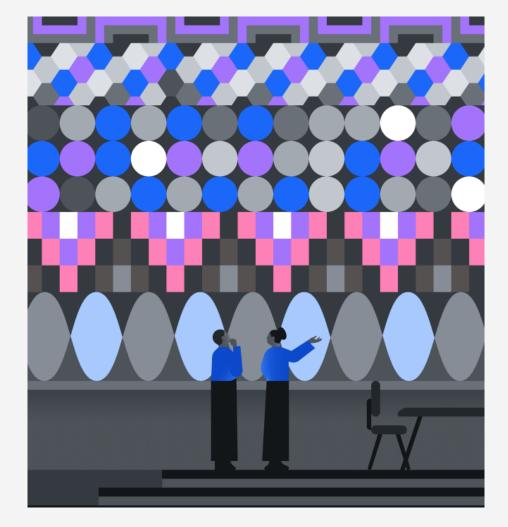
# Understanding quantum information and computation

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  $N \times N$  unitary matrix.

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

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

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  $\lambda_k$ :

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

# 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 quantum state  $|\psi\rangle$ .

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

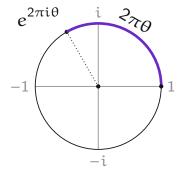
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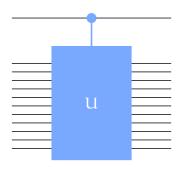
We can approximate  $\theta$  by a fraction

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

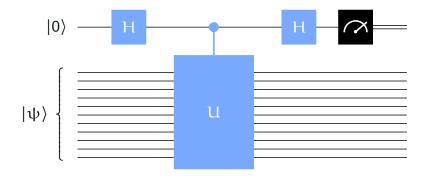
for  $y \in \{0, 1, ..., 2^m - 1\}$ .

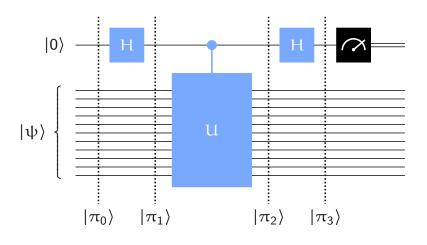
This approximation is taken "modulo 1."

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

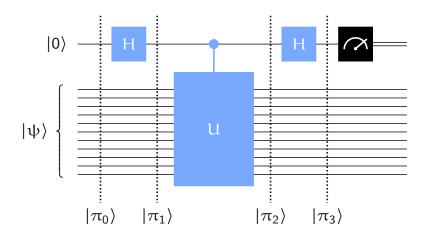


Let's consider this circuit:





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

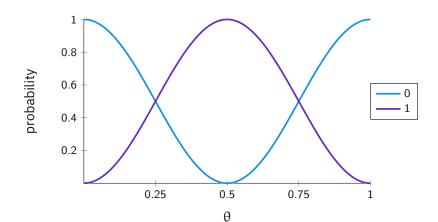


$$\begin{split} |\pi_2\rangle &= |\psi\rangle \otimes \left(\frac{1}{\sqrt{2}}|0\rangle + \frac{e^{2\pi i\theta}}{\sqrt{2}}|1\rangle\right) \\ |\pi_3\rangle &= |\psi\rangle \otimes \left(\frac{1 + e^{2\pi i\theta}}{2}|0\rangle + \frac{1 - e^{2\pi i\theta}}{2}|1\rangle\right) \end{split}$$

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

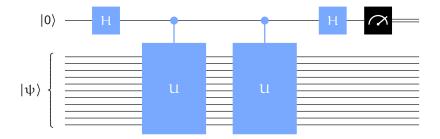
Measuring the top qubit yields the outcomes 0 and 1 with these probabilities:

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

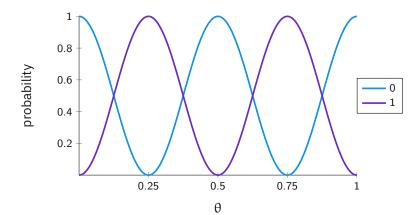


# Iterating the unitary operation

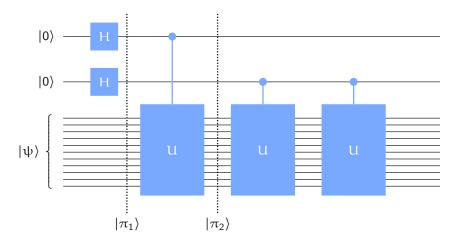
How can we learn more about  $\theta$ ? One possibility is to apply the controlled-U operation twice (or multiple times):



