

Improved Classical and Quantum Algorithms for Subset-Sum

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Outline

- 1 Introduction
- 2 Representations
- 3 Subset-Sum with Quantum Search
- 4 Subset-Sum with Quantum Walks

Introduction

The Subset-Sum Problem

Problem

Given: $\mathbf{a} = (a_1, \dots, a_n)$ a vector of ℓ -bit integers, and an ℓ -bit target S , find $\mathbf{e} = (e_1, \dots, e_n) \in \{0, 1\}^n$ such that $\mathbf{e} \cdot \mathbf{a} = \sum_i e_i a_i = S \pmod{2^\ell}$.

- The decision version is NP-complete
- The low-density case ($\ell \gg n$) is related to lattice SVP
- The high-density case ($\ell \ll n$) is solvable efficiently
- The density-1 case ($\ell \simeq n$) is hard

Subset-sums in (post-quantum) cryptography

- Repeatedly used as a hard problem for post-quantum cryptography ^a
- Similar techniques that we will see in this presentation apply to other problems (generic decoding algorithms) ^b
- Solving subset-sums is also useful in quantum hidden shift algorithms ^c

^aLyubashevsky, Palacio, and Segev, “Public-Key Cryptographic Primitives Provably as Secure as Subset Sum”, TCC 10

^bKachigar and Tillich, “Quantum Information Set Decoding Algorithms”, PQCrypto 17

^cBonnetain, *Improved Low-qubit Hidden Shift Algorithms*, 2019

The **random** Subset-Sum Problem

Problem

Given: $\mathbf{a} = (a_1, \dots, a_n)$ a vector of **n -bit integers**, and an **n -bit target S** , find $\mathbf{e} = (e_1, \dots, e_n) \in \{0, 1\}^n$ such that $\mathbf{e} \cdot \mathbf{a} = \sum_i e_i a_i = S \pmod{2^n}$; **where \mathbf{a}, S are selected uniformly at random**.

- Classical and quantum algorithms run in time $\tilde{O}(2^{\beta n})$: we are interested in the value of β
- In this talk, we optimize the **time exponent** (not the memory)
- We consider w.l.o.g. that \mathbf{e} has Hamming weight $\frac{n}{2}$

Classical algorithms

The time is $\tilde{O}(2^{\beta n})$.

Technique	β	Ref.
MIM	0.5	HS74 (Slide 8)
4-list merge	0.5	SS81 (Slide 9)
$\{0, 1\}$	0.3370	HGJ10 (Slide 18)
$\{-1, 0, 1\}$	0.2909	BCJ11 (Slide 23)
$\{-1, 0, 1\} + \text{NN}$	0.287	Ilya Ozerov's PhD thesis
$\{-1, 0, 1, 2\}$	0.283	Ours

Classical algorithm: meet-in-the-middle

$$\text{Cut the solution } \mathbf{e} = \left(\underbrace{0 \dots 0}_{n/2 \text{ bits}} \mid \underbrace{* \dots *}_{n/2 \text{ bits}} \right) + \left(\underbrace{* \dots *}_{n/2 \text{ bits}} \mid \underbrace{0 \dots 0}_{n/2 \text{ bits}} \right)$$

$$\underbrace{\hspace{10em}}_{2^{n/2} \text{ choices: } L_l} \quad \underbrace{\hspace{10em}}_{2^{n/2} \text{ choices: } L_r}$$

Then find $\mathbf{e}_l \in L_l, \mathbf{e}_r \in L_r$ s.t. $\mathbf{e}_l \cdot \mathbf{a} = -\mathbf{e}_r \cdot \mathbf{a} + S \pmod{2^n}$.

Complexities

Time: $\mathcal{O}(2^{n/2})$ (best worst-case time); **Memory:** $\mathcal{O}(2^{n/2})$

Horowitz and Sahni, "Computing Partitions with Applications to the Knapsack Problem", J. ACM

Schroeppel and Shamir's 4-list merging

By cutting in 4 instead of 2, we can decrease the memory to $2^{n/4}$.

$$\mathbf{e} = \underbrace{(*|0|0|0)}_{\mathbf{e}_0 \in L_0} + \underbrace{(0|*|0|0)}_{\mathbf{e}_1 \in L_1} + \underbrace{(0|0|*|0)}_{\mathbf{e}_2 \in L_2} + \underbrace{(0|0|0|*)}_{\mathbf{e}_3 \in L_3}$$

We now look for $(\mathbf{e}_0, \mathbf{e}_1, \mathbf{e}_2, \mathbf{e}_3) \in L_0 \times L_1 \times L_2 \times L_3$ s.t.

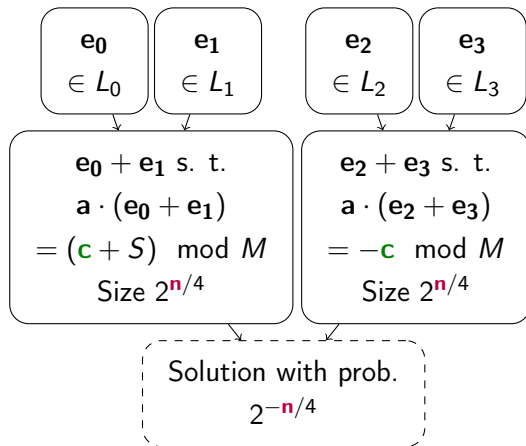
$$\mathbf{a} \cdot (\mathbf{e}_0 + \mathbf{e}_1 + \mathbf{e}_2 + \mathbf{e}_3) = S \pmod{2^n}$$

\Rightarrow there is still one solution among $2^{n/4} \times 2^{n/4} \times 2^{n/4} \times 2^{n/4}$.

Schroeppel and Shamir, "A $T = O(2^{n/2})$, $S = O(2^{n/4})$ Algorithm for Certain NP-Complete Problems", SIAM 81

Schroeppel and Shamir's 4-list merging (ctd.)

