A LECTURE ON

GROVER'S QUANTUM SEARCH ALGORITHM VERSION 1.1

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ABSTRACT. This paper is a written version of a one hour lecture given on Lov Grover's quantum database search algorithm. It is based on [4], [5], and [9].

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1. Problem definition

We consider the problem of searching an unstructured database of $N=2^n$ records for exactly one record which has been specifically marked. This can be rephrased in mathematical terms as an oracle problem as follows:

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Label the records of the database with the integers

$$0,1,2,\ldots,N-1$$
,

and denote the label of the unknown marked record by x_0 . We are given an oracle which computes the n bit binary function

$$f: \{0,1\}^n \longrightarrow \{0,1\}$$

defined by

$$f(x) = \begin{cases} 1 & \text{if } x = x_0 \\ 0 & \text{otherwise} \end{cases}$$

We remind the readers that, as a standard oracle idealization, we have no access to the internal workings of the function f. It operates simply as a blackbox function, which we can query as many times as we like. But with each such a query comes an associated computational cost.

Search Problem for an Unstructured Database. Find the record labeled as x_0 with the minimum amount of computational work, i.e., with the minimum number of queries of the oracle f.

From probability theory, we know that if we examine k records, i.e., if we compute the oracle f for k randomly chosen records, then the probability of finding the record labeled as x_0 is k/N. Hence, on a classical computer it takes $O(N) = O(2^n)$ queries to find the record labeled x_0 .

2. The quantum mechanical perspective

However, as Lov Grover so a stutely observed, on a quantum computer the search of an unstructured data base can be accomplished in $O(\sqrt{N})$ steps, or more precisely, with the application of $O(\sqrt{N} \lg N)$ sufficiently local unitary transformations. Although this is not exponentially faster, it is a significant speedup.

Let \mathcal{H}_2 be a 2 dimensional Hilbert space with orthonormal basis

$$\{|0\rangle, |1\rangle\}$$
;

and let

$$\{|0\rangle, |1\rangle, \ldots, |N-1\rangle\}$$

denote the induced orthonormal basis of the Hilbert space

$$\mathcal{H} = \bigotimes_{0}^{N-1} \mathcal{H}_2$$
.

From the quantum mechanical perspective, the oracle function f is given as a blackbox unitary transformation U_f , i.e., by

$$\mathcal{H}\otimes\mathcal{H}_2 \stackrel{U_f}{\longrightarrow} \mathcal{H}\otimes\mathcal{H}_2$$

$$|x\rangle \otimes |y\rangle \longmapsto |x\rangle \otimes |f(x) \oplus y\rangle$$

where ' \oplus ' denotes exclusive 'OR', i.e., addition modulo $2.^1$

Instead of U_f , we will use the computationally equivalent unitary transformation

$$I_{|x_0\rangle}\left(|x\rangle\right) = (-1)^{f(x)} |x\rangle = \begin{cases} -|x_0\rangle & \text{if } x = x_0 \\ |x\rangle & \text{otherwise} \end{cases}$$

That $I_{|x_0\rangle}$ is computationally equivalent to U_f follows from the easily verifiable fact that

$$U_f\left(|x\rangle\otimes\frac{|0\rangle-|1\rangle}{\sqrt{2}}\right) = \left(I_{|x_0\rangle}\left(|x\rangle\right)\right)\otimes\frac{|0\rangle-|1\rangle}{\sqrt{2}}\ ,$$

and also from the fact that U_f can be constructed from a controlled- $I_{|x_0\rangle}$ and two one qubit Hadamard transforms. (For details, please refer to [10], [11].)

The unitary transformation $I_{|x_0\rangle}$ is actually an **inversion** [1] in \mathcal{H} about the hyperplane perpendicular to $|x_0\rangle$. This becomes evident when $I_{|x_0\rangle}$ is rewritten in the form

$$I_{|x_0\rangle} = I - 2|x_0\rangle \langle x_0| ,$$

where 'I' denotes the identity transformation. More generally, for any unit length ket $|\psi\rangle$, the unitary transformation

$$I_{|\psi\rangle} = I - 2 |\psi\rangle \langle \psi|$$

is an inversion in \mathcal{H} about the hyperplane orthogonal to $|\psi\rangle$.

¹Please note that $U_f = (\nu \circ \iota)(f)$, as defined in sections 10.3 and 10.4 of [12].

3. Properties of the inversion $I_{|\psi\rangle}$

We digress for a moment to discuss the properties of the unitary transformation $I_{|\psi\rangle}$. To do so, we need the following definition.