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

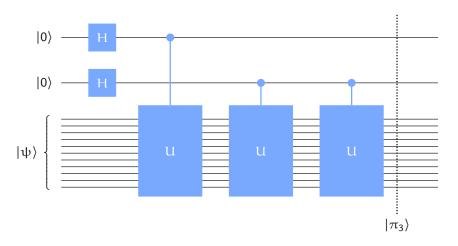


Let's use two control qubits to perform the controlled- $\ensuremath{\mathsf{U}}$  operations — and then we'll see how best to proceed.



$$\begin{split} |\pi_1\rangle &= |\psi\rangle \otimes \frac{1}{2} \sum_{\alpha_0=0}^1 \sum_{\alpha_1=0}^1 |\alpha_1 \alpha_0\rangle \\ |\pi_2\rangle &= |\psi\rangle \otimes \frac{1}{2} \sum_{\alpha_0=0}^1 \sum_{\alpha_1=0}^1 e^{2\pi i \alpha_0 \theta} |\alpha_1 \alpha_0\rangle \end{split}$$

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



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

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

What can we learn about  $\theta$  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.

$$\begin{split} |\varphi_{y}\rangle &= \frac{1}{2} \sum_{x=0}^{3} e^{2\pi i \frac{xy}{4}} |x\rangle \\ |\varphi_{0}\rangle &= \frac{1}{2} |0\rangle + \frac{1}{2} |1\rangle + \frac{1}{2} |2\rangle + \frac{1}{2} |3\rangle \\ |\varphi_{1}\rangle &= \frac{1}{2} |0\rangle + \frac{i}{2} |1\rangle - \frac{1}{2} |2\rangle - \frac{i}{2} |3\rangle \\ |\varphi_{2}\rangle &= \frac{1}{2} |0\rangle - \frac{1}{2} |1\rangle + \frac{1}{2} |2\rangle - \frac{1}{2} |3\rangle \\ |\varphi_{3}\rangle &= \frac{1}{2} |0\rangle - \frac{i}{2} |1\rangle - \frac{1}{2} |2\rangle + \frac{i}{2} |3\rangle \end{split}$$

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

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

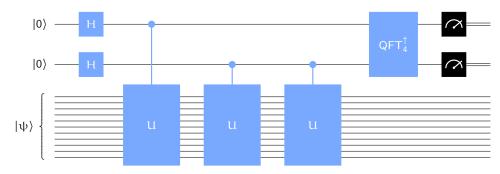
# Two-qubit phase estimation

$$\mathsf{QFT}_4 = \frac{1}{2} \begin{pmatrix} 1 & 1 & 1 & 1 \\ 1 & \mathsf{i} & -1 & -\mathsf{i} \\ 1 & -1 & 1 & -1 \\ 1 & -\mathsf{i} & -1 & \mathsf{i} \end{pmatrix}$$

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

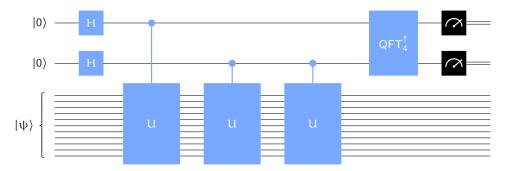
When we think about this matrix as a unitary operation, we call it the quantum Fourier transform.

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

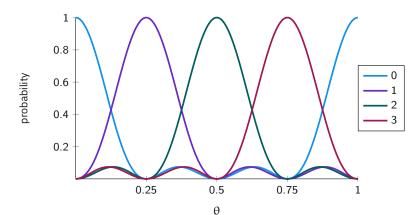


# Two-qubit phase estimation

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  $\theta$ :



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

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

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

### Example

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

$$\mathsf{QFT}_3 = \frac{1}{\sqrt{3}} \begin{pmatrix} 1 & 1 & 1\\ 1 & \frac{-1+\mathrm{i}\sqrt{3}}{2} & \frac{-1-\mathrm{i}\sqrt{3}}{2}\\ 1 & \frac{-1-\mathrm{i}\sqrt{3}}{2} & \frac{-1+\mathrm{i}\sqrt{3}}{2} \end{pmatrix}$$

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

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### Example

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

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

Example

$$\begin{aligned} \mathsf{QFT}_{\mathsf{N}} &= \frac{1}{\sqrt{\mathsf{N}}} \sum_{\mathsf{x}=0}^{\mathsf{N}-1} \sum_{\mathsf{y}=0}^{\mathsf{N}-1} e^{2\pi \mathfrak{i} \frac{\mathsf{x} \mathsf{y}}{\mathsf{N}}} |\mathsf{x}\rangle \langle \mathsf{y}| \\ \\ \mathsf{QFT}_{\mathsf{N}} |\mathsf{y}\rangle &= \frac{1}{\sqrt{\mathsf{N}}} \sum_{\mathsf{x}=0}^{\mathsf{N}-1} e^{2\pi \mathfrak{i} \frac{\mathsf{x} \mathsf{y}}{\mathsf{N}}} |\mathsf{x}\rangle \end{aligned}$$

