

QUANTUM ANNEALING SPEEDUP

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Supervisors: *Oriel Kiss, Michele Grossi, Sofia Vallecora*

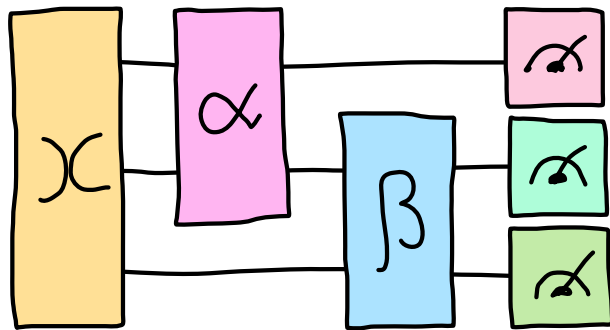
a project for



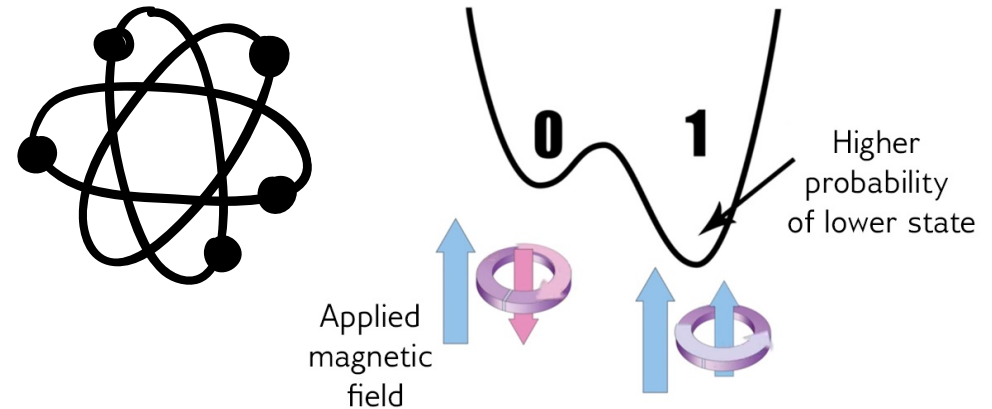
16/08/2023

Quantum Computation

The circuit model is not the only paradigm for quantum computation.



GATE BASED



ADIABATIC COMPUTING and ANNEALING

evolve a system towards low-energy states

Quantum Annealing

we love to describe systems
with **Hamiltonians**

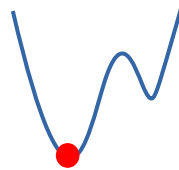
 $\hat{\mathcal{H}}$ 

Seems a good idea to diagonalize it, ya?

$$\hat{\mathcal{H}} |\psi_i\rangle = E_i |\psi_i\rangle$$

E_i is the energy of (eigen)state $|\psi_i\rangle$

The eigenvector associated to the lowest energy eigenvalue is called the **ground state**.

 $|GS\rangle$ 

An Hamiltonian is an operator that describes the quantum system in terms of energy. Thus, **the eigenvalues are the energy of our system.**

KEY IDEA OF QA and QAC



- encode the target problem as \mathcal{H}_p
- each state is a solution to the problem
- the GS is the **optimal solution**

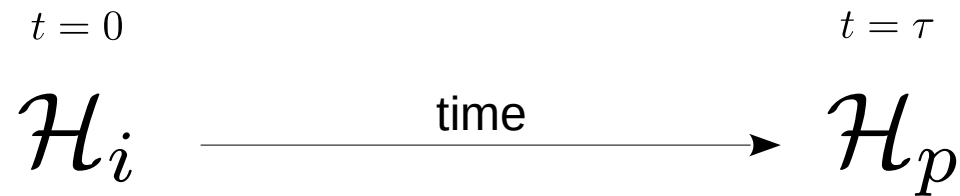
Example problems: last mile resupply problem, layout planning, route planning, molecular unfolding, machine learning, lattice gauge, ...

Further reading

https://docs.dwavesys.com/docs/latest/c_gs_2.html
<https://www.dwavesys.com/learn/featured-applications/>

How to do it?

The system evolves in time, thus $\hat{\mathcal{H}}(t)$:



We can prepare a state here...

