

Fractional quantum Hall effect on an infinite cylinder: **Characterization of topological order using iDMRG**

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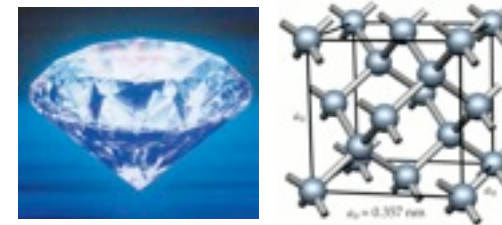
Introduction

- Different phases of matter are usually understood using **local order parameters** (“symmetry breaking”)

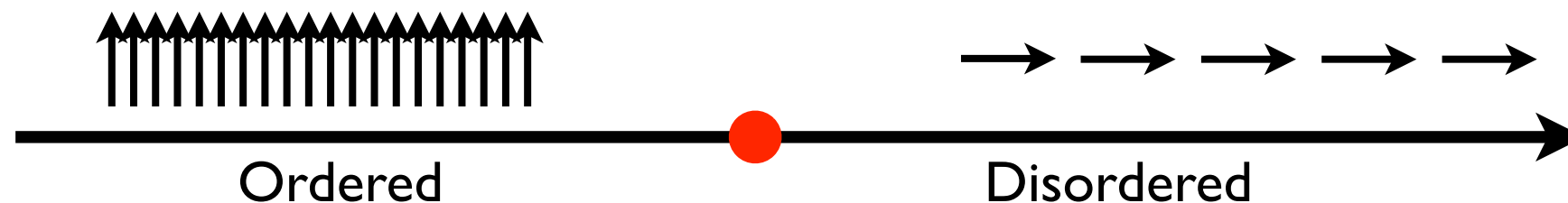
- **Magnets:** spin rotation and TR symmetry broken



- **Crystals:** translation and rotation symmetry broken

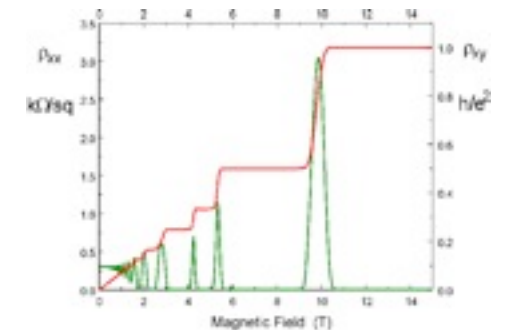


- Phase diagram of an Ising model (magnetization as order parameter)



Introduction

- In the last years several **“new kinds” of phases** have been discovered which **cannot be described by local order parameters**
 - Quantum Hall effect [Klitzing '80]
 - Fractional Quantum Hall effect [Tsui, Stormer '82, Laughlin '83]
 - Topological spin liquids [Anderson '87]
 - ...
- How to characterize topological order?
- How to tell what kind of topological order we have from **numerical simulations?**

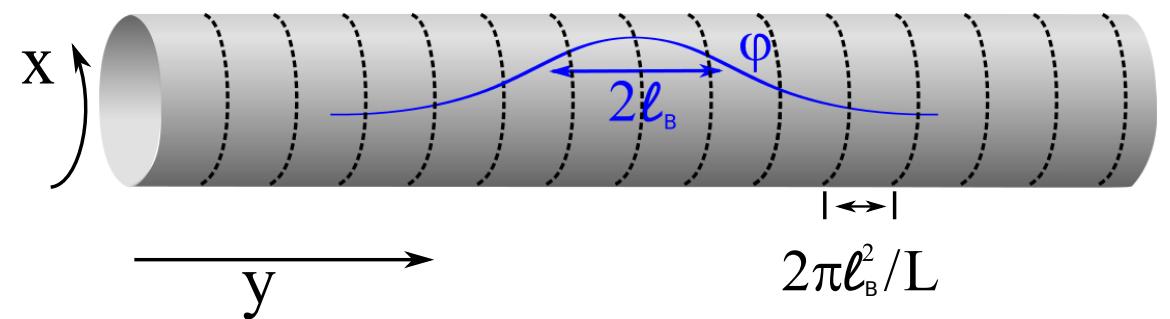


Outline

- **Fractional quantum Hall on a cylinder**
- **Efficient simulations of infinite 1D system:**
Matrix product states (MPS) and density matrix renormalization group (DMRG)
- **Characterization of topological order using MPS and the entanglement spectrum**
 - Charge distribution of a quasi particle
 - Topological entanglement entropy
 - Quantum dimensions
 - Topological spin

Fractional quantum Hall on a cylinder

- Two-dimensional electron gas in a strong magnetic field
- Lowest Landau Level on an infinite cylinder with circumference L



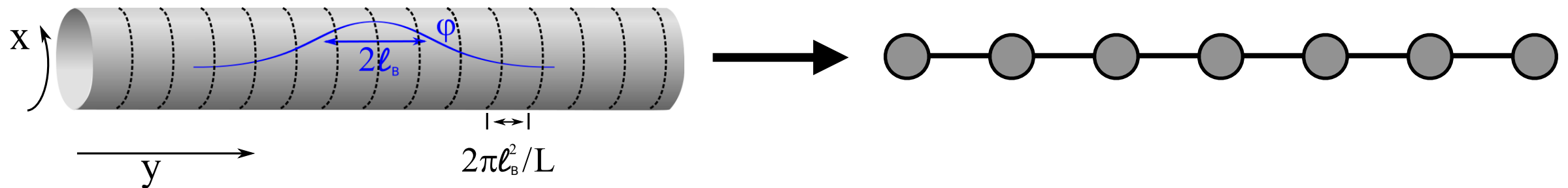
Landau gauge $\mathbf{A} = \ell_B^{-2}(-y, 0)$

