

Title: Entanglement loss properties and classical simulability of quantum many-body systems

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Abstract: In this talk I will expose different results concerning the properties of quantum many-body systems: on the one hand, I will introduce the concept of fine-grained entanglement loss together with its relation with majorization relations along parameter flows and Renormalization Group flows. The machinery of Conformal Field Theory will allow us to derive very general analytical properties, and some examples -like the XY quantum spin chain- will also be considered. On the other hand, I will describe results concerning the classical simulability of quantum many-body systems by means of Matrix Product States. In particular, I will present an approximated classical simulation of a quantum algorithm by adiabatic evolution solving hard instances of an NP-Complete problem up to 100 qubits.

# Entanglement loss properties and classical simulability of quantum many-body systems

**Román Orús**



*Perimeter Institute, 6 of February 2006*

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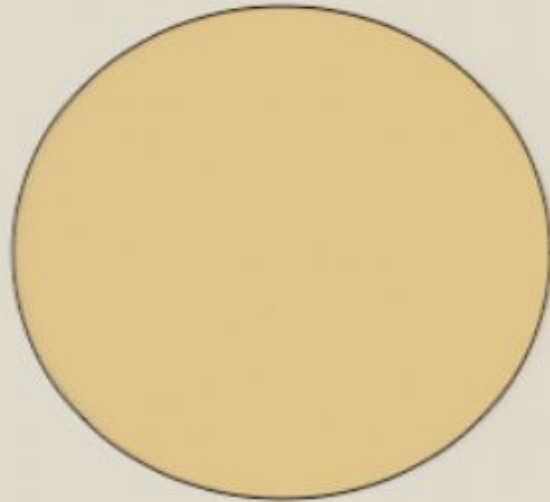
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# The working scenario...

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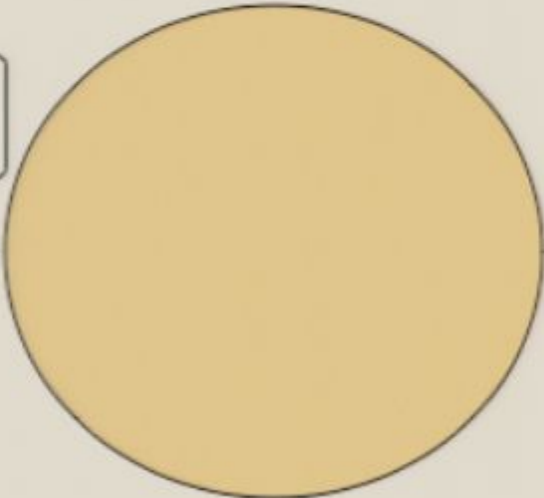
Quantum information and computation



# The working scenario...

## Quantum information and computation

Entanglement theory,  
adiabatic algorithms...

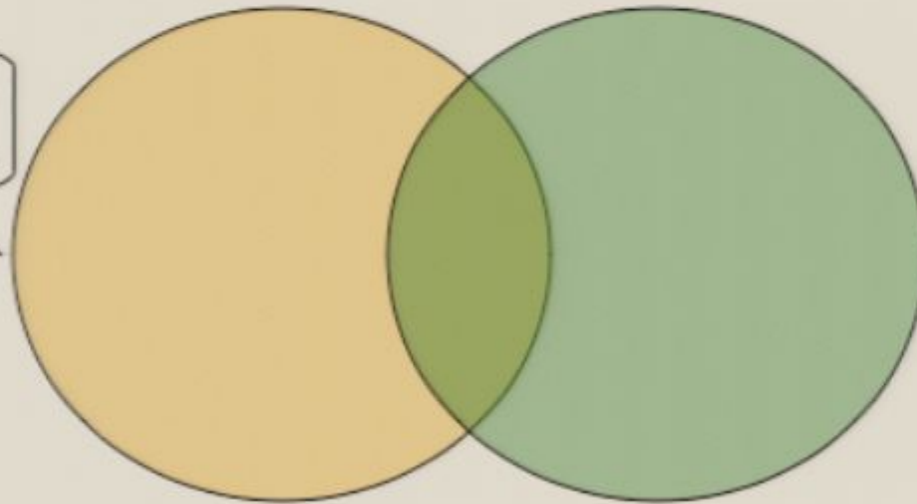


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Quantum information and computation

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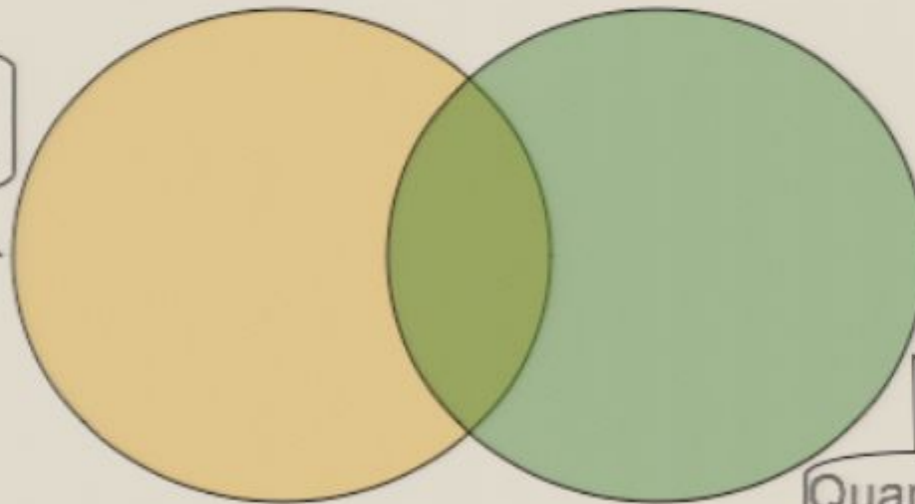


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Quantum phase transitions,  
DMRG...



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High energy physics and quantum field theory

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Condensed matter physics

Entanglement theory,  
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Quantum phase transitions,  
DMRG...

Conformal field theory,  
black-hole entropy...

High energy physics and quantum field theory

# My work at Barcelona



- Robustness of weakly-entangled states

*PRA 70, 050101 (2004)*

- Majorization in quantum algorithms

*QIP 1 (4), 283-302 (2002); EPJD 29, 119-132 (2004)*

- Classical computational complexity of quantum problems

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- Entanglement in quantum algorithms and quantum phase transitions

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PRA 70, 062318 (2004); PRA 71, 064101 (2005); quant-ph/0509023*

- Majorization and entanglement loss along parameter flows and RG-flows

*PRA 71, 052327 (2005), Erratum-ibid 73, 019904 (2006)*

- Classical simulation of quantum many-body systems

*quant-ph/0503174 (to appear in PRA),  
and more to (hopefully) come...*



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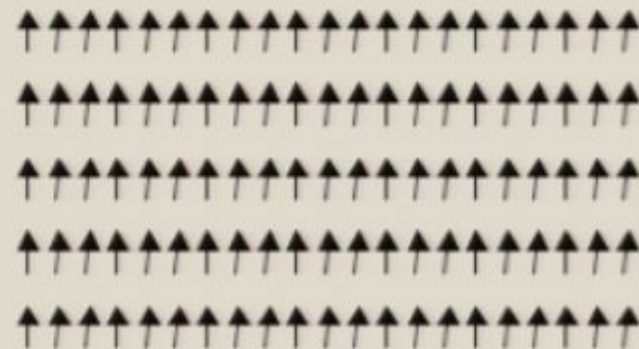
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# Majorization and entanglement loss along parameter flows and RG-flows

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# Majorization theory

$$\vec{x}, \vec{y} \in R^d \text{ probability vectors } \left( \sum_{i=1}^d x_i = \sum_{i=1}^d y_i = 1 \right)$$

- **Definition:**  $\vec{x}$  is majorized by  $\vec{y}$  ( $\vec{x} \prec \vec{y}$ ) if and only if

$$\left\{ \begin{array}{l} \bullet \vec{x} = \sum_j p_j P_j \vec{y} \quad \begin{array}{l} \text{permutation matrices} \\ \text{probabilities} \end{array} \\ \bullet \sum_{i=1}^k x_i \leq \sum_{i=1}^k y_i \quad k = 1, 2, \dots, d-1 \quad \text{(components sorted into decreasing order)} \\ \bullet \vec{x} = D \vec{y} \quad \text{doubly stochastic matrix} \end{array} \right.$$

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**Majorization imposes a very strong sense of order in probability distributions.**




# Global, monotonous and fine-grained entanglement loss

[Latorre et al., 2004]

$\rho = \rho_A = \text{tr}_B(|\Omega\rangle\langle\Omega|)$   $|\Omega\rangle$  ground state (vacuum) of a physical system that depends on a parameter  $t$

Transformation:  $\begin{cases} t_1 \rightarrow t_2 \\ \rho(t_1) \rightarrow \rho(t_2) \end{cases} \quad t_2 > t_1$



- **Global entanglement loss:**  $S(\rho(t_1)) > S(\rho(t_2))$
- **Monotonous entanglement loss:**  $\frac{dS}{dt} < 0$
- **Fine-grained entanglement loss:**  $\bar{\rho}(t) < \bar{\rho}(t + \delta t)$




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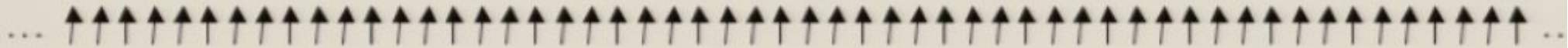
● **Fine-grained entanglement loss:**  $\vec{\rho}(t) \prec \vec{\rho}(t + \delta t)$

# Conformal symmetry in a nutshell

- Physical properties of quantum critical theories are dictated by **scale invariance**.

