# Quantum Algorithm for Linear Systems of Equations

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- 1) Preliminary Concepts and Classical Algorithms.
- 2) Overview of Quantum Algorithm
- 3) Some Details of Quantum Algorithm

### References

- 1) A.W. Harrow, A. Hassidim, S. Lloyd, PRL **103**, 150502 (2009).
- 2) arXiv:0811.3171v3 [quant-ph].
- 3) http://www.multimedia.ethz.ch/conferen ces/2010/qip/?doi=10.3930/ETHZ/AV-37c9a9ce-428f-40c7-964e-6a20c9ec9757
- 4) http://www.youtube.com/watch?v=KtIPA PyaPOg

## Introduction

Let A denote an N by N matrix and  $\mathbf{b}$  be a vector. Then we want to solve for  $\mathbf{x}$ 

$$A \mathbf{x} = \mathbf{b}$$

This is a very important subroutine in all kinds of calculations.

One class of algorithms to solve for x is Gaussian elimination. The complexity is  $O(N^3)$ .

## Introduction

If *A* is hermitian, positive-definite and sparse, then it is much faster to use iterative methods.

Try to iteratively find x that minimizes

$$f(\mathbf{x}) = \frac{1}{2}\mathbf{x}^{\dagger}A\mathbf{x} - \mathbf{b}^{\dagger}\mathbf{x}$$

A unique global minimum is guaranteed to exist.

## Introduction

One of the best algorithm is conjugate gradient method. The complexity is

$$O\left(\kappa \log(1/\epsilon) \cdot Ns\right)$$

The algorithm takes  $O\left(\kappa \log(1/\epsilon)\right)$  iterations. Each iteration takes  $O\left(Ns\right)$ .

- s -- Sparsity of each row.
- $\kappa$  -- Condition number.
- $\epsilon$  -- Error bound.  $\|\mathbf{x} A^{-1}\mathbf{b}\|^2 \leqslant \epsilon$ .

The vector **b** is represented as a ket

$$|b
angle = \sum_{i=1}^N b_i |i
angle$$

The output will be in the form

$$|x\rangle = \sum_{i=1}^{N} x_i |i\rangle$$

Let the eigenvalues and eigenvectors of *A* be denoted as:

$$A|\mu_j\rangle = \lambda_j|\mu_j\rangle$$

The b ket expanded in this eigenbasis is

$$|b\rangle = \sum_{j=1}^{N} \beta_j |\mu_j\rangle$$

Assuming matrix A is sparse and hermitian, and  $1/\kappa \leqslant |\lambda_j| \leqslant 1$ 

Assuming matrix A is sparse and hermitian, the complexity of the quantum algorithm is

$$O\left(\frac{\kappa^2}{\epsilon}\log(N)s^2\right)$$

and recall the output is  $|x\rangle = \sum_{i=1}^{N} x_i |i\rangle$ .

Do you see a problem?

Assuming matrix A is sparse and hermitian, the complexity of the quantum algorithm is

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and recall the output is  $|x\rangle = \sum_{i=1}^{N} x_i |i\rangle$ .

It takes at least N steps to readout the vector coefficients.

To overcome this problem, the actual output of algorithm is  $\langle x|M|x\rangle$ .

The complexity of the quantum algorithm is

$$O\left(\frac{\kappa^2}{\epsilon}\log(N)s^2\right)$$

Compare this to the classical algorithm,

$$O\left(\kappa \log(1/\epsilon) \cdot Ns\right)$$

The classical complexity is the same whether the output is  $\mathbf{x}$  or  $\mathbf{x}^{\dagger}M\mathbf{x}$ .

The complexity of the quantum algorithm is

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Compare this to the classical algorithm,

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In some cases the quantum algorithm can be exponentially faster.

