$2 + \varepsilon$ quantum learning algorithms

Ashley Montanaro

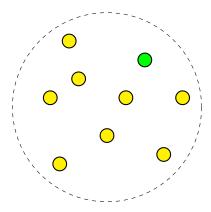
Talk based on joint work with Andris Ambainis and ongoing joint work with Scott Aaronson, David Chen, Daniel Gottesman and Vincent Liew.

20 March 2013





What is learning?



In this talk

Learning a set $S \equiv$ identifying an arbitrary, unknown object picked from *S*.

This talk

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On this principle, I'll talk about three optimal quantum algorithms for learning an unknown...

- ... bit-string, given access to "wildcard" queries;
- ... low-degree multilinear polynomial;
- ... stabilizer state.

Bonus mini-result: A composition theorem for classical decision tree complexity.

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Example

Imagine the hidden string is x = 01101. Then querying...

- 0 * 1 * 1 returns 1;
- *1 * 1* returns 0.

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Contrast: In the standard model, there is a quantum speed-up by about a factor of 2 [van Dam '98], and this is optimal.

Solving SWW

The solution to SWW is based on this claim:

Measurement Lemma

Fix $n \ge 1$ and, for any $0 \le k \le n$, set

$$|\psi_x^k\rangle := \frac{1}{\binom{n}{k}^{1/2}} \sum_{S \subseteq [n], |S|=k} |S\rangle |x_S\rangle,$$

where $|x_S\rangle := \bigotimes_{i \in S} |x_i\rangle$. Then, for any $k = n - O(\sqrt{n})$, there is a quantum measurement (POVM) which, on input $|\psi_x^k\rangle$, outputs \tilde{x} such that the expected Hamming distance $d(x, \tilde{x})$ is O(1).

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Why does this let us solve SWW?

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- ... but after each measurement, an expected *O*(1) bits are incorrect.
- How to fix these?

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In particular, we would like to minimise the dependence on *n*.

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- Many applications known: molecular biology, data streaming algorithms, compressed sensing, pattern matching in strings, ...
- See the book "Combinatorial Group Testing and Its Applications" [Du and Hwang '00] for more.

Quantum algorithms for CGT

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Basic idea:

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- In the CGT problem, we have access to an oracle which computes $f(s) = \bigvee_i x_i s_i$ for arbitrary $s \in \{0, 1\}^n$. But if $|x| \leq 1$, then for any s, $\bigvee_i x_i s_i = x \cdot s$.

Generalising this idea to arbitrary k

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- Following each successful query, we reduce *k* by 1 and exclude the bit that we just learned from future queries.
- In order to learn *x* completely, the expected overall number of queries used is *O*(*k*).

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- A wildcard query corresponding to S ⊆ [n] and the string *x*_S returns 1 iff all bits of *x*_S are correct. Negating the output gives a query that behaves the same as a CGT query.
- So we can use the algorithm for CGT to find, and correct, all incorrect bits using *O*(1) queries.

Summary

• Using an efficient algorithm for CGT as a subroutine, we can solve search with wildcards using $O(\sqrt{n})$ queries.

• This is a square-root speed-up which (apparently) does not come from amplitude amplification or quantum walks.

• Open problem: Determine the quantum query complexity of CGT. We have an upper bound of O(k) and a lower bound of $\Omega(\sqrt{k})$.

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- If *f* is picked from a known set *F*, we need at least $\log_q |\mathcal{F}|$ classical queries.
- On a quantum computer, we have the ability to query *f* in superposition, i.e. to perform the map

 $|x\rangle|z\rangle \mapsto |x\rangle|z+f(x)\rangle.$

 $f : \mathbb{F}_q^n \to \mathbb{F}_q$ is a degree *d* multilinear polynomial:

$$f(x) = \sum_{S \subseteq [n], |S| \leqslant d} \alpha_S \prod_{i \in S} x_i$$

for some coefficients $\alpha_S \in \mathbb{F}_q$, where $[n] := \{1, \ldots, n\}$.

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- Note that for $S = \emptyset$ we define $\prod_{i \in S} x_i = 1$.
- For example, any multilinear polynomial of degree 3 can be written as

$$f(x) = \alpha_{\emptyset} + \sum_{i} \alpha_{\{i\}} x_i + \sum_{i < j} \alpha_{\{i,j\}} x_i x_j + \sum_{i < j < k} \alpha_{\{i,j,k\}} x_i x_j x_k.$$

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- The set of degree *d* polynomials over 𝔽₂ is known as the binary Reed-Muller code of order *d*.

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- Upper bound: It suffices to query f(x) for all strings $x \in \mathbb{F}_q^n$ that contain only 0 and 1, and such that $|x| \leq d$.
- Lower bound: there are $q^{\Theta(n^d)}$ distinct multilinear degree d polynomials of n variables over \mathbb{F}_q ; each classical query to f only provides $\log_2 q$ bits of information.

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Notes:

- The lower bound follows from Holevo's theorem.
- The Bernstein-Vazirani algorithm [Bernstein and Vazirani '97] is the case q = 2, d = 1.
- Rötteler previously gave a bounded-error quantum algorithm for the case q = 2, d = 2 [Rötteler '09].
- A quantum algorithm for estimating a quadratic form over the reals had previously been given by Jordan [Jordan '08].

