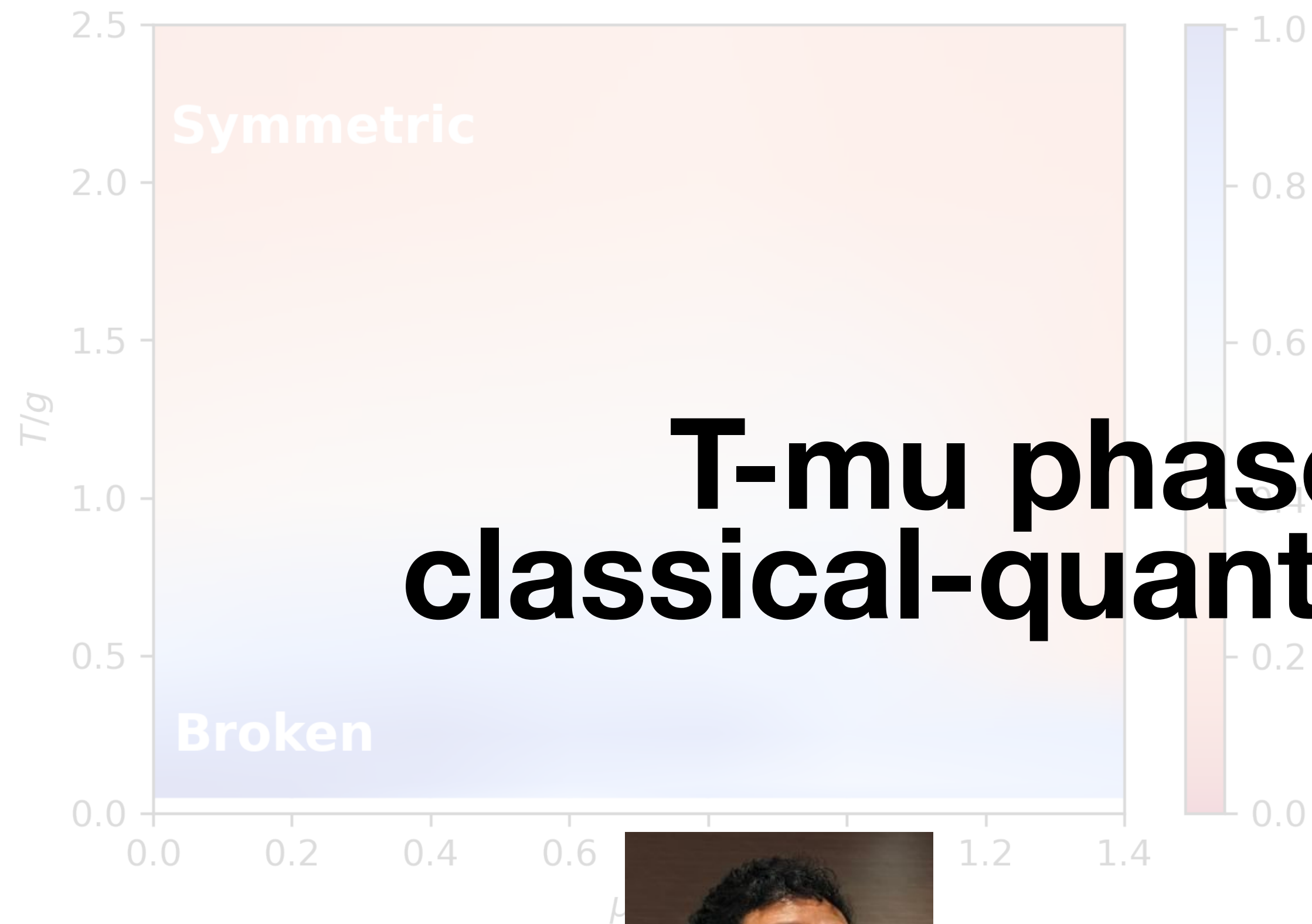


T-mu phase diagram using classical-quantum hybrid algorithm



Akio Tomiya
(Assistant Prof in IPUT Osaka)

Based on arXiv: [2205.08860](https://arxiv.org/abs/2205.08860) →



NEWS: A new research grant (Japan governmental) initiated!

Advertisement

“Machine Learning Physics Initiative”

2022-2027, 10M USD, 70 researchers

Director : K. Hashimoto

A.Tanaka: Math and Application of DL

Y.Kabashima: Statistical data ML

K.Fukushima: Topology and Geometry of ML

A.Tomiya: Computational physics

M.Nojiri: High Energy Physics

T.Ohtsuki: Condensed Matter Physics

K.Hashimoto: Quantum and Gravity Physics

ML
Phys

FY2022-2026 MEXT -KAKENHI- Grant-in-Aid for Transformative Research Areas (A)

科研費
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Assistant professor with a fixed term appointment at University of Tsukuba (Division of Particle Physics) in Japan

Research field, Content of work:

Computational particle physics. In collaboration with Prof. Hiroshi Ohno, the successful candidate will conduct research on lattice field theories incorporating the machine learning approach related to the research project “Fusion of Computational Physics and Machine Learning” supported by Grant-in-Aid for Transformative Research Areas (A)

(Principal Investigator: *Akio Tomiya*). The applicant will be expected to have expertise, skills, and experience on one of the followings:

- 1) *lattice field theory* and related numerical computations,
- 2) *machine learning*, especially generative models, and related numerical computations.

Deadline: Sep, 30

URL: <https://www.ccs.tsukuba.ac.jp/recruit-20220930e/>



WE WANT YOU!

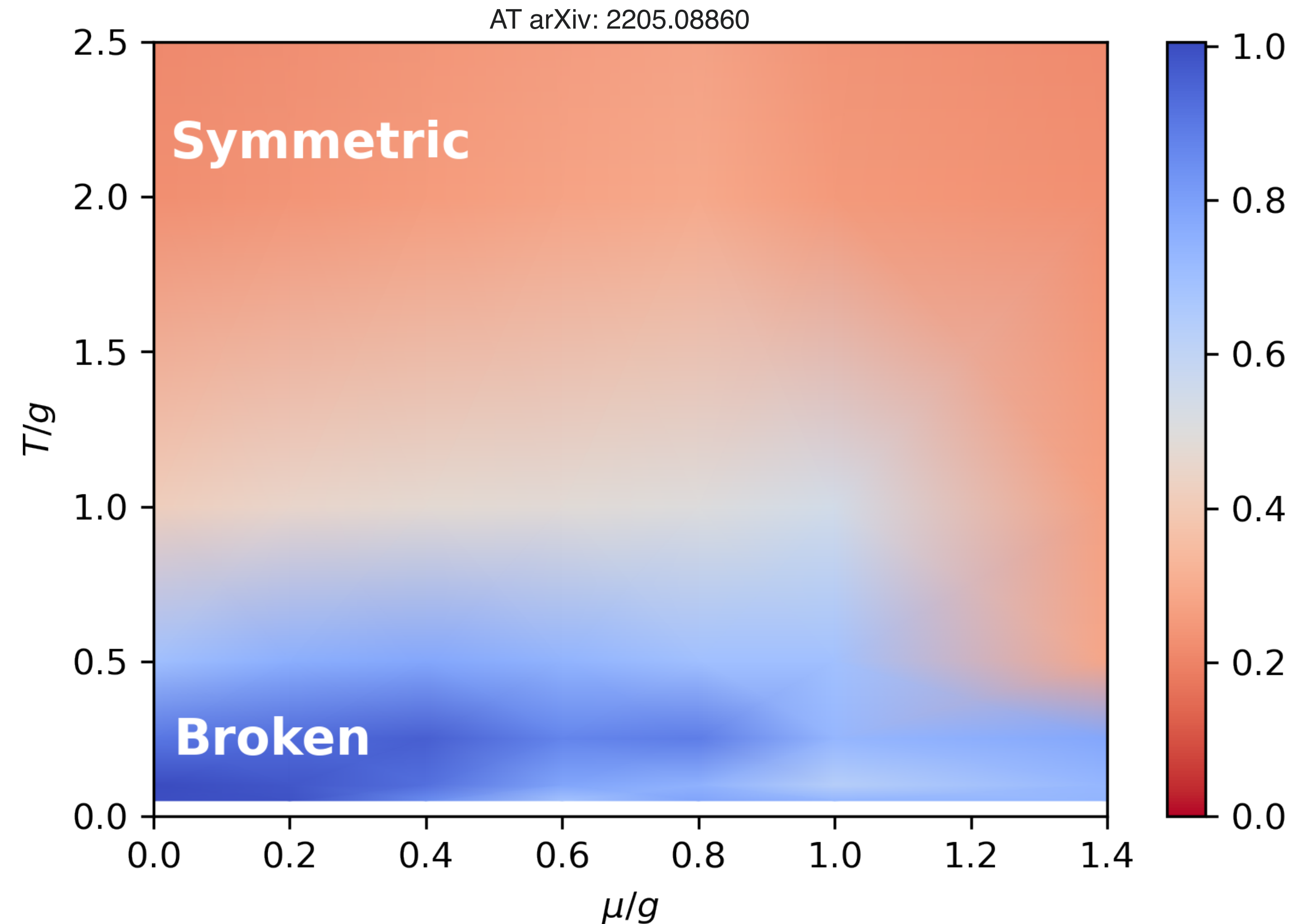
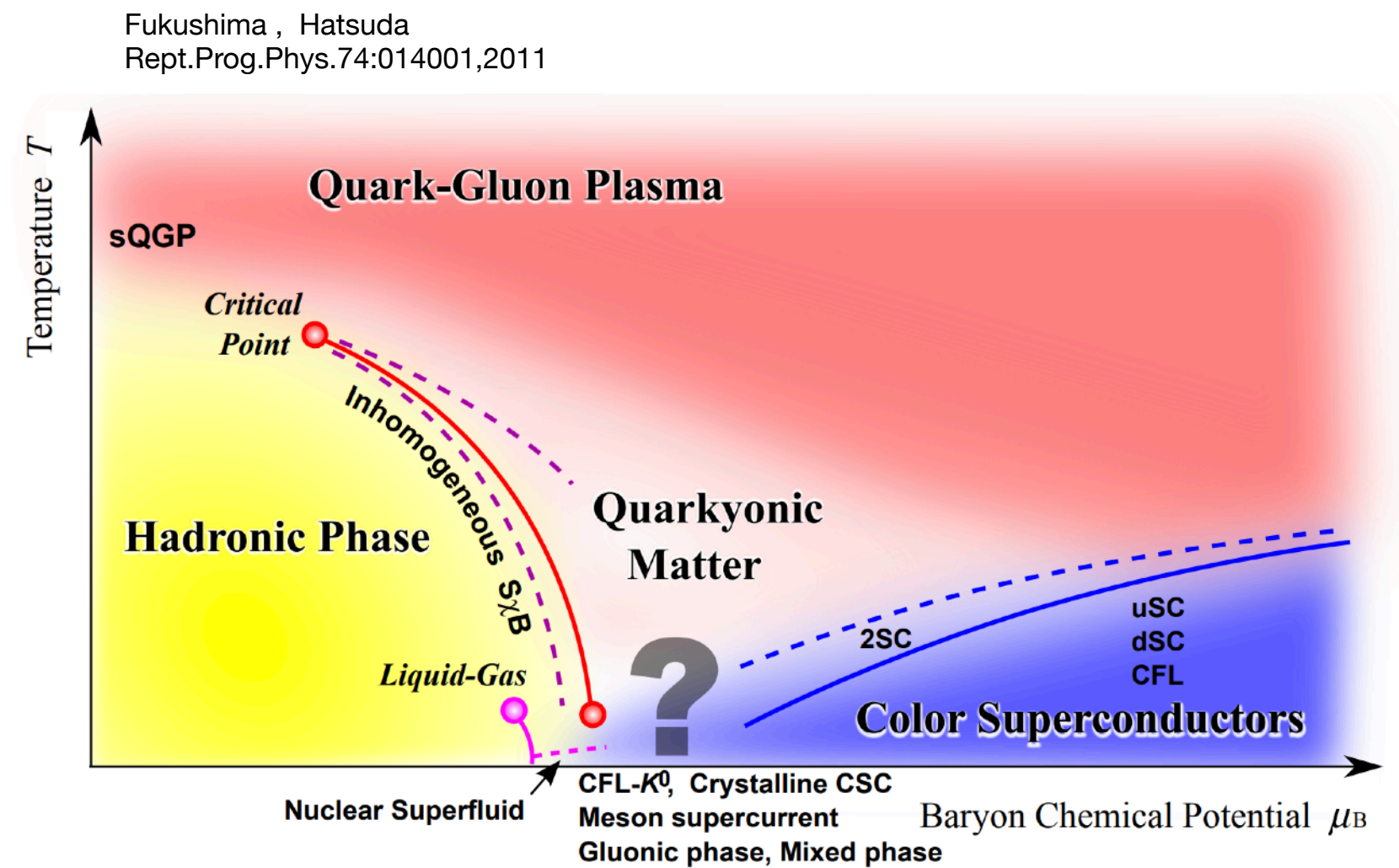


