

Entanglement Partners\_ >





# Introduction to Tensor Network States and Methods





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### Some reviews

- J. Eisert, Modeling and Simulation 3, 520 (2013), arXiv:1308.3318
- N. Schuch, QIP, Lecture Notes of the 44th IFF Spring School 2013, arXiv:1306.5551
- R. Orus, Annals of Physics 349 (2014) 117-158
- R. Orus, Nature Reviews Physics 1, 538-550 (2019)
- J. I. Cirac, F. Verstraete, J. Phys. A: Math. Theor. 42, 504004 (2009)
- F. Verstratete, J. I. Cirac, V. Murg, Adv. Phys. 57,143 (2008)
- J. Jordan, PhD thesis, www.romanorus.com/JordanThesis.pdf
- G. Evenbly, PhD thesis, arXiv:1109.5424
- U. Schollwöck, RMP 77, 259 (2005)
- U. Schollwöck, Annals of Physics 326, 96 (2011)

### Outline

1) Basics 2) 1d MPS 3) 2d PEPS 4) Numerical algorithms 5) MERA 6) Extras

### Outline



# The main problem

• Aim: efficient representation of a quantum many-body state











Entanglement obeys area-law



#### key resource in quantum information

teleportation, quantum algorithms, quantum error correction, quantum cryptography...

#### Entanglement

#### 2d system





### key resource in quantum information

teleportation, quantum algorithms, quantum error correction, quantum cryptography...

 $\rho_{A} = \operatorname{tr}_{E}(|\Psi\rangle\langle\Psi|)$ 

$$S(A) = -\mathrm{tr}(\rho_A \log \rho_A)$$

Reduced density matrix of subsystem A

Entanglement entropy (von Neumann entropy)

For many ground states



#### Entanglement

2d system



### key resource in quantum information

teleportation, quantum algorithms, quantum error correction, quantum cryptography...

 $\rho_A = \operatorname{tr}_E(|\Psi\rangle\langle\Psi|)$ 

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Reduced density matrix of subsystem A

Entanglement entropy (von Neumann entropy)

 $S(A) \sim L$ 

 $(L > \xi)$ 

For many ground states

### In d dimensions

Generic  $S(A) \sim L^d$ state (volume) Ground states S(A) of (most) local Hamiltonians

 $S(A) \sim L^{d-1}$  (area)

Srednicki, Plenio, Eisert, Dreißig, Cramer, Wolf...

Locality of interactions 🔶 area-law

# Many-body Hilbert space is far too large

# Hilbert space is a convenient illusion

Hilbert space of a N-body many-body system



### Hilbert space is a convenient illusion

Hilbert space of a N-body many-body system

> , Set of area-law states Y. Ge, J. Eisert, arXiv:1411.2995

Set of TN states (low-energy eigenstates of local Hamiltonians)

Set of product states (mean field)

### Hilbert space is a convenient illusion

Hilbert space of a N-body many-body system



Most states here are not even reachable by a time evolution with a local Hamiltonian in polynomial time

Poulin, Qarry, Somma, Verstraete, PRL 106 170501 (2011) "Exploration" time ~  $O(10^{10^{23}})$  sec.

Compare to... Age of the universe ~  $O(10^{17})$  sec.

Set of area-law states Y. Ge, J. Eisert, arXiv:1411.2995

Set of TN states (low-energy eigenstates of local Hamiltonians)

Set of product states (mean field)

We need a language to target the relevant corner of quantum states directly

### **Tensor Networks**

### A new language

 $|\Psi\rangle = \sum_{i's} \Psi_{i_1 i_2 \dots i_N} \otimes |i_2\rangle \otimes \dots \otimes |i_N\rangle$ 







Tensors are local building blocks for the quantum state (like a DNA, or LEGO)



### **Tensor Networks**

e.g. RO, Annals of Physics 349 (2014) 117–158











Efficient O(poly(N)), satisfy area-law, low-energy eigenstates of local Hamiltonians

# **Two exact examples**





### Outline



### Outline



1) Basics

2) 1d MPS

3) 2d PEPS

4) Numerical algorithms

5) MERA

6) Extras


e.g., U. Schollwöck, Ann. of Phys. 326, 96 (2011)



 $|\Psi\rangle$ 

e.g., U. Schollwöck, Ann. of Phys. 326, 96 (2011)



