

# *Classical Simulation of Quantum Systems via Tensor Networks*

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# Quantum simulation

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- quantum systems have complex behavior
- want to *simulate* them, i.e. compute the outcome classically without actually building the system
- hard in the worst case, but there are systems for which this is feasible

# Tensors

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- drawn as a creature with a number of legs

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## Tensor networks

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- connecting two legs = contracting a common index  $i$

$$R_{x,y,z}^{a,b,c} = \sum_i P_{a,b,c,i} Q_{x,y,z,i}$$

- requires equal rank
- generalizes matrix multiplication
- can contract more than 1 leg at the same time

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  - an arbitrary quantum state is one fat spider with many legs
  - product states can be drawn as a group of skinnier creatures
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    - ⇒ fewer coefficients are needed!
- is there anything between?
  - for larger family of states
  - stil efficient

# Schmidt decomposition

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- every bipartite quantum state can be written as

$$|\phi\rangle = \sum_{i=1}^r |\psi_{A,i}\rangle |\psi_{B,i}\rangle,$$

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- hence we can slash any creature into two smaller ones connected by just **one** leg
  - notice that these legs may be longer

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- can this description possibly be **efficient**?
  - yes as long as the Schmidt ranks are not too high
  - then we need at most  $n \cdot R^3$  coefficients, where  $R = \max_e r_e$  is the maximal rank
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  - not every possible tensor networks yields efficient ranks!
- can apply unitaries and measurements fast on states with efficient networks
  - the tree structure is not altered much
  - hence we can simulate computation as long as all intermediate states are efficient

## Efficient tensor networks

- can we connect the  $n$  qubits by a 3-regular graph such that the rank of the worst bipartition is not too high?
- that is, optimize *Schmidt-rank width* defined as

$$\text{rwd}(|\psi\rangle) = \log \min_{\text{tree } T} \max_{\text{edge } e \in T} \chi_{A_T^e, B_T^e}(|\psi\rangle),$$

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- [S.-I. Oum, PhD thesis] polynomial time constant approximation algorithm for the width of every sub-modal function  $\chi$  (which is our case)
  - it is polynomial assuming that  $\chi_{A_T^e, B_T^e}$  is an oracle whose computation takes constant time



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  - for example,  $|\psi\rangle$  may be computed by a small quantum circuit  
this is hopeless, as it would solve factoring
  - works when  $|\psi\rangle$  is a cluster state, because then the Schmidt rank of a bipartition equals the  $\mathbb{GF}(2)$  rank of the adjacency matrix of this bipartition

## Cluster states

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- [Raussendorf & Briegel] *one-way quantum computer*
  - start in a highly entangled cluster state
  - perform a sequence of adaptive one-qubit measurements
  - universal for quantum computation

## How does cluster state computation work?

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- if we have a chain (cluster state corresponding to a path), then left-to-right one-qubit measurements in a certain basis *teleport* quantum information to the right **and** one can also perform some unitaries along the way

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- CPHASE gates can also be applied by incorporating them into the underlying cluster state
- this set of gates is universal  $\implies$  every quantum circuit can be efficiently rewritten into this form
  - the cluster state basically resembles the shape of the circuit

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  - on the other hand, there are graphs with constant rank width and tree width  $n$
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- when applied to factoring, the complexity lies in the modular exponentiation and the approximate QFT is easy

# Summary

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1. tensor networks
2. representation of quantum states
3. can find quickly the most efficient tensor network polynomial algorithm for representing cluster states
4. simulating general quantum circuits on cluster states
5. polynomial time simulation of quantum computation when the Schmidt-rank width is at most logarithmic