

# Two Basic Quantum Paradigms: Eigenvalue Estimation & Amplitude Amplification\*

\* But not necessarily in that order

**Richard Cleve**

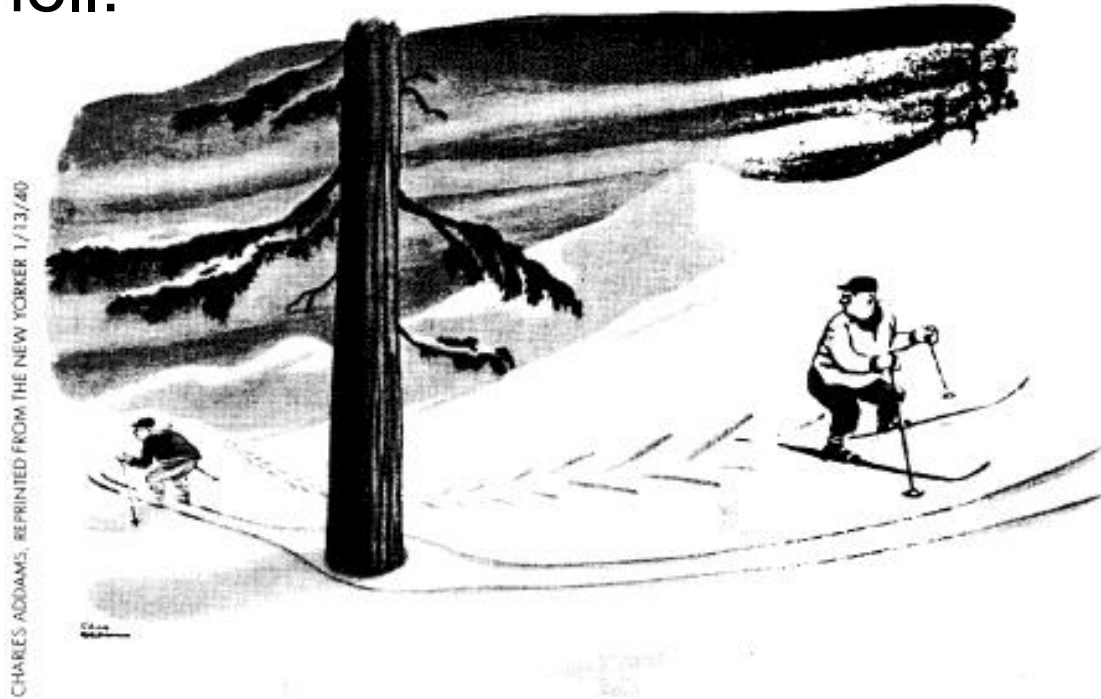
Institute for Quantum Computing & School of Computer Science  
University of Waterloo



Fields Institute, Toronto, October 27, 2014

# How do quantum algorithms work?

In a nutshell:



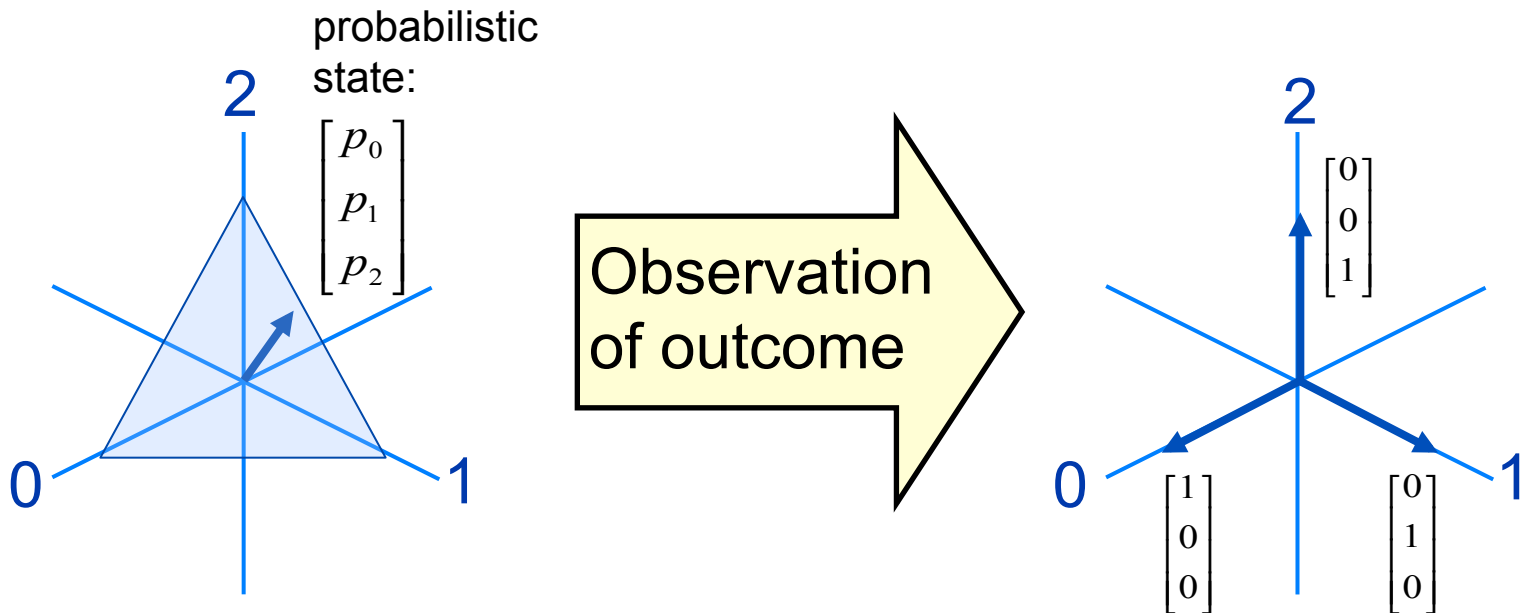
But this isn't "quantum parallelism" in the sense of quantum computers simply performing several computations at once (that approach doesn't work)

# Probability

**Probability vector:**

$$\forall x \in \Omega, p_x \geq 0 \quad \text{and} \quad \sum_x p_x = 1$$

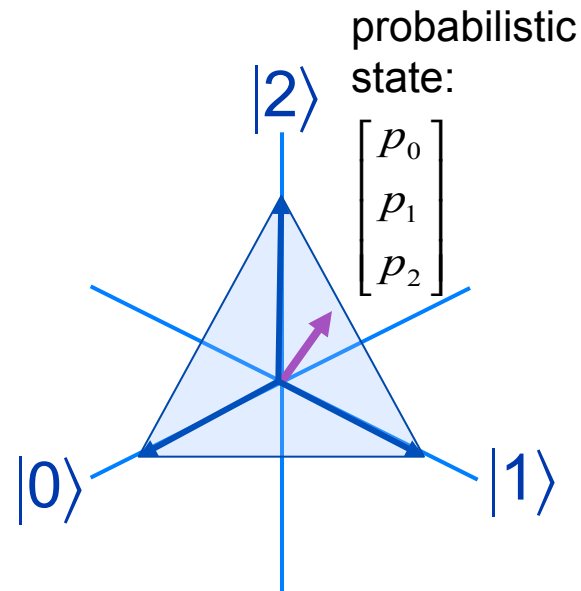
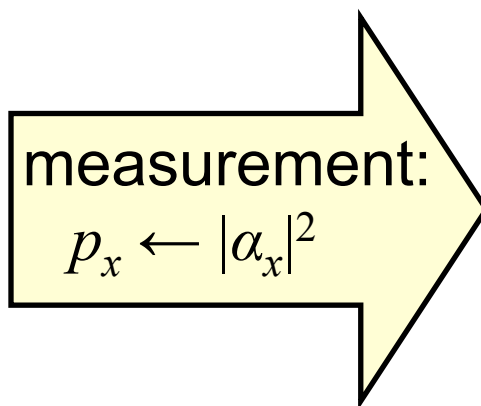
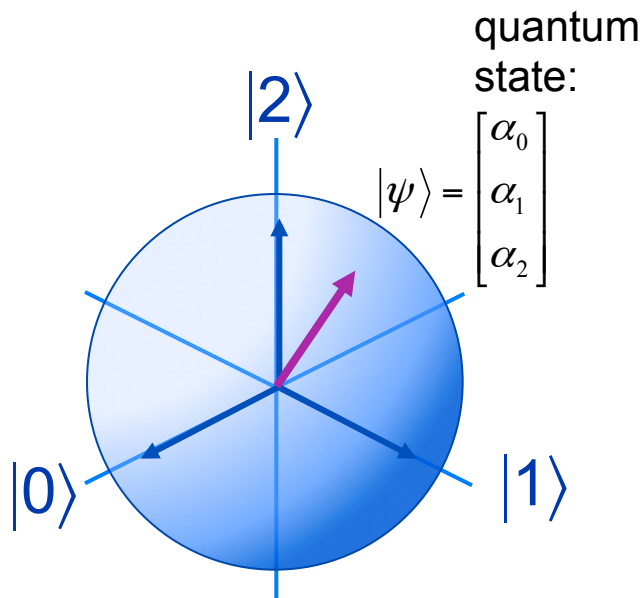
**Example:** for  $\Omega = \{0,1,2\}$