Definition 1. Let $|\psi\rangle$ and $|\chi\rangle$ be two kets in \mathcal{H} for which the bracket product $\langle\psi \mid \chi\rangle$ is a real number. We define

$$\mathcal{S}_{\mathbb{C}} = Span_{\mathbb{C}}(|\psi\rangle, |\chi\rangle) = \{\alpha |\psi\rangle + \beta |\chi\rangle \in \mathcal{H} \mid \alpha, \beta \in \mathbb{C}\}$$

as the sub-Hilbert space of \mathcal{H} spanned by $|\psi\rangle$ and $|\chi\rangle$. We associate with the Hilbert space $\mathcal{S}_{\mathbb{C}}$ a real inner product space lying in $\mathcal{S}_{\mathbb{C}}$ defined by

$$S_{\mathbb{R}} = Span_{\mathbb{R}}(|\psi\rangle, |\chi\rangle) = \{a|\psi\rangle + b|\chi\rangle \in \mathcal{H} \mid a, b \in \mathbb{R}\}$$
,

where the inner product on $S_{\mathbb{R}}$ is that induced by the bracket product on \mathcal{H} . If $|\psi\rangle$ and $|\chi\rangle$ are also linearly independent, then $S_{\mathbb{R}}$ is a 2 dimensional real inner product space (i.e., the 2 dimensional Euclidean plane) lying inside of the complex 2 dimensional space $S_{\mathbb{C}}$.

Proposition 1. Let $|\psi\rangle$ and $|\chi\rangle$ be two linearly independent unit length kets in \mathcal{H} with real bracket product; and let $\mathcal{S}_{\mathbb{C}} = Span_{\mathbb{C}}(|\psi\rangle, |\chi\rangle)$ and $\mathcal{S}_{\mathbb{R}} = Span_{\mathbb{R}}(|\psi\rangle, |\chi\rangle)$. Then

1) Both $S_{\mathbb{C}}$ and $S_{\mathbb{R}}$ are invariant under the transformations $I_{|\psi\rangle}$, $I_{|\chi\rangle}$, and hence $I_{|\psi\rangle} \circ I_{|\chi\rangle}$, i.e.,

- 2) If $L_{|\psi^{\perp}\rangle}$ is the line in the plane $\mathcal{S}_{\mathbb{R}}$ which passes through the origin and which is perpendicular to $|\psi\rangle$, then $I_{|\psi\rangle}$ restricted to $\mathcal{S}_{\mathbb{R}}$ is a reflection in (i.e., a Möbius inversion [1] about) the line $L_{|\psi^{\perp}\rangle}$. A similar statement can be made in regard to $|\chi\rangle$.
- 3) If $|\psi^{\perp}\rangle$ is a unit length vector in $\mathcal{S}_{\mathbb{R}}$ perpendicular to $|\psi\rangle$, then

$$-I_{|\psi
angle}=I_{\left|\psi^{\perp}
ight
angle}$$
 .

(Hence, $\langle \psi^{\perp} | \chi \rangle$ is real.)

Finally we note that, since $I_{|\psi\rangle} = I - 2 |\psi\rangle \langle \psi|$, it follows that

Proposition 2. If $|\psi\rangle$ is a unit length ket in \mathcal{H} , and if U is a unitary transformation on \mathcal{H} , then

$$UI_{|\psi\rangle}U^{-1} = I_{U|\psi\rangle}$$
.

4. The method in Lov's "madness"

Let $H: \mathcal{H} \longrightarrow \mathcal{H}$ be the Hadamard transform, i.e.,

$$H = \bigotimes_{0}^{n-1} H^{(2)} ,$$

where

$$H^{(2)} = \left(\begin{array}{cc} 1 & 1\\ 1 & -1 \end{array}\right)$$

with respect to the basis $|0\rangle$, $|1\rangle$.

We begin by using the Hadamard transform H to construct a state $|\psi_0\rangle$ which is an equal superposition of all the standard basis states $|0\rangle$, $|1\rangle, \ldots, |N-1\rangle$ (including the unknown state $|x_0\rangle$), i.e.,

$$|\psi_0\rangle = H|0\rangle = \frac{1}{\sqrt{N}} \sum_{k=0}^{N-1} |k\rangle$$
.