- Choose an $n/4$ -bit number c mod M
- Repeat for every value of c :



Complexities

Time: $\mathcal{O}(2^{n/4} \times 2^{n/4})$

Memory: $\mathcal{O}(2^{n/4})$

Representations

Breaking the $2^{n/2}$ bound

- When we have **one** solution among 2^n tuples, we don't know of any better time than $2^{n/2}$
- The idea of Howgrave-Graham and Joux (HGJ): cut e with respect to its Hamming weight

Suppose that e is of weight $n/2$ (worst case). Write for example:

$$\underbrace{e}_{\text{Weight } n/2} = \underbrace{e_0}_{\text{Weight } n/8} + \underbrace{e_1}_{\text{Weight } n/8} + \underbrace{e_2}_{\text{Weight } n/8} + \underbrace{e_3}_{\text{Weight } n/8}$$

notice that: $\binom{n}{n/8}^4 \simeq 2^{2.174n} \ggg \binom{n}{n/2} \simeq 2^n$: many solution tuples!

Howgrave-Graham and Joux, "New Generic Algorithms for Hard Knapsacks", EC 10

Notations

We introduce **distributions** and **weight constraints**.

Distributions

$\mathbf{e} \in D^n[\alpha]$ if \mathbf{e} contains αn “1” and $(1 - \alpha)n$ “0”.

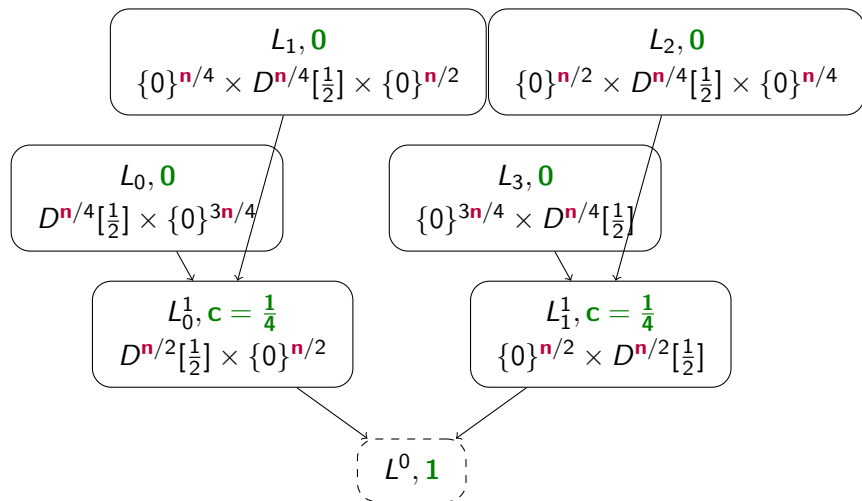
Then if $\mathbf{e}_1 \in D^n[\alpha_1]$, $\mathbf{e}_2 \in D^n[\alpha_2]$ we have $\mathbf{e}_1 + \mathbf{e}_2 \in D^n[\alpha_1 + \alpha_2]$ **with some probability** (to be continued).

Weight constraints

\mathbf{e} has “a cn -bit weight constraint” if we are able to constrain the (knapsack) weight of \mathbf{e} as $\mathbf{e} \cdot \mathbf{a} = s \pmod M$ for (previously) chosen cn -bit integers M and s .

If \mathbf{e}_1 has a cn -bit-weight cons. and \mathbf{e}_2 has a cn -bit-weight cons., then $\mathbf{e}_1 + \mathbf{e}_2$ as well by linearity (but the precise moduli don't matter!)

Example: Schroeppe-Shamir



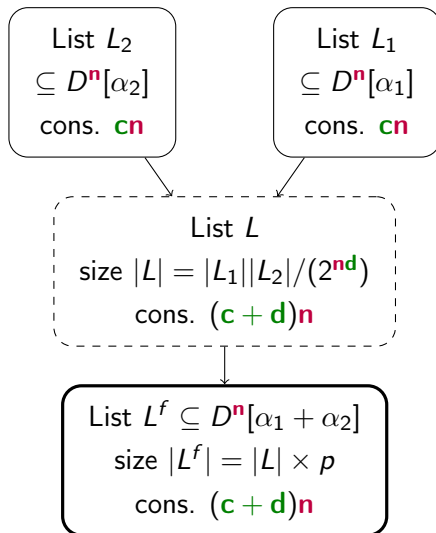
Generic layout

Solution

The solution \mathbf{e} is the only vector with an n -bit weight constraint and $\mathbf{e} \in D^n[1/2]$.

- We start from vectors \mathbf{e} with a 0-bit weight constraint and distributions $D^n[\alpha]$
- We sum them, trying to increase the weight constraint (“merging”)
- Eventually we get to a n -bit weight constraint and a distribution $D^n[1/2]$

Merging and filtering



Step 1: merging

Find pairs with more constrained weights.

We produce L in time $\max(\min(|L_1|, |L_2|), |L|)$.

Step 2: filtering

Keep only the $\mathbf{e}_1 + \mathbf{e}_2$ that conform to the expected distribution.

p is the “filtering probability” for: $D^n[\alpha_1] \times D^n[\alpha_2] \rightarrow D^n[\alpha_1 + \alpha_2]$

Merging and filtering (ctd.)