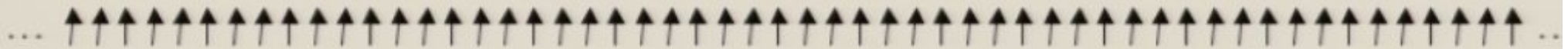
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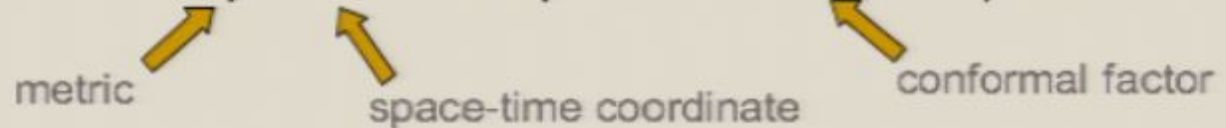
# Conformal symmetry in a nutshell

- Physical properties of quantum critical theories are dictated by **scale invariance**.



- In the continuum limit, when these theories are described in terms of a quantum field theory, one sees that **the symmetry group is even larger: conformal group**. The theory is then a **conformal field theory (CFT)**.

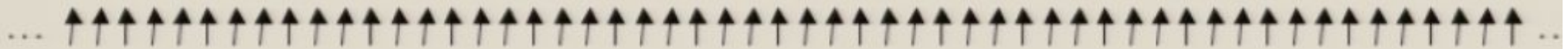
Conformal transformation:  $g_{\mu\nu}(x) \rightarrow g'_{\mu\nu}(x') = \Omega(x)g_{\mu\nu}(x)$





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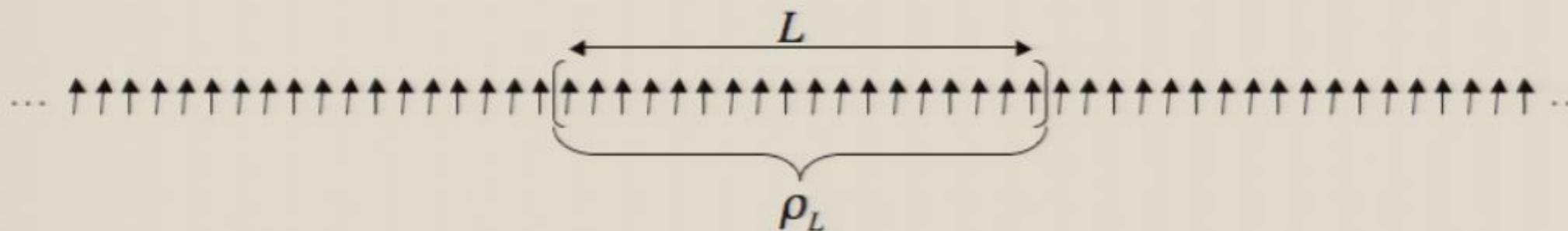
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metric  $\nearrow$   $\nwarrow$  space-time coordinate  $\nwarrow$  conformal factor

- Conformal symmetry completely determines all the physical properties of critical theories in (1+1) dimensions (e.g. some critical spin chains).**

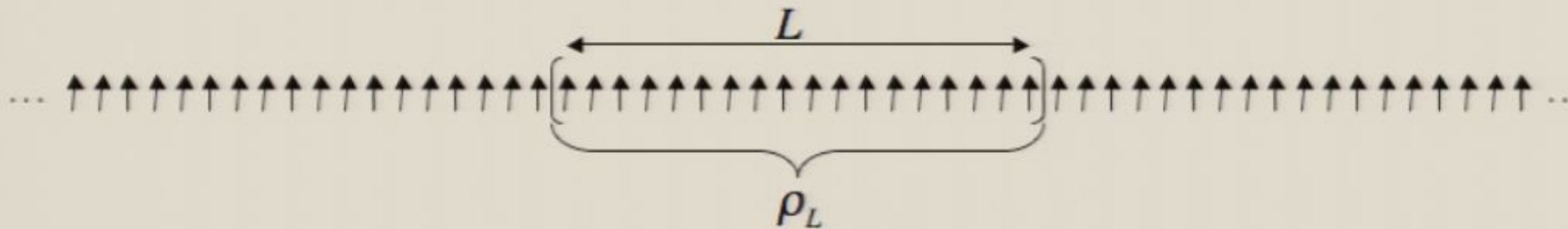
space  $\curvearrowright$   $\curvearrowleft$  time (usually fixed)

# Majorization with $L$ in CFT with (1+1) dimensions





# Majorization with $L$ in CFT with (1+1) dimensions



- **Theorem:**  $\rho_L \prec \rho_{L'}$  if  $L \geq L'$  for all possible (1+1)-dimensional CFT.

$$\rho_L = \frac{1}{Z(q)} q^{-\frac{c}{12}(L_0 + \bar{L}_0)}$$

partition function  $\curvearrowright$   $Z(q)$   $\curvearrowright$  central charge  $c$   
 $\uparrow$  0th Virasoro operators

$$\log(q) = -\frac{2\pi^2}{\log(L/\epsilon)}$$

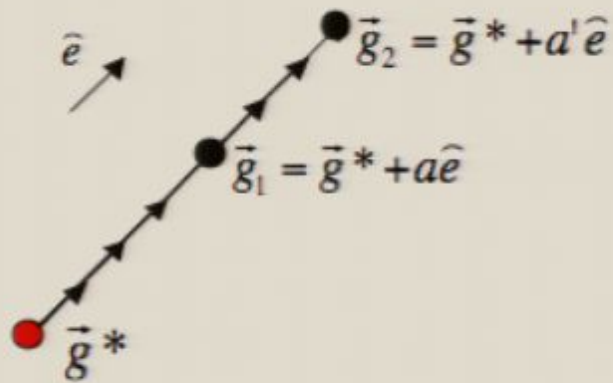
$\nearrow$  short-distance cut-off  
 (=1 for spin chains)

- **Example:** XX quantum spin chain  $H = -\sum_{l=1}^{\infty} (\sigma_l^x \sigma_{l+1}^x + \sigma_l^y \sigma_{l+1}^y)$  [Peschel, 2004]



# Majorization along parameter flows

$\vec{g} = (g_1, g_2, \dots)$  vector of parameters

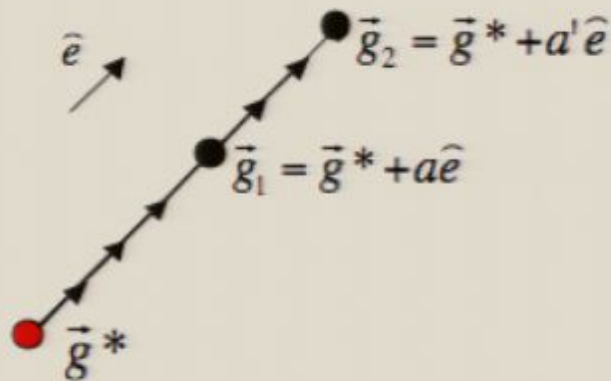


$\vec{g}^*$  conformal point

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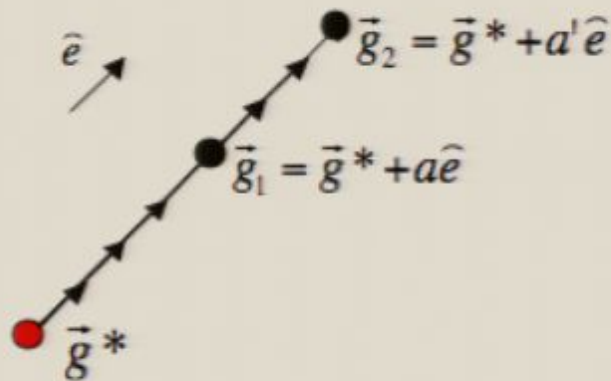
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- **Theorem:** if the reduced density matrices at  $\vec{g} = \vec{g}^* + a\hat{e}, a > 0$  preserve the same structure than at  $\vec{g}^*$ , with a monotonically decreasing  $q(\vec{g})$  along the flow in  $a$ , then  $\rho(\vec{g}_1) \prec \rho(\vec{g}_2)$ .

