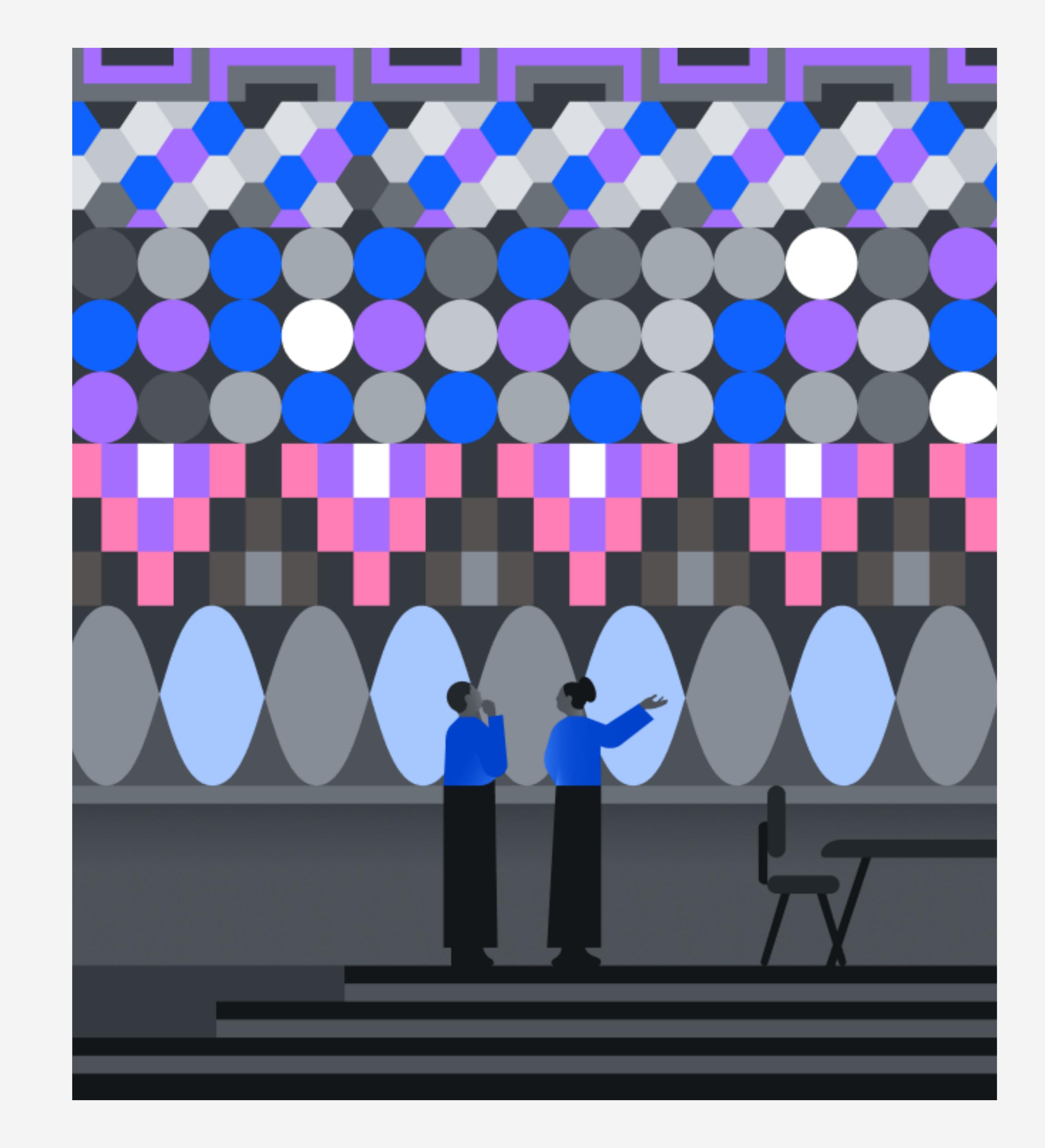
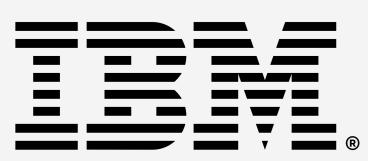
Understanding quantum information and computation

By John Watrous

Lesson 7

Phase estimation and factoring





Spectral theorem for unitary matrices

The *spectral theorem* is an important fact in linear algebra. Here is a statement of a special case of this theorem, for *unitary matrices*.

Spectral theorem for unitary matrices

Suppose U is an N x N unitary matrix.

There exists an orthonormal basis $\{|\psi_1\rangle,\ldots,|\psi_N\rangle\}$ of vectors along with complex numbers

$$\lambda_1 = e^{2\pi i\theta_1}, \dots, \lambda_N = e^{2\pi i\theta_N}$$

such that

$$U = \sum_{k=1}^{N} \lambda_k |\psi_k\rangle\langle\psi_k|$$

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such that

$$U = \sum_{k=1}^{N} \lambda_k |\psi_k\rangle\langle\psi_k|$$

Each vector $|\psi_k\rangle$ is an *eigenvector* of U having *eigenvalue* λ_k :

$$U|\psi_k\rangle = \lambda_k|\psi_k\rangle = e^{2\pi i\theta_k}|\psi_k\rangle$$

Phase estimation problem

In the phase estimation problem, we're given two things:

- 1. A description of a *unitary quantum circuit* on n qubits.
- 2. An n-qubit $\frac{quantum state}{\psi}$.

We're <u>promised</u> that $|\psi\rangle$ is an eigenvector of the unitary operation U described by the circuit, and our goal is to approximate the corresponding eigenvalue.

Phase estimation problem

Input: A unitary quantum circuit for an n-qubit operation U

and an n qubit quantum state $|\psi\rangle$

Promise: $|\psi\rangle$ is an eigenvector of U

Output: An approximation to the number $\theta \in [0, 1)$ satisfying

$$U|\psi\rangle = e^{2\pi i\theta}|\psi\rangle$$

Phase estimation problem

Phase estimation problem

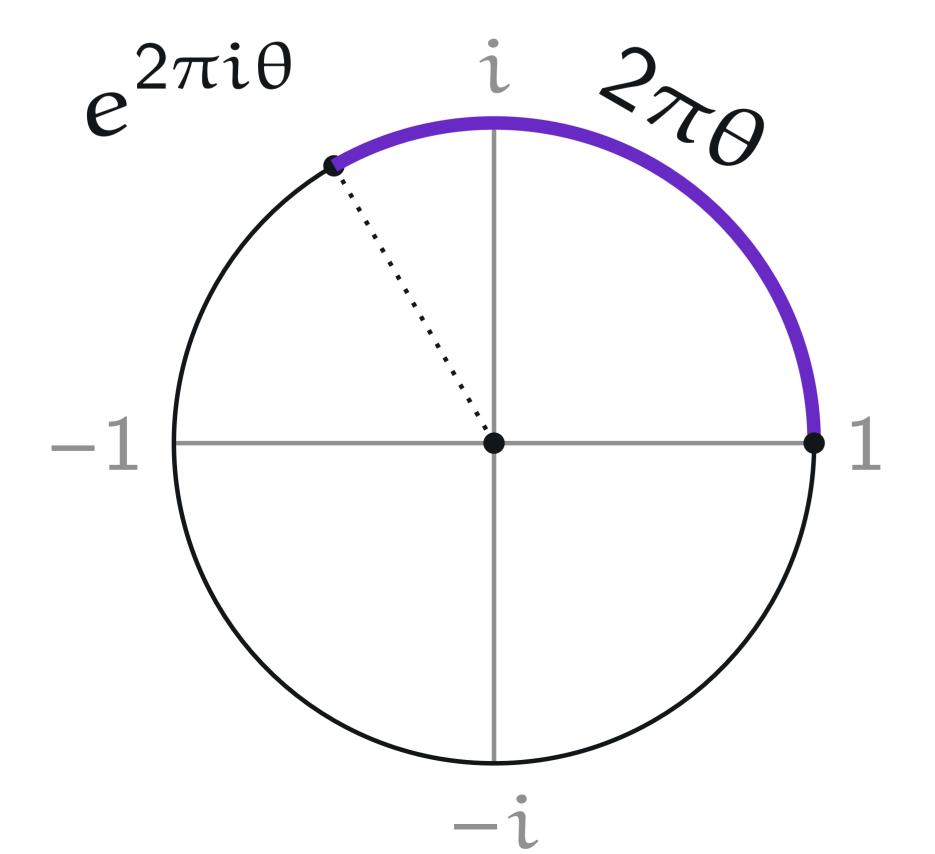
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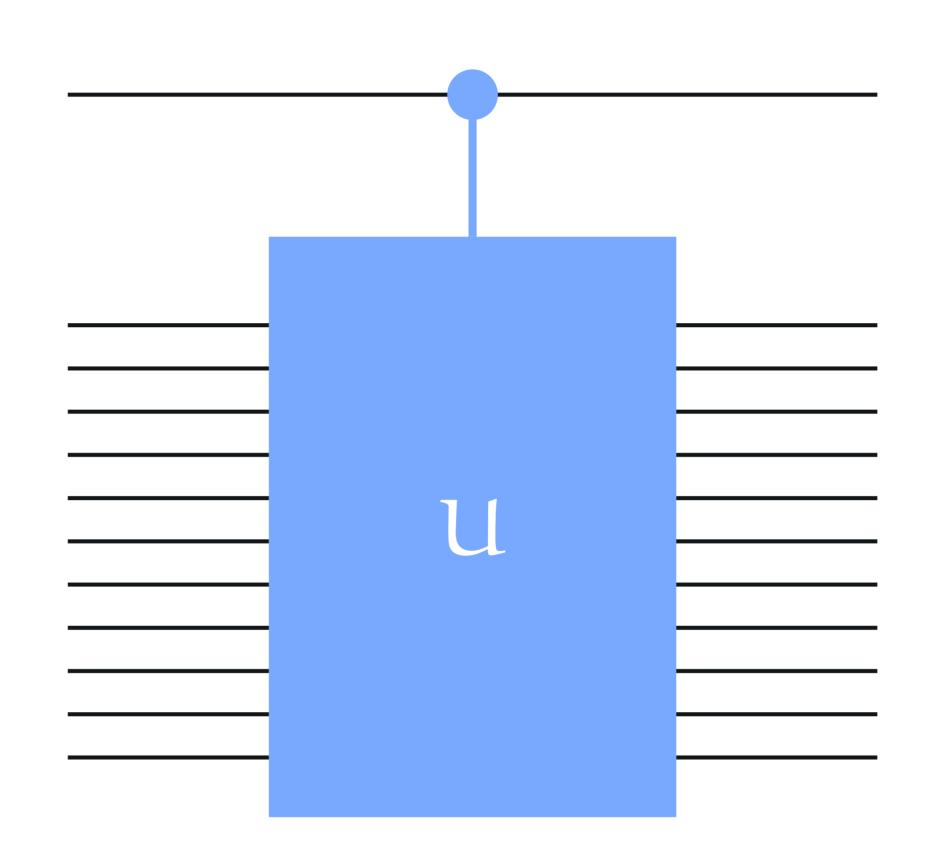
We can approximate θ by a fraction

$$\theta \approx \frac{y}{2^{m}}$$

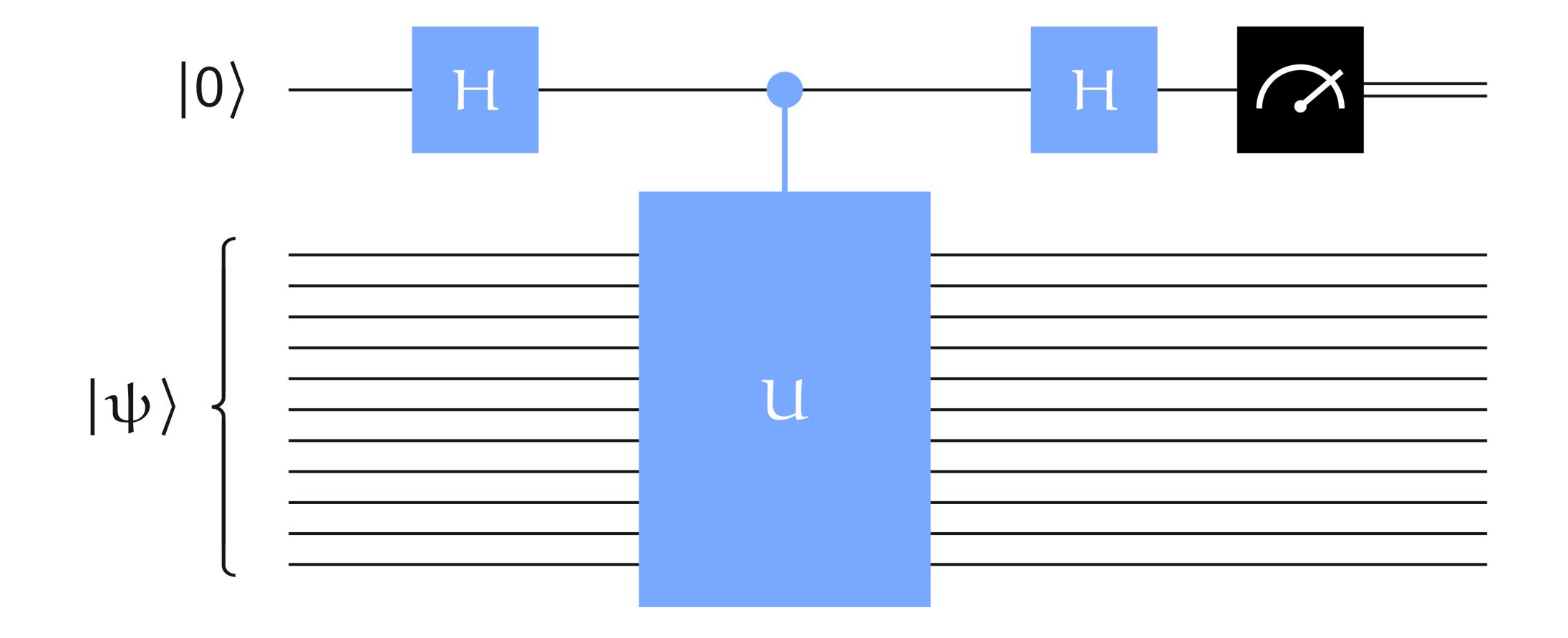
for
$$y \in \{0, 1, ..., 2^m - 1\}$$
.