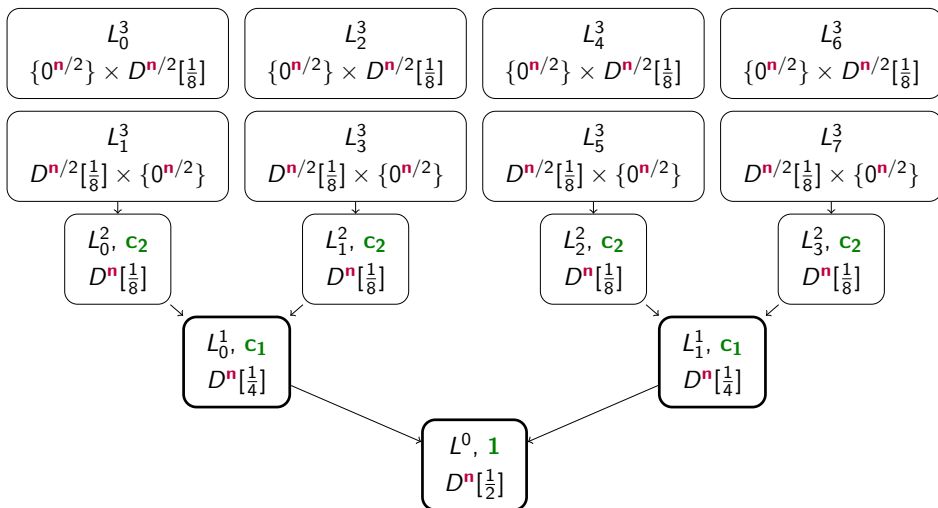
Heuristic

The vectors in L^f are uniformly distributed in $D^{\mathbf{n}}[\alpha_1 + \alpha_2]$.

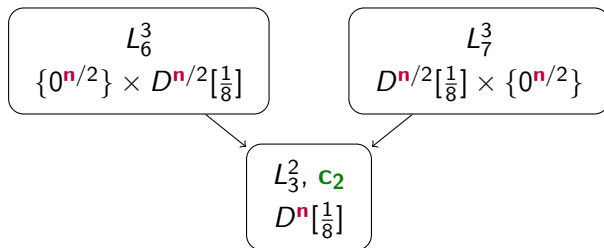
Approximations

- Representation sets have size: $D^{\mathbf{n}}[\alpha] = \binom{\mathbf{n}}{\alpha \mathbf{n}} \simeq 2^{h(\alpha)\mathbf{n}}$
- $h(x) = -x \log x - (1-x) \log(1-x)$
- In general we have a filtering probability
 $D^{\mathbf{n}}[\alpha_1] \times D^{\mathbf{n}}[\alpha_2] \rightarrow D^{\mathbf{n}}[\alpha_1 + \alpha_2]$ of: $\simeq 2^{(h(\alpha_1/(1-\alpha_2)) - h(\alpha_1))\mathbf{n}}$

The HGJ algorithm

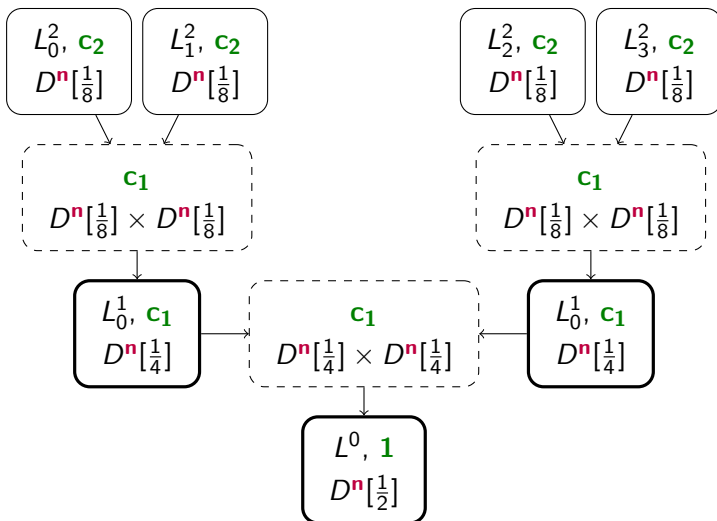


HGJ step 3: left-right split with a modulus



At this level we can afford to **merge without filtering**: find vectors $\mathbf{e} \in D^n[\frac{1}{8}]$ with a **c_2** -weight constraint (on $\mathbf{e} \cdot \mathbf{a}$).

HGJ step 2 and 1: merge and filter



An optimization problem

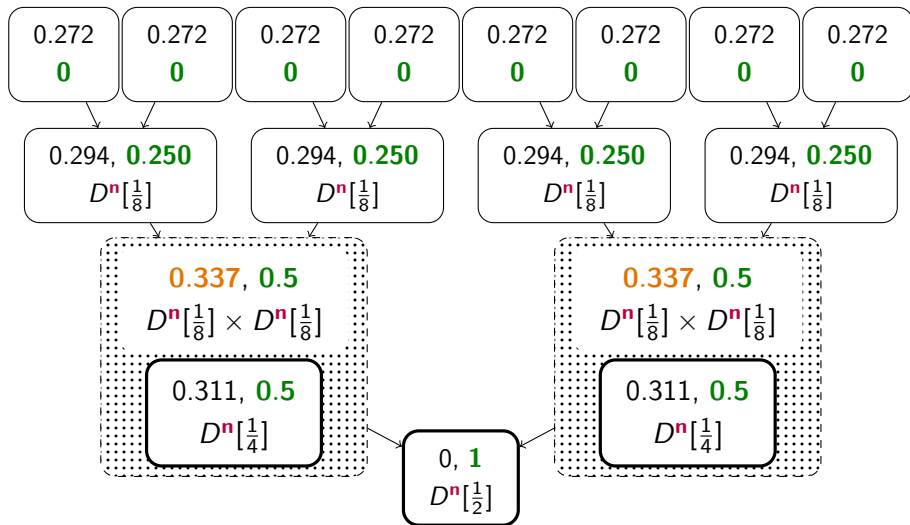
We write all parameters in \log_2 , relative to \mathbf{n} :

$$\left\| \begin{array}{l} |D^{\mathbf{n}/2}[\frac{1}{8}]| \\ |L_i^j| \\ |L_1||L_2|/(2^{\mathbf{n}\mathbf{d}}) \\ |L| \times p \end{array} \right\| \left\| \begin{array}{l} h(1/8)/2 \simeq 0.2718 \\ \ell_i^j \\ \ell_1 + \ell_2 - d \\ \ell + p\mathbf{f} \end{array} \right.$$

We compute the filtering probabilities:

- $D^{\mathbf{n}}[0, \frac{1}{8}] \times D^{\mathbf{n}}[0, \frac{1}{8}] \rightarrow D^{\mathbf{n}}[0, \frac{1}{4}]$: $2^{-0.02585\mathbf{n}}$
- $D^{\mathbf{n}}[0, \frac{1}{4}] \times D^{\mathbf{n}}[0, \frac{1}{4}] \rightarrow D^{\mathbf{n}}[0, \frac{1}{2}]$: $2^{-0.12256\mathbf{n}}$

Optimized parameters for HGJ



Better representations: BCJ

Sample distributions with $\{-1, 0, 1\}$.

