

# Finding many collisions via quantum walks

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# Classical Collision Finding

# Classical collision algorithms

$$f : \{0, 1\}^n \rightarrow \{0, 1\}^n$$

## A first algorithm

- ▶ Create a sorted list of size  $2^{n/2}$
- ▶ Look for collisions

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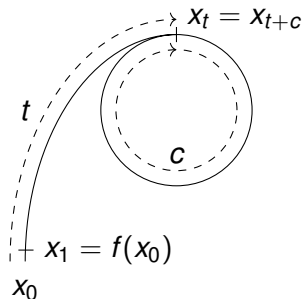
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## Pollard-Rho

- ▶ The cyclic part of the functional graph of  $f$  is of size  $\Theta(2^{n/2})$
- ▶ Seek a collision in the graph
- ▶ Memory-less, parallelizable



## Other cases

$$f : \{0, 1\}^n \rightarrow \{0, 1\}^m, n \leq m \leq 2n$$

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Lower bounds

Matching query lower bound in all cases

# Quantum Collision Finding



# Quantum Search

## Operators

- ▶ Test:  $|x\rangle \mapsto -|x\rangle$  if  $x$  marked, identity otherwise
- ▶ Diffusion:  $\sum_{i=0}^{2^n-1} |i\rangle \mapsto -\sum_{i=0}^{2^n-1} |i\rangle$ , identity otherwise

## Principle

Use the test operator to differentiate marked elements from the others, and use the diffusion operator to make marked elements interfere constructively.

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## Algorithm

- ▶ Test and Diffusion are reflexions in the same plane
- ▶ Composition is a rotation of angle  $2 \arcsin \sqrt{\epsilon}$
- ▶ Need to repeat until angle is  $\sim \pi/2$

# Quantum Collision Finding via Quantum Search

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## What works: a list + an enumeration (BHT algorithm)

- ▶ Take a list  $L = (f(y_0), \dots, f(y_{2^u}))$
- ▶ Search for an  $x$  such that there exists  $i$  with  $f(x) = f(y_i)$  and  $x \neq y_i$
- ▶ Cost  $2^u$  memory,  $2^u + \sqrt{\frac{2^n}{2^u}}$  time

## BHT algorithm

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## Finding $2^k$ collisions

- ▶ Use one larger list of size  $2^{n/3+2k/3}$
- ▶ Do  $2^k$  quantum searches  $\left( \text{cost } 2^k \times \sqrt{\frac{2^n}{2^{n/3-2k/3}}} = 2^{n/3+2k/3} \right)$