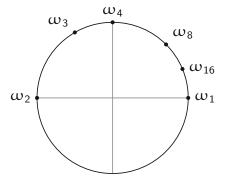
# $\mathsf{QFT}_8 = \frac{1}{2\sqrt{2}} \begin{pmatrix} 1 & 1 & 1 & 1 & 1 & 1 & 1 & 1 \\ 1 & \frac{1+i}{\sqrt{2}} & i & \frac{-1+i}{\sqrt{2}} & -1 & \frac{-1-i}{\sqrt{2}} & -i & \frac{1-i}{\sqrt{2}} \\ 1 & i & -1 & -i & 1 & i & -1 & -i \\ 1 & \frac{-1+i}{\sqrt{2}} & -i & \frac{1+i}{\sqrt{2}} & -1 & \frac{1-i}{\sqrt{2}}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

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$$\mathsf{QFT}_{\mathsf{N}} = \frac{1}{\sqrt{\mathsf{N}}} \sum_{\mathsf{x}=0}^{\mathsf{N}-1} \sum_{\mathsf{y}=0}^{\mathsf{N}-1} e^{2\pi i \frac{\mathsf{x}\,\mathsf{y}}{\mathsf{N}}} |\mathsf{x}\rangle \langle \mathsf{y}| = \frac{1}{\sqrt{\mathsf{N}}} \sum_{\mathsf{x}=0}^{\mathsf{N}-1} \sum_{\mathsf{y}=0}^{\mathsf{N}-1} \omega_{\mathsf{N}}^{\mathsf{x}\,\mathsf{y}} |\mathsf{x}\rangle \langle \mathsf{y}|$$

Useful shorthand notation:

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



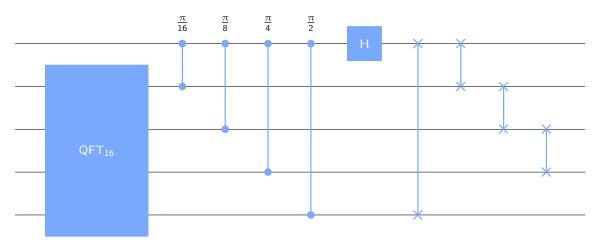
# Circuits for the QFT

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  $\frac{recursive}{recursive}$  in nature. As an example, here is the circuit for QFT<sub>32</sub>:



# Circuits for the QFT

### Cost analysis

Let  $s_m$  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

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

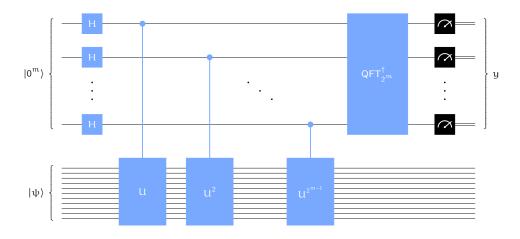
This is a *recurrence relation* with a closed-form solution:

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

### Additional remarks:

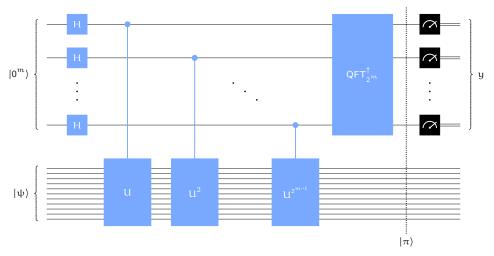
- The number of swap gates can be reduced.
- Approximations to QFT<sub>2</sub><sup>m</sup> can be done at lower cost (and lower depth).

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 *high cost*.



