



Theory and Phenomenology
of Fundamental Interactions

UNIVERSITY AND INFN · BOLOGNA

Workshop
Quantum Computing @ INFN
November 14, 2022

Bayesian Optimization for QAOA

Elisa Ercolessi

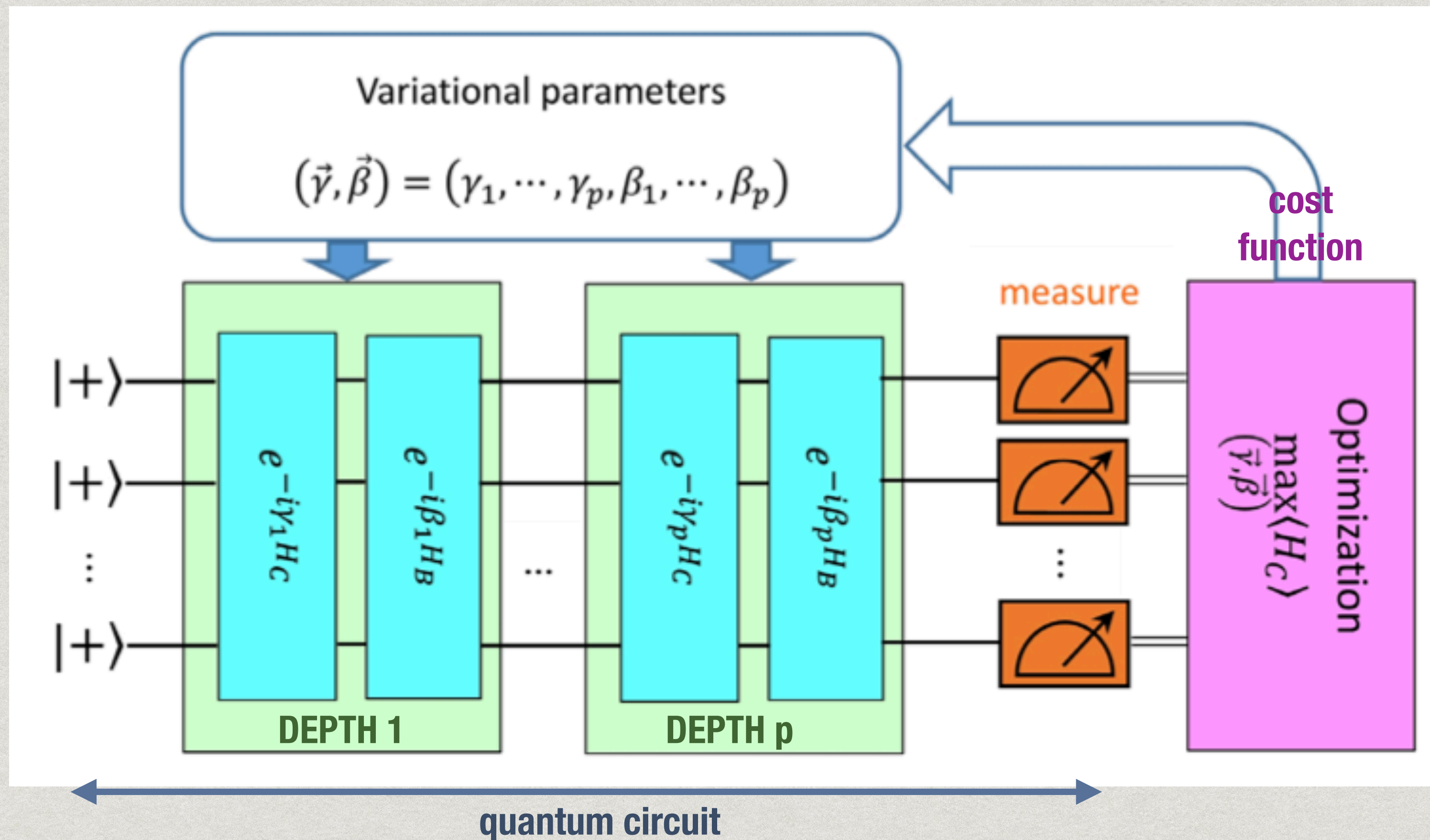
*Dept. Physics & Astronomy - University of Bologna
& INFN - Sezione di Bologna*

with: D. Vodola, C. Sanavio, F. Dell'Anna, S. Tibaldi, C. Degli Esposti Boschi, E. Tignone



• Hybrid classical-quantum protocol: QAOA

(Quantum Approximate Optimisation Algorithm) Farhi et al.- arXiv:1411.4028 (2014)



quantum

- Objective function to be minimised
- Quantum evolution (circuit) with variational parameters
- Measurement scheme for average Hamiltonian to get cost function

classical

- Optimisation algorithm to get the best values of the parameters

• Ground State Preparation for Nontrivial Quantum Hamiltonians

Z2 Lattice (pure) Gauge Theory

$$H(h) = H_E + h H_B$$

Variational Ground State

$$|\psi_P(\gamma, \beta)\rangle = \left(\prod_{m=1}^P e^{-i\beta_m H_E} e^{-i\gamma_m H_B} \right) |\psi_0\rangle$$