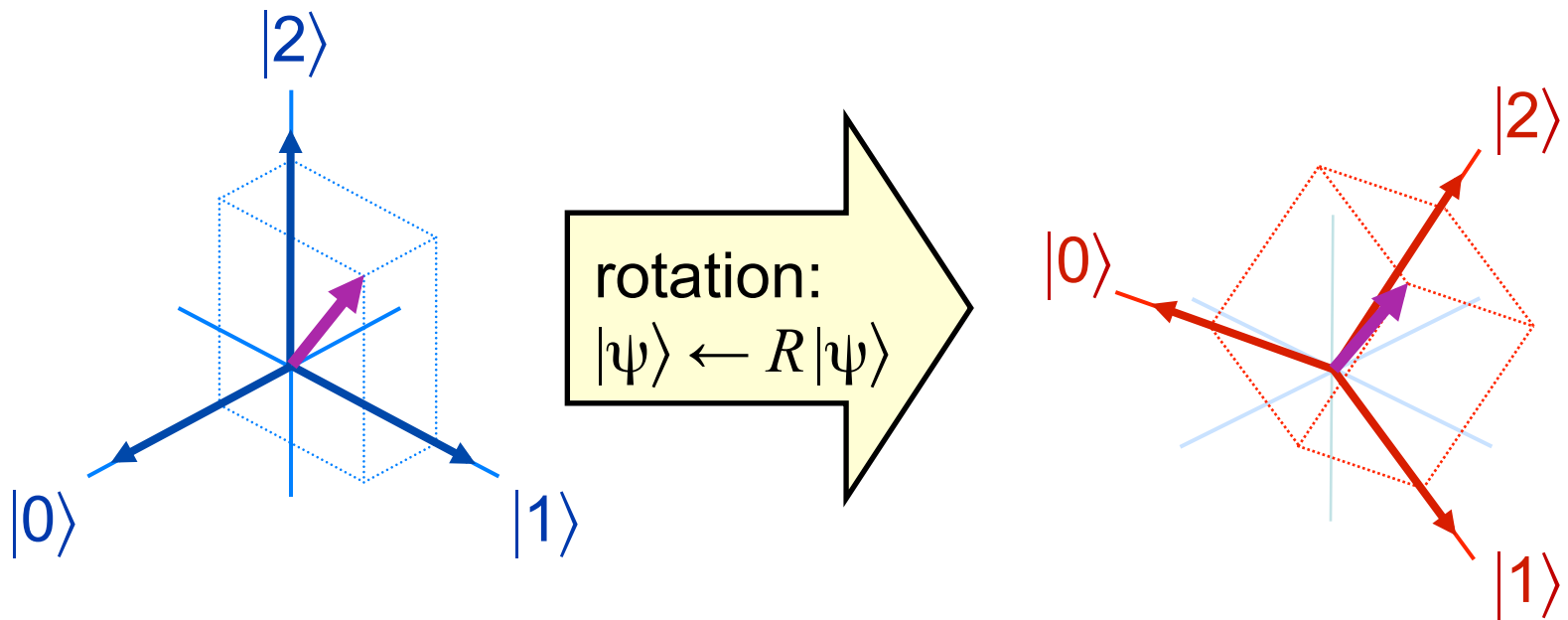
# Quantum information framework

**Probability *amplitudes*:**

$$\forall x \in \Omega, \alpha_x \in \mathbb{C} \text{ and } \sum_x |\alpha_x|^2 = 1$$



# Can rotate the coordinate system!



Different rotations may result in different probability distributions

# $n$ -qubit systems

Probabilistic states:

$$\forall x, p_x \geq 0$$

$$\sum_x p_x = 1$$

$$\begin{bmatrix} p_{000} \\ p_{001} \\ p_{010} \\ p_{011} \\ p_{100} \\ p_{101} \\ p_{110} \\ p_{111} \end{bmatrix}$$

Quantum states:

$$\forall x, \alpha_x \in \mathbb{C}$$

$$\sum_x |\alpha_x|^2 = 1$$

$$\begin{bmatrix} \alpha_{000} \\ \alpha_{001} \\ \alpha_{010} \\ \alpha_{011} \\ \alpha_{100} \\ \alpha_{101} \\ \alpha_{110} \\ \alpha_{111} \end{bmatrix}$$

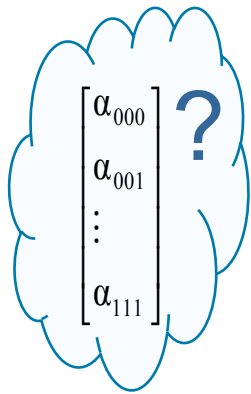
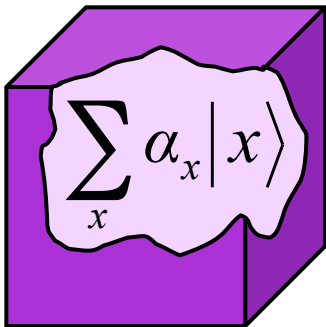
Dirac notation:  $|000\rangle, |001\rangle, |010\rangle, \dots, |111\rangle$  are basis vectors

$$|\psi\rangle = \sum_x \alpha_x |x\rangle$$

# Operations on $n$ -qubit states

**Unitary operations:**  $\sum_x \alpha_x |x\rangle \mapsto U\left(\sum_x \alpha_x |x\rangle\right)$   
( $U^\dagger U = I$ )

**Measurements:**



$$\left\{ \begin{array}{ll} 000 & \text{with prob } |\alpha_{000}|^2 \\ 001 & \text{with prob } |\alpha_{001}|^2 \\ \vdots & \vdots \\ 111 & \text{with prob } |\alpha_{111}|^2 \end{array} \right.$$

... and the quantum state collapses

# Basic operations on qubits (I)

(0) Initialize qubit to  $|0\rangle$  or to  $|1\rangle$

(1) Apply a unitary operation  $U$  (unitary means  $U^\dagger U = I$ )

↑  
conjugate transpose

## Examples:

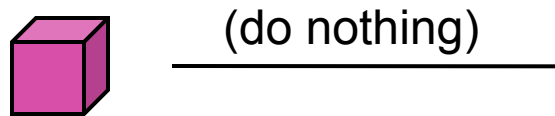
**Rotation:** 
$$\begin{bmatrix} \cos \theta & -\sin \theta \\ \sin \theta & \cos \theta \end{bmatrix}$$

**NOT (bit flip):**  $\sigma_x = X = \begin{bmatrix} 0 & 1 \\ 1 & 0 \end{bmatrix}$

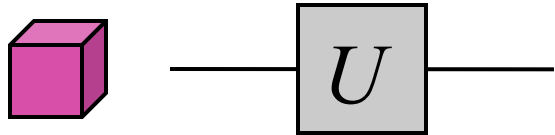
**Hadamard:**  $H = \frac{1}{\sqrt{2}} \begin{bmatrix} 1 & 1 \\ 1 & -1 \end{bmatrix}$