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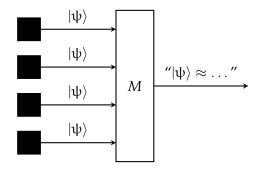
- Claim: Taking the derivative of *f* in (*d*−1) different directions *i* leaves a linear function, which can be learned using one query.
- Claim 2: Computing this derivative can be done using 2^{d-1} queries, and we need to do it for at most $\binom{n}{d-1}$ different sets.

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- For any degree *d* multilinear polynomial *f* : 𝔽ⁿ_q → 𝔽_q, define the discrete derivative of *f* in direction *i* ∈ [*n*] as

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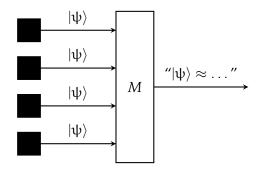
- Claim: Taking the derivative of *f* in (*d*−1) different directions *i* leaves a linear function, which can be learned using one query.
- Claim 2: Computing this derivative can be done using 2^{d-1} queries, and we need to do it for at most $\binom{n}{d-1}$ different sets.
- So we can learn f using $O(n^{d-1})$ queries.

Consider the basic task of quantum state estimation.



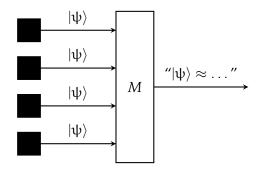
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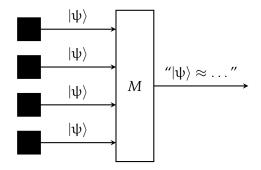
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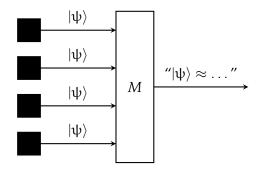
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- Standard quantum state tomography uses $2^{\Theta(n)}$ copies of $|\psi\rangle$ to achieve constant fidelity.
- Can we do better?

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- To achieve constant fidelity between our guess and |ψ⟩, we need 2^{Ω(n)} copies of |ψ⟩.
- In order to determine |ψ⟩ efficiently (using poly(*n*) copies) we must restrict to classes of states which have efficient descriptions, or change the problem.

Some examples where this has been done:

- [Cramer et al '10] give an efficient algorithm for learning matrix product states.
- [Aaronson '06] introduces "pretty good tomography": relax to attempting to predict the outcomes of "most" measurements on the state.
- [Flammia and Liu '11] and [da Silva et al '11] give efficient algorithms for certifying the production of certain states.

Learning stabilizer states

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- Examples include GHZ states, cluster states, states occurring in quantum error-correcting codes, ...

Today I'll talk about a learning algorithm for another important class of quantum states with efficient descriptions: stabilizer states.

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A stabilizer state of *n* qubits is completely specified by a generating set for its stabilizer (*n* Pauli matrices on *n* qubits). There are $2^{\Theta(n^2)}$ stabilizer states of *n* qubits.

Prior work on learning stabilizer states

[Aaronson and Gottesman '08] have previously given quantum algorithms for learning an unknown stabilizer state $|\psi\rangle$:

- An algorithm which uses O(n) copies of $|\psi\rangle$ and runs in time $O(n^4)$;
- An algorithm which uses $O(n^2)$ copies of $|\psi\rangle$, runs in time $O(n^4)$ and uses only single-copy measurements.

Theorem

There is a quantum algorithm which learns an unknown stabilizer state $|\psi\rangle$ given access to O(n) copies of $|\psi\rangle$, and runs in time $O(n^3)$ (or better).

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Notes on this result:

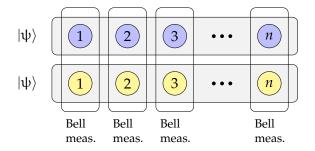
- By Holevo's theorem, this is optimal in terms of the scaling of the number of copies of |ψ⟩ used.
- Any algorithm for learning stabilizer states requires Ω(n²) time just to write the output.
- $\bullet\,$ The algorithm makes measurements on two copies of $|\psi\rangle\,$ at a time.

The algorithm

The algorithm is based on the following subroutine.

Bell sampling

- Create two copies of $|\psi\rangle$.
- **2** Measure each pair of qubits of $|\psi\rangle^{\otimes 2}$ in the Bell basis.



• For
$$z, x \in \{0, 1\}$$
, write $\sigma_{zx} := \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}^z \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}^x$.

• For $s \in \{0, 1\}^{2n}$, write

$$\sigma_s := \sigma_{s_1 s_2} \otimes \cdots \otimes \sigma_{s_{2n-1} s_{2n}}.$$

Fact

Let $|\psi\rangle$ be a state of *n* qubits. Performing Bell sampling on $|\psi\rangle^{\otimes 2}$ returns outcome *s* with probability

 $\frac{|\langle \psi | \sigma_{\scriptscriptstyle S} | \psi^* \rangle|^2}{2^n}$

 $\bullet~$ Up to an overall phase every stabilizer state $|\psi\rangle$ can be written in the form

$$|\psi\rangle = \frac{1}{\sqrt{|A|}} \sum_{x \in A} i^{\ell(x)} (-1)^{q(x)} |x\rangle,$$

where *A* is an affine subspace of \mathbb{F}_2^n , and $\ell, q : \{0, 1\}^n \to \{0, 1\}$ are linear and quadratic (respectively) polynomials over \mathbb{F}_2 [Dehaene and Moor '02].