- Degenerate orbitals take the form
(with $k_j = \frac{2\pi j}{L}$)

$$\varphi_j(x, y) = \frac{e^{ik_j x - \frac{1}{2\ell_B^2}(y - k_j \ell_B^2)^2}}{\sqrt{L\ell_B \pi^{1/2}}}$$

Fractional quantum Hall on a cylinder

- Orbitals are localized at $y_j = k_j \ell_B^2$: **ID model** using an occupation number basis $|N_j\rangle$, $N_j = 0, 1$



- Degenerate states in the orbital basis represented as

$$\dots | \dots 0010101011 \dots \rangle, | \dots 1000101011 \dots \rangle, \dots$$

Fractional quantum Hall on a cylinder

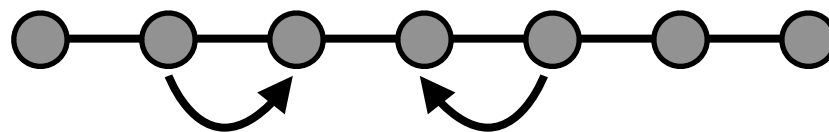
- The most general interaction allowed by symmetry is

$$\hat{H} = \sum_n \sum_{k \geq |m|} V_{km} c_{n+m}^\dagger c_{n+k}^\dagger c_{n+m+k} c_n$$

- We are using the Haldane pseudo potentials V_i
(Trugman-Kivelson $\nabla^2 \delta(r)$: $V_{km} \propto e^{-\frac{1}{2} 4\pi^2 (k^2 + m^2) / L^2}$):

[Haldane '83; Trugman & Kivelson '85]

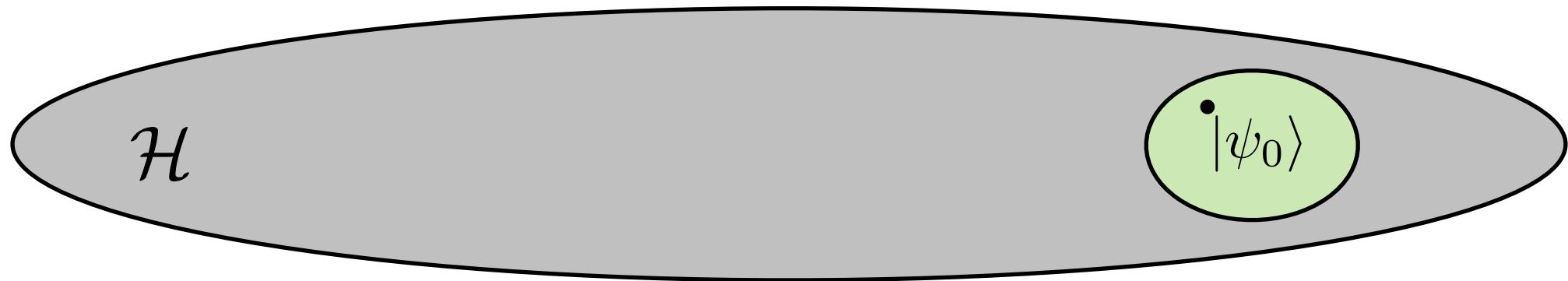
- Total charge $\hat{C} = \sum_j \hat{N}_j$ and momentum $\hat{K} = \sum_j j \hat{N}_j$
are conserved



[Haldane & Rezayi '94; Bergholtz et al.; '05, Seidel et al. '05]

Efficient simulations in 1D

- Consider local Hamiltonians acting on a Hilbert space \mathcal{H} :



- Ground states live in a very small corner of the Hilbert space as they generically fulfill the area law (proven for 1D, gapped systems [Hastings '07])
- Efficiently constrain our simulations to this subspace: In 1D we represent the state as **matrix-product state (MPS)**
- Main idea:** Truncate the Schmidt decomposition