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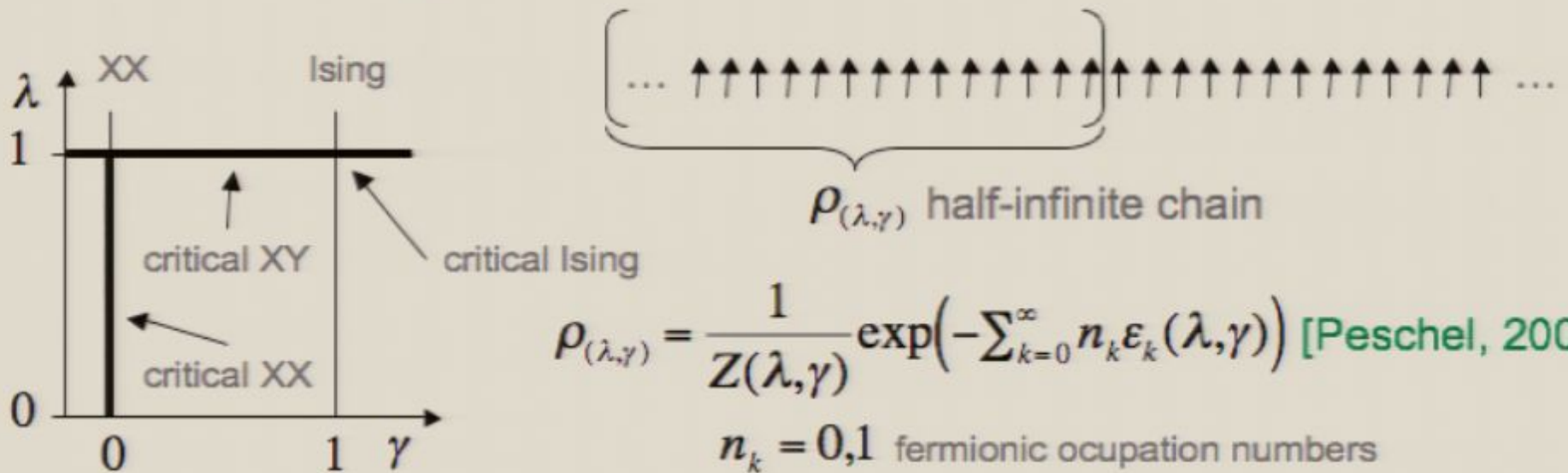
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- If  $a$  coincides with the scale of an RG-flow, then we have proven **continuous majorization along an RG-flow of some integrable theories (by construction)**.
- If the theory has only one parameter, monotonicity of the parameter along RG is trivial, but majorization of the reduced vacuum is not.

# Example: XY quantum spin chain

$$H = -\sum_{l=1}^{\infty} \left( (1+\gamma)\sigma_l^x\sigma_{l+1}^x + (1-\gamma)\sigma_l^y\sigma_{l+1}^y + 2\lambda\sigma_l^z \right)$$



$$\rho_{(\lambda, \gamma)} = \frac{1}{Z(\lambda, \gamma)} \exp\left(-\sum_{k=0}^{\infty} n_k \varepsilon_k(\lambda, \gamma)\right) \text{ [Peschel, 2004]}$$

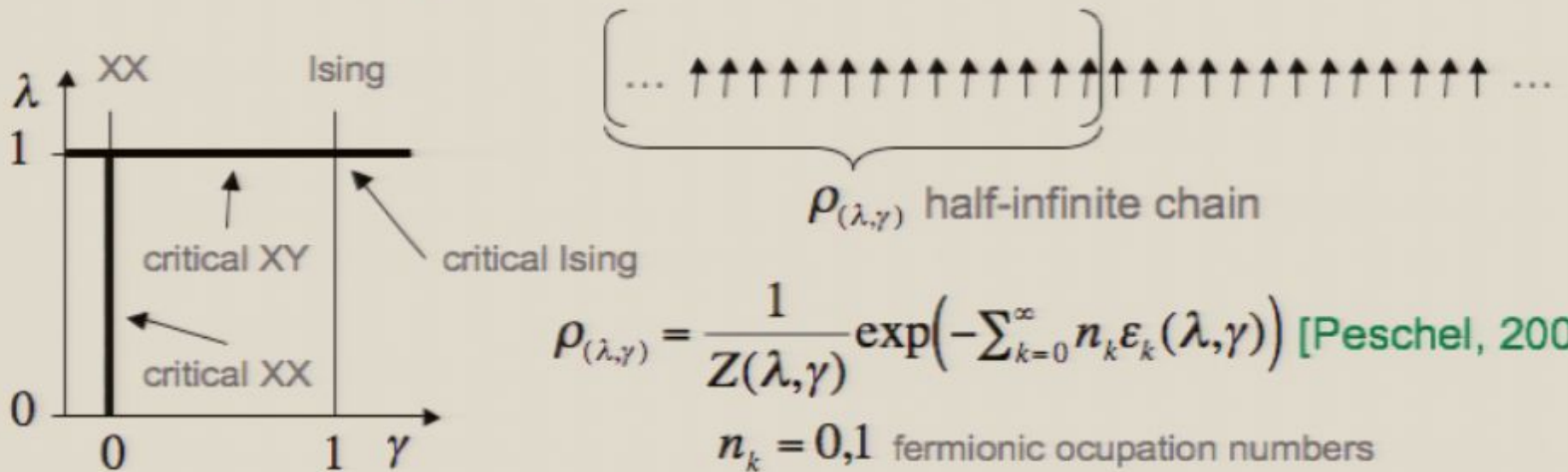
$n_k = 0, 1$  fermionic occupation numbers

$\varepsilon_k(\lambda, \gamma)$  monotonic function of the two parameters



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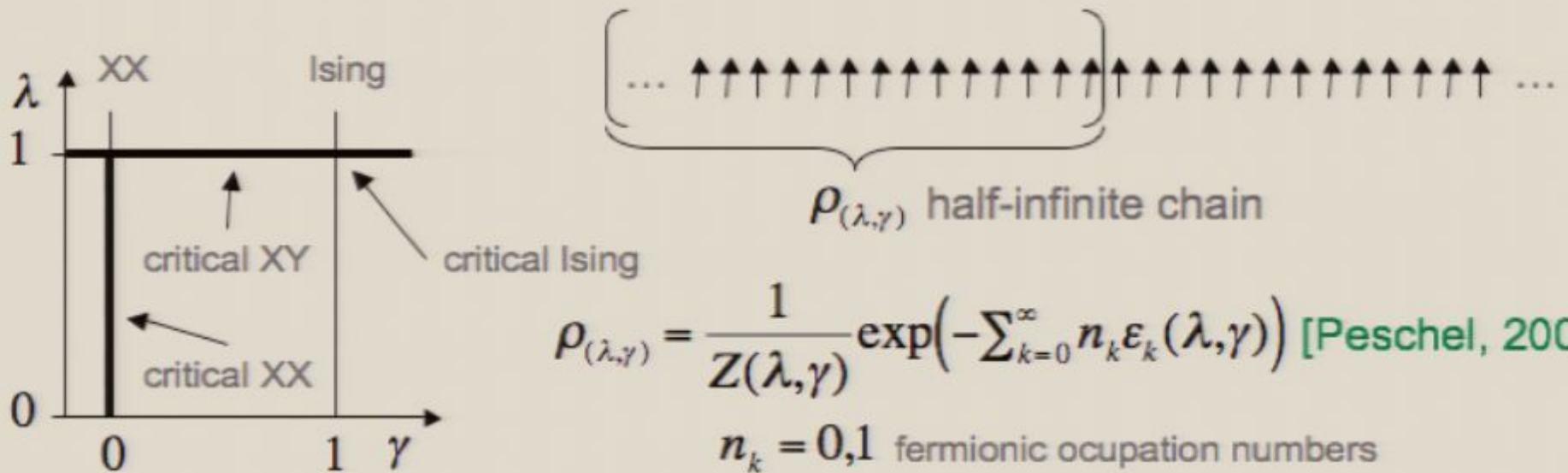
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- Majorization relations hold along the flows in  $\lambda$  and  $\gamma$ . For  $\gamma = 1$ , the flow in  $\lambda$  along the Ising line coincides with an RG-flow.
- Majorization may hint irreversibility of RG-flows in **higher dimensions** also. [Riera, Latorre, *in preparation*]

# Other interesting results with CFT

- Single-copy entanglement scales asymptotically as half the entanglement entropy for conformal field theories in (1+1) dimensions.

[Orús et al., 2005]

$$E_1(\rho_L) = \frac{c}{6} \log L - \frac{c}{6} \frac{\pi^2}{\log L} + O(1/L)$$

$$S(\rho_L) = \frac{c}{3} \log L + O(1/L)$$

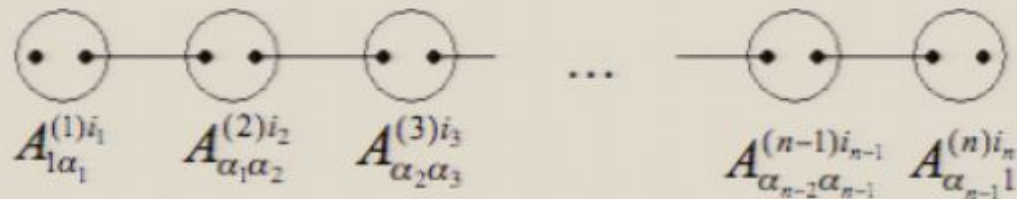
- Entanglement entropy in the Lipkin-Meshkov-Glick model behaves as if the model, though defined on a simplex, were governed by an underlying conformal symmetry in (1+1) dimensions.

[Latorre et al., 2004]

$$S(\rho_L) \approx \frac{c_{\text{eff}}}{3} \log L$$



# Classical simulation of quantum many-body systems





# A “bit” of motivation...

State of an n-qubit quantum system:  $|\psi\rangle = \sum_{\{i\}} c_{i_1 i_2 \dots i_n} |i_1, i_2, \dots, i_n\rangle$

$c_{i_1 i_2 \dots i_n}$    $2^n$  coefficients



Exact classical simulations of general Hamiltonians  
only possible up to ~24 spins (qubits) with exact methods  
(motivation to build a quantum computer [Feynmann, 1982]).