**Phase flip:**  $\sigma_z = Z = \begin{bmatrix} 1 & 0 \\ 0 & -1 \end{bmatrix}$

# Example of a one-qubit gate applied to a two-qubit system



$$U = \begin{bmatrix} u_{00} & u_{01} \\ u_{10} & u_{11} \end{bmatrix}$$



The resulting 4x4 matrix is

Maps basis states as:

$$|0\rangle|0\rangle \rightarrow |0\rangle U|0\rangle$$

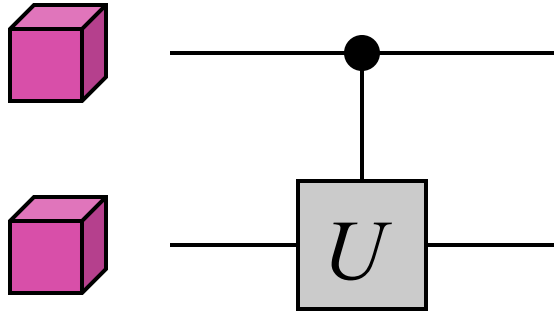
$$|0\rangle|1\rangle \rightarrow |0\rangle U|1\rangle$$

$$|1\rangle|0\rangle \rightarrow |1\rangle U|0\rangle$$

$$|1\rangle|1\rangle \rightarrow |1\rangle U|1\rangle$$

$$I \otimes U = \begin{bmatrix} u_{00} & u_{01} & 0 & 0 \\ u_{10} & u_{11} & 0 & 0 \\ 0 & 0 & u_{00} & u_{01} \\ 0 & 0 & u_{10} & u_{11} \end{bmatrix}$$

# Controlled- $U$ gates



$$U = \begin{bmatrix} u_{00} & u_{01} \\ u_{10} & u_{11} \end{bmatrix}$$

Resulting 4x4 matrix is  
controlled- $U =$

Maps basis states as:

$$|0\rangle|0\rangle \rightarrow |0\rangle|0\rangle$$

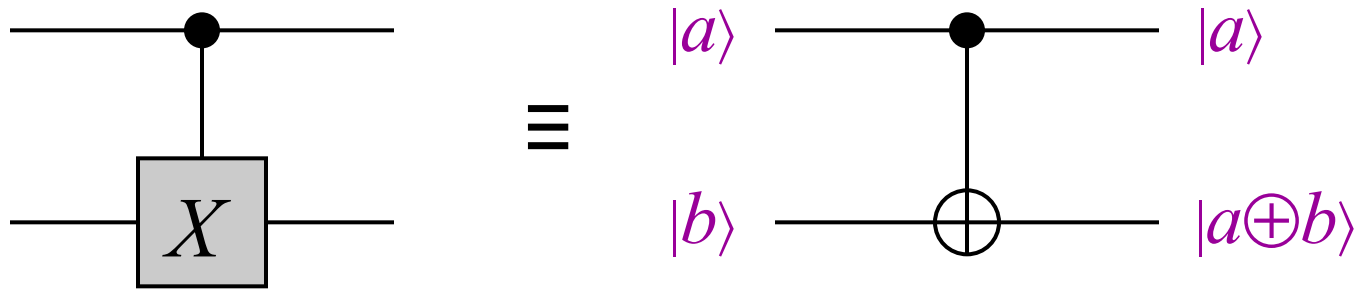
$$|0\rangle|1\rangle \rightarrow |0\rangle|1\rangle$$

$$|1\rangle|0\rangle \rightarrow |1\rangle U|0\rangle$$

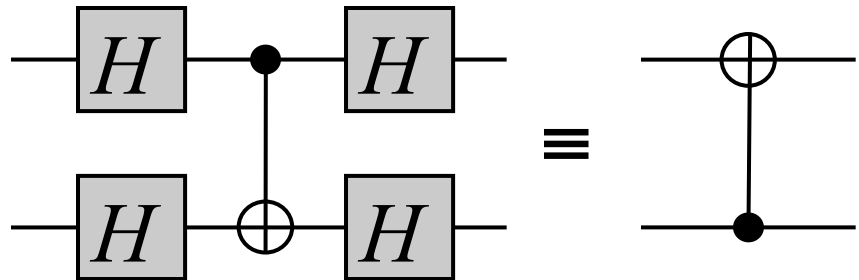
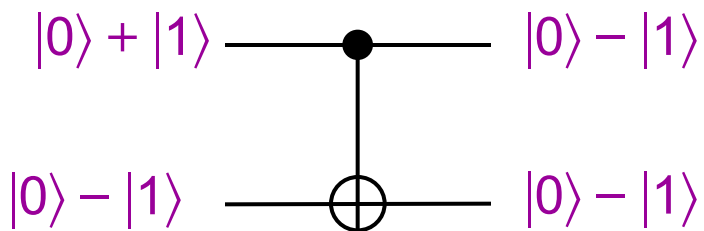
$$|1\rangle|1\rangle \rightarrow |1\rangle U|1\rangle$$

$$\begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & u_{00} & u_{01} \\ 0 & 0 & u_{10} & u_{11} \end{bmatrix}$$

# Controlled-NOT (CNOT)



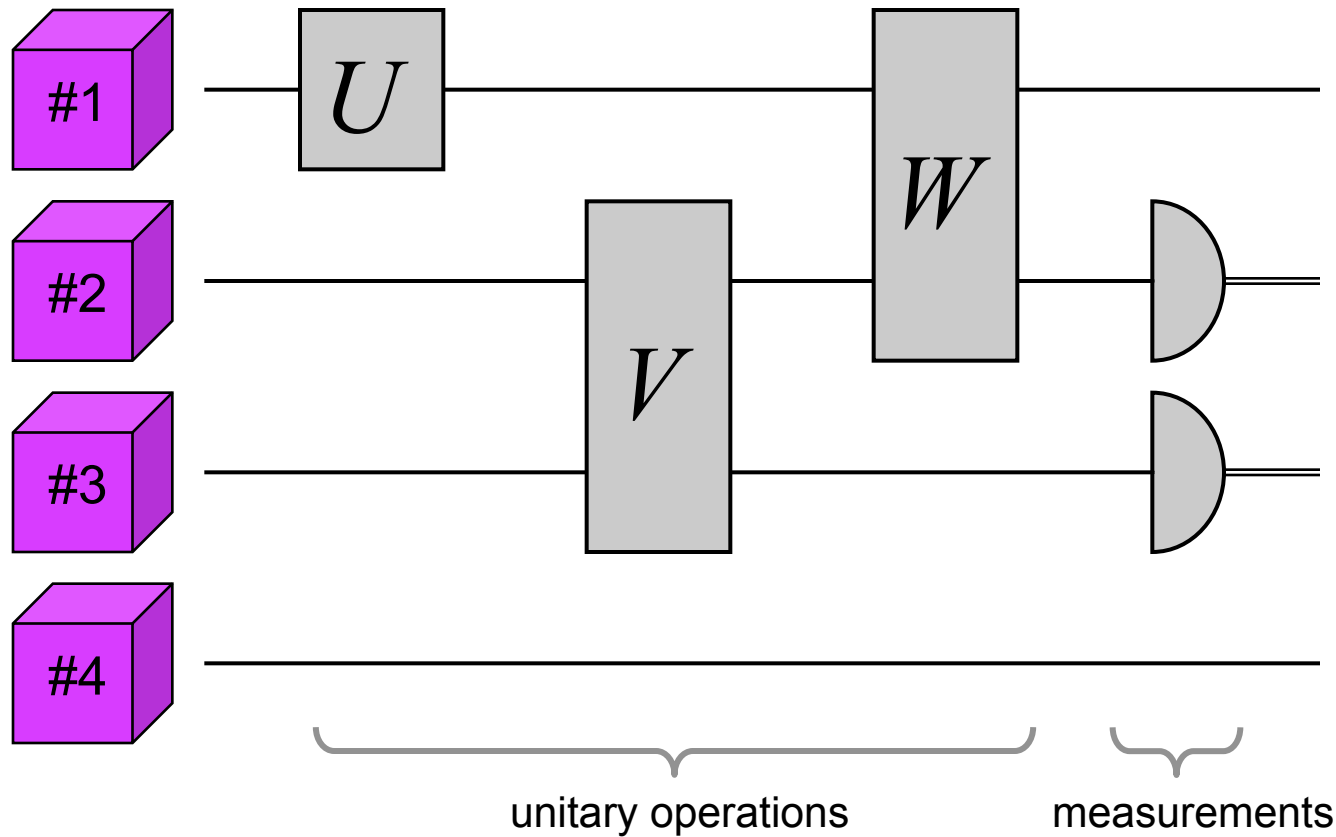
**Note:** “control” qubit may change on some input states!