... and measure the final state

Quantum Annealing (QA)

the initial state is **metaheuristic**, the transverse field allows to explore the hypersurface of solutions

Adiabatic Quantum Computation (AQC)

the initial state is ground-state and easy to prepare, the **system evolves slowly**

Life is not easy for any of us.

Adiabatic theorem(s)

Such theorems bound the **minimum total evolution time**, such as:

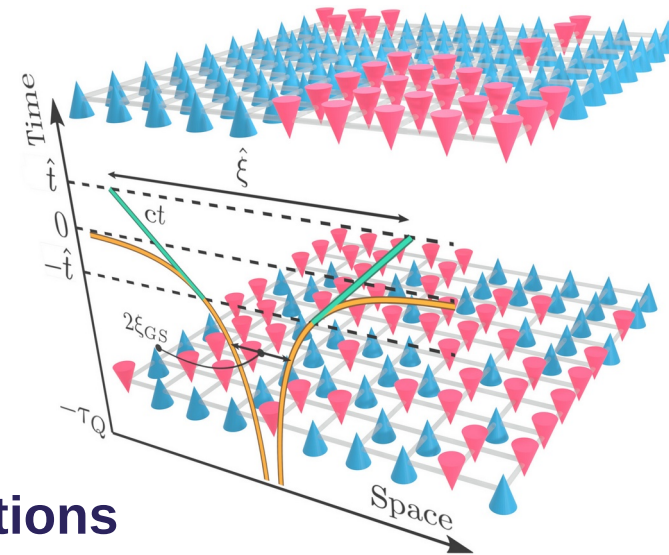
$$\tau \geq \frac{1}{\delta} \left(\int_0^s \left[\frac{\|\partial_s^2 H(\sigma)\|}{\Delta^2(s)} + 7 \frac{\|\partial_s H(\sigma)\|^2}{\Delta^3(s)} \right] d\sigma + B \right)$$

to guarantee that

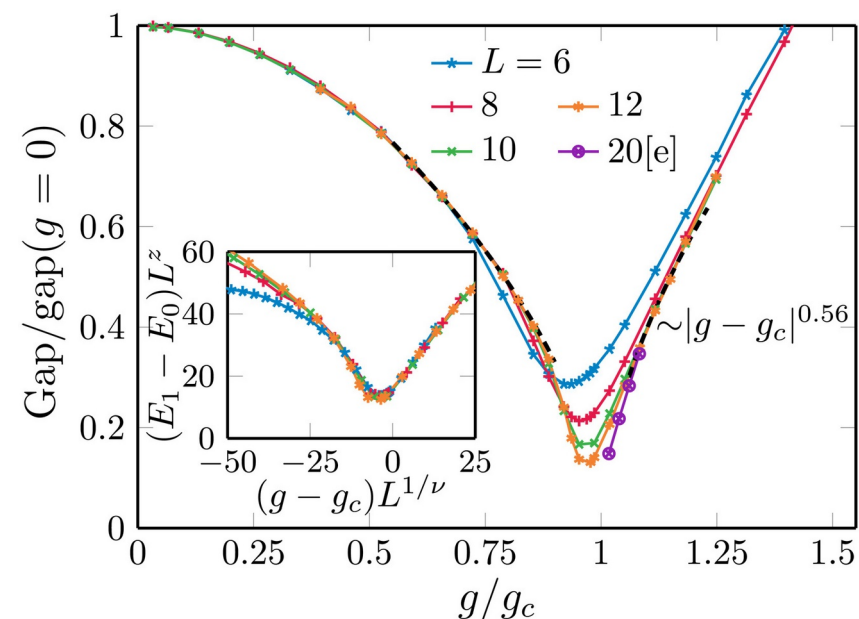
$$|\langle \psi(s) | \psi^T(s) \rangle| \geq 1 - \delta \quad \forall s \in [0, 1]$$

where $|\psi(s)\rangle$ instantaneous GS of $\mathcal{H}(s)$

$\Delta(s)$ gap between GS and first excited eigenstate



Phase transitions



Possible solutions

QUANTUM OPTIMAL CONTROL

Use numerically optimized paths to avoid gap closures.

SHORTCUTS TO ADIABATICITY (STA)

Remove diabatic transitions instantaneously.

