

# **Real-time dynamics of the Schwinger model via variational quantum algorithms**

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Based on:

LN, A. Bapat and C. W. Bauer (LBNL) [arXiv:2302.10933] (to be published in PRD)

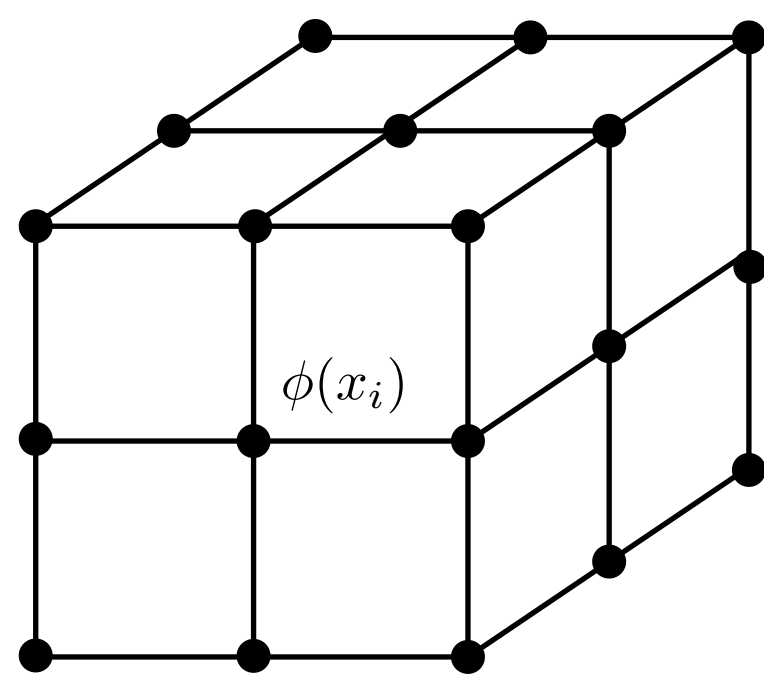
# Lattice gauge theory

- (conventional) lattice gauge theory

- discretize **spacetime**  
→ using Monte Carlo method

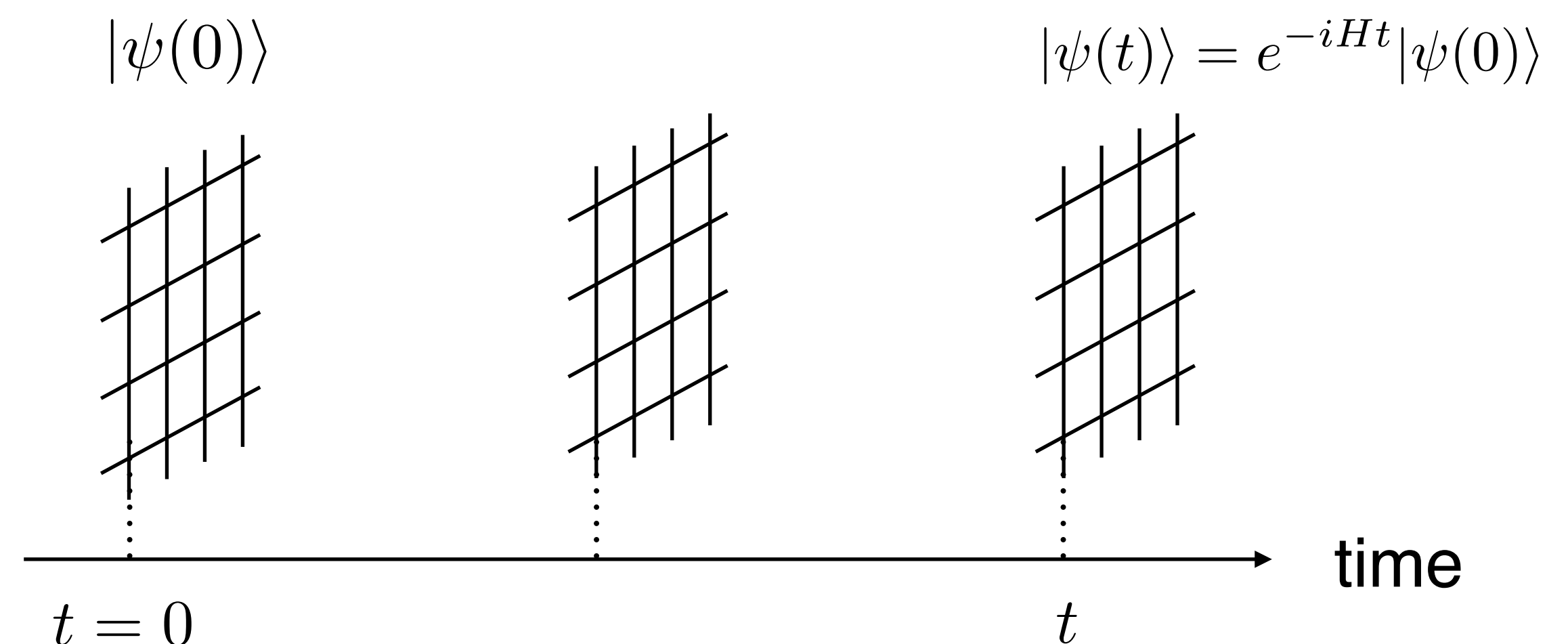
$$Z = \int [d\phi] e^{-S[\phi]} \rightarrow \sum_{\{\phi_i\}} e^{-S(\phi_i)}$$

- infamous **sign problem**
  - topological term
  - real-time dynamics, etc.



- Hamiltonian simulation

- discretize **space**
- no sign problem!
- need exponential resources...
  - quantum simulation
  - tensor network, etc.



# Quantum simulation

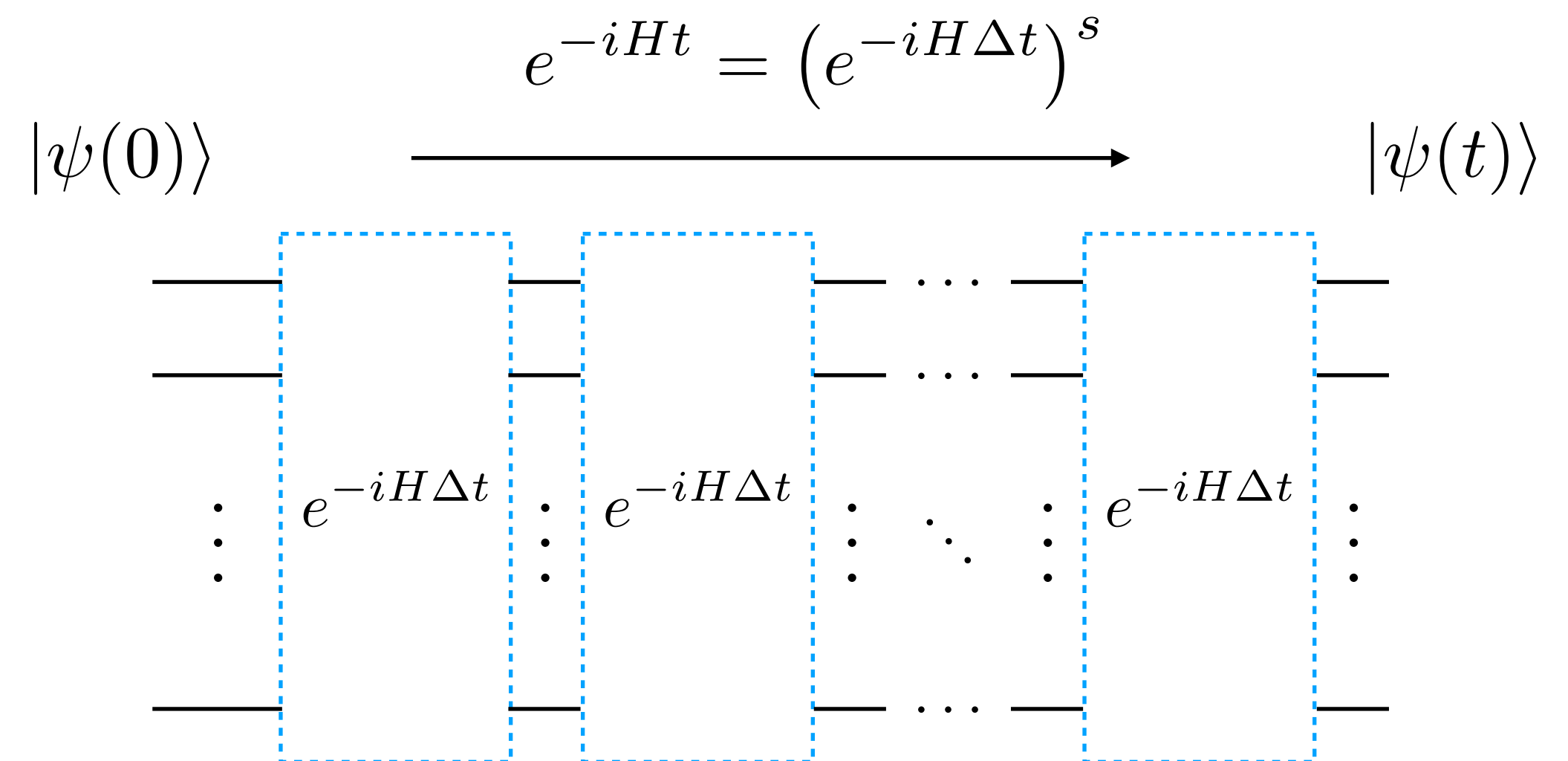
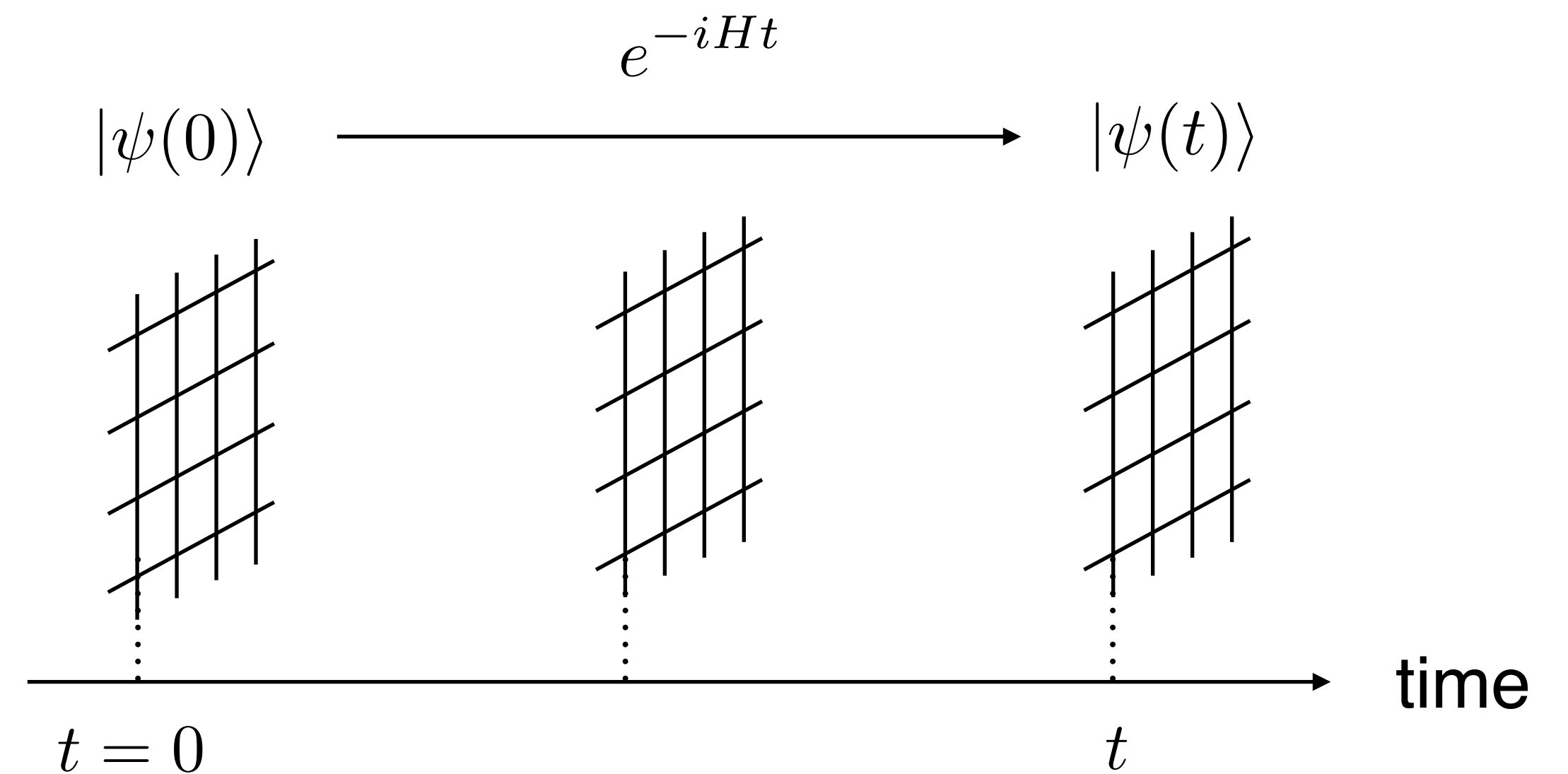
- **quantum simulation:**  
simulation using a quantum computer
  - real-time evolution  $|\psi(t)\rangle = e^{-iHt}|\psi(0)\rangle$
  - adiabatic time evolution

$$|\psi_{\text{GS}}\rangle = e^{-i \int dt H_A(t)t} |\psi_{\text{GS}}^{(0)}\rangle$$

- applications to HEP: e.g. scattering problem

[Jordan, Lee, Preskill, Science 336, 1130-1133 (2012)]

- good: exponential advantage
- bad:
  - still need many resources  
(#gates, #depths, #qubits)
  - near-term (NISQ) applications?