Both $|\psi_0\rangle$ and the unknown state $|x_0\rangle$ lie in the Euclidean plane $\mathcal{S}_{\mathbb{R}} = Span_{\mathbb{R}}(|\psi_0\rangle, |x_0\rangle)$. Our strategy is to rotate within the plane $\mathcal{S}_{\mathbb{R}}$ the state $|\psi_0\rangle$ about the origin until it is as close as possible to $|x_0\rangle$. Then a measurement with respect to the standard basis of the state resulting from rotating $|\psi_0\rangle$, will produce $|x_0\rangle$ with high probability.

To achieve this objective, we use the oracle $I_{|x_0\rangle}$ to construct the unitary transformation

$$Q = -HI_{|0\rangle}H^{-1}I_{|x_0\rangle} ,$$

which by proposition 2 above, can be reexpressed as

$$Q = -I_{|\psi_0\rangle}I_{|x_0\rangle} .$$

Let $|x_0^{\perp}\rangle$ and $|\psi_0^{\perp}\rangle$ denote unit length vectors in $\mathcal{S}_{\mathbb{R}}$ perpendicular to $|x_0\rangle$ and $|\psi_0\rangle$, respectively. There are two possible choices for each of $|x_0^{\perp}\rangle$

and $|\psi_0^{\perp}\rangle$ respectively. To remove this minor, but nonetheless annoying, ambiguity, we select $|x_0^{\perp}\rangle$ and $|\psi_0^{\perp}\rangle$ so that the orientation of the plane $\mathcal{S}_{\mathbb{R}}$ induced by the ordered spanning vectors $|\psi_0\rangle$, $|x_0\rangle$ is the same orientation as that induced by each of the ordered bases $|x_0^{\perp}\rangle$, $|x_0\rangle$ and $|\psi_0\rangle$, $|\psi_0^{\perp}\rangle$. (Please refer to Figure 2.)

Remark 1. The removal of the above ambiguities is really not essential. However, it does simplify the exposition given below.

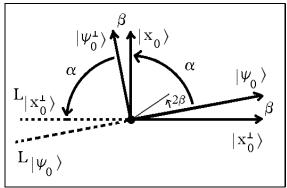


Figure 2. The linear transformation $Q|_{\mathcal{S}_{\mathbb{R}}}$ is reflection in the line $L_{|x_0^{\perp}\rangle}$ followed by reflection in the line $L_{|\psi_0\rangle}$ which is the same as rotation by the angle 2β . Thus, $Q|_{\mathcal{S}_{\mathbb{R}}}$ rotates $|\psi_0\rangle$ by the angle 2β toward $|x_0\rangle$.

We proceed by noting that, by the above proposition 1, the plane $\mathcal{S}_{\mathbb{R}}$ lying in \mathcal{H} is invariant under the linear transformation Q, and that, when Q is restricted to the plane $\mathcal{S}_{\mathbb{R}}$, it can be written as the composition of two inversions, i.e.,

$$Q|_{\mathcal{S}_{\mathbb{R}}} = I_{|\psi_0^{\perp}\rangle} I_{|x_0\rangle}$$
.

In particular, $Q|_{\mathcal{S}_{\mathbb{R}}}$ is the composition of two inversions in $\mathcal{S}_{\mathbb{R}}$, the first in the line $L_{|x_0^{\perp}\rangle}$ in $\mathcal{S}_{\mathbb{R}}$ passing through the origin having $|x_0\rangle$ as normal, the second in the line $L_{|\psi_0\rangle}$ through the origin having $|\psi_0^{\perp}\rangle$ as normal.²

We can now apply the following theorem from plane geometry:

²The line $L_{\left|x_{0}^{\perp}\right\rangle}$ is the intersection of the plane $\mathcal{S}_{\mathbb{R}}$ with the hyperplane in \mathcal{H} orthogonal to $\left|x_{0}\right\rangle$. A similar statement can be made in regard to $L_{\left|\psi_{0}\right\rangle}$.

Theorem 1. If L_1 and L_2 are lines in the Euclidean plane \mathbb{R}^2 intersecting at a point O; and if β is the angle in the plane from L_1 to L_2 , then the operation of reflection in L_1 followed by reflection in L_2 is just rotation by angle 2β about the point O.

