#### Search via quantum walk

Ashwin Nayak University of Waterloo, and Perimeter Institute for Theoretical Physics

Joint work with

Frédéric Magniez<sup>1</sup>, Jérémie Roland<sup>2</sup>, Miklos Santha<sup>1</sup> <sup>1</sup>LRI-CNRS, France, <sup>2</sup>UC Berkeley

#### Abstract search problem

- Input:
  - Set X = {a, b, c, ...}
  - Marked elements M subset of X (say, {a, g})
  - Procedure to answer "x in M?"
- Output:
  - Some element x in M.
- Additional structure: Markov chain *P* on *X*





## Random walk for search

- (*s*,*t*)-Connectivity
  - Input: Graph G on n vertices, two specified vertices s,t
  - Question: is there is a path from *s* to *t*?



- Algorithm: start at u = s, and repeat  $O(n^3)$  times
  - Pick a random vertex *v* adjacent to *u*
  - If v = t, stop. Else, set u = v.

#### Second example

- Element Distinctness (ED)
  - Input: list of *n* numbers  $\{x_1, x_2, x_3, ..., x_n\}$
  - Question: are all the numbers distinct (or is there a collision:  $x_i = x_{j}, i \neq j$ )
- Deterministic Algorithm:
  - Sort elements; check if consecutive numbers are equal
  - Time complexity:  $O(n \log n)$
- Not graph search, but can be recast as one.

Element distinctness as graph search

- Johnson Graph (*n*, *r* )
  - Vertices: size *r* subsets of {1, 2, ..., *n*}
  - Edges:  $\{S, T\}$  is an edge iff they differ by 2 elements
- Example: n = 15, r = 4



• Search for subset with collision

## Randomized algorithm for ED

• Start at a random vertex of the Johnson graph

Pick *r* indices uniformly at random to form a set *S*; sort the elements  $x_i$  for *i* in *S*; check for collisions.

- Repeat for  $T_1$  steps
  - Perform a random walk on the graph for T<sub>2</sub> steps
     In each step, swap random element *i* in *S* and *j* not in *S*;
     remove x<sub>i</sub>, insert x<sub>i</sub> into sorted list
  - check for a collision in *S*
- If no collision is found, output "no collision".
   (Less natural algorithm, but adapts well to quantum)

#### Randomized algorithm for ED **X**<sub>2</sub> **X**5 **X**<sub>2</sub> **X**<sub>3</sub> **X**<sub>13</sub> **X**8 **X**<sub>5</sub> **X**<sub>13</sub> **X**3 $X_2 \quad X_3$ **X**<sub>2</sub> X<sub>8</sub> **X**<sub>13</sub> **X**5 **X**<sub>7</sub>

- Intuition:
  - In  $T_2 = O(r)$  steps of walk, S is nearly uniformly distributed
  - Pr[ collision in random S ]  $\approx$   $(r/n)^2$
  - So in  $T_1 = O((n/r)^2)$  repetitions, a collision will be found
- Runtime:  $r \log r + T_1 (T_2 \log r + 1)$ Set up cost checking cost checking cost

#### Speed-up via quantum walk

- Quantum analogue of randomized algorithm
- Speeds up both  $T_1$  and  $T_2$  quadratically

[Ambainis '04]

• Run time of quantum algorithm for ED  $r \log r + (n/r) (r^{1/2} \log r + 1)$ 

 $n^{2/3} \log n$  (setting  $r = n^{2/3}$ )

- A second algorithm, for symmetric Markov chains
- Quadratic speed-up in detecting marked elements

[Szegedy '04]

## This talk: New search algorithm

- Quantum walk from any irreducible Markov chain
- Algorithm finds a marked element, if any, from any M
- Run time: set-up +  $T_1^{1/2}$  ( $T_2^{1/2}$  update + check) Pr(*M*)<sup>-1/2</sup> singular value gap<sup>-1/2</sup>
- Simple --- conceptually, and to analyze
- Unifies and improves several applications

# Talk outline

- Classical algorithm
- Quantum walk
- Quantum subroutines
  - Amplitude amplification
  - Phase estimation
- Search algorithm

#### **Classical search algorithm**



- Start in some start distribution s
- Repeat for  $T_1$  steps
  - Simulate  $T_2$  steps of the Markov chain P
  - Check if current state is marked
- If no marked element is found, output "none marked".