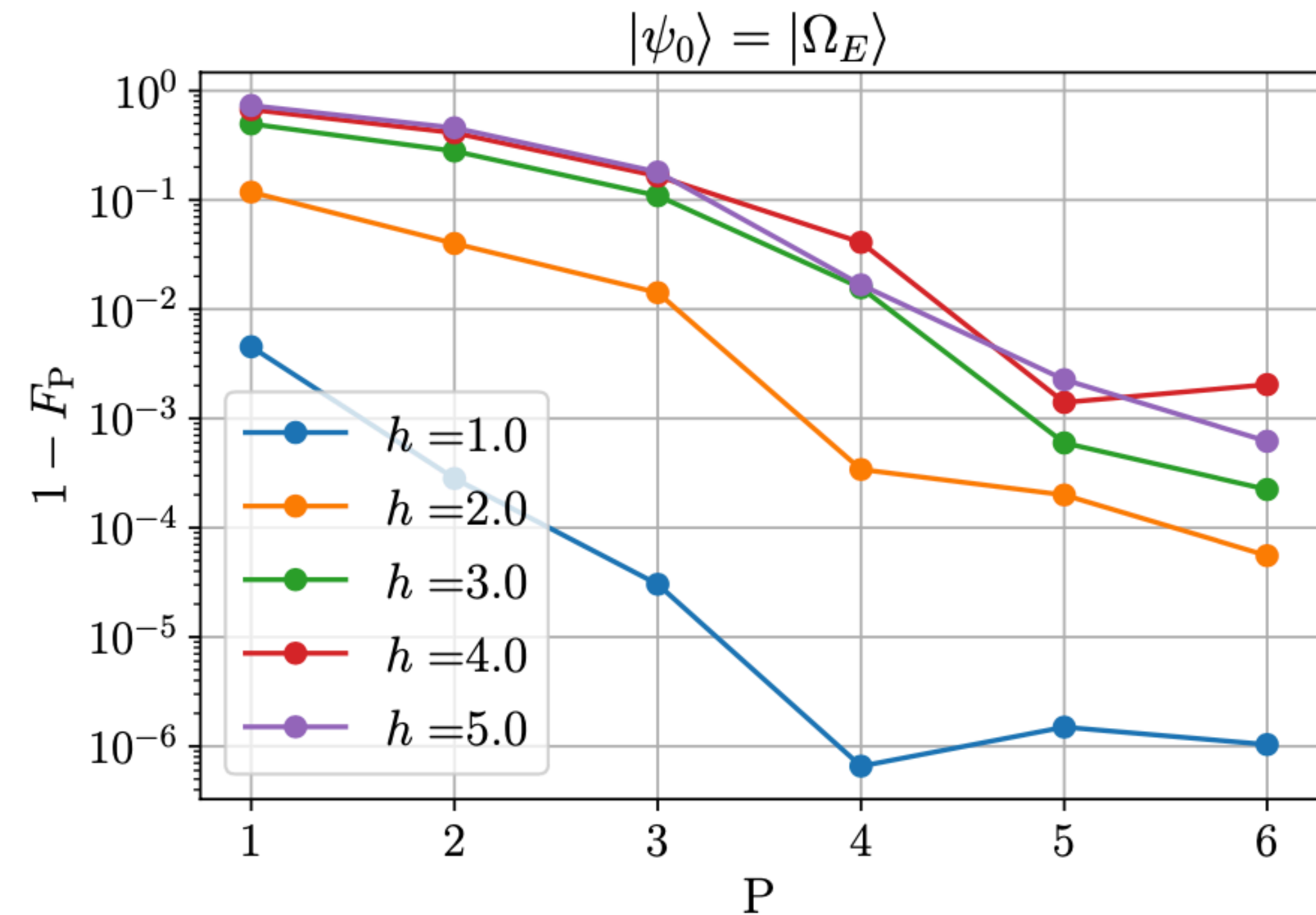
via quantum circuit

Cost function to be minimized
by classical optimiser:

$$E_P(\gamma, \beta) = \langle \psi_P(\gamma, \beta) | H(h) | \psi_P(\gamma, \beta) \rangle$$

(within: QuantERA 2020 Project “QuantHEP”, with BA and PD
& Project in INFN-Quantum Computing HUB CERN-IBM)

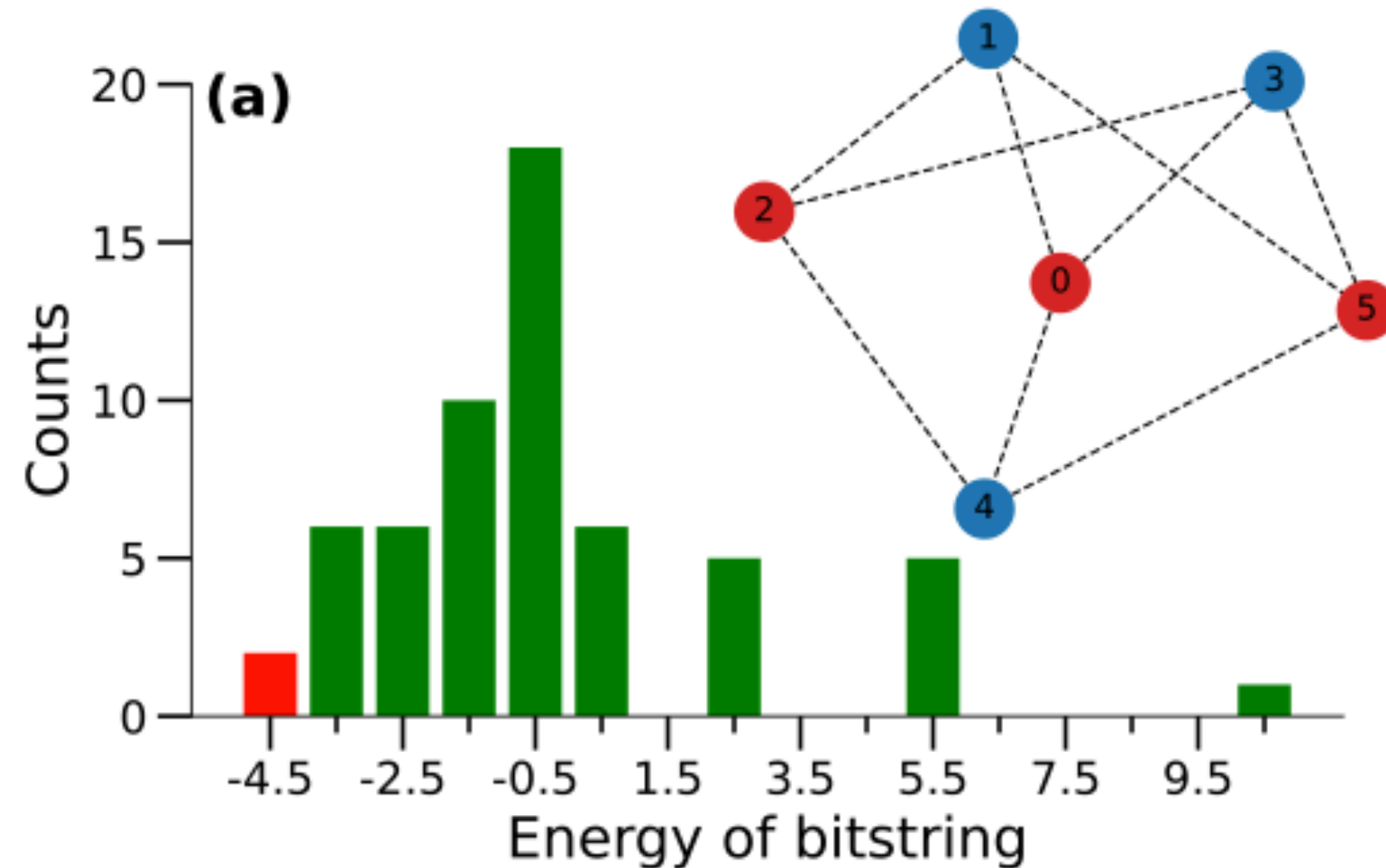
PRX Quantum 3, 020320 (2022)