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• Hence

$$|\psi^*\rangle = \sigma_{10}^{\otimes S} |\psi\rangle.$$

• If we perform Bell sampling on $|\psi\rangle^{\otimes 2}$, we receive outcome *t* with probability

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• Let *G* stabilize $|\psi\rangle$ and let *T* denote the set of strings $t \in \{0, 1\}^{2n}$ such that $\sigma_t \in G$, up to a phase. Then *T* is an *n*-dimensional linear subspace of \mathbb{F}_2^{2n} .

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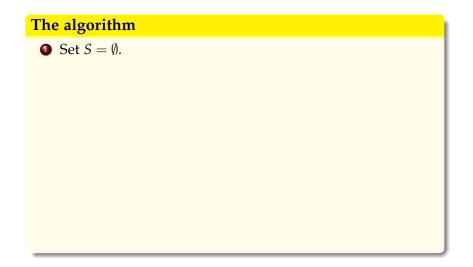
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- Bell sampling gives an outcome *r* which is uniformly distributed on the set {*t* ⊕ *s* : *t* ∈ *T*} for some *s* ∈ {0, 1}²ⁿ.

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 - If the dimension of the subspace of 𝔽₂²ⁿ spanned by these vectors is *n*, any basis of this subspace is a basis for *T*.
- Although *T* does not contain information about phases, determining *T* suffices to uniquely determine $|\psi\rangle$.
 - Once we have found a basis for *T*, we can measure $|\psi\rangle$ in the eigenbasis of each corresponding Pauli matrix *M* to decide whether $M|\psi\rangle = |\psi\rangle$ or $M|\psi\rangle = -|\psi\rangle$.



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- For each element of *B*, measure a copy of |ψ⟩ in the eigenbasis of the corresponding Pauli matrix *M* to determine whether *M*|ψ⟩ = |ψ⟩ or *M*|ψ⟩ = -|ψ⟩.

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 The algorithm fails (i.e. does not identify |ψ⟩) if each of the 2*n* samples *r* ⊕ *r*₀ lies in a subspace of *T* of dimension at most *n* − 1. This occurs with probability at most 2^{−n}.

Bonus: a composition theorem for decision tree complexity

Imagine we want to compute a function of the form

$$h(x) = g(f^1(x^1), \ldots, f^n(x^n)),$$

where $x^i \in \{0, 1\}^{m_i}$, using the minimal number of classical queries to *x*.

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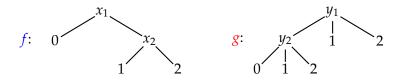
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"Theorem": The x^i inputs are independent, so this is the most efficient way to compute g.

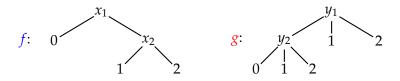
Counterexample to "theorem"

Let $f : \{0, 1\}^2 \rightarrow \{0, 1, 2\}$ and $g : \{0, 1, 2\}^2 \rightarrow \{0, 1, 2\}$ be defined by the decision trees below (where edges correspond to elements of $\{0, 1\}$ or $\{0, 1, 2\}$ in ascending order from left to right).

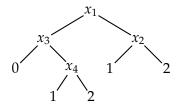


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Set $h(x_1, x_2, x_3, x_4) = g(f(x_1, x_2), f(x_3, x_4))$. Then *h* can be computed using only 3 queries:



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- Implies various corollaries, e.g. a direct sum theorem for decision tree complexity (a special case of a result of [Jain, Klauck and Santha '10]) and optimal bounds for iteratively defined functions.
- The quantum equivalent of this result was proven by [Høyer, Lee and Špalek '07] and [Reichardt '09].

Summary

We can learn...

- ... *n*-bit strings with $O(\sqrt{n})$ wildcard queries;
- ... degree *d n*-variate multilinear polynomials with $O(n^{d-1})$ queries;
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Open problems:

- Determine the quantum query complexity of CGT.
- Other applications of SWW! A possible example: testing juntas.
- What about testing stabilizer states?

Thanks!

Some further reading:

- The algorithm for search with wildcards: **arXiv:1210.1148** (joint work with Andris Ambainis)
- The algorithm for learning multilinear polynomials: arXiv:1105.3310
- The algorithm for learning stabilizer states: arXiv:13??.???? (joint work with Scott Aaronson, David Chen, Daniel Gottesman and Vincent Liew)
- The composition theorem for decision tree complexity: arXiv:1302.4207

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- *G_{xy}* depends only on *x* ⊕ *y*, so *G* is diagonalised by the Fourier transform over Zⁿ₂ and D_k does not depend on *x*.
- *D_k* can be upper bounded using Fourier duality and some combinatorics.