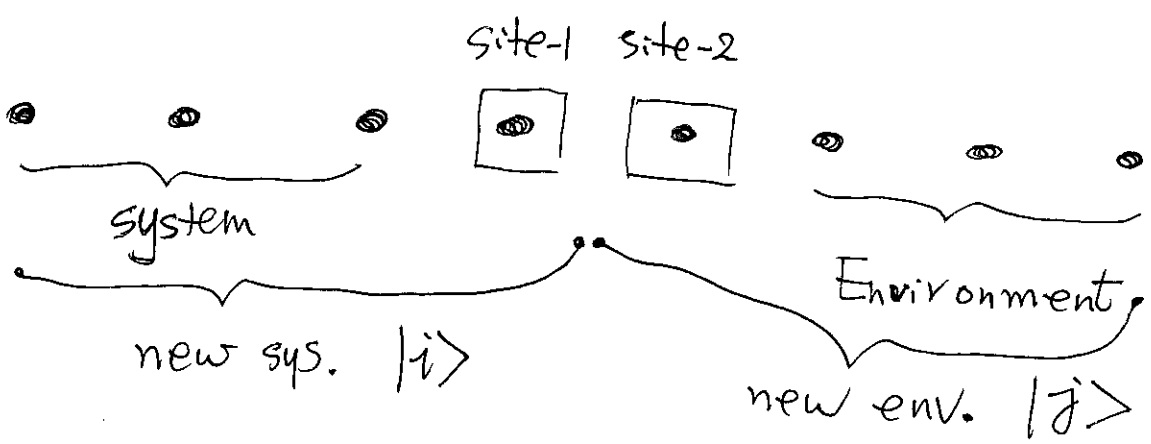


Notes on "DMRG for interacting systems"

- I. DMRG set up and symmetries (e.g. SU2)
- II. Generalize to fermionic systems
- III. Momentum as additional quantum number (or momentum space DMRG) - applying to FQHE
- IV. Examples



- use tensor product between system and environment.

$$|\psi\rangle = \sum_{ij} \psi_{ij} |i\rangle \otimes |j\rangle$$

① Diagonalize H_{full} to obtain $|\psi\rangle$.

② Construct the reduced density matrix

$$\rho_{ii'} = \sum_k \psi_{ik} \psi_{i'k}^*$$

(matrix representation $\psi \sim M, \rho = M M^\dagger$)

diagonalize ρ and keep m eigenvectors corresponding to m largest eigenvalues. $\text{Tr} \rho = 1$.

control truncation error $\epsilon = 1 - \sum_{i=1}^m \rho_i$

③ Construct m new basis vectors from 2^m states

$$|\alpha_l\rangle |S_{l+1}\rangle \Rightarrow |\alpha_{l+1}\rangle$$

$$|\alpha_{l+1}\rangle = \sum_{S_{l+1}, \alpha_l} A_{\alpha_{l+1}, \alpha_l}^{l+1}(S_{l+1}) |\alpha_l\rangle \otimes |S_{l+1}\rangle$$

↑ each row is the eigenvector of \hat{J} .

④ update operators and truncate them, eg.

$$\tilde{H}_{sys} = \underbrace{H'_{sys}}_{\text{previous}} \otimes I_{l+1} + \sum c_{ij} B'_i \otimes B_j^{l+1}$$

$$H_{sys} = O H O^\dagger$$

O is $m \times 2m$ matrix, the rows of O are the kept states.

$$\begin{pmatrix} A^\uparrow_{2l+1, \alpha_l} & A^\downarrow_{2l+1, \alpha_l} \end{pmatrix}$$

* SU2 symmetry

common eigenstates of \hat{S}^2 , S_z and H .

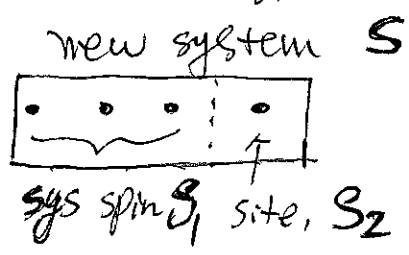
basis states $|s, m, \alpha\rangle$.