- Of course we still need to obtain $D^n[\frac{1}{2}]$ in the end
- The “-1” need to be canceled out by “1”
- The “-1” shouldn’t sum up to “-2”!
- More parameters, new filtering probabilities
- Improvement on the time exponent: $0.291 < 0.337$

Becker, Coron, and Joux, “*Improved Generic Algorithms for Hard Knapsacks*”, EC 11

New results

Idea 1: do not saturate the lists

Starting lists are **not equal** to $D^{n/2}[*]$ but **sampled** u.a.r. from it.

BCJ without saturation: $0.289 < 0.291$

Idea 2: still better representations

Why stop at “-1”? Add some “2”.

- Of course we still need to obtain $D^n[\frac{1}{2}]$ in the end
- Some “1” can sum up to “2” (but not too much)
- A “-1” and a “2” give a “1”

BCJ without saturation and with “2”: $0.283 < 0.289$

Bonnetain et al., *Improved Classical and Quantum Algorithms for Subset-Sum*, ePrint 2020/168

Going Quantum

Classical search

$$\text{Let } \underbrace{X}_{\substack{\text{Search space,} \\ \text{size } N}} = \underbrace{G}_{\substack{\text{Good ones,} \\ \text{size } T}} \cup \underbrace{B}_{\substack{\text{Bad ones, size} \\ N - T}}$$

Let *Sample* and *Test* be functions to *sample* x from X and *test* if $x \in G$, in time t_{Sample} and t_{Test} .

There exists a function Sample_G that samples from G in time:

$$\frac{N}{T} (t_{\text{Sample}} + t_{\text{Test}})$$

\Rightarrow we transform a sampling procedure for the “search space” into a sampling procedure for the “solution space”.

Quantum search (amplitude amplification)

$$\underbrace{X}_{\substack{\text{Search space,} \\ \text{size } N}} = \underbrace{G}_{\substack{\text{Good ones,} \\ \text{size } T}} \cup \underbrace{B}_{\substack{\text{Bad ones, size} \\ N - T}}$$

Let QSample and QTest be quantum algorithms to **quantumly sample** X and **quantumly test if** $x \in G$, in time t_{Sample} and t_{Test} .

There exists an algorithm QSample_G that samples G in time:

$$\sqrt{\frac{N}{T}} (t_{\text{QSample}} + t_{\text{QTest}})$$

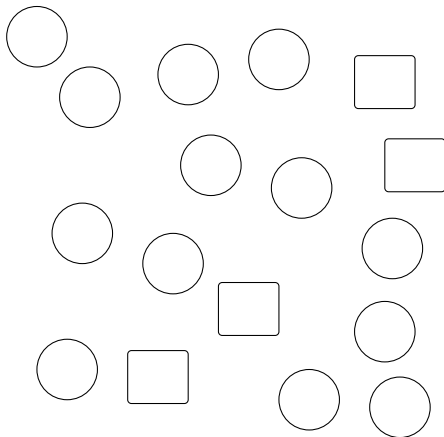
Quantum test: means testing **any** $x \in X$ **in superposition**

Quantum sample: means producing a uniform superposition of X

Brassard et al., “Quantum amplitude amplification and estimation”, 2002

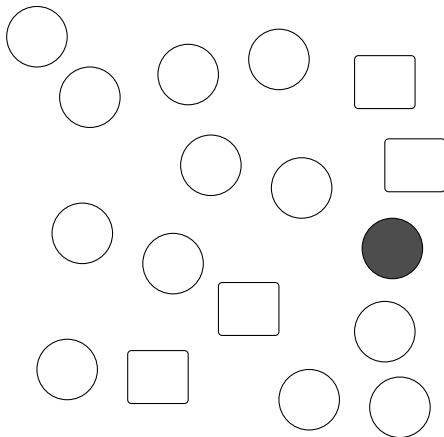
Classical search vs. quantum search

In the classical realm, we test elements x at random until we have found (a random) $x \in G$.



