

# Quantum Algorithms for Simulating Quantum Field Theories

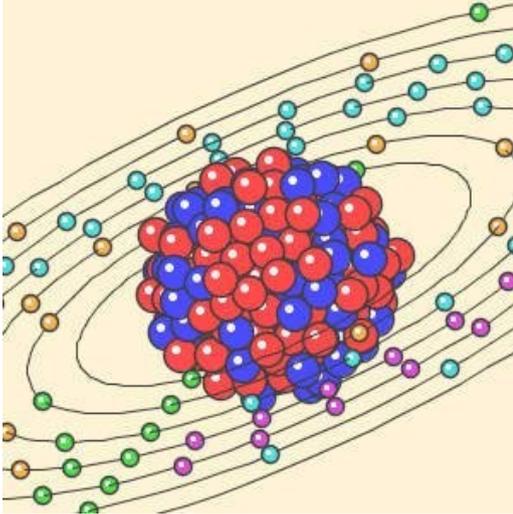
Stephen Jordan

joint work with:

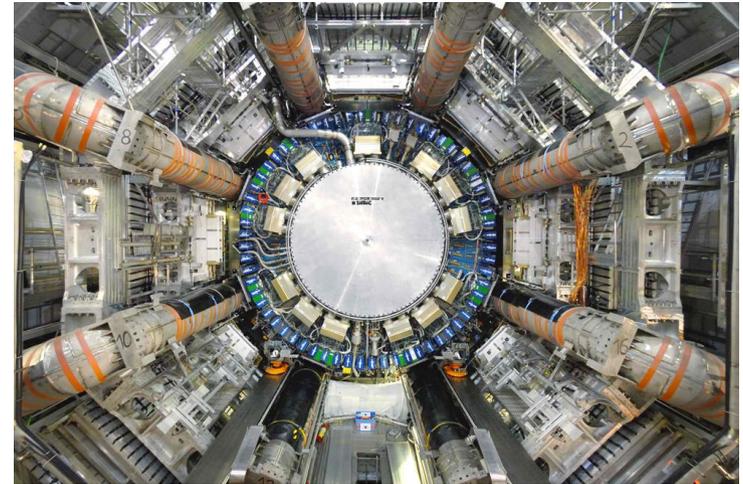
Keith Lee

John Preskill

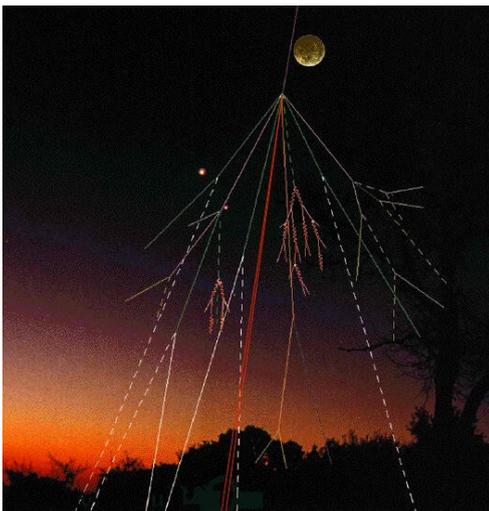
# When do we need QFT?