The matrix element can be obtained using Wigner-Eckart theorem.

$$\langle s', m', \alpha' | T_{q,i}^{(k)} | s, m, \alpha \rangle = (-1)^{s-m'} \begin{pmatrix} s' & k & s \\ -m' & q & m \end{pmatrix} \underbrace{\langle s', \alpha' || T_i^{(k)} || s, \alpha \rangle}_{\text{reduced element}}$$

$T_i^{(k)}$ the irreducible tensor operator. i -other index, site for example. $T_{q,i}^{(k)}$ it is q component ($q = -k, -k+1, \dots, k$).

1. Set basis



$$|SM\alpha\rangle = \sum_{m_1, m_2} |s_1 m_1 s_2 m_2, \alpha\rangle \langle s_1 m_1 s_2 m_2 | SM \rangle$$

Clebsch-Gordan coefficients representing an additional Unitary transformation to basis set of U1.

2. Identifying irreducible tensors

definition $[\mathbf{S}_i^z, T_{q,i}^{(k)}] = \hbar q T_{q,i}^{(k)}$

$$[S_{\pm i}, T_{q,i}^{(k)}] = \sqrt{(k \mp q)(k \pm q + 1)} \hbar T_{q \pm 1, i}^{(k)}$$

$$T_1^{(1)} = -\frac{1}{\sqrt{2}} S^+$$

$$T_0^{(1)} = S^z$$

$$T_{-1}^{(1)} = \frac{1}{\sqrt{2}} S^-$$

$$\langle S || T^{(1)} || S \rangle = \sqrt{(2S+1)(S+1)S}$$

Two irreducible tensor operator $T^{(k_1)}, U^{(k_2)}$ can form a new irreducible operator $X^{(k)}$

$$X_q^{(k)} = \sum_{q_1 q_2} (-1)^{-k_1 + k_2 + q} \sqrt{2k+1} \begin{pmatrix} k_1 & k_2 & k \\ q_1 & q_2 & -q \end{pmatrix} T_{q_1}^{(k_1)} U_{q_2}^{(k_2)}$$

$$\vec{S}_i \cdot \vec{S}_j = -T_{1i}^{(1)} T_{-1j}^{(1)} - T_{-1i}^{(1)} T_{1j}^{(1)} + T_{0i}^{(1)} T_{0j}^{(1)} \quad \text{rank-0}$$

similarly, $T^{(1)} = \begin{cases} T_1^{(1)} = \frac{i}{\sqrt{2}} \left(-\frac{1}{\sqrt{2}}\right) [(\vec{S}_i \times \vec{S}_j)_x + i(\vec{S}_i \times \vec{S}_j)_y] \\ T_0^{(1)} = \frac{i}{\sqrt{2}} (\vec{S}_i \times \vec{S}_j)_z \\ T_{-1}^{(1)} = \frac{i}{\sqrt{2}} \cdot \frac{1}{\sqrt{2}} [(\vec{S}_i \times \vec{S}_j)_x - i(\vec{S}_i \times \vec{S}_j)_y] \end{cases}$

is useful for systems with three vector chiral terms.

5

3. Beside the transformation to total S basis, and using the reduced matrix elements for irreducible tensor operators, everything else can be carried out similar to U(1)-DMRG.

* Example, t-J model

$$H = \sum_{\langle ij \rangle} -t_{ij} (c_{i\sigma}^{\dagger} c_{j\sigma} + h.c.) + \sum_{\langle ij \rangle} J_{ij} (\vec{S}_i \cdot \vec{S}_j - n_i n_j / 4)$$

Irreducible tensor - rank $-\frac{1}{2}$

$$\begin{cases} T_{-\frac{1}{2}, i}^{(\frac{1}{2})} = c_{i\uparrow} \\ T_{\frac{1}{2}, i}^{(\frac{1}{2})} = -c_{i\downarrow} \end{cases} \quad \begin{cases} T'_{-\frac{1}{2}, i}^{(\frac{1}{2})} = c_{i\downarrow} \\ T'_{\frac{1}{2}, i}^{(\frac{1}{2})} = c_{i\uparrow} \end{cases}$$

Thus that the hopping term H_t can be expressed as

$$H_t = \sum_{\langle ij \rangle} -t_{ij} (T_{\frac{1}{2}, i}^{(\frac{1}{2})} T_{-\frac{1}{2}, j}^{(\frac{1}{2})} + T_{\frac{1}{2}, j}^{(\frac{1}{2})} T_{-\frac{1}{2}, i}^{(\frac{1}{2})} - T'_{-\frac{1}{2}, i}^{(\frac{1}{2})} T'_{\frac{1}{2}, j}^{(\frac{1}{2})} - T'_{-\frac{1}{2}, j}^{(\frac{1}{2})} T'_{\frac{1}{2}, i}^{(\frac{1}{2})})$$

