Quantum Algorithms: An overview of techniques

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Outline

Main quantum tricks and techniques

- Quantum Fourier Transform
- ► The SWAP test
- Unitaries as representations
- Quantum simulation
- Dissipative & stochastic state preparation
- Quantum walks, Grover search

Quantum Fourier Transform

QFT over \mathbb{Z}_N

$$DFT_{N} = QFT_{N} = \frac{1}{\sqrt{N}} \begin{pmatrix} 1 & 1 & 1 & \dots & 1 \\ 1 & \omega & \omega^{2} & \dots & \omega^{N-1} \\ 1 & \omega^{2} & \omega^{4} & \dots & \omega^{2(N-1)} \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ 1 & \omega^{N-1} & \omega^{2(N-1)} & \dots & \omega^{(N-1)(N-1)} \end{pmatrix}, \text{ where } \omega = e^{\frac{2\pi i}{N}}.$$

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- Given a Boolean function *f*: {0, 1}ⁿ → {0, 1} decide whether it is constant (0 or 1) or balanced (50% 0 and 1).
- The function is given as an oracle $O_f : |x\rangle|b\rangle \mapsto |x\rangle|b \oplus f(x)\rangle$.

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Take away message

- Constructive interference can be used as a computational resource
- Studying problems in a black-box setting gives useful insights

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Take away message

- Shows the power of Fourier transform (over the group \mathbb{Z}_{2}^{n})
- (+1 Phase kickback is a surprising and useful quantum effect)

- Given a function $f : \mathbb{Z}_{K}^{n} \to \mathbb{Z}_{K}$ so that $f(x) = s \cdot x \pmod{K}$; find *s*.
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Recall: $QFT_{K} : |j\rangle \mapsto \frac{1}{\sqrt{K}} \sum_{\ell=0}^{K-1} e^{2\pi i \frac{j\ell}{K}} |\ell\rangle$

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Jordan's algorithm ($\mathbb{Z}_{\mathcal{K}} \rightsquigarrow \mathbb{R}$)

- For a differentiable function $f \colon \mathbb{R}^n \to \mathbb{R}$ we have $f(x_0 + \delta_x) \approx f(x_0) + \nabla f \cdot \delta_x$
- Discretize \mathbb{R} and run the above algorithm for large enough K (resolution is $\approx \frac{1}{K}$)
- ► Implement $U_f: |\delta_x\rangle \mapsto e^{\frac{2\pi i f}{K}f(x_0+\delta_x)} |\delta_x\rangle \approx e^{\frac{2\pi i f(x_0)}{K}} e^{\frac{2\pi i (\nabla f \cdot \delta_x)}{K}} |\delta_x\rangle$ with one evaluation of f

Generalizations and applications of Jordan's algorithm

Convex functions

- Have at least one subgradient at every point
- Around most points can be well approximated by a linear function

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Separating hyperplanes

Exponential speed-up for finding separating hyperplanes (2018):

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Gradient computation for variational qauntum circits (QAOA)

• $\frac{1}{\epsilon}$ Quadratic speed-up for computing the gradient (**G**, Arunachalam, Wiebe 2017)

Phase estimation ($\mathbb{Z}_2^n \rightsquigarrow \mathbb{Z}_{2^n}$)

Phase estimation problem

Given $U = \sum_{\lambda} e^{2\pi i \lambda} |\psi_{\lambda}\rangle \langle \psi_{\lambda}|$ and an eigenstate $|\psi_{\lambda}\rangle$ output λ .