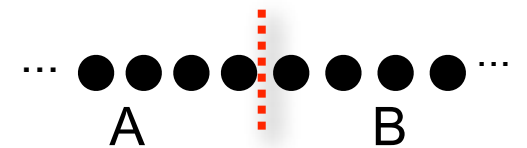
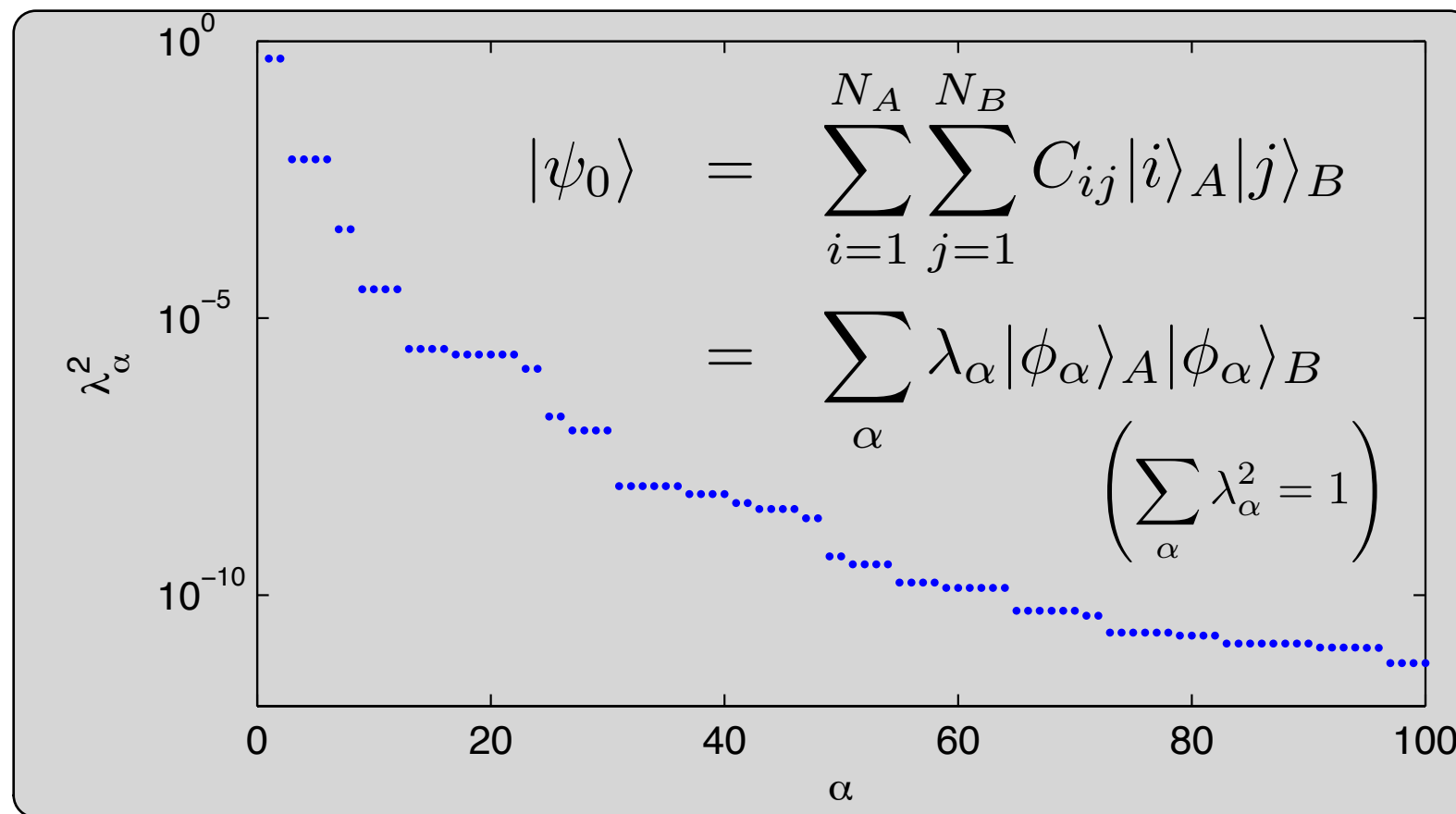
$$|\psi_0\rangle = \sum_{\alpha} \lambda_{\alpha} |\phi_{\alpha}\rangle_A |\phi_{\alpha}\rangle_B \quad \cdots \bullet \bullet \bullet \bullet \bullet \bullet \cdots$$

A B

The diagram shows a sequence of black dots representing sites in a 1D chain. A vertical dashed red line is placed between two dots, dividing the chain into two parts, labeled A and B. The dots are arranged in a slightly wavy line, and there are ellipses at both ends of the sequence.

Efficient simulations in 1D

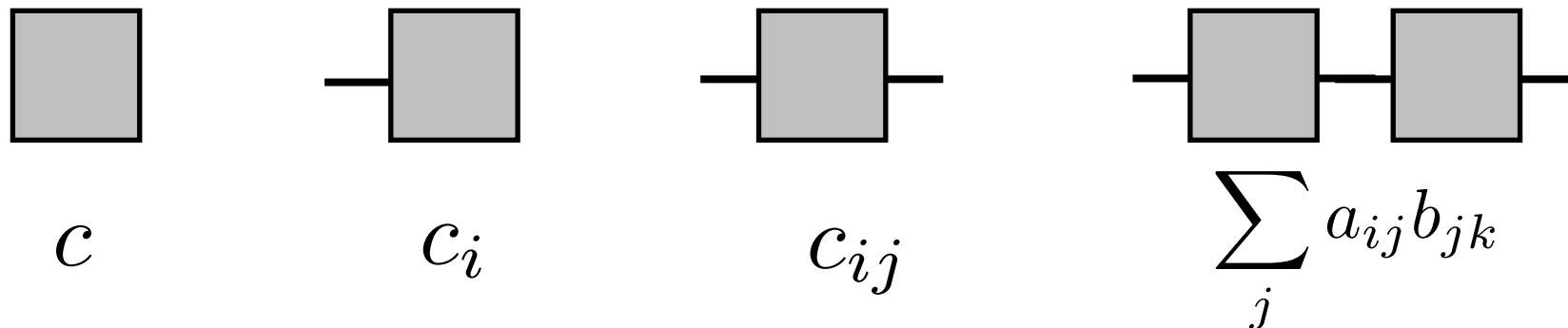
- Example: Entanglement spectrum (Schmidt spectrum) of the Spin-1 Heisenberg chain $H = \sum_j \vec{S}_j \cdot \vec{S}_{j+1}$



- **Schmidt values decay rapidly:** Almost the entire weight is contained in only **few important states (AREA LAW!)**