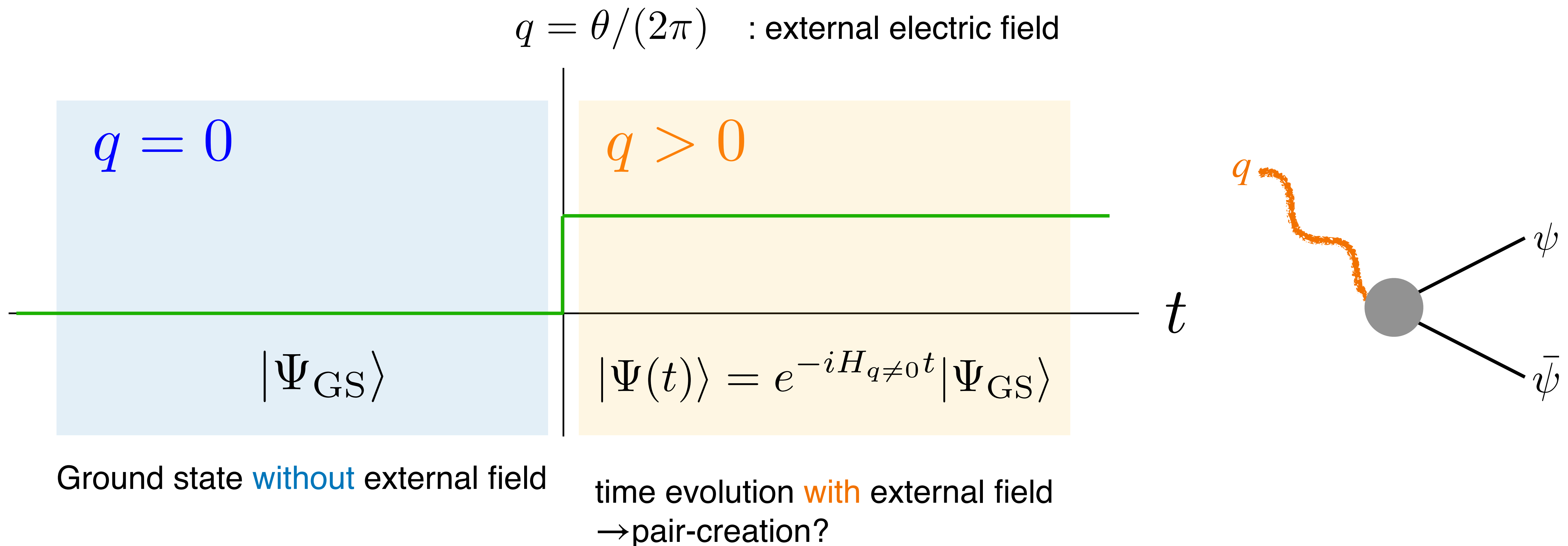
# Schwinger model

$$\mathcal{L} = -\frac{1}{4}F_{\mu\nu}F^{\mu\nu} + \frac{g\theta}{4\pi}\epsilon^{\mu\nu}F_{\mu\nu} + i\bar{\psi}\gamma^\mu(\partial_\mu + igA_\mu)\psi - m\bar{\psi}\psi$$

- simple toy model: 1+1d U(1) gauge theory = **Schwinger model** [Schwinger, *Phys. Rev.* 128, 2425]
  - exactly solvable for  $m = 0$
  - mass perturbation is available for small mass regime
- simple but still non-trivial
  - screening/confinement phenomena
  - we can include **the topological term** (cannot be treated in the MC method)
    - the effects of the external field (constant  $\theta$ )
    - the effects of probe charges (position dependent  $\theta$ )

# Quench dynamics in the Schwinger model

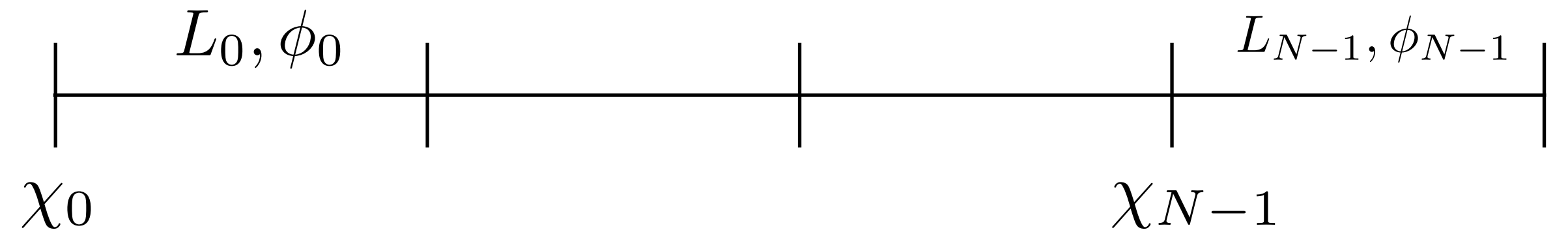
- Schwinger effect: particle pair creation due to strong **external electric field** [Schwinger, Phys. Rev. 82, 664, (1951)]
- Method: variational quantum algorithm (variational quantum eigensolver/simulation)



# Method

# Lattice Hamiltonian of Schwinger model

- $\chi_n$ : staggered fermion [Kogut, Susskind, Phys. Rev. D **11**, 395]
- $L_n, \phi_n$ : link variables (gauge field)



$$H_{\text{lat}} = J \sum_{n=0}^{N-2} \left( L_n + \frac{\theta}{2\pi} \right)^2 - i w \sum_{n=0}^{N-2} \left( \chi_n^\dagger e^{i\phi_n} \chi_{n+1} - \text{c.c.} \right) + m \sum_{n=0}^{N-1} (-)^n \chi_n^\dagger \chi_n$$

- gauge invariance: **Gauss's law constraint**

$$L_n - L_{n-1} = \chi_n^\dagger \chi_n - \frac{1 - (-)^n}{2}$$

- we can **eliminate** gauge fields (open boundary condition)!
  - automatically gauge invariant, no boson fields
  - cannot be used in higher dimension

# Spin description of the Schwinger model

- continuum Hamiltonian → lattice Hamiltonian (Kogut-Susskind formulation)
- lattice Hamiltonian + Gauss's law → fermionic representation of lattice Hamiltonian
- fermionic lattice Hamiltonian → spin system (Jordan-Wigner transformation)

$$H_{\text{spin}} = J \sum_{n=0}^{N-2} \left( \sum_{k=0}^n \frac{Z_k + (-)^k}{2} + q \right)^2 + \frac{w}{2} \sum_{n=0}^{N-2} (X_n X_{n+1} + Y_n Y_{n+1}) + \frac{m}{2} \sum_{n=0}^{N-1} (-)^n Z_n$$

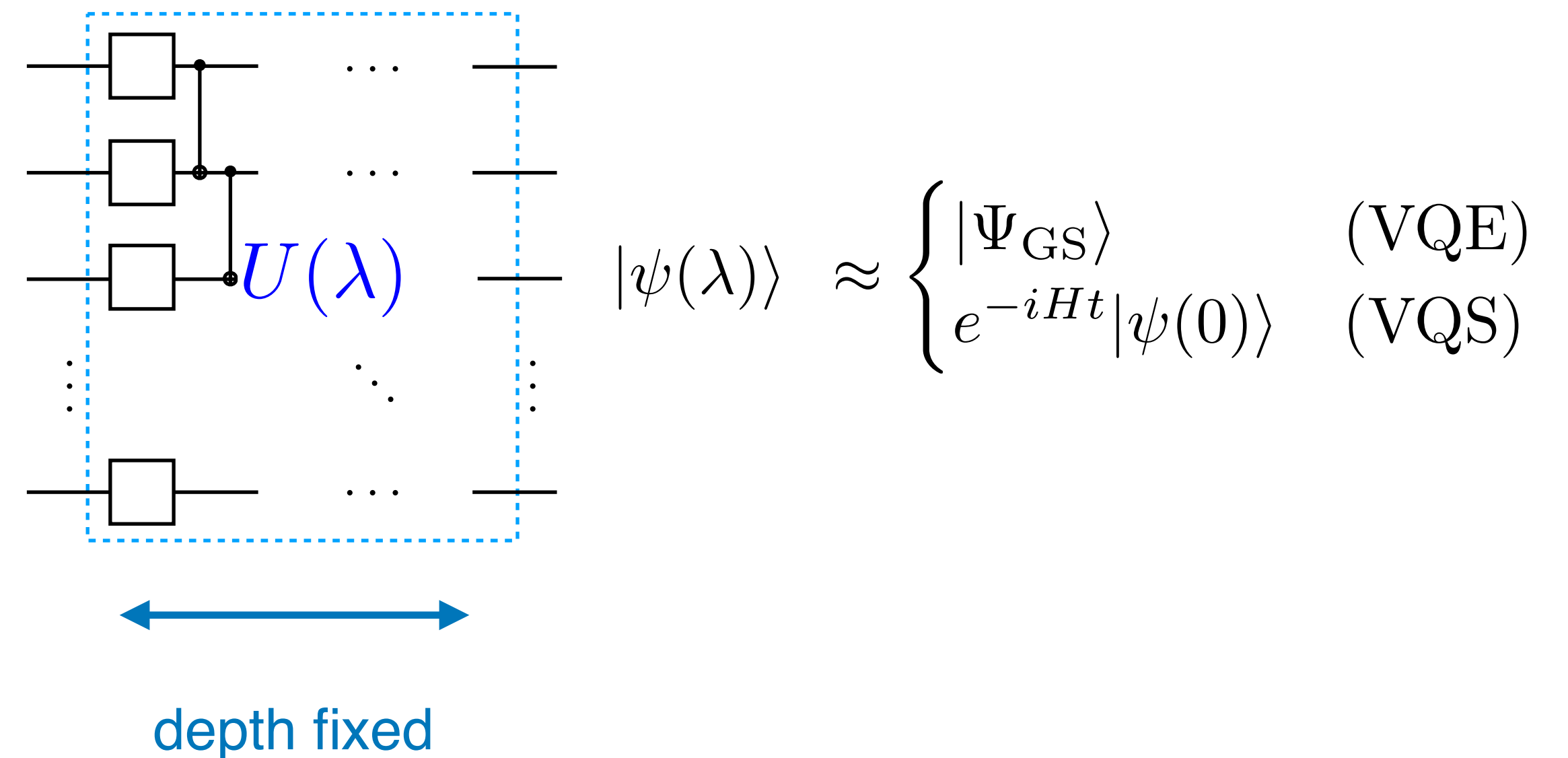
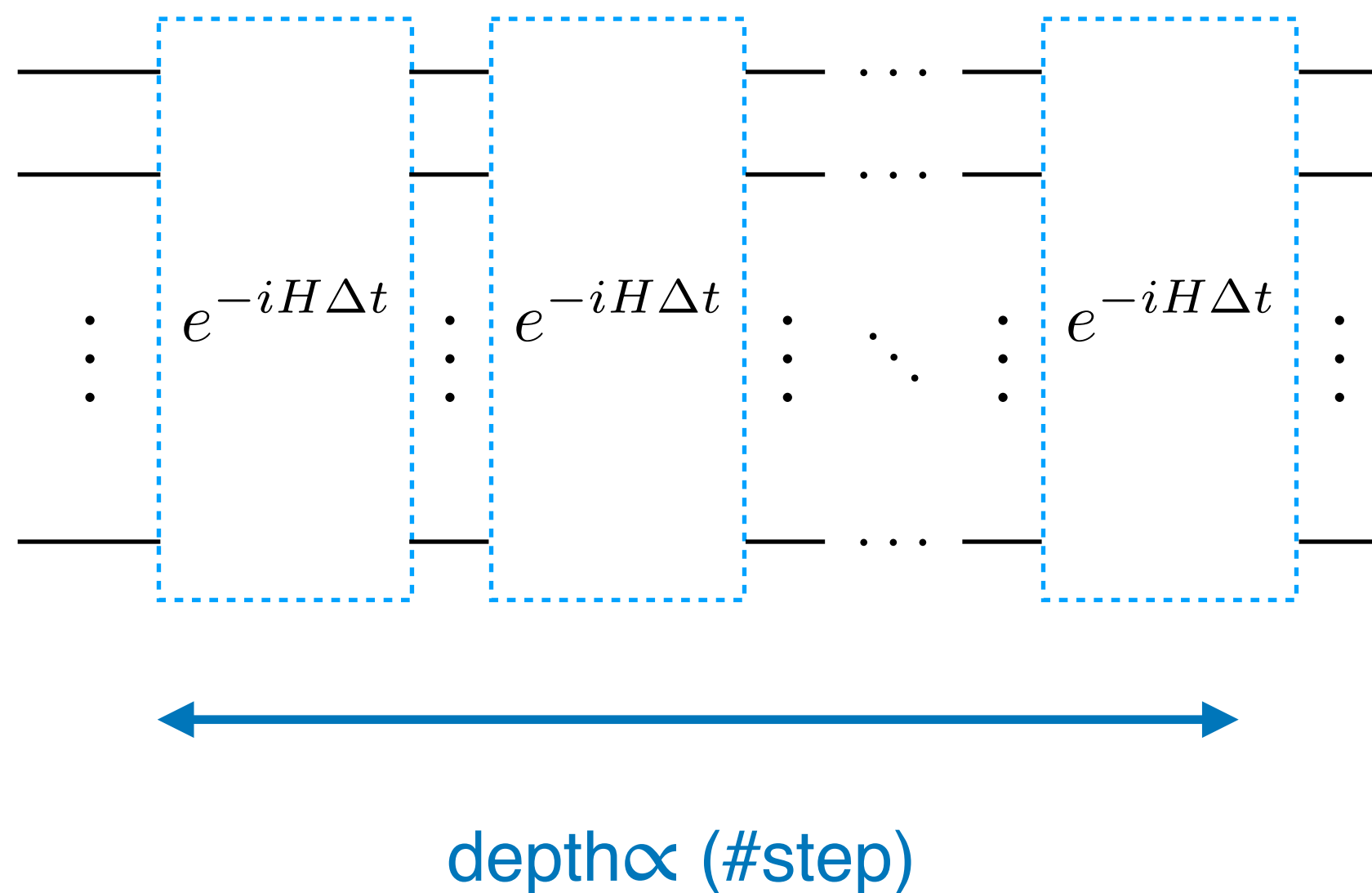
electric field term                      fermion kinetic term                      fermion mass term