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The Hidden subgroup problem (HSP) ($\mathbb{Z}_{2^n} \rightsquigarrow G$)

- ▶ Input: Oracle access to a function $f: G \rightarrow S$ for some group G and (finite) set S
- ▶ **Promise:** There is a subgroup $H \le G$ such that f(x) = f(y) iff $x^{-1}y \in H$
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Algorithm for solving the problem – Kitaev (1995)

$$\begin{array}{c} \frac{1}{\sqrt{|G|}} \sum_{g \in G} |g\rangle \\ |0\rangle \end{array} \begin{array}{c} - O_f \\ - O_f \end{array}$$

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Works well for Abelian groups

- Samples a uniformly random character / irrep. of G that is trivial on H
- One can find a generator system of H after a few repetitions
- ▶ We can implement *QFT*_G efficiently

Some examples of the Abelian HSP

Simon's problem

- Function: $f: \{0, 1\}^n \rightarrow \{0, 1\}$ (the group is \mathbb{Z}_2^n)
- ▶ Subgroup: $\{0, s\}$, i.e., f(x) = f(y) iff $x y \in \{0, s\}$
- ► Output: s

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Period finding (and Shor's algorithm)

- ▶ **Function:** $f: \mathbb{Z} \to \mathbb{Z}_N$ (in Shor's algorithm $f(x) = a^x \mod N$ for some *a*)
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Discrete log (for given γ , A find a such that $A = \gamma^a$)

- ▶ **Function:** $f: \mathbb{Z}_N \times \mathbb{Z}_N \to \mathbb{Z}_N$ mapping $(x, y) \mapsto \gamma^x A^{-y} \mod N$
- ▶ Subgroup: $\langle (a,1) \rangle$, i.e., f(x,y) = f(x',y') iff $\exists c \in \mathbb{Z}_N : (x x', y y') = (ac, c)$
- Output: a

For more info see, e.g., Ronald de Wolf's lecture notes: arXiv:1907.09415

More advanced algorithms based on Abelian HSPs

Solving Pell's equation (Hallgren 2002)

 $x^2 - dy^2 = 1$

- Solving the principal ideal problem (Hallgren 2002)
- Period finding over \mathbb{R} and \mathbb{R}^n

▶ :

- Computing the unit group of number fields
- Breaking elliptic curve based cryptography

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See Sean Hallgren's talk on Thursday for more on this direction!

The non-Abelian HSP

What works and what does not

- ► *QFT_G* is somewhat harder to define and implement
- Unclear how to efficiently recover the subgroup
- ► However, the same algorithm is actually query efficient (Barnum & Knill 2002)
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Important example: Graph isomorphism (i.e., deciding whether $G \simeq G'$)

- Group: S_{2n} , Function: permute the vertices of $G \cup G'$
- Subgroup: Automorphisms of $G \cup G'$
- ▶ Output: whether there is a generator interchanging vertices of G and G'

The SWAP test

A simpler algorithm for graph isomorphism

Prepare a uniform superposition

- Let $|\psi_0\rangle \propto \sum_{s\in S_n} |s(G)\rangle$
- Let $|\psi_1
 angle \propto \sum_{s \in S_n} \left| s(G') \right\rangle$
- Observe that

$$\langle \psi_0 | \psi_1
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The probability of getting outcome + is

$$\frac{1}{2} + \frac{1}{2} |\langle \psi_0 | \psi_1 \rangle|^2$$

Unitaries as representations

Towards approximating the Jones polynomial

The Hadamard test



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13/20

Towards approximating the Jones polynomial

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The Jones polynomial – a link invariant

A *link* is a collection of loops embedded into \mathbb{R}^3 , in a possibly intertwined way. A *link invariant* is a quantity associated to links that is invariant under smooth transformations of the embedding.



$$|\psi\rangle - U$$

13/20

Towards approximating the Jones polynomial

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Links from braids

A *braid* is a collection of parallel strands, where adjacent strands are allowed to cross under or over each other. One can get a link by connecting the bottom and top ends of the strands.



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Quantum algorithms and connections to field theory

For a root of unity e^{2πi/k}, the relevant representation is unitary; the corresponding value of the Jones polynomial can be approx. evaluated via estimating (ψ|U|ψ). This (BQP-complete) algorithm is due to Aharonov, Jones, and Landau (2006).

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- Witten showed that the Jones polynomial is closely related to topological quantum field theory (TQFT).
- Friedman, Kitaev, Larsen, and Wang (2001) showed that quantum computers can efficiently simulate TQFTs.