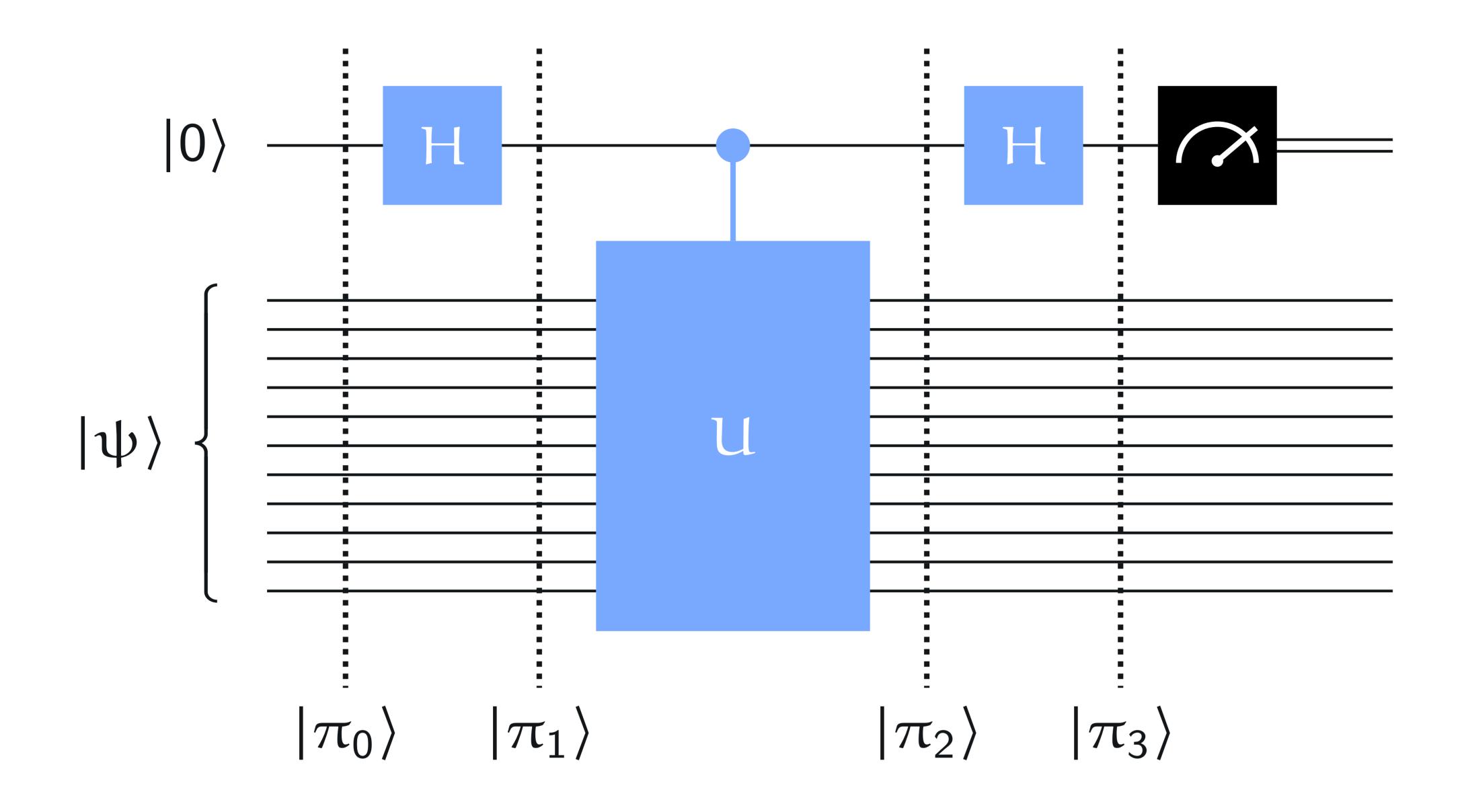
This approximation is taken "modulo 1."

Given a circuit for U, we can create a circuit for a controlled-U operation:

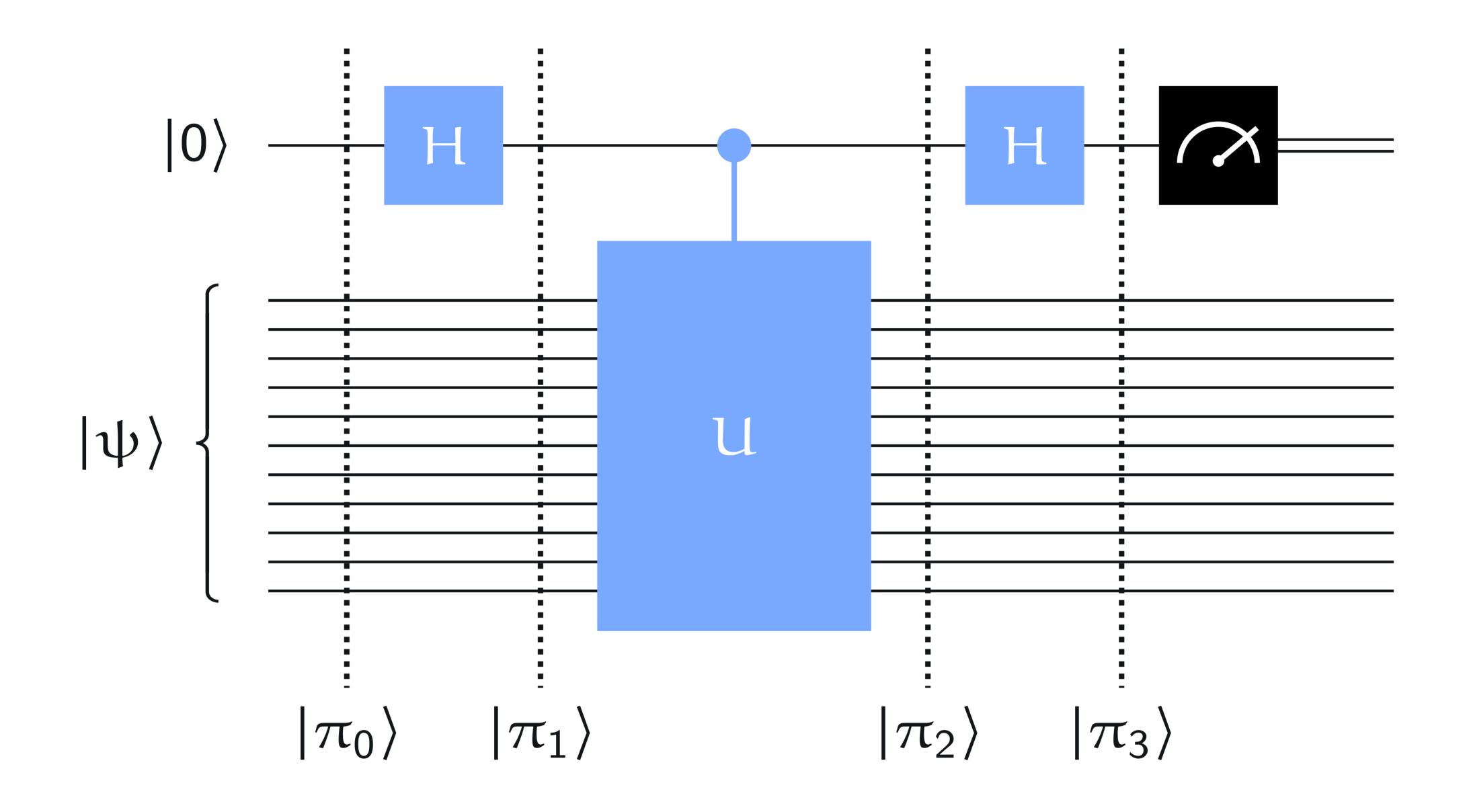


Let's consider this circuit:





$$\begin{split} |\pi_{0}\rangle &= |\psi\rangle|0\rangle \\ |\pi_{1}\rangle &= \frac{1}{\sqrt{2}}|\psi\rangle|0\rangle + \frac{1}{\sqrt{2}}|\psi\rangle|1\rangle \\ |\pi_{2}\rangle &= \frac{1}{\sqrt{2}}|\psi\rangle|0\rangle + \frac{1}{\sqrt{2}}(U|\psi\rangle)|1\rangle = |\psi\rangle \otimes \left(\frac{1}{\sqrt{2}}|0\rangle + \frac{e^{2\pi i\theta}}{\sqrt{2}}|1\rangle\right) \end{split}$$



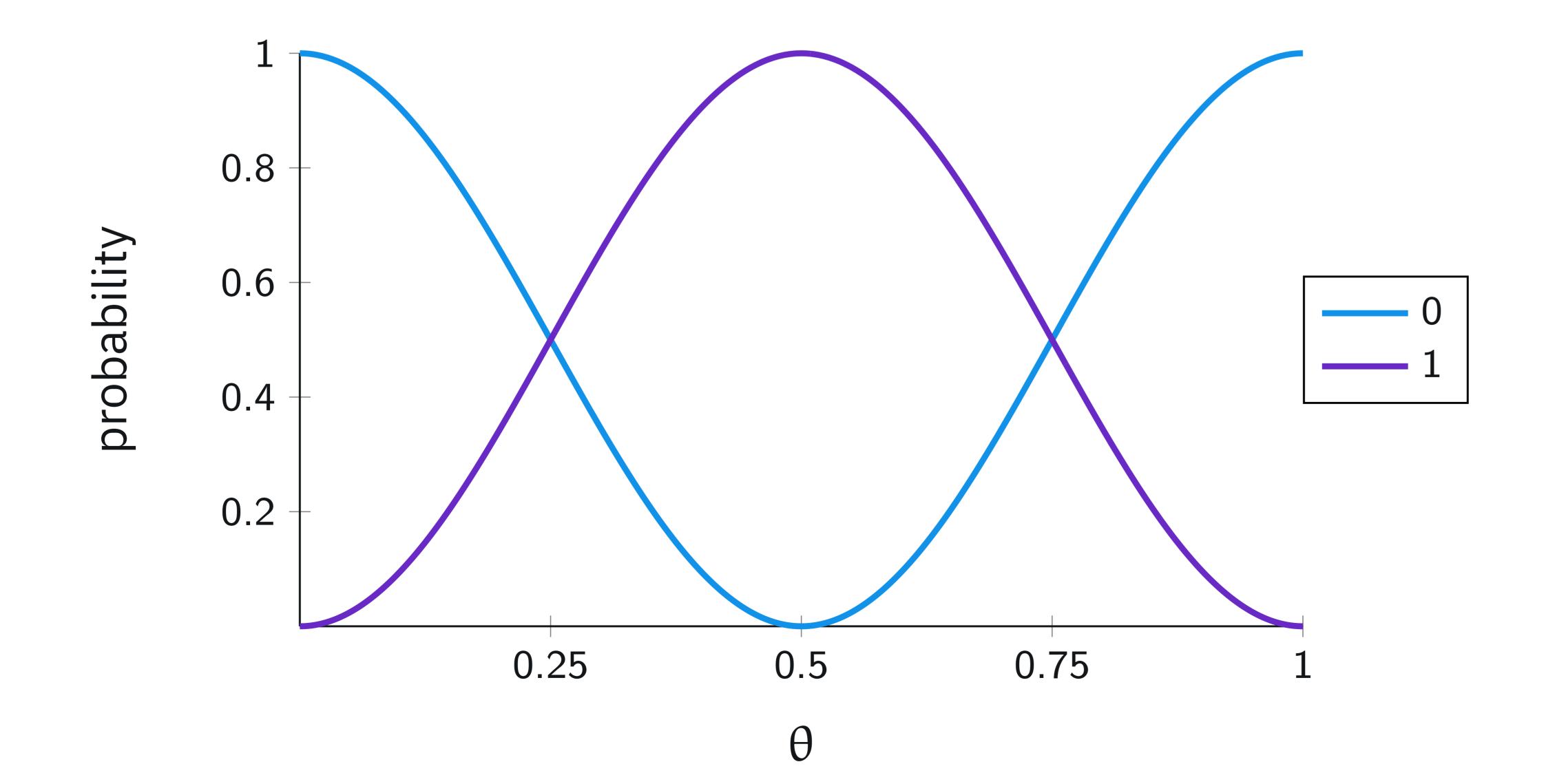
$$|\pi_{2}\rangle = |\psi\rangle \otimes \left(\frac{1}{\sqrt{2}}|0\rangle + \frac{e^{2\pi i\theta}}{\sqrt{2}}|1\rangle\right)$$

$$|\pi_{3}\rangle = |\psi\rangle \otimes \left(\frac{1 + e^{2\pi i\theta}}{2}|0\rangle + \frac{1 - e^{2\pi i\theta}}{2}|1\rangle\right)$$

$$|\psi\rangle\otimes\left(\frac{1+e^{2\pi\mathrm{i}\theta}}{2}|0\rangle+\frac{1-e^{2\pi\mathrm{i}\theta}}{2}|1\rangle\right)$$