 $|\Psi\rangle$ 







 $\left|out(\bar{\alpha})\right\rangle$  $\bar{\alpha} = 1, 2, ..., D^2$ 



$$\Psi \rangle = \sum_{\overline{\alpha}=1}^{D^2} |in(\overline{\alpha})\rangle |out(\overline{\alpha})\rangle$$
$$\overline{\alpha} = 1, 2, \dots, D^2$$



$$\Psi \rangle = \sum_{\overline{\alpha}=1}^{D^2} |in(\overline{\alpha})\rangle |out(\overline{\alpha})\rangle$$
$$\overline{\alpha} = 1, 2, ..., D^2$$

$$\rho_{in} = \operatorname{tr}_{out}\left(\left|\Psi\right\rangle\left\langle\Psi\right|\right) = \sum_{\overline{\alpha},\overline{\alpha}'} X_{\overline{\alpha},\overline{\alpha}'}\left|in(\overline{\alpha})\right\rangle\left\langle in(\overline{\alpha}')\right| \qquad X_{\overline{\alpha},\overline{\alpha}'} = \left\langle out(\overline{\alpha}')\right|out(\overline{\alpha})\right\rangle$$

$$rank(\rho_{in}) \le D^2$$
  $S(L) = -tr(\rho_{in} \log \rho_{in}) \le \log(D)2$ 



$$\Psi \rangle = \sum_{\overline{\alpha}=1}^{D^2} |in(\overline{\alpha})\rangle |out(\overline{\alpha})\rangle$$
$$\overline{\alpha} = 1, 2, ..., D^2$$

$$\rho_{in} = \operatorname{tr}_{out}\left(\left|\Psi\right\rangle \left\langle \Psi\right|\right) = \sum_{\overline{\alpha},\overline{\alpha}'} X_{\overline{\alpha},\overline{\alpha}'} \left|in(\overline{\alpha})\right\rangle \left\langle in(\overline{\alpha}')\right| \qquad X_{\overline{\alpha},\overline{\alpha}'} = \left\langle out(\overline{\alpha}') \left|out(\overline{\alpha})\right\rangle \right\rangle$$

$$rank(\rho_{in}) \le D^{2} \qquad S(L) = -tr(\rho_{in} \log \rho_{in}) \le \log(D)^{2}$$

$$prefactor \qquad size of the 0d boundary$$









 $O(pD^2)$ 







 $O(pD^2) + O(pD^3)$ 





 $O(pD^2) + O(pD^3)$ 





e.g., U. Schollwöck, Ann. of Phys. 326, 96 (2011)

 $O(pD^2) + O(pD^3) + O(pD^3)$ 

 $\langle \Psi | \Psi \rangle$ 



e.g., U. Schollwöck, Ann. of Phys. 326, 96 (2011)

 $\langle \Psi | \Psi \rangle$ 

 $O(pD^{2}) + O(pD^{3}) + O(pD^{3}) + ...$ 



... and so on...

Exact contraction in  $O(NpD^3)$  time, and the same for expectation values of local observables

### MPS corr. length is finite

 $C(r) = \left\langle O_i O_{i+r} \right\rangle - \left\langle O_i \right\rangle \left\langle O_{i+r} \right\rangle$ 

### MPS corr. length is finite











# MPS corr. length is finite

e.g., U. Schollwöck, Ann. of Phys. 326, 96 (2011)

$$C(r) = \left\langle O_i O_{i+r} \right\rangle - \left\langle O_i \right\rangle \left\langle O_{i+r} \right\rangle$$









 $C(r) \approx f(r)ae^{-r/\xi}$  $\xi \equiv -1/\log|\lambda_2/\lambda_1|$ 

#### MPS transfer matrix



# **MPS** canonical form

RO, G. Vidal, Phys. Rev. B 78, 155117 (2008)



# **Finding MPS canonical form**

RO, G. Vidal, Phys. Rev. B 78, 155117 (2008)



Bond indices correspond to orthonormal Schmidt basis

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### From MPS to PEPS

Matrix Product States (MPS)

 $\mathbf{e} \quad \mathbf{e} \quad$ 

1d systems

### But we want to go beyond 1d systems!!!





Very painful for DMRG...





### PEPS are not your friends...

### MPS



### PEPS are not your friends...

### MPS









### PEPS are not your friends...

### MPS













### PEPS obey 2d area-law



# PEPS obey 2d area-law L out $|\Psi\rangle$ in 1..D

### PEPS obey 2d area-law



### **PEPS obey 2d area-law**
























## **Critical correlation functions**

F. Verstraete et al, PRL 96, 220601 (2006)

$$|\Psi(\beta)\rangle = \frac{1}{\sqrt{Z(\beta)}} \exp\left(\frac{\beta}{2} \sum_{\langle i,j \rangle} \sigma_z^i \sigma_z^j\right)|+,+...+\rangle$$

Expectation values are those of the classical 2d Ising model

$$\left\langle \sigma_{z}^{r} \sigma_{z}^{r'} \right\rangle_{\beta} = \frac{1}{Z(\beta)} \sum_{\{s\}} s^{r} s^{r'} \exp\left(\beta \sum_{\langle i,j \rangle} s^{i} s^{j}\right) \qquad s = \pm 1$$