Nuclear Physics



Accelerator Experiments

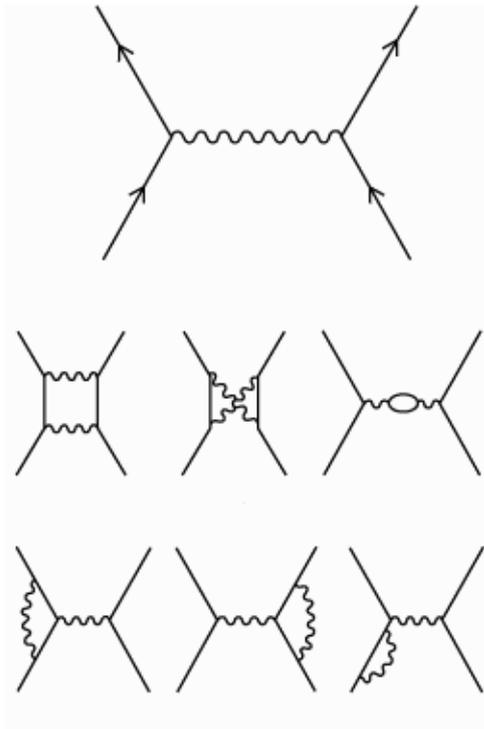


Cosmic Rays

→Whenever quantum mechanical and relativistic effects are both significant.

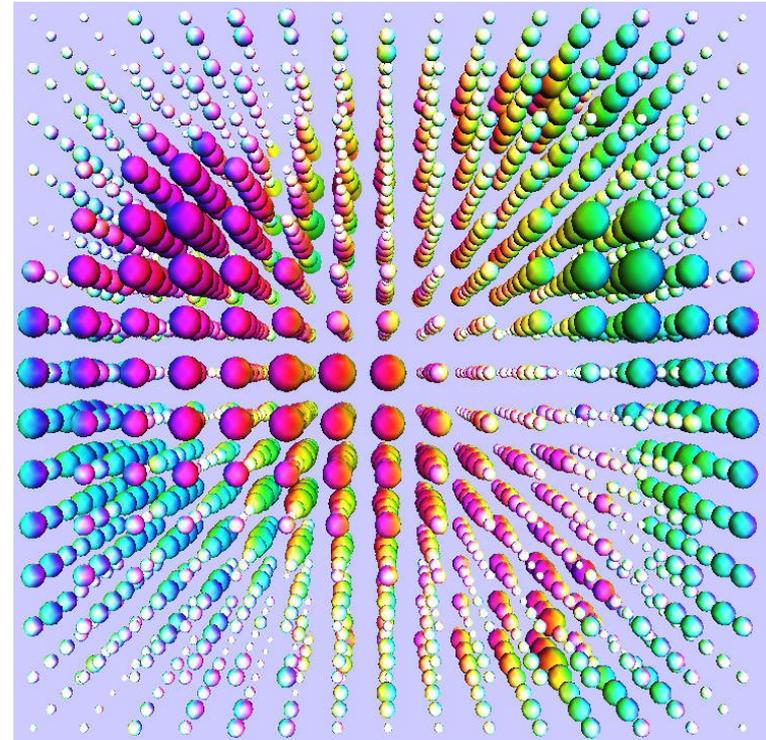
# Classical Algorithms

## Feynman Diagrams



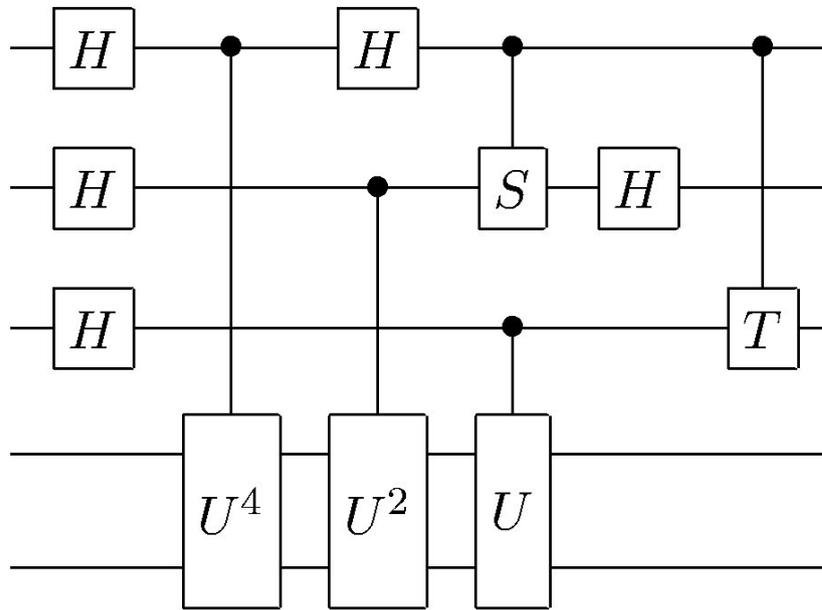
break down at strong  
coupling or high precision

