DMRG: Basics

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Exponential Wall H_{H}

- Size of the Hilbert space grows exponentially with system size $\sim d^N$
- Size of the Hilbert space occupied the ground state grows much slower $\sim dN$

Density Matrix

Probability p_i in the pure state $|\psi_i\rangle$

$$
\rho = \sum_{i} p_{i} | \psi_{i} \rangle \langle \psi_{i} |
$$
\n
$$
\begin{array}{c}\n\text{for } \langle \psi_{i} \rangle \langle \psi_{i} | \\
\text{for } \rho = 1 \\
\text{for } \langle \psi | \rho | \psi \rangle \geq 0 \\
\text{for } \langle \psi | \rho | \psi \rangle \geq 0\n\end{array}
$$

Observable

$$
\langle \mathcal{O} \rangle = \text{tr}(\rho \mathcal{O}) = \sum_{i} p_i \langle \psi_i | \mathcal{O} | \psi_i \rangle
$$

Reduced Density Matrix educed Density Matrix

$$
\rho_{A} = \text{tf}_{B}(\rho_{AB})
$$
\n
$$
|\psi\rangle = \frac{1}{\sqrt{2}}(|\uparrow \downarrow \rangle + |\downarrow \uparrow \rangle)
$$
\n
$$
|\psi\rangle = \frac{\sqrt{2}}{\sqrt{2}}(|\uparrow \downarrow \rangle + |\downarrow \uparrow \rangle)
$$
\n
$$
\rho_{A} = \text{tr}_{B}(|\psi\rangle\langle\psi|) = \frac{1}{2} \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}
$$
\nBest description of region A

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$$
\nBest description of region A

Schmidt Decomposition

• If $|\psi\rangle$ is a pure state, it can always be decomposed into

$$
|\psi\rangle = \sum_{i}^{A} \lambda_{i} \begin{vmatrix} i \\ i \end{vmatrix} \begin{vmatrix} \lambda_{B} \\ i \end{vmatrix}
$$

where
 $\lambda_{i} \ge 0$ and
 $\{ |i_{A} \rangle \}, \{ |i_{B} \rangle \}$ are orthonormal basis of A, B

$$
\rho_A = \text{tr}_B(|\psi\rangle\langle\psi|) = \sum_i^{N_\lambda} \lambda_i^2 |i_A\rangle\langle i_A|
$$

B

Subsystem states Subsyste

• What are the most important subsystem states ?

Hamiltonian
\n
$$
H = H_S + H_E + H_{SE}
$$
\nWavefunction

\n
$$
|\psi\rangle = \sum_{i,\alpha} \psi_{i,\alpha} |i\rangle_S |\alpha\rangle_E
$$

Best approximation with m subsystem states . Best approximation with *m* subsystem states:

$$
|\tilde{\psi}\rangle = \sum_{n=1}^{m} \sum_{\alpha} \tilde{\psi}_{n,\alpha} |\phi_n\rangle_S |\alpha\rangle_E
$$

Minimize the distance between states:

$$
S = \left| |\tilde{\psi} \rangle - |\psi \rangle \right|^2
$$

Eigenstates of reduced DM

Controlled Approximation

$$
|\psi\rangle \approx |\psi_{AB}^m\rangle \equiv \sum_{i}^{m} \lambda_i |i_A\rangle |i_B\rangle, \quad m < N_\lambda
$$
\n
$$
\epsilon = 1 - \sum_{i=m+1}^{N_\lambda} \lambda_i^2
$$

• The accutacy (of the)
approximation depends on how the state of A approximation depends on how fast λ_i decays.