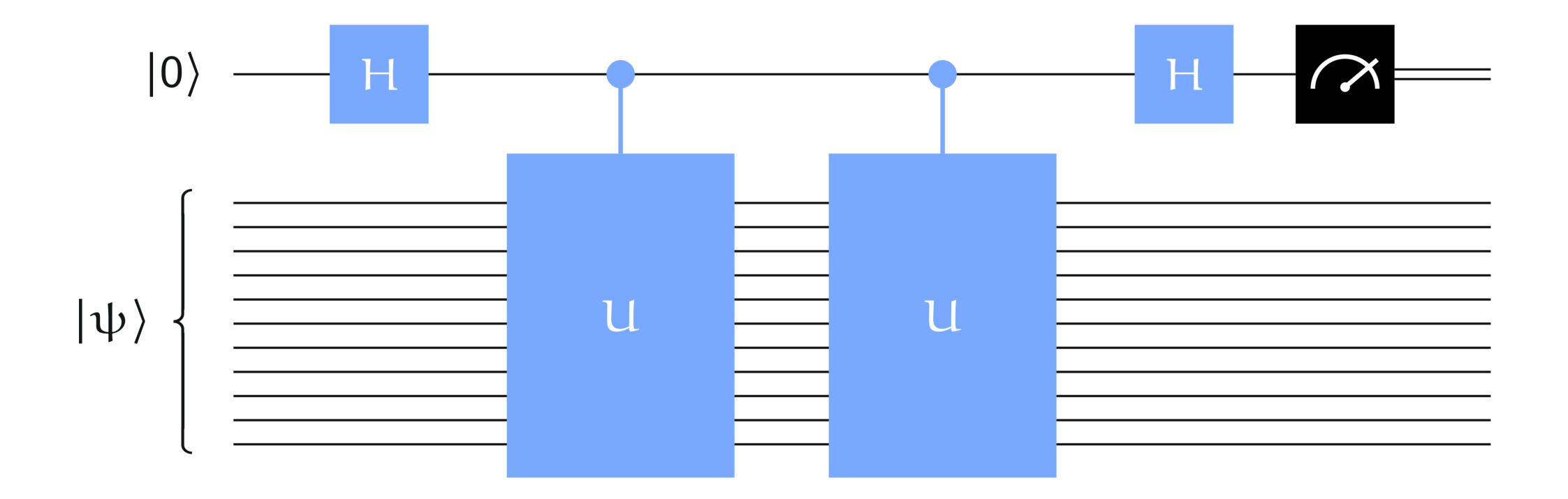
Measuring the top qubit yields the outcomes 0 and 1 with these probabilities:

$$p_0 = \left| \frac{1 + e^{2\pi i\theta}}{2} \right|^2 = \cos^2(\pi\theta)$$
 $p_1 = \left| \frac{1 - e^{2\pi i\theta}}{2} \right|^2 = \sin^2(\pi\theta)$

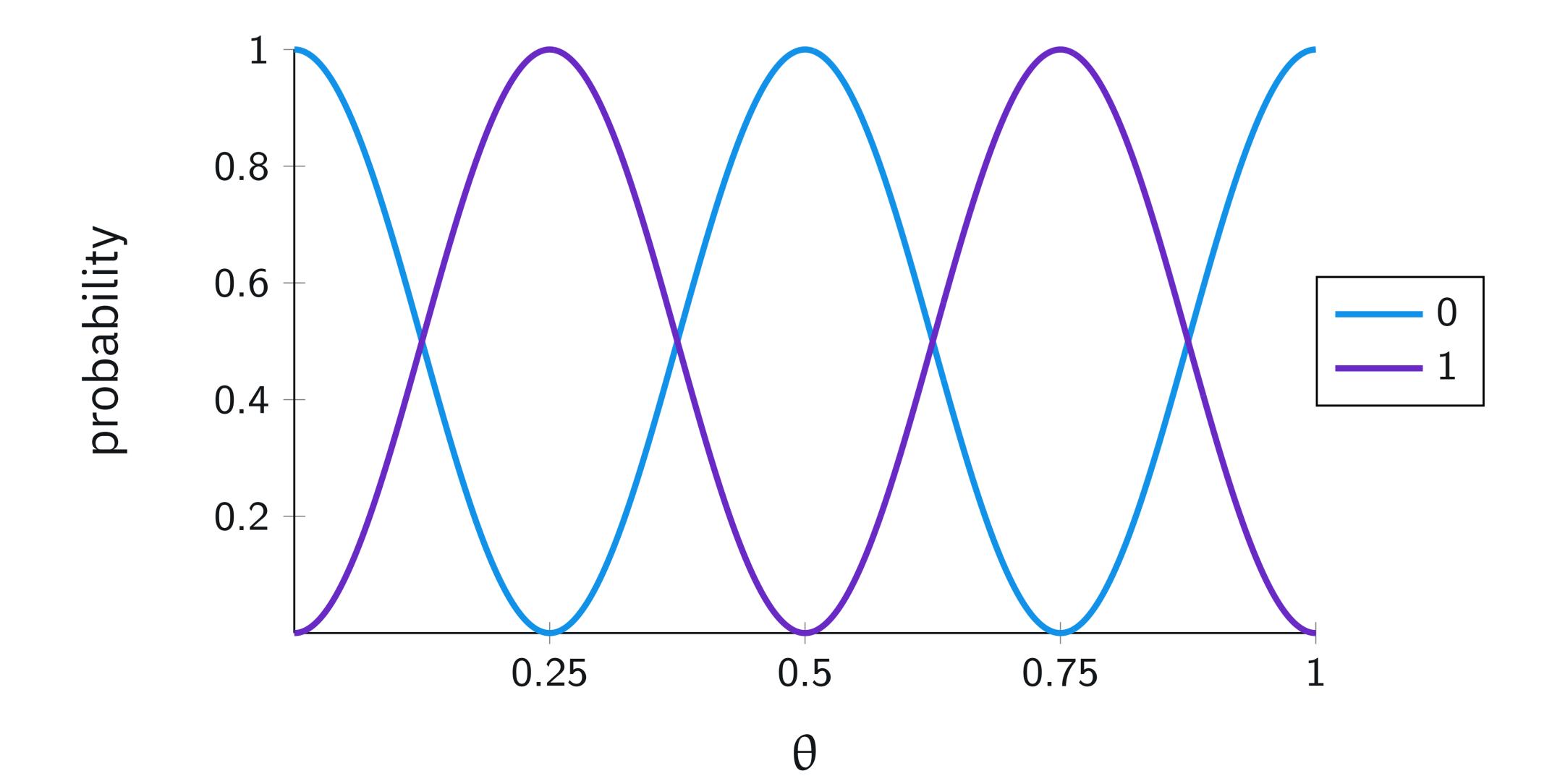


Iterating the unitary operation

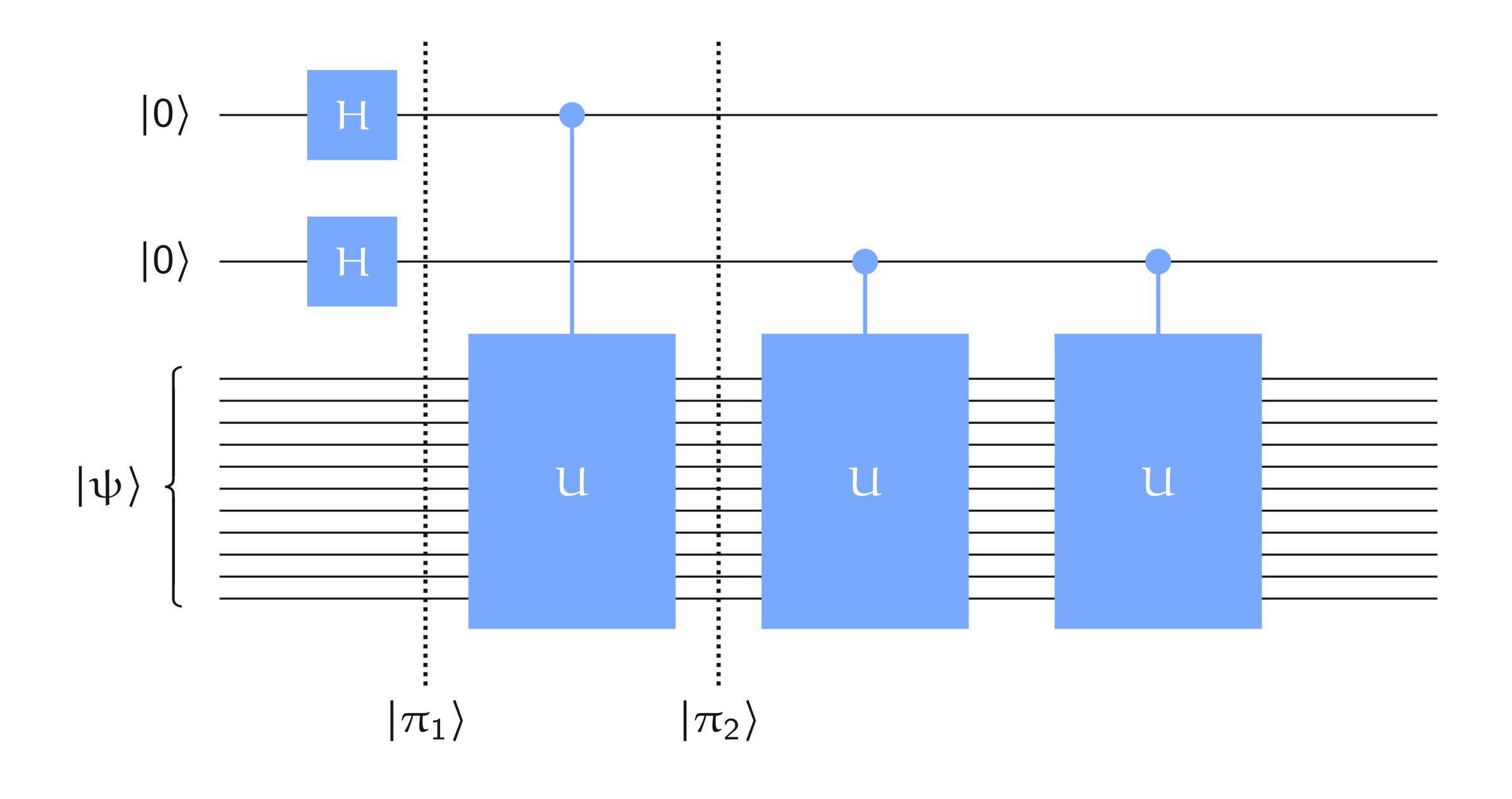
How can we learn more about θ ? One possibility is to apply the controlled-U operation twice (or multiple times):



Performing the controlled-U operation twice has the effect of squaring the eigenvalue:



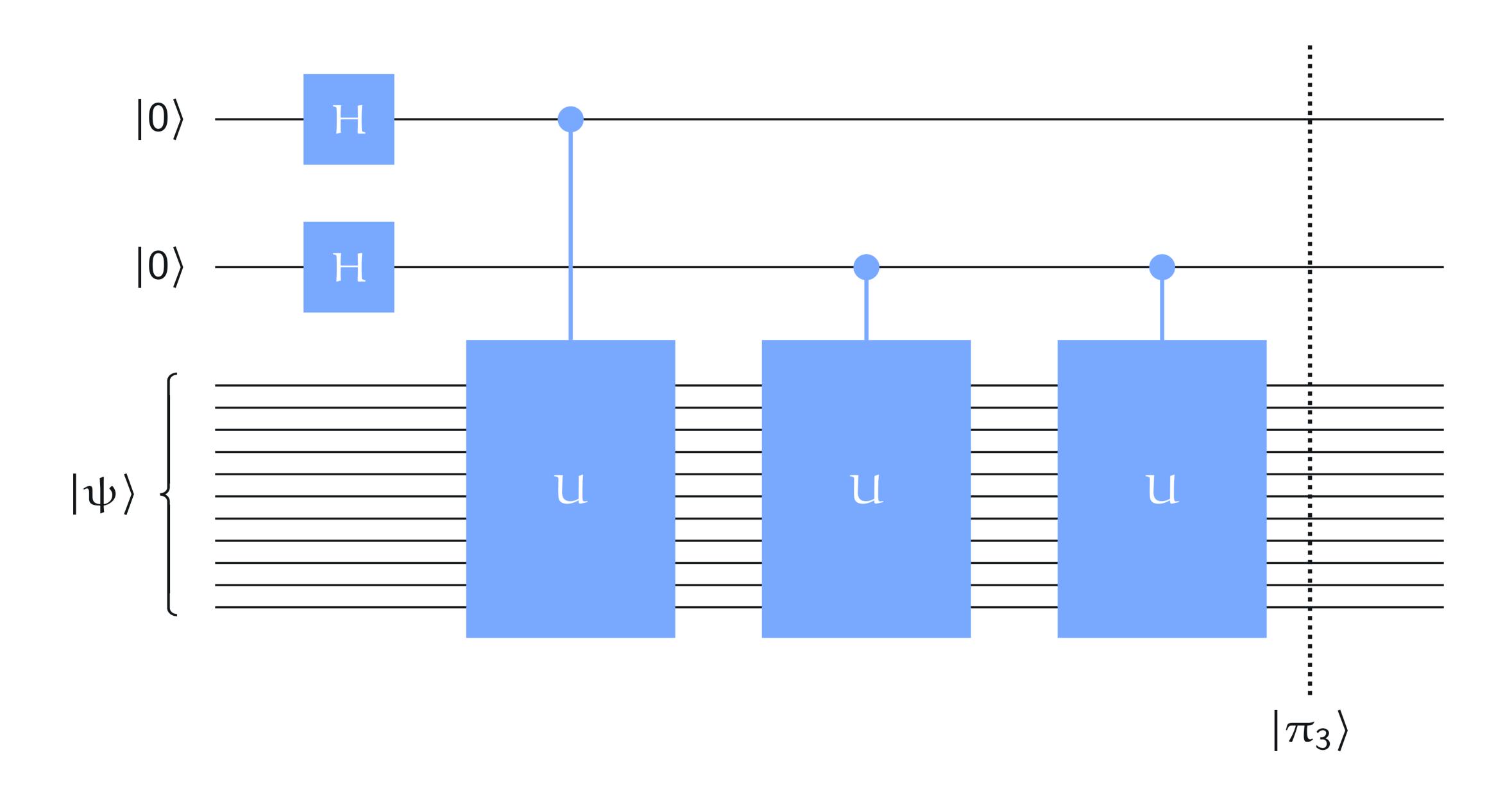
Let's use two control qubits to perform the controlled-U operations — and then we'll see how best to proceed.