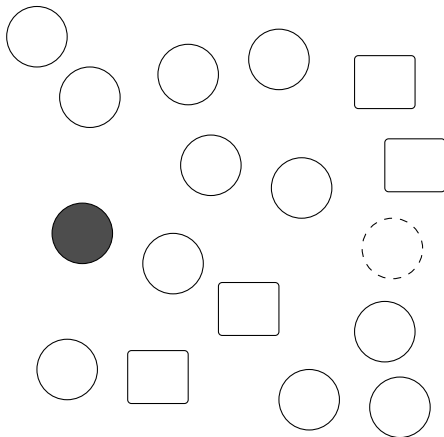
Classical search vs. quantum search

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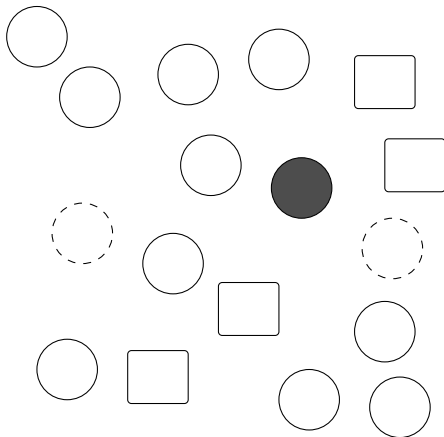
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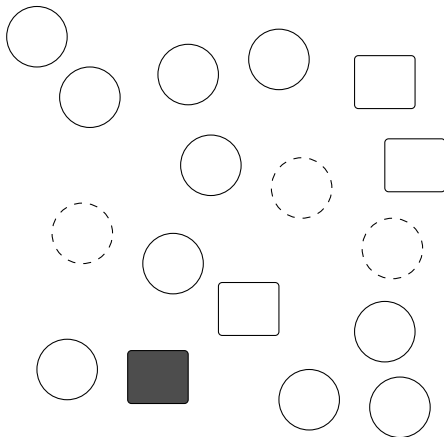
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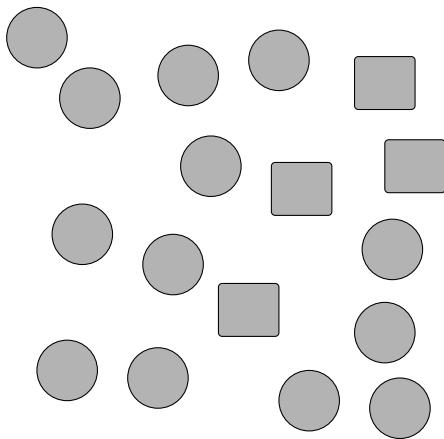
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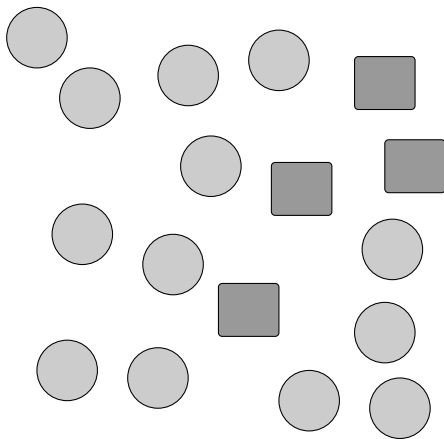
Classical search vs. quantum search

In the quantum realm, we move globally from X to G .



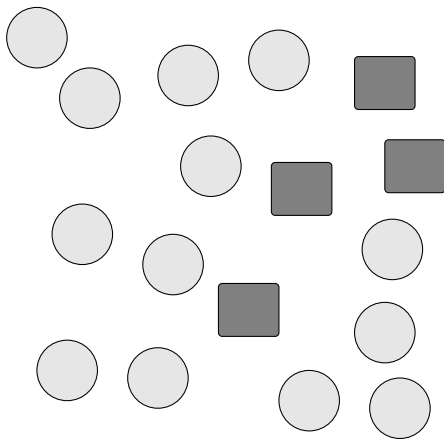
Classical search vs. quantum search

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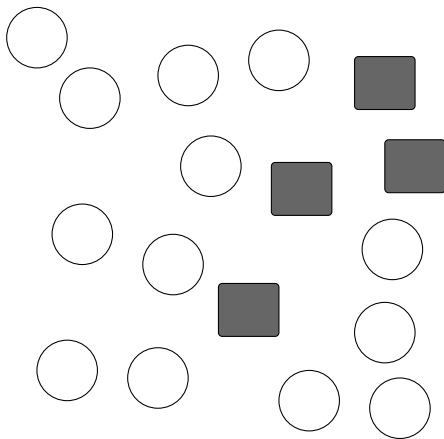
Classical search vs. quantum search

In the quantum realm, we move globally from X to G .



Classical search vs. quantum search

In the quantum realm, we move globally from X to G .



Interlude: quantum memory models

What happens if X is in memory?

Classical sample: only reads a single $x \in X$ (easy)

Quantum sample: must read all of X in superposition (maybe not easy): this is **quantum random access**

	Quantum random access	
Classical write	Classical memory quantum random access QRACM	\Rightarrow This section
Quantum write	Quantum memory quantum random access QRAQM	Previous quantum subset-sum algos

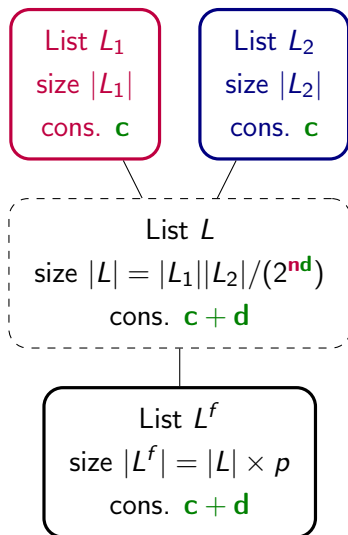
Quantum algorithms for subset-sum

The time is $\tilde{O}(2^{\beta n})$.

Technique	β	Ref.	Classical version	Model
MIM	0.3334	BHT98	HS74	QRACM
4-list merge	0.3	BJLM13	SS81	QRAQM
$\{0, 1\}$	0.241	BJLM13	HGJ10	QRAQM + conj.
$\{0, 1\}$	0.2356	Ours	HGJ10	QRACM
$\{-1, 0, 1\}$	0.226	HM18	BCJ11	QRAQM + conj.
$\{-1, 0, 1, 2\}$	0.2156	Ours		QRAQM + conj.
$\{-1, 0, 1, 2\}$	0.2182	Ours		QRAQM

Subset-Sum with Quantum Search

“Sampling-and-filtering”



Let's separate:

- **The sampled list**
- **The intermediate list**

We turn samples from L_1 into:

- Samples from L :

$$t_{\text{Sample}(L)} = \max\left(\frac{2^{nd}}{|L_2|}, 1\right) t_{\text{Sample}(L_1)}$$

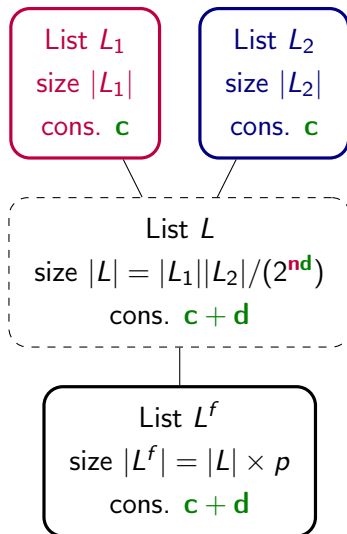
(find an element with same modulus)

- Samples from L^f :

$$t_{\text{Sample}(L^f)} = \frac{1}{p} t_{\text{Sample}(L)}$$

(wait until the filter is passed)

Quantum “sampling-and-filtering”



Assume that we have quantum samples from L_1 .