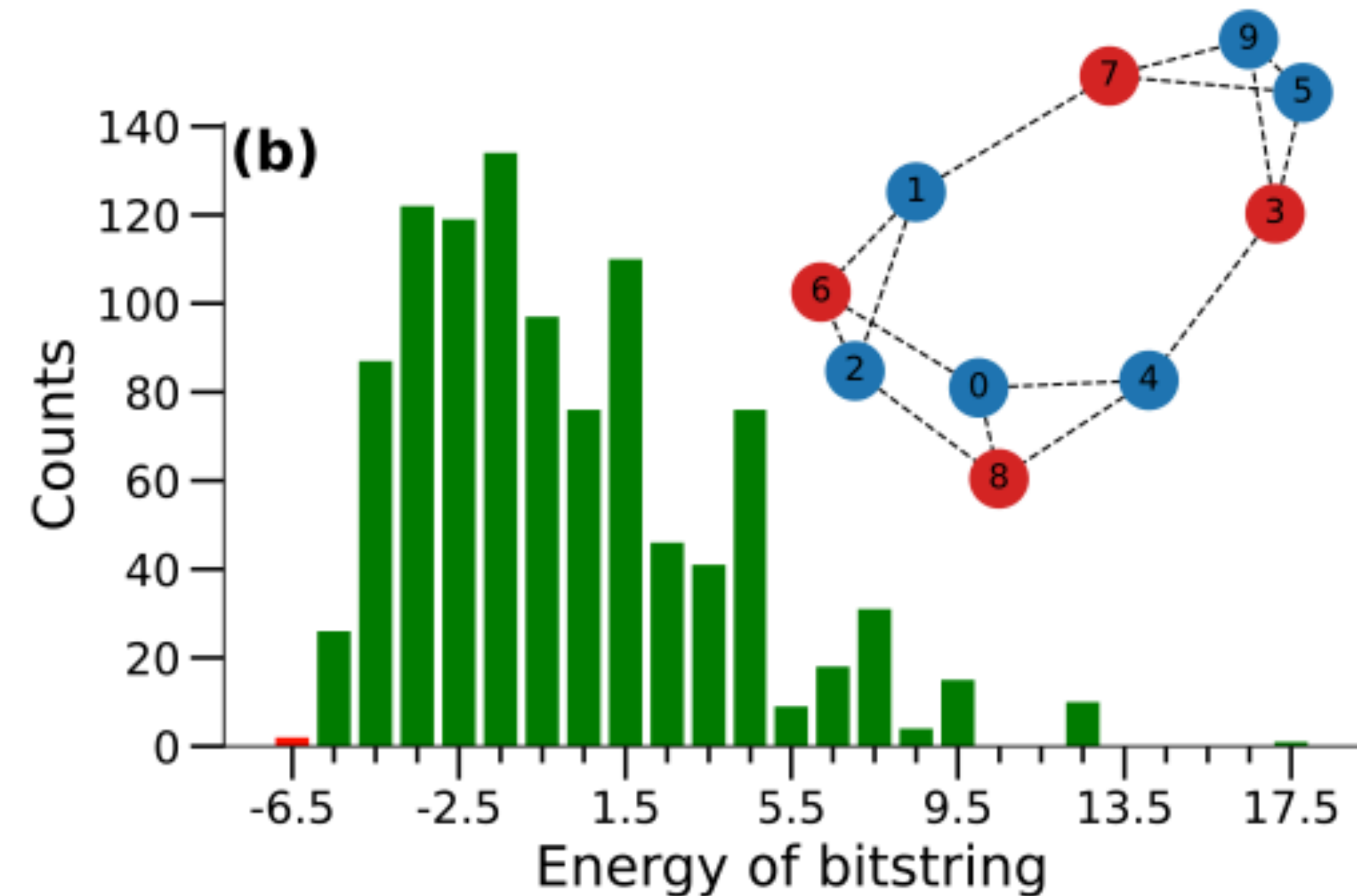
• Solution of Combinatorial problems on Graphs

MIS problem: find the largest set of nodes not adjacent

Max-Cut problem: partition the graph in two sets of nodes interconnected by the largest number of links



GS of $H_C^{\text{MIS}} = \sum_i Z_i + \omega \sum_{(i,j) \in E} Z_i Z_j,$



GS of $H_C^{\text{MC}} = - \sum_{(i,j) \in E} (1 - Z_i Z_j)/2.$

$$H_M = \sum_i X_i$$

where Z_j, X_j are the Pauli matrices acting on the qubit representing node j

The Quantum circuit now prepares the state

$$|\boldsymbol{\theta}\rangle = \prod_{l=1}^p e^{-i\beta_l H_M} e^{-i\gamma_l H_C} |+\rangle,$$

with $\theta = (\gamma_1\beta_1, \dots, \gamma_p\beta_p)$

The energy cost has a minimum for γ^*, β^*
to be found by means of a classical protocol

$$E(\boldsymbol{\theta}) = \langle \boldsymbol{\theta} | H_C | \boldsymbol{\theta} \rangle$$

Issue:

