

# **CPEN 400Q Lecture 18**

## **Hamiltonian simulation**

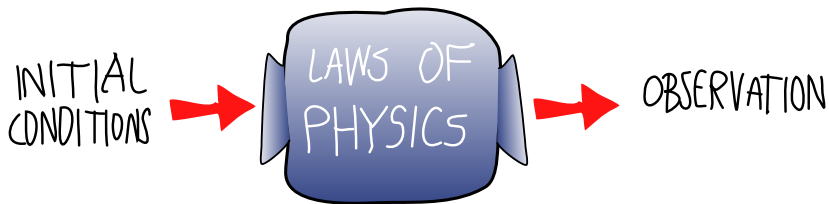
Wednesday 13 March 2024

# Announcements

- Hands-on 3 due Monday (18 March) at 23:59
- Assignment 3 due Wednesday (20 March) 23:59
- Quiz 8 on Monday (on today's lecture)

- Define a Hamiltonian and describe the relationship between its structure and the energy of a physical system
- Perform time evolution of simple systems in PennyLane

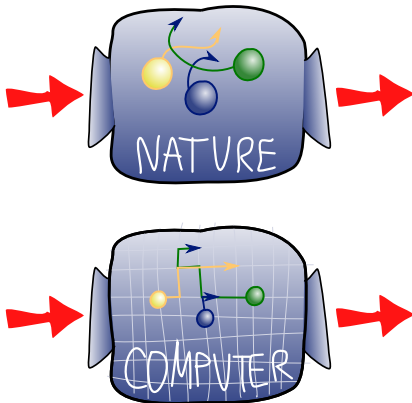
# Motivation



The role of a physicist is to work out what happens in the middle.

# Motivation

We can use computers to approximately model or simulate how a system evolves with time. This requires *discretizing* the systems.

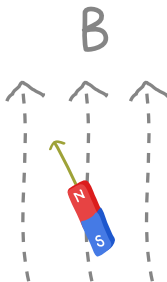


To do this on a quantum computer we need:

1. A meaningful way of describing physical systems using qubits
2. A way of manipulating qubits that corresponds to time evolution of that system
3. A way to extract relevant physical quantities

# Hamiltonians

Example: a bar magnet in a magnetic field along the  $Z$  direction.

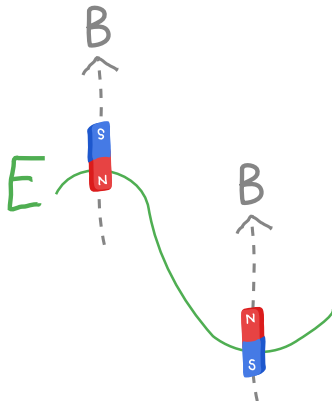


Every orientation of the magnet has an energy associated with it.

Image credit: Xanadu Quantum Codebook node H.3

# Hamiltonians

Bar magnets like to *align* with the external field.



Aligned configuration is the *ground state* (lowest energy state).



# Hamiltonians

Magnets have some orientation represented by a vector in 3-dimensional space. We can map this to a qubit.

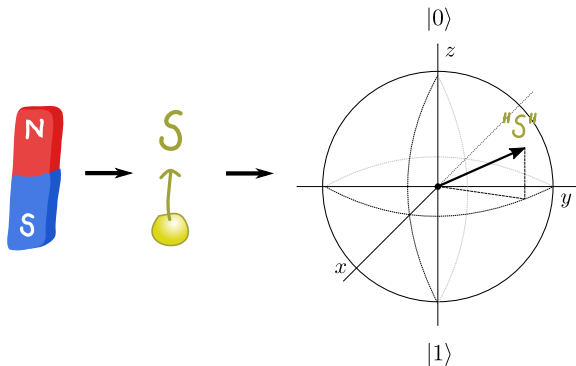


Image credit: Xanadu Quantum Codebook node H.5

# Hamiltonians

Can you think of a way to represent these magnets as qubit states?

