

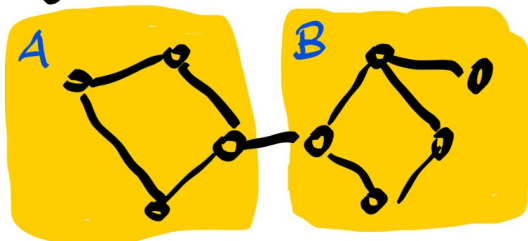
# A literature review of quantum parallelisation

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LIP6

2 OCT 2024

Divide and Conquer...  
Saleem, Tomesh, +4  
2022

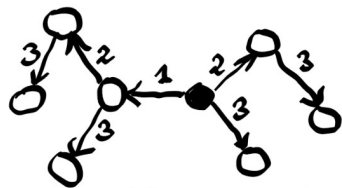
Figure I don't understand



It says: vertex = qubit  
edge = entanglement  
A, B = parallel qu. computers

In my understanding we capture only part of the picture. Ideally we need knot theory to describe entanglement.

Example  
○ = initial  $|0\rangle$   
● = initial  $\frac{1}{\sqrt{2}}(|0\rangle + |1\rangle)$   
○ → ● controlled NOT  
 $v := v \oplus u$   
 $t =$  time of operation



now qubits are all  $|0\rangle$  w. ampl.  $\frac{1}{\sqrt{2}}$   
and all  $|1\rangle$  w. ampl.  $\frac{1}{\sqrt{2}}$   
entanglement graph = clique  $\neq$  interaction graph

## Maximum Independent Set

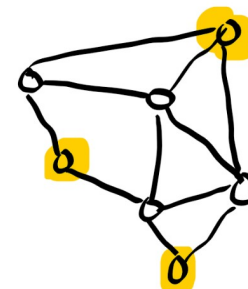
input  $G = (V, E)$

output  $S$  s.t.  $\forall u, v \in E$

$u \notin S$  or

$v \notin S$

objective  $\max |S|$



## State of the art for classical computers

easy for bipartite graphs  
interval graphs

hard to approximate (cannot decide if  $OPT \leq n^\epsilon$  or  $OPT \geq n^{1-\epsilon}$ )  
is also hard for quantum, so don't expect much

exact algorithm in  $1.1926^n \cdot n^{O(1)}$  [Xiao, Nagamochi 2017]

My suggestion: combine with quantum computing

## Quantum Algorithm

qubit = vertex

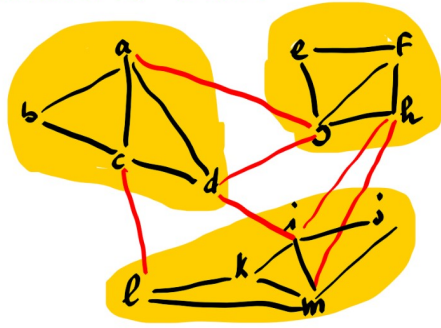
$|1\rangle =$  belong to solution set

constraint

$|0\rangle \rightarrow |1\rangle$  only possible if all neighbors are  $|0\rangle$

- 1) partition graph, each part  $\rightarrow$  dedicated quantum comp.
- 2) run quantum optimisation algorithm on each qu. comp.

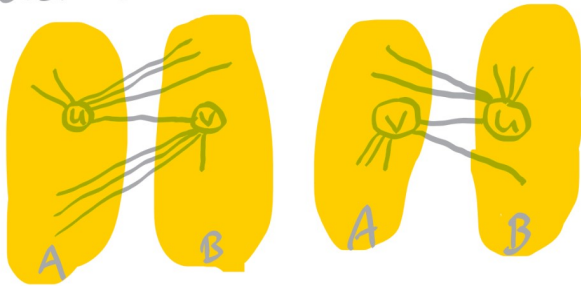
# Partition The Graph



Size of each part  $\leq$  # qbits per machine  
 Minimize # crossing edges

## Their choice

Kernighan-Lin 1970  
 partition  $V = A \cup B, |A|, |B| = \frac{|V|}{2}$   
 local search



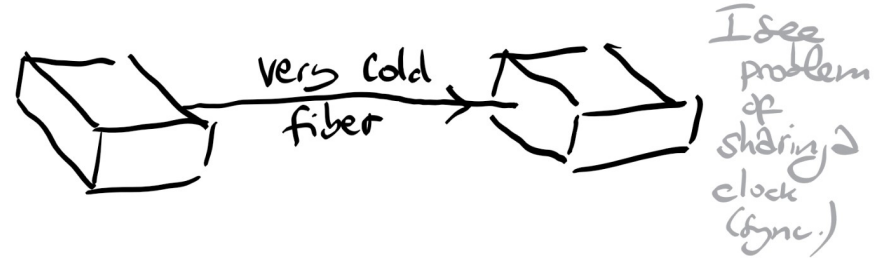
## My suggestion

Use integer linear programming formulation  
 Appropriate for small instances.  
 Easier than actual maximum independent set problem?