> COUNTERDIABATIC DRIVING (CD) *or transitionless driving*

Suppress losses that arise due to fast deformations of the system via **analytical** compensation.

$$\mathcal{H}_{CD} = \mathcal{H}_0(\lambda) + \dot{\lambda} \hat{A} \quad \text{st} \quad \hat{A} = i\hbar \sum_{m \neq l} \frac{\langle \epsilon_m(\lambda) | \partial_\lambda \mathcal{H}_0(\lambda) | \epsilon_l(\lambda) \rangle}{\epsilon_l(\lambda) - \epsilon_m(\lambda)} |\epsilon_m(\lambda)\rangle \langle \epsilon_l(\lambda)|$$

> APPROXIMATE COUNTERDIABATIC DRIVING

$$\mathcal{A} \simeq \alpha \sum_i \sigma_i^y + \gamma \sum_i (\sigma_i^x \sigma_{i+1}^y + \sigma_i^y \sigma_{i+1}^x) + \xi \sum_i (\sigma_i^y \sigma_{i+1}^z + \sigma_i^z \sigma_{i+1}^y)$$

$$\text{!} \min_{\mathcal{A}} \text{Tr}[G_t(\mathcal{A})^2]$$

$$G_t = -\partial_t \mathcal{H}_0 + \frac{i}{\hbar} [\mathcal{A}, \mathcal{H}_0]$$

Target solution

Counterdiabatic Optimized Local Driving

Ieva Čepaitė, Anatoli Polkovnikov, Andrew J. Daley, Callum W. Duncan

$$\mathcal{H}_{\text{COLD}}(t, \beta) = H_0(t) + \underbrace{\alpha(t, \beta)\mathcal{O}_{\text{LCD}}}_{\text{CD}} + \underbrace{f(t, \beta)\mathcal{O}_{\text{opt}}}_{\text{QOC}}$$

My contribution

A library to manage the simulation of arbitrary spin systems annealing in the COLD formulation.

\mathcal{O}_{LCD} : can be local (first order) or non-local (higher orders)

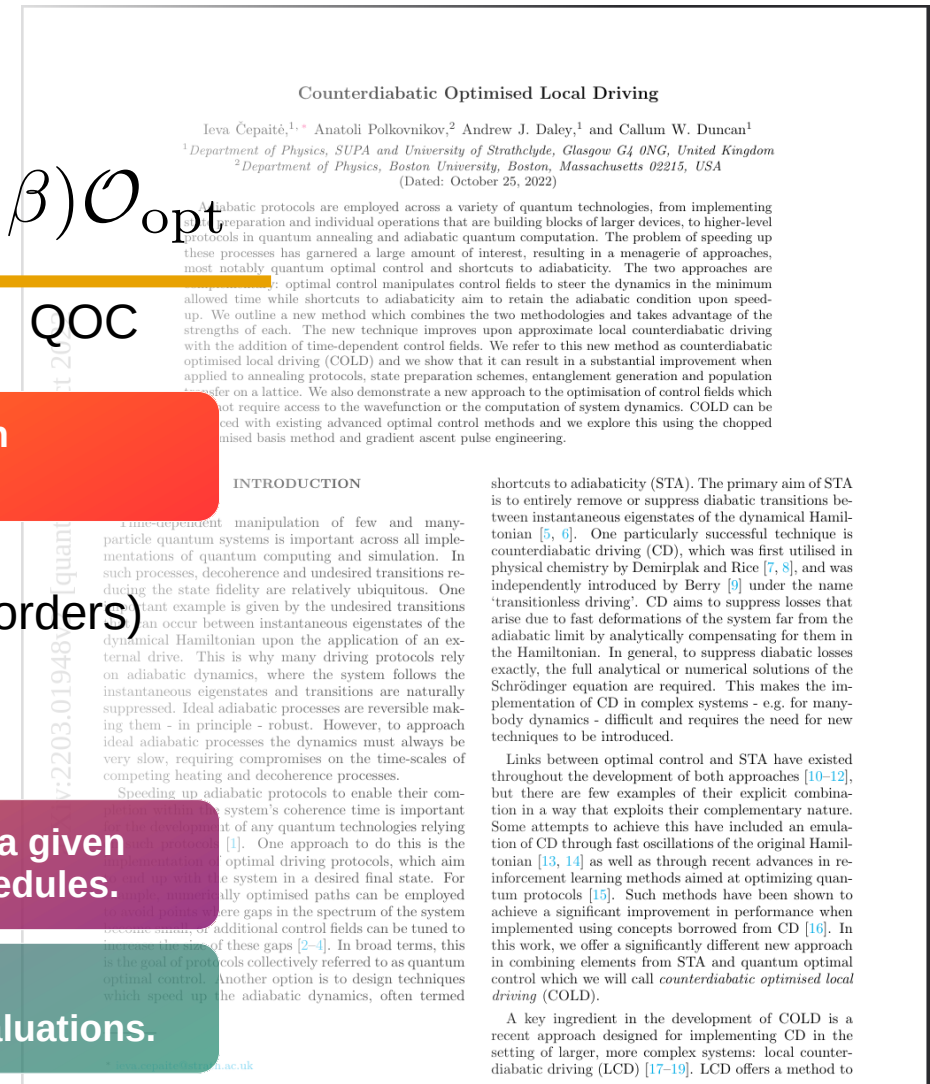
\mathcal{O}_{opt} : additional degrees of freedom