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

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

### Best approximations

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

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

Then the probability to measure y will relatively high:

$$p_{y} \ge \frac{4}{\pi^2} \approx 0.405$$

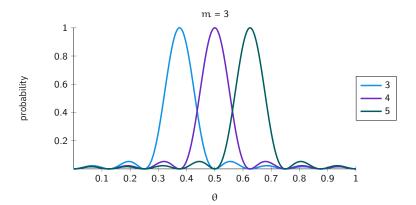
### Worse approximations

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

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$$p_y \leq \frac{1}{4}$$



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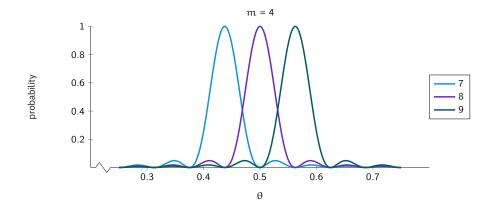
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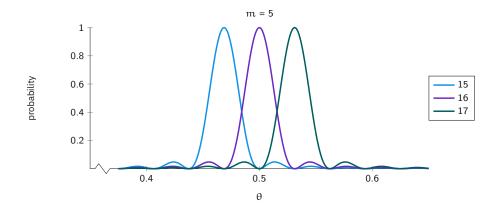
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Then the probability to measure y will be relatively low:

$$p_y \leq \frac{1}{4}$$

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

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

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.

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

### Example

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

# The order-finding problem

### Fact

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

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

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(a, N) = 1!

### Example

Let N = 15 and  $\alpha$  = 2. The operation  $M_{\alpha}$  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 <u>eigenvalues</u> of  $M_a$  are closely connected with the <u>order</u> of 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  $M_a$ :

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

The associated eigenvalue is 1:

$$M_{\alpha}|\psi_{0}\rangle = \frac{|\alpha\rangle + |\alpha^{2}\rangle + \dots + |\alpha^{r}\rangle}{\sqrt{r}} = \frac{|\alpha\rangle + \dots + |\alpha^{r-1}\rangle + |1\rangle}{\sqrt{r}} = |\psi_{0}\rangle$$

To identify more eigenvectors, first recall that

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

This is another eigenvector of  $M_a$ :

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

# 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  $M_a$ :

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

# A convenient eigenvector

$$\begin{split} |\psi_1\rangle &= \frac{|1\rangle + \omega_r^{-1}|\alpha\rangle + \dots + \omega_r^{-(r-1)}|\alpha^{r-1}\rangle}{\sqrt{r}} \\ M_\alpha |\psi_1\rangle &= \omega_r |\psi_1\rangle = e^{2\pi i \frac{1}{r}} |\psi_1\rangle \end{split}$$

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  $j\in\{0,\ldots,r-1\}$ . 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  $j\in\{0,\ldots,r-1\}$ . 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.

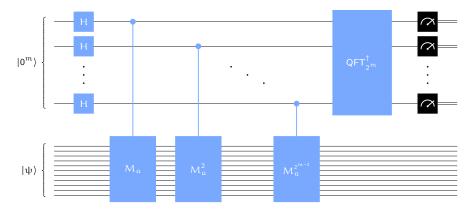
How much precision do we need to correctly determine u/v = j/r?

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

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  $M_\alpha$ . 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  $M_a$  at cost  $O(n^2)$ .

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

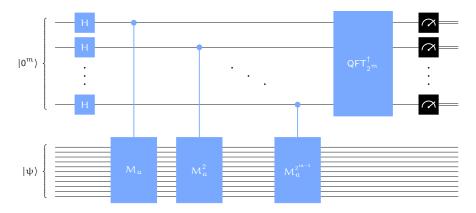
Compute  $b = a^k \pmod{N}$ .

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  $M_\alpha$ . Let's measure the cost as a function of  $n = \lg(N)$ .

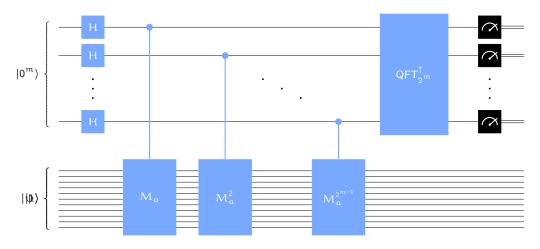


### Cost for phase estimation

- m Hadamard gates: cost O(n)
- m controlled unitary operations: cost  $O(n^3)$
- Quantum Fourier transform: cost  $O(n^2)$

Total cost:  $O(n^3)$ 

# Implementation



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

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

This works because of the following equation:

$$\left|1\right\rangle = \frac{\left|\psi_{0}\right\rangle + \cdots + \left|\psi_{r-1}\right\rangle}{\sqrt{r}}$$

The outcome is the same as if we chose  $j \in \{0, 1, \dots, r-1\}$  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  $d = \gcd(a^{r/2} 1, N)$ . 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  $a^r \equiv 1 \pmod{N}$ .

$$a^r \equiv 1 \pmod{N}$$
  $\iff$  N 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  $(a^{r/2} + 1)$  or  $(a^{r/2} - 1)$ . For a random a, at least one of the prime factors of N is likely to divide  $(a^{r/2} - 1)$ .