Then we have:

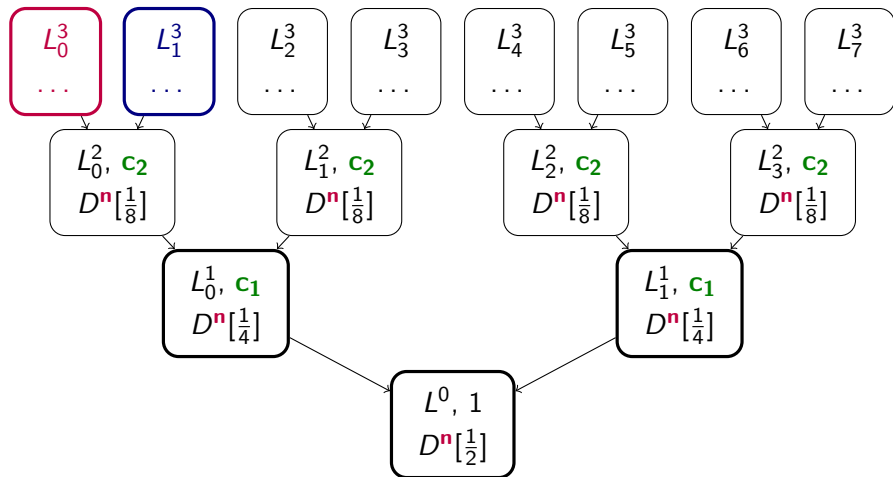
- Quantum samples from L :

$$t_{\text{QSample}(L)} = \max \left(\sqrt{\frac{2^{nd}}{|L_2|}}, 1 \right) t_{\text{QSample}(L_1)}$$

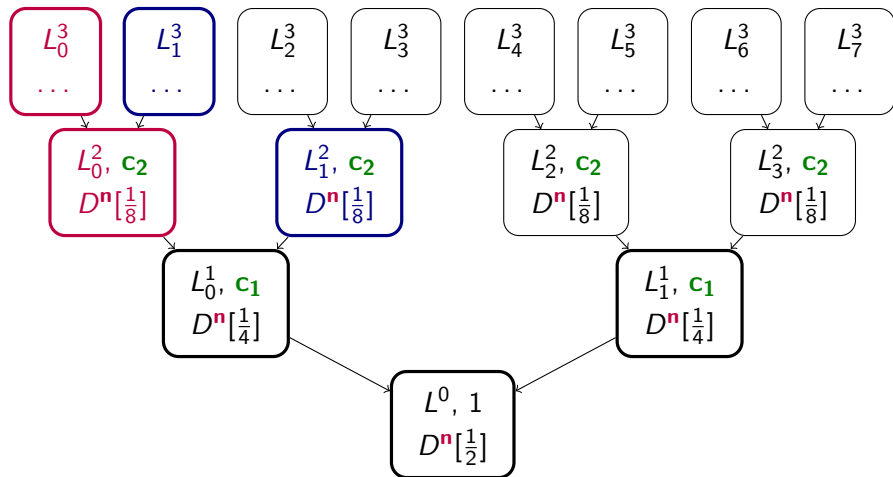
- Quantum samples from L^f :