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**We have efficiently simulated *approximated* quantum computations up to 100 qubits with matrix product states.**

# Matrix product states (MPS)

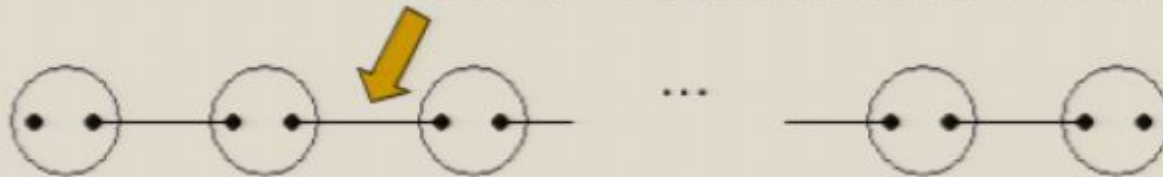
[Affleck et al., 1987]



# Matrix product states (MPS)

[Affleck et al., 1987]

maximally entangled pair of ancillas of dimension  $\chi$

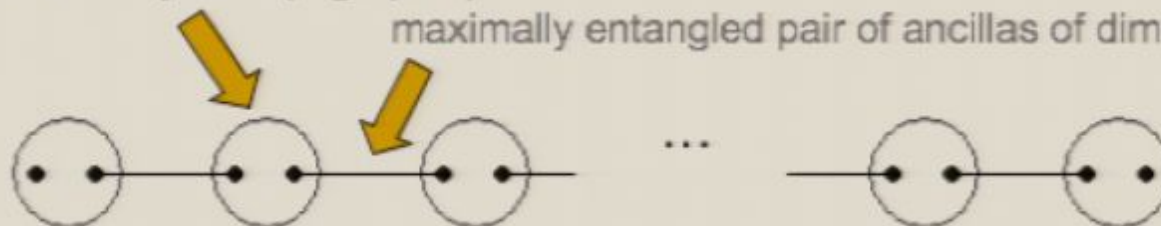


# Matrix product states (MPS)

[Affleck et al., 1987]

projection on physical local system (e.g. qubit)

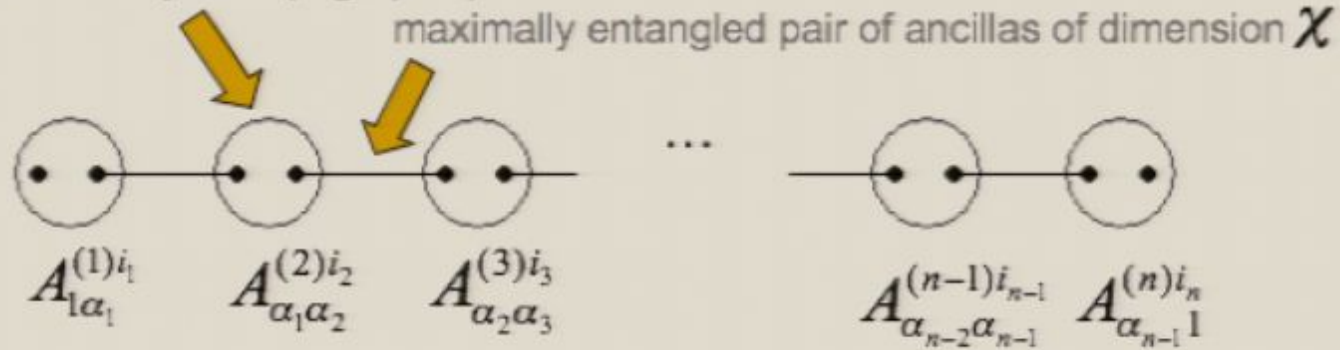
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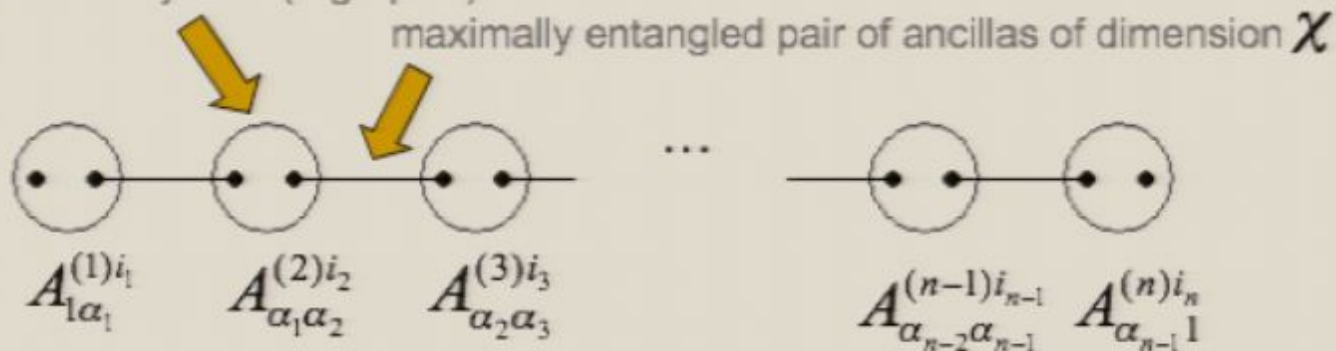
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# Matrix product states (MPS)

[Affleck et al., 1987]

projection on physical local system (e.g. qubit)



$$|\psi\rangle = \sum_{\{i\}} \sum_{\{\alpha\}} A_{1\alpha_1}^{(1)i_1} A_{\alpha_1\alpha_2}^{(2)i_2} \dots A_{\alpha_{n-1}1}^{(n)i_n} |i_1, i_2, \dots, i_n\rangle \quad i_a = 1, 2, \dots, d \quad \alpha_a = 1, \dots, \chi$$

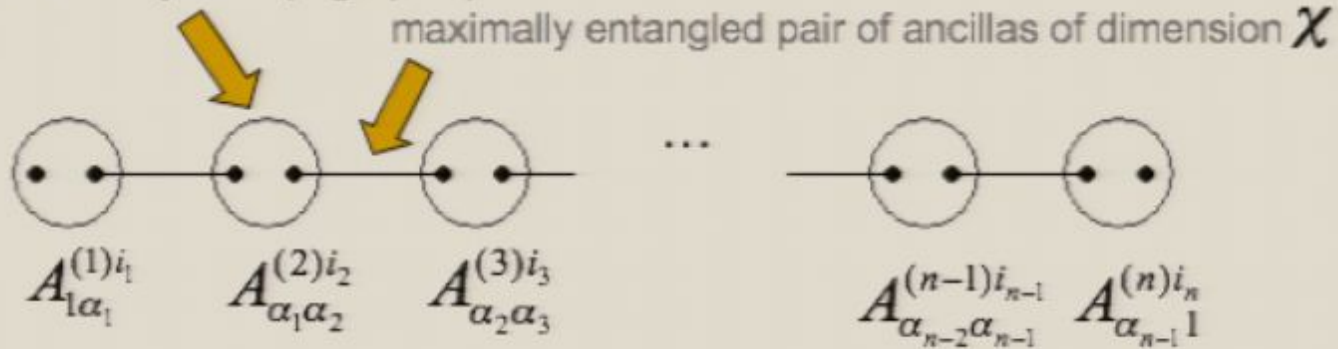
$$A_{\alpha_{a-1}\alpha_a}^{(a)i_a} = \Gamma_{\alpha_{a-1}\alpha_a}^{(a)i_a} \lambda_{\alpha_a}^{(a)}$$



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$$A_{\alpha_{a-1}\alpha_a}^{(a)i_a} = \Gamma_{\alpha_{a-1}\alpha_a}^{(a)i_a} \boxed{\lambda_{\alpha_a}^{(a)}} \longrightarrow \chi \text{ Schmidt coefficients [Vidal, 2003]}$$

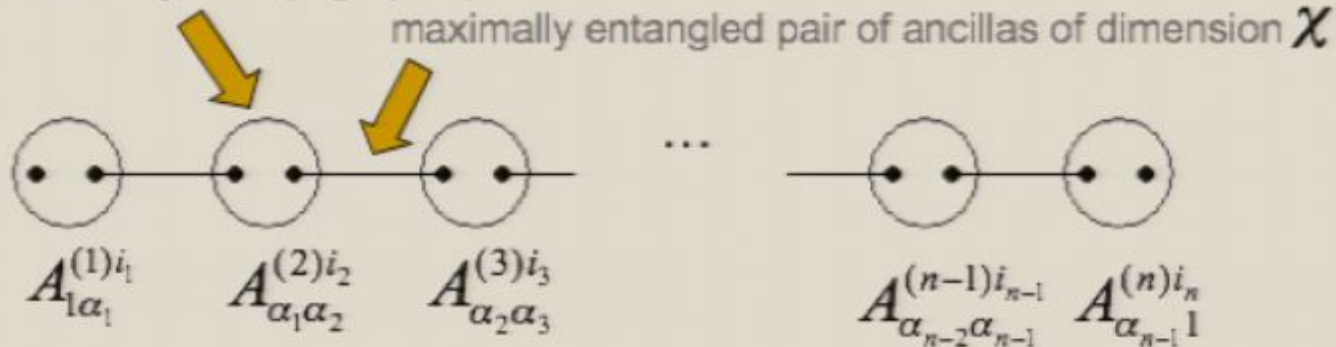
(measure of entanglement;  $\chi \approx 2^{S(\rho)}$ )



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**Generalizations** { To more than 1 dimension  $\longrightarrow$  PEPS [Verstraete, Cirac, 2004]  
With disentangling unitaries  $\longrightarrow$  MERA [Vidal, 2005]

# Why are MPS “good”?

- The state is represented with  $nd\chi^2$  parameters, instead of  $d^n$ .
- Any quantum state can be represented as an MPS, with large enough  $\chi$ .
- Physical observables (e.g. correlators) can be computed in  $O(\text{poly}(\chi))$  time.
- They allow to compute low-energy properties of Hamiltonians (DMRG, euclidean time evolution...).
- They allow to compute the dynamics of quantum states.