( $X, Y, Z$ : Pauli matrices)



# Suzuki-Trotter vs variational method

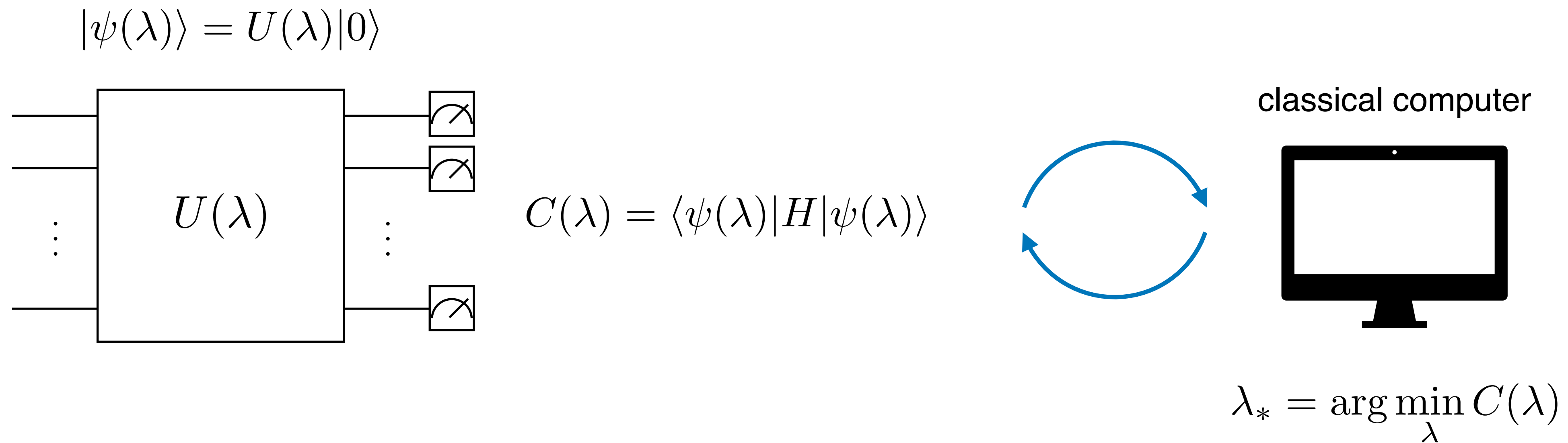
- Suzuki-Trotter method
  - #depth grows with #steps
  - decoherence problem on NISQ devices
- variational quantum algorithm (VQA)
  - approximate states by ansatz with **fixed depth**
  - state preparation: variational quantum eigensolver (VQE)
  - time-evolution: variational quantum simulation (VQS)



# Variational quantum eigensolver

[Peruzzo, A. *et al.* Nat. Commun. 5:4213 (2014)]

- goal: obtain the ground state
- approximate the ground state by ansatz  $|\psi(\lambda)\rangle$
- optimize cost function  $C(\lambda) = \langle \psi(\lambda) | H | \psi(\lambda) \rangle$  via classical computer  
→ ground state is given by  $|\psi(\lambda_*)\rangle$

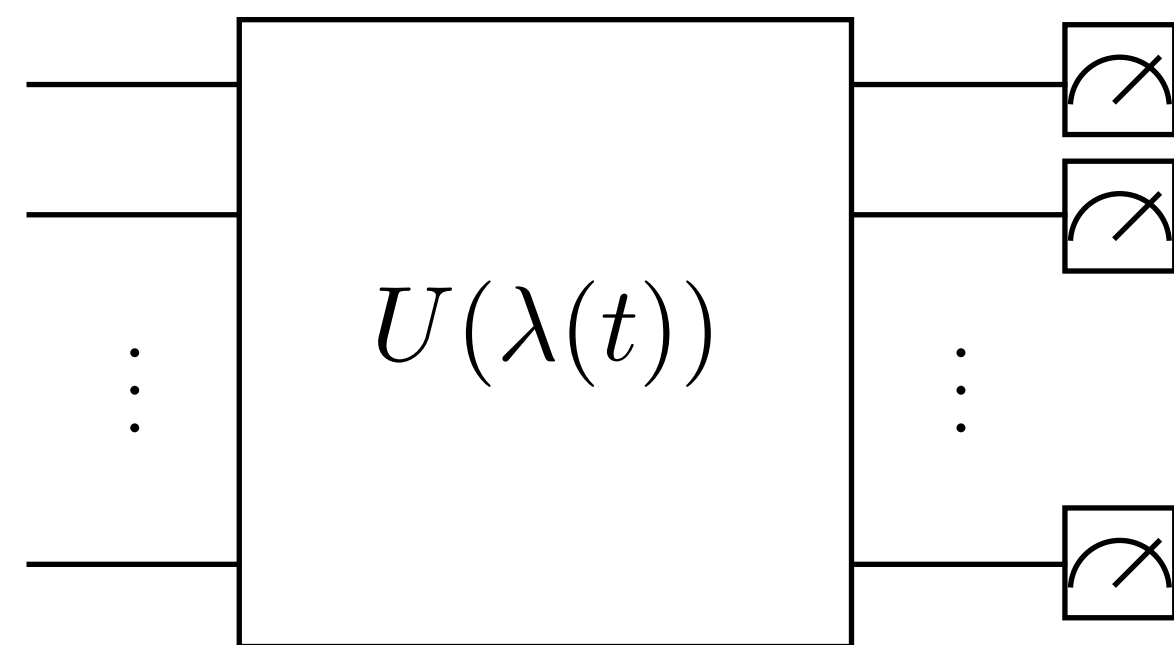


# Variational quantum simulation

[Li, Benjamin, Phys. Rev. X 7, 021050, (2017)]

- goal: obtain time-evolved state  $|\Psi(t)\rangle = e^{-iHt} |\Psi(0)\rangle$
- approximate  $|\Psi(t)\rangle$  by ansatz  $|\psi(\lambda(t))\rangle$  with time-dependent parameters
- evolution of states  $\rightarrow$  evolution of parameters  $\lambda(t)$  via McLacran's variational principle
- we use the same ansatz (Hamiltonian variational ansatz) for both VQE and VQS  
 $\rightarrow$  quench dynamics: set  $\lambda(0) = \lambda_*$  (obtained by VQE)

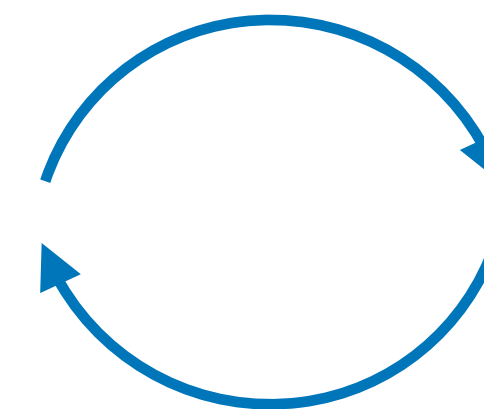
$$|\psi(\lambda(t))\rangle = U(\lambda(t))|0\rangle$$



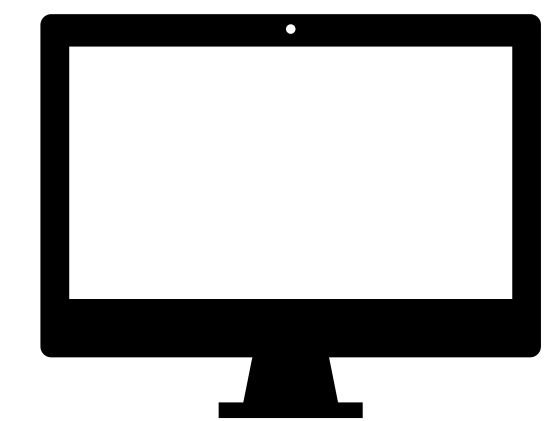
$$M_{ij} = \text{Re} \frac{\partial \langle \psi(\lambda) |}{\partial \lambda_i} \frac{\partial | \psi(\lambda) \rangle}{\partial \lambda_j}$$

$$V_i = \text{Im} \frac{\partial \langle \psi(\lambda) |}{\partial \lambda_i} H | \psi(\lambda) \rangle$$

(+correction terms)



classical computer



$$\sum_j M_{ij} \dot{\lambda}_j = V_i$$

# Hamiltonian variational ansatz (HVA)

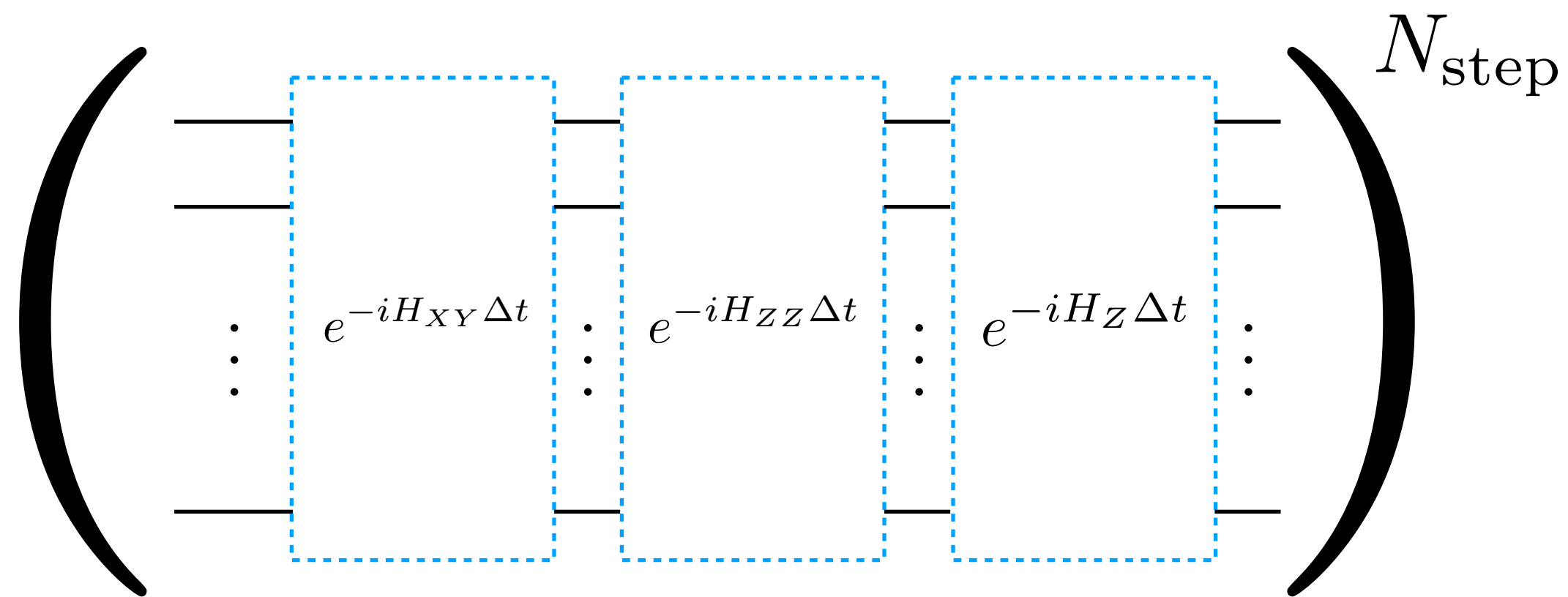
[Wecker, Hastings, Troyer,  
Phys. Rev. A, 92, 042303 (2015)]  
[Ho, Hsieh, SciPost Phys. 6, 029 (2019)]  
[Wiersema, *et. al.*  
PRX Quantum 1, 020319 (2020)]