My contribution

The Gauge potential is computed symbolically for a given configuration, allowing fast prototyping of QA schedules.

My contribution

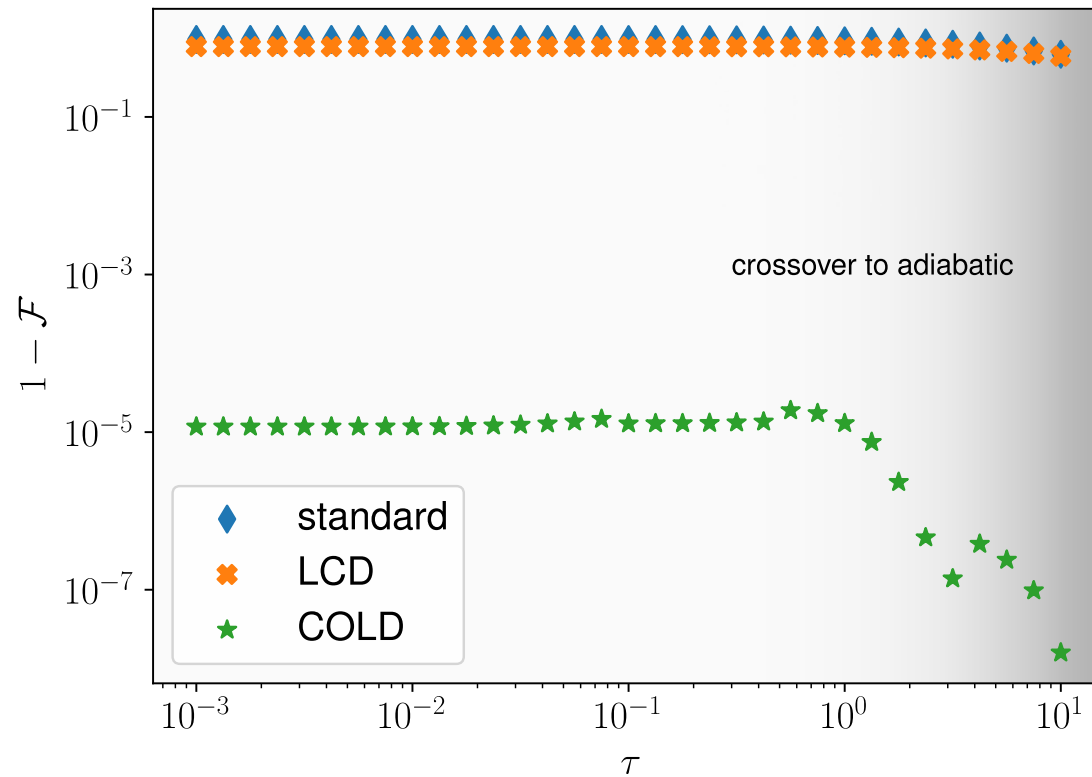
The QOC optimization is carried out with bayesian sampling methods, to reduce the loss function evaluations.