## Lower bound [LZ19]

General query lower bound  $\Omega(2^{m/3+2k/3})$

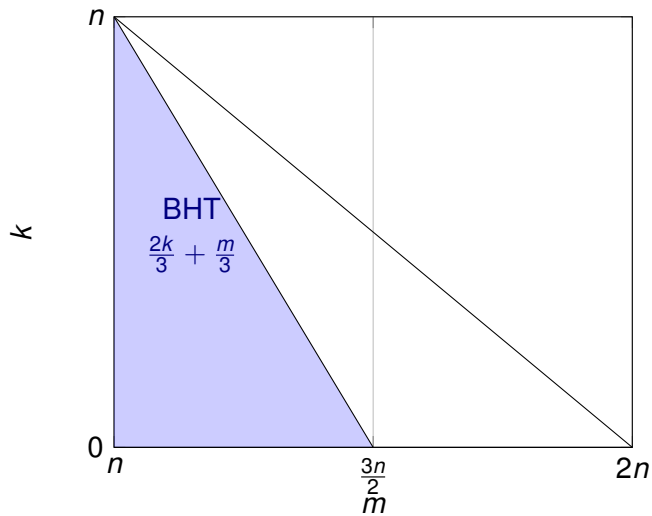
# With larger $m$

## BHT algorithm

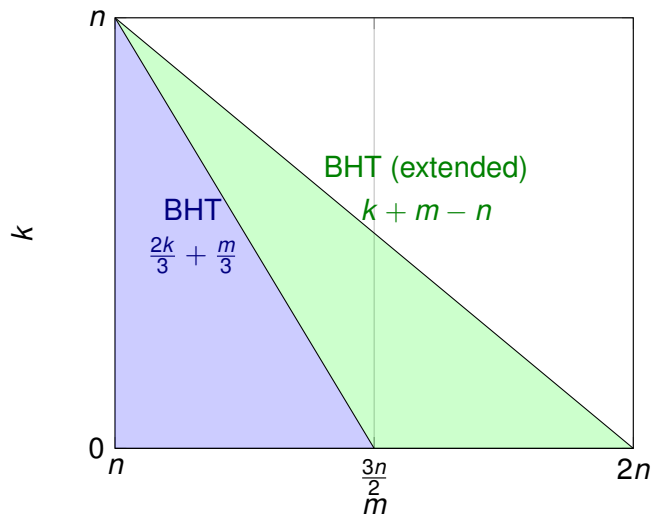
Query a list of size  $2^\ell$

- ▶ Only  $2^{2n-m}$  inputs are part of a collision
- ▶ Each element has probability  $2^{n-m}$  to be in a collision pair
- ▶  $\mathcal{O}(2^{\ell-m+n})$  collisions pairs for the initial list.
- ▶ Output  $2^k$  collision pairs, need  $\ell - m + n \geq k$ , otherwise the list might contain no relevant input  $\leadsto m \leq \frac{3}{2}n$

# Summary



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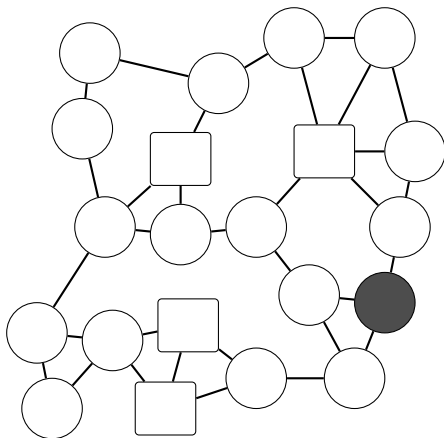


# Collision Finding

via Random Walks

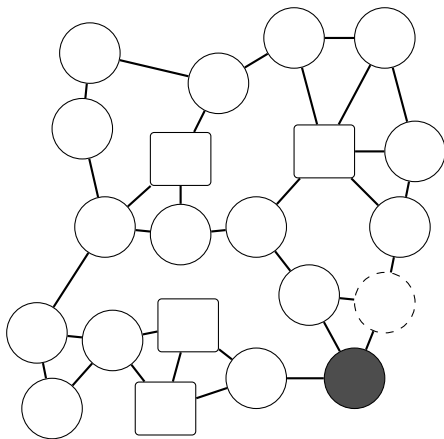
# Classical random walk

We move to random neighbors until we find a marked vertex.



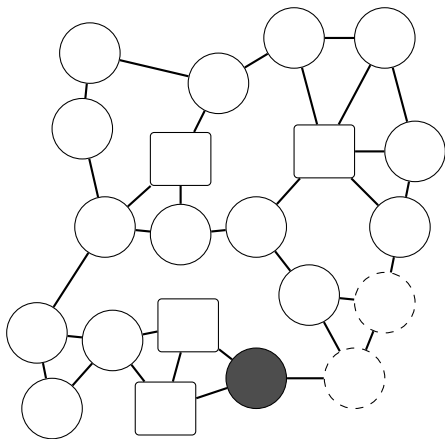
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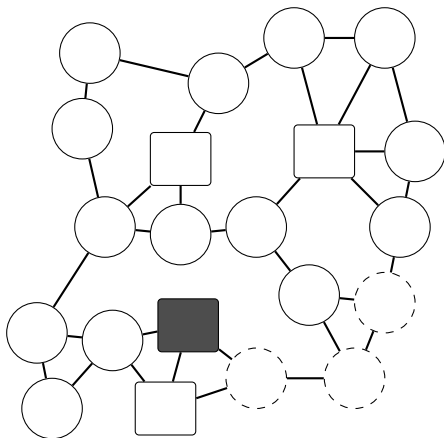
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# Cost of a classical random walk

We need procedures:

- ▶ To **setup** a starting arbitrary vertex (S)
- ▶ To **move** from one vertex to one of its neighbors (U)
- ▶ To **check** if a vertex is marked (trivial) (C)

We will find a marked vertex in time:

$$S + \underbrace{\frac{1}{\epsilon}}_{\epsilon \text{ proportion of marked vertices}} \left( \underbrace{\frac{1}{\delta}}_{\delta \text{ spectral gap of the graph}} (U + C) \right)$$

where  $\frac{1}{\delta}$  is the number of updates before we reach a new uniformly random vertex.