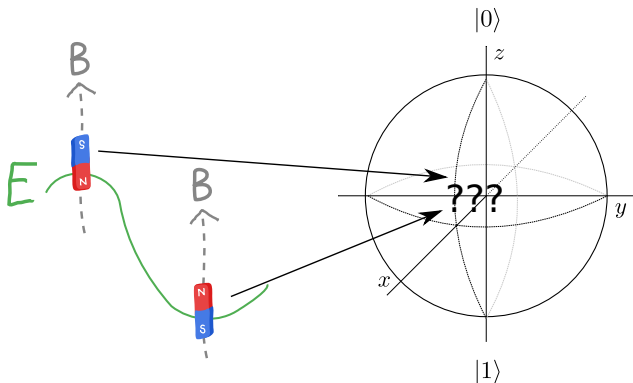


Image credit: Xanadu Quantum Codebook node H.3

# Hamiltonians

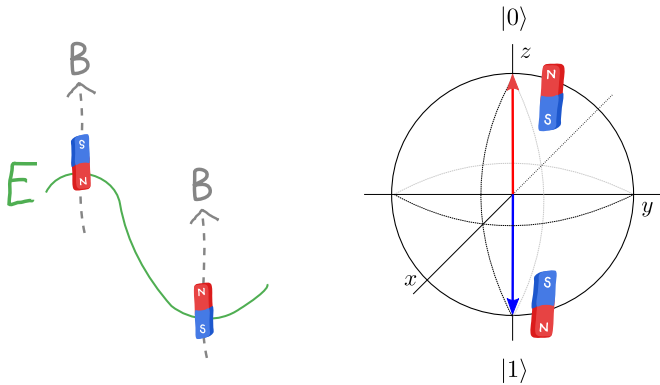


Image credit: Xanadu Quantum Codebook node H.3

# Hamiltonians

What can we use to denote the *energy* of the magnet?

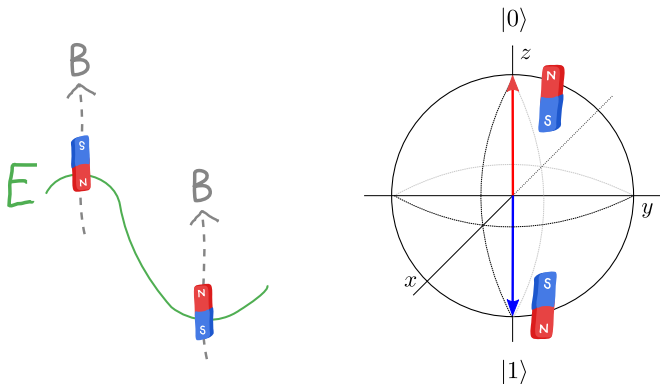
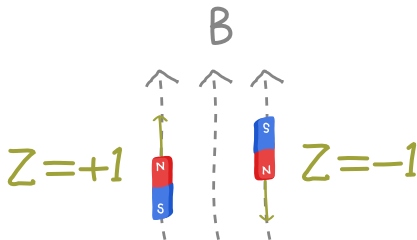


Image credit: Xanadu Quantum Codebook node H.3

# Hamiltonians

A Hermitian operator called a *Hamiltonian* describes the system:

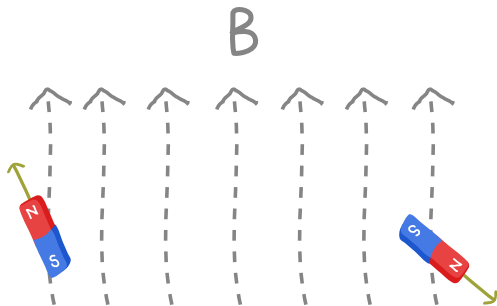
where  $\alpha$  is a coefficient that depends on physical values.



Computing the energy of the system corresponds to *measuring the expectation value of  $\hat{H}$*  (which in this case, is just  $-\alpha\langle Z \rangle$ ).

# Hamiltonians

Hamiltonians can describe systems with multiple parts. For example, two magnets that are far apart:



The energy of this system is

# Hamiltonians

Question: when is the energy of this system minimized?