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$$\left\langle \sigma_{z}^{r} \sigma_{z}^{r'} \right\rangle_{\beta} = \frac{1}{Z(\beta)} \sum_{\{s\}} s^{r} s^{r'} \exp\left(\beta \sum_{\langle i,j \rangle} s^{i} s^{j}\right) \qquad s = \pm 1$$

It is a PEPS with D=2:

$$\frac{1}{|+\rangle} \frac{1}{1} = \left(\cosh(\beta/2)\right)^4 \qquad \frac{1}{|-\rangle} \frac{1}{1} = \left(\cosh(\beta/2)\right)^3 \left(\sinh(\beta/2)\right)$$

$$\frac{1}{|+\rangle} \frac{1}{1} = \left(\cosh(\beta/2)\right)^2 \left(\sinh(\beta/2)\right)^2 \qquad \frac{2}{|-\rangle} \frac{1}{1} = \left(\cosh(\beta/2)\right) \left(\sinh(\beta/2)\right)^3$$

$$\frac{2}{|+\rangle} \frac{1}{2} = \left(\sinh(\beta/2)\right)^4 \qquad + \text{ permutations}$$

At  $\beta_c = \left(\log(1+\sqrt{2})\right)/2$  the correlation length is infinite:  $\left<\sigma_z^r \sigma_z^r\right>_{\beta}$ 

$$_{\beta_c} \approx \frac{a}{\left|r-r'\right|^{1/4}}$$

#### One more thing about PEPS

#### They have no exact canonical form...



#### ...but there are approximate versions

*M. P. Zelatel, F. Pollmann, arXiv:1902.05100 R. Haghshenas, M. J. O'Rourke, G. K.-Lic Chan, arXiv:1903.03843* 

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#### **Tensor** Networks as an ansatz

#### Variational optimization

(e.g. DMRG)

$$\min_{|\Psi\rangle\in TN} \langle \Psi | H | \Psi \rangle / \langle \Psi | \Psi \rangle$$
ground states

e.g., S. White, PRL 69, 2863 (1992)

#### **Real/Imaginary time evolution**

(e.g. TEBD)

 $e^{-iHt}|\Psi
angle$ 

 $e^{-Ht}|\Psi
angle$ 

 dynamics
 ground states

 e.g., G. Vidal, PRL 91, 147902 (2003)

#### **Tensor** Networks as an ansatz

#### Variational optimization

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 $\min_{|\Psi\rangle \in TN} \langle \Psi | H | \Psi \rangle / \langle \Psi | \Psi \rangle$ ground states

e.g., S. White, PRL 69, 2863 (1992)

#### **Real/Imaginary time evolution**

(e.g. TEBD)

$$e^{-iHt} |\Psi
angle$$



 dynamics
 ground states

 e.g., G. Vidal, PRL 91, 147902 (2003)

- MPS methods for 1d are very efficient (e.g. DMRG obc  $\rightarrow$  D<sup>3</sup>)
- 2d PEPS ~ D<sup>10</sup>. But low D expected because of high connectivity, or entanglement monogamy. D=2 can be critical.
- Infinite lattices for translation-invariant systems
   (thermodynamic limit)
- Internal symmetries, fermionic systems, continuum limit (cMPS → quantum field theories)
- Limitation: amount and structure of entanglement



#### **Examples: DMRG and TEBD**

VOLUME 69, NUMBER 19

#### PHYSICAL REVIEW LETTERS

**9 NOVEMBER 1992** 

#### **Density Matrix Formulation for Quantum Renormalization Groups**

Steven R. White

Department of Physics, University of California, Irvine, California 92717 (Received 22 May 1992)

A generalization of the numerical renormalization-group procedure used first by Wilson for the Kondo problem is presented. It is shown that this formulation is optimal in a certain sense. As a demonstration of the effectiveness of this approach, results from numerical real-space renormalization-group calculations for Heisenberg chains are presented.

PACS numbers: 75.10.Jm, 02.70.+d, 05.30.-d

Efficient simulation of one-dimensional quantum many-body systems

Guifré Vidal<sup>1</sup>

<sup>1</sup>Institute for Quantum Information, California Institute of Technology, Pasadena, CA 91125, USA (Dated: February 1, 2008)

We present a numerical method to simulate the time evolution, according to a Hamiltonian made of local interactions, of quantum spin chains and systems alike. The efficiency of the scheme depends on the amount of the entanglement involved in the simulated evolution. Numerical analysis indicate that this method can be used, for instance, to efficiently compute time-dependent properties of lowenergy dynamics of sufficiently regular but otherwise arbitrary one-dimensional quantum many-body systems.

