

Lecture 2: Collision finding

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1 Quantum walk search

We can summarize the (query) cost of the QW search algorithm using the following quantities:

- “Setup cost” \mathcal{S} : the number of queries needed to create the initial state $|\pi\rangle$ in step 1.
- “Checking cost” \mathcal{C} : the number of queries needed to check whether an element is marked in step 2.(a).
- “Update cost” \mathcal{U} : the number of queries needed to implement a step of the QW.

The total cost (up to constants) is then

$$\mathcal{S} + \sqrt{\frac{n}{m}} \left(\mathcal{C} + \frac{1}{\sqrt{\delta}} \mathcal{U} \right).$$

This algorithm is called the “MNRS algorithm”, after Magniez-Nayak-Roland-Santha [MNRS07].

2 Collision finding with quantum walk search

Assume that we are given an array of integers x_1, x_2, \dots, x_N . A *collision* is a pair of distinct i, j such that $x_i = x_j$. How many elements do we have to query in order to find a collision (or decide that no collision exists)? Classically this essentially requires to query the full array, and so the classical query complexity is $\Omega(N)$. In contrast, using an algorithm based on *quantum walk search* we can find a collision with a *sublinear* number of queries.¹

The quantum walk algorithm for collision finding was proposed by Ambainis [Amb07]. The algorithm runs quantum walk search over elements or “words” $\mathcal{Y} = (Y, x_Y)$, consisting of (i) a size- k subset $Y \subseteq [N]$, and (ii) the list x_Y of integers x_j with index $j \in Y$. We call an element \mathcal{Y} marked if Y contains both indices of a collision (equivalently, x_Y must contain a collision).

Exercise 1. Let $n = \binom{N}{k}$ denote the number of elements and m the number of marked elements. Show that $m/n \in \Omega(k^2/N^2)$.

To use quantum walk search, we consider a graph G with vertex set V indexed by the elements \mathcal{Y} . There is an edge between $\mathcal{Y} = (Y, x_Y)$ and $\mathcal{Y}' = (Y', x_{Y'})$ if the subsets Y and Y' differ in exactly one element (i.e., we can obtain Y' from Y by replacing one index). The resulting graph G has $n = \binom{N}{k}$ vertices and is $k(n - k)$ -regular. It corresponds to a so-called *Johnson graph*, and one can show that its spectral gap is $\delta \in \Omega(1/k)$ when $k \ll n$.

¹While we focus on query complexity for ease of exposition, all algorithms can be implemented with a similar runtime.

A star state $|\psi_{\mathcal{Y}}\rangle$ centered on a vertex \mathcal{Y} of G is given by the state

$$|\psi_{\mathcal{Y}}\rangle = \frac{1}{\sqrt{k(n-k)}} \sum_{\mathcal{Y}' \sim \mathcal{Y}} |\mathcal{Y}, \mathcal{Y}'\rangle,$$

where the sum runs over neighboring elements \mathcal{Y}' of \mathcal{Y} . The quantum walk search algorithm then starts from the uniform superposition

$$|\pi\rangle = \frac{1}{\sqrt{n}} \sum_{\mathcal{Y}} |\psi_{\mathcal{Y}}\rangle,$$

and the algorithm has cost

$$\mathcal{S} + \frac{N}{k}(\sqrt{k}\mathcal{U} + \mathcal{C}),$$

where \mathcal{U} is the *update cost* or cost of implementing a single quantum walk step on the Johnson graph G . We now bound the different costs.

For the checking cost \mathcal{C} , note that we can check whether a given state $|\psi_{\mathcal{Y}}\rangle$ is marked simply by checking whether the list $x_{\mathcal{Y}}$ contains a collision. Since this list is given explicitly in the description of $|\psi_{\mathcal{Y}}\rangle$, this requires no queries and so $\mathcal{C} = 0$.

The setup cost \mathcal{S} amounts to creating the state $|\pi\rangle = \frac{1}{\sqrt{n}} \sum_{\mathcal{Y}} |\psi_{\mathcal{Y}}\rangle$. We do this in a few steps. First, we prepare the state $\frac{1}{\sqrt{n}} \sum_{\mathcal{Y}} |\mathcal{Y}\rangle |0\rangle$. This takes k queries. Then, we construct the mapping U_{ψ} defined by $U_{\psi} |\mathcal{Y}\rangle |0\rangle = |\psi_{\mathcal{Y}}\rangle$. We do this in two steps:

$$\begin{aligned} |\mathcal{Y}\rangle |0\rangle &= |Y, x_Y\rangle |0\rangle \xrightarrow{(i)} \frac{1}{\sqrt{k(n-k)}} \sum_{\mathcal{Y}' \sim \mathcal{Y}} |Y, x_Y\rangle |Y', 0\rangle \\ &\xrightarrow{(ii)} \frac{1}{\sqrt{k(n-k)}} \sum_{\mathcal{Y}' \sim \mathcal{Y}} |Y, x_Y\rangle |Y', x_{Y'}\rangle = |\psi_{\mathcal{Y}}\rangle. \end{aligned}$$

Step (i) requires no queries. Step (ii) amounts to gathering the elements $x_{Y'}$ with index in Y' . Since $x_{Y'}$ contains exactly one element not in x_Y , this requires exactly one query. The setup cost \mathcal{S} is hence roughly k .

Finally, we bound the cost \mathcal{U} of a single call to the quantum walk operator W . Recall from last lecture that $W = S \cdot C$ where S is a simple swap (i.e., $S |\mathcal{Y}, \mathcal{Y}'\rangle = |\mathcal{Y}', \mathcal{Y}\rangle$), requiring no queries, and $C = 2(\sum_{\mathcal{Y}} |\psi_{\mathcal{Y}}\rangle \langle \psi_{\mathcal{Y}}|) - I$ is a reflection around the star subspace. Using a similar trick as before, we can implement the reflection C by making two calls to the preparation operator U_{ψ} , which requires a single query. This proves that the update cost \mathcal{U} is $O(1)$.

Exercise 2. Verify that $C = U_{\psi} R_0 U_{\psi}^{\dagger}$.

Combining these different arguments, we can bound the total cost by

$$\mathcal{S} + \sqrt{\frac{n}{m}} \left(\frac{1}{\sqrt{\delta}} \mathcal{U} + \mathcal{C} \right) \approx k + \frac{N}{\sqrt{k}}.$$

If we set $k = N^{2/3}$ then this yields a quantum algorithm for collision finding with complexity $\tilde{O}(N^{2/3})$. This is essentially optimal by the $\Omega(N^{2/3})$ lower bound of Aaronson and Shi [AS04].

References

- [Amb07] Andris Ambainis. Quantum walk algorithm for element distinctness. *SIAM Journal on Computing*, 37(1):210–239, 2007.
- [AS04] Scott Aaronson and Yaoyun Shi. Quantum lower bounds for the collision and the element distinctness problems. *Journal of the ACM (JACM)*, 51(4):595–605, 2004.
- [MNRS07] Frédéric Magniez, Ashwin Nayak, Jérémie Roland, and Miklos Santha. Search via quantum walk. In *Proceedings of the thirty-ninth annual ACM symposium on Theory of computing*, pages 575–584, 2007.