$$|\pi_1\rangle = |\psi\rangle \otimes \frac{1}{2} \sum_{\alpha_0=0}^{1} \sum_{\alpha_1=0}^{1} |\alpha_1 \alpha_0\rangle$$

$$|\pi_2\rangle = |\psi\rangle \otimes \frac{1}{2} \sum_{\alpha_0=0}^{1} \sum_{\alpha_1=0}^{1} e^{2\pi i \alpha_0 \theta} |\alpha_1 \alpha_0\rangle$$

Let's use two control qubits to perform the controlled-U operations — and then we'll see how best to proceed.



$$\begin{aligned} |\pi_3\rangle &= |\psi\rangle \otimes \frac{1}{2} \sum_{\alpha_0=0}^1 \sum_{\alpha_1=0}^1 e^{2\pi i (2\alpha_1 + \alpha_0)\theta} |\alpha_1 \alpha_0\rangle \\ &= |\psi\rangle \otimes \frac{1}{2} \sum_{x=0}^3 e^{2\pi i x\theta} |x\rangle \end{aligned}$$

$$\frac{1}{2} \sum_{x=0}^{3} e^{2\pi i x \theta} |x\rangle$$

What can we learn about θ from this state? Suppose we're promised that $\theta = \frac{y}{4}$ for $y \in \{0, 1, 2, 3\}$. Can we figure out which one it is?

Define a two-qubit state for each possibility:

$$|\phi_{y}\rangle = \frac{1}{2} \sum_{x=0}^{3} e^{2\pi i \frac{xy}{4}} |x\rangle$$

$$|\phi_{0}\rangle = \frac{1}{2} |0\rangle + \frac{1}{2} |1\rangle + \frac{1}{2} |2\rangle + \frac{1}{2} |3\rangle$$

$$|\phi_{1}\rangle = \frac{1}{2} |0\rangle + \frac{i}{2} |1\rangle - \frac{1}{2} |2\rangle - \frac{i}{2} |3\rangle$$

$$|\phi_{2}\rangle = \frac{1}{2} |0\rangle - \frac{1}{2} |1\rangle + \frac{1}{2} |2\rangle - \frac{1}{2} |3\rangle$$

$$|\phi_{3}\rangle = \frac{1}{2} |0\rangle - \frac{i}{2} |1\rangle - \frac{1}{2} |2\rangle + \frac{i}{2} |3\rangle$$

These vectors are *orthonormal* — so they can be discriminated perfectly by a projective measurement.

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The unitary matrix V whose columns are $|\phi_0\rangle$, $|\phi_1\rangle$, $|\phi_2\rangle$, $|\phi_3\rangle$ has this action:

$$V|y\rangle = |\phi_y\rangle$$
 (for every $y \in \{0, 1, 2, 3\}$)

We can identify y by performing the inverse of V then a standard basis measurement.

$$V^{\dagger}|\phi_{y}\rangle=|y\rangle$$
 (for every $y\in\{0,1,2,3\}$)

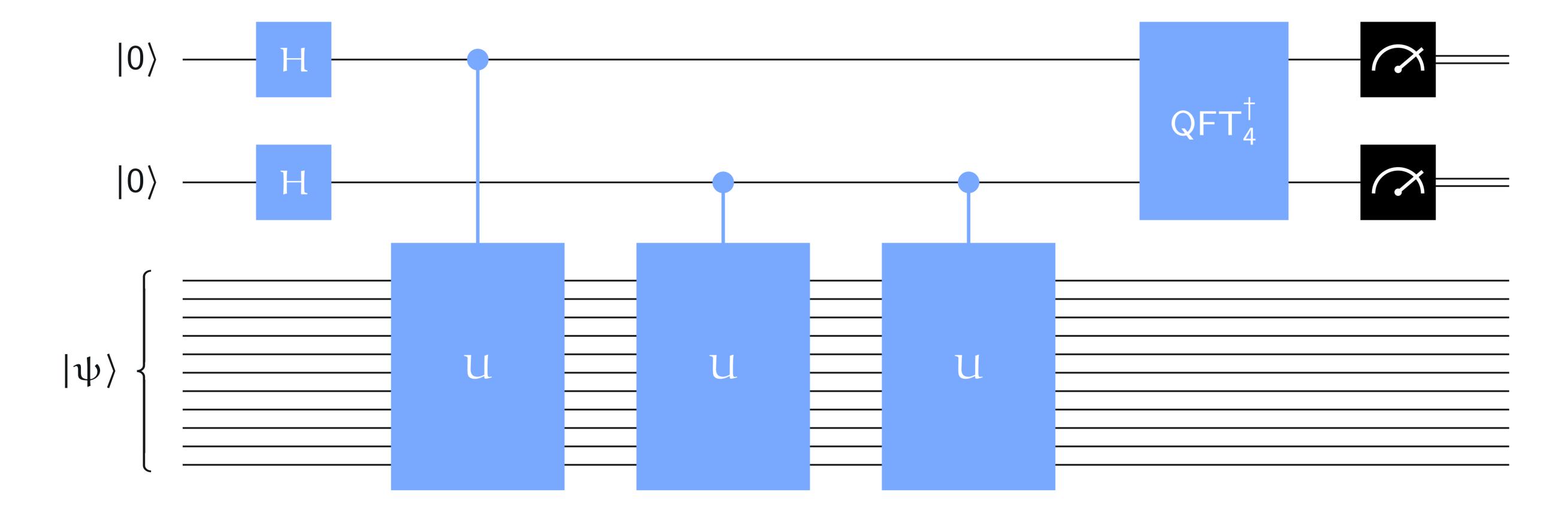
Two-qubit phase estimation

$$\mathsf{QFT_4} = \frac{1}{2} \begin{pmatrix} 1 & 1 & 1 & 1 \\ 1 & i & -1 & -i \\ 1 & -1 & 1 & -1 \\ 1 & -i & -1 & i \end{pmatrix}$$

This matrix is associated with the discrete Fourier transform (for 4 dimensions).

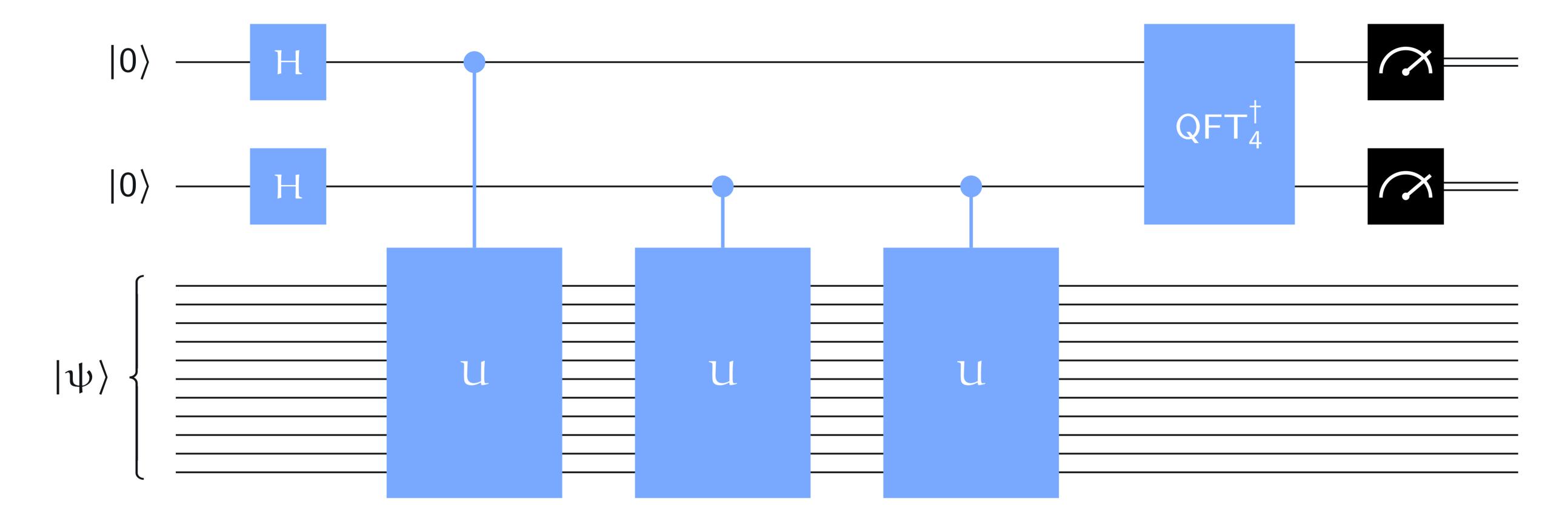
When we think about this matrix as a unitary operation, we call it the quantum Fourier transform.

The complete circuit for learning $y \in \{0, 1, 2, 3\}$ when $\theta = y/4$:

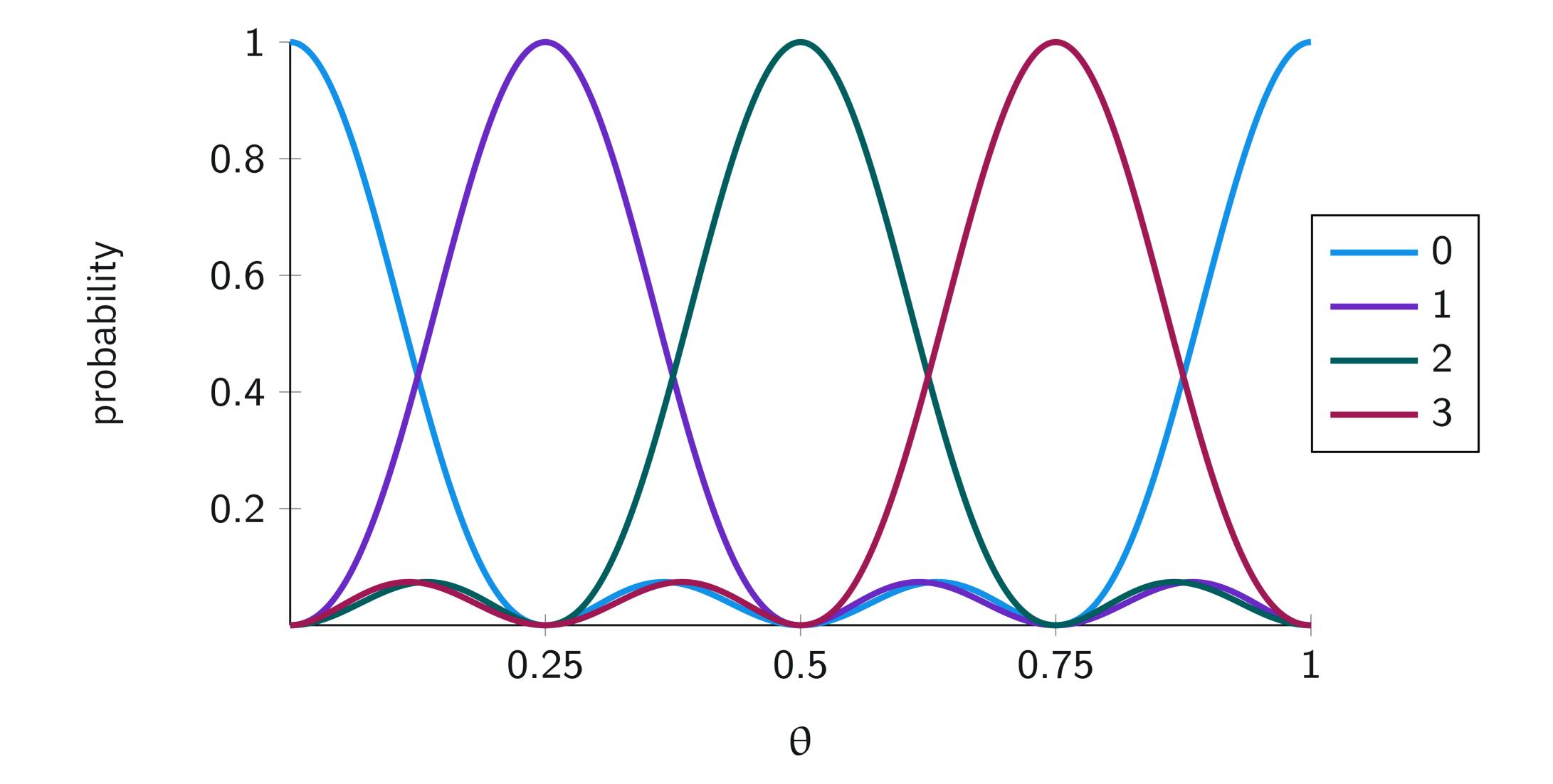


Two-qubit phase estimation

The complete circuit for learning $y \in \{0, 1, 2, 3\}$ when $\theta = y/4$:



The outcome probabilities when we run the circuit, as a function of θ :



The quantum Fourier transform is defined for each positive integer N as follows.

$$QFT_{N} = \frac{1}{\sqrt{N}} \sum_{x=0}^{N-1} \sum_{y=0}^{N-1} e^{2\pi i \frac{xy}{N}} |x\rangle\langle y|$$

$$QFT_{N}|y\rangle = \frac{1}{\sqrt{N}} \sum_{x=0}^{N-1} e^{2\pi i \frac{xy}{N}} |x\rangle$$

Example

$$QFT_2 = \frac{1}{\sqrt{2}} \begin{pmatrix} 1 & 1 \\ 1 & -1 \end{pmatrix} = H$$

$$QFT_{3} = \frac{1}{\sqrt{3}} \begin{pmatrix} 1 & 1 & 1\\ 1 & \frac{-1+i\sqrt{3}}{2} & \frac{-1-i\sqrt{3}}{2}\\ 1 & \frac{-1-i\sqrt{3}}{2} & \frac{-1+i\sqrt{3}}{2} \end{pmatrix}$$

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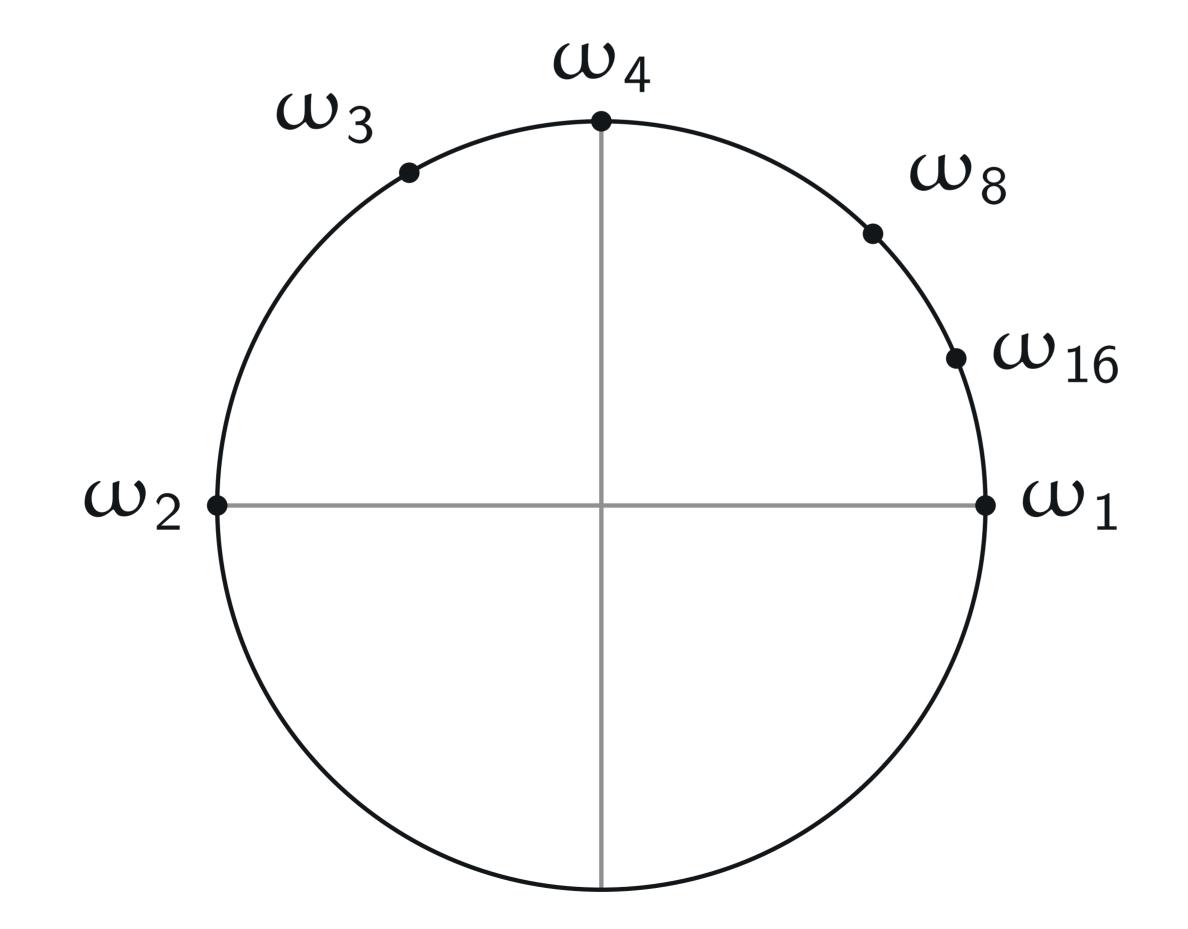
$$\mathsf{QFT}_8 = \frac{1}{2\sqrt{2}} \begin{pmatrix} 1 & 1 & 1 & 1 & 1 & 1 & 1 & 1 \\ 1 & \frac{1+i}{\sqrt{2}} & i & \frac{-1+i}{\sqrt{2}} & -1 & \frac{-1-i}{\sqrt{2}} & -i & \frac{1-i}{\sqrt{2}} \\ 1 & i & -1 & -i & 1 & i & -1 & -i \\ 1 & \frac{-1+i}{\sqrt{2}} & -i & \frac{1+i}{\sqrt{2}} & -1 & \frac{1-i}{\sqrt{2}} & i & \frac{-1-i}{\sqrt{2}} \\ 1 & -1 & 1 & -1 & 1 & -1 & 1 & -1 \\ 1 & \frac{-1-i}{\sqrt{2}} & i & \frac{1-i}{\sqrt{2}} & -1 & \frac{1+i}{\sqrt{2}} & -i & \frac{-1+i}{\sqrt{2}} \\ 1 & -i & -1 & i & 1 & -i & -1 & i \\ 1 & \frac{1-i}{\sqrt{2}} & -i & \frac{-1-i}{\sqrt{2}} & -1 & \frac{-1+i}{\sqrt{2}} & i & \frac{1+i}{\sqrt{2}} \end{pmatrix}$$

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$$QFT_{N} = \frac{1}{\sqrt{N}} \sum_{x=0}^{N-1} \sum_{y=0}^{N-1} e^{2\pi i \frac{xy}{N}} |x\rangle\langle y| = \frac{1}{\sqrt{N}} \sum_{x=0}^{N-1} \sum_{y=0}^{N-1} \omega_{N}^{xy} |x\rangle\langle y|$$

Useful shorthand notation:

$$\omega_{N} = e^{\frac{2\pi i}{N}} = \cos\left(\frac{2\pi}{N}\right) + i\sin\left(\frac{2\pi}{N}\right)$$



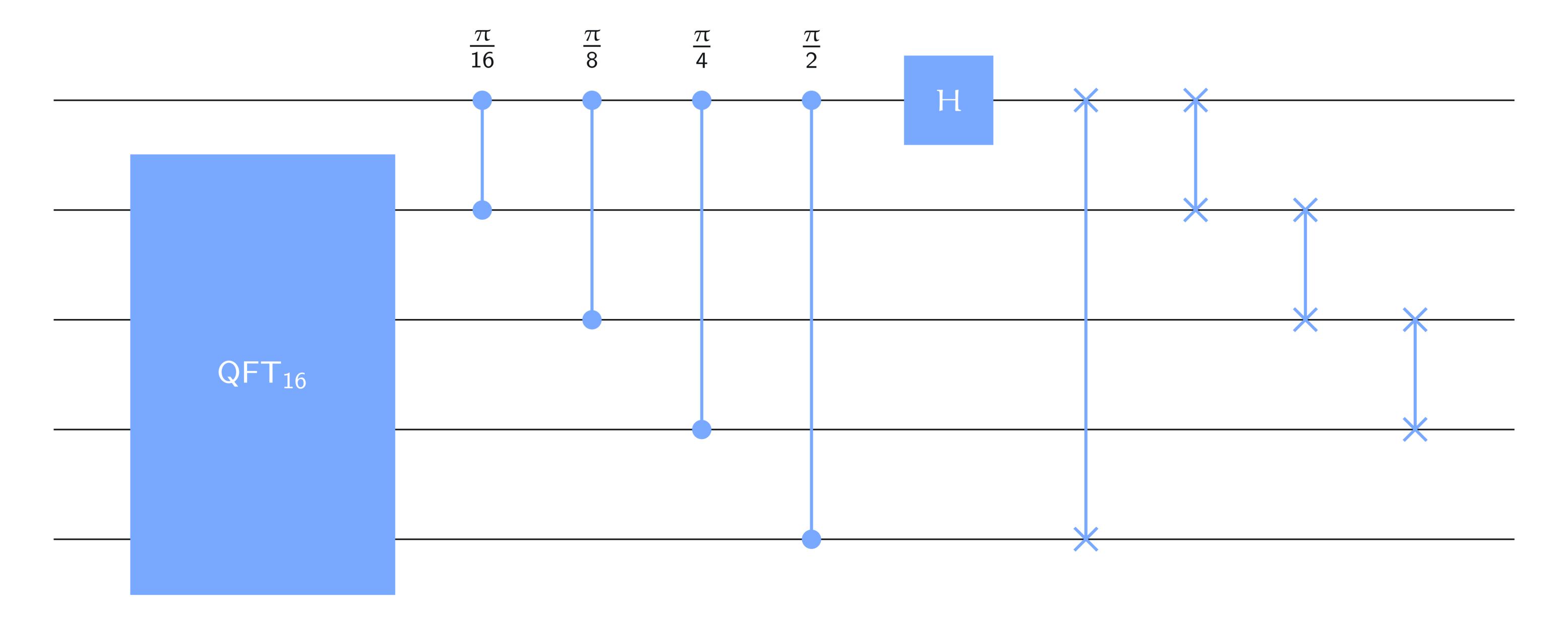
Circuits for the QFT

We can implement QFT_N efficiently with a quantum circuit when N is a power of 2.

The implementation makes use of controlled-phase gates:



The implementation is recursive in nature. As an example, here is the circuit for QFT₃₂:



Circuits for the QFT

Cost analysis

Let $s_{\rm m}$ denote the number of gates we need for m qubits.

- For m = 1, a single Hadamard gate is required.
- For $m \ge 2$, these are the gates required:

 s_{m-1} gates for the QFT on m-1 qubits

m – 1 controlled phase gates

m – 1 swap gates

1 Hadamard gate

$$s_{m} = \begin{cases} 1 & m = 1 \\ s_{m-1} + 2m - 1 & m \ge 2 \end{cases}$$

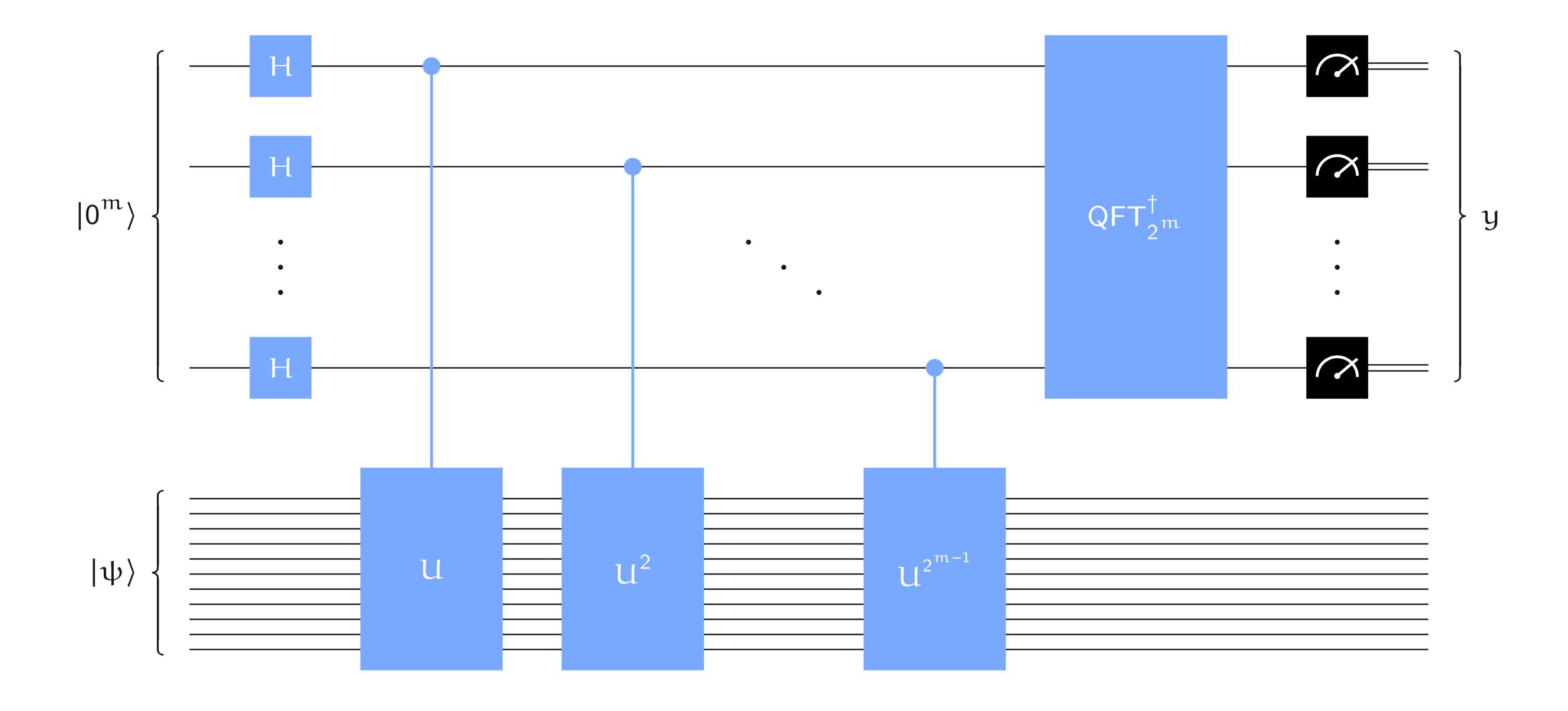
This is a *recurrence relation* with a closed-form solution:

$$s_{\rm m} = \sum_{k=1}^{\rm m} (2k - 1) = {\rm m}^2$$

Additional remarks:

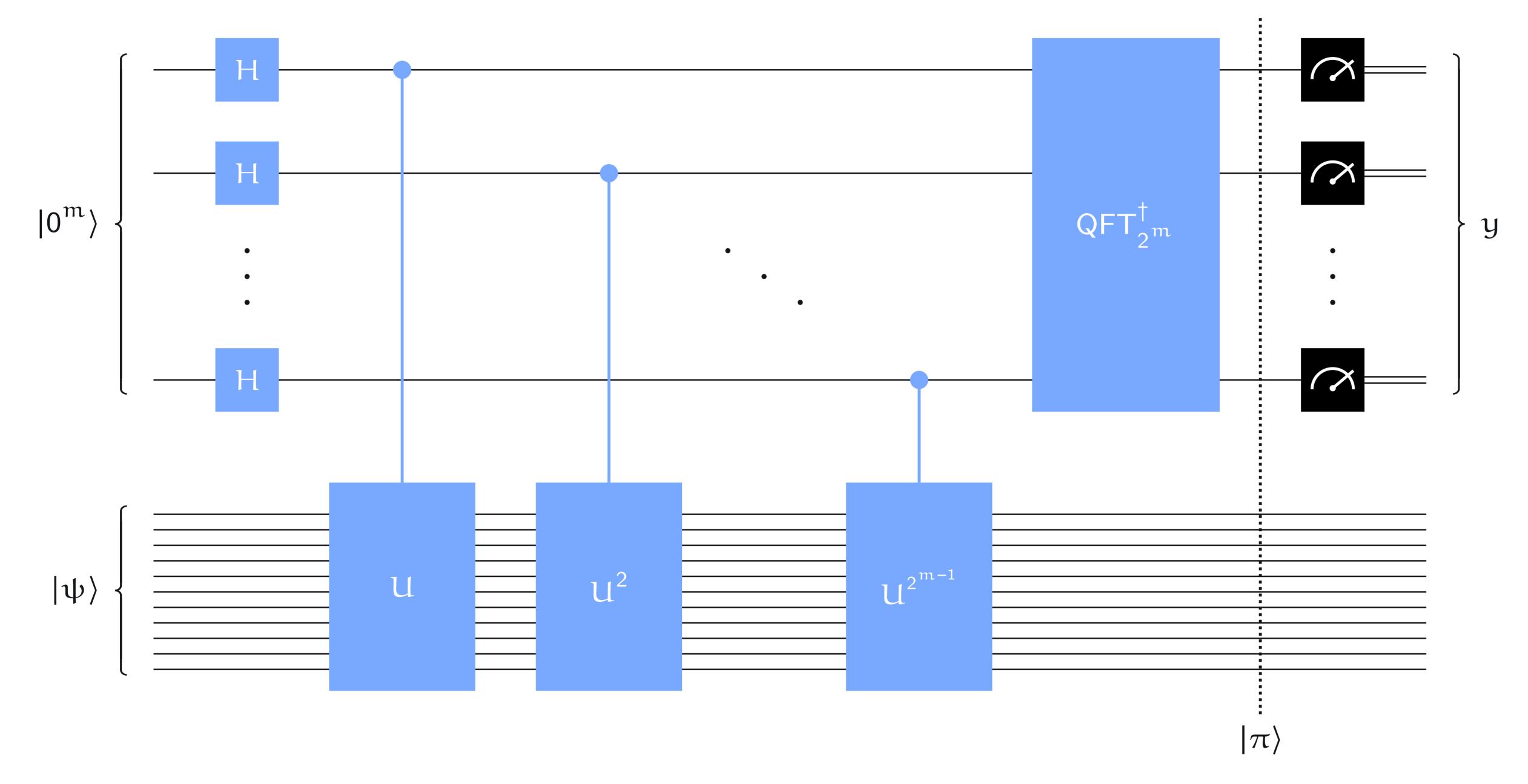
- The number of swap gates can be reduced.
- Approximations to QFT_{2m} can be done at lower cost (and lower depth).

The general phase-estimation procedure, for any choice of m:



— Warning

If we perform each U^k -operation by repeating a controlled-U operation k times, increasing the number of control qubits m comes at a <u>high cost.</u>



$$|\pi\rangle = |\psi\rangle \otimes \frac{1}{2^{m}} \sum_{y=0}^{2^{m}-1} \sum_{x=0}^{2^{m}-1} e^{2\pi i x (\theta - y/2^{m})} |y\rangle$$

$$p_{y} = \left| \frac{1}{2^{m}} \sum_{x=0}^{2^{m}-1} e^{2\pi i x (\theta - y/2^{m})} \right|^{2}$$

Best approximations

Suppose $y/2^m$ is the *best approximation* to θ :

$$\left|\theta - \frac{y}{2^{m}}\right|_{1} \le 2^{-(m+1)}$$

Then the probability to measure y will relatively high:

$$p_y \geq \frac{4}{\pi^2} \approx 0.405$$

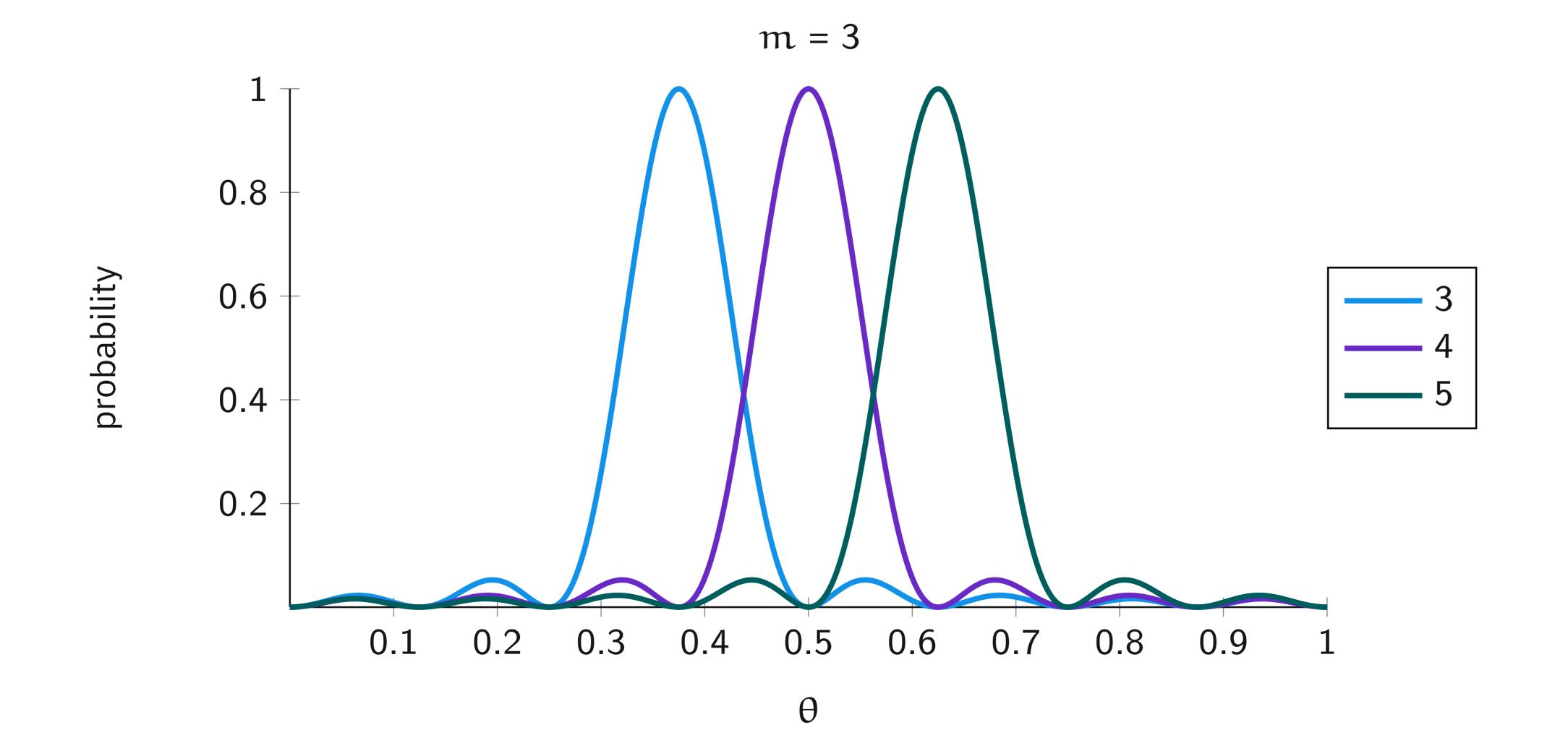
Worse approximations

Suppose there's a *better approximation* to θ between $y/2^m$ and θ :