PACS numbers: 03.67.-a, 03.65.Ud, 03.67.Hk



Optimize over each tensor individually and sweep over the entire system (e.g., DMRG)



 $\min\left(\frac{\langle \Psi | H | \Psi \rangle}{\langle \Psi | \Psi \rangle}\right)$  Optimize over each tensor individually and sweep over the entire system (e.g., DMRG)





 $\min\left(\frac{\langle \Psi | H | \Psi \rangle}{\langle \Psi | \Psi \rangle}\right) \qquad \text{Optimize over each tensor individually and} \\ \text{sweep over the entire system (e.g., DMRG)}$ 



$$\frac{\partial}{\partial A^{*i}} \left( \langle \Psi | H | \Psi \rangle - \lambda \langle \Psi | \Psi \rangle \right) = 0$$

Minimization of quadratic function



 $\min\left(\frac{\langle \Psi | H | \Psi \rangle}{\langle \Psi | \Psi \rangle}\right)$  Optimize over each tensor individually and sweep over the entire system (e.g., DMRG)

mın



$$\Rightarrow \left| \Psi^{N} \right\rangle \xrightarrow{A^{N-1}} \left| \Psi^{N+1} \right\rangle \xrightarrow{A^{N-2}} \\ E^{N} \ge E^{N+1} \ge$$

$$\frac{\partial}{\partial A^{*i}} \left( \langle \Psi | H | \Psi \rangle - \lambda \langle \Psi | \Psi \rangle \right) = 0$$

Minimization of quadratic function

mın

 $\mathbf{H}_{eff}^{i} \vec{A}^{i} = \lambda \mathbf{N}^{i} \vec{A}^{i}$  Generalized eigenvalue problem Once  $\mathbf{H}_{eff}^{i}$  and  $\mathbf{N}^{i}$  are known, we can solve this problem efficiently  $\mathbf{H}_{_{
m eff}}^{i}$  and  $\mathbf{N}^{i}$  are exact in 1d, and approximate in 2d

# e.g. 2d calculation of $\mathbf{N}^i \vec{A}^i$

*F. Verstraete, I. Cirac, cond-mat/0407066* 

$$\frac{\partial}{\partial A^{*i}} \left( \left\langle \Psi \middle| H \middle| \Psi \right\rangle - \lambda \left\langle \Psi \middle| \Psi \right\rangle \right) = 0 \qquad \blacksquare$$



 $\mathbf{H}_{eff}^{i}\vec{A}^{i} = \lambda \mathbf{N}^{i}\vec{A}^{i}$ 









# **e.g.** 2d calculation of $\mathbf{N}^{i}\vec{A}^{i}$ F. Verstraete, I. Cirac, $\frac{\partial}{\partial A^{*i}}(\langle \Psi|H|\Psi\rangle - \lambda\langle \Psi|\Psi\rangle) = 0 \implies \mathbf{H}_{eff}^{i}\vec{A}^{i} = \lambda\mathbf{N}^{i}\vec{A}^{i}$



 $\frac{\partial}{\partial A^{*i}} \langle \Psi | \Psi \rangle = \mathbf{N}^{i} \vec{A}^{i}$ 

**e.g.** 2d calculation of 
$$\mathbf{N}^{i}\vec{A}^{i}$$
 F. Verstraete, I. Cirac,  
 $\frac{\partial}{\partial A^{*i}}(\langle \Psi|H|\Psi\rangle - \lambda\langle \Psi|\Psi\rangle) = 0$ 

$$\mathbf{MPS}$$

$$\frac{1..D^{2}}{MPO}$$

$$\frac{\partial}{\partial A^{*i}}\langle \Psi|\Psi\rangle = \mathbf{N}^{i}\vec{A}^{i}$$

1d problem: use a 1d method for MPS (e.g., DMRG or TEBD)

**e.g.** 2d calculation of 
$$\mathbf{N}^{i} \vec{A}^{i}$$
 F. Verstraete, I. Cirac,  
 $\partial = \partial A^{*i} (\langle \Psi | H | \Psi \rangle - \lambda \langle \Psi | \Psi \rangle) = 0$ 

$$\mathbf{H}^{i}_{eff} \vec{A}^{i} = \lambda \mathbf{N}^{i} \vec{A}^{i}$$

$$\frac{\partial}{\partial A^{*i}} \langle \Psi | \Psi \rangle = \mathbf{N}^{i} \vec{A}^{i}$$