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**Works great for quantum many-body systems in 1 spatial dimension, because of the logarithmic scaling of the entanglement entropy.**

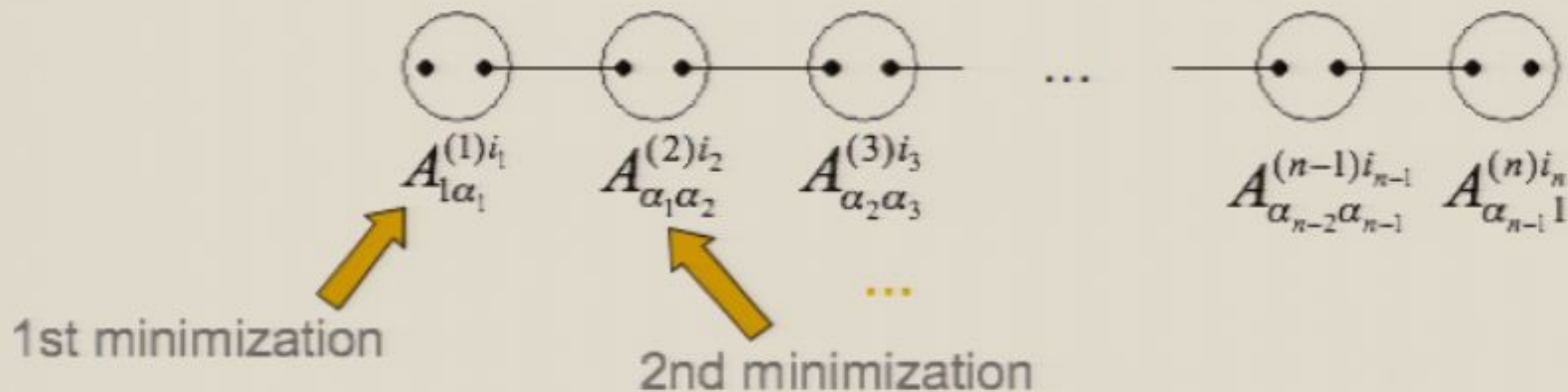
[Vidal et al., 2002]

# Computing ground states with MPS

- **Aim:** compute the ground state properties of a given Hamiltonian  $H$ .

# Computing ground states with MPS

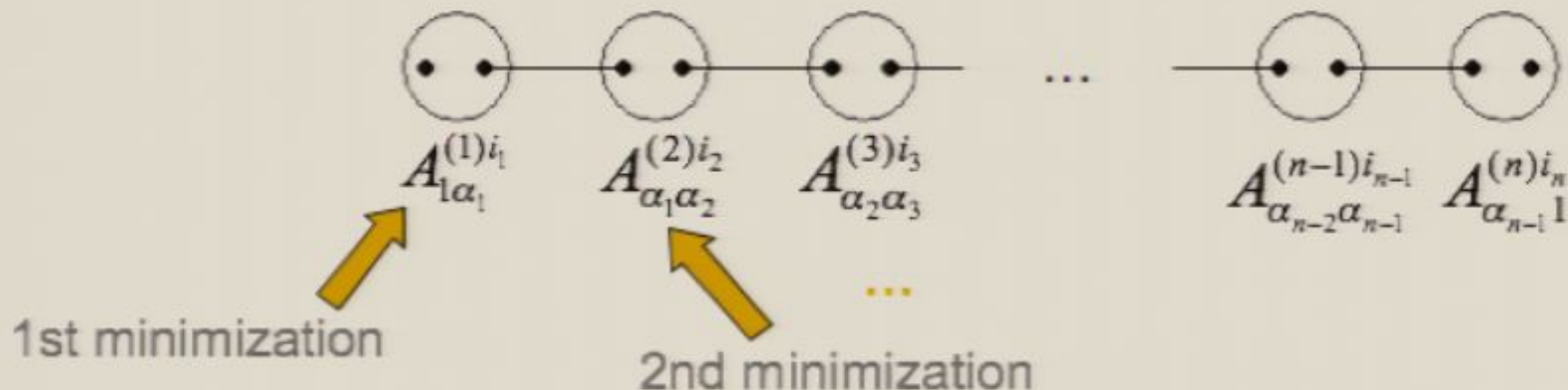
- **Aim:** compute the ground state properties of a given Hamiltonian  $H$ .
- **Basic idea:** minimize the quantity  $\langle \psi | H | \psi \rangle$  sequentially sweeping back and forth over all the matrices  $A_{\alpha_{a-1}\alpha_a}^{(a)i_a}$  for a fixed  $\chi$ .





# Computing ground states with MPS

- **Aim:** compute the ground state properties of a given Hamiltonian  $H$ .
- **Basic idea:** minimize the quantity  $\langle \psi | H | \psi \rangle$  sequentially sweeping back and forth over all the matrices  $A_{\alpha_{a-1}\alpha_a}^{(a)i_a}$  for a fixed  $\chi$ .



**This procedure is equivalent to the DMRG algorithm. [White, 1992]**



# Example: harmonic chain ( $\phi^2$ theory)

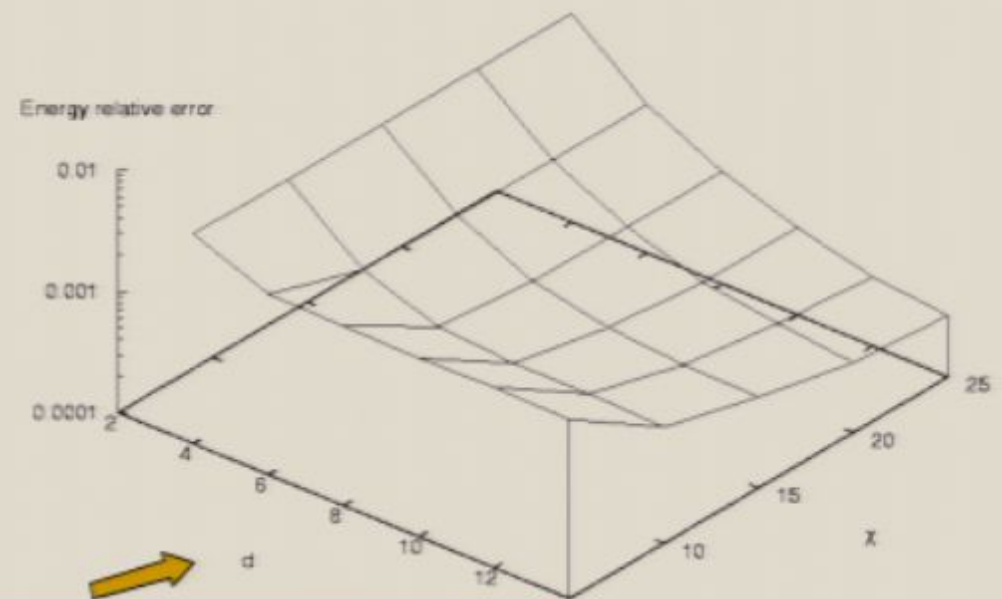
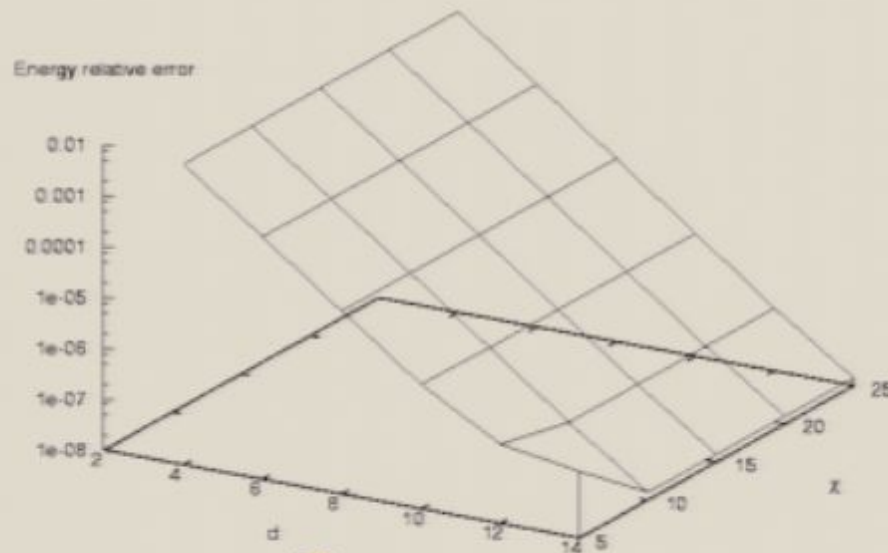
[Iblisdir, Orús, Latorre, *in preparation*]

$$H = \frac{1}{2} \sum_{i=1}^N p_i^2 + \frac{1}{2} \sum_{i=1}^{N-1} (x_i - x_{i+1})^2 + \frac{\mu}{2} \sum_{i=1}^N x_i^2 \approx \int \left( \frac{1}{2} \dot{\phi}^2 + \frac{1}{2} (\nabla \phi)^2 + \frac{\mu}{2} \phi^2 \right) dx$$

Off-critical ( $\mu = 1$ )

( $N = 30$ )

Critical ( $\mu = 0$ )



Truncation in the local dimension of the Hilbert space  
(Fock basis of free oscillators).