Quantum simulation

(Dynamical) Hamiltonian simulation

Time-independent Hamiltonians

Schrödinger's equation ($\hbar = 1$) for time-independent quantum systems:

$$rac{d}{dt}|\psi
angle=-iH|\psi
angle \Longrightarrow |\psi(t)
angle=e^{-itH}|\psi(0)
angle$$

Recap – matrix functions

Any Hermitian matrix *H* can be diagonalised using some unitary *V* such that $H = V^{\dagger}DV = \sum_{\lambda} \lambda |\lambda \rangle \langle \lambda |$.

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Wait a minute, don't we build quantum computers using Hamiltonian simulation???

Product formula approach (Lloyd 1996)

Time-independent local Hamiltonians

Let $H = \sum_{k=1}^{K} H_k$, where each term H_k acts on a constant (say 2) number of qubits.

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$$e^{-itH} = (e^{-\frac{itH}{r}})^r = (e^{-\frac{itH_1}{r}}e^{-\frac{itH_2}{r}}\cdots e^{-\frac{itH_K}{r}})^r + O\left(\frac{(tK)^2}{r}\right).$$

Choosing $r = \Theta((tK)^2/\varepsilon)$ guarantees an ε -approximation.

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(Query) Optimal Hamiltonian simulation of sparse matrices

Quantum Signal Processing (QSP): (Low & Chuang 2016)

 $O(t \|H\|_{\max} \mathbf{s} + \log(1/\varepsilon))$

For a recent survey see: Childs, Maslov, Nam, Ross, Su - arXiv: 1711.10980

More generalizations and improvements

A few more recent generic results (without being exhaustive)

► Time-dependent sparse Hamiltonians: (Berry, Child, Su, Wang, Wiebe 2019)

$$\widetilde{O}\left(s\int_{0}^{t}\left\|H(\tau)\right\|_{\max}d au
ight)$$

Quantum chemistry: (Babbush, Berry, McClean, Neven 2019)

 $\widetilde{O}(N^{\frac{1}{3}}\eta^{\frac{8}{3}})$, with N : #plane wave orbitals, η : #electrons

Lattice Hamiltonians: (Haah, Hastings, Kothari, Low: QIP'19)

 $\widetilde{O}(nt)$

..., multi-product formulas, interaction picture simulation, ...

More generalizations and improvements

A few more recent generic results (without being exhaustive)

► Time-dependent sparse Hamiltonians: (Berry, Child, Su, Wang, Wiebe 2019)

$$\widetilde{O}\left(s\int_{0}^{t}\left\|H(\tau)\right\|_{\max}d au
ight)$$

Quantum chemistry: (Babbush, Berry, McClean, Neven 2019)

 $\widetilde{O}(N^{\frac{1}{3}}\eta^{\frac{8}{3}})$, with N : #plane wave orbitals, η : #electrons

Lattice Hamiltonians: (Haah, Hastings, Kothari, Low: QIP'19)

 $\widetilde{O}(nt)$

..., multi-product formulas, interaction picture simulation, ...

Simulating quantum field theory? See Preskill's recent survey: arXiv:1811.10085

Dissipative & stochastic state preparation

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 - pick an unchecked constraint and check (measure) it
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A loosely related result

Quant. Metropolis samp. (Temme, Osborne, Vollbrecht, Poulin, Verstraete 2009)

Quantum walks

Continuous-time quantum / random walks

Laplacian of a weighted graph

Let G = (V, E) be a finite simple graph, with non-negative edge-weights $w : E \to \mathbb{R}_+$. The Laplacian is defined as

$$u \neq v$$
: $L_{uv} = w_{uv}$, and $L_{uu} = -\sum_{v} w_{uv}$.

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Evolution of the state:

$$\frac{d}{dt}p_{u}(t) = \sum_{v \in V} L_{uv}p_{v}(t) \qquad \Longrightarrow \qquad p(t) = e^{tL}p(0)$$

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$$irac{d}{dt}\psi_u(t)=\sum_{v\in V}L_{uv}\psi_v(t) \implies \psi(t)=e^{-itL}\psi(0)$$
Exponential speedup by a quantum walk

Childs, Cleve, Deotto, Farhi, Gutmann, and Spielman: quant-ph/0209131