- 1.Introduction, motivation
- 2.VQE (a classical-quantum hybrid algorithm)
- 3.Density matrix, KL-U divergence
- 4.Beta VQE (VQE for $T>0$)
- 5.Results
- 6.Summary

Introduction

Summary of this talk

Hybrid algorithm = Quantum + “machine learning”



I investigated T- μ phase diagram using a quantum algorithm & neural network (β -VQE, No sign problem) for Schwinger model (toy model of QCD)

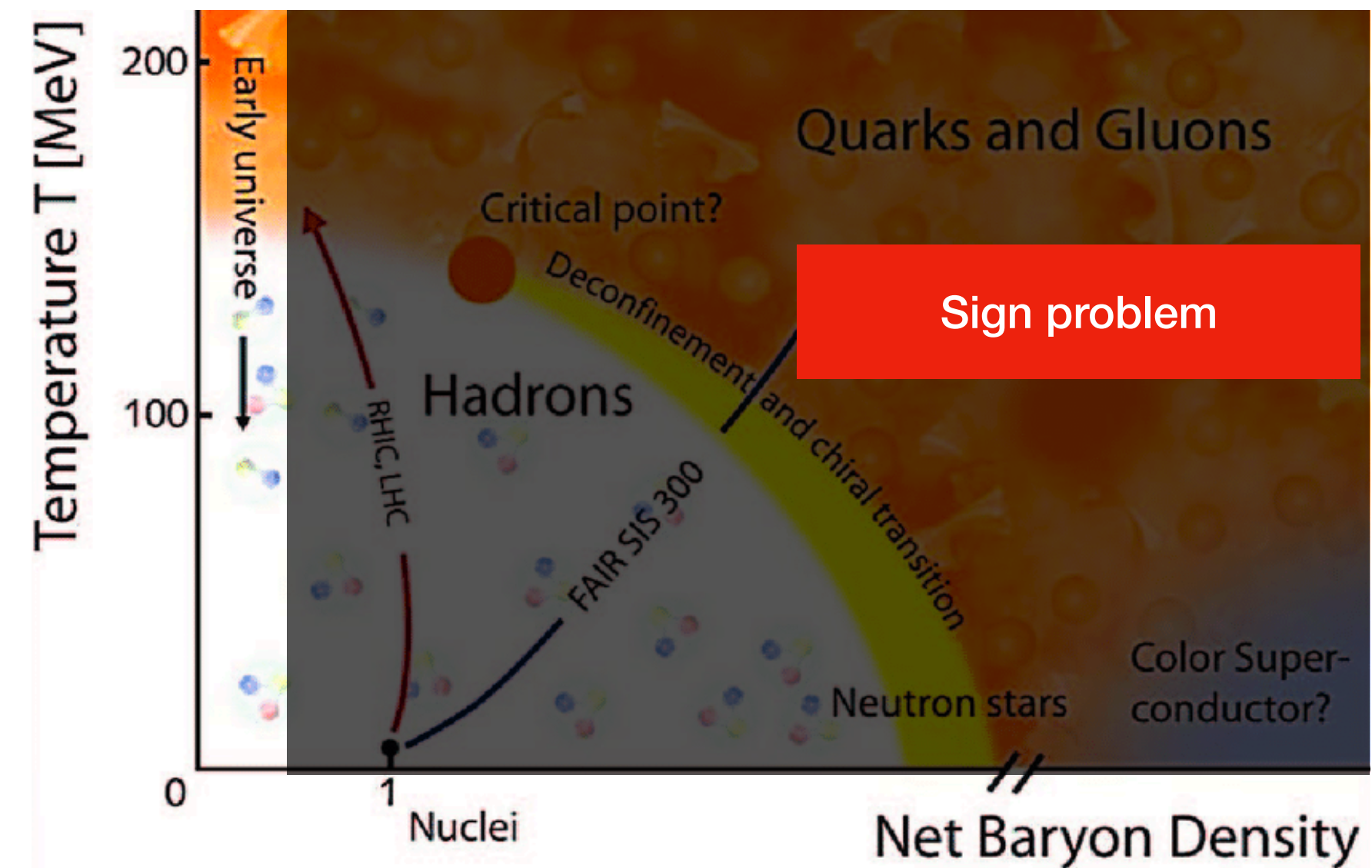
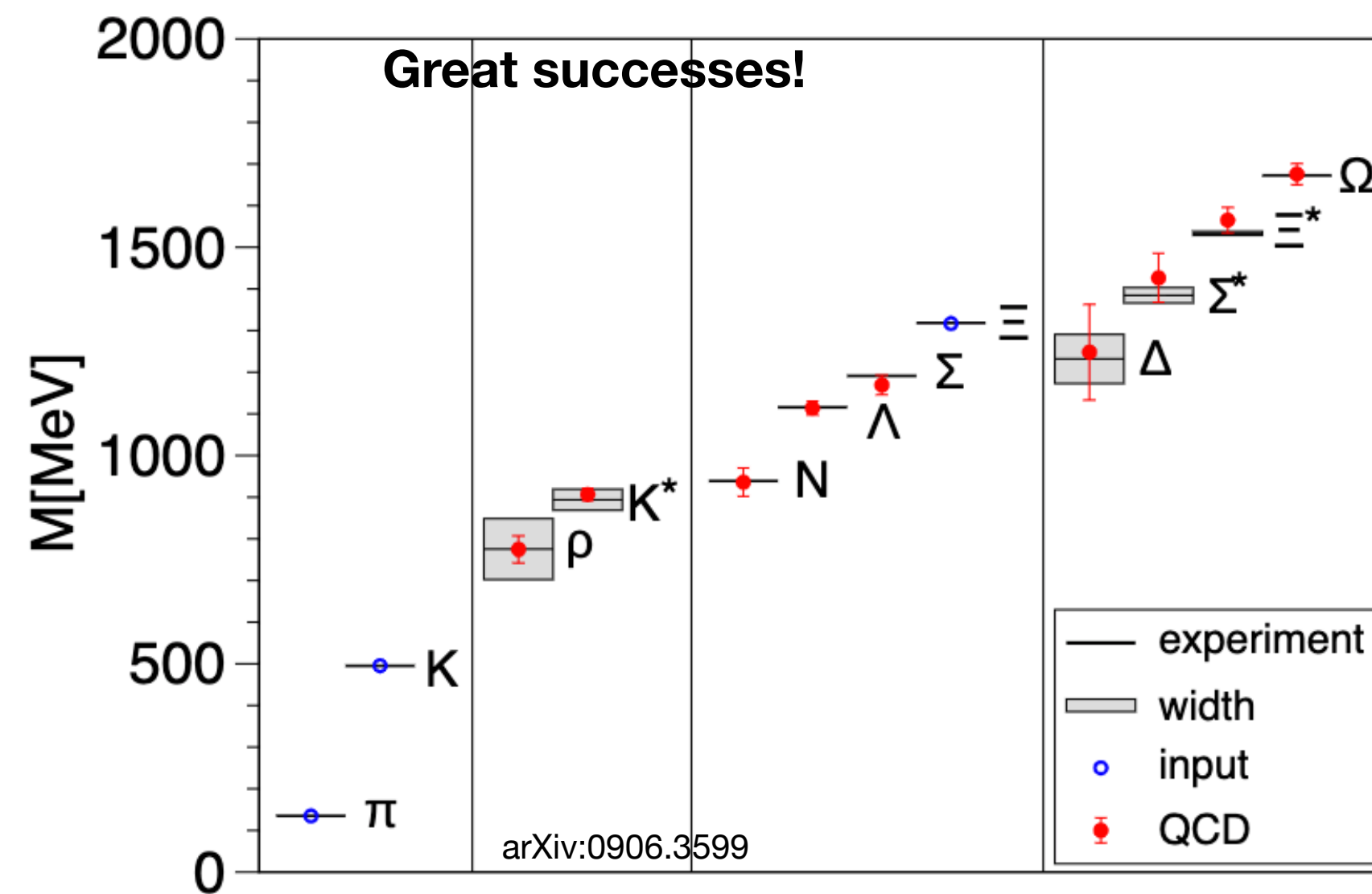
Introduction