Efficient simulations in 1D

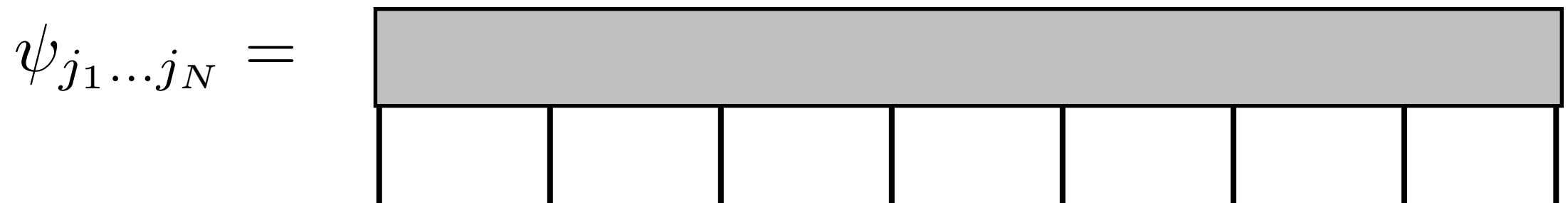
- Useful way to represent tensors



- A generic quantum state for a N site system has a d^N dimensional Hilbert space ($j_n = 1 \dots d$)

$$|\psi\rangle = \sum_{i_1 \dots i_N} \psi_{j_1 \dots j_N} |j_1\rangle \cdots |j_N\rangle$$

- Coefficients are given by a rank N tensor:



Efficient simulations in 1D

- In an **MPS** representation we write the rank-N tensor as a product of rank-3 tensors

[Fannes et al. '92]

$$\psi_{j_1 \dots j_N} = \begin{array}{cccccccc} \boxed{} & \boxed{} & \boxed{} & \boxed{} & \boxed{} & \boxed{} & \boxed{} & \boxed{} \\ | & | & | & | & | & | & | & | \end{array}$$

$$B_{\alpha, \beta}^{[n] j_n} = \alpha \begin{array}{c} \boxed{} \\ | \\ j_n \end{array} \beta \quad \text{with dimensions} \quad \begin{array}{l} j_n = 1 \dots d \\ \alpha, \beta = 1 \dots \chi_n \end{array}$$

- Dimension χ generically grows exponentially with system size: But If we represent $|\psi\rangle$ in terms of the **Schmidt basis at each bond** we know how to truncate it!
- **Variational wave function** for which the number of parameters depends only on entanglement ($N d \chi^2$)

Efficient simulations in 1D

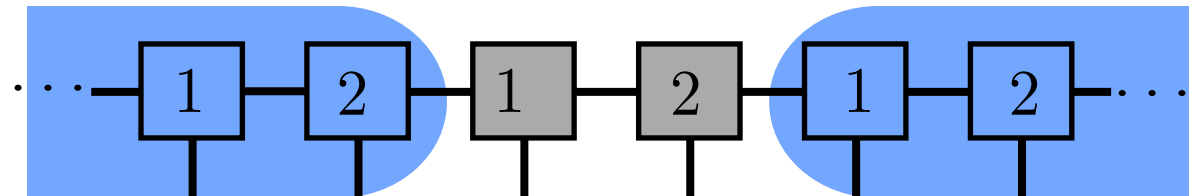
- Local expectation values calculated efficiently ($\propto N\chi^3d$)
- Operators can be expressed in terms of **matrix-product operators (MPO)** as a generalization of MPS
- For infinite systems we assume a finite unit cell such that $B^{[n]} = B^{[n+M]}$

$$\psi_{\dots j_1 \dots j_N \dots} = \dots \text{---} \begin{array}{c} \boxed{1} \\ | \end{array} \text{---} \begin{array}{c} \boxed{2} \\ | \end{array} \text{---} \begin{array}{c} \boxed{3} \\ | \end{array} \text{---} \begin{array}{c} \boxed{1} \\ | \end{array} \text{---} \begin{array}{c} \boxed{2} \\ | \end{array} \text{---} \begin{array}{c} \boxed{3} \\ | \end{array} \text{---} \dots$$

- **FQH**: Matrix dimension χ needed to represent the ground state grows $\propto \exp(L)$: This is bad - but not as bad as a $\propto \exp(L^2)$ growth :-)
- Same scaling with L as in finite system DMRG...
[Shibata '01; Feiguin '08; Hu '12; Zhao '12;...]

Efficient simulations in 1D

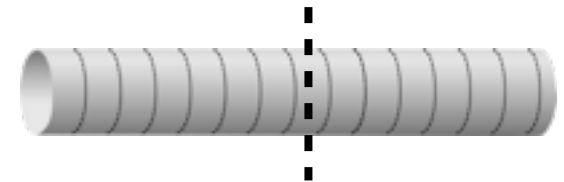
- The ground state is found using an infinite system variant of the DMRG algorithm [White '92; McCulloch '08]
- **Iterative minimization of the energy**
 - Optimizing some tensors while keeping the environment fixed (e.g., Lanczos)



- After each optimization, the wave function is truncated back to the original variational space of MPS's (SVD)
- Sufficient number of iterations yields a fixed point and we can read off the tensors $B^{[1]} \dots B^{[M]}$ which approximate the ground state

iDMRG results: Quasiparticles

- **How can we characterize topological order using iDMRG?**
- Topologically ordered states are characterized by their quasiparticle (QP) excitations: **GS degeneracy equal to the number of QP types** a [Wen '90; Kitaev '05]
- We choose a basis such that each GS corresponds to a specific QP type $\mathcal{H} = \bigoplus_a \mathcal{V}_a$: **minimally entangled states** (MES) for the bipartition of the cylinder $|\Xi_a\rangle$
[Zhang et al. '12; Grover et al. '12; Cincio & Vidal '12]
- Choose a specific “**pattern of zeros**” (e.g., ...0101...) as initial state [Wen & Wang '08; Bernevig '08]
Ergodicity: Two-site update not sufficient



iDMRG results: Quasiparticles

- Example: **Moore-Read phase** (six different quasi-particle sectors with different initial configurations)