Step 1 – Prepare the ancilla

$$|\Psi_0\rangle = \sqrt{\frac{2}{T}} \sum_{\tau=0}^{T-1} \sin \frac{\pi(\tau + \frac{1}{2})}{T} |\tau\rangle$$

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$$\left(\sum_{\tau=0}^{T-1} [\tau] \otimes e^{iA\tau t_0/T}\right) |\Psi_0\rangle \otimes \sum_{j=1}^N \beta_j |\mu_j\rangle$$

Ref: D.W. Berry at al, Efficient Quantum Algorithms for Simulating Sparse Hamiltonians, Comm. Math. Phys. **270**, 359-371 (2007).

Step 1 – Prepare the ancilla and act the following controlled-unitary to it

$$\left(\sum_{\tau=0}^{T-1} [\tau] \otimes e^{iA\tau t_0/T}\right) |\Psi_0\rangle \otimes \sum_{j=1}^N \beta_j |\mu_j\rangle$$

produces

$$\sum_{\tau=0}^{T-1} \sum_{j=1}^{N} \sin \frac{\pi(\tau + \frac{1}{2})}{T} |\tau\rangle \otimes e^{i\lambda_j \tau t_0/T} \beta_j |\mu_j\rangle$$

Step 2 – Fourier transform the first qudit

$$\sum_{\tau=0}^{T-1} \sum_{j=1}^{N} \sin \frac{\pi(\tau + \frac{1}{2})}{T} |\tau\rangle \otimes e^{i\lambda_j \tau t_0/T} \beta_j |\mu_j\rangle$$

given by

$$|\tau\rangle \to \frac{1}{\sqrt{T}} \sum_{k=0}^{T-1} e^{2\pi i \tau k/T} |k\rangle$$

and we get

$$\sum_{k=0}^{T-1} \sum_{j=1}^{N} \alpha_{k|j} \beta_j |k\rangle |\mu_j\rangle$$

Step 3 – Add ancilla qubit and perform controlled-rotation

$$\sum_{k=0}^{T-1} \sum_{j=1}^{N} \alpha_{k|j} \beta_j |k\rangle |\mu_j\rangle |0\rangle$$

$$\rightarrow \sum_{k=0}^{T-1} \sum_{j=1}^{N} \alpha_{k|j} \beta_j |k\rangle |\mu_j\rangle \left(\sqrt{1 - \frac{C^2}{k^2}} |0\rangle + \frac{C}{k} |1\rangle\right)$$

Step 4 – Undo phase estimation and uncompute  $|k\rangle$ 

$$\sum_{k=0}^{T-1} \sum_{j=1}^{N} \alpha_{k|j} \beta_{j} |k\rangle |\mu_{j}\rangle \left( \sqrt{1 - \frac{C^{2}}{k^{2}}} |0\rangle + \frac{C}{k} |1\rangle \right)$$

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The coefficients  $\alpha_{k|j}$  have the property

$$\alpha_{k|j} = \begin{cases} 1 & \text{if } \lambda_j = 2\pi k/t_0 \\ 0 & \text{otherwise} \end{cases}$$

Step 4 – Undo phase estimation and uncompute  $|k\rangle$ 

$$\sum_{k=0}^{T-1} \sum_{j=1}^{N} \alpha_{k|j} \beta_{j} |\mu_{j}\rangle \left( \sqrt{1 - \frac{C^{2}}{k^{2}}} |0\rangle + \frac{C}{k} |1\rangle \right)$$

which simplifies to

$$\sum_{j=1}^{N} \beta_j |\mu_j\rangle \left( \sqrt{1 - \frac{C'^2}{\lambda_j^2}} |0\rangle + \frac{C'}{\lambda_j} |1\rangle \right)$$

Step 5 – Measure ancilla qubit

$$\sum_{j=1}^{N} \beta_j |\mu_j\rangle \left( \sqrt{1 - \frac{C'^2}{\lambda_j^2}} |0\rangle + \frac{C'}{\lambda_j} |1\rangle \right)$$

If 1 is measured, we have the state

$$\sum_{j=1}^{N} \lambda_j^{-1} \beta_j |\mu_j\rangle = A^{-1} |b\rangle$$

## References

- 1) A.W. Harrow, A. Hassidim, S. Lloyd, PRL **103**, 150502 (2009).
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- 4) http://www.youtube.com/watch?v=KtIPA PyaPOg