Two irreducible $T^{(k_1)}, T^{(k_2)}$ can compose a new tensor $X^{(k)}$:

$$X_{q_1 q_2}^{(k)} = \sum_{q_1 q_2} T_{q_1}^{(k_1)} T_{q_2}^{(k_2)} S_{q_1 q_2, k, q} = \sum_{q_1 q_2} (-1)^{-k_1 + k_2 + q} \sqrt{2k+1} \begin{pmatrix} k_1 & k_2 & k \\ q_1 & q_2 & q \end{pmatrix} T_{q_1}^{(k_1)} T_{q_2}^{(k_2)}$$

$$\text{So } X_0^{(0)} = \sum_{q_1 q_2} T_{q_1}^{(\frac{1}{2})} T_{q_2}^{(\frac{1}{2})} \begin{pmatrix} \frac{1}{2} & \frac{1}{2} & 0 \\ q_1 & q_2 & 0 \end{pmatrix} = \frac{1}{\sqrt{2}} (T_{\frac{1}{2}}^{(\frac{1}{2})} T_{-\frac{1}{2}}^{(\frac{1}{2})} - T_{-\frac{1}{2}}^{(\frac{1}{2})} T_{\frac{1}{2}}^{(\frac{1}{2})})$$

Therefore, we have $H_t = \sum_{\langle i, j \rangle} -\sqrt{2} t_{ij} (X_{0, (ij)}^{(0)} + X_{0, (ji)}^{(0)})$

II. DMRG for fermionic system

In DMRG, basis states usually form block structure according to their quantum numbers.

basis for system $|q_{n1}, q_{n2}, \dots\rangle$

$q_{n1} \dots$ maybe S_z^{tot} of system

$q_{n2} \dots$ maybe electron number N_e^{tot} of system

operators are sparse matrices, which connect certain basis states with fixed change of quantum numbers.

$c_{i\uparrow}^+$ gives $\Delta q_{n1} = 1/2, \Delta q_{n2} = 1$. etc.

(In DMRG, you match the quantum number between operators and basis.)

Jordan-Wigner transformation is easy to implement in such a system

$$c_{i\sigma}^+ = e^{i\pi \sum_{l < i} n_l} b_{i\sigma}^+$$

$b_{i\sigma}^+$ is the corresponding bosonic creation operator, which adds one particle at site i to the basis state (its a hardcore boson operator).

For $C_{i\sigma}^+$ in the system, the phase $e^{i\pi \sum_{l < i} n_l}$ will be combined with $b_{i\sigma}^+$ when i -site is being included into system. But for $C_{i\sigma}^+$ in the environment, we incorporate part of sign

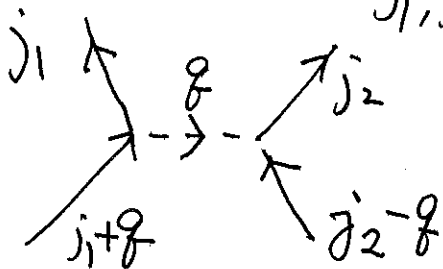
$$C_{i\sigma}^+(\text{env}) \sim e^{i\pi \sum_{\substack{l < i \\ l \in \text{env}}} n_l}$$

When dealing with H_{full} , all remaining signs for operators from environment or sys/env sites should be taken into account. They are easy to get since each basis has a good quantum number for electron number.

III Momentum space DMRG set up

$$H = \sum_j \epsilon_j C_j^+ C_j + \sum_{j_1 j_2 j_3 j_4} A_{j_1 j_2 j_3 j_4} C_{j_1}^+ C_{j_2}^+ C_{j_3} C_{j_4}$$

$$= \sum_j \epsilon_j C_j^+ C_j + \sum_{j_1, j_2, \delta} A_{j_1, j_2, j_2 - \delta, j_1 + \delta} C_{j_1}^+ C_{j_2}^+ C_{j_2 - \delta} C_{j_1 + \delta}$$



C_j^+ creates an electron with momentum j . The total momentum is conserved.