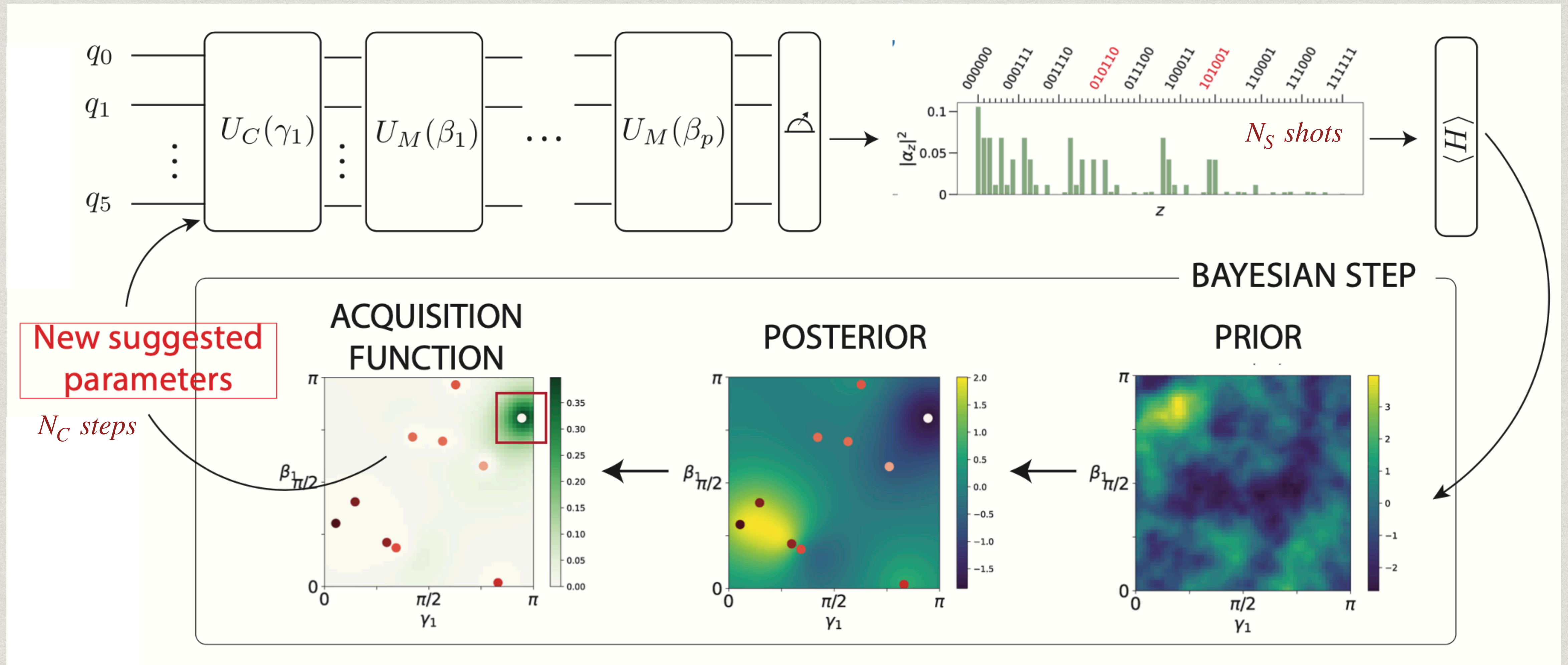
- the landscape of the energy can become very rough and develop barren plateaus
- convergence can become hard (global techniques are needed) and a large number of points are to be tested

but, on a real device, we have to consider also problems related to

- the measure of the cost function, which requires many measurements (N_S)
to give a good estimate of the average, a procedure that has to be repeated for all points (N_C) necessary to reach the minimum with the desired precision

• Bayesian algorithm for classical optimization

warm up: we sample $N_W = 10$ values of the parameters θ_j and measure the cost function $\langle \theta_j | H_C | \theta_j \rangle$



subroutine: a prior distribution is updated by means of the observed data & the posterior distribution is acquired through an acquisition function whose maximum indicates the new value of the parameter where to perform the run (Gaussian process)