Approximate Wavefunctions Approximate wavefunctions N_{Sch}

m-dimensional MPS m-dimensional MPS

1D ground state

General, incl. 2D General, including 2D

Entanglement Entropy

Von Neumann Entanglement Entropy

$$
S(A) = -\operatorname{tr}\left[\rho_A \ln\left(\rho_A\right)\right] = -\sum_i p_i \ln p_i = S(B)
$$

Reduced density matrix

- Measures how entangled subregions A and B are.
- The number of states to keep, m , $\begin{array}{cc} \sqrt{A+1} & \sqrt{A+1} \\ \hline \end{array}$ scales with S :

$$
\boxed{m \sim e^{S(A)}}
$$

Scaling of entanglement entropy

RG transformation

- Diagonalization of the reduced density matrix gives you the RG transformation.
- Truncation is done by truncating the transformation matrix.

$$
\boxed{\rho_A^{dia} = U \rho_A U^{-1}}
$$

$$
U = \begin{pmatrix} 1_{A} & 1_{A} & 1_{A} \\ u_{11} & u_{12} & \cdots & u_{1N_{\lambda}} \\ u_{21} & u_{22} & \cdots & u_{2N_{\lambda}} \\ \vdots & \vdots & \ddots & \vdots \\ u_{N_{\lambda}1} & u_{N_{\lambda}2} & \cdots & u_{N_{\lambda}N_{\lambda}} \end{pmatrix}
$$

RG transformation

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$$
U \rightarrow U_m = \begin{pmatrix} |1_A\rangle & |2_A\rangle & & |m_A\rangle \\ u_{11} & u_{12} & \cdots & u_{1m} \\ u_{21} & u_{22} & \cdots & u_{2m} \\ \vdots & \vdots & \ddots & \vdots \\ u_{N_A 1} & u_{N_A 2} & \cdots & u_{N_A m} \end{pmatrix} \qquad | \psi \rangle \rightarrow | \psi_m \rangle
$$

$$
S_{i} S_{i+1}, \t S = 1/2
$$

\nHeisenberg model
\n
$$
H = \sum_{i} S_{i} S_{i+1} = \sum_{i} S_{i}^{z} S_{i+1}^{z} + \frac{1}{2} \left(S_{i}^{+} S_{i+1}^{-} + S_{i}^{-} S_{i+1}^{+} \right)
$$
\n
$$
S^{z} = \begin{pmatrix} 1/2 & 0 \\ 0 & -1/2 \end{pmatrix} S^{+} = \begin{pmatrix} 0 & 1 \\ 0 & 0 \end{pmatrix} S^{-} = \begin{pmatrix} 0 & 0 \\ 1 & 0 \end{pmatrix}
$$
\n
$$
S(l) = \frac{1}{6} \ln \left[\frac{2L}{\pi} \sin \left(\frac{\pi l}{L} \right) \right] + \frac{1}{2} c' + \ln g
$$
\n
$$
m \sim e^{S(L/2)} \approx L^{1/6}
$$

i

Split chain into blocks

$$
H = H_{e_{i-1}} + \mathbf{S}_{i-1} \cdot \mathbf{S}_i + \mathbf{S}_i \cdot \mathbf{S}_{i+1} + \mathbf{S}_i \cdot \mathbf{S}_{i+2} + H_{b_{i+2}}
$$

Block Hamiltonian Block Hamiltonian

$$
|\psi\rangle = \sum_{\substack{e_{i-1}, \sigma_i, \\ \sigma_{i+1}, \sigma_{i+2}}} c_{e_{i-1}, \sigma_i, \sigma_{i+1}, b_{i+2}} \sqrt{\frac{e_{i-1}}{e_{i-1}} \otimes \sigma_i} \otimes \sqrt{\sigma_{i+1}} \otimes \sqrt{\sigma_{i+2}}
$$

$$
\bullet\hspace{-.7em}-\hspace{-.7em}\bullet\hspace
$$

$$
H_b^{(s)} = \frac{1}{2} \left(S_{s,1}^+ \otimes S_{s,2}^- + S_{s,1}^- \otimes S_{s,2}^+ \right) + S_{s,1}^z \otimes S_{s,2}^z
$$

Single spin operator in the block *I* ⊗ **S**

$$
H_b^{(e)} = \frac{1}{2} \left(S_{s,3}^+ \otimes S_{s,4}^- + S_{s,3}^- \otimes S_{s,4}^+ \right) + S_{s,3}^z \otimes S_{s,4}^z
$$