* Adding q_{n3} ~ total momentum to basis quantum numbers.

$$|q_{n1}, q_{n2}, q_{n3}, \dots\rangle$$

$$\begin{matrix} \uparrow & \uparrow & \uparrow & \dots \\ S^z & N_e & q & \dots \end{matrix}$$

* Order the summation to speed up

$$H = \sum_{\substack{j_1, j_2, q \\ \{j_1 < j_2 \\ j_1 + q < j_2 - q\}}} R_{j_1 j_2, j_2 - q, j_1 + q} C_{j_1}^+ C_{j_2}^+ C_{j_2 - q} C_{j_1 + q}$$

$$R_{j_1 j_2 j_3 j_4} = A_{j_1 j_2 j_3 j_4} + A_{j_2 j_1 j_4 j_3} - A_{j_2 j_1 j_3 j_4} - A_{j_1 j_2 j_4 j_3}$$

$$A_{j_1 j_2 j_3 j_4} = \int \psi_{j_1}^*(\vec{r}_1) \psi_{j_2}^*(\vec{r}_2) V(\vec{r}_1 - \vec{r}_2) \psi_{j_3}(r_2) \psi_{j_4}(r_1)$$

* Combining similar terms to reduce the number of operators, e.g.

define $\sum_{j_2, q} R_{j_1 j_2, j_2 - q, j_1 + q} C_{j_1}^+ C_{j_2 - q} C_{j_1 + q} \equiv C_3(j_1)$

$H \rightarrow \sum_{j_1} C_{j_1}^+ C_3(j_1)$ Only when 3 C operators are all in system or environment!
(keep remaining terms unchanged).

- * carefully including all terms to your H_{full} (so many terms...)
- * benefit is that we can target states with different q^{tot} . It works for torus, infinite cylinder (or finite) or sphere.

* Exact diagonalization program for FQHE (torus example)

The single particle wavefunction: (N -Landau level index)

$$\psi_{N,j}(x,y) = \left(\frac{1}{2^N N! \pi^{1/2} L_y l} \right)^{1/2} e^{i \frac{x_j}{l^2} y - \left(\frac{x_j - x}{2l^2} \right)^2} H_N \left(\frac{x_j - x}{l} \right)$$

(infinite system $L_x = \infty$).

$$x_j = \frac{2\pi l^2}{L_y} j, \quad j = 0, 1, 2, \dots, N_s - 1.$$

But, we should use the finite one: $L_x L_y = 2\pi l^2 N_s$

↑ number of flux

$$\sum_{k=-\infty}^{+\infty} e^{i \frac{x_{jk} y}{l^2}} e^{-\frac{(x - x_{jk})^2}{2l^2}} H_N \left(\frac{x_{jk} - x}{l} \right)$$

$$x_{jk} = \frac{2\pi l^2 j}{L_y} + k L_x \quad (\text{forcing the periodic condition}).$$

$$V_{(j_1 j_2 j_3 j_4)} = \frac{1}{2} \int d^2 \vec{r}_1 d^2 \vec{r}_2 \psi_{N j_1}^* \psi_{N j_2}^* V(\vec{r}_1 - \vec{r}_2) \psi_{N j_3} \psi_{N j_4}$$

set up ED code for $N_s = 6$ (flux number) and $N_e = 2$ (electron number).

totally $\frac{N_s!}{N_e! (N_s - N_e)!}$ states = $\frac{6 \cdot 5}{2} = 15$

110000
101000 ----- too many, wait, use k -total momentum!

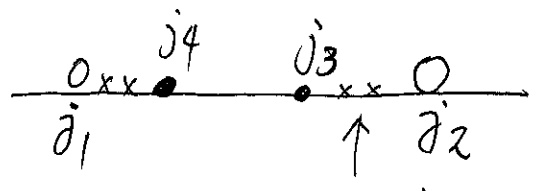
$k=0$ sector (mod N_s).

000, 001, 1+5=0
 001, 010, 2+4=0

$$\begin{pmatrix} R_{j_1 j_2 j_3 j_4} = A_{j_1 j_2 j_3 j_4} + A_{j_2 j_1 j_4 j_3} \\ -A_{j_2 j_1 j_3 j_4} - A_{j_1 j_2 j_4 j_3} \end{pmatrix}$$

only 2 states $H = \begin{pmatrix} R_{1551} & f R_{1542} \\ f R_{1542}^* & R_{2442} \end{pmatrix}$

f - fermion sign ($f=1$ here)
 check (if $N_e > 2$).