# Structure among subsystems

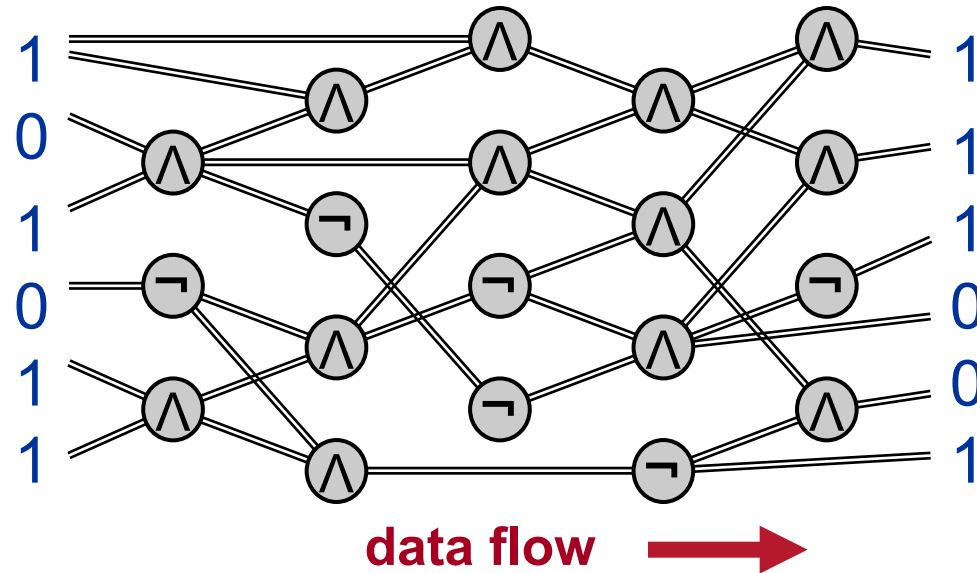
qubits:

time  $\longrightarrow$

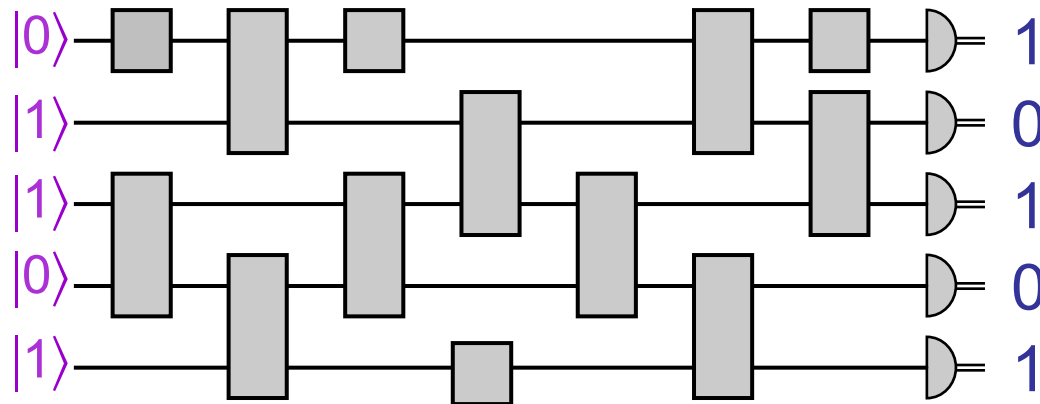


# Models of computation

**Classical circuits:**



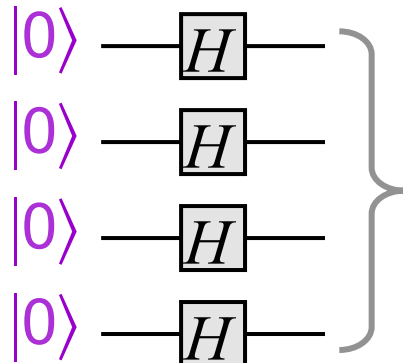
**Quantum circuits:**



Circuit families of size  $O(n^2 \text{polylog } n)$  exist for prime factorization, whereas the best known classical circuit family size is  $\approx 2^{n^{1/3}}$

# Equally weighted superpositions

**Hadamard:**  $H = \frac{1}{\sqrt{2}} \begin{bmatrix} 1 & 1 \\ 1 & -1 \end{bmatrix}$        $H|0\rangle = \frac{1}{\sqrt{2}}|0\rangle + \frac{1}{\sqrt{2}}|1\rangle$



$H \otimes H \otimes \dots \otimes H |00 \dots 0\rangle$   
 $= \left(\frac{1}{\sqrt{2}}|0\rangle + \frac{1}{\sqrt{2}}|1\rangle\right) \otimes \left(\frac{1}{\sqrt{2}}|0\rangle + \frac{1}{\sqrt{2}}|1\rangle\right) \otimes \dots \otimes \left(\frac{1}{\sqrt{2}}|0\rangle + \frac{1}{\sqrt{2}}|1\rangle\right)$   
 $= \frac{1}{\sqrt{2^n}} \sum_{x \in \{0,1\}^n} |x\rangle$

# How do quantum algorithms work?

Given a polynomial-time classical algorithm for  $f : \{0,1\}^n \rightarrow T$ , it is straightforward to construct a quantum algorithm that creates the state

$$\frac{1}{\sqrt{2^n}} \sum_x |x\rangle |f(x)\rangle$$

at the cost of about **one** evaluation of  $f$

Is this exponentially many computations at polynomial cost?

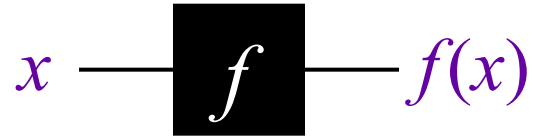
No! — the most straightforward way of extracting information from the state yields just  $(x, f(x))$  for a random  $x \in \{0,1\}^n$

But we can make some interesting **tradeoffs**:

instead of learning about any  $(x, f(x))$  point, one can learn something about a **global property** of  $f$

# Query scenario

**Input:** a function  $f$ , given as a black box (a.k.a. oracle)



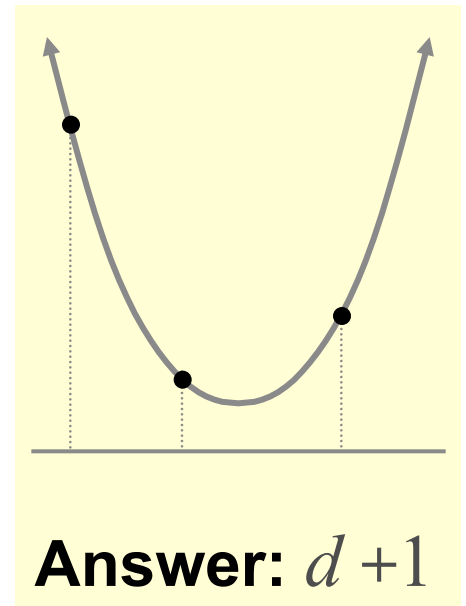
**Goal:** determine some information about  $f$  making as few queries to  $f$  (and other operations) as possible

**Example:** polynomial interpolation

**Let:**  $f(x) = c_0 + c_1x + c_2x^2 + \dots + c_dx^d$

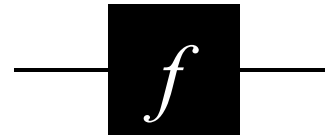
**Goal:** determine  $c_0, c_1, c_2, \dots, c_d$