## Complexity of classical strategy

- *P* symmetric (for simplicity), ergodic
- Uniform stationary distribution (1-eigenvector)
- Say we start in s = uniform distribution
- Run-time characterized by
  - Spectral gap  $\delta(P) = 1 \text{second largest |eigenvalue|}$
  - Probability of marked elements  $\varepsilon = \Pr(M) = |M| / |X|$
- Proposition

Run-time of the classical strategy is set-up + (1/ $\epsilon$ ) ((1/ $\delta$ ) update + check)  $T_1 \rightarrow T_2$ 

## Talk outline

- Classical algorithm Run time =  $1/\epsilon\delta$
- Quantum walk
- Quantum subroutines
  - Amplitude amplification
  - Phase estimation
- Search algorithm



[Watrous '01, Szegedy '04]





- The quantum walk W(P)
  - State space: pairs of neighbouring vertices  $|x\rangle |y\rangle$
  - Step of walk: diffuse y over neighbours of x, new nbr. y' then, diffuse x over neighbours of y'
  - Diffusion: analogous to Grover search operator (reflection about state  $|x\rangle \sum_{y} \sqrt{p_{x,y}} |y\rangle$ , for each x)

## Spectrum of W(P)

[Szegedy '04]

- W(P) = product of two reflection operators
- Assume *P* is symmetric, ergodic Has uniform stationary distribution
- Spectrum of W(P) related to that of P
- For every singular value of P,  $\sigma = \cos \theta$  in (0,1) W(P) has eigenvalues  $\exp(\pm 2i \theta)$
- The remaining eigenvalues are ±1

## Spectral gap

- Largest singular value of *P* = 1, and is unique
   W(*P*) has unique eigenvalue 1 (in walk subspace)
- Eigenvector of W(P) with eigenvalue 1 is  $|\pi\rangle = (1/n^{1/2}) \sum |x\rangle |n\rangle$  where

$$|\Pi| = (\Pi \Pi) \sum_{x} |x| |P_x|$$

$$|p_x\rangle = \sum_{y} p_{xy}^{1/2} |y\rangle$$

• If  $\sigma = \cos \theta < 1$  is second largest singular value, eigenvalue gap of W(P) is  $|1 - \exp(2i\theta)| \ge 2(1 - \sigma)^{1/2} = 2\delta(P)^{1/2}$ square-root of spectral gap of P

# Talk outline

- Classical algorithm Run time =  $1/\epsilon \delta$
- Quantum walk Spectral gap =  $\delta^{1/2}$
- Quantum subroutines
  - Amplitude amplification
  - Phase estimation
- Search algorithm

### Amplitude amplification [Grover '96, BBHT '98, ...]

- Search for one out of n states
- Start state:  $|\pi\rangle = (1/n^{1/2}) \Sigma_x |x\rangle$
- Desired final state:  $|a\rangle$
- Alternately reflect through  $|a^{\perp}\rangle$  and  $|\pi\rangle$



## Complexity of amplitude amplification

- Angle of rotation =  $2 \varphi$  (sin  $\varphi$  =  $1/n^{1/2}$ )
- Number of iterations  $\approx (\pi/2) / (2\varphi) \approx n^{1/2}$
- Required reflection operators have small circuits
- Multiple marked states
  - Fraction of marked states  $\varepsilon = m/n$
  - target state =  $(1/m)^{1/2} \Sigma_{x \text{ in } M} |x\rangle$
  - Angle of rotation =  $2 \varphi$  (sin  $\varphi = (m/n)^{1/2} = \varepsilon^{1/2}$ )
  - Number of iterations  $\approx$  1/ $\varepsilon^{1/2}$
  - Quadratic speed-up over classical