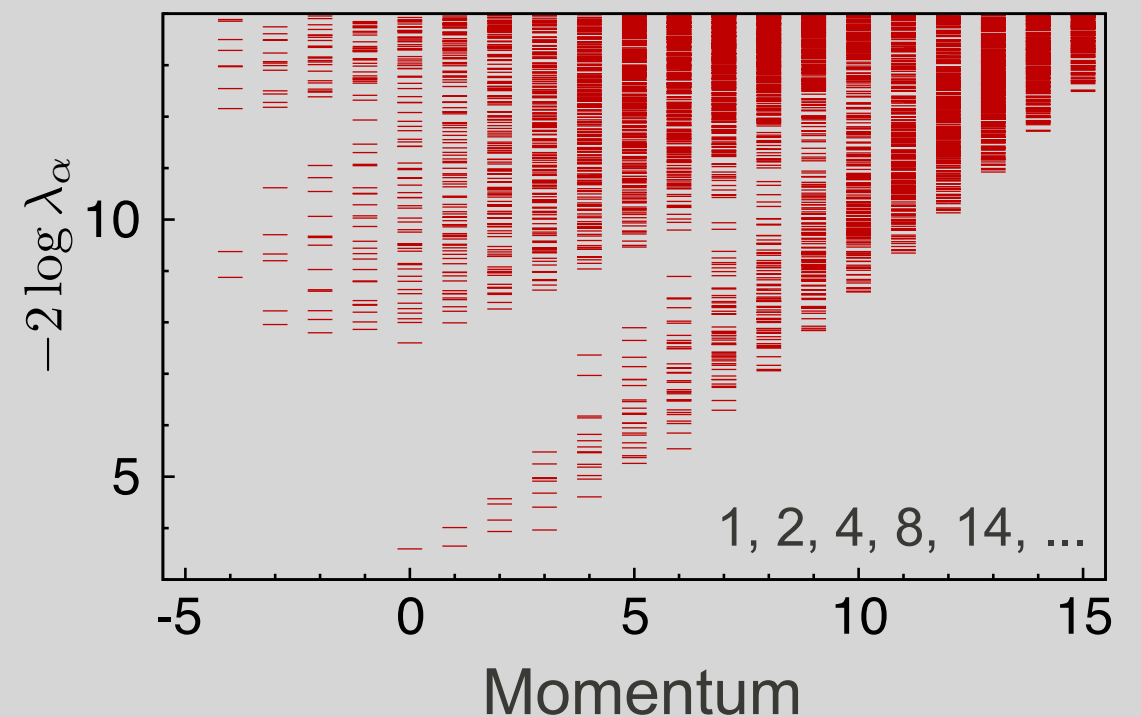
[Moore & Read '91; Fendley '06; Bergholtz '06; Fradkin '08; ...]

QP type	Charge	Seed
1	0	0110
$V_{+1/\sqrt{2}}$	$e/2$	0011
$V_{-1/\sqrt{2}}$	$e/2$	1100
ψ	0	1001
$\sigma V_{+1/2\sqrt{2}}$	$-e/4$	0101
$\sigma V_{-1/2\sqrt{2}}$	$e/4$	1010



Real space entanglement spectrum
in the $\nu = 1/2$ phase:

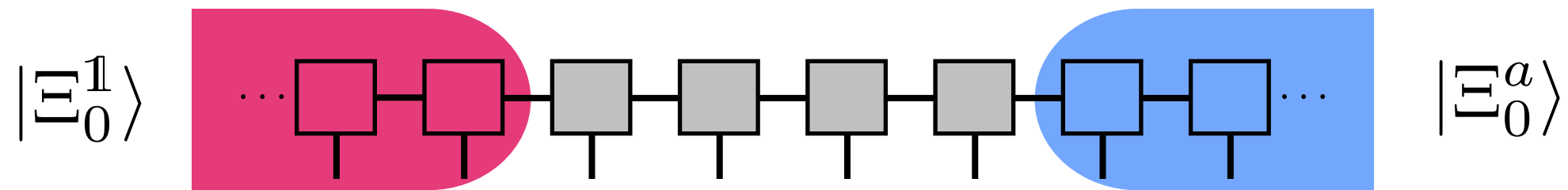
$$L = 21.5\ell_B, V_1 = 1, V_3 = 0.65$$



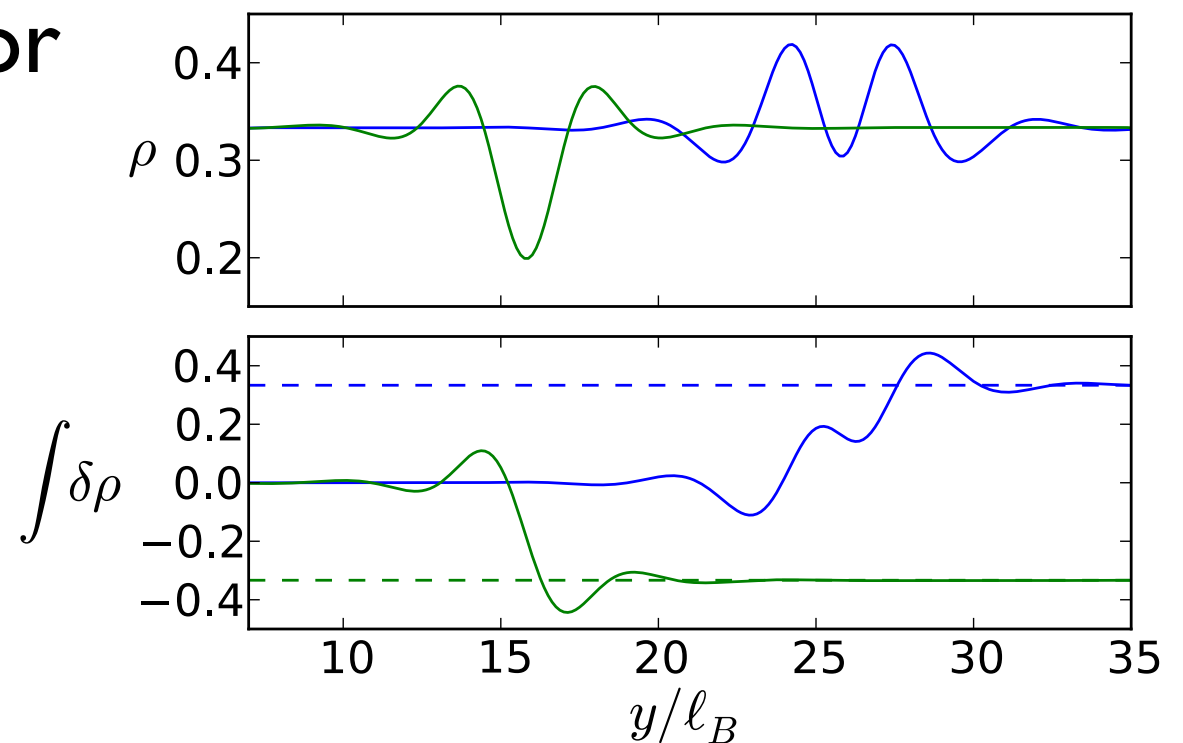
[Read '96; Zaletel '12; Dubail '12;
Sterdyniak '12; ...]

iDMRG results: Quasiparticles

- Creating a QP by glueing together two ground states with different charges at the boundary



- Example: Charge density for the $\nu = 1/3$ state
- QP charge by measuring expectation values of the charge in Schmidt states

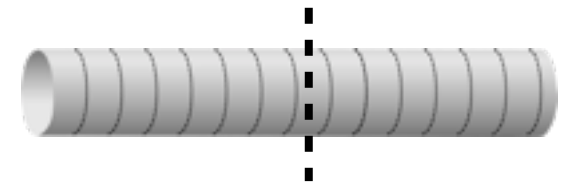


iDMRG results: Entanglement

- The entanglement entropy for a **topologically ordered ground state** $|\Xi_a\rangle$

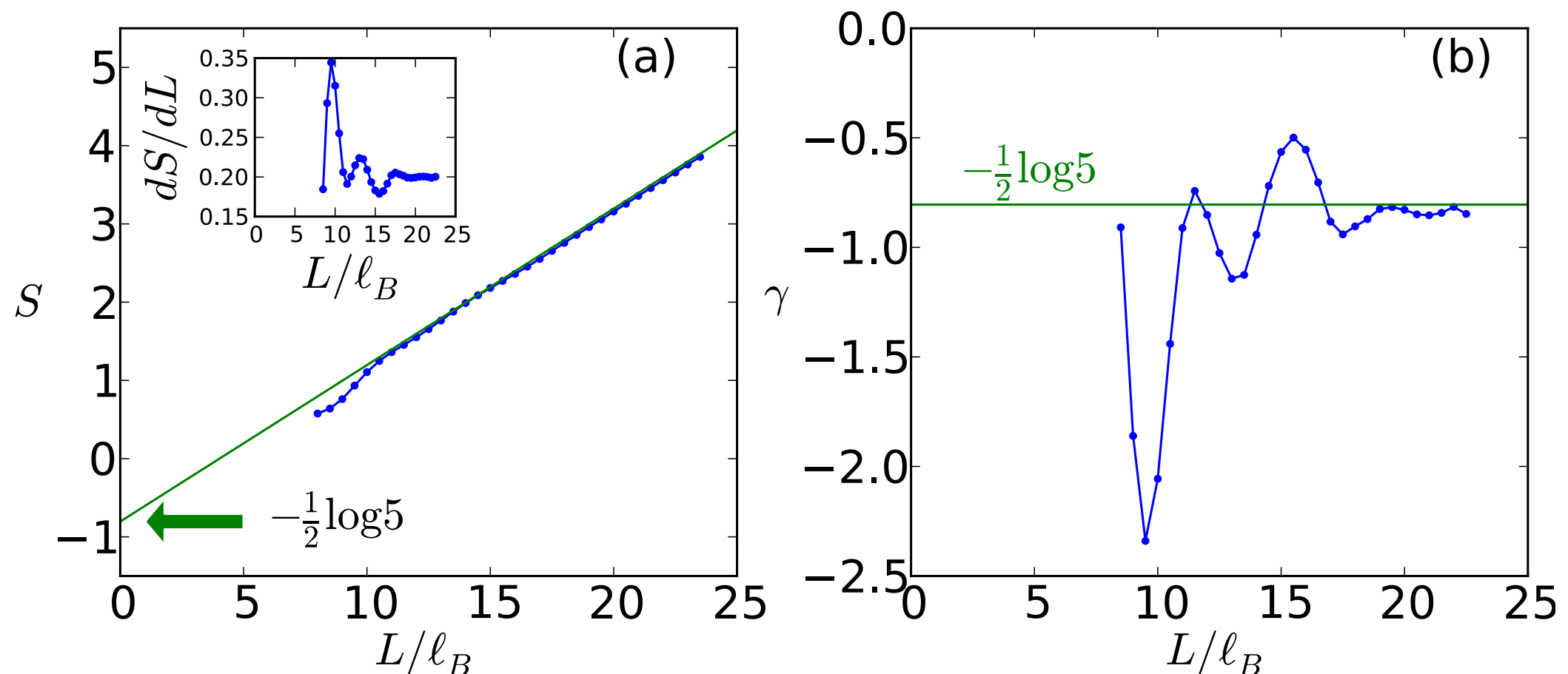
$$S = sL - \gamma_a \quad \text{with} \quad \gamma_a = -\log \left(d_a / \sqrt{\sum_b d_b^2} \right)$$

