# University of Jyväskylä - Course TIEJ6003 Introduction to Quantum Computing

#### Day-05a

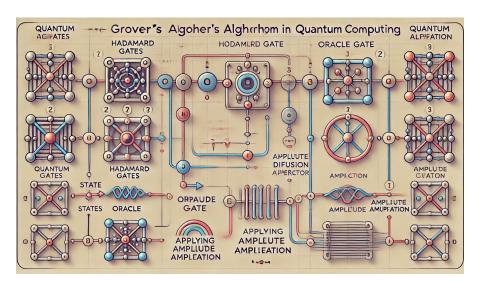
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Summer 2024 Jyväskylä, Finland

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#### GROVER SEARCH ALGORITHM



Summer 2024

## the challenge: unstructured search

We would like to devise a method to locate known objects ('needles') in a large unordered set ('haystack') of N objects.

Suppose that f is a function from  $\{0, 1, ..., N-1\} \rightarrow \{0, 1\}$ , and the target is to locate some x for which f(x) = 1.

For simplicity, assume a single winner w for which f(w) = 1.

Evaluating f(x) for random x will locate w in an expected N/2 steps (alternatively,  $\mathcal{O}(\frac{N}{2})$  steps to find the object with probability 50%).

Grover developed a quantum algorithm to this end and published it in 1996. His method achieves a success-probability greater than 50% in  $\mathcal{O}(\sqrt{N})$  steps!

# introducing the oracle

We assume that we are provided with a black-box that has the ability to *recognize* solutions to the problem.

This assumption is not as strong as it may sound at first - there are many black-box problems in reality whose solutions may be recognized once located. We shall call it the *oracle*.

Intuitively, the oracle enables us to define a Hilbert subspace of winners versus non-winners, and to define a rotation in this subspace toward the winners.

It will be beneficial to work with the uniform superposition state  $|h\rangle = \frac{1}{\sqrt{N}} \sum_{0 \leq x < N} |x\rangle$  (prepared by  $H^{\otimes n} |0\rangle^{\otimes n}$ ).

Then, starting with  $|h\rangle$ , we will repeatedly apply rotations for T iterations and measure!

#### the oracle formalized

The oracle's recognition is signalled via a qubit  $|q\rangle$ , and altogether the oracle's operation may be formally described as follows:

$$\mathcal{U}_{\text{oracle}} |x\rangle |q\rangle = |x\rangle |q \oplus f(x)\rangle.$$
 (1)

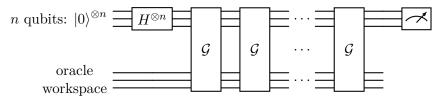
Importantly, it pays off to apply this operator on the superposition state  $(|0\rangle - |1\rangle)/\sqrt{2}$  (similar to Deutsch-Jozsa):

$$\mathcal{U}_{\text{oracle}} |x\rangle \left(\frac{|0\rangle - |1\rangle}{\sqrt{2}}\right) = (-1)^{f(x)} |x\rangle \left(\frac{|0\rangle - |1\rangle}{\sqrt{2}}\right)$$

In practice, the oracle marks the solution by shifting its phase.

Given an unstructured search with M solutions, it turns out that a solution can be located by calling the oracle  $\mathcal{O}(\sqrt{N/M})$  times.

### quantum search: schematic circuit



The Grover operator  $\mathcal{G}$  represents an iteration, which subsequently applies the Oracle followed by the so-called Diffusion operator  $\mathcal{D}$ :

$$\mathcal{G} := \mathcal{U}_{\mathrm{oracle}} \mathcal{D}$$

#### the Grover iteration

The Grover operator  $\mathcal{G}$  is applied repeatedly as a subroutine in each iteration. It can be broken down to four steps:

- 1 Call the oracle  $\mathcal{U}_{oracle}$
- 2 Apply the Hadamard transform  $H^{\otimes n}$
- 3 Conditionally apply phase shift:

if 
$$x > 0$$
 then  $|x\rangle \to -|x\rangle$  else  $|0\rangle \to |0\rangle$ 

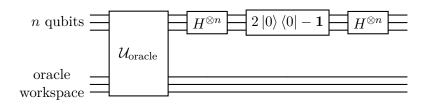
4 Apply the Hadamard transform  $H^{\otimes n}$ 

The combined effect of steps 2+3+4 is the reflection on the uniform superposition  $|h\rangle$  (the Diffusion operator  $\mathcal{D}$ ):

$$H^{\otimes n}\left(2\left|0\right\rangle^{\otimes n}\left\langle 0\right|^{\otimes n}-I^{\otimes n}\right)H^{\otimes n}=2\left|h\right\rangle\left\langle h\right|-I^{\otimes n}=\mathcal{D}$$

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# Grover iteration: concrete circuit per $\mathcal{G}$



$$\begin{split} \mathcal{G} &:= \quad \mathcal{D}\mathcal{U}_{\text{oracle}} \\ \mathcal{U}_{\text{oracle}} &:= \quad \mathbf{1} - 2 \left| w \right\rangle \left\langle w \right| \\ \mathcal{D} &:= \quad 2 \left| h \right\rangle \left\langle h \right| - \mathbf{1} \end{split}$$

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## subspace framework

Let S denote the set of objects we are searching for, and let  $M \geq 1$  be its cardinality. S is the solution set ("winners"), and its elements are called the solutions.