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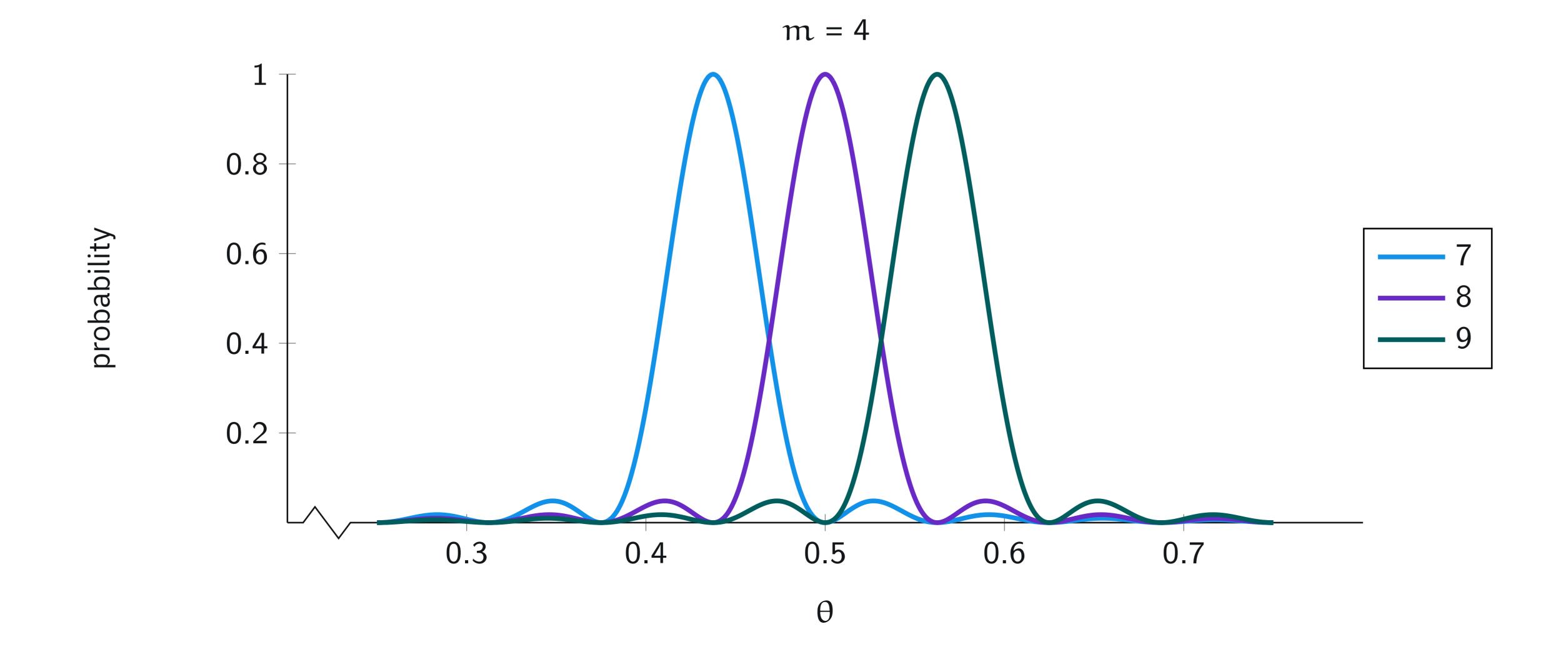
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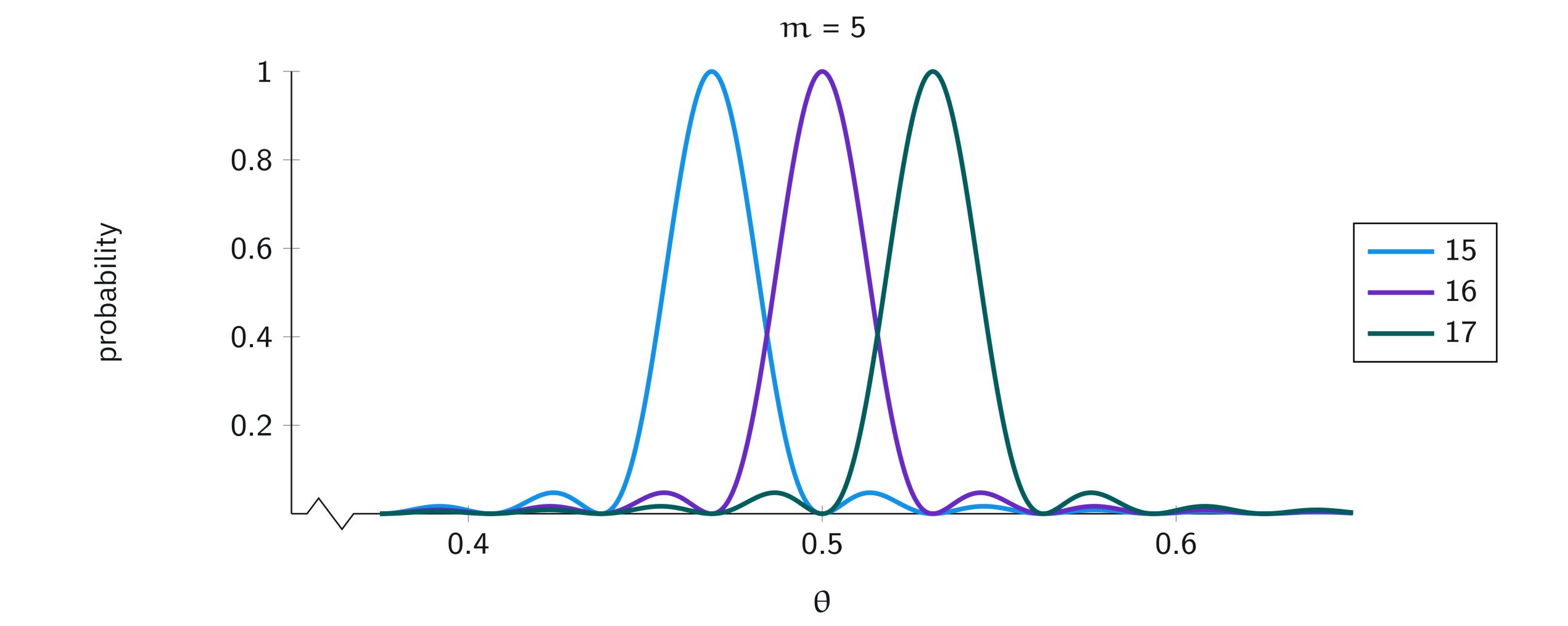
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To obtain an approximation $y/2^m$ that is very likely to satisfy

$$\left|\theta-\frac{y}{2^{m}}\right|_{1}<2^{-m}$$

we can run the phase estimation procedure using \mathfrak{m} control qubits several times and take \mathfrak{y} to be the mode of the outcomes.

(The eigenvector $|\psi\rangle$ is unchanged by the procedure and can be reused as many times as needed.)

The order-finding problem

For each positive integer N we define

$$\mathbb{Z}_{N} = \{0, 1, \dots, N-1\}$$

For instance, $\mathbb{Z}_1 = \{0\}$, $\mathbb{Z}_2 = \{0, 1\}$, $\mathbb{Z}_3 = \{0, 1, 2\}$, and so on.

We can view arithmetic operations on \mathbb{Z}_{N} as being defined modulo N.

Example

Let N = 7. We have $3 \cdot 5 = 15$, which leaves a remainder of 1 when divided by 7.

This is often expressed like this:

$$3 \cdot 5 \equiv 1 \pmod{7}$$

We can also simply write $3 \cdot 5 = 1$ when it's clear we're working in \mathbb{Z}_7 .

The elements $\alpha \in \mathbb{Z}_N$ that satisfy $gcd(\alpha, N) = 1$ are special.

$$\mathbb{Z}_{N}^{*} = \{\alpha \in \mathbb{Z}_{N} : \gcd(\alpha, N) = 1\}$$

$$\mathbb{Z}_{21}^* = \{1, 2, 4, 5, 8, 10, 11, 13, 16, 17, 19, 20\}$$

The order-finding problem

Fact

For every $\alpha \in \mathbb{Z}_N^*$ there must exist a positive integer k such that $\alpha^k = 1$. The smallest such k is called the <u>order</u> of α in \mathbb{Z}_N^* .

Example

For N=21, these are the smallest powers for which this works:

$$1^{1} = 1$$
 $5^{6} = 1$ $11^{6} = 1$ $17^{6} = 1$ $2^{6} = 1$ $8^{2} = 1$ $13^{2} = 1$ $19^{6} = 1$ $4^{3} = 1$ $10^{6} = 1$ $16^{3} = 1$ $20^{2} = 1$

Order-finding problem

Input: Positive integers α and N with $gcd(\alpha, N) = 1$.

Output: The smallest positive integer r such that $\alpha^r \equiv 1 \pmod{N}$

No efficient classical algorithm for this problem is known — an efficient algorithm for order-finding implies an efficient algorithm for integer factorization.

Order-finding by phase-estimation

To connect the order-finding problem to phase estimation, consider a system whose classical state set is \mathbb{Z}_N .

For a given element $\alpha \in \mathbb{Z}_N^*$, define an operation as follows:

$$M_{\alpha}|x\rangle = |\alpha x\rangle$$
 (for each $x \in \mathbb{Z}_{N}$)

This is a *unitary operation* — but only because $gcd(\alpha, N) = 1!$

Example

Let N=15 and $\alpha=2$. The operation M_{α} has this action:

$$\begin{array}{lll} M_2|0\rangle = |0\rangle & M_2|5\rangle = |10\rangle & M_2|10\rangle = |5\rangle \\ M_2|1\rangle = |2\rangle & M_2|6\rangle = |12\rangle & M_2|11\rangle = |7\rangle \\ M_2|2\rangle = |4\rangle & M_2|7\rangle = |14\rangle & M_2|12\rangle = |9\rangle \\ M_2|3\rangle = |6\rangle & M_2|8\rangle = |1\rangle & M_2|13\rangle = |11\rangle \\ M_2|4\rangle = |8\rangle & M_2|9\rangle = |3\rangle & M_2|14\rangle = |13\rangle \end{array}$$

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Main idea

The eigenvalues of $\mathcal{M}_{\mathfrak{a}}$ are closely connected with the order of \mathfrak{a} .

By approximating certain eigenvalues with enough precision using phase estimation, we'll be able to compute the order.

Eigenvectors and eigenvalues

This is an eigenvector of M_{α} :

$$|\psi_0\rangle = \frac{|1\rangle + |\alpha\rangle + \dots + |\alpha^{r-1}\rangle}{\sqrt{r}}$$

The associated eigenvalue is 1:

$$\mathcal{M}_{\alpha}|\psi_{0}\rangle = \frac{|\alpha\rangle + |\alpha^{2}\rangle + \cdots + |\alpha^{r}\rangle}{\sqrt{r}} = \frac{|\alpha\rangle + \cdots + |\alpha^{r-1}\rangle + |1\rangle}{\sqrt{r}} = |\psi_{0}\rangle$$

To identify more eigenvectors, first recall that

$$\omega_{\rm r} = e^{2\pi i/r}$$

This is another eigenvector of M_{α} :

$$|\psi_1\rangle = \frac{|1\rangle + \omega_r^{-1}|\alpha\rangle + \dots + \omega_r^{-(r-1)}|\alpha^{r-1}\rangle}{\sqrt{r}}$$

Eigenvectors and eigenvalues

$$\begin{split} M_{\alpha}|\psi_{1}\rangle &= \frac{|\alpha\rangle + \omega_{r}^{-1}|\alpha^{2}\rangle + \cdots + \omega_{r}^{-(r-1)}|\alpha^{r}\rangle}{\sqrt{r}} \\ &= \frac{\omega_{r}|1\rangle + |\alpha\rangle + \omega_{r}^{-1}|\alpha^{2}\rangle + \cdots + \omega_{r}^{-(r-2)}|\alpha^{r-1}\rangle}{\sqrt{r}} \\ &= \omega_{r} \left(\frac{|1\rangle + \omega_{r}^{-1}|\alpha\rangle + \omega_{r}^{-2}|\alpha^{2}\rangle + \cdots + \omega_{r}^{-(r-1)}|\alpha^{r-1}\rangle}{\sqrt{r}}\right) \\ &= \omega_{r}|\psi_{1}\rangle \end{split}$$

Additional eigenvectors can be identified by similar reasoning...

For each $j \in \{0, ..., r-1\}$, this is an eigenvector of M_a :

$$|\psi_{j}\rangle = \frac{|1\rangle + \omega_{r}^{-j}|\alpha\rangle + \dots + \omega_{r}^{-j(r-1)}|\alpha^{r-1}\rangle}{\sqrt{r}}$$

$$M_{\alpha}|\psi_{j}\rangle = \omega_{r}^{j}|\psi_{j}\rangle$$

A convenient eigenvector

$$|\psi_{1}\rangle = \frac{|1\rangle + \omega_{r}^{-1}|\alpha\rangle + \dots + \omega_{r}^{-(r-1)}|\alpha^{r-1}\rangle}{\sqrt{r}}$$

$$M_{\alpha}|\psi_{1}\rangle = \omega_{r}|\psi_{1}\rangle = e^{2\pi i \frac{1}{r}}|\psi_{1}\rangle$$

Suppose we're given $|\psi_1\rangle$ as a quantum state. We can attempt to learn r as follows:

- 1. Perform phase estimation on the state $|\psi_1\rangle$ and a quantum circuit implementing M_a . The outcome is an approximation $y/2^m \approx 1/r$.
- 2. Output $2^{m}/y$ rounded to the nearest integer:

$$\operatorname{round}\left(\frac{2^{m}}{y}\right) = \left\lfloor \frac{2^{m}}{y} + \frac{1}{2} \right\rfloor$$

How much precision do we need to correctly determine r?

$$\left|\frac{y}{2^{m}} - \frac{1}{r}\right| \le \frac{1}{2N^2} \implies \text{round}\left(\frac{2^{m}}{y}\right) = r$$

Choosing $m = 2 \lg(N) + 1$ in phase estimation makes such an approximation likely.

A random eigenvector

$$|\psi_{j}\rangle = \frac{|1\rangle + \omega_{r}^{-j}|\alpha\rangle + \dots + \omega_{r}^{-j(r-1)}|\alpha^{r-1}\rangle}{\sqrt{r}}$$

$$M_{\alpha}|\psi_{j}\rangle = \omega_{r}^{j}|\psi_{1}\rangle = e^{2\pi i \frac{j}{r}}|\psi_{1}\rangle$$

Suppose we're given $|\psi_j\rangle$ as a quantum state for a random choice of $j \in \{0, ..., r-1\}$. We can attempt to learn j/r as follows:

- 1. Perform phase estimation on the state $|\psi_j\rangle$ and a quantum circuit implementing M_a . The outcome is an approximation $y/2^m \approx j/r$.
- 2. Among the fractions u/v in lowest terms satisfying $u, v \in \{0, ..., N-1\}$ and $v \neq 0$, output the one closest to $y/2^m$. This can be done efficiently using the continued fraction algorithm.

How much precision do we need to correctly determine u/v = j/r?

$$\left|\frac{y}{2^m} - \frac{j}{r}\right| \le \frac{1}{2N^2} \implies \frac{u}{v} = \frac{j}{r}$$

Choosing $m = 2 \lg(N) + 1$ for phase estimation makes such an approximation likely. We might get unlucky: j could have common factors with r.