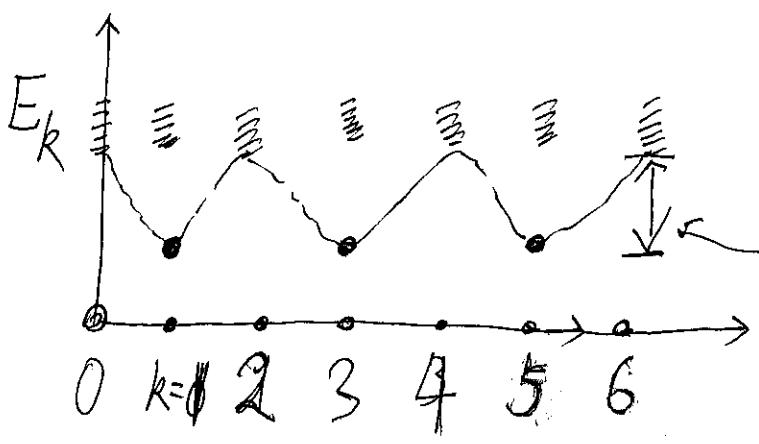


are there particles in between to give fermionic signs?

$k=1$ sector

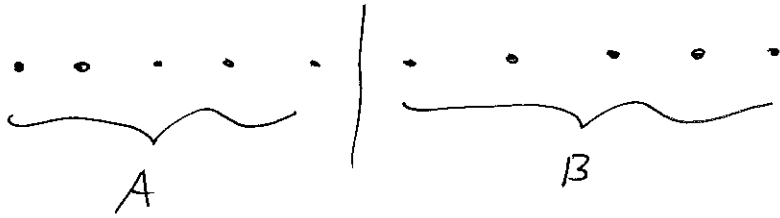
10, 000
 001, 001
 000, 110

$$\begin{pmatrix} R_{0110} & f_1 R_{0152} & f_2 R_{0143} \\ * & R_{2552} & f_3 R_{2543} \\ * & * & R_{3443} \end{pmatrix}$$



3-fold degenerating Laughlin FQHE
 excitation gap is finite and robust

more electrons, separating orbits to A & B



$|i_A\rangle \equiv |n_0, n_1, n_2, \dots, n_{N_s/2}\rangle \dots$
 occupation number for a set of orbitals in A.

a table for finding i_A from $\{n_j\}$.

so once you use $C_{j_1}^\dagger C_{j_2}^\dagger \dots |i_A\rangle$ etc
 you can get new index $|i_A'\rangle$ quickly.