We denote the set of objects that are not a solution ("non-winners") by

$$S^{\perp} := \{0, 1, \dots, N-1\} \setminus S$$

and accordingly construct two vectors:

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$$|\Psi_S\rangle := \frac{1}{\sqrt{M}} \sum_{x \in S} |x\rangle \qquad |\Psi_{S^{\perp}}\rangle := \frac{1}{\sqrt{N-M}} \sum_{x \in S^{\perp}} |x\rangle \qquad (2)$$

#### projections

By considering the two projection operators onto the two subspaces

$$P_S := \sum_{x \in S} |x\rangle \langle x| \qquad P_{S^{\perp}} := \sum_{x \in S^{\perp}} |x\rangle \langle x| = \mathbf{1} - P_S, \tag{3}$$

and by recalling that  $S \cup S^{\perp} = \{0, 1, \dots, N-1\}$ , every state can be expanded in the computational basis as follows

$$|\psi\rangle = (P_S + P_{S^{\perp}})|\psi\rangle = \sum_{x \in S} \psi_x |x\rangle + \sum_{x \in S^{\perp}} \psi_x |x\rangle = \alpha |\Psi_S\rangle + \beta |\Psi_{S^{\perp}}\rangle$$
(4)

In particular, the uniform superposition state  $|h\rangle$  can be expanded as:

$$|h\rangle = \sqrt{\frac{M}{N}} |\Psi_S\rangle + \sqrt{\frac{N-M}{N}} |\Psi_{S^{\perp}}\rangle$$
 (5)

### interpretation: reflections

By observing the 2D subspace spanned by  $|\Psi_S\rangle$  ("the good") and  $|\Psi_{S^{\perp}}\rangle$  ("the bad"), Grover's iteration may be interpreted as two reflection operations when starting with  $|h\rangle$ :

- 1  $\mathcal{U}_{\text{oracle}}$  performs a reflection about the vector  $|\Psi_{S^{\perp}}\rangle$ .
- 2  $\mathcal{D}$  performs a reflection about the vector  $|h\rangle$ .

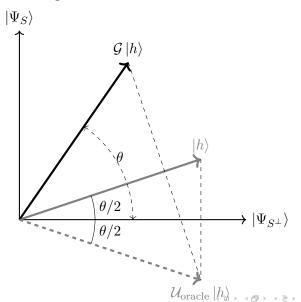
Importantly, the repeated application of  $\mathcal{G}$  always keeps the states in the plane defined by  $|\Psi_S\rangle$  and  $|\Psi_{S^{\perp}}\rangle$ .

Altogether,  $\mathcal{G}$  yields a rotation defined as

$$\mathcal{G} := \begin{bmatrix} \cos \theta & -\sin \theta \\ \sin \theta & \cos \theta \end{bmatrix},\tag{6}$$

with  $\theta$  satisfying  $\sin \theta = 2\sqrt{M(N-M)}/N$ .

# geometric visualization



## Grover formalized per M=1

**Inputs:** (1) a black-box oracle  $\mathcal{U}_{\text{oracle}}$  which performs the operation  $\mathcal{U}_{\text{oracle}} |x\rangle |q\rangle = |x\rangle |q \oplus f(x)\rangle.$ 

(2) (n+1) qubits in the state  $|0\rangle$ .

Outputs:  $x_0$ .

**Runtime:**  $\mathcal{O}(\sqrt{2^n})$  operations. Succeeds with probability  $\mathcal{O}(1)$ .

Procedure: 
$$|\psi_0\rangle = |0\rangle^{\otimes n} |0\rangle$$

$$|\psi_1\rangle = \frac{1}{\sqrt{2^n}} \sum_{x=0}^{2^n - 1} |x\rangle \left[ \frac{|0\rangle - |1\rangle}{\sqrt{2}} \right]$$

$$\rightarrow \left[ \mathcal{D}\mathcal{U}_{\text{oracle}} \right]^T |\psi_1\rangle \approx |x_0\rangle \left[ \frac{|0\rangle - |1\rangle}{\sqrt{2}} \right]$$

$$\rightarrow x_0$$

# Grover: complexity and aftermath

Now that the search is reduced to rotations in the plane of  $\{|\Psi_S\rangle, |\Psi_{S^{\perp}}\rangle\}$ , the algorithm's complexity is reduced to the question

"how many radians are needed to approach  $|\Psi_S\rangle$ ?"

Given the starting point  $|h\rangle = \sqrt{\frac{M}{N}} |\Psi_S\rangle + \sqrt{\frac{N-M}{N}} |\Psi_{S^{\perp}}\rangle$ , rotating it through  $\arccos(\sqrt{M/N})$  radians will drive the system to  $|\Psi_S\rangle$ :

$$T = \text{round}\left(\frac{\arccos(\sqrt{M/N})}{\theta}\right) \le \left[\frac{\pi}{4}\sqrt{\frac{N}{M}}\right].$$
 (7)

 $\mathcal{O}(\sqrt{N/M})$  Grover iterations (and thus oracle calls) are required, versus  $\mathcal{O}(N/M)$  oracle calls that are required classically.