**Question:** How many  $f$ -queries does one require for this?



# Example 1: Deutsch's problem

Let  $f: \{0,1\} \rightarrow \{0,1\}$



There are **four** possibilities:

$x$	$f_1(x)$
0	0
1	0

$x$	$f_2(x)$
0	1
1	1

$x$	$f_3(x)$
0	0
1	1

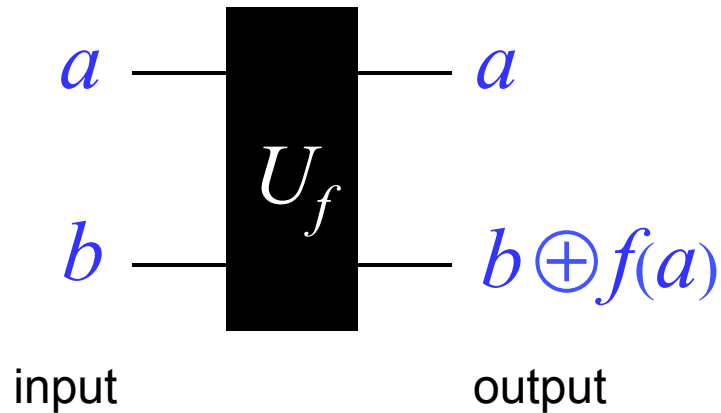
$x$	$f_4(x)$
0	1
1	0

**Goal:** determine whether or not  $f(0) = f(1)$  (i.e.  $f(0) \oplus f(1)$ )

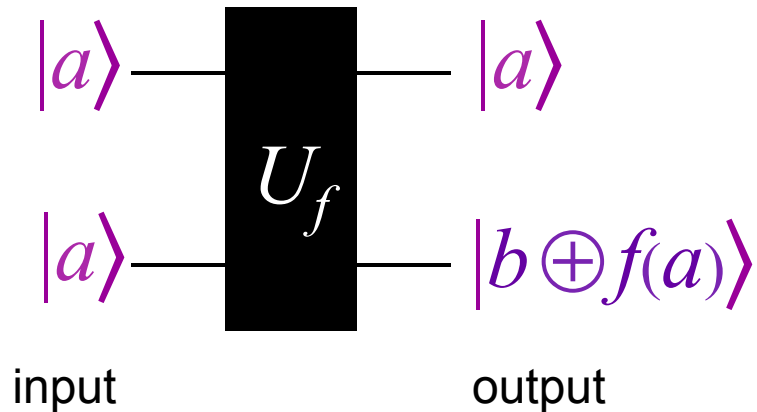
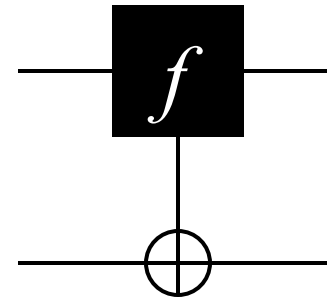
Any classical method must make **two** queries

There is a quantum algorithm that makes only **one** query

# Reversible black box for $f$



alternate  
notation:



# Example 2: search problem

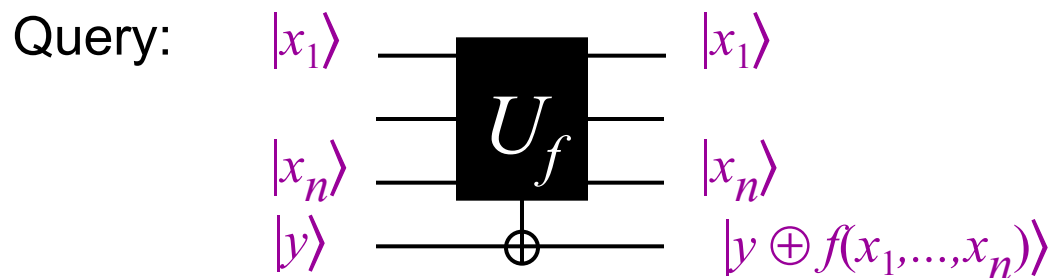
**Given:** a black box computing  $f: \{0,1\}^n \rightarrow \{0,1\}$

**Goal:** determine if  $f$  is **satisfiable** (if  $\exists x \in \{0,1\}^n$  s.t.  $f(x) = 1$ )

In positive instances, it makes sense to also **find** such a satisfying assignment  $x$

**Classically**, using probabilistic procedures, order  $2^n$  queries are necessary to succeed—even with probability  $\frac{3}{4}$  (say)

Grover's **quantum** algorithm that makes only  $O(\sqrt{2^n})$  queries

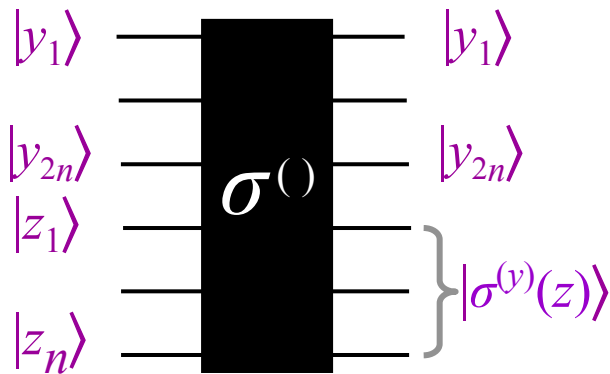


[Grover '96]

# Example 3: period finding

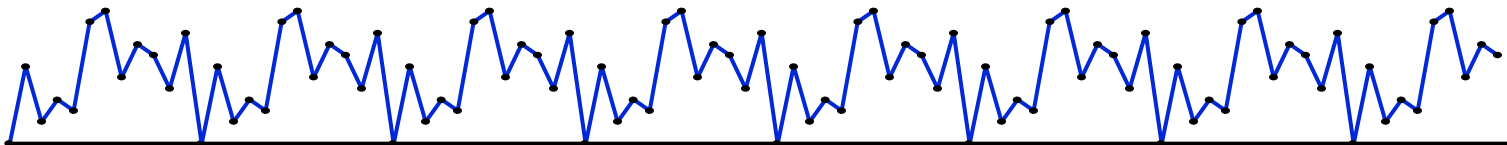
Let  $\sigma : \{0,1\}^n \rightarrow \{0,1\}^n$  be a(n unknown) permutation, and let  $\sigma^{(y)}(z)$  denote  $y$  iterations of  $\sigma$ , namely  $\sigma(\sigma(\sigma(\dots \sigma(z) \dots)))$

**Given:** a black box computing  $\sigma^{(y)}(z)$ , where  $y \in \{0,1\}^{2n}$  and  $z \in \{0,1\}^n$



(enables fast forwarding of large iterations of  $\sigma$ )