# Computing dynamics with MPS

[Vidal, 2003]

- Local gate on system  $a$ : 
$$A'_{\alpha\beta}^{(a)i'_a} = U_{i_a i'_a}^{(a)} A_{\alpha\beta}^{(a)i_a}$$

# Computing dynamics with MPS

[Vidal, 2003]

Local gate on system  $a$ : 
$$A'_{\alpha\beta}^{(a)i'_a} = U_{i_a i'_a}^{(a)} A_{\alpha\beta}^{(a)i_a}$$

Non-local gate on adjacent systems  $a$  and  $a+1$ :

diagonalization of 
$$\rho_{\alpha\gamma}^{ij} = |\lambda_{\beta}^{(a-1)}|^2 \Theta_{\beta\alpha}^{ki} \Theta_{\beta\gamma}^{*kj} \quad (\text{rank } d\chi)$$

$$\Theta_{\alpha\gamma}^{i'_a i'_{a+1}} = U_{i'_a i'_{a+1}, i_a i_{a+1}}^{(a,a+1)} A_{\alpha\beta}^{(a)i_a} A_{\beta\gamma}^{(a+1)i_{a+1}} = A'_{\alpha\beta}^{(a)i'_a} A'_{\beta\gamma}^{(a+1)i'_{a+1}}$$

$1, \dots, \chi$

$1, \dots, d\chi$

creation of entanglement

(exponential growth!  
Some truncation  
is needed...)

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$1, \dots, \chi$

$1, \dots, d\chi$

creation of entanglement

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Some truncation  
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- Non-local gate on non-adjacent systems: SWAP systems until they are together.

# Truncation schemes

- **Local procedure** [Vidal, 2003]: retain only a fixed number  $\chi$  of terms that carry most of the entanglement in the decomposition.



**Entanglement-inspired** truncation method



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- **Non-local procedure** [Verstraete, Cirac, 2004]:

compute *all* the matrices  $\tilde{A}_{\alpha\beta}^{(a)i_a}$  with indices up to a fixed  $\chi$  such that

$$\|\tilde{|\psi\rangle} - |\psi'\rangle\|^2 \text{ is minimum.}$$



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**Precision-inspired** truncation method

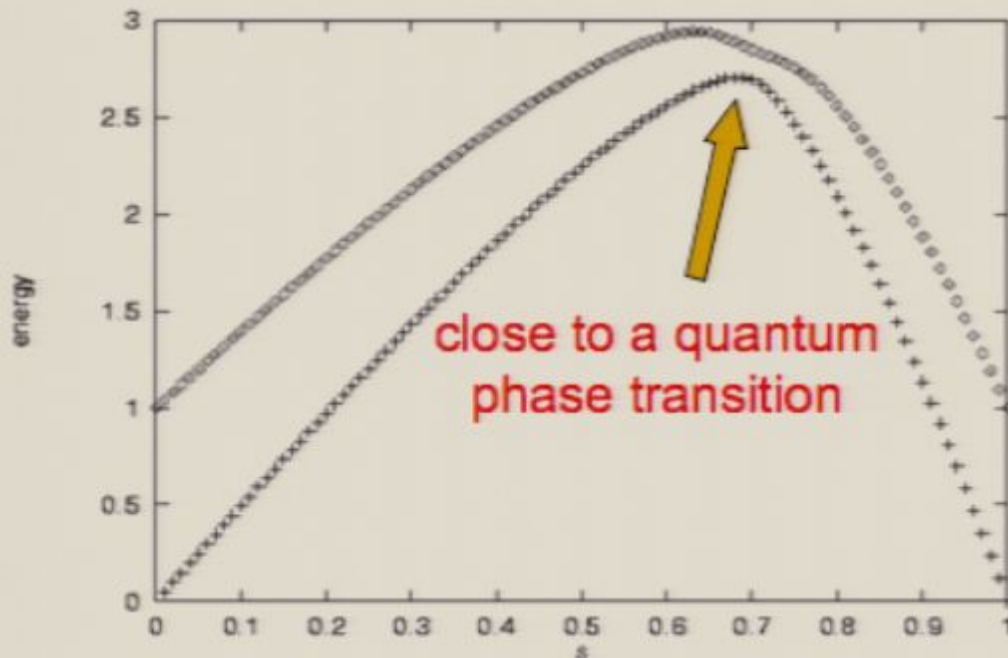
**We applied the first local procedure in our time-evolution simulations.**

# Adiabatic quantum algorithms

[Farhi et al., 2000]

$H_p$  codifies the solution to a problem in its ground state

$H_0$  with an easy-to-prepare ground state



At  $t = 0$  prepare the ground state of  $H_0$  and evolve adiabatically with the time-dependent Hamiltonian

$$H(t/T) = (1-t/T)H_0 + (t/T)H_p$$

Adiabatic theorem

$$T > \frac{\Delta}{g_{\min}^2}$$

$$\Delta = \max \left| \frac{dH_{10}(t/T)}{dt} \right| \quad s = t/T$$

$$g_{\min} = \min(E_1(t) - E_0(t))$$

$$|\langle \psi(T) | E_0(T) \rangle|^2 \approx 1$$

# The Exact Cover NP-Complete problem

**Instance:** set of  $m$  clauses (constraints) over  $n$  bits

Clause for bits  $i, j, k$ :

	$i, j, k$	
	100	} Satisfies the clause
	010	
	001	

**Problem:** is there an assignment of the  $n$  bits that satisfies *all* the clauses?

Hard classical instances  $\longrightarrow \frac{m}{n} \approx 0.8$

# Adiabatic algorithm solving Exact Cover

$$H_0 = \sum_{i=1}^n \frac{d_i}{2} (1 - \sigma_i^x) \quad H_p = \sum_{c(i,j,k)}^m (z_i + z_j + z_k - 1)^2$$

$d_i = \# \text{ clauses for } i$        $z_i = \frac{1}{2}(1 - \sigma_i^z)$

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Highly non-local  
interacting spin Hamiltonian



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Average numerical scalings for ~20 qubits

$$\left\{ \begin{aligned}
 T &\approx n^2 \quad [\text{Farhi et al., 2001}] \\
 \chi_{\max} &\approx 2^{0.1n} \quad [\text{Orús, Latorre, 2004}]
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 \end{aligned} \right.$$



Small! (but exponential).  
Possibility of a good MPS simulation  
for "many" qubits?

# Discretization in gates

Time evolution  $\longrightarrow U_{T,0} = T\left(e^{-i\int_0^T H(t/T)dt}\right) \approx U_{T,T-\Delta} \dots U_{2\Delta,\Delta} U_{\Delta,0}$

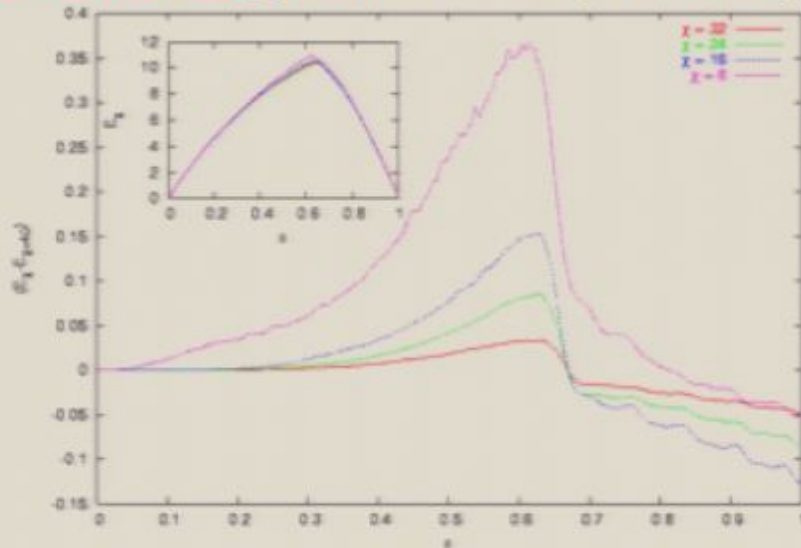
$$U_{(l+1)\Delta,l\Delta} = e^{-i\Delta H(s)} \approx \left( e^{-i\frac{\delta}{2}(1-s)H_0} e^{-i\delta s H_p} e^{-i\frac{\delta}{2}(1-s)H_0} \right)^{\frac{\Delta}{\delta}} \quad s = \frac{l\Delta}{T}$$

2nd order Trotter expansion

$$\left\{ \begin{aligned} e^{-i\frac{\delta}{2}(1-s)H_0} &= \prod_{i=1}^n e^{-i\frac{\delta}{4}(1-s)d_i(1-\sigma_i^x)} \\ e^{-i\delta s H_p} &= \prod_{c(i,j,k)} e^{-i\delta s(z_i^2 - 2z_i)} e^{-i\delta s(z_j^2 - 2z_j)} e^{-i\delta s(z_k^2 - 2z_k)} e^{-i\delta s} \\ &\quad e^{-i2\delta s z_i z_j} e^{-i2\delta s z_i z_k} e^{-i2\delta s z_j z_k} \end{aligned} \right.$$