total basis $|\mathbf{k}\rangle \equiv |i_A\rangle \otimes |j_B\rangle$

new index, adding one basis when

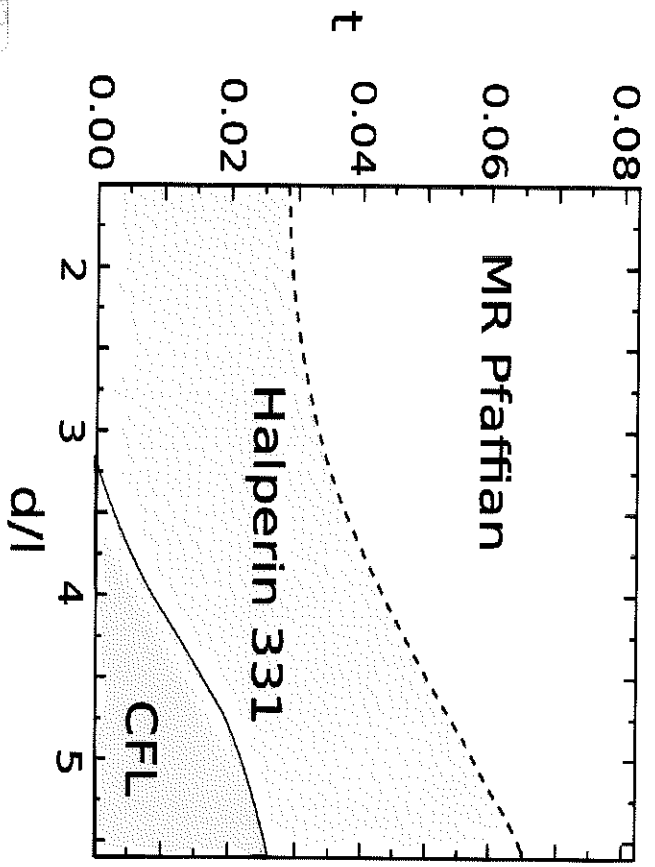
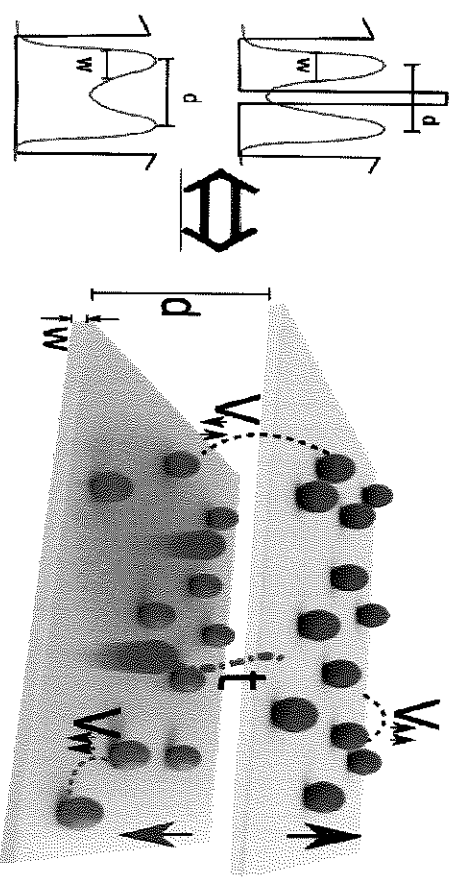
$$J_n(i_A) + J_n(j_B) = \text{desired momentum } k.$$

a table for finding i quickly from (i_A, i_B) .

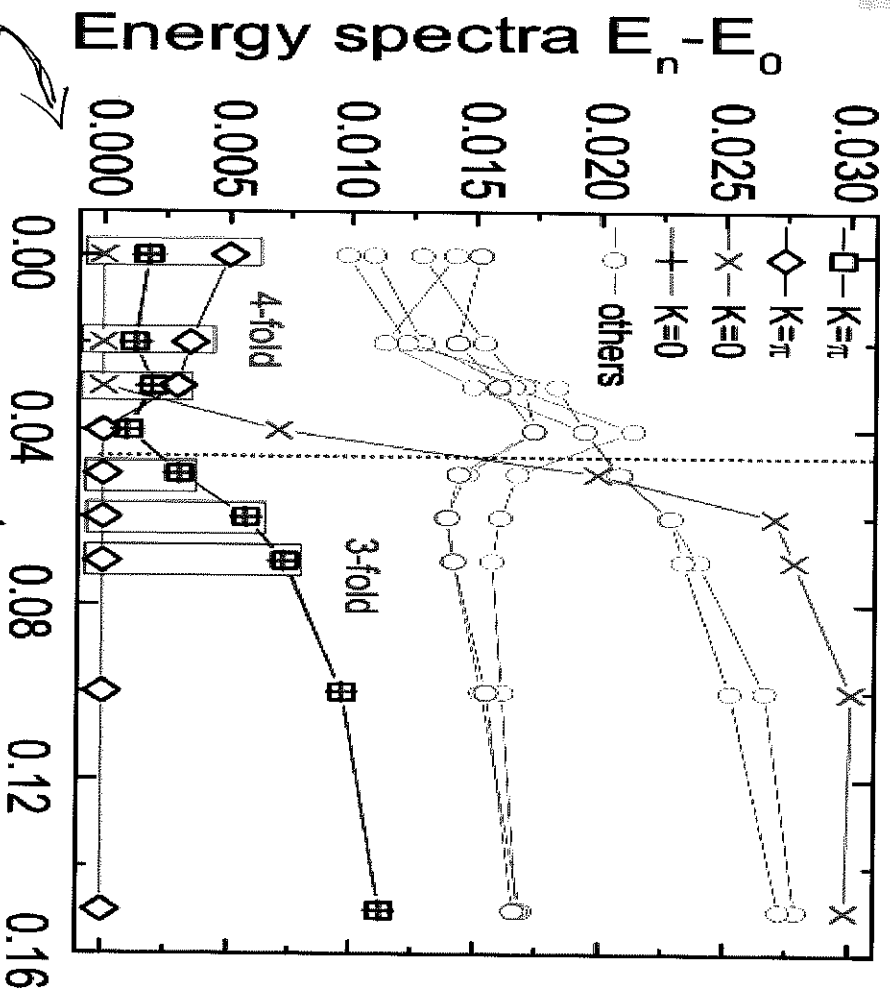
Set up matrix elements, then we can diagonalize H
 for a chosen total momentum k . $|\Psi\rangle = \sum_{ij} \Psi_{ij} |i_A\rangle \otimes |j_B\rangle$

$E(k)$ gives information of topological degeneracy.