- For abelian QP $d_a = 1$ and the topological entanglement entropy does not depend on a
- Sum rule allows us to check if we found all ground states $\sum_i e^{-2\gamma_i} = 1$



iDMRG results: Entanglement

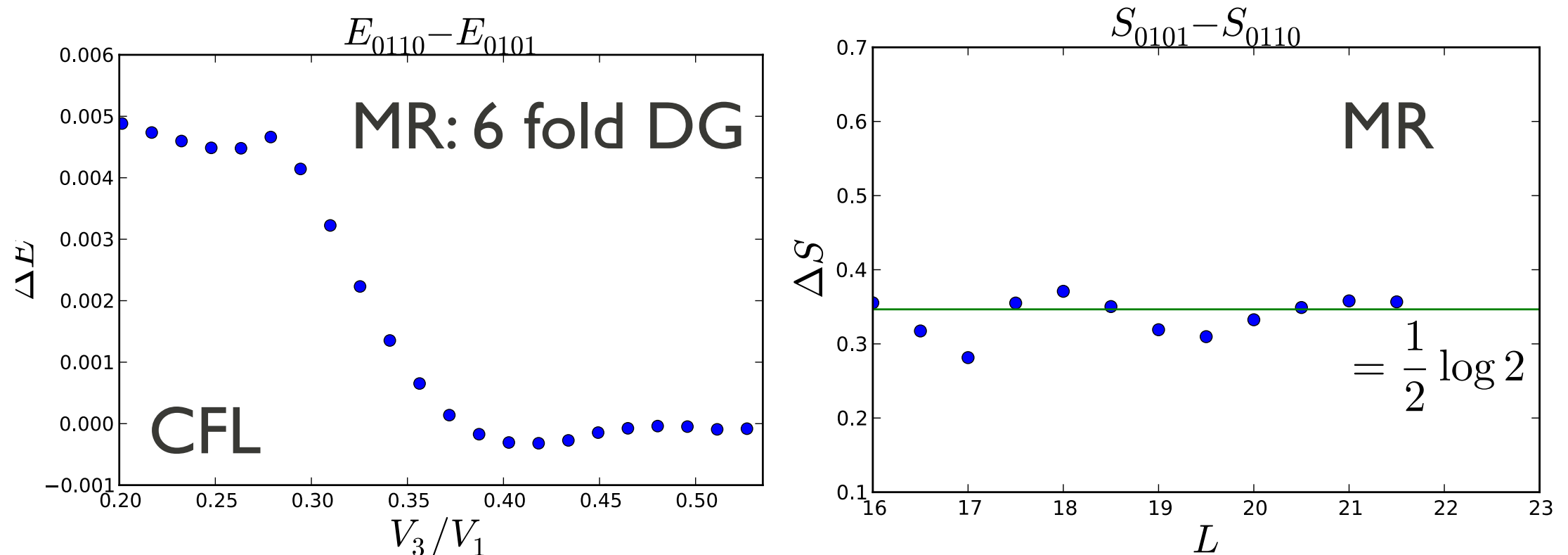
- Example: Entanglement entropy for an (abelian) $\nu = 2/5$ state