Let β denote the angle in $S_{\mathbb{R}}$ from $L_{\left|x_{0}^{\perp}\right\rangle}$ to $L_{\left|\psi_{0}\right\rangle}$, which by plane geometry is the same as the angle from $\left|x_{0}^{\perp}\right\rangle$ to $\left|\psi_{0}\right\rangle$, which in turn is the same as the angle from $\left|x_{0}\right\rangle$ to $\left|\psi_{0}^{\perp}\right\rangle$. Then by the above theorem $Q|_{\mathcal{S}_{\mathbb{R}}}=I_{\left|\psi_{0}^{\perp}\right\rangle}I_{\left|x_{0}\right\rangle}$ is a rotation about the origin by the angle 2β .

The key idea in Grover's algorithm is to move $|\psi_0\rangle$ toward the unknown state $|x_0\rangle$ by successively applying the rotation Q to $|\psi_0\rangle$ to rotate it around to $|x_0\rangle$. This process is called **amplitude amplification**. Once this process is completed, the measurement of the resulting state (with respect to the standard basis) will, with high probability, yield the unknown state $|x_0\rangle$. This is the essence of Grover's algorithm.

But how many times K should we apply the rotation Q to $|\psi_0\rangle$? If we applied Q too many or too few times, we would over- or undershoot our target state $|x_0\rangle$.

We determine the integer K as follows:

Since

$$|\psi_0\rangle = \sin\beta |x_0\rangle + \cos\beta |x_0^{\perp}\rangle$$
,

the state resulting after k applications of Q is

$$|\psi_k\rangle = Q^k |\psi_0\rangle = \sin\left[(2k+1)\beta\right] |x_0\rangle + \cos\left[(2k+1)\beta\right] |x_0^{\perp}\rangle$$
.

Thus, we seek to find the smallest positive integer K = k such that

$$\sin\left[\left(2k+1\right)\beta\right]$$

is as close as possible to 1. In other words, we seek to find the smallest positive integer K=k such that

$$(2k+1)\beta$$

is as close as possible to $\pi/2$. It follows that³

$$K = k = round \left(\frac{\pi}{4\beta} - \frac{1}{2}\right) ,$$

where "round" is the function that rounds to the nearest integer.

We can determine the angle β by noting that the angle α from $|\psi_0\rangle$ and $|x_0\rangle$ is complementary to β , i.e.,

$$\alpha + \beta = \pi/2$$
,

and hence,

$$\frac{1}{\sqrt{N}} = \langle x_0 \mid \psi_0 \rangle = \cos \alpha = \cos(\frac{\pi}{2} - \beta) = \sin \beta .$$

Thus, the angle β is given by

$$\beta = \sin^{-1}\left(\frac{1}{\sqrt{N}}\right) \approx \frac{1}{\sqrt{N}}$$
 (for large N),

and hence,

$$K = k = round \left(\frac{\pi}{4 \sin^{-1} \left(\frac{1}{\sqrt{N}} \right)} - \frac{1}{2} \right) \approx round \left(\frac{\pi}{4} \sqrt{N} - \frac{1}{2} \right) \text{ (for large } N \text{)}.$$

5. Summary of Grover's algorithm

In summary, we provide the following outline of Grover's algorithm:

 $^{^3}$ The reader may prefer to use the floor function instead of the round function.

Grover's Algorithm

STEP 0. (Initialization)

$$|\psi\rangle \longleftarrow H |0\rangle = \frac{1}{\sqrt{N}} \sum_{j=0}^{N-1} |j\rangle$$

STEP 1. Loop until $k = round\left(\frac{\pi}{4\sin^{-1}(1/\sqrt{N})} - \frac{1}{2}\right) \approx round\left(\frac{\pi}{4}\sqrt{N} - \frac{1}{2}\right)$

$$|\psi\rangle \longleftarrow Q |\psi\rangle = -HI_{|0\rangle}HI_{|x_0\rangle} |\psi\rangle$$

 $k \longleftarrow k+1$

STEP 2. Measure $|\psi\rangle$ with respect to the standard basis $|0\rangle, |1\rangle, \ldots, |N-1\rangle$ to obtain the marked unknown state $|x_0\rangle$ with probability $\geq 1 - \frac{1}{N}$.

We complete our summary with the following theorem:

Theorem 2. With a probability of error⁴

$$Prob_E \leq \frac{1}{N}$$
,

Grover's algorithm finds the unknown state $|x_0\rangle$ at a computational cost of

$$O\left(\sqrt{N}\lg N\right)$$

Proof.