## Example: Walk-based collision finding

### Definition (Johnson graph)

- ▶ Nodes are sets of  $2^r$  elements among  $2^n$
- ▶  $N_1$  and  $N_2$  are adjacent is  $|N_1 \cap N_2| = 2^r - 1$
- ▶  $\frac{1}{\delta} = \frac{2^r(2^n - 2^r)}{2^n} \simeq 2^r$  (We need to replace all elements.)

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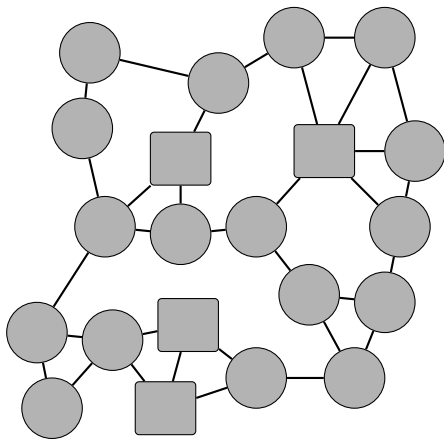
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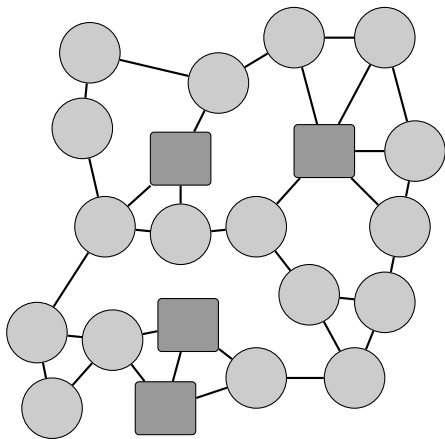
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As in quantum search, the walk transforms a uniform superposition over the whole graph into a superposition over marked vertices.



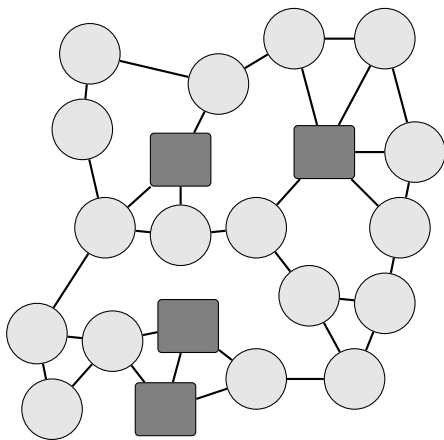
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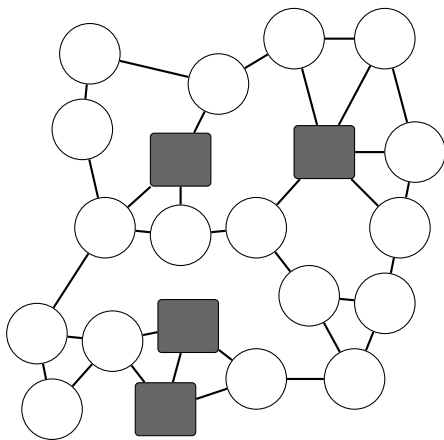
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# Time of a quantum walk (MNRS framework)

- ▶ The **setup** now requires to create a superposition over **all** vertices
- ▶ As in quantum search, we perform  $\sqrt{\frac{1}{\epsilon}}$  steps instead of  $\frac{1}{\epsilon}$
- ▶ But the mixing is also accelerated!

