger systems
- Explicit circuit costs and dynamics

Open Questions:

- Given a set of coupling operators, can we provably converge to ground states of LGTs?
- Are there *optimal* choices of coupling operators for LGTs?
- Can we use dissipative state preparation in conjunction with other state preparation methods?

Acknowledgements



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Zohreh Davoudi

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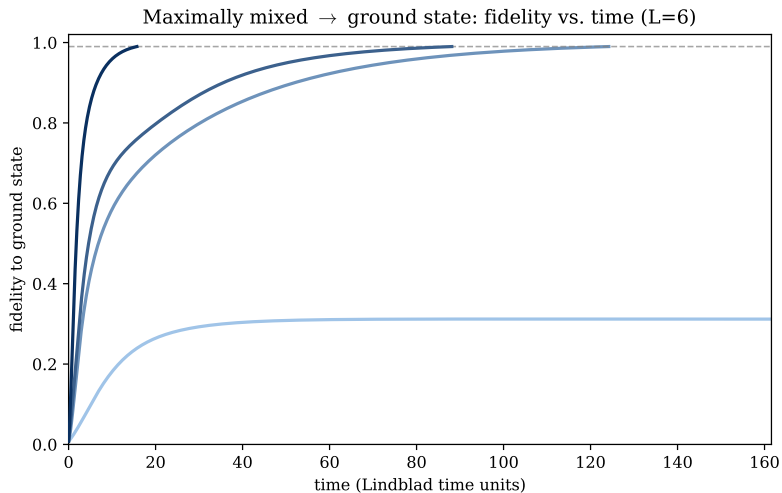


Questions?

Backup Slides

Operator Sets

Vacuum fidelity given different choices of coupling operators.



Filter Function

$$\hat{f}(\omega) = \frac{1}{2} \left(\operatorname{erf} \left(\frac{\omega + a}{\delta_a} \right) - \operatorname{erf} \left(\frac{\omega + b}{\delta_b} \right) \right)$$

Where

$$\operatorname{erf}(\omega) = \frac{2}{\sqrt{\pi}} \int_0^\omega e^{-x^2} dx$$

The inverse Fourier transform of the filter function is given by:

$$f(s) = \frac{\exp \left(-\frac{\delta_a^2 s^2}{4} \right) \exp(ias) - \exp \left(-\frac{\delta_b^2 s^2}{4} \right) \exp(ibs)}{2\pi i s}$$

We choose $a = 2.5\|H\|$, $\delta_a = .5\|H\|$, and $b = \delta_b = E_1 - E_0$.

$\mathbb{Z}_2$ LGT Hamiltonian

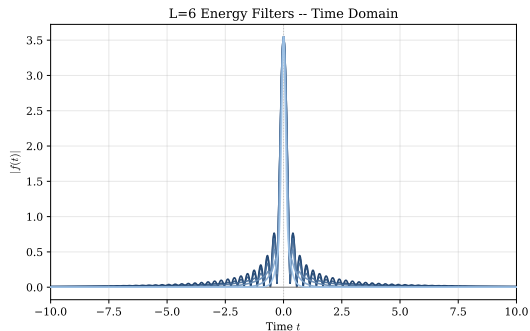
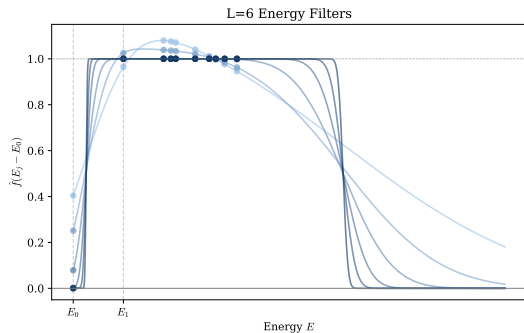
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After gauge-fixing and Jordan-Wigner transforming onto qubits:

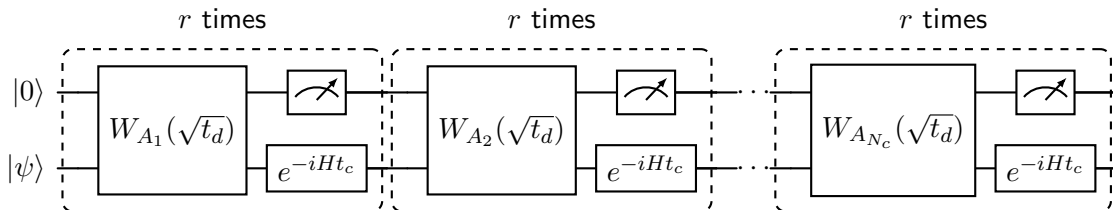
$$\begin{aligned} H_{\text{qubit}} = & -\frac{1}{2} \sum_{j=1}^{L-1} \left(\sigma_j^+ \sigma_{j+1}^- + \text{h.c.} \right) - \frac{1}{2} \left(\sigma_L^+ X_g \left(\prod_{k=2}^{L-1} Z_k \right) \sigma_1^- + \text{h.c.} \right) \\ & + m \sum_{j=1}^L (-1)^{j-1} \frac{1 - Z_j}{2} + 2\varepsilon Z_g \sum_{J=1}^L \prod_{k=1}^J [(-1)^{k-1} Z_k] \end{aligned}$$

Filter Function



Quantum Circuit

- N_c coupling operators
- Dissipative time step t_d
- Coherent time step t_c



Quantum Circuit