## Lattice Methods

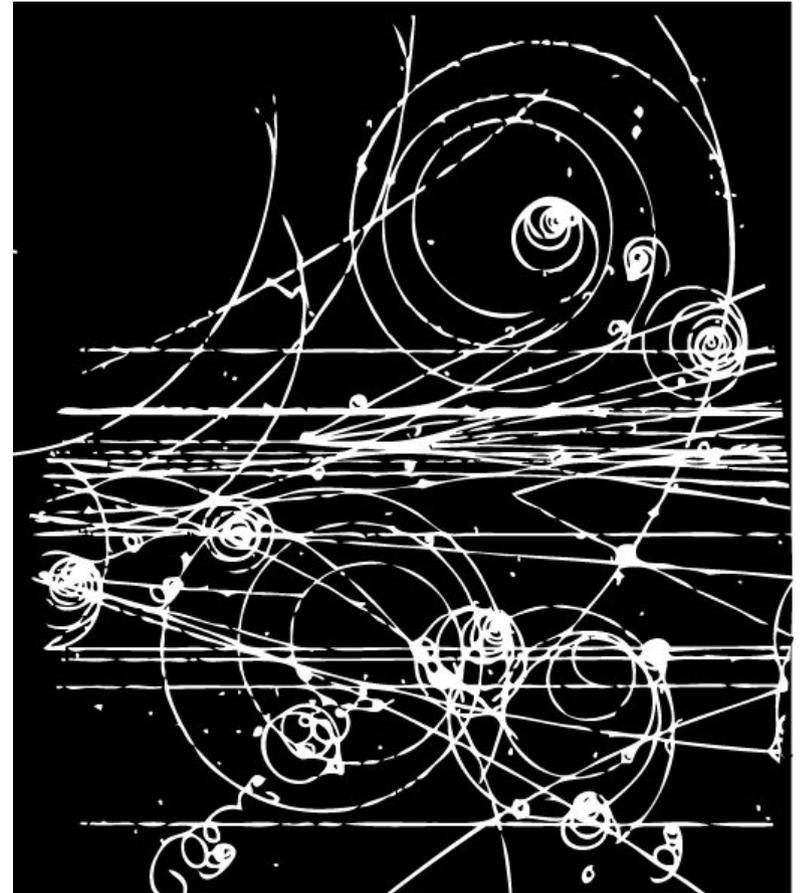


cannot calculate dynamical  
quantities

# What is the computational power of our universe?



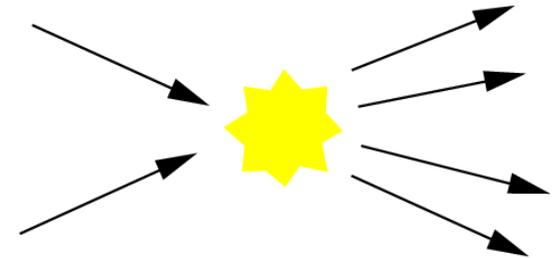
→  
simulate  
←



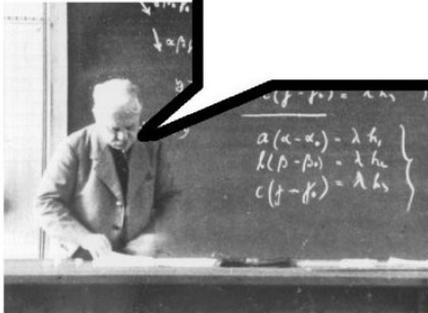
# A QFT Computational Problem

**input:** a list of momenta of incoming particles

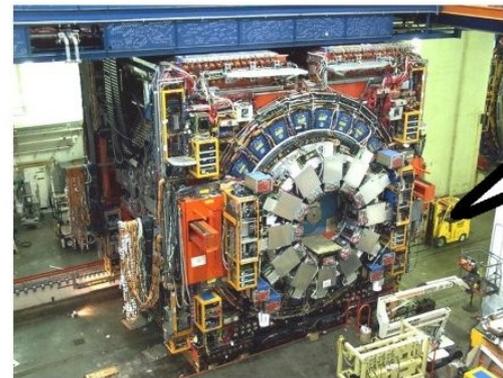
**output:** a list of momenta of outgoing particles time  $t$  later



S-matrix



Particle  
accelerator



I will present a polynomial-time quantum algorithm to compute scattering amplitudes in  $\phi^4$ -theory with nonzero mass.

$\phi^4$ -theory is a simple toy model, but it illustrates some of the main difficulties of simulating QFTs

Discretizing spacetime

Preparing initial states

Measuring observables

Time evolution is the easy part (use Trotter)

# $\phi^4$ -theory

Lagrangian density:

$$\mathcal{L} = \frac{1}{2} \partial^\mu \phi \partial_\mu \phi - m^2 \phi^2 - \frac{\lambda}{4!} \phi^4$$

For quantum simulation, we prefer Hamiltonian formulation (equivalent):

$$H = \int d^d x \left[ \pi^2 + (\nabla \phi)^2 + m^2 \phi^2 + \lambda \phi^4 \right]$$

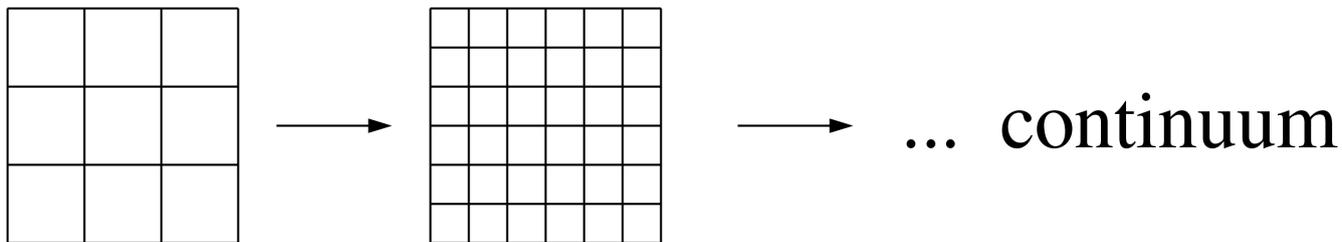
$$[\phi(x), \pi(y)] = i\delta^{(d)}(x - y)$$

# Lattice Cutoff

To get finite answers one needs a cutoff. We use a spatial lattice:

$$H(a) = \frac{1}{2} \sum_{x \in \Omega} a^d [\pi^2 + (\nabla \phi)^2 + m^2 \phi^2 + \lambda \phi^4]$$

Continuum QFT = limit of a sequence of theories on successively finer lattices:



***m* and  $\lambda$  are functions of *a* !**

After imposing a spatial lattice we have a many-body quantum system with a local Hamiltonian

Simulating the time evolution in polynomial time is a **solved problem**

Standard methods scale as  $N^2$  We can do  $N$  .)