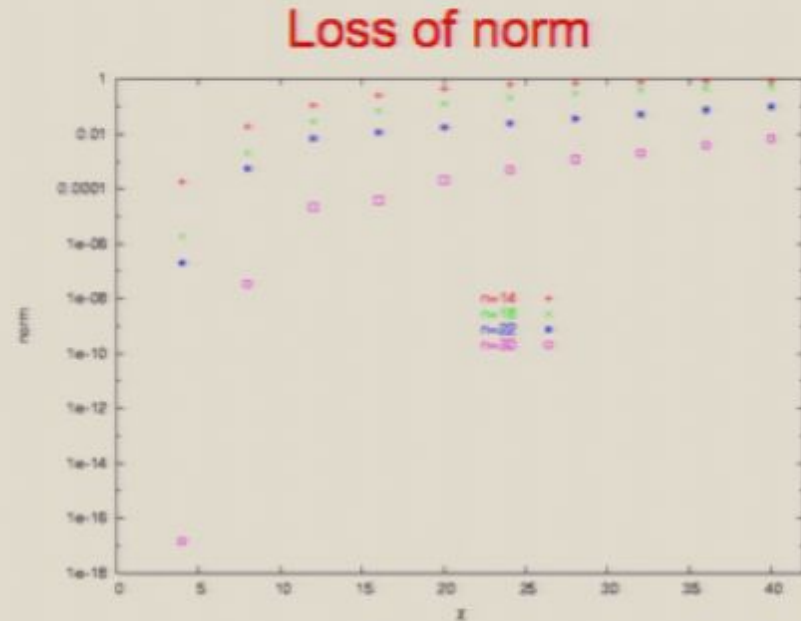
# Simulation results (1)

Expected energy for  $n = 30, m = 24, T = 100$



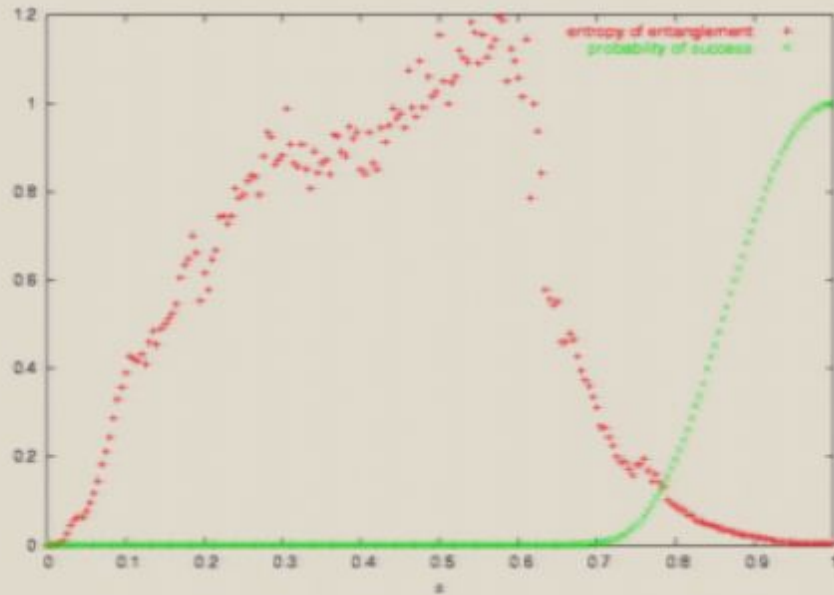
Good accuracy in  
"small" computation time (as  
compared to the exact calculation)

Increasing precision  
with increasing  $\chi$



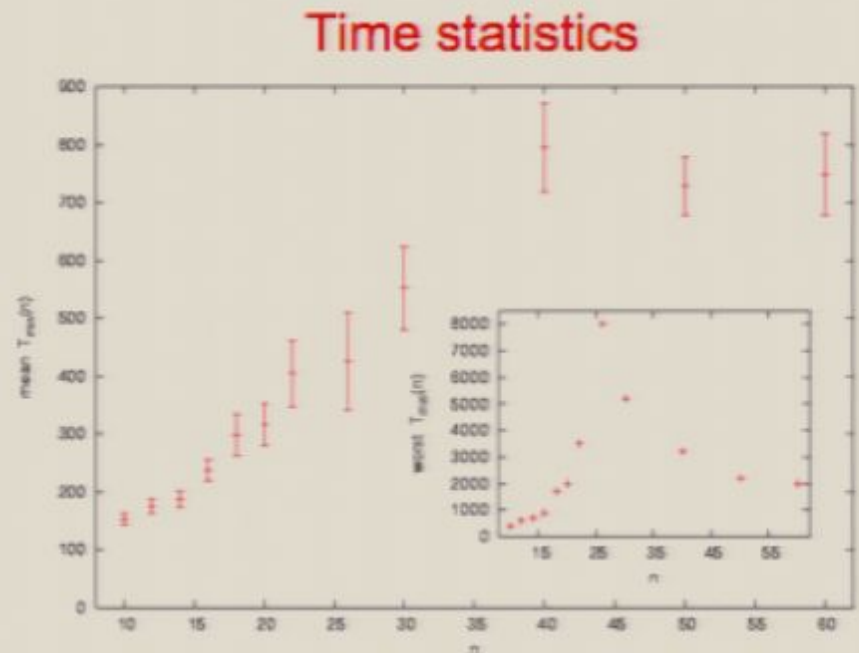
# Simulation results (2)

100-qubit instance with  $n = 100, m = 84, T = 2000, \chi = 14 \ll 2^{50}$



← Finds the correct solution among  $2^{100} \approx 10^{30}$  possibilities!

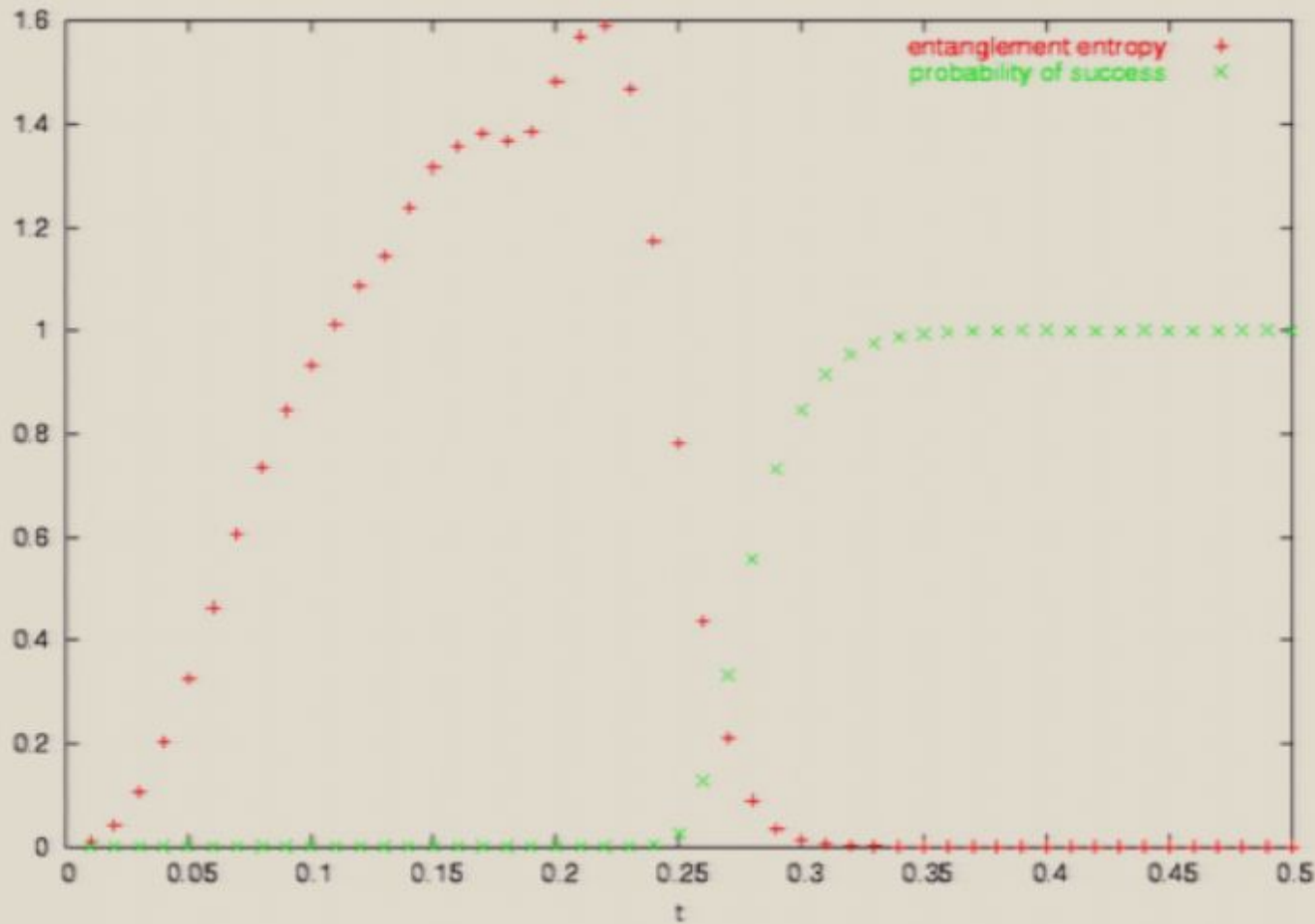
Average appears to grow very slowly with  $n$  →