Results

ISING MODEL

$$\mathcal{H} = -J(t) \sum_i \sigma_i^z \sigma_{i+1}^z + X(t) \sum_i \sigma_i^x + Z(t) \sum_i \sigma_i^z$$



$$-\sum_i \sigma_i^x \longrightarrow \mathcal{H}_{\text{Ising}}$$

1D lattice: 5 spins, nearest neighbor interactions

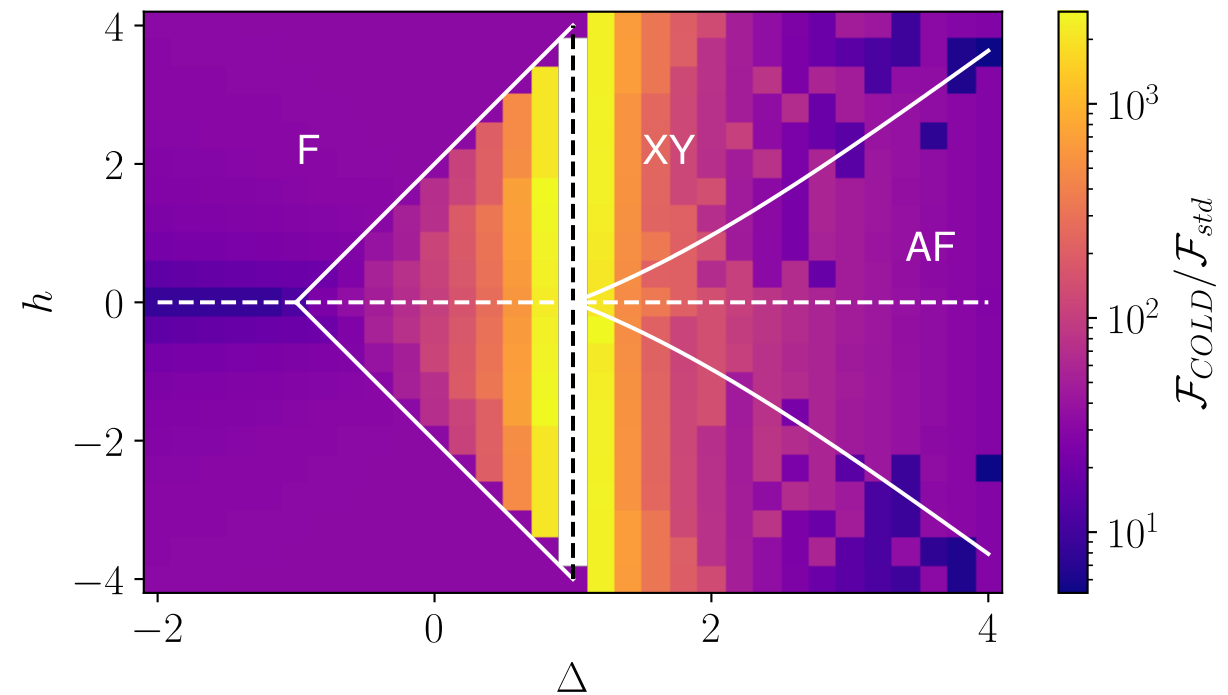
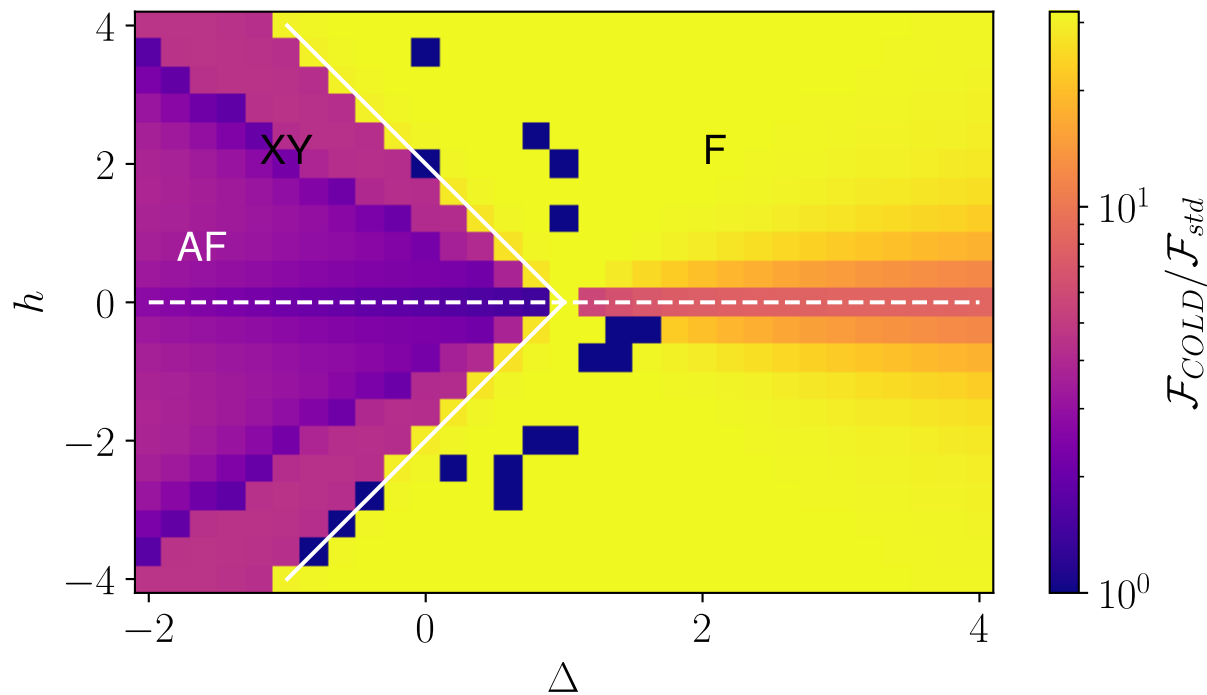
Simulated annealing: > 2000 steps, min timestep constant in magnitude