**Goal:** determine length of cycle containing 0 (say)



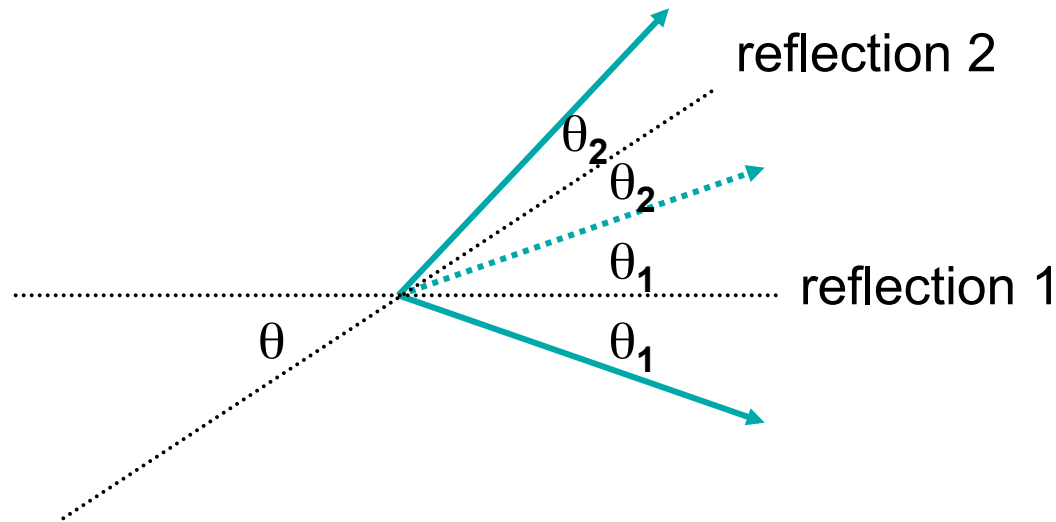
$$f(y) = \sigma^{(y)}(00\dots 0)$$

**Query complexity:** classically:  $\Omega(2^{n/3})$  queries; quantumly:  $O(1)$

# Prelude to Grover's search algorithm:

## two reflections = a rotation

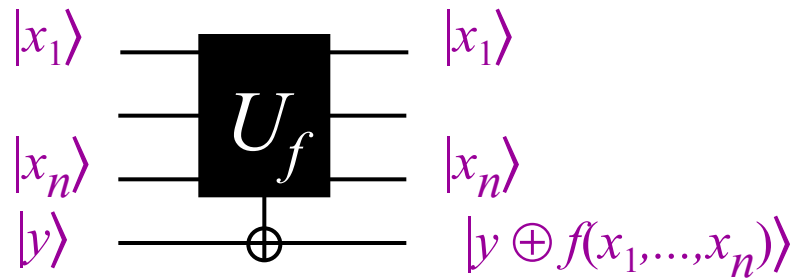
Consider two lines with intersection angle  $\theta$ :



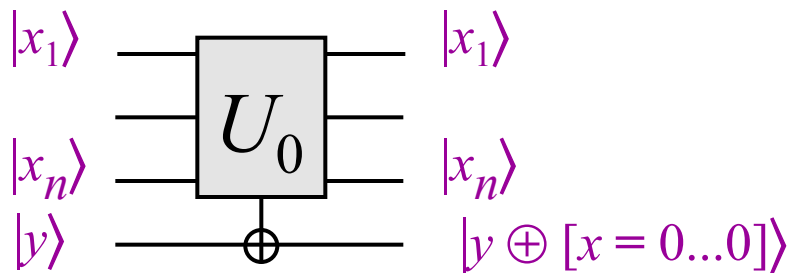
Net effect: rotation by angle  $2\theta$ , *regardless of starting vector*

# Grover's algorithm: description I

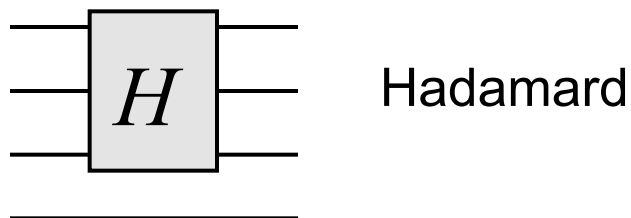
Basic operations used:



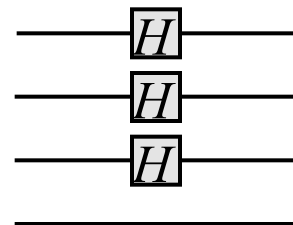
$$U_f |x\rangle(|0\rangle - |1\rangle) = (-1)^{f(x)} |x\rangle(|0\rangle - |1\rangle)$$



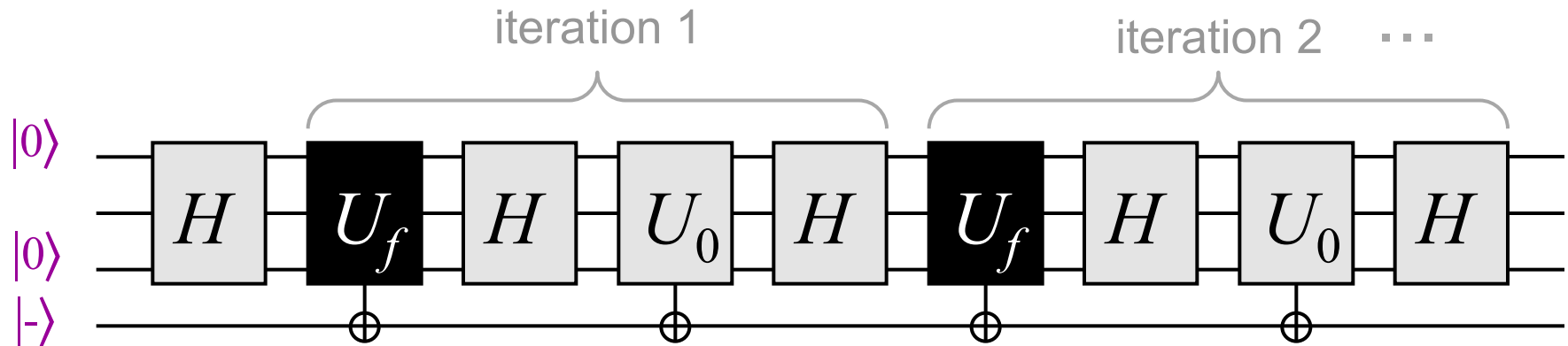
$$U_0 |x\rangle(|0\rangle - |1\rangle) = (-1)^{[x = 0\dots 0]} |x\rangle(|0\rangle - |1\rangle)$$



Hadamard



# Grover's algorithm: description II



1. construct state  $H|0\dots 0\rangle(|0\rangle-|1\rangle)$
  2. repeat  $k$  times:  
     apply  $-HU_0HU_f$  to state
  3. measure state, to get  $x \in \{0,1\}^n$ , and check if  $f(x)=1$
- (The setting of  $k$  will be determined later)

# Grover's algorithm: analysis I

Let  $A = \{x \in \{0,1\}^n : f(x) = 1\}$  and  $B = \{x \in \{0,1\}^n : f(x) = 0\}$

and  $N = 2^n$  and  $a = |A|$  and  $b = |B|$

Let  $|A\rangle = \frac{1}{\sqrt{a}} \sum_{x \in A} |x\rangle$  and  $|B\rangle = \frac{1}{\sqrt{b}} \sum_{x \in B} |x\rangle$

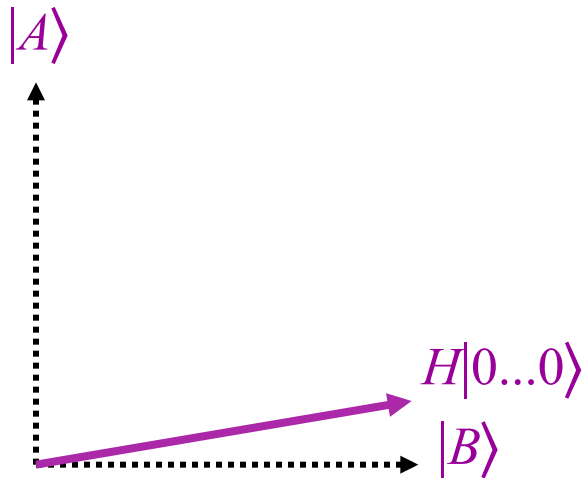
Consider the space spanned by  $|A\rangle$  and  $|B\rangle$

$|A\rangle$  ← goal is to get close to this state

$$H|0\dots 0\rangle = \frac{1}{\sqrt{N}} \sum_{x \in \{0,1\}^n} |x\rangle = \sqrt{\frac{a}{N}} |A\rangle + \sqrt{\frac{b}{N}} |B\rangle$$

Interesting case:  $a \ll N$

# Grover's algorithm: analysis II



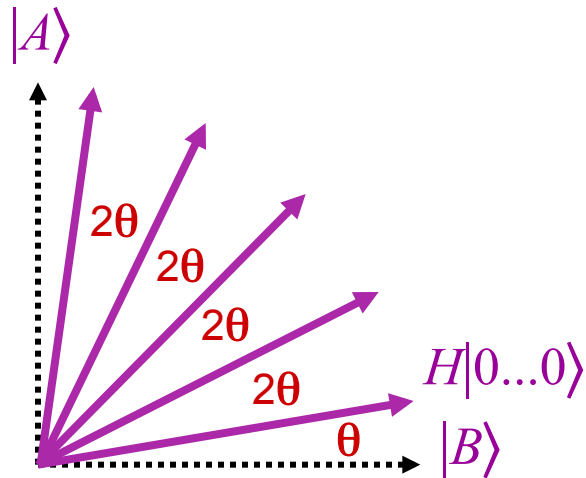
$$\text{Algorithm: } (-HU_0HU_f)^k H|0\dots 0\rangle$$