MCMC is powerful, if $\mu = 0$

- Monte-Carlo enables us to evaluate expectation values for “statistical system”, like lattice QCD in imaginary time

$$\langle O[U] \rangle = \frac{1}{N_{\text{conf}}} \sum_c^{N_{\text{conf}}} O[U_c] + \mathcal{O}\left(\frac{1}{\sqrt{N_{\text{conf}}}}\right)$$

$$U_c \leftarrow P(U) = \frac{1}{Z} e^{-S[U]} \in \mathbb{R}_+$$



- If we turn on **the baryon number density μ** , **Monte-Carlo methods do not work** because $e^{-S[U]}$ becomes complex. This is no more probability. (sign problem)
- Operator formalism does not have such problem! But it requires huge memory to store quantum states, which cannot be realized even on supercomputer.
- Quantum states should not be realized on classical computer but on quantum computers, as Feynman said.**

Introduction

Finite T is good for classical machine. Finite mu is good for quantum

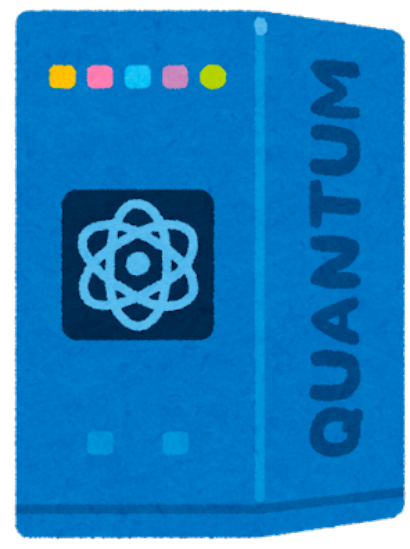


Classical machine: Lattice field theory calculations rely on

$$P(U) = \frac{1}{Z} e^{-S[U]} \det(D[U] + m)^2 \in \mathbb{R}_+$$

Since 1980 (M. Creutz)~

- This P cannot be regarded as probability if $\mu \neq 0$ (sign problem)



Quantum machines can realize (any) unitary evolutions (Solovay Kitaev theorem),

$$U(t) = e^{-i\hat{H}t} \quad : \text{unitary}$$

B. Chakraborty, AT+ *Phys.Rev.D* 105 (2022) 9, 094503 and references therein

- No problem for $\mu \neq 0$ because we can only use unitary gates (operators)
- “Short time evolution” (shallow circuit) is preferred for near-term devices
- (Efficient way of) calculation of non-unitary cases (i.e. Boltzmann weight)?

	Classical Computers	Quantum Computers
Finite Density	Sign Problem	✓
Finite Temperature	✓	Challenging

Credit: Connor Powers (U. of Maryland)*

We need a method to calculate $T>0$ and $\mu \neq 0$ for QCD and for near-term quantum devices!

*<https://indico.hiskp.uni-bonn.de/event/40/contributions/484/attachments/358/630/Powers%20Talk%20Final%20Draft%20Updated.pdf>

State preparation, VQE

State preparation, VQE

Realization of the exact ground state is difficult

Operator formalism

H : Hamiltonian in QFT $[A_\mu, E_\nu] = i\hbar\delta_{\mu\nu}\dots$

Minkowski in M^{d+1}

$$U(t) = e^{-itH}$$

AT+ *Phys.Rev.D* 105 (2022) and a lot!



Real time

Euclid($t \rightarrow \tau$)

$t = -i\tau$

Minkowski($\tau \rightarrow t$)

Finite temperature/imaginary time

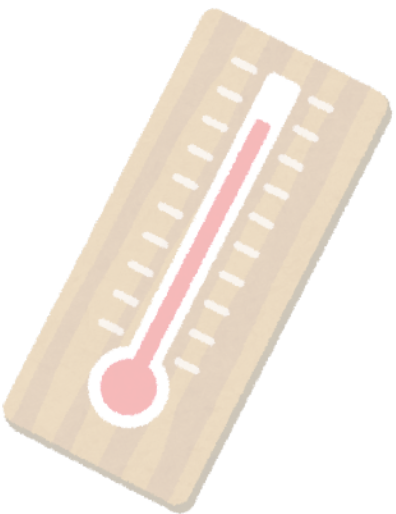
$$U(\tau) = e^{-\tau H}$$

Euclid in $S^1 \times M^d$

(This work)

$$\langle OO(\tau) \rangle = \text{Tr}[O(0)O(\tau)\rho]$$

$$\rho = U(\tau)/Z$$



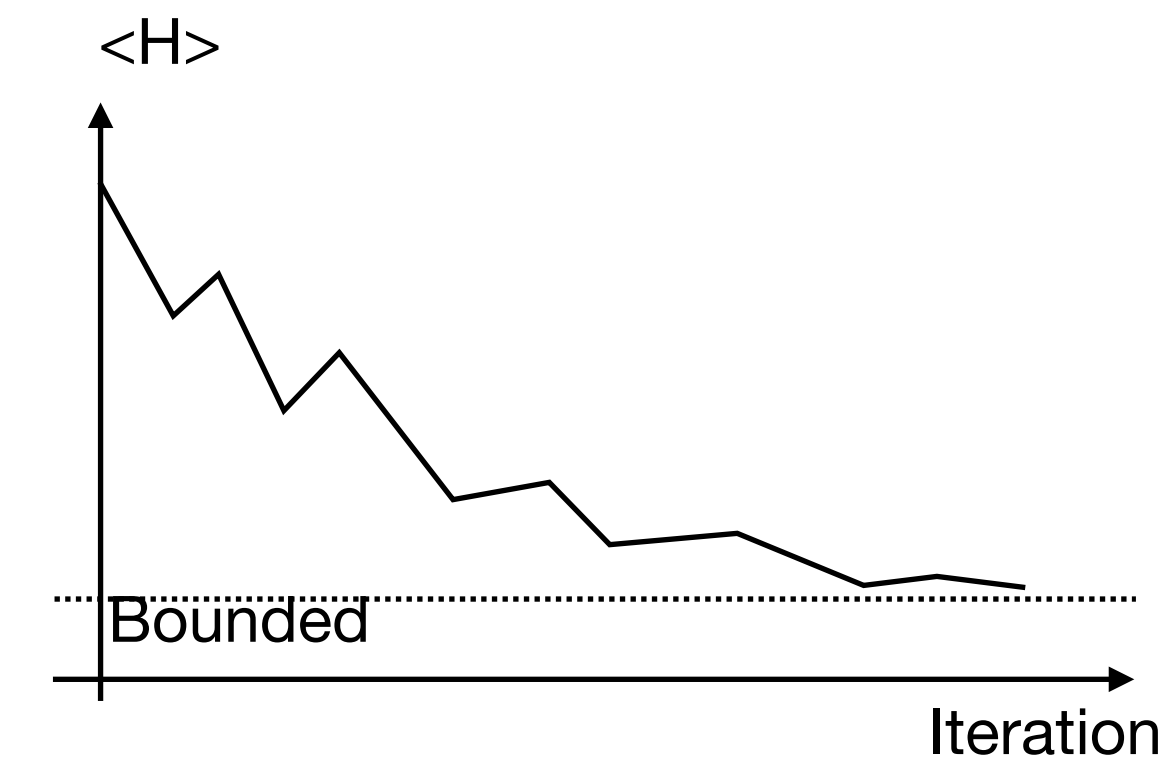
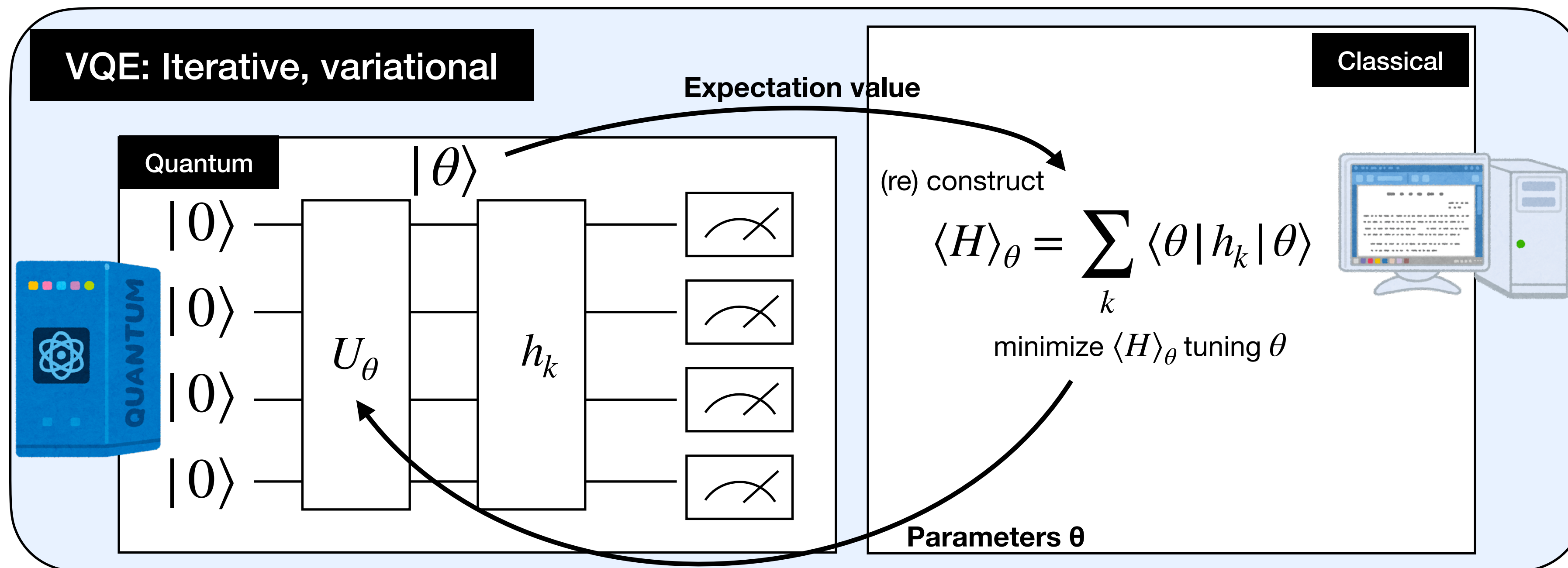
- Typical use of quantum algorithm is real-time simulation (\because Unitary).
 - Main interest: $\langle \Omega | O | \Omega \rangle$, where $|\Omega\rangle$ is the exact ground state
- Difficulty: State preparation* (as ground state extraction in conventional LQCD)
 - > Variational ansatz (next)
- Our work uses the operator formalism. No problem for $\mu > 0$
- Variational calculations with thermal state? -> Beta VQE

* very good summary in a talk by Alexei Bazavov <https://indico.hiskp.uni-bonn.de/event/40/contributions/469/>

State preparation, VQE

Mimic the ground state with a parametrized state

- Quantum machine: Exact ground state preparation is hard. In particular, it is difficult on near term devices
- **Variational method for a pure state** with a short circuit (VQE, variational quantum eigen-solver).
 - Quantum/Classical hybrid algorithm, iterative, variational
 - Parametrized **unitary** circuit (~parametrized wave-function $|\theta\rangle$, θ : a set of parameters)
- Basically, it mimics the ground state (pure state)



- Pros: Cheap. NISQ
- Cons: Systematic error (opposite to adiabatic ones)

State preparation, VQE

Thermal ... Boltzmann weight is not unitary...