Results

XXZ MODEL
$$\mathcal{H} = J(t) \left(\sum_i \sigma_i^x \sigma_{i+1}^x + \sum_i \sigma_i^y \sigma_{i+1}^y \right) + \Delta(t) \sum_i \sigma_i^z \sigma_{i+1}^z + h(t) \sum_i \sigma_i^z$$

$$-\sum_i \sigma_i^x \rightarrow \mathcal{H}(J = 1, -h, -\Delta)$$

$$-\sum_i \sigma_i^x \rightarrow -\mathcal{H}(J = 1)$$



i Simulation with annealing time 0.01, 200 discrete steps on 5 spin systems.

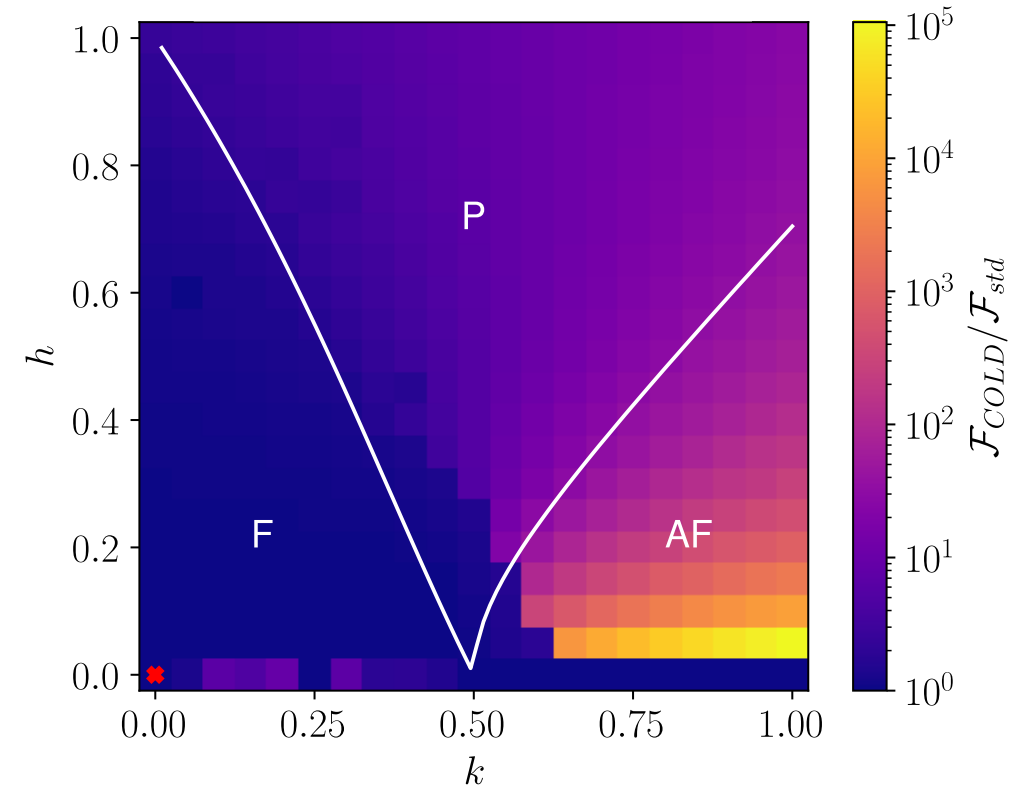
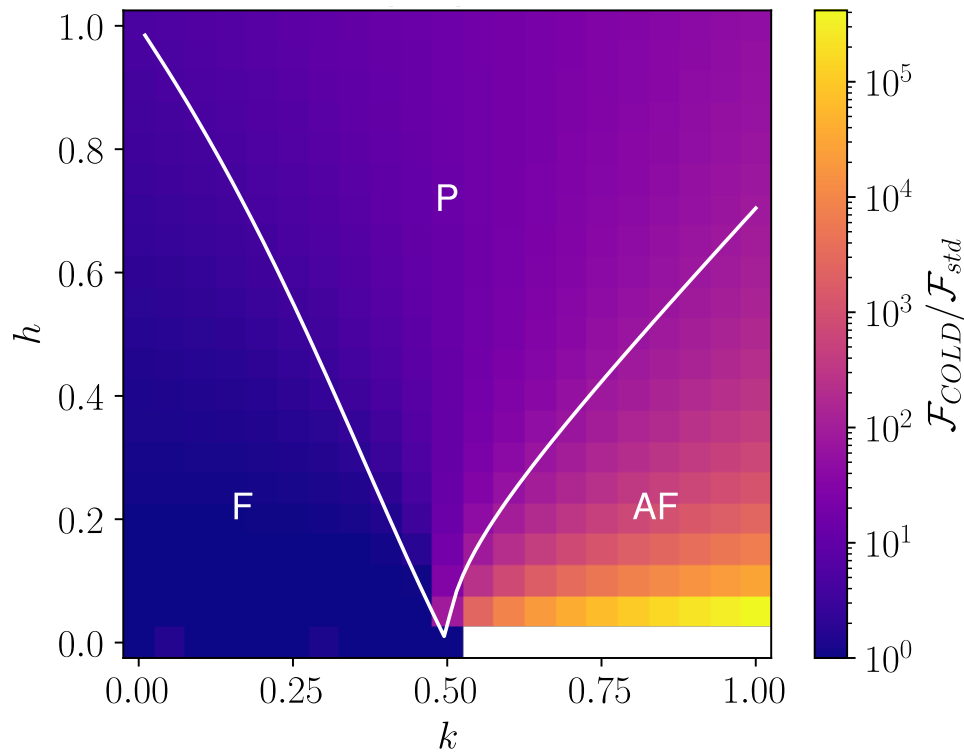
Results

ANNNI MODEL

$$\mathcal{H} = -J(t) \sum_i \sigma_i^z \sigma_{i+1}^z + k(t) \sum_i \sigma_i^z \sigma_{i+2}^z + h(t) \sum_i \sigma_i^z$$

$$-\sum_i \sigma_i^x \longrightarrow \mathcal{H}_{\text{ANNNI}}$$

$$\mathcal{H}_{\text{ANNNI}}^{(\text{ferro})} \longrightarrow \mathcal{H}_{\text{ANNNI}}$$



i Simulation with annealing time 0.01, 200 discrete steps on 5 spin systems.

Goal: prepare a good initial state
for QPE algorithms



QUANTUM
TECHNOLOGY
INITIATIVE



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I would like to say

*Thanks to my supervisors, Oriel, Michele and Sofia.
Thanks to the openlab staff for this opportunity.
Thanks to all the students for being here this summer.*



Bibliography

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- [Schmitt2022] **Quantum phase transition dynamics in the two-dimensional transverse-field Ising model** (doi.org/10.1126/sciadv.abl6850)
- **D-Wave Systems** (<https://www.dwavesys.com/>)

Attributions

Slide 2 – PennyLane illustration of quantum circuit

Slide 3 – Einstein illustration, credits to Albert Vectors by Vecteezy