## Observations:

$U_f$  is a reflection about  $|B\rangle$ :  $U_f|A\rangle = -|A\rangle$  and  $U_f|B\rangle = |B\rangle$

$-HU_0H$  is a reflection about  $H|0\dots 0\rangle$

# Grover's algorithm: analysis III



Algorithm:  $(-HU_0HU_f)^k H|0\dots 0\rangle$

Since  $-HU_0HU_f$  is a composition of two reflections, it is a rotation by  $2\theta$ , where  $\sin(\theta) = \sqrt{a/N} \approx \sqrt{a/N}$

When  $a = 1$ , we want  $(2k+1)(1/\sqrt{N}) \approx \pi/2$ , so  $k \approx (\pi/4)\sqrt{N}$

More generally, it suffices to set  $k \approx (\pi/4)\sqrt{N/a}$

(Can be adapted to work even if  $a$  is not known in advance)

# Applications of quantum search

The function  $f$  could be realized as a **3-CNF formula**:

$$f(x_1, \dots, x_n) = (x_1 \vee \bar{x}_3 \vee x_4) \wedge (\bar{x}_2 \vee x_3 \vee \bar{x}_5) \wedge \dots \wedge (\bar{x}_1 \vee x_5 \vee \bar{x}_n)$$

Running time:  $O(\sqrt{2^n} \text{ polylog } n)$  gates

But there are better **classical** SAT solvers, such as [Schöning '99]:

$O(\sqrt{2^n} \text{ polylog } n)$  gates

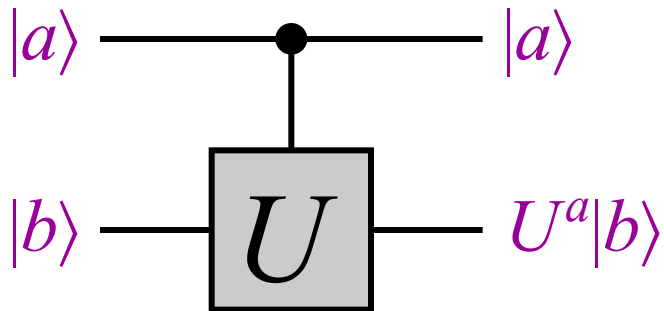
We can recast Schöning's algorithm as a search on a space of size

$O(\sqrt{2^n} \text{ polylog } n)$  and then use Grover's algorithm to attain

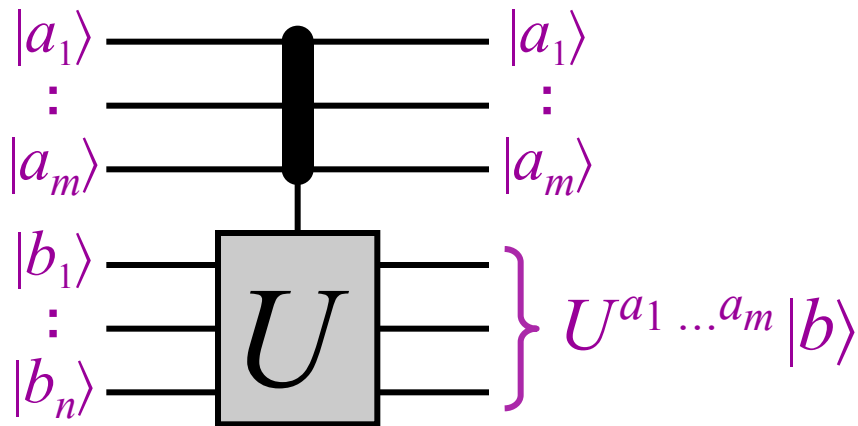
algorithm with  $O((2^n)^{1/4} \text{ polylog } n)$  gates

That's essentially what is called **amplitude amplification**

# Generalized controlled- $U$ gates



$$\begin{bmatrix} I & 0 \\ 0 & U \end{bmatrix}$$



$$\begin{bmatrix} I & 0 & 0 & \dots & 0 \\ 0 & U & 0 & \dots & 0 \\ 0 & 0 & U^2 & \dots & 0 \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & 0 & \dots & U^{2^m-1} \end{bmatrix}$$

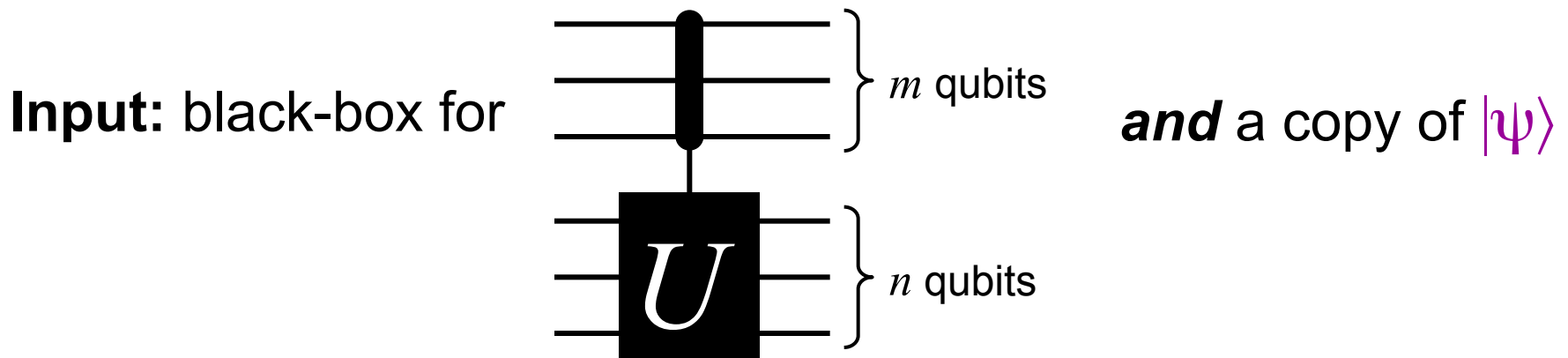
**Example:**  $|1101\rangle|0101\rangle \rightarrow |1101\rangle U^{1101}|0101\rangle$

# Eigenvalue estimation problem

$U$  is a unitary operation on  $n$  qubits

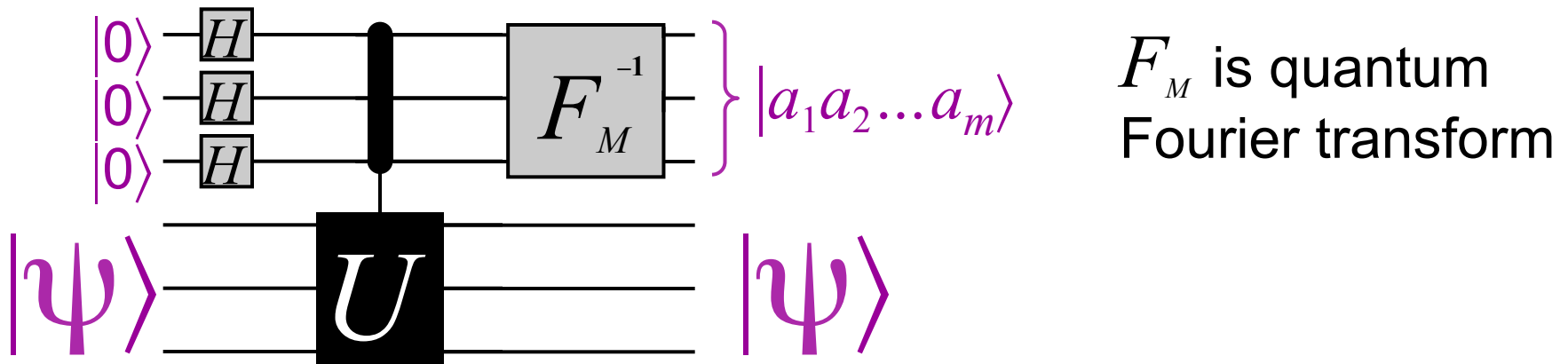
$|\psi\rangle$  is an eigenvector of  $U$ , with eigenvalue  $e^{2\pi i\phi}$