Questions:

Convergence as  $a \rightarrow 0$

How to prepare initial wavepackets

How to measure particle momenta

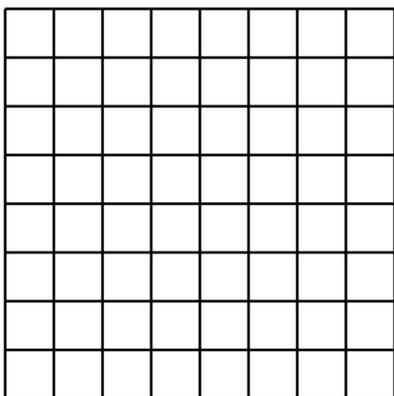
# Convergence as $a \rightarrow 0$

Particles of energy  $E$  have associated

lengthscale  $\frac{hc}{E}$

Intuition: if  $a \ll \frac{hc}{E}$  discretization errors should be small

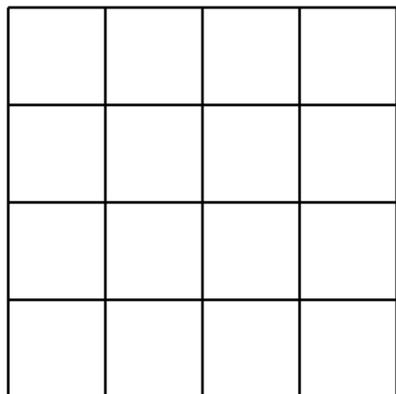
We can formalize this by analyzing a sequence of effective Hamiltonians (RG flow)



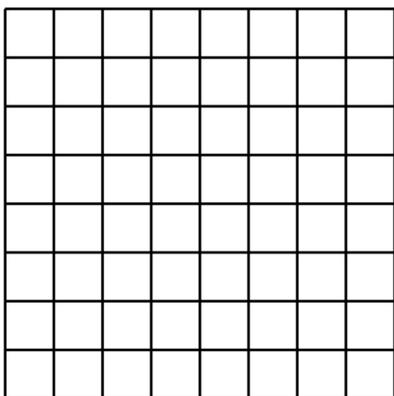
$$H = \frac{1}{2} \sum_{x \in \Omega} a^d [\pi^2 + (\nabla \phi)^2 + m^2 \phi^2 + \lambda \phi^4]$$



Coarse grain



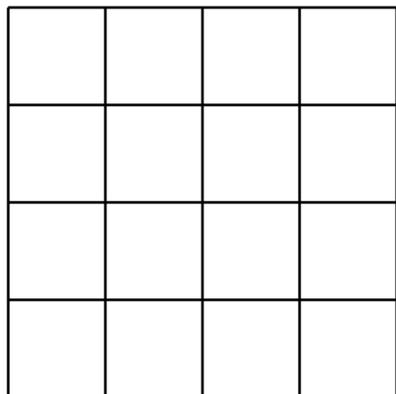
$$H_{\text{eff}} = \frac{1}{2} \sum_{x \in \Omega'} (2a)^d [\pi^2 + (\nabla' \phi)^2 + m_{\text{eff}}^2 \phi^2 + \lambda_{\text{eff}} \phi^4 + g \phi^6 + \dots]$$



$$H = \frac{1}{2} \sum_{x \in \Omega} a^d [\pi^2 + (\nabla \phi)^2 + m^2 \phi^2 + \lambda \phi^4]$$



Coarse grain



$$H_{\text{eff}} = \frac{1}{2} \sum_{x \in \Omega'} (2a)^d [\pi^2 + (\nabla' \phi)^2 + m_{\text{eff}}^2 \phi^2 + \lambda_{\text{eff}} \phi^4 + g \phi^6 + \dots]$$

How does simulation convergence as  $a \rightarrow 0$ ?

Answer: as  $a^2$

TODO:

Efficiently prepare initial wavepackets

Efficiently measure particle momenta

# Creating a particle

$a_p^\dagger$  creates a particle of momentum  $p$  and

energy  $\omega_p = \sqrt{p^2 + m^2}$

We can define single-particle states in the interacting theory by adiabatic continuation from the free theory.