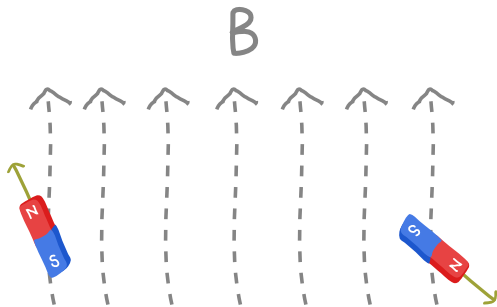
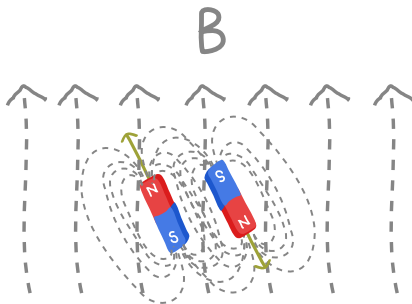


Image credit: Xanadu Quantum Codebook node H.5

# Hamiltonians

If the two magnets are close, we need parts of the Hamiltonian that describe the *interaction*.





# Hamiltonians

Question: What is the energy? What is the ground state (energy)?

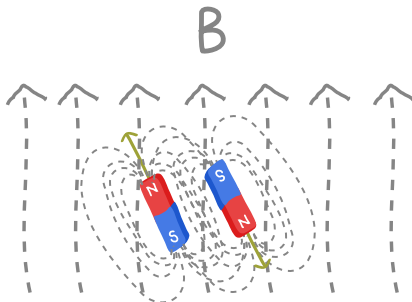


Image credit: Xanadu Quantum Codebook node H.5

More generally, a Hamiltonian can be expressed as

i.e., as a *linear combination of Pauli operators*.

Key points:

- An  $n$ -qubit Hamiltonian is a giant  $2^n \times 2^n$  Hermitian matrix
- The Paulis are a group and form a basis for Hermitian matrices

Computation of expectation values is linear:

**Exercise:** Consider two qubits interacting under the Hamiltonian

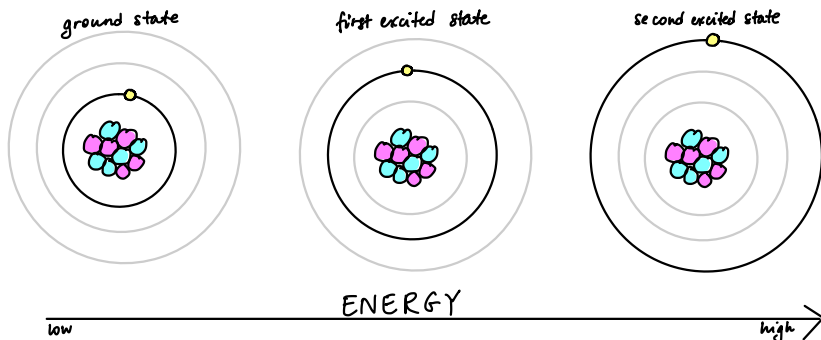
$$\hat{H} = -3(Z_0 + Z_1) + 2Z_0Z_1 + 4(X_0 + X_1)$$

What is the energy of the system if the qubits are in state

$$|\psi\rangle = |+\rangle \otimes |0\rangle$$

# Energy of a physical system

Determining energy levels of physical systems is relevant for applications of quantum computing like physics and chemistry.



The energies (and associated configurations) are given by the *eigenvalues* and *eigenvectors* of the system Hamiltonian.

# Computing the energy of a Hamiltonian

Given an arbitrary  $\hat{H}$ , how can we compute its energy?

Option: *exact diagonalization* to find its eigenvalues.

$$\hat{H} = \begin{pmatrix} * & * & * \\ * & * & * \\ * & * & * \end{pmatrix} \xrightarrow{\text{np.linalg.eig}} \{E_i, |E_i\rangle\}$$

Many other purely classical methods too. But we're more interested in the quantum ones.

# Time evolution

The Schrödinger equation from physics describes how the state of a system evolves with time.

The solution is:

$e^{-i\hat{H}t/\hbar}$  is the *time evolution operator*, and  $\hbar$  is a fundamental physical constant (we set it to 1 for simplicity).

Fun fact: if  $\hat{H}$  is Hermitian,

$$e^{-i\hat{H}t}$$

is unitary.

Then

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

describes how a system evolves with time.

If we can implement  $U$  on a quantum computer, we can compute its energy eigenvalues...