$(0 \leq \phi < 1)$



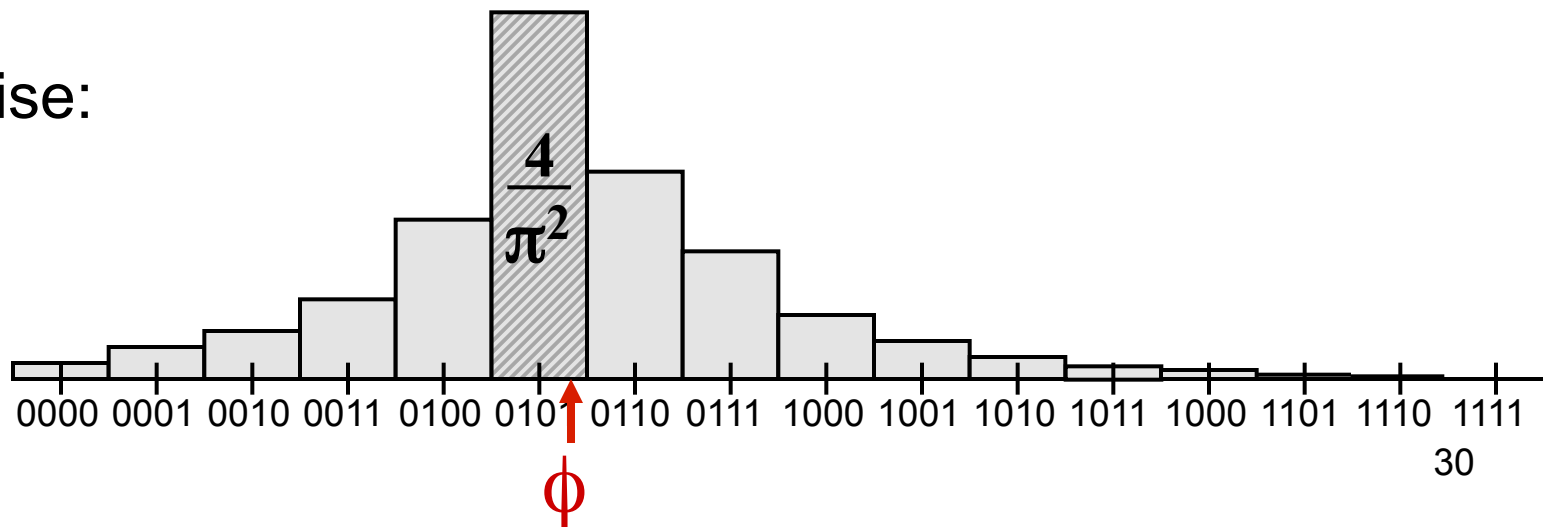
**Output:**  $\phi$  ( $m$ -bit approximation)

# Algorithm for eigenvalue estimation



If  $\phi = 0.a_1a_2\dots a_m$  then the above procedure yields  $|a_1a_2\dots a_m\rangle$  (from which  $\phi$  can be deduced exactly)

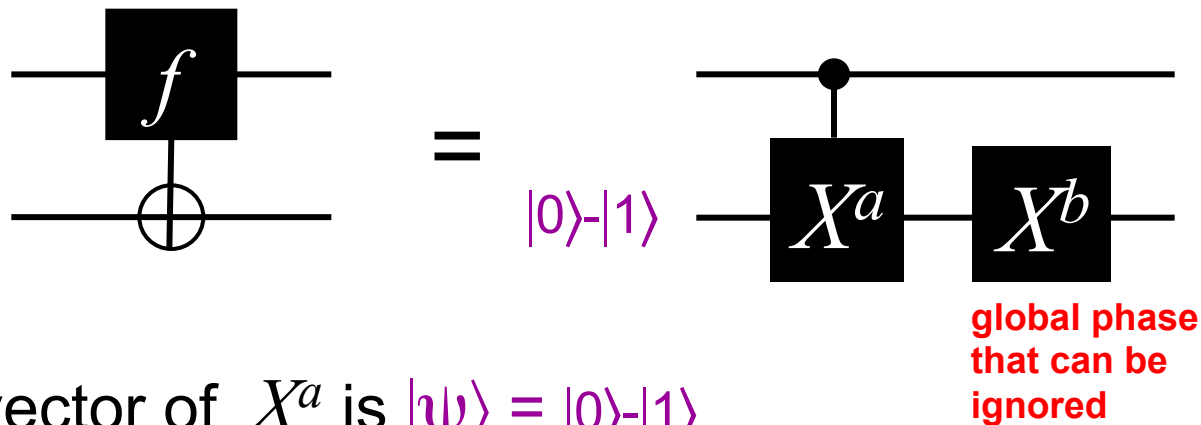
Otherwise:



# Simple example: Deutsch's problem

Recall  $f: \{0,1\} \rightarrow \{0,1\}$       $f(z) = a \cdot z + b \pmod 2$

Note that  $f$  is constant iff  $a = 0$



Eigenvector of  $X^a$  is  $|\psi\rangle = |0\rangle - |1\rangle$

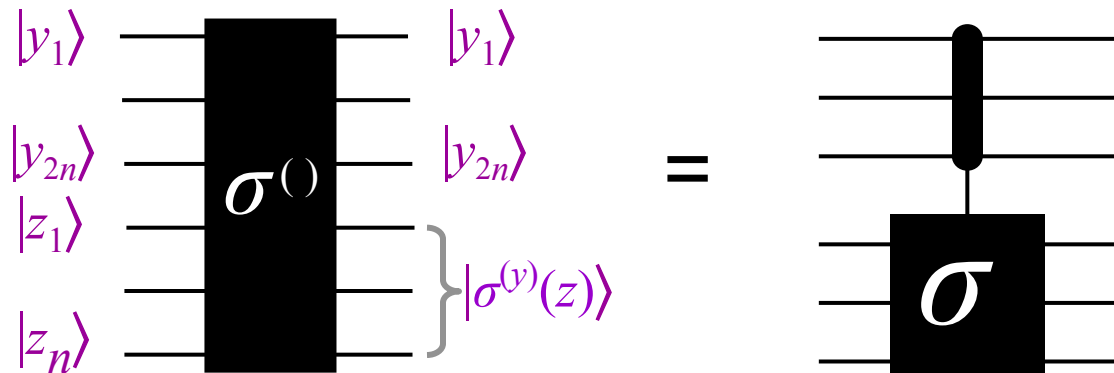
Eigenvalue of  $X^a$  is: +1 if constant; -1 if balanced

Eigenvalue can be determined with a single query to  $f$

# Another example: period finding

Recall that  $\sigma : \{0,1\}^n \rightarrow \{0,1\}^n$  is an unknown permutation, and  $\sigma^{(y)}(z)$  denotes  $y$  iterations of  $\sigma$ , namely  $\sigma(\sigma(\sigma(\dots \sigma(z) \dots)))$

Eigenvectors of  $\sigma$ :  $|\psi_1\rangle = \frac{1}{\sqrt{r}} \sum_{j=0}^{r-1} |\sigma^{(j)}(0)\rangle$      $|\psi_k\rangle = \frac{1}{\sqrt{r}} \sum_{j=0}^{r-1} e^{-i2k\pi j/r} |\sigma^{(j)}(0)\rangle$



Eigenvalue is  $k/r$  from which  $r$  can be deduced by continued fractions

Trick for bypassing construction of  $|\psi_k\rangle$  by using  $|0\rangle$

**Order-finding** (min  $r$  such that  $a^r \bmod m = 1$ ) reduces to **period finding**  
(fast forward  $\sigma(z) = az \bmod m$  by repeated squaring)

# Thank you