Quantum State-Preparation Algorithm:

1. Build the free vacuum (Gaussian)
2. Excite wavepackets
3. Adiabatically turn on interaction

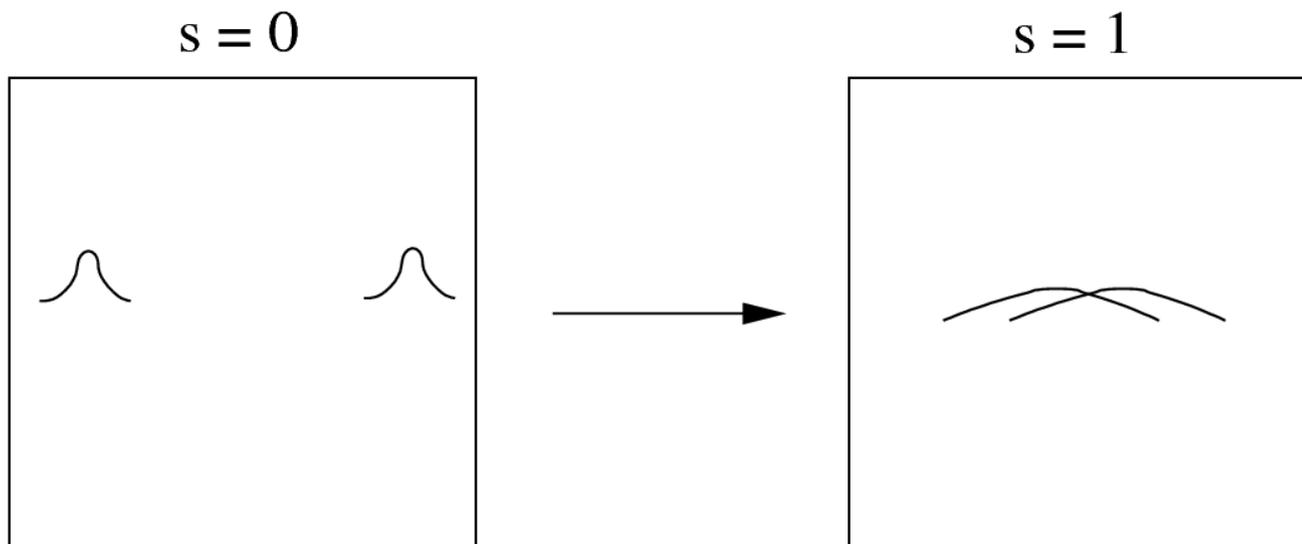
# Adiabatically Turn on Interaction

$$H(s) = \sum_{x \in \Omega} \left[ \pi^2 + (\nabla \phi)^2 + m^2 \phi^2 + s \lambda \phi^4 \right]$$

Use standard Trotter technique to simulate  $H(s)$   
with  $s$  slowly varying from 0 to 1

This **almost** works...