● Results of the simulations

for the Max-Cut problem
of the 10-node graph

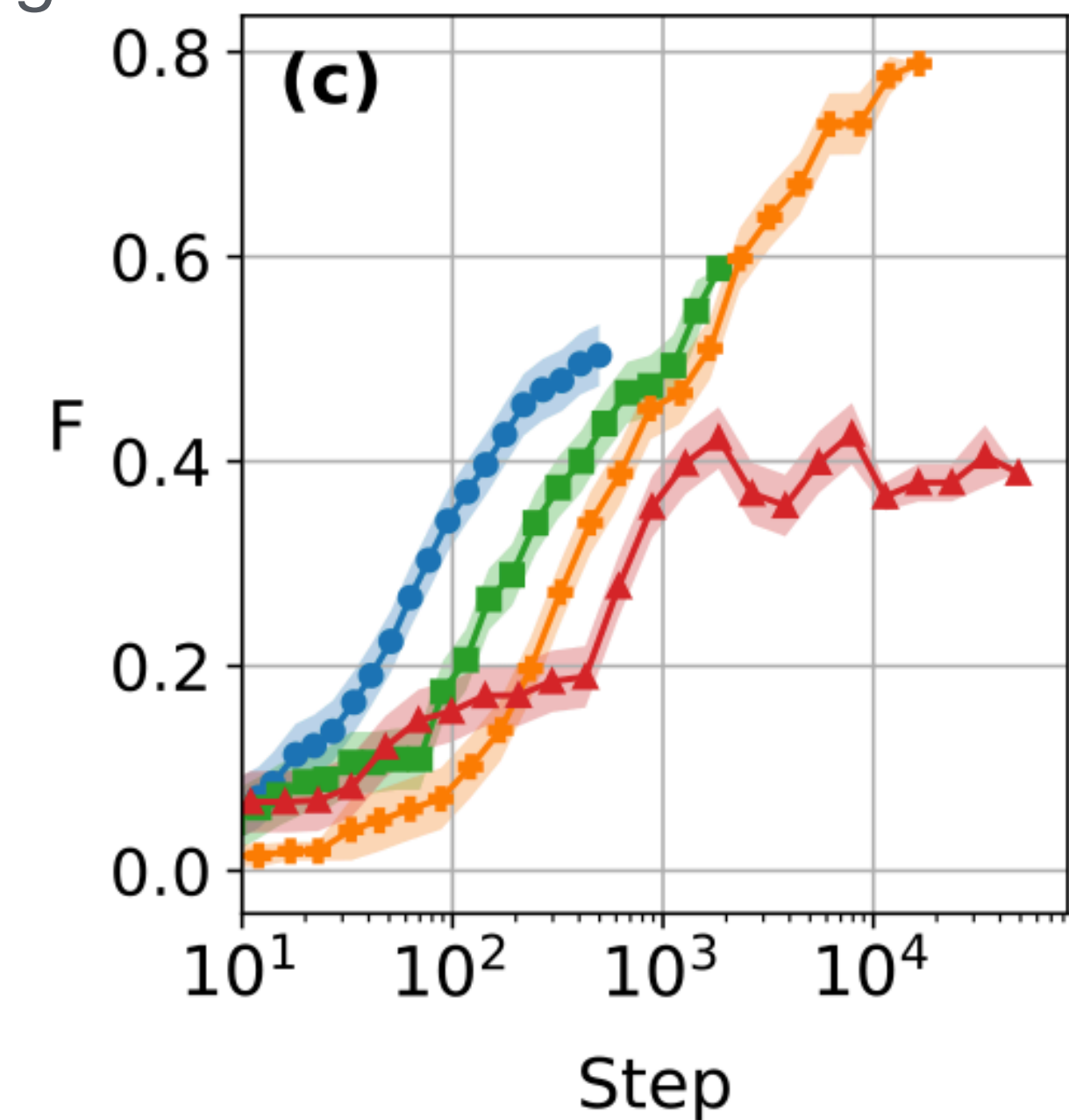
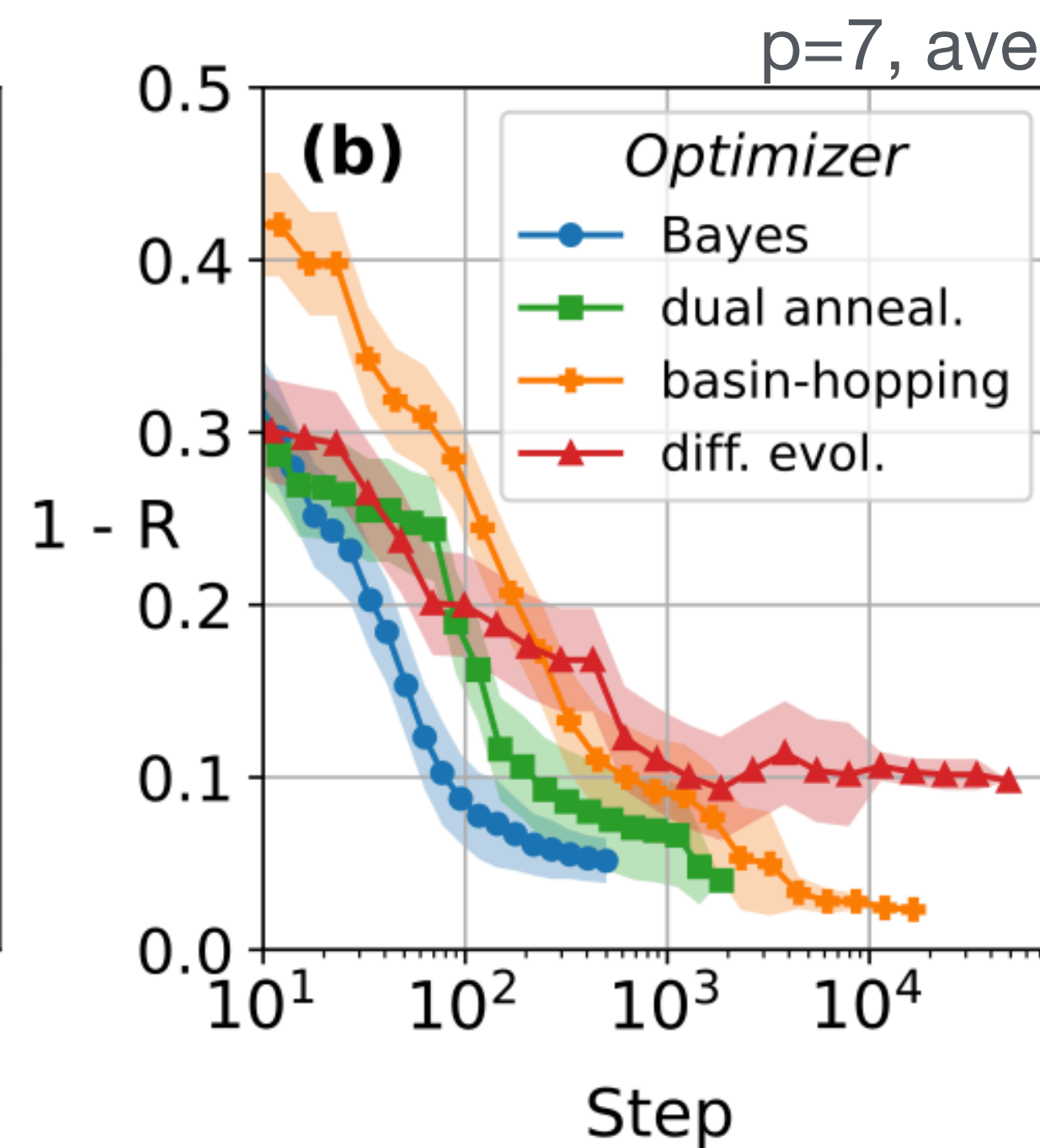
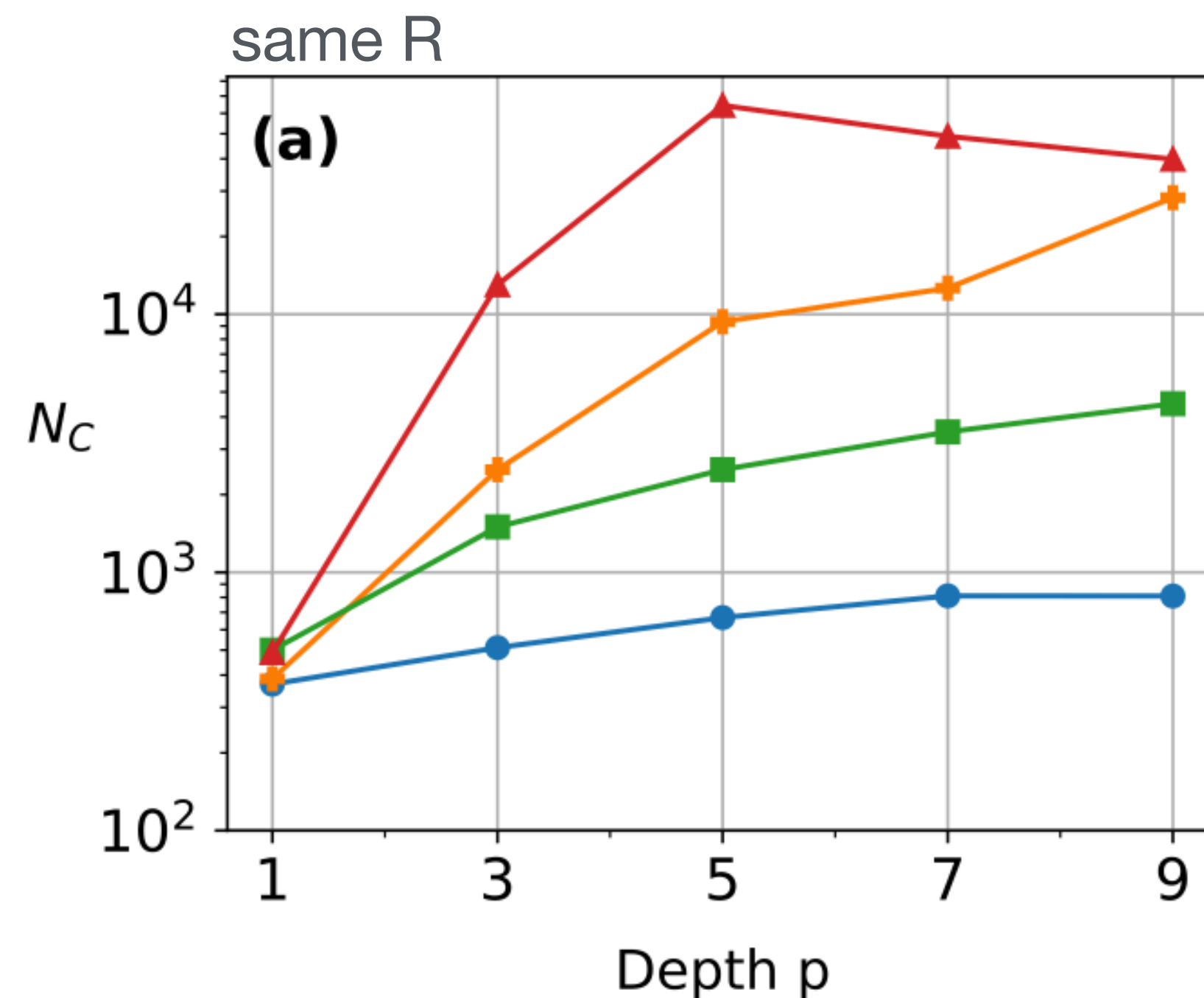
approximation ratio