➡ Extracted topological entanglement entropy: $\gamma \approx 0.83$

iDMRG results: Entanglement

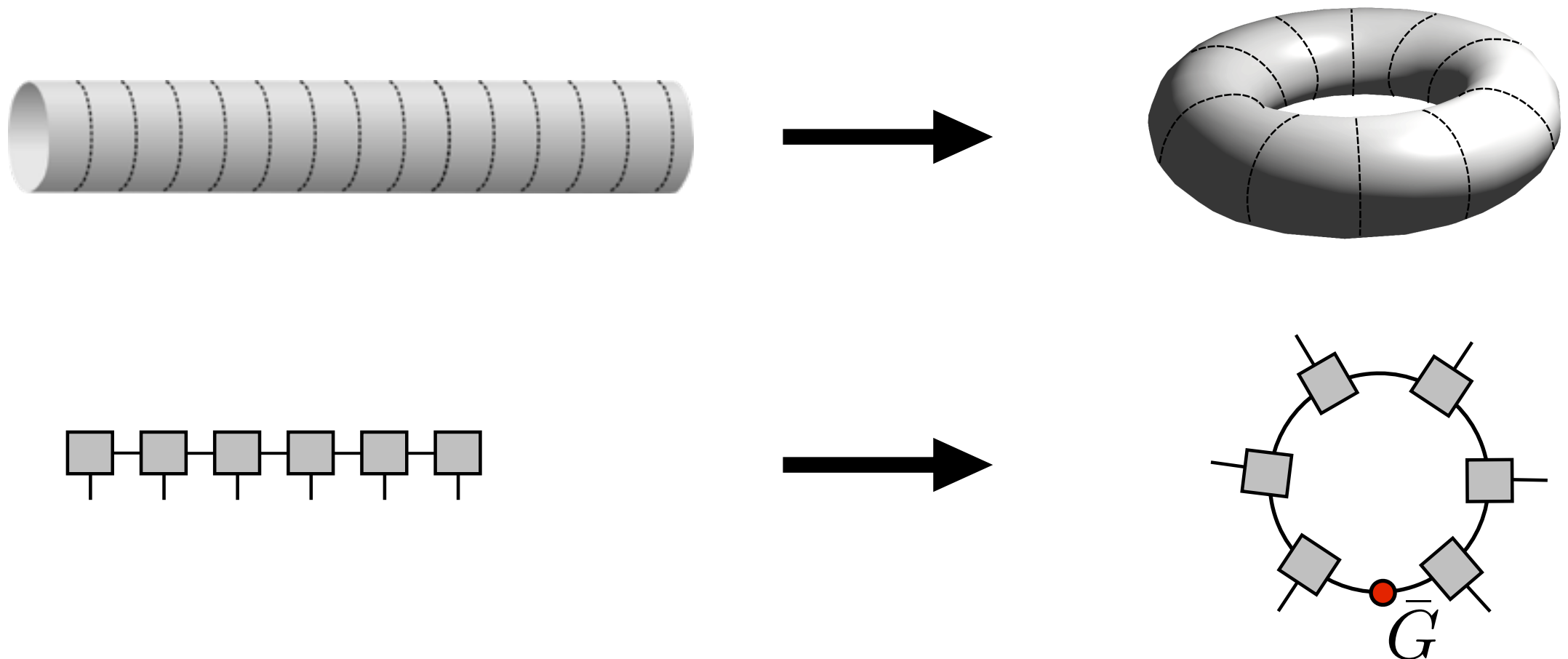
- Phase diagram of the $\nu = 1/2$ state by varying the Haldane pseudo potentials



- Difference in topological entropy $\Delta S = \gamma_{\sigma V_{-1/2\sqrt{2}}} - \gamma_1$
 ➡ Non-abelian anyons with $d_a \approx 1.43$

iDMRG results: From a cylinder to a torus

- Put the ground state of the infinite cylinder onto a torus (correlation length ξ is short compared to the length)

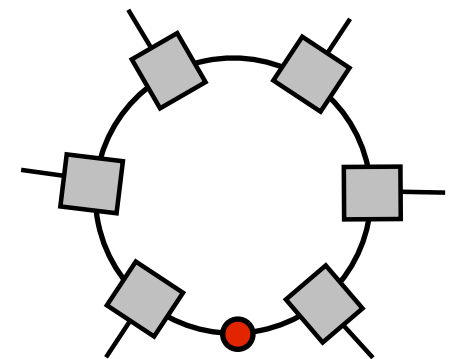


- The operator \bar{G} determines the boundary conditions

iDMRG results: From a cylinder to a torus

- Fluxes and modular parameter can be accounted for by inserting a diagonal matrix • when connecting the two edge auxiliary bonds

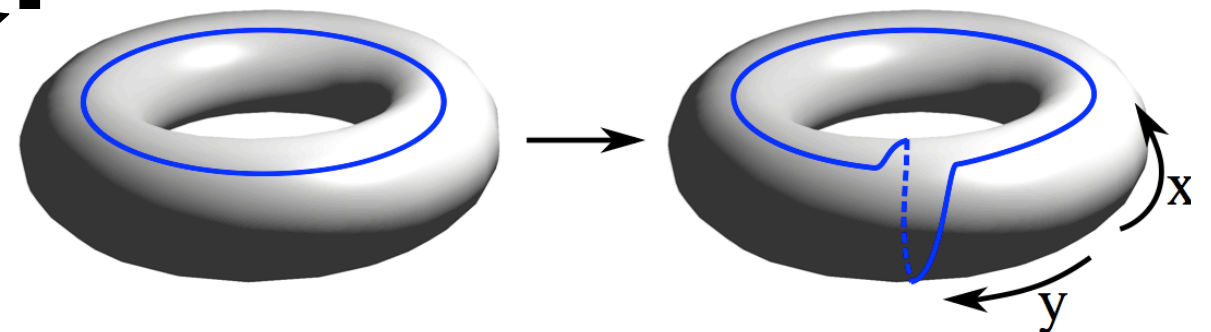
$$\bar{G} = (-1)^{(N_e-1)\frac{\bar{C}}{q}} \exp \left[-2\pi i \tau_x \frac{\bar{K}}{q} + i\Phi_y \frac{\bar{C}}{q} \right]$$



- Inserting a flux Φ_y (Shifting between GS's): **QP charge**

- Dehn Twist (Modular T):
Topological spin of QP

- Numerical verifications in progress...



Summary

- Simulating FQH systems using infinite MPS allows us to extract many relevant quantities
 - Clean entanglement spectra (both in real and orbital space)
 - Topological entanglement entropy and individual quantum dimensions
 - Quasiparticle charges
 - Topological spin