Operator formalism

H : Hamiltonian in QFT

$$[A_\mu, E_\nu] = i\hbar\delta_{\mu\nu}\dots$$

Real time

Finite temperature/imaginary time

(This work)

Minkowski in M^{d+1}

Euclid in $S^1 \times M^d$

$$U(t) = e^{-itH}$$

$$U(\tau) = e^{-\tau H}$$

Euclid($t \rightarrow \tau$)

$t = -i\tau$

Minkowski($\tau \rightarrow t$)

AT+ *Phys.Rev.D* 105 (2022) and a lot!



$$\langle OO(\tau) \rangle = \text{Tr}[O(0)O(\tau)\rho]$$

$$\rho = U(\tau)/Z$$



- Typical use of quantum algorithm is real-time simulation (\because Unitary).
 - Main interest: $\langle \Omega | O | \Omega \rangle$, where $|\Omega\rangle$ is the exact ground state
- Difficulty: State preparation* (as ground state extraction in conventional LQCD)
 - > Variational ansatz
- Our work uses the operator formalism. No problem for $\mu > 0$
- But how can we realized thermal “state”? (Next)

Density matrix, KL- U divergence

Density matrix

can describe statistical mechanics

Pure states: System is purely quantum

$$\rho_{\text{pure}} = |\Psi\rangle\langle\Psi|$$

$$\langle O \rangle = \text{Tr}[O\rho_{\text{pure}}] = \langle\Psi|O|\Psi\rangle$$

Mixed states: States are classically mixed (\neq superposition)

$$\rho_{\text{mixed}} = \sum_i w_i |\psi_i\rangle\langle\psi_i|$$

$$\langle O \rangle = \text{Tr}[O\rho_{\text{mixed}}] = \sum_i w_i \langle\psi_i|O|\psi_i\rangle$$

$w_i \in \mathbb{R}$ represents probability to find a pure state $|\psi_i\rangle$

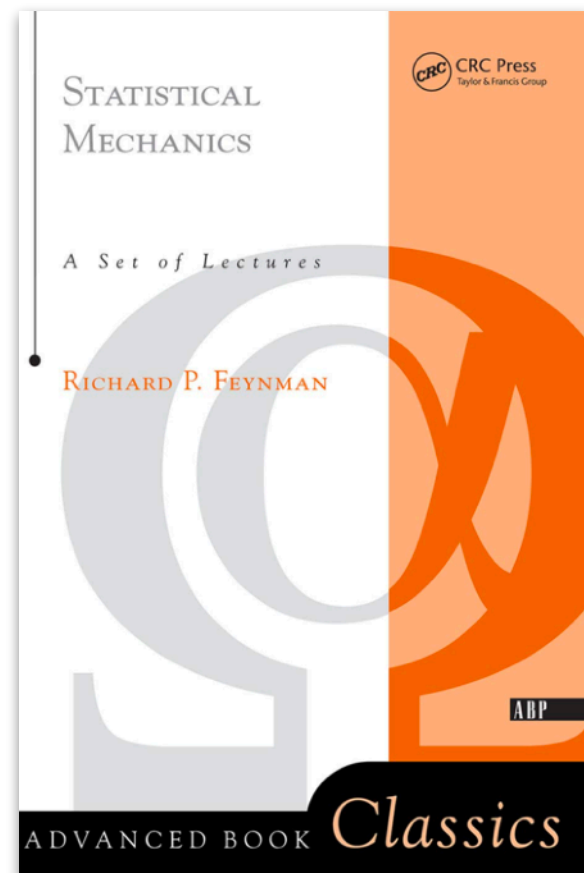
Thermal states (Grand-canonical):

$$\rho_{T,\mu} = \frac{1}{Z} e^{-\frac{1}{T}(\hat{H} - \mu\hat{N})}$$

$$\langle O \rangle_{T,\mu} = \text{Tr}[O\rho_{T,\mu}]$$

Thermal-quantum average in general

$$\langle O \rangle = \text{Tr}[O\rho]$$



Density matrix

Contain full information of the system, quantum version of probability distributions

*P. Caputa, AT + Physics Letters B 772, 53-57 (and a lot!)

Thermal-quantum average in general

$$\langle O \rangle = \text{Tr}[O\rho]$$

General Properties of density matrix ρ

- It unifies discretion of pure states and mixed states
- Normalized as $\text{Tr}[\rho] = 1$
- This can be used to investigate entanglement/thermalization*
- Density matrix ρ can be regarded as quantum version of probability distribution $p(x)$
 - e.g.) $S = - \int dx p(x) \log p(x)$ (Shannon entropy)
 - $\longleftrightarrow S = - \text{Tr}[\rho \log \rho]$ (Von-Neumann entropy)
- Distance between two density matrices = quantum relative entropy (next)

KL-*U* divergence

“Distance” between two density matrices

- KL divergence for ρ is Kullback–Leibler *Umegaki* divergence (Pseudo-distance for ρ)
- Classical version: $D(p \mid q) = \int dx \, p(x) \log p(x)/q(x)$ (KL divergence)
 - Relative entropy, quantifies difference of two distributions (~distance)
 - Positive definite, Used in machine learning
 - $D=0$ if and only if p, q are equal
- **Quantum version:** $D(\rho_1 \mid \rho_2) = \text{Tr}[\rho_1 \log \rho_1 / \rho_2]$ (KL-*Umegaki* divergence ~ distance)
 - Positive definite
 - $D=0$ if and only if ρ_1, ρ_2 are equal
- Kullback–Leibler *Umegaki* divergence can be used for variational approaches (as the flow models do)

Ansatz (model) for ρ ?

KL- U divergence

Beta VQE = VQE for $T>0$ = VQE + Neural net

J. -Guo Liu+ 1902.02663

*M. Germain+ 1502.03509

- Variational method for mixed states: Variational method on ρ

$$\bullet \rho_{\Theta}^{\text{ansatz}} = \sum_{\{\vec{x}\}} p_{\phi}[\vec{x}] U_{\theta} |\vec{x}\rangle \langle \vec{x}| U_{\theta}, \quad \Theta = \theta \cup \phi \text{ (parameters for state/distribution)}$$

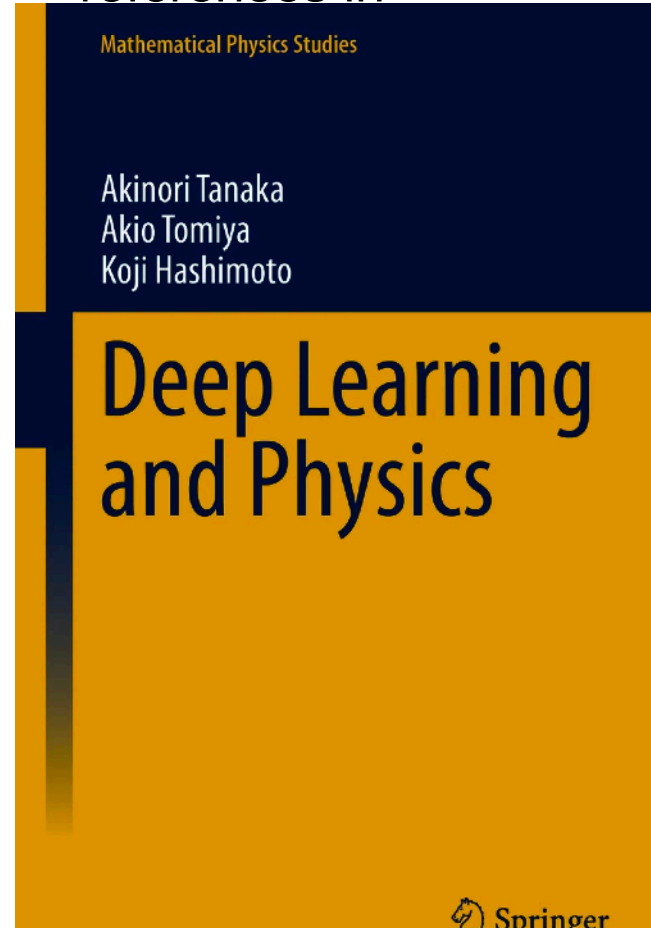
- $\vec{x} = (x_1, x_2, x_3, \dots, x_k, \dots)^T$, and $x_k \in \{0, 1\}$: Array of binary numbers

- $U_{\theta} |\vec{x}\rangle$: parametrized pure states, similar to the conventional VQE, θ = variational parameters in a circuit

- $p_{\phi}[\vec{x}]$: Classically approximated distribution for a configuration of \vec{x} ,
Neural network (Generative model, MADE*) is used.
 ϕ = parameters in the neural network

- **Neural network = Universal approximator of functions****
 Here, it approximate thermal distribution for fermions (generative model)