- motivation: imitating Suzuki-Trotter decomposition of adiabatic or real-time evolution
- we use U(1) preserving decomposition

$$H = H_{XY}^{(\text{even})} + H_{XY}^{(\text{odd})} + H_{ZZ} + H_Z$$

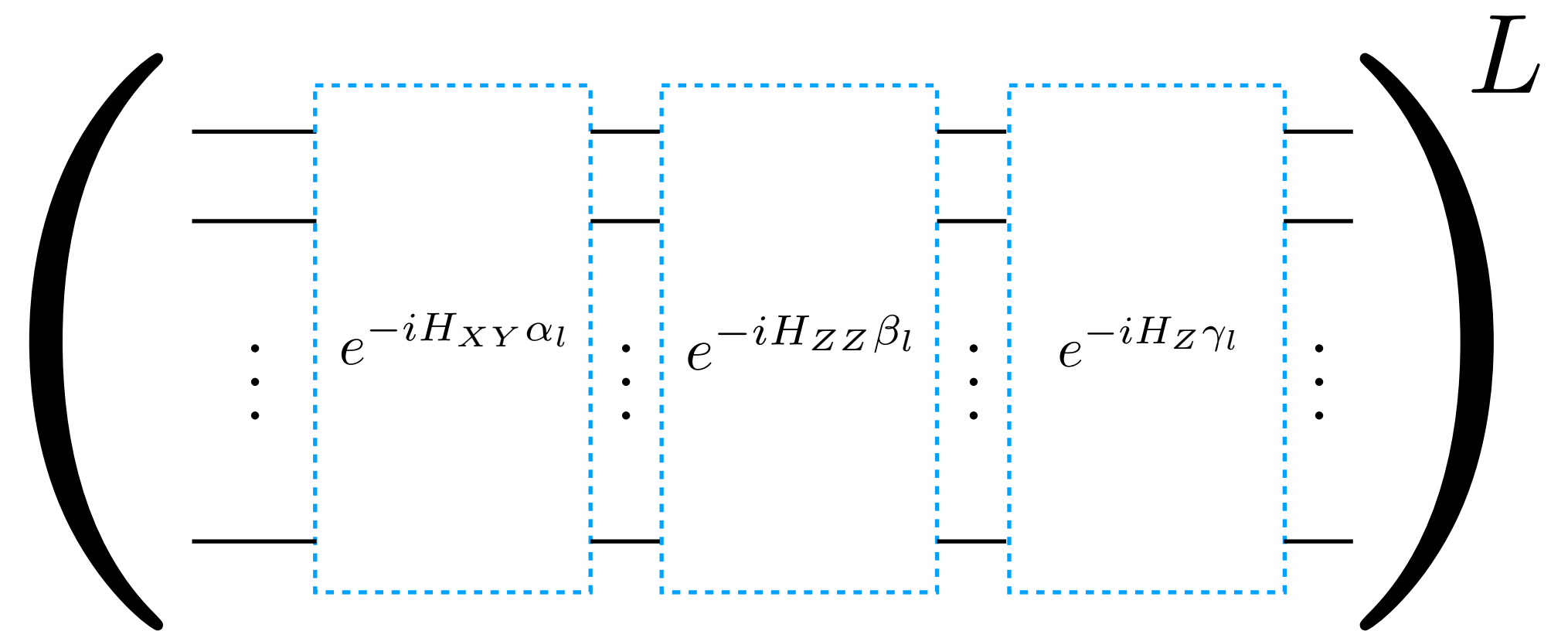
- replace  $\Delta t$  with variational parameters  $(\alpha_l, \beta_l, \gamma_l)$
- parameters  $(\alpha, \beta, \gamma)$  can depend on sites

Suzuki-Trotter evolution



$$\Delta t = T_{\text{max}}/N_{\text{step}}$$

HVA



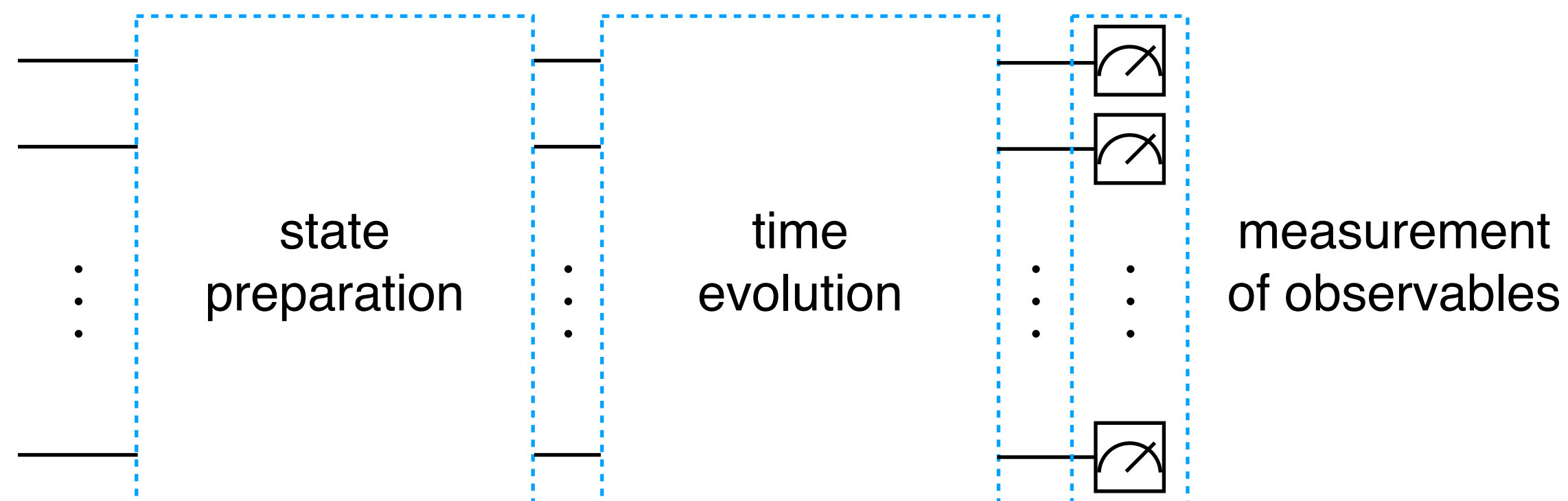
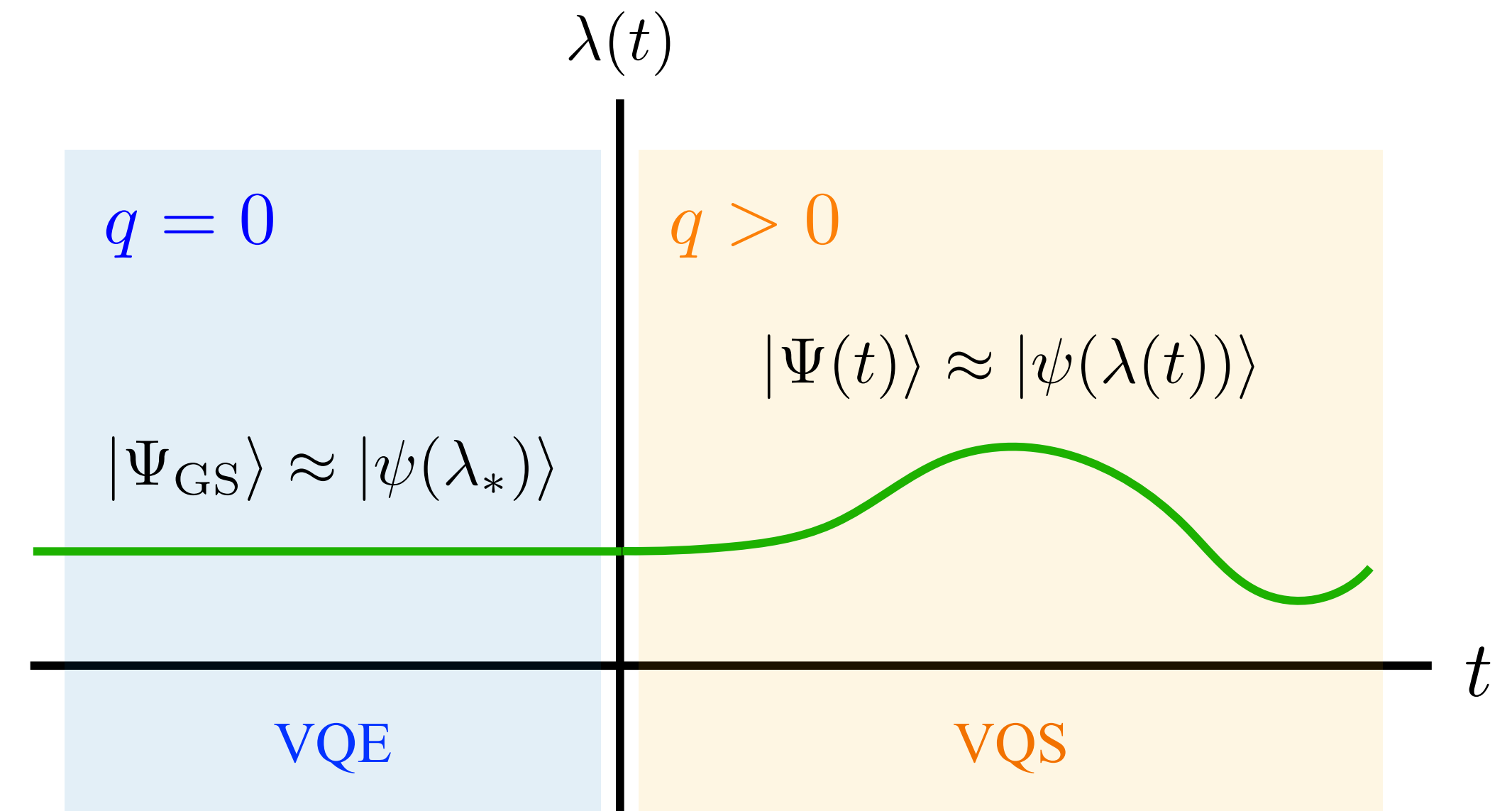
$(\alpha_l, \beta_l, \gamma_l)$  : variational parameters

# Summary of our protocol

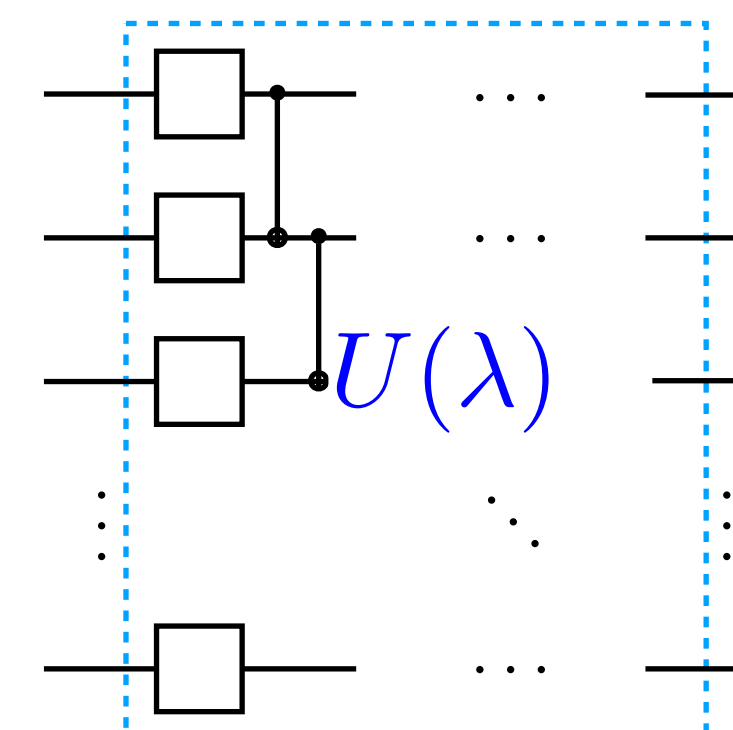
- Quench dynamics in the Schwinger model
  - ground state without external field  $q$ :  $|\Psi_{\text{GS}}\rangle$
  - time evolution via Hamiltonian with external field  $q$ :

$$|\Psi(t)\rangle = e^{-iH_{q\neq 0}t}|\Psi_{\text{GS}}\rangle$$

- perform VQE and VQS using the **same** ansatz  $|\psi(\lambda)\rangle$ 
  - reduce overall circuit depth
  - simulation with fixed depth