$$S + \underbrace{\sqrt{\frac{1}{\epsilon}}}_{\text{Walk steps}} \left( \underbrace{\sqrt{\frac{1}{\delta}} U}_{\text{Mixing time}} + C \right)$$

- ▶ The **Update** handles all vertices and all edges in superposition

# Ambainis's algorithm for Collision Finding

## Problem

$f : \{0, 1\}^n \rightarrow \{0, 1\}^m$ ,  $n \leq m \leq 2n$ , find a collision

## MNRS walk in a Johnson graph

- ▶ Setup : Create the uniform superposition of all lists of  $2^r$  elements.
- ▶ Fraction of marked nodes :  $\epsilon = 2^{2r-m}$
- ▶ Mixing time:  $\sqrt{\frac{1}{\delta}} = 2^{r/2}$
- ▶ Assume Update and Test polynomial time
- ▶ Cost  $2^r + 2^{m/2-r} \times 2^{r/2} \simeq \max(2^r, 2^{m/2-r/2})$



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 $\leadsto$  optimal for  $r = m/3$

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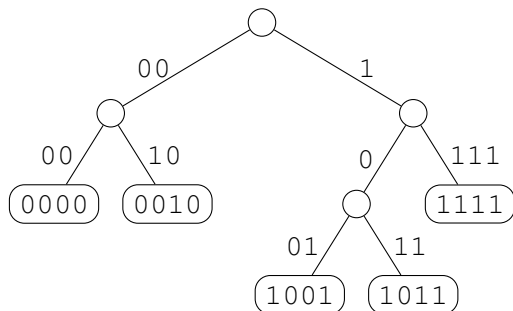
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## History-free quantum data structures: step 1

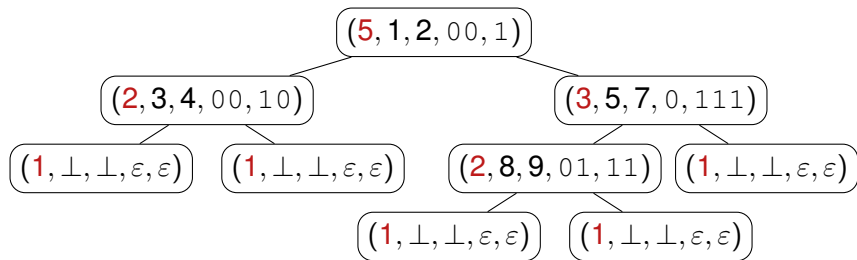
Take an efficient classical data structure (ex. radix tree)



Tree representing  $\{0000, 0010, 1001, 1011, 1111\}$ .

# Quantum data structures

Example for set  $S = \{0000, 0010, 1001, 1011, 1111\}$ :



Format: (# leaves, addr left, addr right, label left, label right)

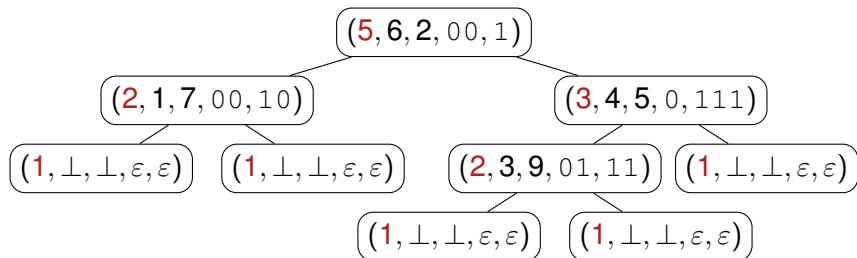
Actual memory content:

0	(5, 1, 2, 00, 1)	1	(2, 3, 4, 00, 10)	2	(3, 5, 7, 10, 11)
3	(1, ⊥, ⊥, ε, ε)	4	(1, ⊥, ⊥, ε, ε)	5	(2, 8, 9, 01, 11)
6	Empty cell	7	(1, ⊥, ⊥, ε, ε)	8	(1, ⊥, ⊥, ε, ε)
9	(1, ⊥, ⊥, ε, ε)				

Valid sequence: (0, 1, 3, 4, 2, 5, 8, 9, 7)

# Quantum data structures

Example for set  $S = \{0000, 0010, 1001, 1011, 1111\}$ :