Fractional Quantum Hall Bilayers at Half-Filling: Tunneling-driven Non-Abelian Phase



Predicted by Xiao-Gang Wen & other works



Energy spectrum for torus system with varying tunneling t

* Calculating topological Chern number

adding generalized boundary phase θ (α, β)

$$\Upsilon(L_x \hat{x}) \psi_{Nj}(x, y) = e^{i\alpha L_x} \psi_{Nj}(x, y)$$

$$\Upsilon(L_y \hat{y}) \psi_{Nj}(x, y) = e^{i\beta L_y} \psi_{Nj}(x, y)$$

Υ is the single-particle magnetic translational operator in the x or y directions.

$$\psi_{Nj}(x, y) = \left(\frac{1}{2^N N! \pi^{1/2} L_y l} \right)^{1/2} e^{i \frac{x_{jk}'}{l^2} - (x - x_{jk}')^2 / 2l^2} \cdot H_N \left(\frac{x_{jk}' - x}{l} \right)$$

\uparrow \nwarrow
 L index momentum index
in Landau gauge

$$x_{jk}' = \frac{2\pi l^2 j}{L_y} + k_x + \beta l^2$$

Nontrivial: changing $\beta = 0$ to $\frac{2\pi}{L_y}$, will increase $j \rightarrow j+1$ (representing the center of the orbital).

Obtain many-body wavefunction as before

$$|\Psi(\alpha, \beta)\rangle = \sum \psi_{ij}(\alpha, \beta) |i_A(\alpha, \beta)\rangle \otimes |j_B(\alpha, \beta)\rangle$$

(One should also do unitary transformation, so that

$$\psi \rightarrow e^{-i\alpha(x_1 + \dots + x_N)} e^{-i\beta(y_1 + \dots + y_N)} \psi$$

$$\text{Hamiltonian } \left(-i \frac{\partial}{\partial x_i} \rightarrow -i \frac{\partial}{\partial x_i} + \alpha, -i \frac{\partial}{\partial y} \rightarrow -i \frac{\partial}{\partial y} + \beta \right)$$

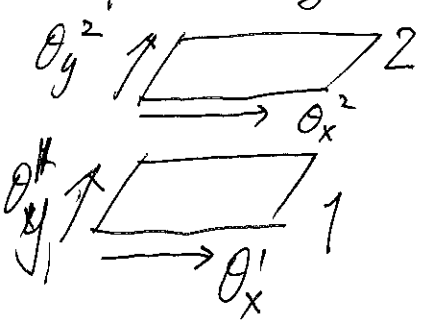
$$C = \frac{i}{4\pi} \iint d\theta_x d\theta_y \left(\left\langle \frac{\partial \Psi}{\partial \theta_x} \middle| \frac{\partial \Psi}{\partial \theta_y} \right\rangle - \left\langle \frac{\partial \Psi}{\partial \theta_y} \middle| \frac{\partial \Psi}{\partial \theta_x} \right\rangle \right) \text{ with } \begin{cases} \theta_x = \alpha L_x \\ \theta_y = \beta L_y \end{cases}$$

For one component FQHE with no disorder, if there is a spectrum gap between ground states and excited states, we have $\Psi_{ij}(\alpha, \beta) \equiv \Psi_{ij}(0, 0)$

(interaction does not depend on boundary phases although $\Psi_{N,ij}$ (orbitals) does).

All Berry phase comes from the rotation of the basis vectors, the Chern number = filling number = N_e/N_s .