$$R = E(\boldsymbol{\theta})/E_{GS}$$

fidelity

$$F = |\langle \boldsymbol{\theta} | z^* \rangle|^2$$

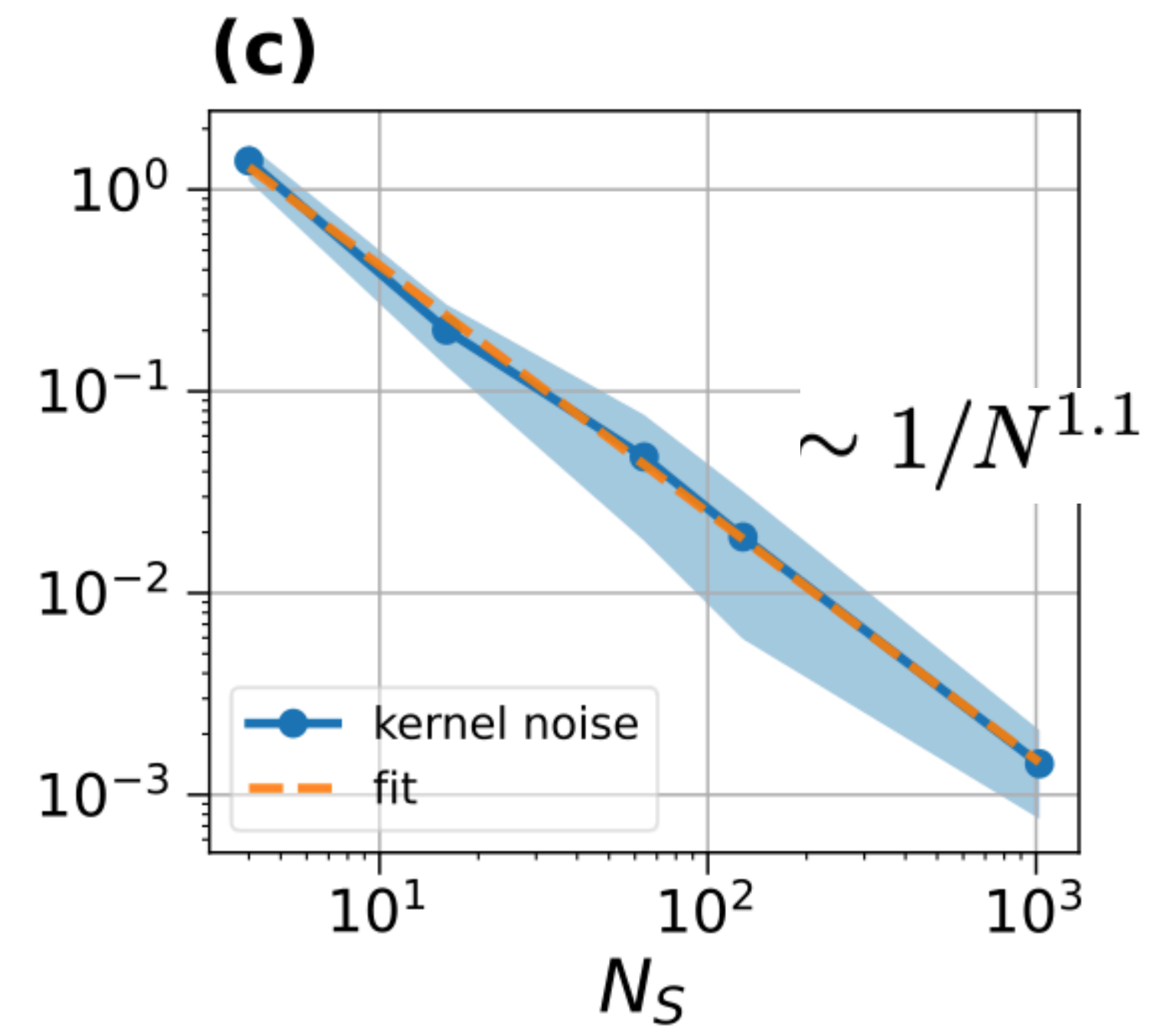