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e.g. 2d calculation of 
$$\mathbf{N}^{i}\vec{A}^{i}$$
 F. Verstraete, I. Cirac,  
 $\partial_{\partial A^{*i}}(\langle \Psi|H|\Psi\rangle - \lambda \langle \Psi|\Psi\rangle) = 0$   $\longrightarrow$   $\mathbf{H}^{i}_{eff}\vec{A}^{i} = \lambda \mathbf{N}^{i}\vec{A}^{i}$   
 $\int_{\partial A^{*i}} \langle \Psi|\Psi\rangle = \mathbf{N}^{i}\vec{A}^{i}$   
 $\frac{\partial}{\partial A^{*i}} \langle \Psi|\Psi\rangle = \mathbf{N}^{i}\vec{A}^{i}$   
 $MPS_{up}$ 





#### **Canonical form helps!**

MPS with obc There is canonical form

 $\bigcap_{i=1}^{\Gamma_1} \sum_{\lambda_1} \sum_{\lambda_2} \sum_{\lambda_2} \sum_{i=1}^{\Gamma_3} \sum_{\lambda_3} \sum_{i=1}^{\Gamma_4} \sum_{i=1}^{$ 

$$\mathbf{H}^{i}_{eff}\vec{A}^{i} = \lambda \vec{A}^{i}$$

Normal eigenvalue problem. Very stable. PEPS, and TNs with loops There is no canonical form



$$\mathbf{H}_{eff}^{i}\vec{A}^{i} = \lambda \mathbf{N}^{i}\vec{A}^{i}$$

Generalized eigenvalue problem. Less stable. Approximate canonical forms.

### **Time evolution**

(real, imaginary)

### Time evolution (e.g. imaginary)

e.g. J. Jordan et al, PRL 101, 250602 (2008)

 $\left|\Psi_{0}\right\rangle = \lim_{\tau \to \infty} \frac{e^{-\tau H} \left|\Psi\right\rangle}{\left\|e^{-\tau H} \left|\Psi\right\rangle\right\|}$ 

Divide into small time-steps  $\delta \tau << 1$ 

#### Time evolution (e.g. imaginary)

e.g. J. Jordan et al, PRL 101, 250602 (2008)



Divide into small time-steps  $\delta \tau << 1$ 




e.g. J. Jordan et al, PRL 101, 250602 (2008)

Divide into small time-steps  $\delta \tau << 1$ 



Split the Hamiltonian (e.g. 2-body n.n.)

 $H = H_{hor}^{even} + H_{hor}^{odd} + H_{ver}^{even} + H_{ver}^{odd}$ 





e.g. J. Jordan et al, PRL 101, 250602 (2008)

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Divide into small time-steps  $\delta \tau << 1$ 



 $H = H_{hor}^{even} + H_{hor}^{odd} + H_{ver}^{even} + H_{ver}^{odd}$ 



e.g. J. Jordan et al, PRL 101, 250602 (2008)

 $e^{-\delta\tau H} \approx e^{-\delta\tau H_{hor}^{even}} e^{-\delta\tau H_{hor}^{odd}} e^{-\delta\tau H_{ver}^{even}} e^{-\delta\tau H_{ver}^{odd}} + O(\delta\tau^2)$ 

 $e^{-\delta au H_{hor}^{even}}$ 



e.g. J. Jordan et al, PRL 101, 250602 (2008)

$$e^{-\delta\tau H} \approx e^{-\delta\tau H_{hor}^{even}} e^{-\delta\tau H_{hor}^{odd}} e^{-\delta\tau H_{ver}^{even}} e^{-\delta\tau H_{ver}^{odd}} + O(\delta\tau^2)$$

$$e^{-\delta\tau H_{hor}^{even}} = \bigotimes_{\langle i,j \rangle} e^{-\delta\tau h_{ij}} = \bigotimes_{\langle i,j \rangle} g_{ij}$$

2-body gates



e.g. J. Jordan et al, PRL 101, 250602 (2008)

$$e^{-\delta\tau H} \approx e^{-\delta\tau H_{hor}^{even}} e^{-\delta\tau H_{hor}^{odd}} e^{-\delta\tau H_{ver}^{odd}} e^{-\delta\tau H_{ver}^{odd}} + O(\delta\tau^2)$$

$$e^{-\delta \tau H_{hor}^{even}} \left| \Psi \right\rangle = \left| \tilde{\Psi} \right\rangle$$

evolved state





Different approaches to this problem: (fast) full update, simplified update, TPVA... **Full update:**  $\min \left\| \tilde{\Psi} \right\rangle - \left| \Psi' \right\rangle \right\|^2$ 

Finite systems: optimize over all tensors in the PEPS (as before)

Infinite systems: optimize over tensors in the PEPS unit cell (iPEPS)

Require *calculations of environments*, like the one shown before.