# Euclidean time evolution: $e^{-H_p t}$

$$n = 26, \chi = 6$$



Unphysical  
(not unitary)

Finds the correct  
solution much  
faster than the  
adiabatic simulations



# Conclusions



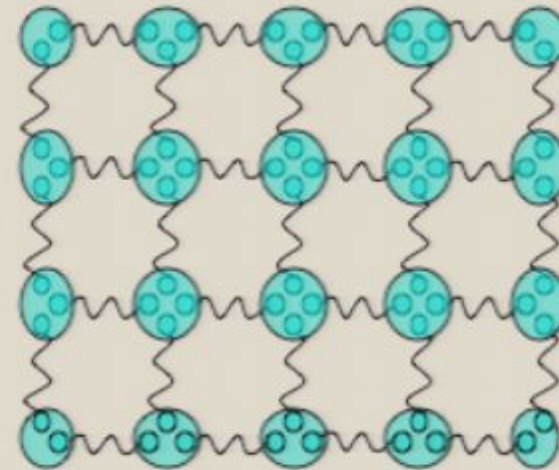
- **Majorization along  $L$  and parameter/RG-flows** can be proven for some integrable theories in  $(1+1)$  dimensions under appropriate assumptions using CFT. This proves some conjectures previously raised for quantum spin chains, and hints a way to prove irreversibility of RG-flows in more dimensions.
- **MPS are a useful computational tool** for calculating ground states and dynamical properties of many interesting physical systems. On top, MPS-simulations of quantum computations can handle a relatively large number of qubits with reasonable approximation schemes.

# Future work



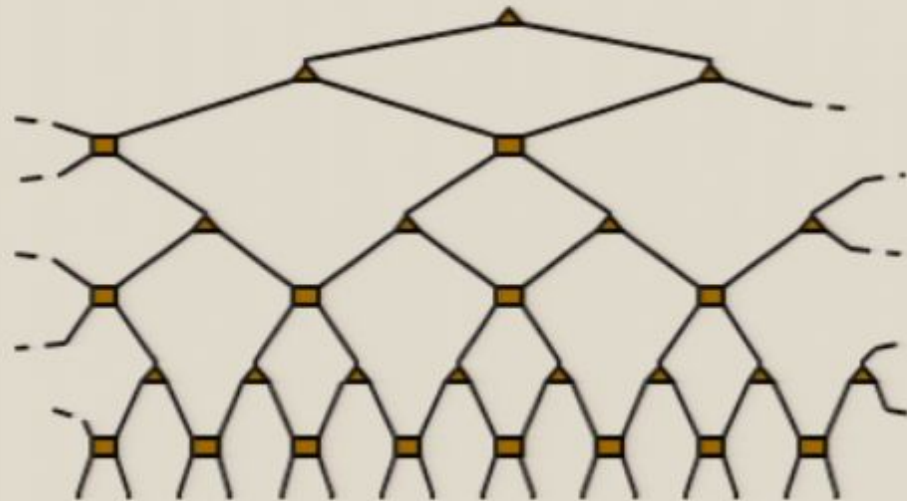
- **PEPS** in more than 1 dimension.

[Verstraete, Cirac, 2004]



- **MERA** (disentangling unitaries) in 1 and more dimensions.

[Vidal, 2005]

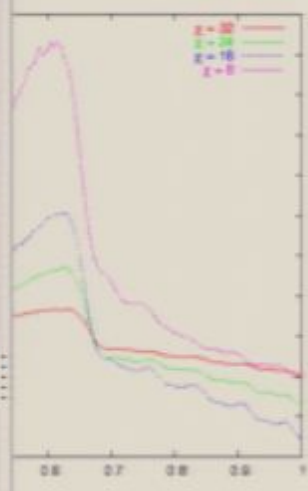




RG\_MPS.ppt

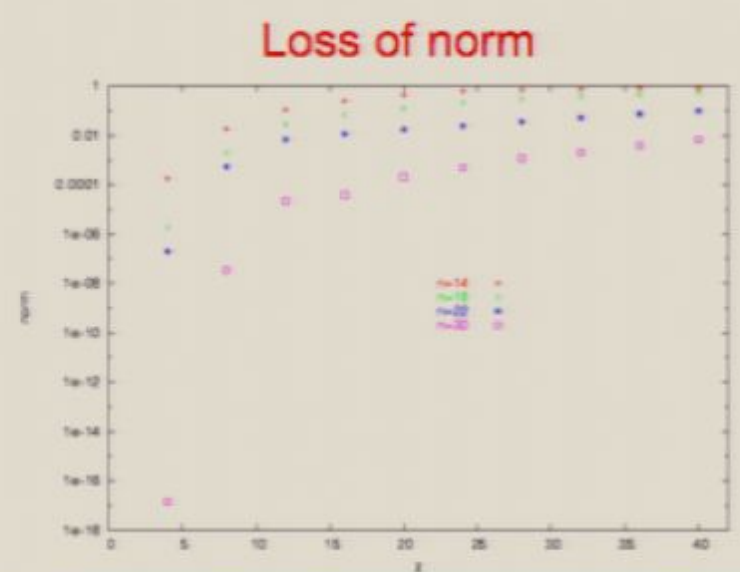
# Simulation results (1)

$n = 30, m = 24, T = 100$



← Good accuracy in "small" computation time (as compared to the exact calculation)

Loss of norm



precision sing  $\chi$  →

Diapositiva en miniatura

Simulation results (1)



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Gráficos

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Fuente

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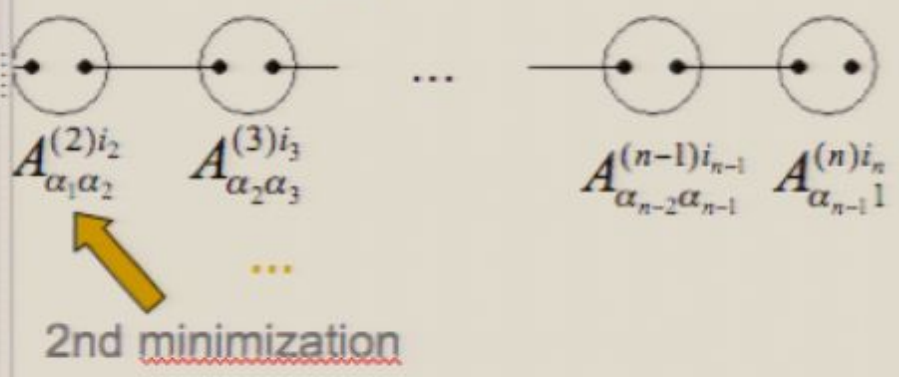
Tamaño: 18

N K S

# Computing ground states with MPS

Computing ground state properties of a given Hamiltonian  $H$ .

Minimize the quantity  $\langle \psi | H | \psi \rangle$  sequentially sweeping back and forth through the matrices  $A^{(a)}_{\alpha_{a-1}\alpha_a}$  for a fixed  $\chi$ .



equivalent to the DMRG algorithm. [White, 1992]

Diapositiva en miniatura

Computing ground states with MPS

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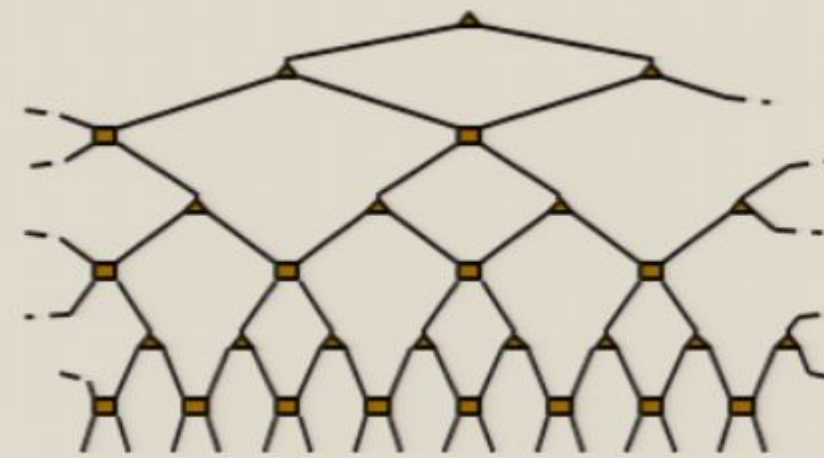
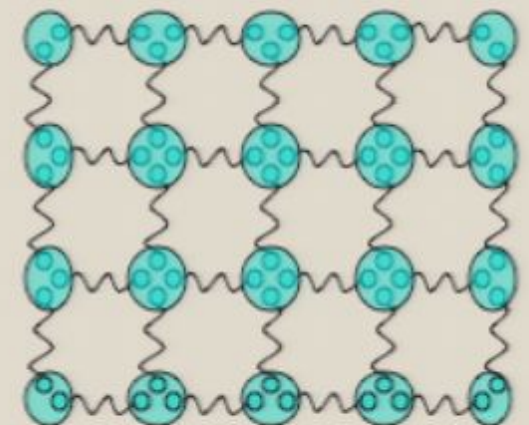

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RG\_MPS.ppt

# Future work



dimension.

[2004]

unitaries)

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Diapositiva en miniatura

Future work



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Gráficos

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Fuente

Nombre: Garan

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