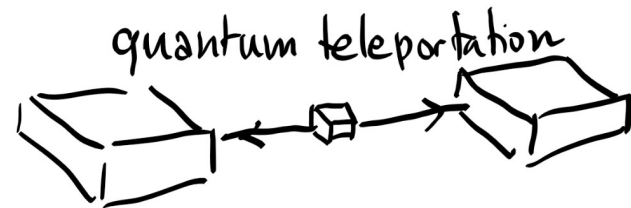
# Communication between Parts

technical aspect

either



or



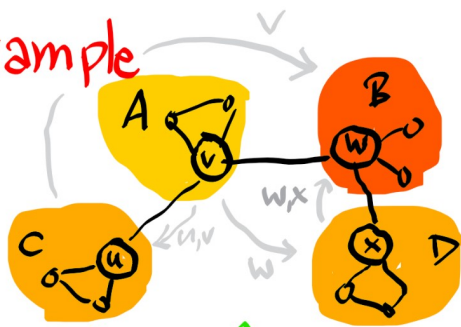
Need one additional qubit on each side.  
 I think we need **sequential** computation

- 1) run QC1 on vertices  $a, b, c, d, g, i, l$
- 2) teleport  $a, d$  from QC1 to QC2
- 3) run QC2 on  $a, d, e, f, h, i, m$
- 4) teleport  $c$  from QC1 to QC3
- 5) teleport  $d, h, i, m$  from QC2 to QC3
- 6) run QC3 on  $c, d, h, i, k, l, m$

I suspect, rather minimizing #cross edges,  
 we would like to minimize a different objective

- $d$  is teleported twice (degree)
- take cross edges into account for QC capacity 2/3

example



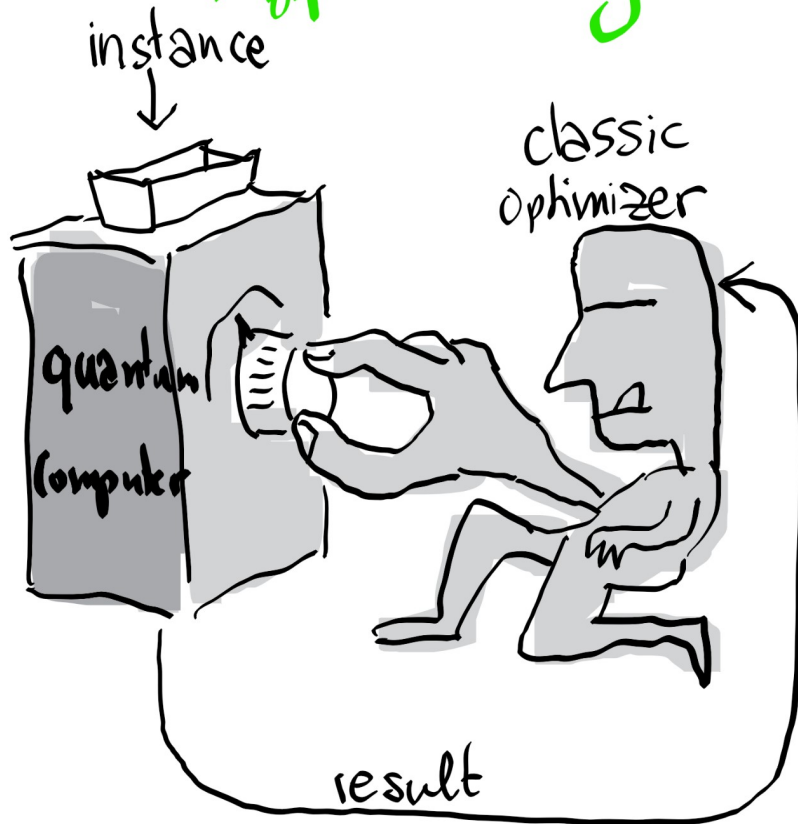
- execution order
- 1)  $AU\{u,v\}$
  - 2)  $CU\{v\} \parallel DU\{w\}$
  - 3)  $BW\{v,x\}$

## Quantum Alternating Operator Ansatz (Approach)

- invariant: all states with nonzero amplitude are feasible
- objective = Hamming Weight.
- Mixer operator on specific qubit  $|0\rangle \rightarrow |1\rangle$  only if all neighbors in state  $|0\rangle$ .
- Apply sequentially to all vertices in some arbitrary order

[...]

## Quantum Approximate Algorithm



## Their Experiments

- random graphs
- 3-regular
- 2-community
- 16 and 26 nodes
- high low high edge probability
- Simulated quantum computer
- Qiskit
- Compared different methods for partitioning the graph
- Compare with classical alg. for max. indep. set but not with state of the art

- alternate between
  - phase separation step (constraint in obj)
  - mixing step (pb indep.)

classic optimize time spend on each step