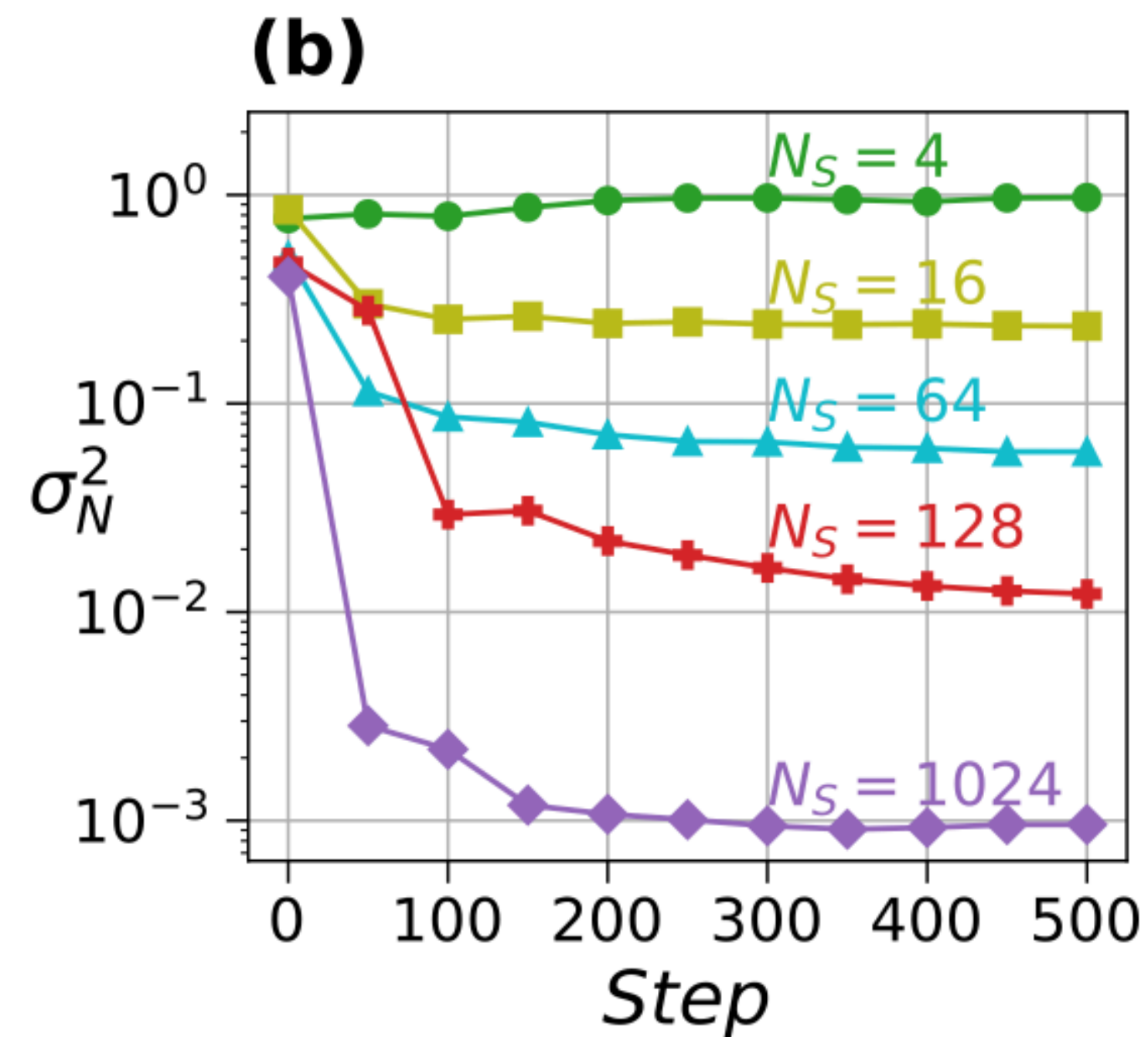
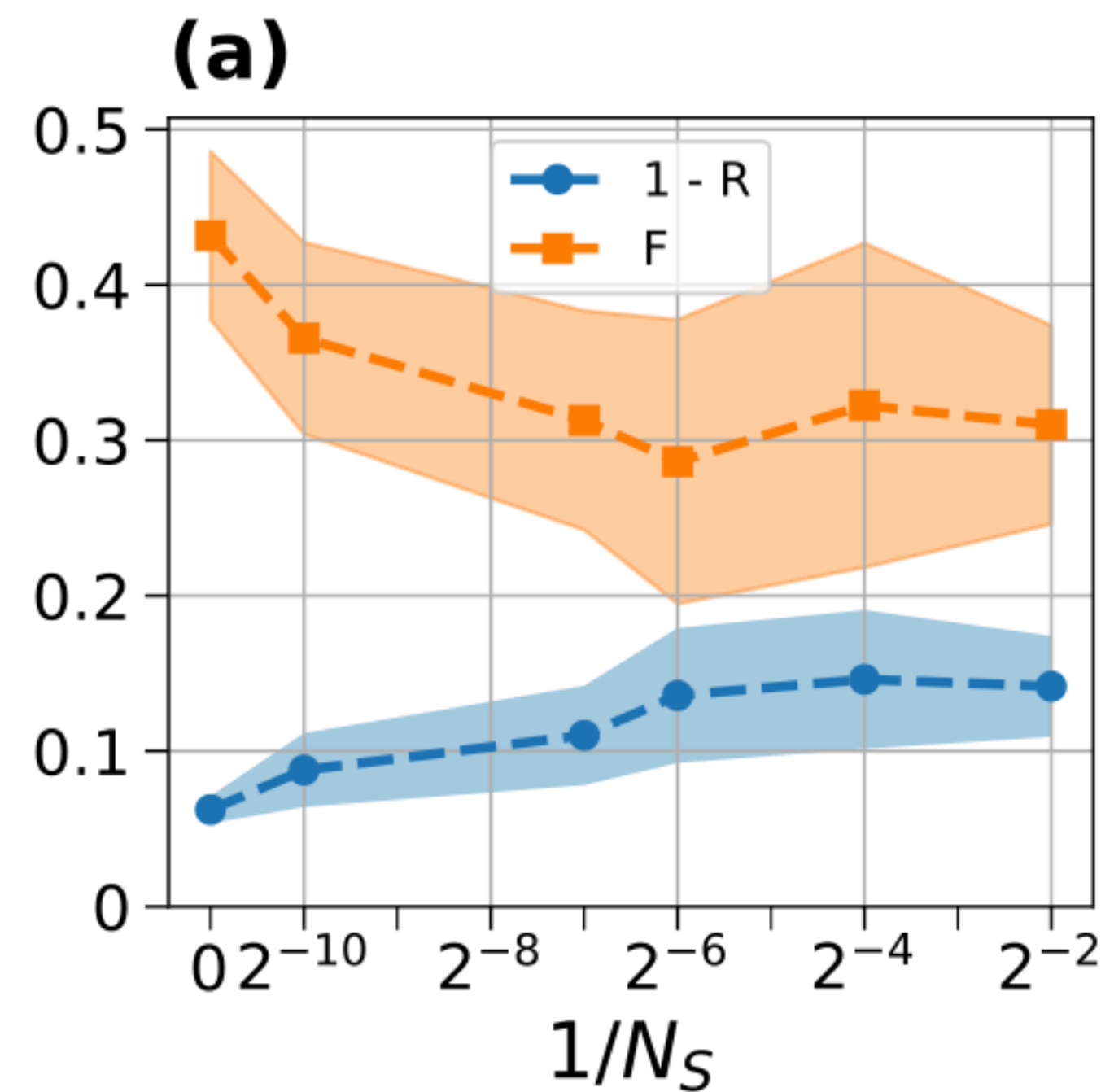
comparison among optimisers