$$t_{\text{Sample}(L_f)} = \sqrt{\frac{1}{p}} t_{\text{Sample}(L)}$$

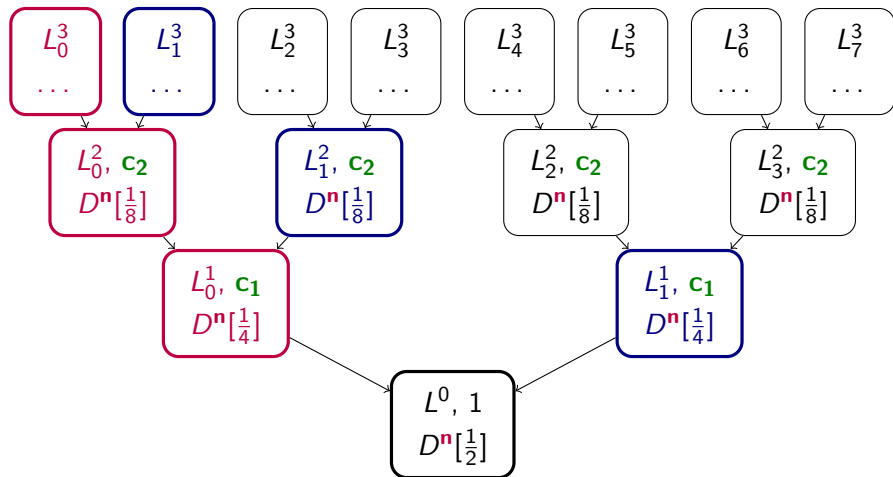
HGJ in the “sampling” framework



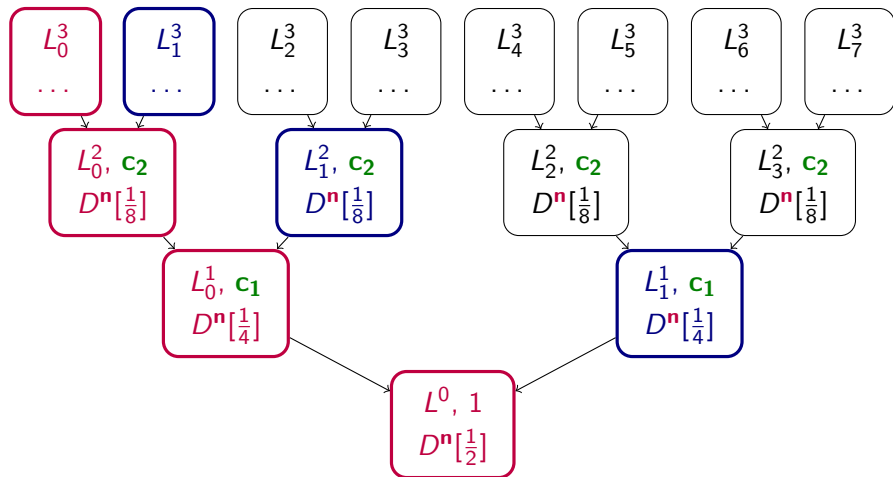
HGJ in the “sampling” framework



HGJ in the “sampling” framework



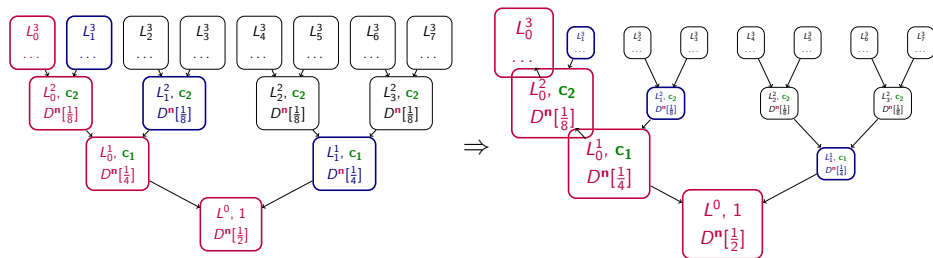
HGJ in the “sampling” framework



Using quantum search

Quantum search will square-root the sampling time of L^0 . But it's useless if the intermediate lists cost the same as before.

⇒ we make the tree unbalanced.



Details and result

- Unbalanced left-right split of L_0^3 and L_1^3 , unbalanced weights for the lists
- L_1^3, L_1^2, L_1^1 are intermediate lists stored in QRACM (classical data with quantum random access)
- **only $\text{poly}(n)$ quantum storage needed**

$0.226 \text{ (HM18)} <$

We use only $\{0, 1\}$
representations

0.2356

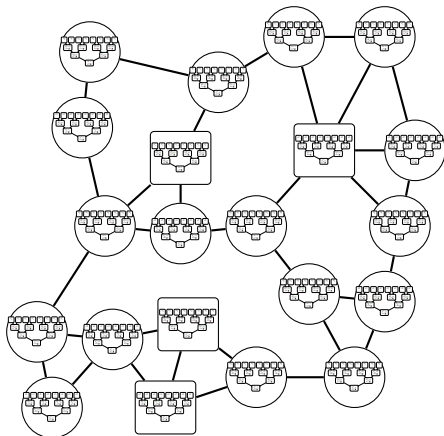
$< 0.241 \text{ (BJLM13)}$

We filter more efficiently

Subset-Sum with Quantum Walks

A classical walk for HGJ

Reduce the HGJ merging tree to a smaller tree, with smaller starting lists. Now L^0 does not always contain a solution.



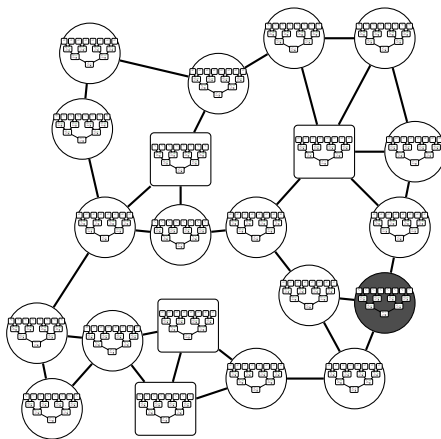
Walking on the graph

We use a (regular, undirected) Johnson graph $J(D, L)$.

- A vertex contains a product of 8 small lists $L_i^3, 0 \leq i \leq 7$, of size $|L_i^3| = L$, chosen among distributions $|D^i| = D$, **and** the whole tree built from these lists.
- There are $\binom{D}{L}^8$ vertices.
- Some vertices are **marked**: they contain the knapsack solution.
- We go from one to another by changing an element in a list L_i^3 **and** updating the tree.

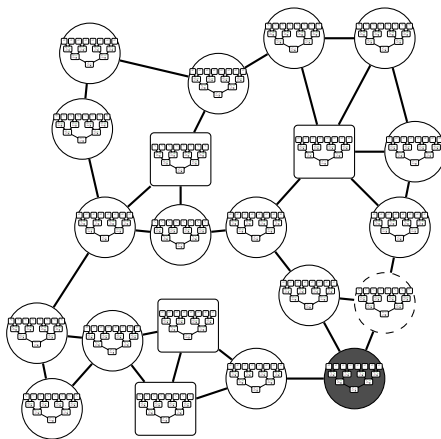
Classical random walk

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



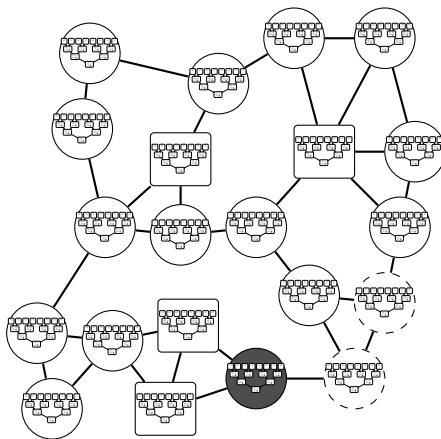
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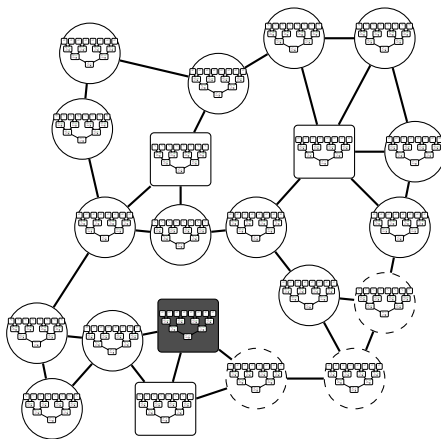
Classical random walk