** references in



KL- U divergence

Optimized parameters enables us calculate $\langle O \rangle$ at T and μ

Jin-Guo Liu+ 1902.02663

- We **approximate** $\rho_{T,\mu}^{\text{exact}} = \frac{1}{Z} e^{-\frac{1}{T}(\hat{H} - \mu \hat{N})}$ by $\rho_{\Theta}^{\text{ansatz}} = \sum_{\{\vec{x}\}} p_{\phi}[\vec{x}] U_{\theta} |\vec{x}\rangle \langle \vec{x}| U_{\theta}$ by tuning/training parameters $\Theta = \theta \cup \phi$ with minimizing $D(\rho_{\Theta}^{\text{ansatz}} | \rho) \geq 0$
- $\langle O \rangle_{T,\mu} \approx \text{Tr}[\rho_{\Theta} O]$, if and only if $\rho_{\Theta} \approx \rho$
- Intuitively:
 - Quantum machine stores a state $U_{\theta} |\vec{x}\rangle$
 - Classical machine makes thermal distribution $p_{\phi}[\vec{x}]$ (neural net, generative model)

KL- U divergence

Shifted KLU can be used if $\ln Z$ is not available

J. -Guo Liu+ 1902.02663

*M. Germain+ 1502.03509

- Variational method for mixed states: Variational method on ρ

- $\rho_{\Theta}^{\text{ansatz}} = \sum_{\{\vec{x}\}} p_{\phi}[\vec{x}] U_{\theta} |\vec{x}\rangle \langle \vec{x}| U_{\theta}, \quad \Theta = \theta \cup \phi \text{ (parameters)}$

- $\vec{x} = (x_1, x_2, x_3, \dots, x_k, \dots)^T$, and $x_k \in \{0, 1\}$: (roughly) fermion excitation

- $U_{\theta} |\vec{x}\rangle$: parametrized pure states, similar to the conventional VQE

- $p_{\phi}[\vec{x}]$: Classically approximated distribution for a configuration of \vec{x} ,

Neural network (MADE*) is used. ϕ = parameters

- Minimizing $D(\rho_{\Theta}^{\text{ansatz}} | \rho_{T,\mu}^{\text{exact}})$, we get approximated a set of states (= thermal)

- Shifted D (by a constant, minimization is not affected) is used in real applications:

- $$\mathcal{L}(\Theta) \equiv D(\rho_{\Theta}^{\text{ansatz}} | \rho_{T,\mu}^{\text{exact}}) - \underbrace{\ln Z}_{\text{const}} = \text{Tr}[\rho_{\Theta}^{\text{ansatz}} \ln \rho_{\Theta}] + \frac{1}{T} \text{Tr}[\rho_{\Theta}^{\text{ansatz}} (\hat{H} - \mu \hat{N})]$$

BUT, in this work, we use exact diagonalization to check the validity of the algorithm and see D

Results

Results

Schwinger model at finite μ and T . Code is written in Julia


- We apply beta-VQE for Schwinger model = QED in 1+1d.

AT. 2205.08860

- Toy model for QCD in 4d (good for testbed). Common features: Confinement, chiral anomaly, topology ...

$$S = \int d^2x \left[-\frac{1}{4} F_{\mu\nu} F^{\mu\nu} + \bar{\psi} (i \not{\partial} - g A - m) \psi \right] \longleftrightarrow H = \int dx \left[-i \bar{\psi} \gamma^1 (\partial_1 + i g A_1) \psi + m \bar{\psi} \psi + \frac{1}{2} \Pi^2 \right]$$

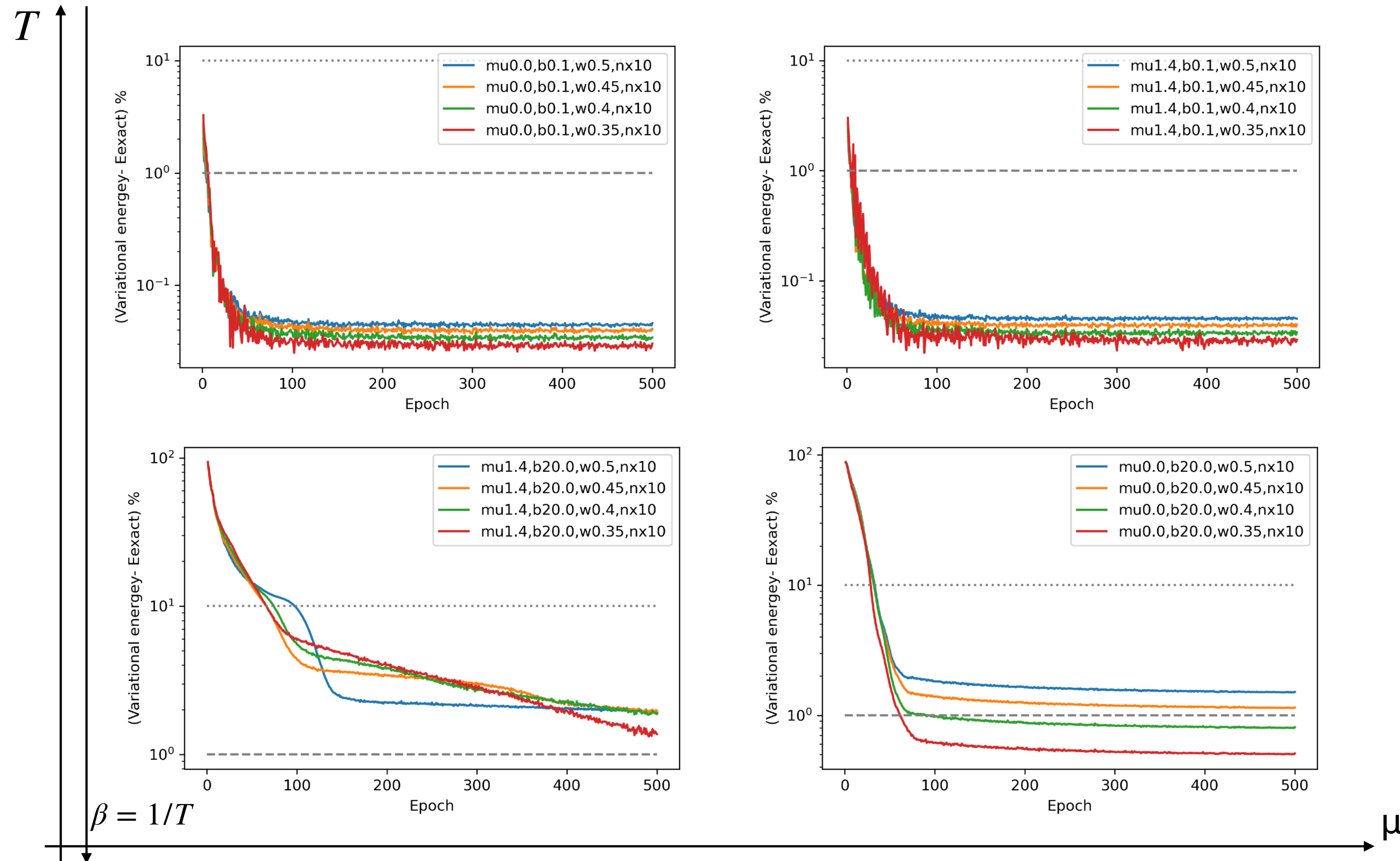
$$\partial_x E = g \bar{\psi} \gamma^0 \psi$$

- Staggered discretization, Jordan-Wigner transformation, and open BC are used
- $g = 1$, $N_x = (4, 6), 8, 10$, $1/T = [0.5-20.0]$, $\mu = [0-1.4]$, 4 lattice spacings $1/2a = [0.5-0.35]$
- We do not take large volume limit but take continuum limit
 - (Practically, $N_x > 10$ cannot be calculated on our numerical resources)
 - (My previous work shows data from $N_x > 12$ are essential to take stable large volume limit though)
- Setup for beta VQE:
 - Unitary part = SU(4) ansatz
 - Classical weight = Masked Auto-Encoder for Distribution Estimation (MADE)
- Training epoch is 500. Sampling = 5000 for classical distribution
- The code is implemented in  . Calculations on Yukawa21 cluster at Kyoto University
- **Observables**
 - Variational free energy (exact = lower-bound, and variational one)
 - (Translationally invariant) Chiral condensate
- **Check point: Dependence of variational error on temperature**

Results

Compare the variational energy ~ D. Looks good

AT. 2205.08860



Vertical axis:
Deviation of D
from the exact one(%)