### But... does it work?

### But... does it work?

# YES, it does

### + lots of other examples

J. Jordan, RO, G. Vidal, F. Verstraete, I. Cirac, PRL **101** 250602 (2008) P. Corboz, RO, B. Bauer, G. Vidal, PRB **81** 165104 (2010)

#### P. Corboz, PRB **93** 045116 (2016)



FIG. 4. (Color online) iPEPS energy of a period-5 stripe in the doped case in the strongly correlated regime (U/t = 8, n = 0.875) in comparison with other methods.

### Simulating a Quantum Computer

### With MPS

### Efficient classical simulation of slightly entangled quantum computations

Guifré Vidal<sup>1</sup>

<sup>1</sup>Institute for Quantum Information, California Institute of Technology, Pasadena, CA 91125, USA (Dated: February 1, 2008)

We present a scheme to efficiently simulate, with a classical computer, the dynamics of multipartite quantum systems on which the amount of entanglement (or of correlations in the case of mixed-state dynamics) is conveniently restricted. The evolution of a pure state of n qubits can be simulated by using computational resources that grow linearly in n and exponentially in the entanglement. We show that a pure-state quantum computation can only yield an exponential speed-up with respect to classical computations if the entanglement increases with the size n of the computation, and gives a lower bound on the required growth.

PACS numbers: 03.67.-a, 03.65.Ud, 03.67.Hk

## Simulating a Quantum Computer

### With PEPS

#### PHYSICAL REVIEW LETTERS 123, 190501 (2019)

#### General-Purpose Quantum Circuit Simulator with Projected Entangled-Pair States and the Quantum Supremacy Frontier

Chu Guo,<sup>1,\*</sup> Yong Liu<sup>(a)</sup>,<sup>2,\*</sup> Min Xiong,<sup>2</sup> Shichuan Xue,<sup>2</sup> Xiang Fu<sup>(a)</sup>,<sup>2</sup> Anqi Huang,<sup>2</sup> Xiaogang Qiang<sup>(a)</sup>,<sup>2</sup> Ping Xu,<sup>2</sup> Junhua Liu,<sup>3,4</sup> Shenggen Zheng,<sup>5</sup> He-Liang Huang,<sup>1,6,7</sup> Mingtang Deng,<sup>2</sup> Dario Poletti,<sup>8,†</sup> Wan-Su Bao,<sup>1,7,‡</sup> and Junjie Wu<sup>(a)</sup>,<sup>2,§</sup>

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<sup>8</sup>Science and Math Cluster and EPD Pillar, Singapore University of Technology and Design, 8 Somapah Road, 487372 Singapore

(Received 30 July 2019; published 4 November 2019)

Recent advances on quantum computing hardware have pushed quantum computing to the verge of quantum supremacy. Here, we bring together many-body quantum physics and quantum computing by using a method for strongly interacting two-dimensional systems, the projected entangled-pair states, to realize an effective general-purpose simulator of quantum algorithms. The classical computing complexity of this simulator is directly related to the entanglement generation of the underlying quantum circuit rather than the number of qubits or gate operations. We apply our method to study random quantum circuits, which allows us to quantify precisely the memory usage and the time requirements of random quantum circuits. We demonstrate our method by computing one amplitude for a  $7 \times 7$  lattice of qubits with depth (1 + 40 + 1) on the Tianhe-2 supercomputer.

DOI: 10.1103/PhysRevLett.123.190501

## **Simulating a Quantum Annealer**

### With PEPS

#### Heuristic optimization and sampling with tensor networks

Marek M. Rams,<sup>1</sup> Masoud Mohseni,<sup>2</sup> and Bartłomiej Gardas<sup>1,3</sup>

<sup>1</sup> Jagiellonian University, Marian Smoluchowski Institute of Physics, Lojasiewicza 11, 30-348 Kraków, Poland <sup>2</sup> Google AI Quantum, Venice, CA 90291 <sup>3</sup> University of Silesia, Institute of Physics, Bankowa 12, 40-007 Katowice, Poland



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### From MPS to MERA

Matrix Product States (MPS) 9
9
9
9
9
9
9
1
d systems

### But we want to do critical systems!!!



Also very painful for DMRG...











holographic dimension (RG)



### **Tensors obey constraints**







### Reason:

## entanglement is built locally at all length scales

# 

L



coarse-grain entangle locally



entangle locally coarse-grain entangle locally









Extra dimension defines an RG flow: Entanglement Renormalization



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## Norm of MERA

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## Entropy of 1d MERA





































# Only tensors inside of the causal cone contribute to the expectation value

# Comparison

	MPS in 1d ●-●-●-●	PEPS in 2d	MERA in 1d
Ent. entropy	S(L) = O(1)	S(L) = O(L)	$S(L) = O(\log L)$
Exact contraction	efficient	inefficient	efficient
Corr. length	finite	finite & infinite	finite & infinite
To/from	1d Ham.	2d Ham.	1d Ham.
Tensors	arbitrary	arbitrary	constrained