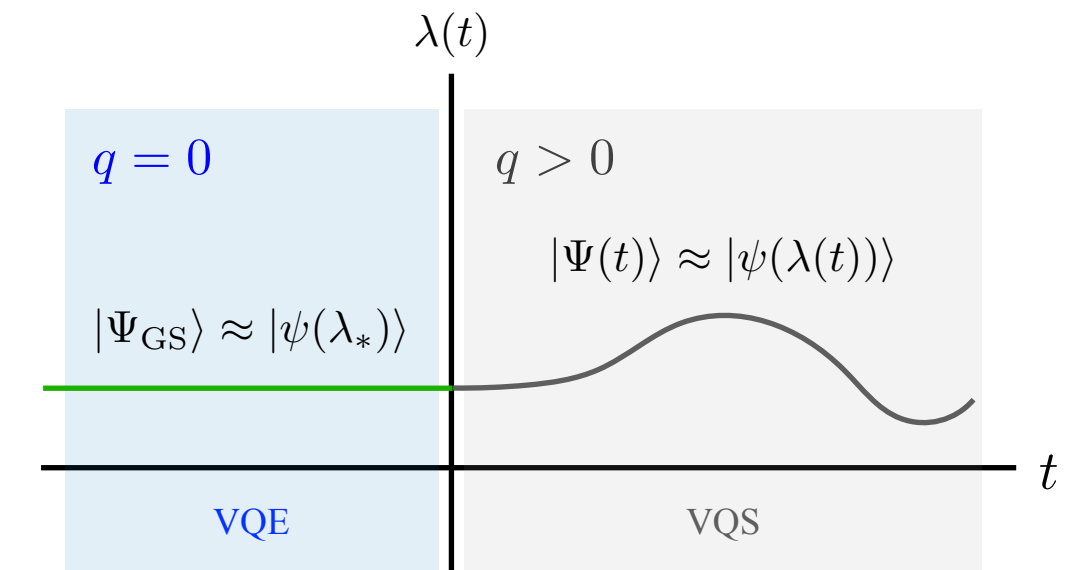
state prep./ time evolution



# Results

(classical) statevector simulation

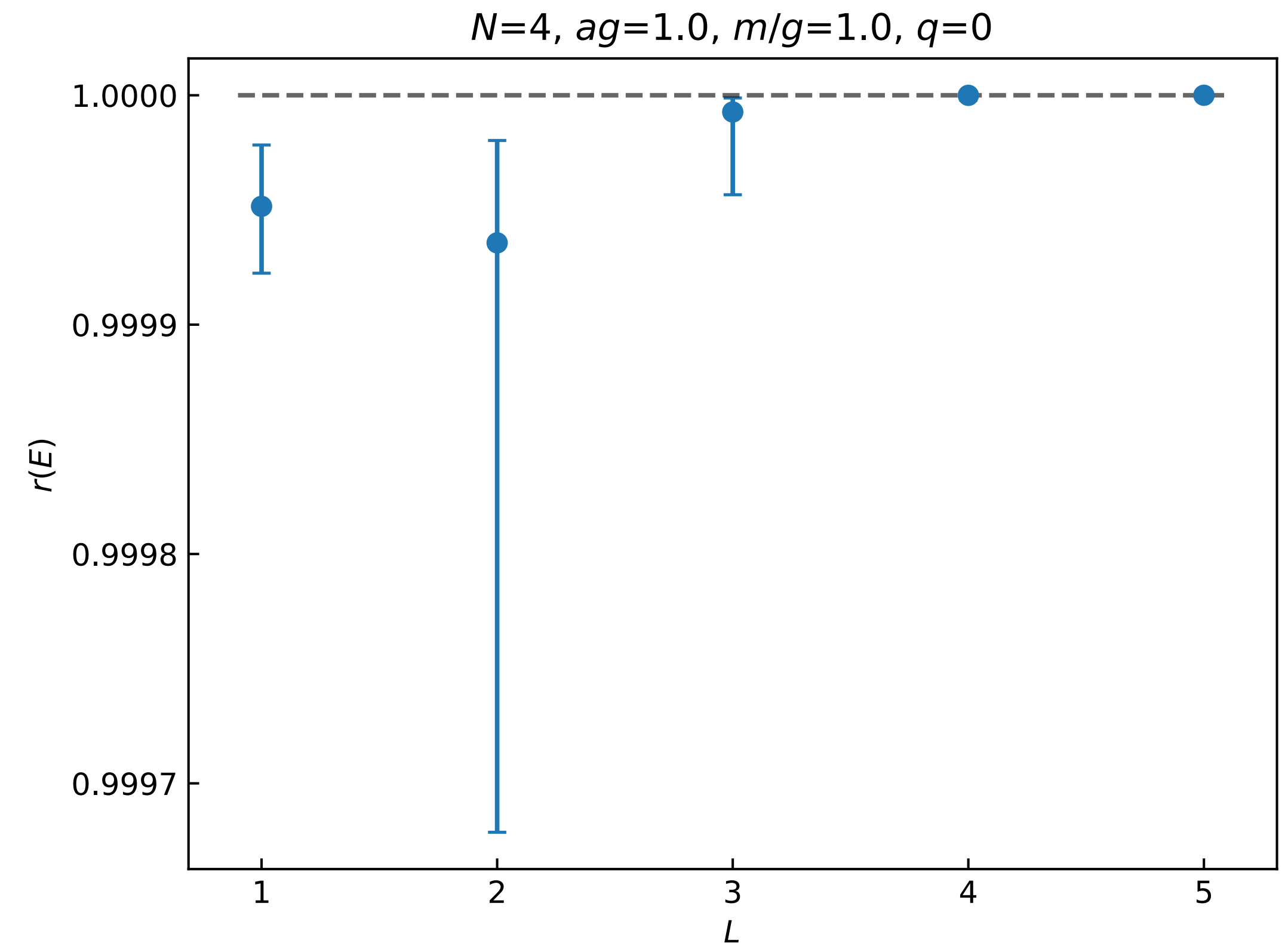
# Ground state preparation via VQE



- compare VQE results with exact diagonalization (ED)

- a metric of accuracy: 
$$r(E) = \frac{E_{\max} - E_{\text{VQE}}}{E_{\max} - E_{\min}}$$

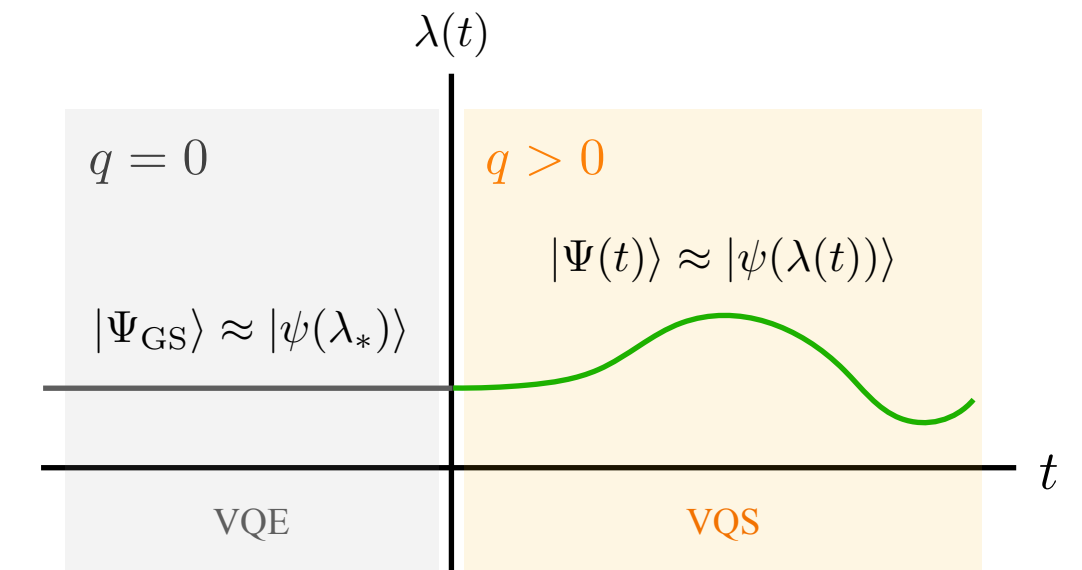
- $E_{\max}, E_{\min}$  : max/min energy obtained by ED
- $r(E) = 1$  for the best case
- $r(E) = 0$  for the worst case
- $L$  : depth of ansatz
- quality drastically improves for  $L \geq 4$



- 20 samples with different initialization
- dots/bars represent medians and 25-75 percentiles

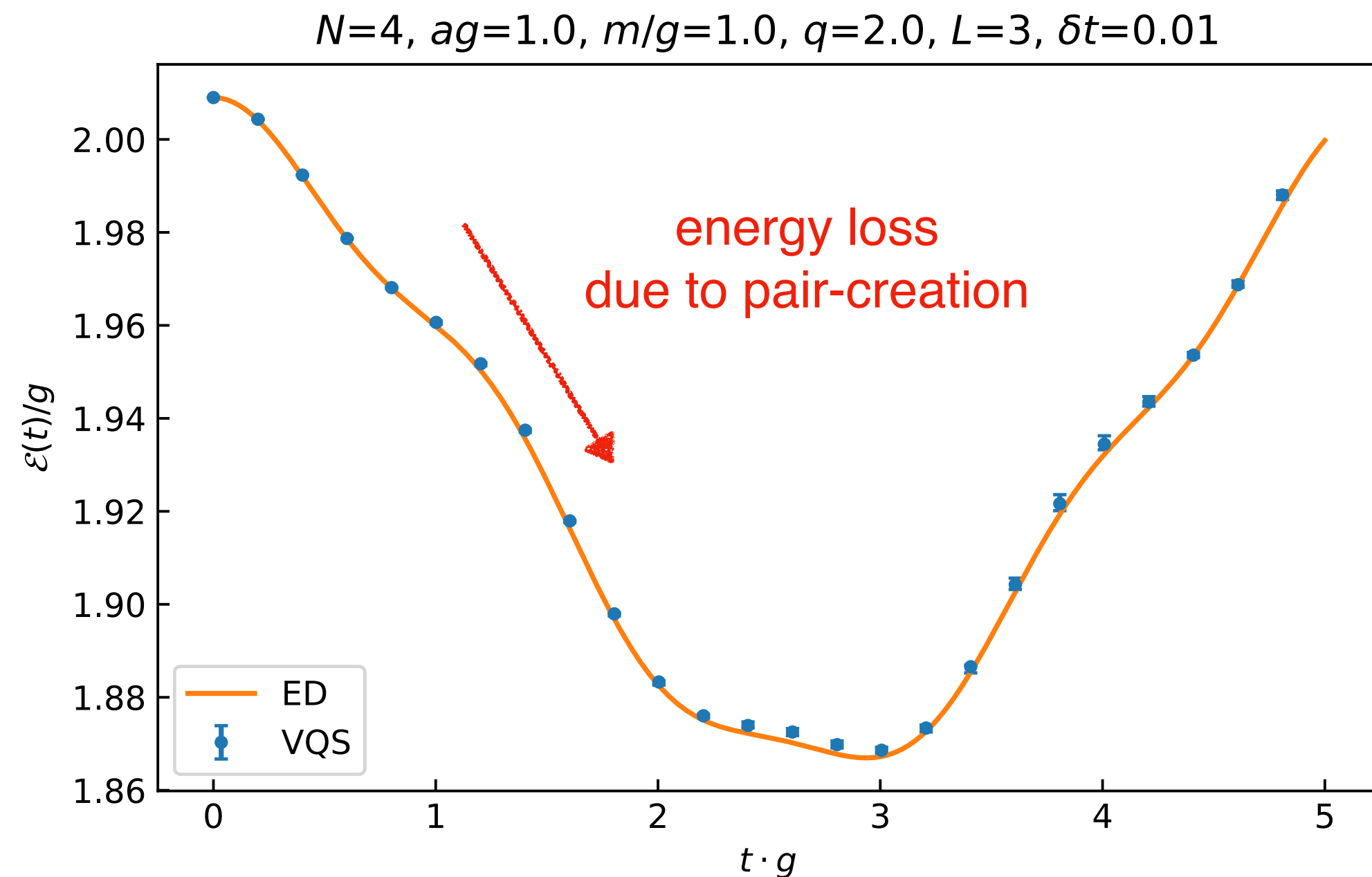
# Real-time evolution via VQS

- two observables:
  - total electric field  $\mathcal{E}$
  - chiral condensation  $\langle \bar{\psi} \psi \rangle$  ( $\sim$ particle number density)
- observing energy loss and pair-creation!

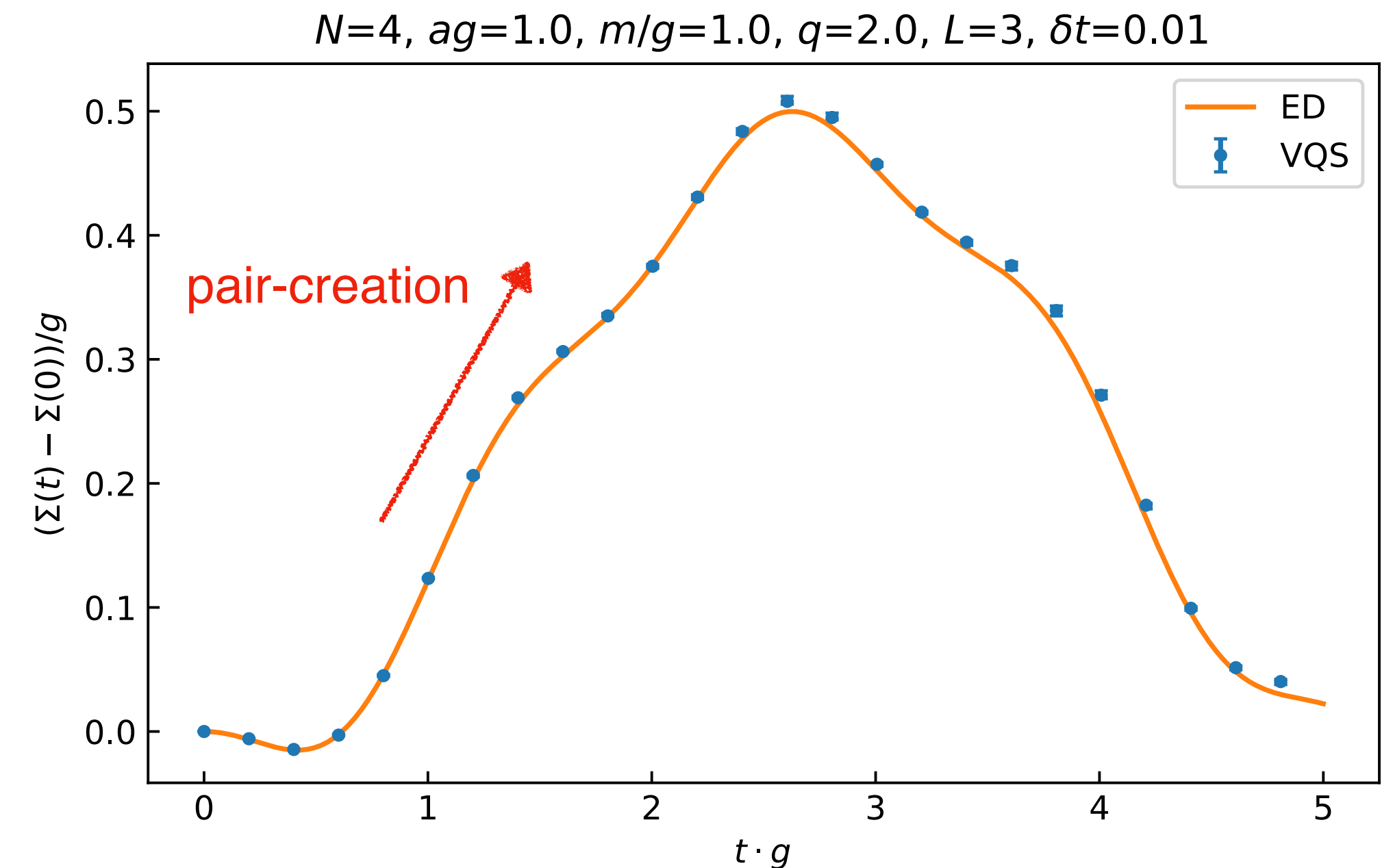


- 20 samples with different initialization
- dots/bars represent medians and 25-75 percentiles