Horizontal axis:
Training epoch

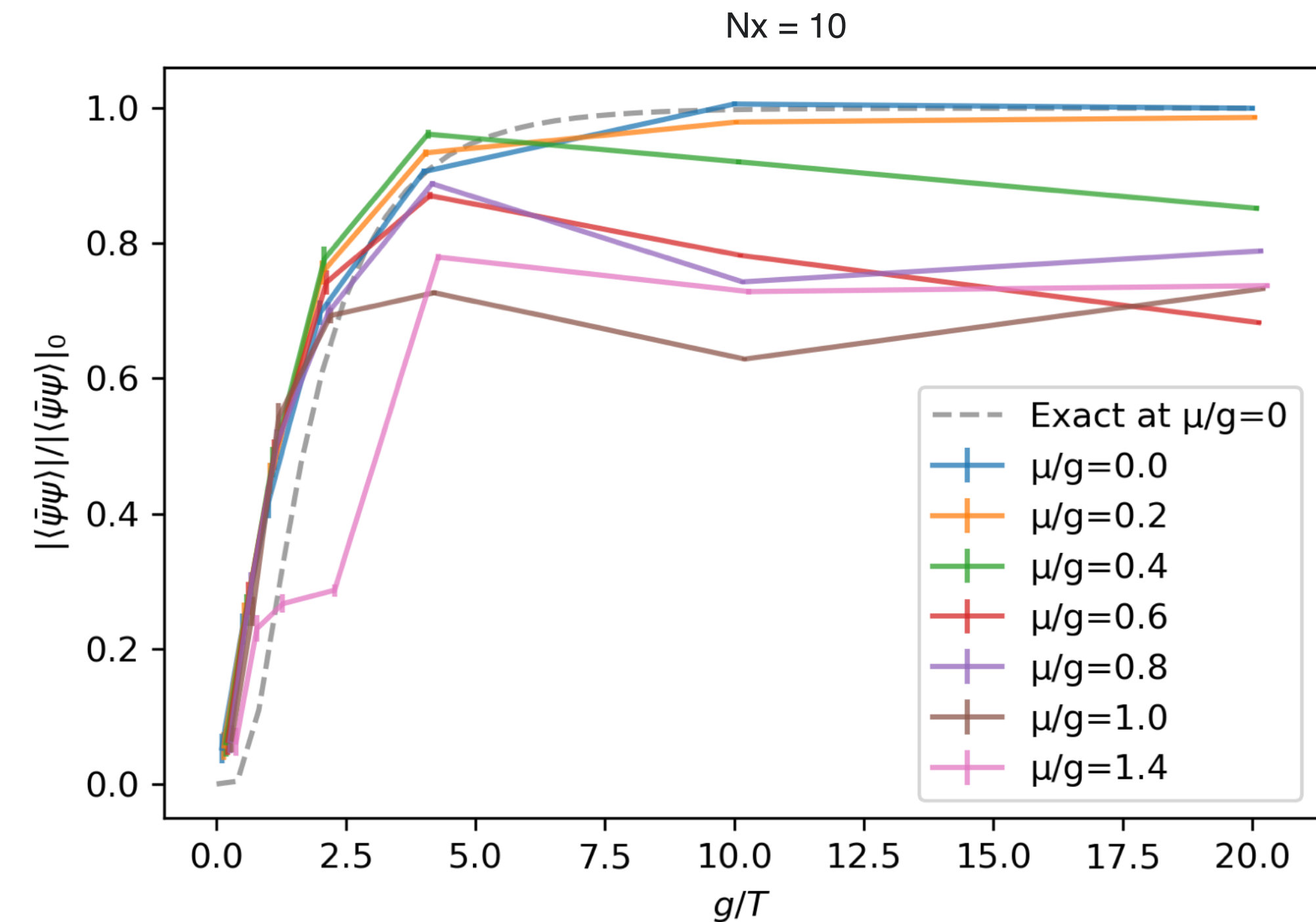
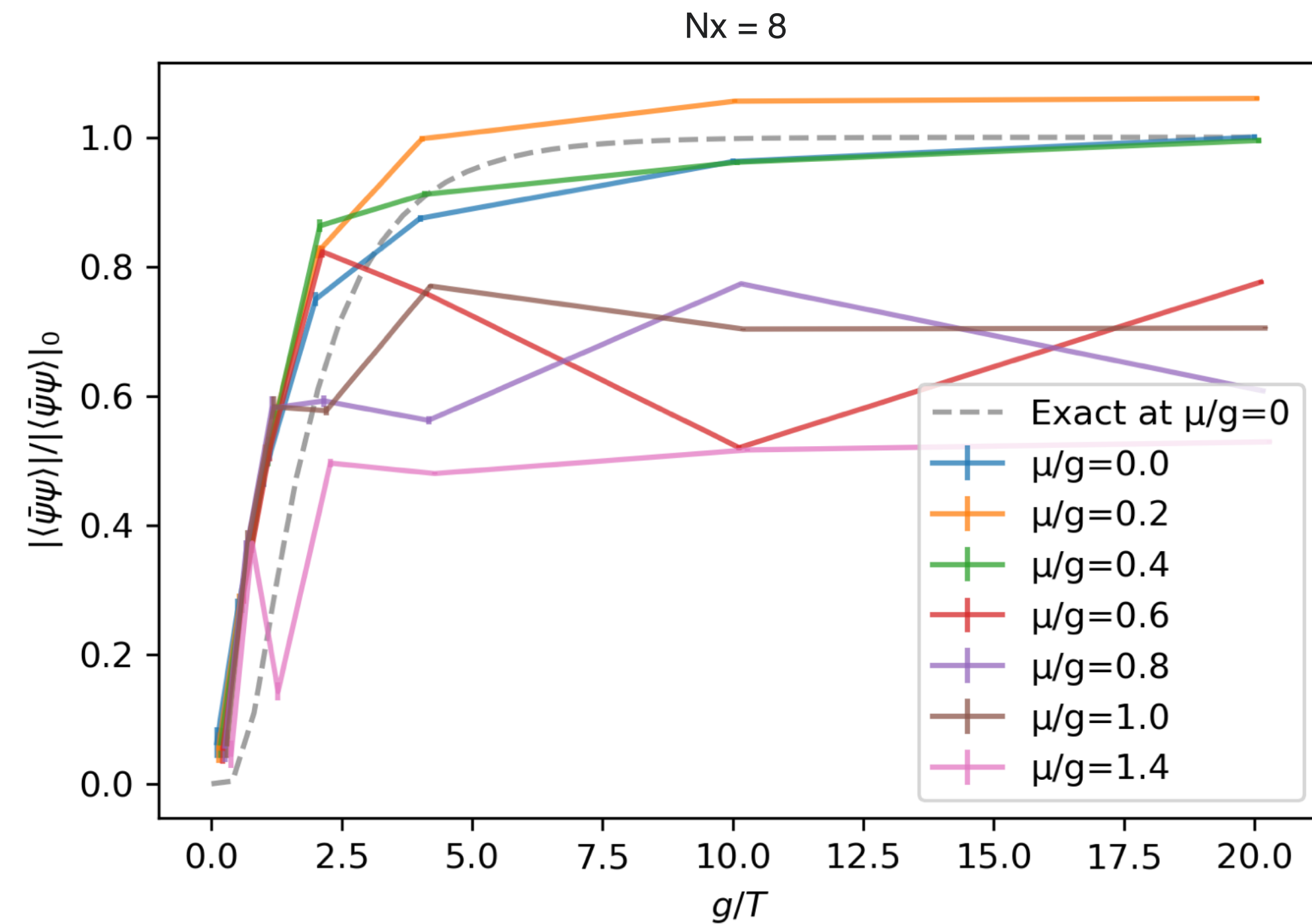
1. Mild dependence on μ
2. Hard for $T \rightarrow 0$ (large deviation) as expected

Results

Chiral condensate as a function of T and μ

AT arXiv: 2205.08860

Qualitatively consistent at $\mu = 0$ (*), and results seems reasonable



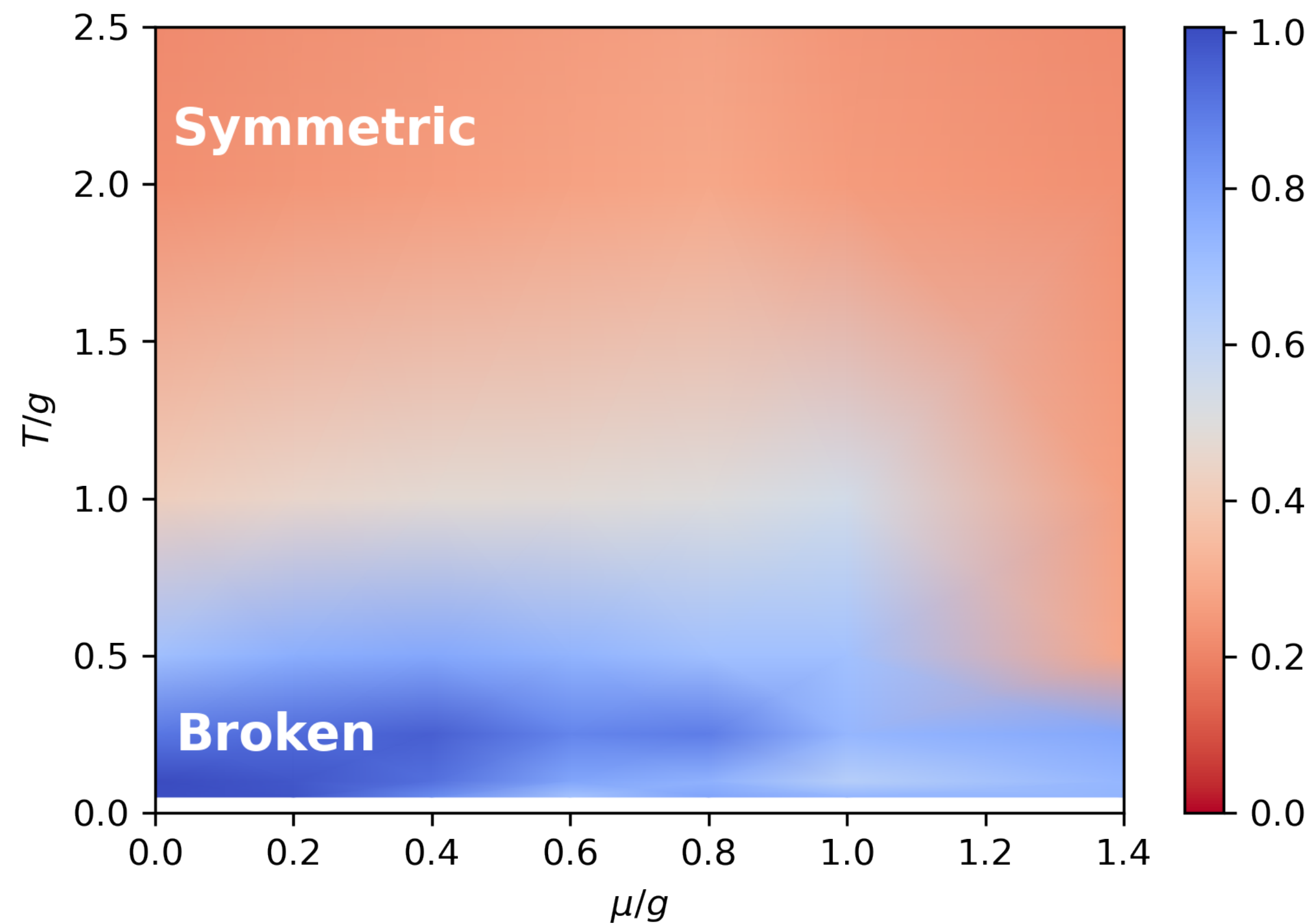
*(I did not include additive mass shift (Ross Dempsey+ arXiv: 2206.05308).

I thank to Takis Angelides (DESY) and Etsuko Itou (RIKEN) for letting me know this important reference!)

We use $N_x = 10$ results for the phase diagram

Summary

Classical-quantum hybrid algorithm: $T > 0$ and $\mu > 0$



AT. 2205.08860



- We investigate T - μ phase diagram for Schwinger model. **This algorithm works other than this model.**
- Continuum extrapolation has been evaluated (w/o mass shift by Ross Dempsey+ arXiv: 2206.05308)
- The variational approach does not show difficulty for our parameter regime
- Future works: Towards to go large volume, optimization of code, GPU version, tensor network. (noise-free) real device!. Finite temperature + finite density + real time??
Pure state ansatz? (relation to TPQ/QITE?), scaling? $SU(N)$?

Thanks!