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



Classical random walk

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



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)

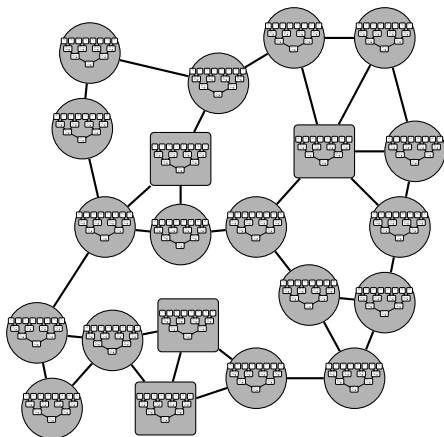
We will find a marked vertex in time:

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

where $\frac{1}{\delta}$ is the number of updates before we reach a new uniformly random vertex. In a Johnson graph $J(D, L)$, $\frac{1}{\delta} \simeq L$. (We need to replace all elements.)

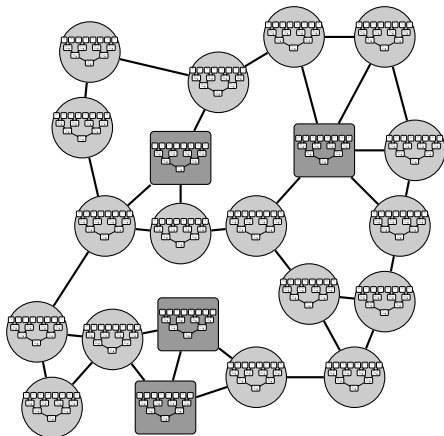
Quantum walk

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



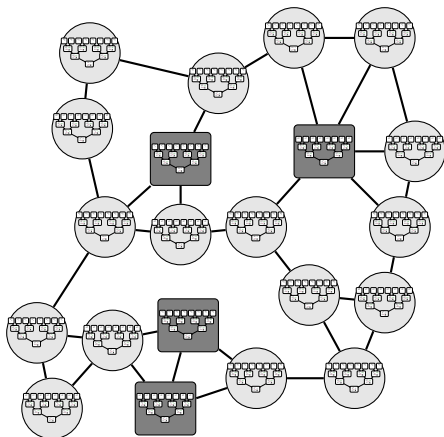
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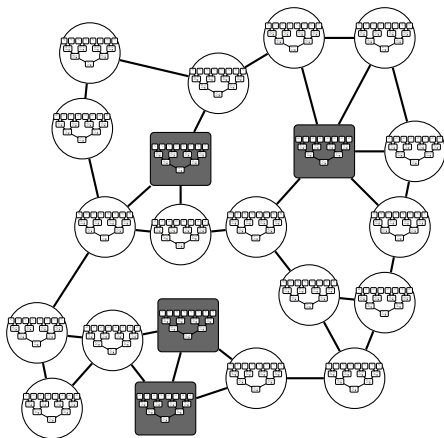
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As in quantum search, the walk transforms a uniform superposition over the whole graph into a superposition over marked vertices.



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!

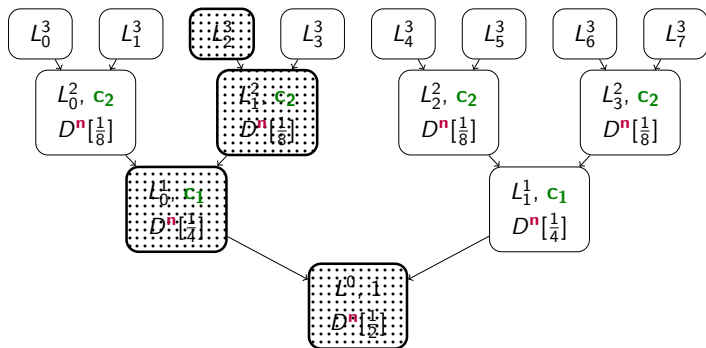
$$S + \underbrace{\sqrt{\frac{1}{\epsilon}}}_{\text{Walk steps}} \left(\underbrace{\sqrt{\frac{1}{\delta}} U}_{\text{Mixing time}} + C \right)$$

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

Magniez et al., “Search via quantum walk”, SIAM 11

Tracking the updates

- The update U must replace one element in a lower-level list **and** update the merging tree data structure.



- On average**, there is a single replacement to make at each level (no problem classically).

Superposition updates

- The update needs to take a fixed time.
- Since we are handling all vertices **and all edges** in superposition, there are cases when updating the tree would cost an exponential time.

Can we abort the bad cases?

Not in the MNRS framework: the data structure (the tree) must depend only on the vertex (the initial lists).

The quantum walk conjecture of Helm and May (TQC18)

With an update of expected time $\mathcal{O}(1)$, we can still do the runtime analysis as if it had an exact time $\mathcal{O}(1)$.

Helm and May, “*Subset Sum Quantumly in 1.17^n* ”, TQC 18

New results

- We have modified the **data structure** to guarantee the update time
- This reduces (marginally) the number of marked vertices

Fact

Previous quantum walks for subset-sum do not require the update conjecture.

- However, this data structure is not enough for our best algorithms ...

Summary of quantum walk results

Technique	Time	Ref.	Classical version	Model
$\{0, 1\}$	0.241	BJLM13	HGJ10	QRAQM + conj.
$\{-1, 0, 1\}$	0.226	HM18	BCJ11	QRAQM + conj.
$\{-1, 0, 1, 2\}$	0.2156	Ours		QRAQM + conj.
$\{-1, 0, 1, 2\}$	0.2182	Ours		QRAQM

Conclusion and open questions

On classical algorithms

More symbols and nearest-neighbor techniques should improve the exponent \Rightarrow how far could we go?

On quantum algorithms with quantum search

Better representations should improve the exponent ... if we manage to make the optimization converge.

Conclusion and open questions (ctd.)

On quantum walks

- The update conjecture can be removed from previous works ... but not completely from ours
- It seems that we still need to adapt the MNRS Quantum Walk framework

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Thank you!