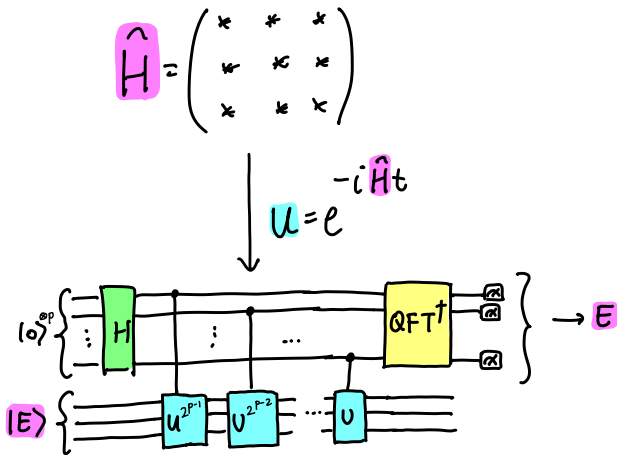


## Computing the energy of a Hamiltonian

Let  $\hat{H}$  be a Hamiltonian and  $|E\rangle$  an eigenstate with energy  $E$ .

# Computing the energy of a Hamiltonian

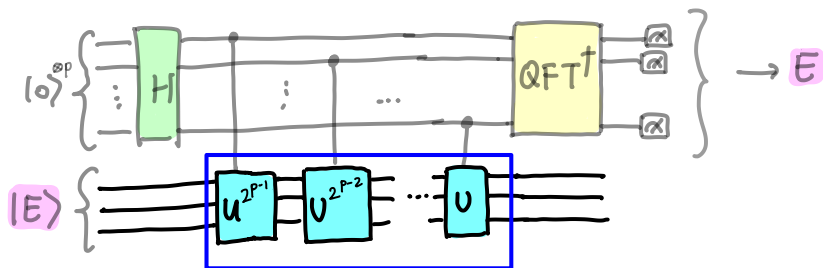
We can use *quantum phase estimation*!



(In hands-on 4, you will do this with the *variational eigensolver*)

# Computing the energy of a Hamiltonian

The next challenge: implementing  $U$



## Example: time evolution

Consider just one qubit for now.

$$\hat{H} = -\alpha Z_0$$

**Exercise:** A qubit starts in state  $|\psi(0)\rangle$  at time  $t = 0$ . What is  $|\psi(t)\rangle$  for this  $\hat{H}$ ? What quantum operation is this?

## Example: time evolution

Recall how we expressed  $RZ$  as a matrix exponential:

We have

## Example: time evolution

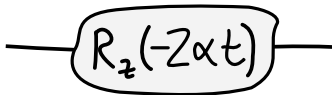
What about two non-interacting qubits?

$$\hat{H} = -\alpha Z_0 - \alpha Z_1$$

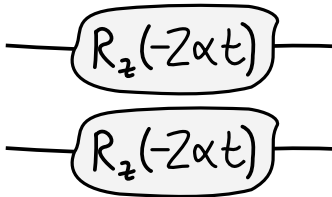
**Exercise:** What do you think the quantum circuit for time evolution under this  $\hat{H}$  looks like?

## Example: time evolution

One qubit:



Two qubits:



This makes Hamiltonian simulation looks pretty easy!

# Hamiltonian simulation in general

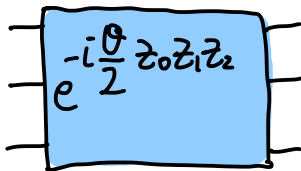
What happens when

1. we have Paulis that aren't just  $Z$ ?
2. we have interaction terms?
3. we have combinations of all Paulis and arbitrary interactions?



## Circuits for individual Pauli terms

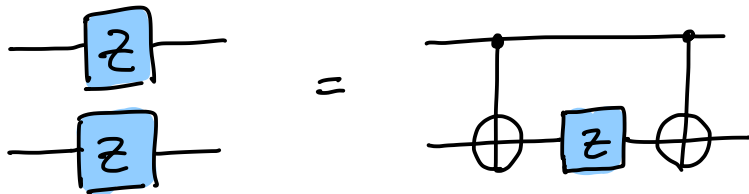
Example:


$$e^{-i\frac{\theta}{2} z_0 z_1 z_2}$$

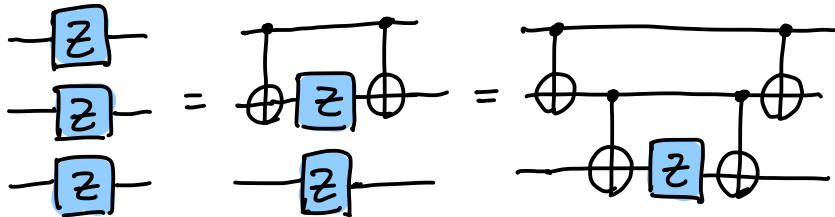
Idea: transform this into an exponential using only single-qubit Pauli Z, then implement those.