Format: (# leaves, addr left, addr right, label left, label right)

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0	(5, 6, 2, 00, 1)	1	(1, ⊥, ⊥, ε, ε)	2	(3, 4, 5, 10, 11)
3	(1, ⊥, ⊥, ε, ε)	4	(2, 3, 9, 01, 11)	5	(1, ⊥, ⊥, ε, ε)
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# Quantum memory allocation

We work with

$$|T(S)\rangle = \sum_{\text{valid sequences } I} |T_I(S)\rangle .$$

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## Approach 2 : Memory tree

Maintain a tree with each memory cell as a leaf. Count for each node the number of free cells.

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## Aim

At the end of each walk, extract collisions and preserve a useful quantum data structure  $\leadsto$  new starting state of the next quantum walk.

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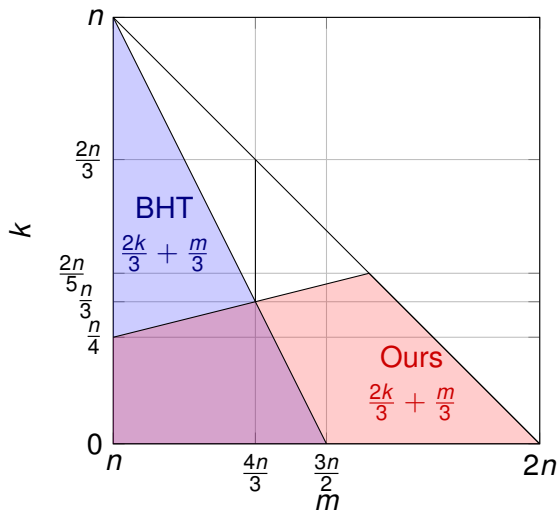
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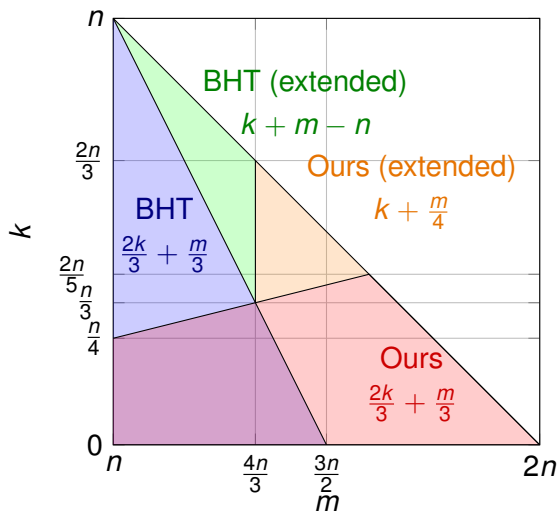
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**Complexity:**  $\tilde{O}(2^\ell + 2^k 2^{m/2 - \ell/2}) = \tilde{O}(2^{k+m/2 - \ell/2})$  where  $\ell \leq \min(2k/3 + m/3, m/2)$ .

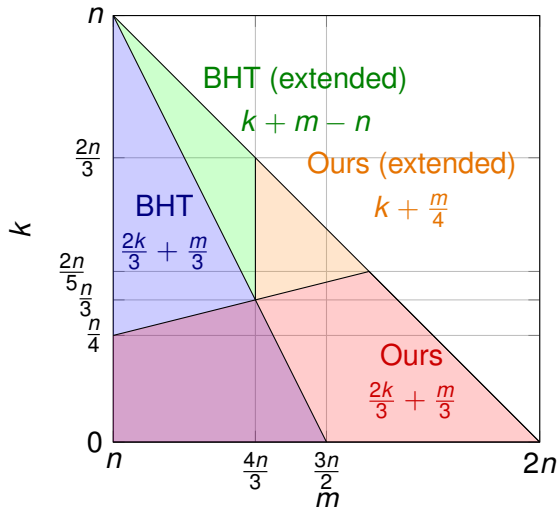
# Results



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Many applications:

- ▶ Quantum impossible differentials
- ▶ Quantum lattice sieving  $2^{0.2570d} \rightarrow 2^{0.2563d}$