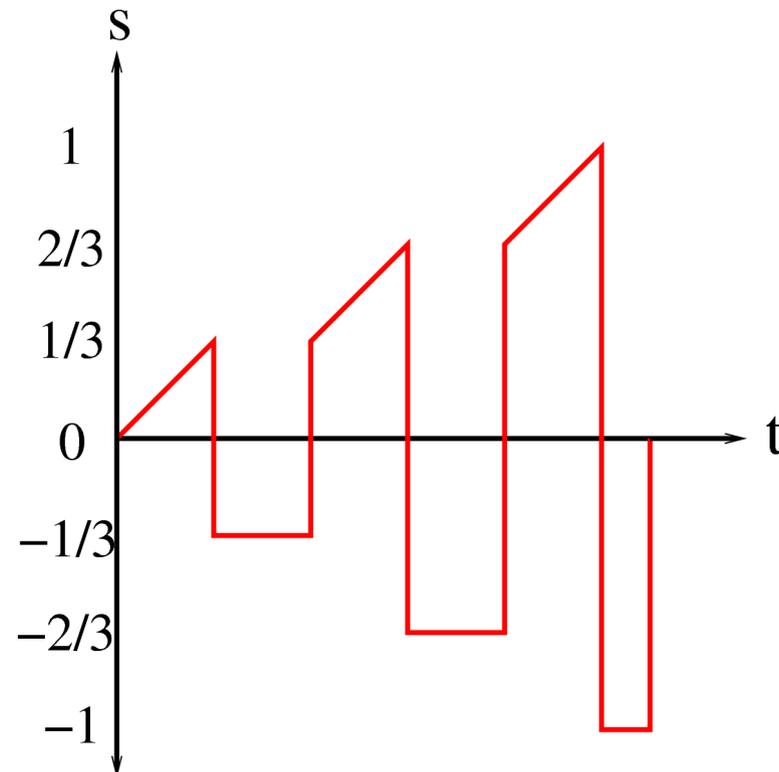
During all this slow time-evolution



the wavepackets propagate and broaden!

**Solution:** intersperse backwards time-evolutions with time-independent Hamiltonians

This winds back the dynamical phase on each eigenstate, without undoing the adiabatic change of eigenbasis.



# Adiabaticity

Q. How slowly must we vary  $s$ ?

A. By the adiabatic theorem, this is determined by the energy gap against excitations. There are two important excitation processes:

particle creation from vacuum:  $\gamma = m_{\text{phys}}$

one particle splits into three:  $\gamma \simeq \frac{m_{\text{phys}}^2}{p}$

# Measurements

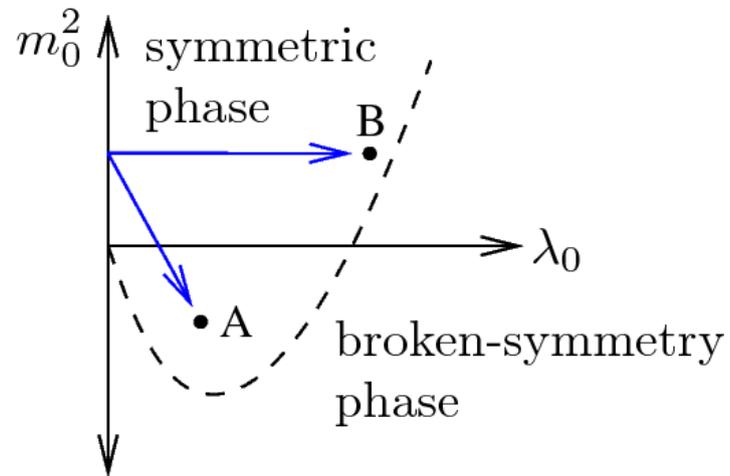
In the free theory, use the number operators:  $a_p^\dagger a_p$

Defining particle number is hard in the interacting theory!

Adiabatically go back to the free theory and measure the number operators using Kitaev's phase estimation technique.

# Strong Coupling

$\phi^4$ -theory in 1+1 and 2+1 dimensions has phase transition in which  $\phi \rightarrow -\phi$  symmetry is spontaneously broken



Near phase transition perturbation theory breaks down and mass gap vanishes:

$$m_{\text{phys}} \sim (\lambda_c - \lambda)^\nu \quad \nu = \begin{cases} 1 & d = 1 \\ 0.63 \dots & d = 2 \end{cases}$$

# Strong Coupling

The only part of the quantum algorithm affected by strong coupling is adiabatic state preparation.

Splitting: 
$$\gamma \simeq \frac{m_{\text{phys}}^2}{p} \sim \frac{(\lambda_c - \lambda)^{2\nu}}{p}$$

Creation from vacuum: 
$$\gamma = m_{\text{phys}} \sim (\lambda_c - \lambda)^\nu$$

Complexity is polynomial in  $(\lambda_c - \lambda)$  and  $p$

## Eventual goal:

Simulate the Standard model in BQP

## In progress:

Fermions

## Open problems:

Massless particles

Gauge symmetries

Phases inaccessible from free theory

# Conclusion

Quantum computers can efficiently calculate scattering amplitudes for  $\phi^4$  theory in 3 or fewer spatial dimensions.

Work remains to be done regarding more complicated theories.

I thank my collaborators:



Thank you for your attention.