## A numerical example: 1d critical systems

G. Evenbly, G. Vidal, in "Strongly Correlated Systems. Numerical Methods", Springer, Vol. 176 (2013)



## Outline



1) Basics

2) 1d MPS

3) 2d PEPS

4) Numerical algorithms

5) MERA

6) Extras



## Outline



## **Machine Learning**

E.g., E. M. Stoudenmire, D. J. Schwab, Advances in Neural Information Processing Systems 29, 4799 (2016)

### **Supervised Kernel Learning**

#### Supervised Learning With Quantum-Inspired Tensor Networks

E. Miles Stoudenmire<sup>1,2</sup> and David J. Schwab<sup>3</sup>

<sup>1</sup>Perimeter Institute for Theoretical Physics, Waterloo, Ontario, N2L 2Y5, Canada <sup>2</sup>Department of Physics and Astronomy, University of California, Irvine, CA 92697-4575 USA <sup>3</sup>Dept. of Physics, Northwestern University, Evanston, IL (Dated: May 22, 2017)

Tensor networks are efficient representations of high-dimensional tensors which have been very successful for physics and mathematics applications. We demonstrate how algorithms for optimizing such networks can be adapted to supervised learning tasks by using matrix product states (tensor trains) to parameterize models for classifying images. For the MNIST data set we obtain less than 1% test set classification error. We discuss how the tensor network form imparts additional structure to the learned model and suggest a possible generative interpretation.

$$f^\ell(\mathbf{x}) = W^\ell \cdot \Phi(\mathbf{x})$$

$$C = rac{1}{2} \sum_{n=1}^{N_T} \sum_{\ell} (f^{\ell}(\mathbf{x}_n) - \delta^{\ell}_{L_n})^2$$

0

Minimize using TN methods

### **Anomaly detection**

### **Anomaly Detection with Tensor Networks**

Jinhui Wang\* Stanford University Stanford, CA 94305, USA wangjh97@stanford.edu

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#### Abstract

Originating from condensed matter physics, tensor networks are compact representations of high-dimensional tensors. In this paper, the prowess of tensor networks is demonstrated on the particular task of one-class anomaly detection. We exploit the memory and computational efficiency of tensor networks to learn a linear transformation over a space with dimension exponential in the number of original features. The linearity of our model enables us to ensure a tight fit around training instances by penalizing the model's global tendency to a predict normality via its Frobenius norm—a task that is infeasible for most deep learning models. Our method outperforms deep and classical algorithms on tabular datasets and produces competitive results on image datasets, despite not exploiting the locality of images.
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Originating from condensed matter physics, tensor networks are compact representations of high-dimensional tensors. In this paper, the prowess of tensor networks is demonstrated on the particular task of one-class anomaly detection. We exploit the memory and computational efficiency of tensor networks to learn a linear transformation over a space with dimension exponential in the number of original futures. The linearity of our model enables us to ensure a tight fit around training unces by penalizing the model's global tendency to a predict normality via obenius norm—a task that is infeasible for most deep learning models. Our method outperforms deep and classical algorithms on tabular datasets ind produces competitive results on image datasets, despite not exploiting the locality of images.

### **Beyond MPS**

#### Machine Learning by Unitary Tensor Network of Hierarchical Tree Structure

Ding Liu,<sup>1,2</sup> Shi-Ju Ran,<sup>3,2,\*</sup> Peter Wittek,<sup>4,5,6,7,†</sup> Cheng Peng,<sup>8</sup> Raul Blázquez García,<sup>2</sup> Gang Su,<sup>8,9</sup> and Maciej Lewenstein<sup>2,10</sup>

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The resemblance between the methods used in quantum-many body physics and in machine learning has drawn considerable attention. In particular, tensor networks (TNs) and deep learning architectures bear striking similarities to the extent that TNs can be used for machine learning. Previous results used one-dimensional TNs in image recognition, showing limited scalability and flexibilities. In this work, we train two-dimensional hierarchical TNs to solve image recognition problems, using a training algorithm derived from the multi-scale entanglement renormalization ansatz. This approach introduces mathematical connections among quantum many-body physics, quantum information theory, and machine learning. While keeping the TN unitary in the training phase, TN states are defined, which encode classes of images into quantum many-body states. We study the quantum features of the TN states, including quantum entanglement and fidelity. We find these quantities could be properties that characterize the image classes, as well as the machine learning tasks.