## Circuits for individual Pauli terms

There is a useful circuit identity that relates  $Z$  and CNOT:

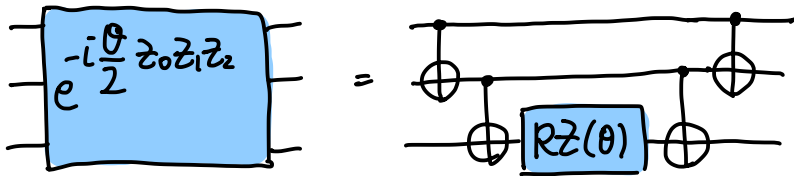


We can apply this multiple times



## Circuits for individual Pauli terms

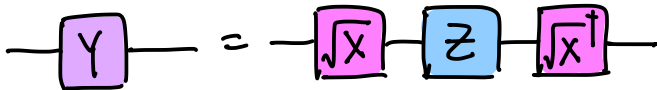
$Z$  is just a special case of  $RZ$  though; so this works in general:



We can deal with any product of Pauli  $Z$  in this way.

## Circuits for individual Pauli terms

$X$  and  $Y$  take a little more work; but they are related to  $Z$  in special ways:



## Circuits for individual Pauli terms

$RX$  has the same relationship:

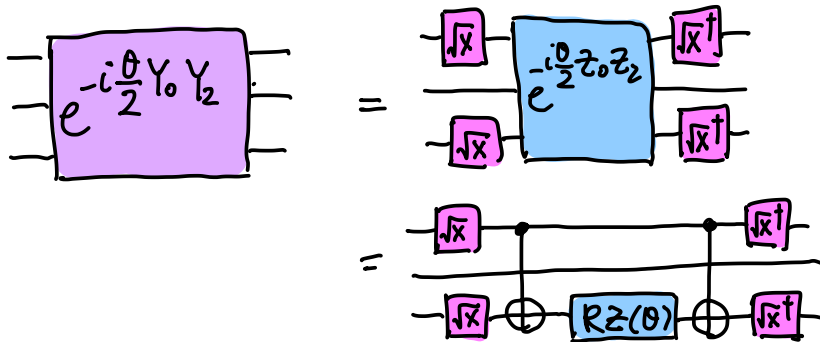
$$\text{---} \boxed{RX(\theta)} \text{---} = \text{---} \boxed{H} \boxed{RZ(\theta)} \boxed{H} \text{---}$$

So when we exponentiate...

$$\begin{aligned} \text{---} \boxed{e^{-i\frac{\theta}{2} X_0 X_1}} \text{---} &= \text{---} \boxed{H} \boxed{e^{-i\frac{\theta}{2} Z_0 Z_1}} \boxed{H} \text{---} \\ &= \text{---} \boxed{H} \text{---} \text{---} \boxed{H} \text{---} \\ &\quad \text{---} \boxed{H} \oplus \boxed{RZ(\theta)} \oplus \boxed{H} \text{---} \end{aligned}$$

## Circuits for individual Pauli terms

Similarly for  $Y$



We can do this for arbitrary Paulis with  $X$ ,  $Y$ , and  $Z$  terms.

Fun fact: the operators applied on either side of the  $Z$  are elements of the **Clifford group**. Cliffords send Paulis to Paulis.

## Dealing with sums of Pauli terms

Now for the hard part. What about:

$$e^{-i\alpha Z_0 X_1 - i\beta Z_0 Y_1}$$

More generally,

$$e^{-i\alpha P - i\beta Q}, \quad P, Q \in \mathcal{P}_n$$

Depends whether  $P, Q$  commute, i.e., if  $[P, Q] = PQ - QP = 0$ .

More next time!

# Next time

## Content:

- Pauli commutation relation
- Hamiltonian simulation with Trotterization
- Error in Hamiltonian simulation

## Action items:

1. Assignment 3 and hands-on 3
2. Work on project (remember to fill out weekly surveys)

## Recommended reading:

- Codebook module H