# Talk outline

- Classical algorithm Run time =  $1/\epsilon \delta$
- Quantum walk Spectral gap =  $\delta^{1/2}$
- Quantum subroutines
  - Amplitude amplification
     Cost = 1/ε<sup>1/2</sup>
  - Phase estimation
- Search algorithm

Phase estimation

- Input: circuit for unitary Usuperposition  $|v\rangle$ , eigenvector with unknown eigenvalue  $\exp(2\pi i\theta)$
- **Output:** approximation to  $\theta$
- Proposition [Kitaev '95, Cleve, Ekert, Macchiavello, Mosca '98] Can compute an approximation to  $\theta$  within  $\eta$ with  $1/\eta$  repetitions of U, one copy of  $|v\rangle$ with probability 3/4

## **Reflection using phase estimation**



- *U* unitary operator
- v isolated eigenvector
- $\varphi$  spectral gap

#### Reflection through $|v\rangle$

- Run phase estimation algorithm on the current state, with U
- If approximate phase is "far" from  $\theta$ , flip sign
- Undo phase estimation

Precision required  $\approx \varphi/2$ Repetitions of  $U \approx 1/\varphi = 1$ / spectral gap

## Reflection via quantum walk W(P)

- $|\pi\rangle$  1-eigenvector of W(*P*)
- $\delta^{1/2}$  spectral gap of W(*P*)
- Reflection through  $|\pi\rangle$ Use phase estimation, as described Repetitions of W(P)  $\approx$  1/ spectral gap  $\approx$  1/ $\delta^{1/2}$

# Talk outline

- Classical algorithm Run time =  $1/\epsilon \delta$
- Quantum walk Spectral gap =  $\delta^{1/2}$
- Quantum subroutines
  - Amplitude amplification
     Cost = 1/ε<sup>1/2</sup>
  - Phase estimation
    - $Cost = 1/\delta^{1/2}$
- Search algorithm

### The search algorithm

• Start state:

$$|\pi\rangle$$
 = (1/n<sup>1/2</sup>)  $\Sigma_x$   $|x\rangle|p_x\rangle$ 

• Desired final state:

$$|\mu\rangle = (1/m^{1/2}) \Sigma_{x \text{ in } M} |x\rangle |p_x\rangle$$

• Alternately reflect through  $|\mu^{\perp}\rangle$  and  $|\pi\rangle$  à la Grover



#### Implementing the reflections

• Reflection through  $|\mu^{\perp}\rangle$ 

If vertex *x* in first register is marked, and second register is in state  $|p_x\rangle$ , then flip sign

• Reflection through  $|\pi
angle$ 

Use phase estimation algorithm, as described

## Complexity of the algorithm

• Angle between  $|\mu^{\perp}\rangle$  and  $|\pi\rangle$ :

 $\sin \varphi = (m/n)^{1/2} = \varepsilon^{1/2},$ 

- $\varepsilon = \Pr(M) = \text{probability of } M \text{ under stationary}$ distribution
- Number of rotations à *la* Grover:  $1/\varepsilon^{1/2}$
- Cost of reflection through  $|\mu^{\perp}\rangle$

check + update cost

• Cost of reflection through  $|\pi\rangle$ :

update cost times  $1/\delta^{1/2}$ 

 $\delta^{1/2}$  = spectral gap of W(*P*)

Complexity

set-up +  $(1/\epsilon^{1/2})$  ( $(1/\delta^{1/2})$  update + check)

#### Final remarks

- Error due to imperfect phase estimation algorithm handled with a recursive search algorithm à la [Hoyer, Mosca, de Wolf '04]
- Algorithm extends to any irreducible Markov chain
- Unified and improved algorithms for Element Distinctness, Triangle Finding, Matrix Product verification, Group Commutativity
- Better algorithms for applications in which checking cost is higher than update cost