- Quantum machine: Exact ground state preparation is hard. In particular, it is difficult on near term devices
- **Variational method for a pure state** with a short circuit (VQE, variation quantum eigen-solver).
 - Quantum/Classical hybrid algorithm, iterative
 - **Parametrized unitary circuit (~parametrized state $|\theta\rangle$, θ : a set of parameters)**
 - $|\theta\rangle = \hat{U}(\theta)(|0\rangle_1|0\rangle_2|0\rangle_3\cdots)$, and $\hat{U}(\theta)$ is a short circuit (entanglement + rotations)
 - If $\langle\theta|H|\theta\rangle = 0$, $|\theta\rangle \approx |\Omega\rangle$, where $|\Omega\rangle$ is the exact ground state
- Systematic error since $|\theta\rangle \neq |\Omega\rangle$ (cannot be exact in practice)
 - How about thermal states? Thermal evolution is not unitary
 - (TPQ/QITE on a classical emulator is an option but short circuit rep is not known)

Beta VQE

O(0.1)% for T>>0. T ~0 is difficult

μ/g	g/T	N_x	$\sim 1/a$ w/g	Approx $\mathcal{L} - \ln Z$	Exact $-\ln Z$	Diff (%)				$\sim 1/a$	Approx	Exact	
0.0	0.1	4	0.5	-27.779	-27.781	0.00804	1.4	0.1	4	0.5	-28.021	-28.023	0.00697
0.0	0.1	4	0.35	-27.807	-27.808	0.005	1.4	0.1	4	0.35	-27.989	-27.991	0.00755
0.0	0.1	10	0.5	-70.686	-70.718	0.0459	1.4	0.1	10	0.5	-70.842	-70.874	0.0453
0.0	0.1	10	0.35	-71.744	-71.765	0.0302	1.4	0.1	10	0.35	-71.742	-71.763	0.0291
0.0	0.5	4	0.5	-5.792	-5.802	0.185	1.4	0.5	4	0.5	-6.784	-6.789	0.0609
0.0	0.5	4	0.35	-5.885	-5.891	0.105	1.4	0.5	4	0.35	-6.644	-6.647	0.0327
0.0	0.5	10	0.5	-17.133	-17.25	0.68	1.4	0.5	10	0.5	-17.989	-18.104	0.636
0.0	0.5	10	0.35	-18.849	-18.934	0.448	1.4	0.5	10	0.35	-19.445	-19.534	0.456
0.0	10.0	4	0.5	-1.748	-1.75	0.161	1.4	10.0	4	0.5	-3.708	-3.71	0.0728
0.0	10.0	4	0.35	-1.829	-1.829	0.0184	1.4	10.0	4	0.35	-3.63	-3.669	1.07
0.0	10.0	10	0.5	-8.218	-8.341	1.48	1.4	10.0	10	0.5	-10.067	-10.243	1.71
0.0	10.0	10	0.35	-9.98	-10.03	0.496	1.4	10.0	10	0.35	-11.763	-11.862	0.837
0.0	20.0	4	0.5	-1.492	-1.739	14.2	1.4	20.0	4	0.5	-3.673	-3.681	0.218
0.0	20.0	4	0.35	-1.653	-1.806	8.46	1.4	20.0	4	0.35	-3.621	-3.669	1.31
0.0	20.0	10	0.5	-8.202	-8.328	1.51	1.4	20.0	10	0.5	-10.028	-10.224	1.92
0.0	20.0	10	0.35	-9.955	-10.006	0.509	1.4	20.0	10	0.35	-11.699	-11.862	1.37

- 1.Mild dependence on μ
- 2.Hard for T -> 0 (large deviation) as expected

state preparation

state

Jin-Guo Liu+ 1902.02663

- We approximate $\rho_{T,\mu}^{\text{exact}} = \frac{1}{Z} e^{-\frac{1}{T}(\hat{H} - \mu \hat{N})}$ by $\rho_{\Theta}^{\text{ansatz}} = \sum_{\{\vec{x}\}} p_{\phi}[\vec{x}] U_{\theta} |\vec{x}\rangle \langle \vec{x}| U_{\theta}$
by tuning/training parameters with minimizing $D(\rho_{\Theta}^{\text{ansatz}} | \rho)$
 - $\langle O \rangle_{T,\mu} \approx \text{Tr}[\rho_{\Theta} O]$, if and only if $\rho_{\Theta} \approx \rho$
 - Quantum machine can store a state $U_{\theta} |\vec{x}\rangle$
 - Classical machine can sample thermal distribution from $p_{\phi}[\vec{x}]$ (neural net)

state preparation

state

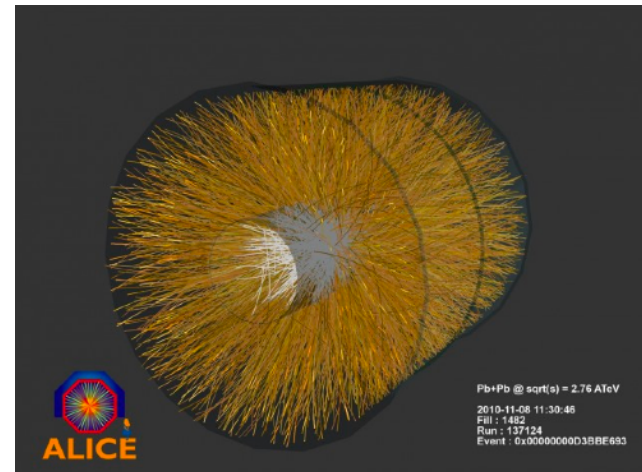
QCD (Quantum Chromo-dynamics) in 3 + 1 dimension

$$S = \int d^4x \left[-\frac{1}{2} \text{tr} F_{\mu\nu} F^{\mu\nu} + \bar{\psi} (\mathbf{i}\not{\partial} + g\mathbf{A} - m) \psi \right]$$

$$F_{\mu\nu} = \partial_\mu A_\nu - \partial_\nu A_\mu - ig[A_\mu, A_\nu]$$

$A_\mu(x) \in su(3)$, 3x3 traceless, hermitian

$$|\psi(t)\rangle = e^{-iHt} |\psi(0)\rangle \quad H: \text{Hamiltonian from } S$$



- Generalization of QED, $A_\mu(x)$ is a matrix (Yang-Mills-Uchiyama)
- Action above enables us to calculate followings:
 - Tc of Quark-Hadron, Matrix elements of QCD
 - Forces between nuclei ... etc!

Lattice QCD in 4 dimension

F. Wegner 1971
K. Wilson 1974

$$U_\mu = e^{igA_\mu}$$

Lattice regulation

$$S[U, \psi, \bar{\psi}] = a^4 \sum_n \left[-\frac{1}{g^2} \text{Re tr} U_{\mu\nu} + \bar{\psi} (\not{D} + m) \psi \right]$$

$$\langle \mathcal{O} \rangle = \frac{1}{Z} \int \mathcal{D}U \mathcal{D}\bar{\psi} \mathcal{D}\psi e^{-S} \mathcal{O}(U)$$

$$|\psi(t)\rangle = e^{-H\tau} |\psi(0)\rangle$$

- Lattice QCD has same long-distance physics with continuum QCD
- Euclidean signature, statistic physics

My related talks

U(1)A at fin. temp by Yu Zhang, 28 Jul 2021, 05:45(EDT)

QCD + magnetic field by Xiaodang Wang, 28 Jul 2021, 22:00 (EDT)

Schwinger model

=2D QED: Solvable at $m=0$, similar to QCD in 4D.

Schwinger model = QED in 1+1 dimension

$$S = \int d^2x \left[-\frac{1}{4} F_{\mu\nu} F^{\mu\nu} + \bar{\psi} (i\partial - gA - m) \psi + \frac{g\theta}{4\pi} \epsilon_{\mu\nu} F^{\mu\nu} \right]$$

Similarities to QCD in 3+1

- Confinement
- Chiral symmetry breaking (different mechanism), gapped even $m=0$

$$\langle \bar{\psi} \psi \rangle = -\frac{e' g}{\pi^{3/2}} = -g 0.16 \dots$$

- Topological term can be included as in QCD
- Vacuum decay by external electric field (Schwinger effect)

Hamiltonian of Schwinger model

Akio Tomiya

=2D QED: Solvable at $m=0$, similar to QCD in 4D.

Schwinger model = QED in 1+1 dimension

$$S = \int d^2x \left[-\frac{1}{4} F_{\mu\nu} F^{\mu\nu} + \bar{\psi} (i\partial\!\!\!/ - gA - m) \psi \right]$$

- Strategy
 1. Derive Hamiltonian with gauge fixing
 2. Rewrite gauge field to fermions using Gauss' law
 3. Use Jordan-Wigner transformation \rightarrow Spin system

Why? next page

Hamiltonian of Schwinger model

Schwinger model in spin language

Schwinger model = QED in 1+1 dimension

$$S = \int d^2x \left[-\frac{1}{4} F_{\mu\nu} F^{\mu\nu} + \bar{\psi} (i\not{\partial} - g\not{A} - m) \psi \right]$$



- Strategy (1 gauge fix, 2 Gauss' law, 3 Jordan-Wigner trf)

Schwinger model on the lattice (staggered fermion, OBC, Spin rep.)

$$H = \frac{1}{4a} \sum_n \left[X_n X_{n+1} + Y_n Y_{n+1} \right] + \frac{m}{2} \sum_n (-1)^n Z_n + \frac{g^2 a}{2} \sum_n \left[\sum_{j=1}^n \left(\frac{Z_j + (-1)^j}{2} \right) + \epsilon_0 \right]^2$$

X_j : Pauli matrix of x on site j

Y_j : Pauli matrix of y on site j

Z_j : Pauli matrix of z on site j

- Spin representation is necessary to use quantum device
(Analogous to floating point rep. in classical machine)
- (QCD + QC also requires this strategy)

Hamiltonian of Schwinger model

Akio Tomiya

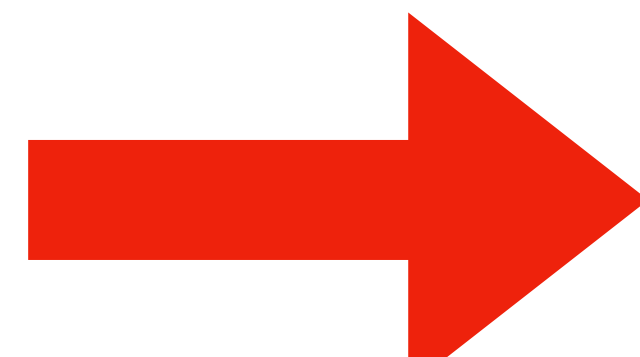
=2D QED: Solvable at $m=0$, similar to QCD in 4D.