slow measurements: $N_s = 1024, 128, 64, 16, 4$

reducing the numbers of measurements, decreases the precision of the energy landscape reconstruction

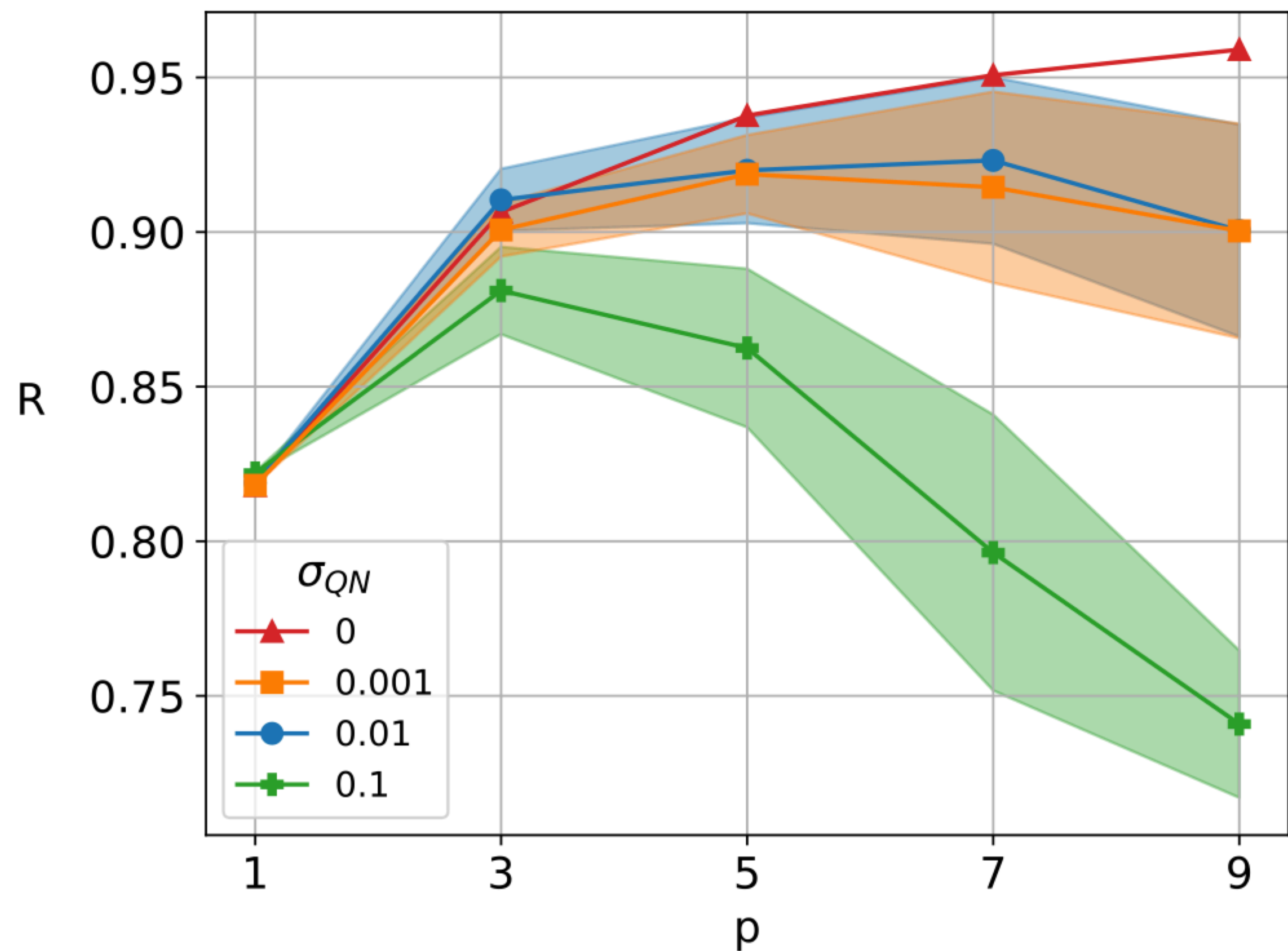
this is a multinomial sampling problem, thus variance of energy should depend on N_s^{-1}



quantum noise:

quantum circuits are subject to different kinds of noise (device dependent)

$$|\boldsymbol{\theta}\rangle = \prod_{l=1}^p e^{-i \sum_i \beta_l^i X_i} e^{i \sum_{\langle i,j \rangle} \gamma_l^{(i,j)} Z_i Z_j} |+\rangle,$$




R is expected to decrease with depth p (that increases the number of evolutions) and by increasing the q. noise variance σ_{QN}

However, up to p=3, R increases, signalling that Bayesian optimization is robust

S. Tibaldi, D. Vodola, E. Tignone. E.E.
arXiv:2209.03824

AND NOW?

- Collaboration, via CINECA, with  to test our Bayesian approach on their Rydberg atom platform
 - ✓ Run simulations on Pulser with their real optical pulses to generate gates and with (different kinds of) noise
 - ★ To be run on real atomic platform
- Applications to different problems:
quantum chemistry, many body hamiltonians

**Thank you for
your attention**