Part 1. The probability of error $Prob_E$ of finding the hidden state $|x_0\rangle$ is given by

$$Prob_E = \cos^2\left[\left(2K+1\right)\beta\right] ,$$

where

$$\begin{cases} \beta = \sin^{-1}\left(\frac{1}{\sqrt{N}}\right) \\ K = round\left(\frac{\pi}{4\beta} - \frac{1}{2}\right) \end{cases},$$

⁴If the reader prefers to use the *floor* function rather than the *round* function, then probability of error becomes $Prob_E \leq \frac{4}{N} - \frac{4}{N^2}$.

where "round" is the function that rounds to the nearest integer. Hence,

$$\frac{\pi}{4\beta} - 1 \le K \le \frac{\pi}{4\beta} \implies \frac{\pi}{2} - \beta \le (2K+1)\beta \le \frac{\pi}{2} + \beta$$

$$\implies \sin \beta = \cos \left(\frac{\pi}{2} - \beta\right) \ge \cos \left[\left(2K+1\right)\beta\right] \ge \cos \left(\frac{\pi}{2} + \beta\right) = -\sin \beta$$

Thus,

$$Prob_E = \cos^2\left[\left(2K + 1\right)\beta\right] \le \sin^2\beta = \sin^2\left(\sin^{-1}\left(\frac{1}{\sqrt{N}}\right)\right) = \frac{1}{N}$$

Part 2. The computational cost of the Hadamard transform $H = \bigotimes_0^{n-1} H^{(2)}$ is $O(n) = O(\lg N)$ single qubit operations. The transformations $-I_{|0\rangle}$ and $I_{|x_0\rangle}$ each carry a computational cost of O(1). STEP 1 is the computationally dominant step. In STEP 1 there are $O\left(\sqrt{N}\right)$ iterations. In each iteration, the Hadamard transform is applied twice. The transformations $-I_{|0\rangle}$ and $I_{|x_0\rangle}$ are each applied once. Hence, each iteration comes with a computational cost of $O(\lg N)$, and so the total cost of STEP 1 is $O(\sqrt{N} \lg N)$.

6. An example of Grover's algorithm

As an example, we search a database consisting of $N = 2^n = 8$ records for an unknown record with the unknown label $x_0 = 5$. The calculations for this example were made with OpenQuacks, which is an open source quantum simulator Maple package developed at UMBC and publically available.

We are given a blackbox computing device

$$\operatorname{In} \to \boxed{I_{\mid ? \rangle}} \to \operatorname{Out}$$

that implements as an oracle the unknown unitary transformation

We cannot open up the blackbox $\to I_{|?\rangle}$ \to to see what is inside. So we do not know what $I_{|x_0\rangle}$ and x_0 are. The only way that we can glean some information about x_0 is to apply some chosen state $|\psi\rangle$ as input, and then make use of the resulting output.

Using of the blackbox \to $\boxed{I_{|?\rangle}}$ \to as a component device, we construct a computing device \to $\boxed{-HI_{|0\rangle}HI_{|?\rangle}}$ \to which implements the unitary operator

We do not know what unitary transformation Q is implemented by the device $\rightarrow \boxed{-HI_{|0\rangle}HI_{|?\rangle}}$ \rightarrow because the blackbox $\rightarrow \boxed{I_{|?\rangle}}$ \rightarrow is one of its essential components.

 $\mathbb{STEP} 0$. We begin by preparing the known state

$$|\psi_0\rangle = H|0\rangle = \frac{1}{\sqrt{8}}(1, 1, 1, 1, 1, 1, 1, 1)^{transpose}$$

STEP 1. We proceed to loop

$$K = round\left(\frac{\pi}{4\sin^{-1}\left(1/\sqrt{8}\right)} - \frac{1}{2}\right) = 2$$

times in STEP 1.

ITERATION 1. On the first iteration, we obtain the unknown state

$$|\psi_1\rangle = Q|\psi_0\rangle = \frac{1}{4\sqrt{2}}(1,1,1,1,5,1,1,1)^{transpose}$$

ITERATION 2. On the second iteration, we obtain the unknown state

$$|\psi_2\rangle = Q |\psi_1\rangle = \frac{1}{8\sqrt{2}} (-1, -1, -1, -1, -1, -1, -1, -1)^{transpose}$$

and branch to $\mathbb{STEP}\ 2$

STEP 2. We measure the unknown state $|\psi_2\rangle$ to obtain either

 $|5\rangle$

with probability

$$Prob_{Success} = \sin^2((2K+1)\beta) = \frac{121}{128} = 0.9453$$

or some other state with probability

$$Prob_{Failure} = \cos^2((2K+1)\beta) = \frac{7}{128} = 0.0547$$

and then exit.

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