electric field



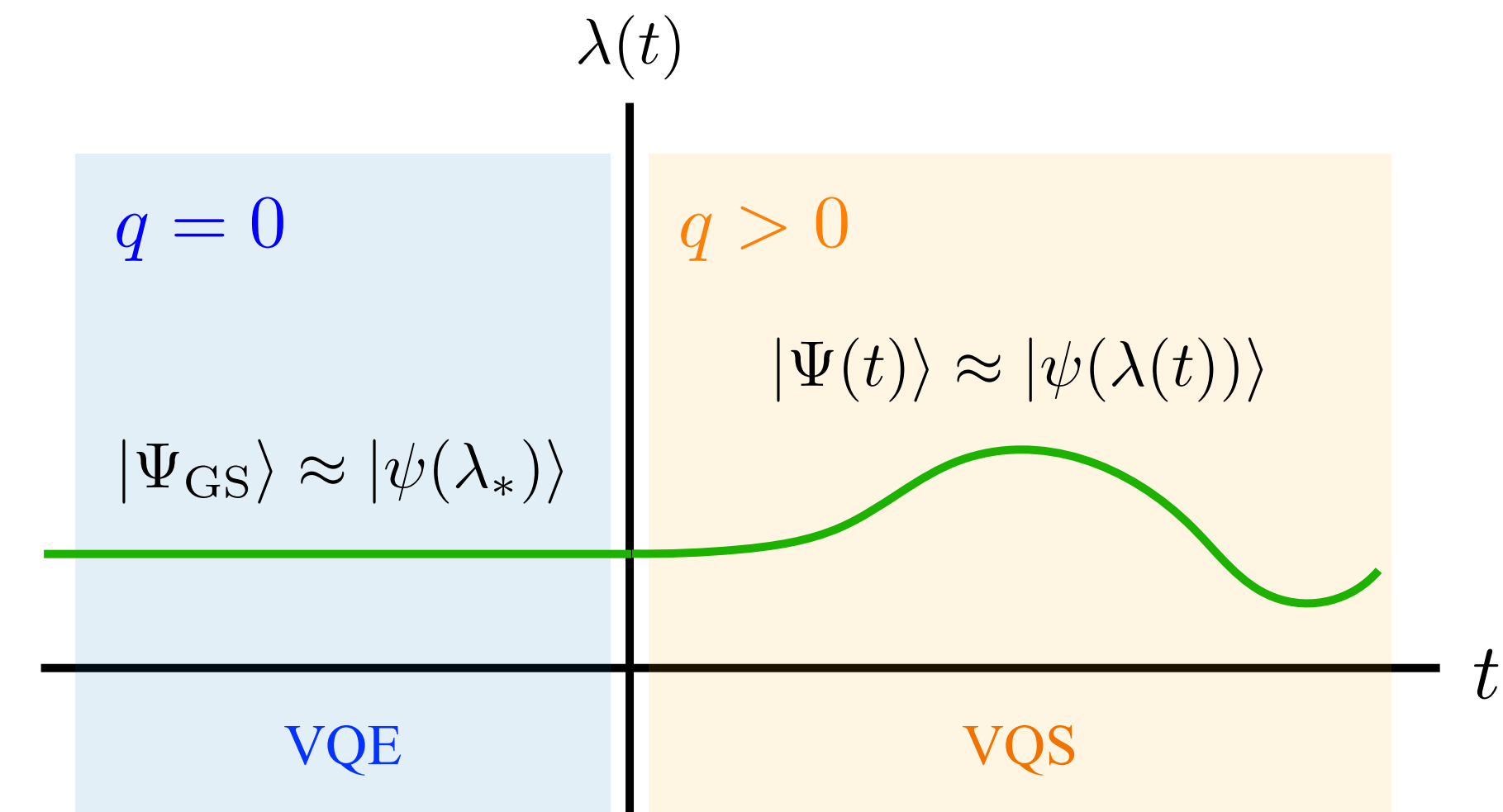
chiral condensation





# Summary and outlooks

- quench dynamics in the Schwinger model via VQAs
  - ground state w/o external field  $q$  via VQE
  - time evolution via Hamiltonian w/ external field  $q$  via VQS
  - we can reduce circuit depth
  - VQA results agree well with ED
- **future directions:**
  - understanding scaling of resources
  - effects of errors (shot noise, quantum noise)
  - reducing sampling cost [in progress]
  - extension to higher dimensional and/or non-Abelian theories



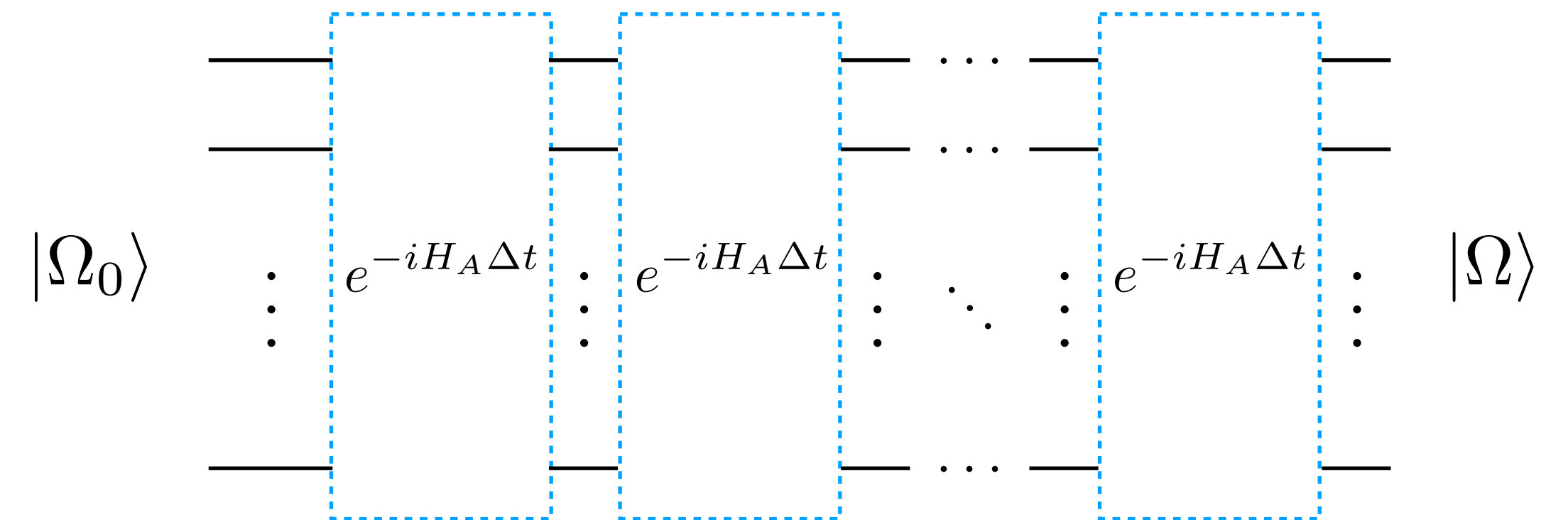
**Backups (including preliminary results)**

# Suzuki-Trotter decomposition

- adiabatic state preparation

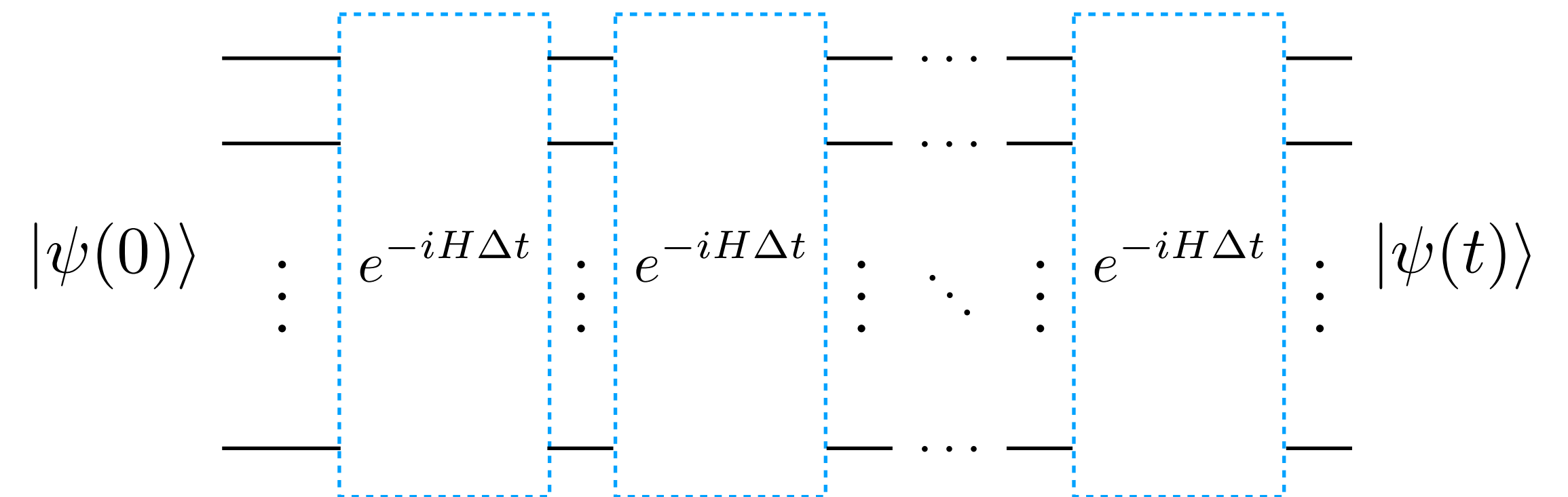
$$|\Omega\rangle = \lim_{T \rightarrow \infty} \text{T exp} \left( -i \int_0^T dt H_A(t) \right) |\Omega_0\rangle$$

$$\simeq \prod_s e^{-iH_A(s\Delta t)\Delta t} |\Omega_0\rangle$$



- real-time evolution

$$|\psi(t)\rangle = e^{-iHt} |\psi(0)\rangle = (e^{-iH\Delta t})^s |\psi(0)\rangle$$



- drawback: #depth grows with #steps

# McLachlan's variational principle

$$\delta \left\| \left( \frac{d}{dt} + iH \right) |\psi(\lambda)\rangle \right\| = 0$$

$$\Rightarrow \sum_j M_{ij} \dot{\lambda}_j = V_i$$

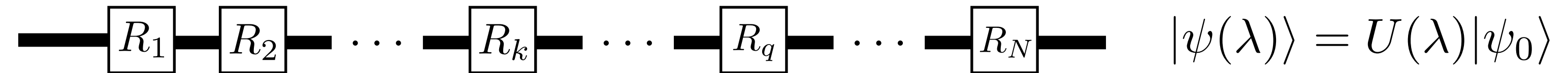
$$M_{ij} = \operatorname{Re} \frac{\partial \langle \psi(\lambda) |}{\partial \lambda_i} \frac{\partial |\psi(\lambda)\rangle}{\partial \lambda_j}$$

$$V_i = \operatorname{Im} \frac{\partial \langle \psi(\lambda) |}{\partial \lambda_i} H |\psi(\lambda)\rangle$$

# Quantum circuit for VQS

[Li, Benjamin, Phys. Rev. X 7, 021050 (2017)]  
 [Yuan et al., Quantum 3, 191 (2019)]

$$U(\lambda) = R_N(\lambda_N) \cdots R_1(\lambda_1)$$



- evaluation of matrix elements  $M_{kq} = \text{Re} \frac{\partial \langle \psi(\lambda) |}{\partial \lambda_k} \frac{\partial |\psi(\lambda)\rangle}{\partial \lambda_q}$

- derivative of each component w.r.t. parameters  $\frac{\partial}{\partial \lambda_k} R_k(\lambda) = U_k R_k(\lambda)$

- quantum circuit:

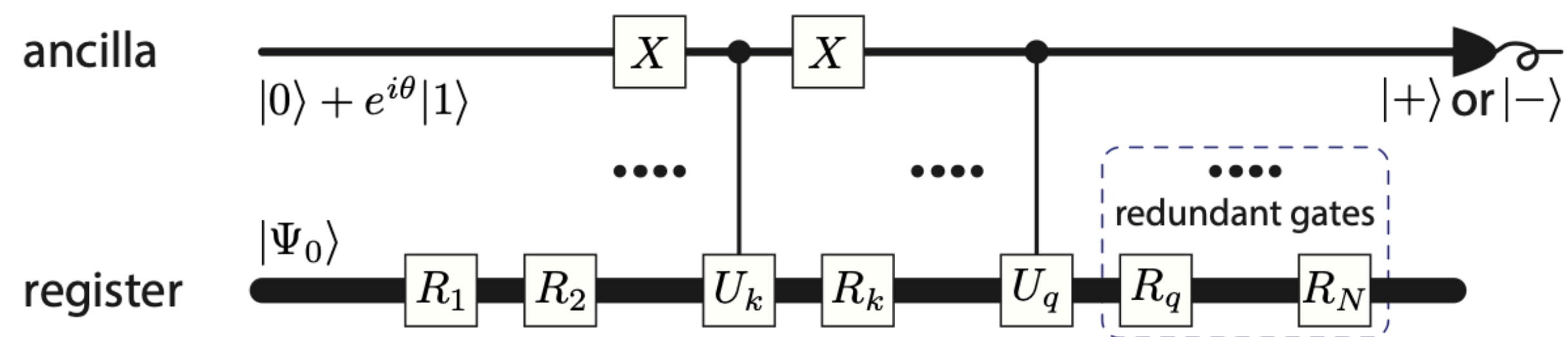
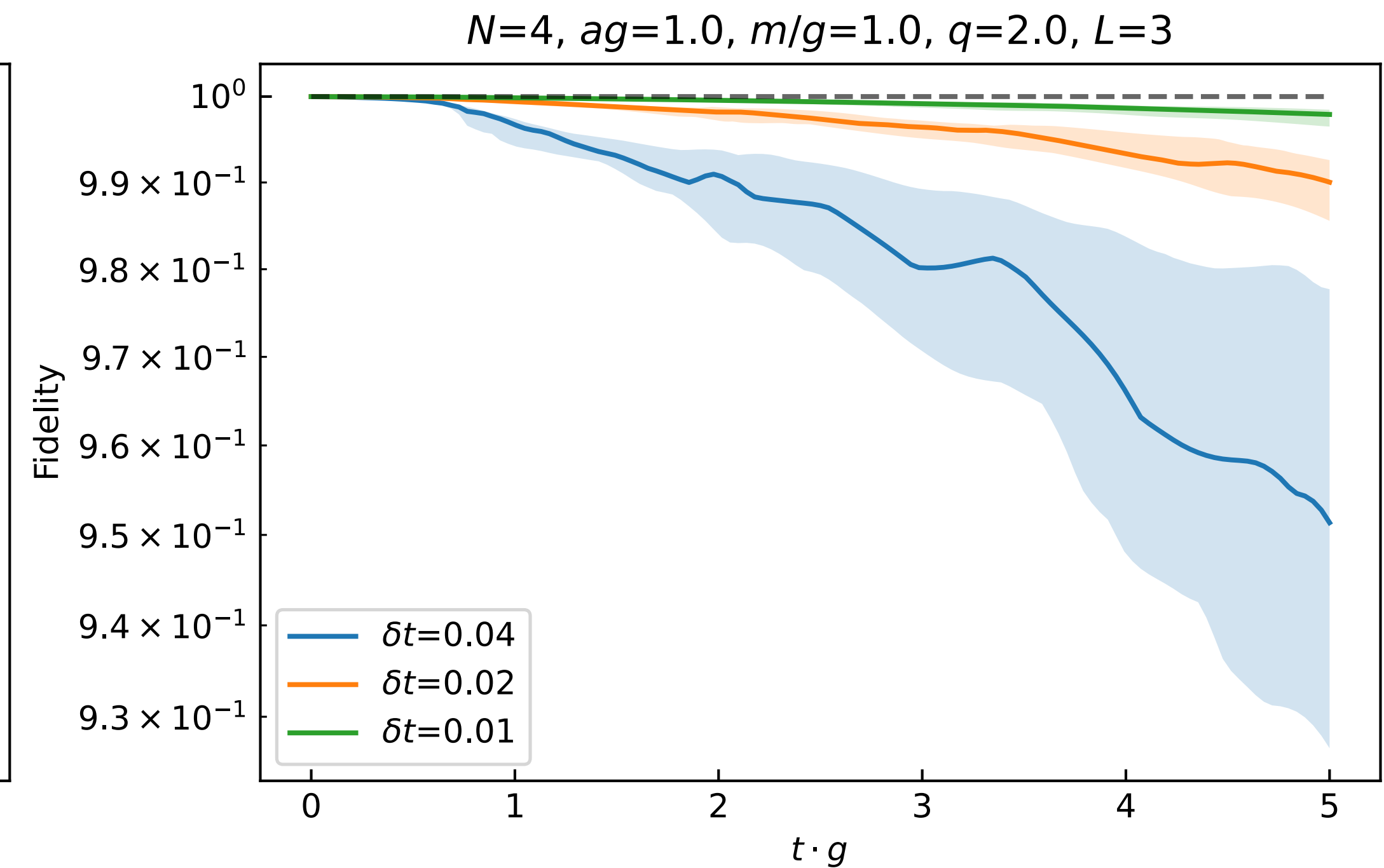
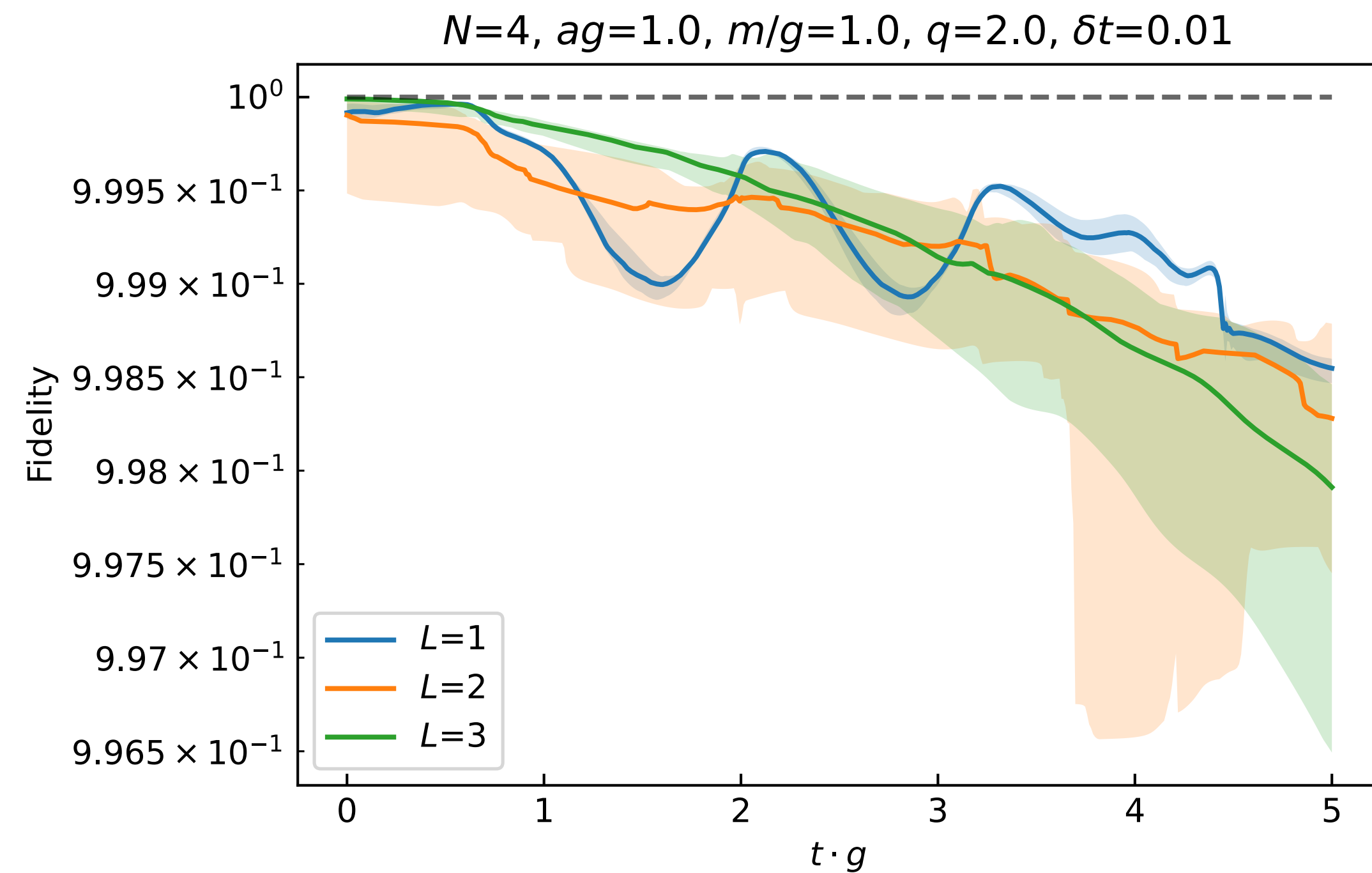


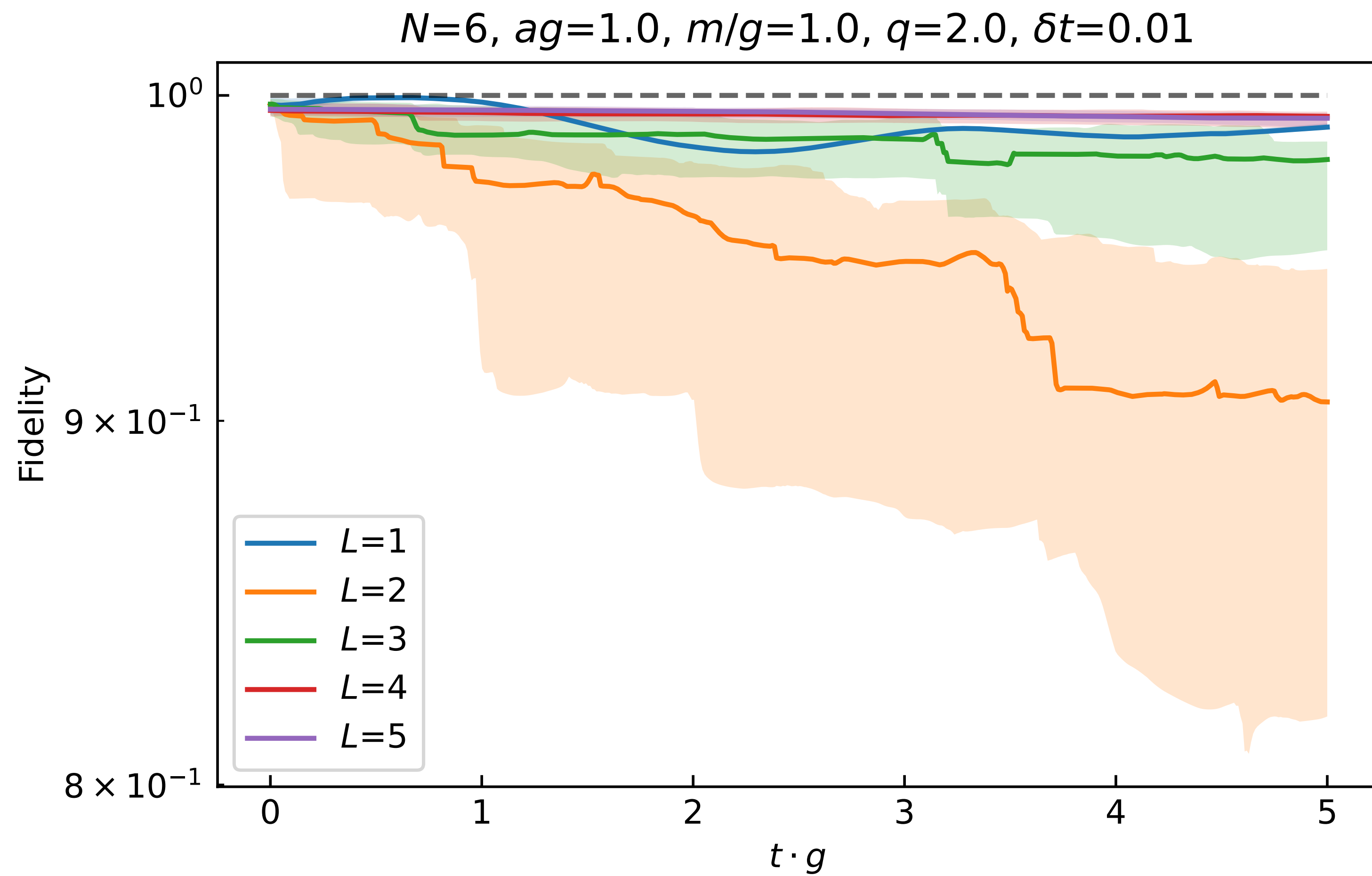
Figure from Yuan et al.