#### TN structure of some neural networks

Y. Levine et al, Phys. Rev. Lett. 122, 065301 (2019)

Recurrent neural networks are MPS (hidden Markov models)



#### Deep convolutional networks are TTNs



## **Symmetries**

e.g., P. Schmoll et al, Annals of Physics 419 (2020) 168232

### Symmetric tensors and Schur's lemma

e.g., S. Singh, R. N. C. Pfeifer, G. Vidal, PRA 82, 050301 (2010)



Structural part depends only on the group properties (intertwiners)

## **Emergent spin networks**

e.g., S. Singh, R. N. C. Pfeifer, G. Vidal, PRA 82, 050301 (2010)



Global and gauge symmetries are handled naturally



#### Symmetries are useful for numerics!

e.g., P. Schmoll, S. Singh, M. Rizzi, RO, arXiv:1809.08180

Canonical example: spin-1/2 Heisenberg quantum spin chain with SU(2)-iDMRG

$$H = \frac{1}{2} \sum_{i} \sigma_i^x \sigma_{i+1}^x + \sigma_i^y \sigma_{i+1}^y + \sigma_i^z \sigma_{i+1}^z$$





#### Fermionic systems

e.g. P. Corboz, R. Orús, B. Bauer, G. Vidal, PRB 81, 165104 (2010)

Rules (ask me later if interested in technical details or derivations)



The leading order of the computational cost is the same as in the bosonic case

#### fermionic order ~ graphical projection



physics is independent of the order physics is independent of graphical projection

*(different choices of Jordan-Wigner transformation, if mapping to a spin system)* 

#### **PEPS & Entanglement Hamiltonians**

e.g. I. Cirac et al, PRB 83, 245134 (2011), N. Schuch et al, PRL 111, 090501 (2013)





Boundary How is physics described here?



### **Emergent Hamiltonians**

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Remember it has double indices...

# **Emergent Hamiltonians**



Particles and energies from Hamiltonians, and Hamiltonians from networks of entanglement + bulk-boundary correspondence

## "branching" MERA

G. Evenbly, G. Vidal, PRL 112, 240502 (2014)





















Exact in many cases Variational ansatz for numerical simulations (e.g. DMRG)

#### MERA & AdS/CFT

e.g. B. Swingle, PRD 86, 065007 (2012), G. Evenbly, G. Vidal, JSTAT 145:891-918 (2011)

### **Emergent space-time**



Picture from M. Nozaki, S. Ryu, T. Takayanagi, JHEP10(2012)193

#### MERA entropy ~ Ryu-Takayanagi prescription







For a scale-invariant MERA, the tensors of a critical model with a CFT limit correspond to a "gravitational" description in a discretized AdS space: "lattice" realization of AdS/CFT correspondence





Finite correlation length (gapped systems) = finite number of layers



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$$\rho_{in} = tr_{out} \left( |\Psi\rangle \langle \Psi| \right) \right]$$
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Same **thermal** spectrum (entanglement Hamiltonian) finite temperature, scale invariance broken



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e.g., T. Hartman, J. Maldacena, JHEP05(2013)014

Thermofield double state

Eternal AdS black-hole

$$|TFD
angle = rac{1}{\sqrt{Z(eta)}}\sum_n e^{-eta E_n/2}|n
angle_1|n
angle_2$$



e.g., T. Hartman, J. Maldacena, JHEP05(2013)014

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#### Entanglement connects upper and lower spacetimes



M. Van Raamsdonk, arXiv:0907.2939

ER=EPR, Maldacena & Susskind





### cMERA

(continuum)

$$\left|\psi\right\rangle = Pe^{-i\int_{u^{2}}^{u^{1}} \left(K(u)+L\right)du} \left|\Omega\right\rangle$$

J. Haegeman et al, Phys. Rev. Lett. 110, 100402 (2013)

K(u) Disentangler generator

*L* Isommetry generator



### cMERA

(continuum)

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J. Haegeman et al, Phys. Rev. Lett. 110, 100402 (2013)

K(u) Disentangler generator

*L* Isommetry generator

$$g_{uu}(u)du^{2} = \mathcal{N}^{-1}\left(1 - \left|\langle \Psi(u)|e^{iL \cdot du}|\Psi(u+du)\rangle\right|^{2}\right)$$
  
Measures the density of strength of disentanglers.  
Compatible with AdS metric

M. Nozaki, S. Ryu, T. Takayanagi, JHEP10(2012)193

curvature ~ change of entanglement at every length scale

# Lots of other things:

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**Chiral PEPS** Entanglement measures Time evolution, TDVP Infinite systems **Classical systems Corner Transfer Matrices Tensor Renormalization Group** Variational updates PESS, gPEPS, TgPEPS Phases of matter 3d PEPS **Topological systems Excited states** Thermal states & dissipation Continuous tensor networks TNs & MonteCarlo, DFT, FRG... Practical implementations **Entanglement Hamiltonians Symmetries** 

Blablablablablabla...