Single spin operator in the block **S** ⊗ *I*

 $H_{se} =$ 1 $\frac{1}{2}$ $\left(S_{s,2}^+ \otimes S_{s,3}^- + S_{s,2}^- \otimes S_{s,3}^+\right) + S_{s,2}^z \otimes S_{s,3}^z$ $S^z = \left($ $1/2$ 0 $\begin{pmatrix} 0 & 0 \\ 0 & -1/2 \end{pmatrix}$ $S^+ = \begin{pmatrix} 0 & 0 \\ 0 & 0 \end{pmatrix}$ 0 1 $\begin{pmatrix} 0 & 1 \\ 0 & 0 \end{pmatrix} \quad S^- = \begin{pmatrix} 0 & 0 \\ 0 & 0 \end{pmatrix}$ 0 0 $1 \quad 0)$

 $H = H_b^{(s)} \otimes I + I \otimes H_b^{(e)} +$ $(S_x \otimes S_x + S_y \otimes S_y + S_z \otimes S_z)$

Find the ground state

- Find the ground state $|\psi_0\rangle$ of H
- Construct the density matrix $\rho = |\psi_0\rangle \langle \psi_0|$
- Construct the reduced density matrix $\rho_s = \sum_{l} \langle k \, | \, \langle l \, | \, \psi_0 \rangle \langle \psi_0 \, | \, k \rangle \, | \, l \rangle$ *kl*
- Keeping m eigenstates $\{ | \phi_i \rangle \}$ with largest eigenvalues $\{ \Lambda_i \}$ of ρ_s

RG transformation

• Construct transformation matrix

$$
U_m = (| \phi_1 \rangle | \phi_2 \rangle \dots | \phi_m \rangle)
$$

• Transform the block Hamiltonian and operators

$$
\tilde{H}_{b}^{(s)} = U_{m}^{\dagger} H_{b}^{(s)} U_{m}, \tilde{\mathbf{S}} = U_{m}^{\dagger} (I \otimes \mathbf{S}) U_{m}
$$
\n
$$
U_{m} = \phi_{1} \phi_{2} \dots \phi_{m} \qquad \qquad \boxed{\tilde{H}} = U_{m}^{\dagger} \qquad H \qquad U_{m}
$$

Building the Hamiltonian Building the Hamiltonian Building the Hamiltonian

$$
H_{b,3}^{(s)} = \tilde{H}_{b,2}^{(s)} + (\tilde{S} \otimes I) \cdot (I \otimes S)
$$

$$
H_{b,3}^{(e)} = \tilde{H}_{b,2}^{(e)} + (I \otimes S) \cdot (\tilde{S} \otimes I)
$$

Building the Hamiltonian Building the Hamiltonian Building the Hamiltonian

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$$

Infinite-size DMRG

Finite-size DMRG

- Grow the chain to the desired size
- Improve ground state (energy) by sweeping

Sweeping

Finite-size DMRG

Sweeping

Measurements

$$
\left\langle \psi \left| S_i^z S_j^z \right| \psi \right\rangle \approx \left\langle \psi_{L/2}^m \left| \tilde{S}_i^z \tilde{S}_j^z \right| \psi_{L/2}^m \right\rangle
$$

$$
S_i^z = O(i, L/2) S_i^z O^t(i, L/2),
$$

$$
O(i, L/2) \equiv U_{trunc}(i) U_{trunc}(i+1) \cdots U_{trunc}(L/2)
$$

$$
O(i, L/2) = U_m(i) U_m(i+1) \cdots U_m(L/2)
$$

Fermionic sign

$$
S_j^z = c_j^{\dagger} c_j - \frac{1}{2}
$$

\n
$$
S_j^+ = c_j^{\dagger} e^{i\pi \sum_{l < j} n_l}
$$

\n
$$
S_j^- = c_j e^{-i\pi \sum_{l < j} n_l}
$$

$$
\underline{1} \qquad \qquad \cdot s_{j-1} \tilde{c}_j, \qquad s_i = e^{i \pi n_i}
$$

$$
c_i^{\dagger} c_j = S_i^+ e^{-i\pi \sum_{l=i+1}^{j-1} n_l} S_j
$$

Jordan-Wigner transformation

Optimization

- Use symmetries
- Guess for Lanczos (wave function transformation)
- Everything under m³
- DGEMM should be your best friend