# Fidelity and algorithmic errors

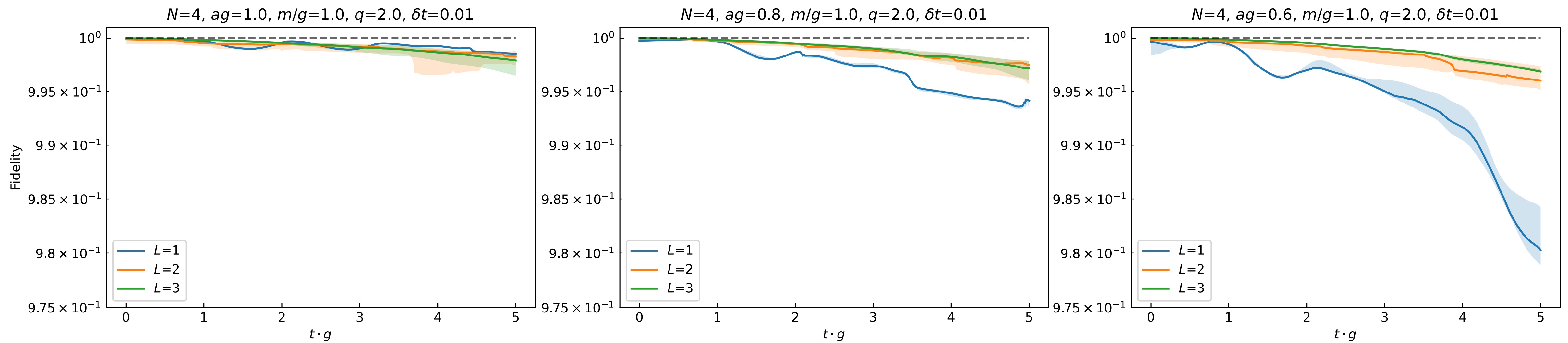
- (averaged) fidelity improves as increasing  $L$  and/or decreasing  $\delta t = T_{\max}/N_{\text{step}}$
- effects from  $\delta t$  is significant



# $L$ -dependence ( $N = 6$ )



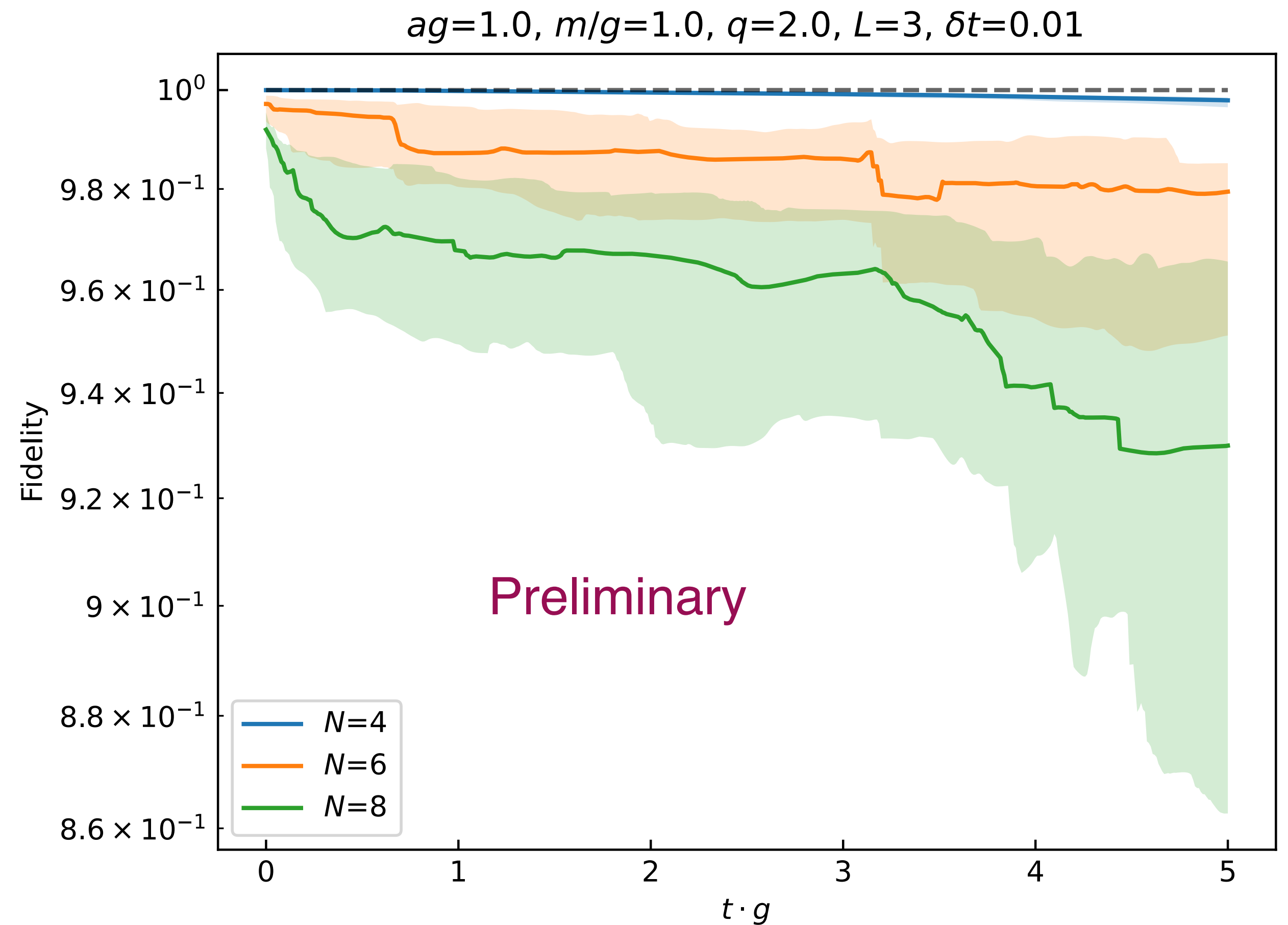
# Lattice spacing dependence



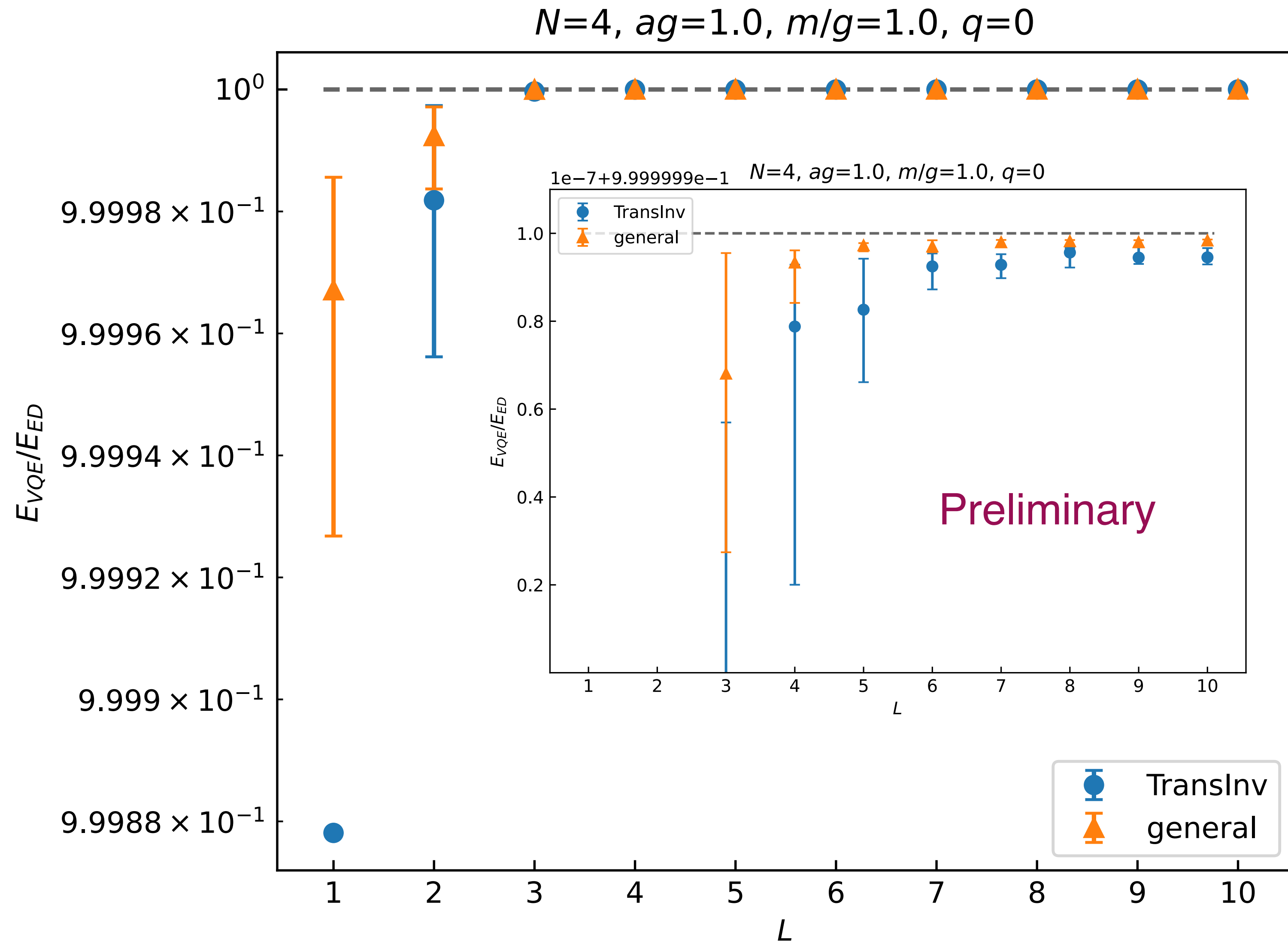


# Comments on scalability

- results get worse with increasing  $N$ ...
  - VQE accuracy (#iteration, optimizer)
  - VQS accuracy (choice of  $\delta t$ ,  $L$ )
- at least  $N \leq 8$ ,  $F > 0.9$  with  $L = 3$

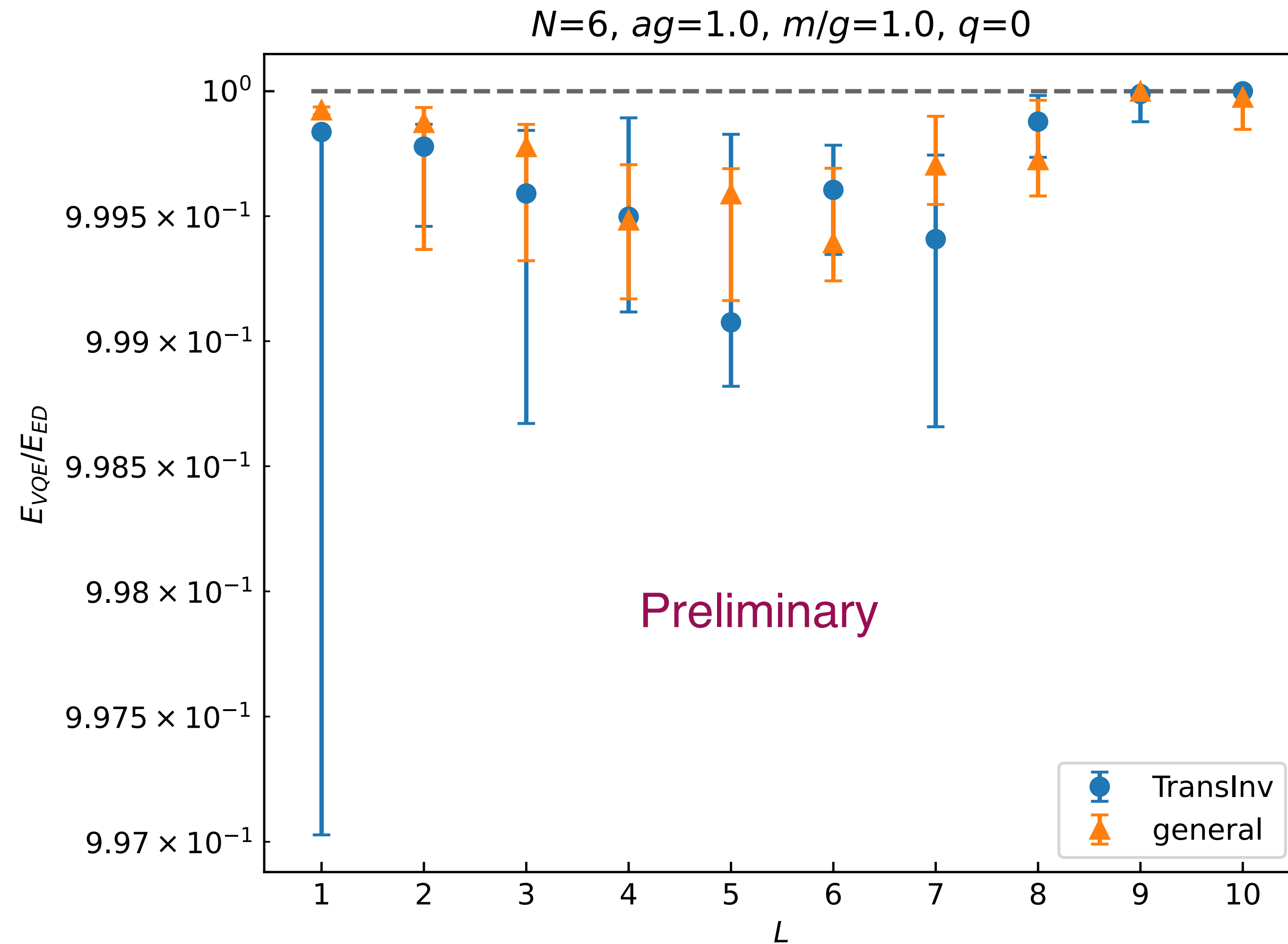


# Translational invariant parameters ( $N = 4$ )



#iter=5,000

# Translational invariant parameters ( $N = 6$ )



#iter=5,000