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Schwinger model = QED in 1+1 dimension

$$S = \int d^2x \left[-\frac{1}{4} F_{\mu\nu} F^{\mu\nu} + \bar{\psi} (i\not{\partial} - g\not{A} - m) \psi \right]$$

$$\Pi(x) = \frac{\partial \mathcal{L}}{\partial \dot{A}^1(x)} = \dot{A}(x) = E(x)$$


$$A_0 = 0 \quad \left\{ \begin{array}{l} H = \int dx \left[-i\bar{\psi}\gamma^1(\partial_1 + igA_1)\psi + m\bar{\psi}\psi + \frac{1}{2}\Pi^2 \right] \\ \partial_x E = g\bar{\psi}\gamma^0\psi \end{array} \right.$$

(Gauss' law constraint)

This constrains time evolution to be gauge invariant

Lattice Hamiltonian formalism

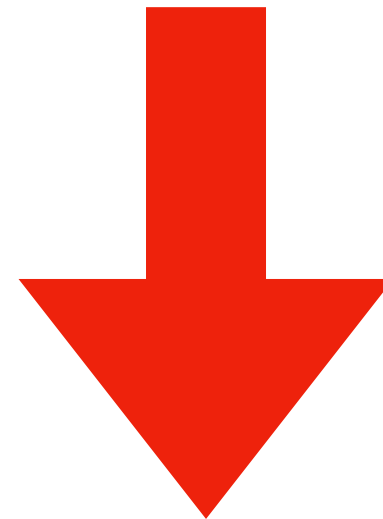
Hamiltonian on a discrete space

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Schwinger model in continuum

$$H = \int dx \left[-i\bar{\psi}\gamma^1(\partial_1 + igA_1)\psi + m\bar{\psi}\psi + \frac{1}{2}\Pi^2 \right]$$

Gauss' law $\partial_x E = g\bar{\psi}\gamma^0\psi$



$$-\frac{1}{g}\Pi(x) \rightarrow L_n$$

upper component of $\psi \rightarrow \chi_{\text{even-site}}$

$$-agA_1(x) \rightarrow \phi_n$$

lower component of $\psi \rightarrow \chi_{\text{odd-site}}$

Schwinger model on the lattice (staggered fermion)

$$H = -\frac{i}{2a} \sum_{n=1}^{N-1} \left[\chi_{n+1}^\dagger e^{-i\phi_n} \chi_n - \chi_n^\dagger e^{i\phi_n} \chi_{n+1} \right] + m \sum_{n=1}^N (-1)^n \chi_n^\dagger \chi_n + \frac{g^2 a}{2} \sum_{n=1}^{N-1} L_n^2$$

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

Lattice Schwinger model = spin system

Akio Tomiya

Gauge trf, open bc, Gauss law \rightarrow pure fermionic system

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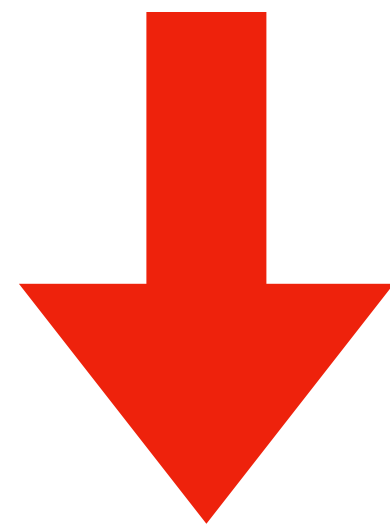
Schwinger model on the lattice (staggered fermion)

$$H = -\frac{i}{2a} \sum_{n=1}^{N-1} \left[\chi_{n+1}^\dagger e^{-i\phi_n} \chi_n - \chi_n^\dagger e^{i\phi_n} \chi_{n+1} \right] + m \sum_{n=1}^N (-1)^n \chi_n^\dagger \chi_n + \frac{g^2 a}{2} \sum_{n=1}^{N-1} L_n^2$$

Gauss' law

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

$L_0 = \epsilon_0 \in \mathbb{R}$ (open B.C.), and insert "Gauss' law"



$$\begin{cases} U_n = \prod_{j=1}^{n-1} e^{-i\phi_j} \\ \chi_n \rightarrow U_n \chi_n \\ e^{-i\phi_{n-1}} \rightarrow U_{n-1} e^{-i\phi_{n-1}} U_n^\dagger \end{cases}$$

remnant gauge transformation

Schwinger model on the lattice (staggered fermion, OBC)

$$H = -\frac{i}{2a} \sum_n \left[\chi_{n+1}^\dagger \chi_n - \chi_n^\dagger \chi_{n+1} \right] + m \sum_n (-1)^n \chi_n^\dagger \chi_n + \frac{g^2 a}{2} \sum_n \left[\sum_j^n \left(\chi_j^\dagger \chi_j - \frac{1 - (-1)^j}{2} \right) + \epsilon_0 \right]^2$$

Lattice Schwinger model

We requires anticommutations to fermions

(skip)

Schwinger model on the lattice (staggered fermion, OBC)

$$H = -\frac{i}{2a} \sum_n \left[\chi_{n+1}^\dagger \chi_n - \chi_n^\dagger \chi_{n+1} \right] + m \sum_n (-1)^n \chi_n^\dagger \chi_n + \frac{g^2 a}{2} \sum_n \left[\sum_j^n \left(\chi_j^\dagger \chi_j - \frac{1 - (-1)^j}{2} \right) + \epsilon_0 \right]^2$$

System is quantized by assuming the canonical anti-commutation relation

$$\{\chi_j^\dagger, \chi_k\} = i\delta_{jk} \quad j, k = \text{site index}$$

On the other hand, Pauli matrices satisfy anti-commutation as well

$$\{\sigma^\mu, \sigma^\nu\} = 2\delta_{\mu\nu} \mathbf{1} \quad \mu, \nu = 1, 2, 3$$

Quantum spin-chain case, each site has Pauli matrix, but they are “commute”.

We can absorb difference of statistical property using Jordan Wigner transformation

Jordan-Wigner transformation:

$$\chi_n = \frac{X_n - iY_n}{2} \prod_{j < n} (iZ_j)$$

X_j : Pauli matrix of x on site j

Y_j : Pauli matrix of y on site j

Z_j : Pauli matrix of z on site j

← This guarantees the statistical property

This (re)produces correct Fock space.

We can rewrite the Hamiltonian in terms of spin-chain

Lattice Schwinger model = spin system Akio Tomiya

Jordan-Wigner transformation: Fermions ~ Spins

(skip)

Schwinger model on the lattice (staggered fermion, OBC)

$$H = -\frac{i}{2a} \sum_n \left[\chi_{n+1}^\dagger \chi_n - \chi_n^\dagger \chi_{n+1} \right] + m \sum_n (-1)^n \chi_n^\dagger \chi_n + \frac{g^2 a}{2} \sum_n \left[\sum_j^n \left(\chi_j^\dagger \chi_j - \frac{1 - (-1)^j}{2} \right) + \epsilon_0 \right]^2$$



$$\begin{cases} \chi_n = \frac{X_n - iY_n}{2} \prod_{j<n} (iZ_j) \\ \chi_n^\dagger = \frac{X_n + iY_n}{2} \prod_{j<n} (-iZ_j) \end{cases}$$

Jordan-Wigner transformation

X_j : Pauli matrix of x on site j

Y_j : Pauli matrix of y on site j

Z_j : Pauli matrix of z on site j

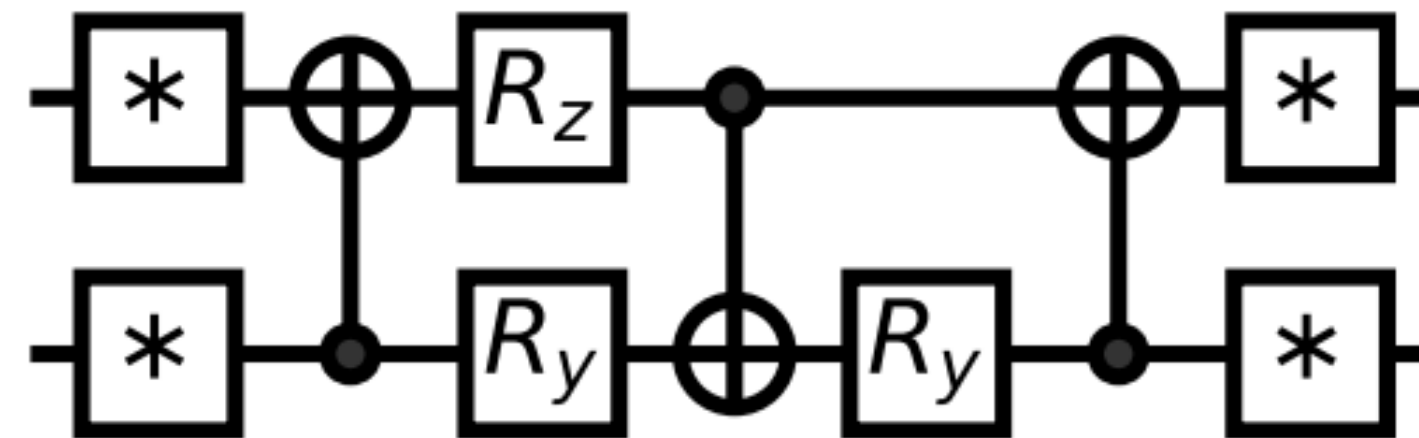
Schwinger model on the lattice (staggered fermion, OBC, Spin rep.)

$$H = \frac{1}{4a} \sum_n \left[X_n X_{n+1} + Y_n Y_{n+1} \right] + \frac{m}{2} \sum_n (-1)^n Z_n + \frac{g^2 a}{2} \sum_n \left[\sum_{j=1}^n \left(\frac{Z_j + (-1)^j}{2} \right) + \epsilon_0 \right]^2$$

[Y. Hosotani 9707129]

SU(4)

Variational ansatz



$$\boxed{*} = \boxed{R_z} \boxed{R_y} \boxed{R_z}$$

The general gate consists of 15 single qubit gates and 3 CNOT gates.
Each two qubit unitary is parametrized by 15 parameters in the rotational gates, which parametrizes the SU(4) group.