A random eigenvector

$$|\psi_{j}\rangle = \frac{|1\rangle + \omega_{r}^{-j}|\alpha\rangle + \dots + \omega_{r}^{-j(r-1)}|\alpha^{r-1}\rangle}{\sqrt{r}}$$

$$M_{\alpha}|\psi_{j}\rangle = \omega_{r}^{j}|\psi_{1}\rangle = e^{2\pi i \frac{j}{r}}|\psi_{1}\rangle$$

Suppose we're given $|\psi_j\rangle$ as a quantum state for a random choice of $j \in \{0, ..., r-1\}$. We can attempt to learn j/r as follows:

- 1. Perform phase estimation on the state $|\psi_j\rangle$ and a quantum circuit implementing M_a . The outcome is an approximation $y/2^m \approx j/r$.
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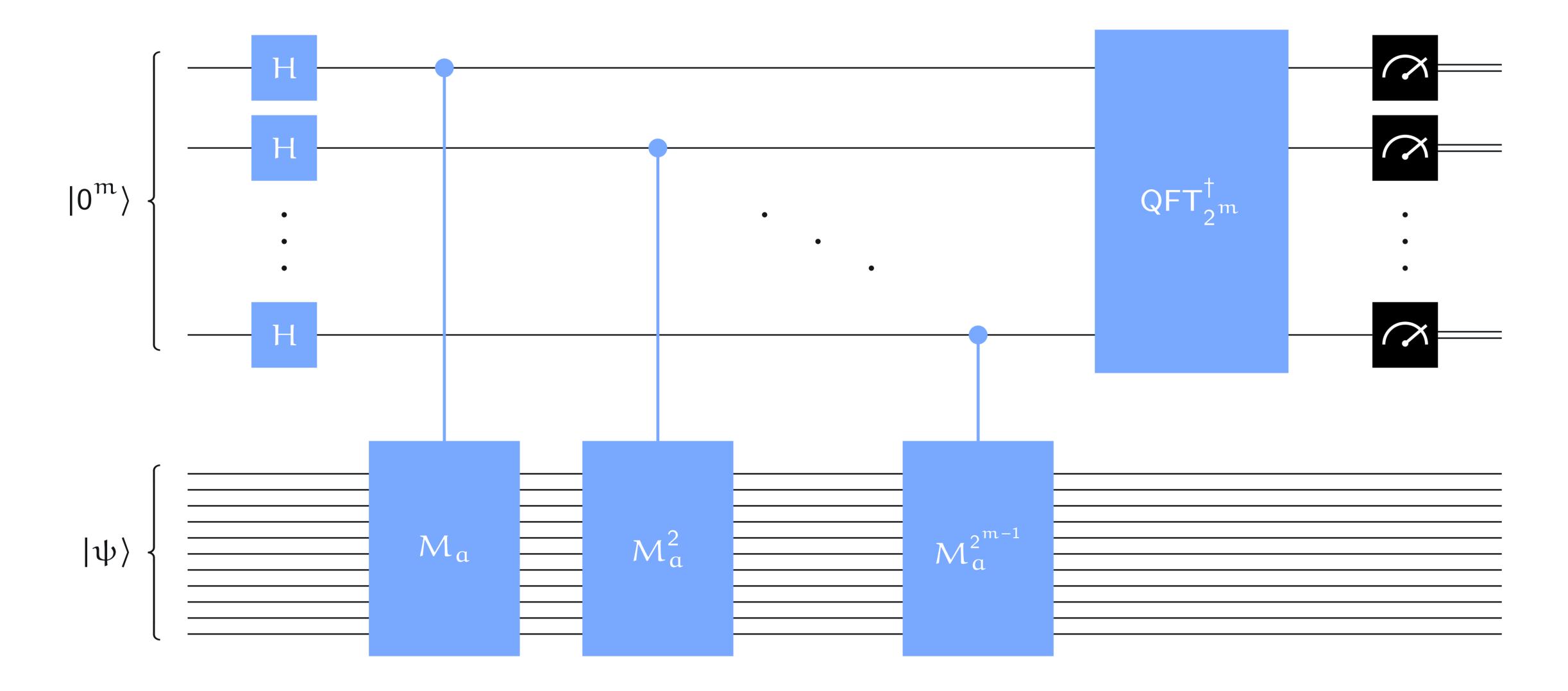
How much precision do we need to correctly determine u/v = j/r?

$$\left|\frac{y}{2^m} - \frac{j}{r}\right| \le \frac{1}{2N^2} \implies \frac{u}{v} = \frac{j}{r}$$

If we can draw *independent samples,* for $j \in \{0, ..., r-1\}$ is chosen uniformly, we can recover r with high probability by computing the *least common multiple* of the values of v we observed.

Implementation

To find the order of $\alpha \in \mathbb{Z}_N^*$, we apply phase estimation to the operation \mathcal{M}_α . Let's measure the cost as a function of $n = \lg(N)$.



Cost for each controlled unitary

Using the techniques from Lesson 6, we can implement $\mathcal{M}_{\mathfrak{a}}$ at cost $O(\mathfrak{n}^2)$.

We need to implement M_a^k for each $k = 1, 2, 4, 8, ..., 2^{m-1}$. Each M_a^k can be implemented as follows:

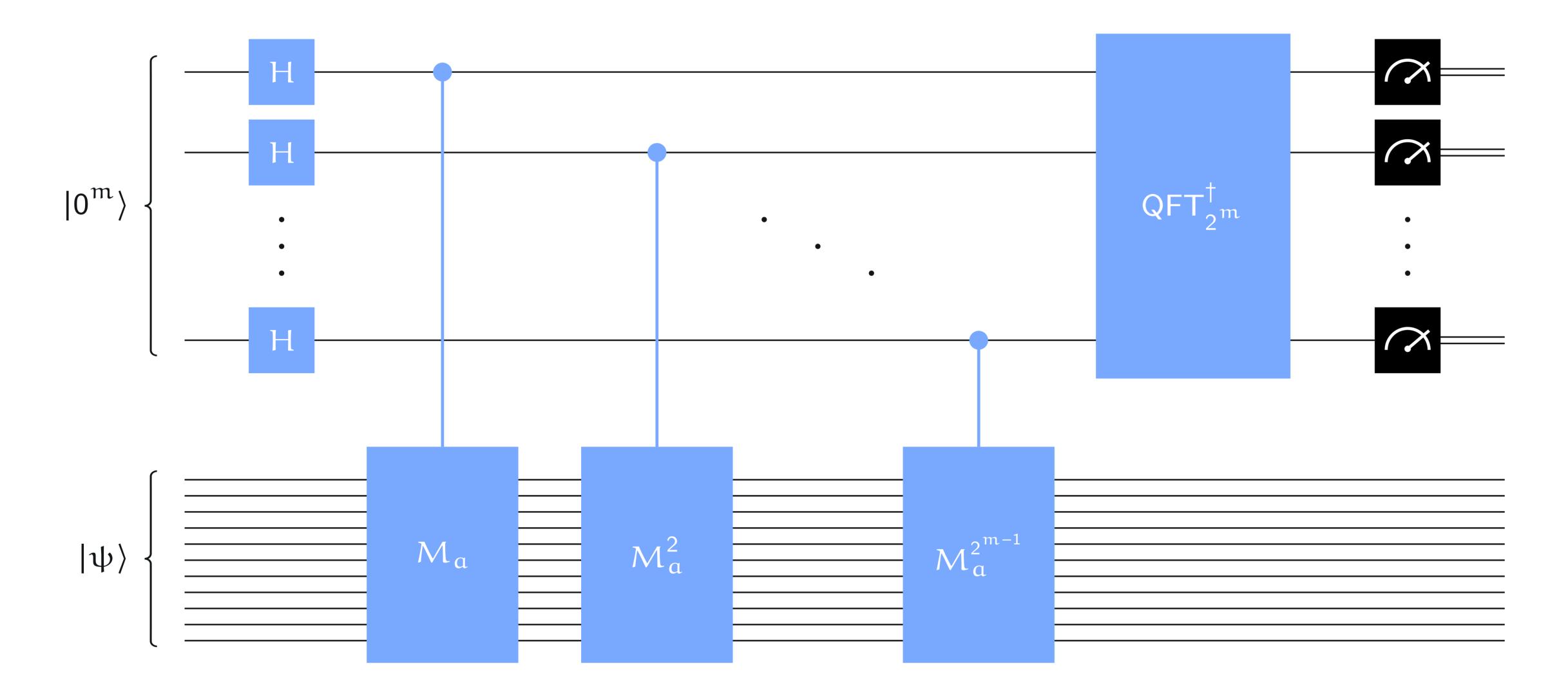
Compute $b = a^k \pmod{N}$.

Use a circuit for M_b .

The cost to implement $M_b = M_a^k$ is $O(n^2)$.

Implementation

To find the order of $\alpha \in \mathbb{Z}_N^*$, we apply phase estimation to the operation \mathcal{M}_{α} . Let's measure the cost as a function of $n = \lg(N)$.

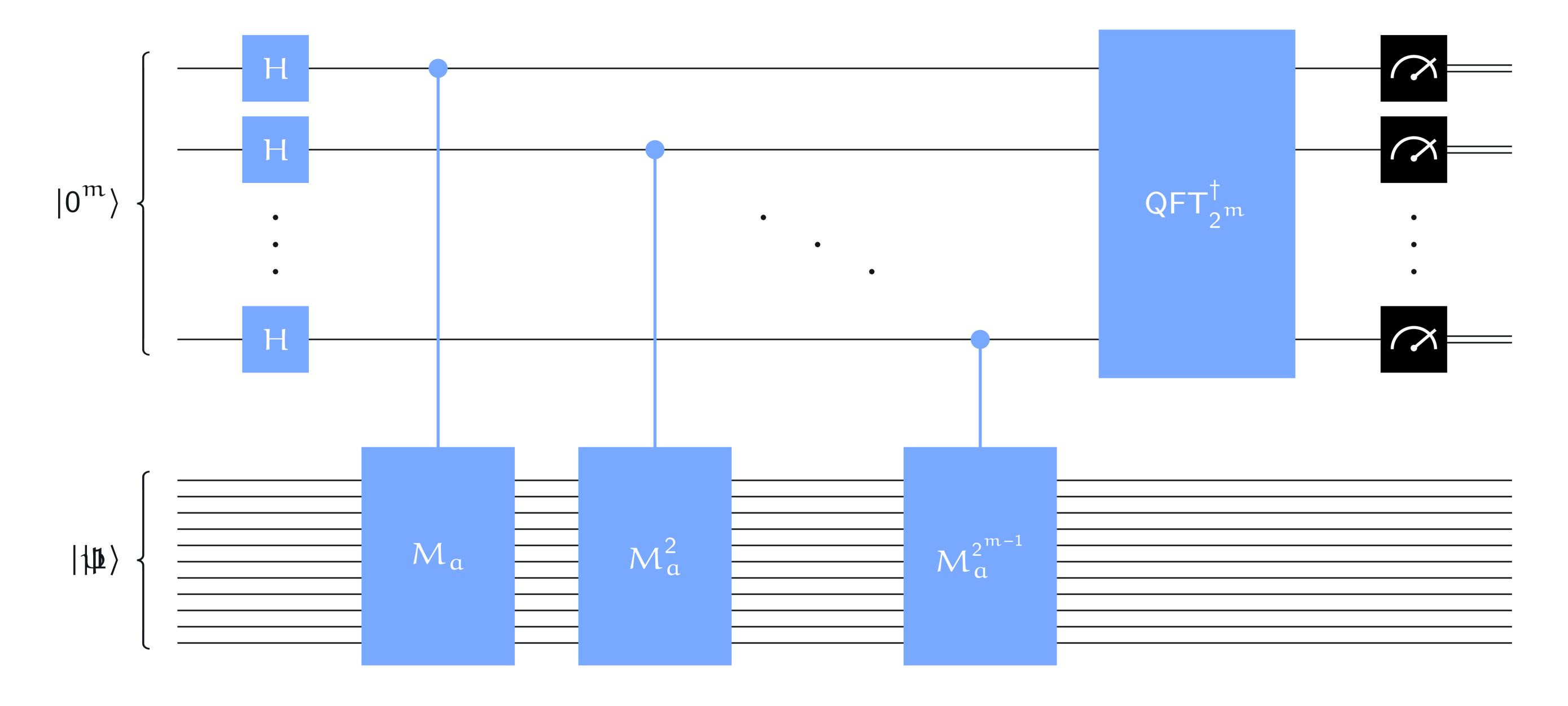


Cost for phase estimation

- m Hadamard gates: cost O(n)
- m controlled unitary operations: cost O(n³)
 Quantum Fourier transform: cost O(n²)

Total cost: $O(n^3)$

Implementation



Remaining issue: getting one of the eigenvectors $|\psi_0\rangle$, . . . , $|\psi_{r-1}\rangle$.

Solution: replace the eigenvector $|\psi\rangle$ with the state $|1\rangle$.

This works because of the following equation:

$$|1\rangle = \frac{|\psi_0\rangle + \dots + |\psi_{r-1}\rangle}{\sqrt{r}}$$

The outcome is the same as if we chose $j \in \{0, 1, ..., r-1\}$ uniformly and used $|\psi\rangle = |\psi_j\rangle$.

Factoring through order-finding

The following method succeeds in finding a factor of N with probability at least 1/2, provided N is odd and not a prime power.

Factor-finding method

- 1. Choose $\alpha \in \{2, ..., N-1\}$ at random.
- 2. Compute $d = \gcd(\alpha, N)$. If $d \ge 2$ then output d and stop.
- 3. Compute the order r of a modulo N.
- 4. If r is even, then compute $d = \gcd(\alpha^{r/2} 1, N)$. If $d \ge 2$, output d and stop.
- 5. If this step is reached, the method has failed.

Main idea

1. By the definition of the order, we know that $\alpha^r \equiv 1 \pmod{N}$.

$$a^{r} \equiv 1 \pmod{N} \iff N \text{ divides } a^{r} - 1$$

2. If r is even, then

$$a^{r} - 1 = (a^{r/2} + 1)(a^{r/2} - 1)$$

Each prime dividing N must therefore divide either $(\alpha^{r/2} + 1)$ or $(\alpha^{r/2} - 1)$. For a random α , at least one of the prime factors of N is likely to divide $(\alpha^{r/2} - 1)$.