The bilayer case is nontrivial.



if (θ_x^1, θ_y^1) change, but $(\theta_x^2 = \theta_y^2 = 0)$,

the interlayer interaction will depend on boundary phases, we can obtain ν drag Hall conductance.

$$C^{\alpha\beta} = \frac{i}{4\pi} \iint d\theta_x^\alpha d\theta_y^\beta \left\{ \left\langle \frac{\partial \Psi}{\partial \theta_x^\alpha} \middle| \frac{\partial \Psi}{\partial \theta_y^\beta} \right\rangle - \left\langle \frac{\partial \Psi}{\partial \theta_y^\beta} \middle| \frac{\partial \Psi}{\partial \theta_x^\alpha} \right\rangle \right\}$$

forms 2x2 Chern number matrix.

for 331 Halperin state: $C = \frac{1}{8} \begin{pmatrix} 3 & -1 \\ -1 & 3 \end{pmatrix}$ inverse of

K matrix. $K = \begin{pmatrix} 3 & 1 \\ 1 & 3 \end{pmatrix}$.

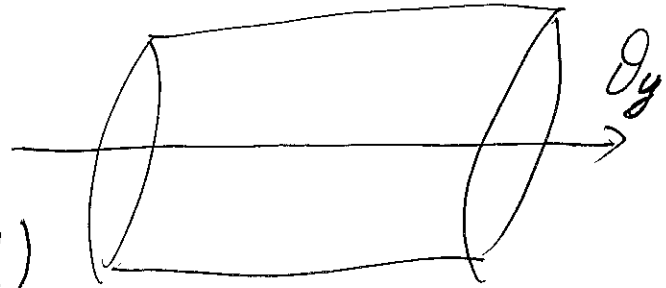
IV Examples

* DMRG inserting flux to detect topological degeneracy

$$S_i^+ S_j^- \rightarrow e^{i\theta_y} S_i^+ S_j^-$$

if the bond crosses the

y -boundary (e.g. $i_y = L_y, j_y = 1$)



using $\psi(\theta_y - \Delta\theta)$ as the initial wavefunction for

new θ_y . Updating $H(\theta_y)$ locally when we sweep pass

different points.

measuring $\langle S_i^z \rangle$ for different θ_y .

comparing entanglement spectrum as function of θ_y .

* Spontaneous Quantum Hall Effect

Using complex code to work with minimum entangled state

if time-reversal symmetry is spontaneously broken

(In real code, ground state would have additional double degeneracy).

* Topological entanglement spectrum and entropy for detecting topological order (FQHE)

root configuration

• 0 0 0, • 0 0 for $1/3$ FQHE
 1 0 0, 1 0 0

{ 1 1 1 0 0, 1 1 1 0 0 for Read-Rezayi $13/5$
 { 1 0 1 0 1, 1 0 1 0 1, 1 0 1 0 1

TABLE I: In this table, we analyze the counting rule of the edge excitations in the "... 11100111001110|00 ..." sector, which has multiplicities 1, 1, 3, 6, ... at $\Delta L = 0, 1, 2, 3, \dots$

$\Delta L = 0$	$\Delta L = 1$	$\Delta L = 2$	$\Delta L = 3$	$\Delta L = 4$
11100111001110 0000	11100111001101 0000	11100111001100 1000	11100111001100 0100	11100111001100 0010
	✓	11100111001011 0000	11100111001010 1000	11100111001010 0100
		11100110101101 0000	11100111000111 0000	11100110101100 0100
		✓	11100110101011 0000	11100111001001 1000
		✓	11100110101100 1000	11100111000110 1000
			11010110101101 0000	11100110101010 1000
				11100110100111 0000
				11100101101011 0000
				11010110101011 0000
				11100110011100 1000
				11010110101100 1000
				11010110101101 0000

TABLE II: In this table, we analyze the counting rule of the edge excitations in the ... 111001110011|00 ... sector, which has multiplicities 1, 2, 5, 9, ... at $\Delta L = 0, 1, 2, 3, \dots$

$\Delta L = 0$	$\Delta L = 1$	$\Delta L = 2$	$\Delta L = 3$
111001110011 0000	111001110010 1000	111001110010 0100	111001110010 0010
	111001101011 0000	111001110001 1000	111001110001 0100
		111001101010 0000	111001101010 0100
		111001100111 0000	111001101001 1000
		110101101011 0000	111001100110 1000
			111001011010 1000
			110101101010 1000
			110101100111 0000
			110100110101101011 0000

TABLE III: In this table, we analyze the counting rule of the edge excitations in the ... 101011010110101| ... sector, which has multiplicities 1, 3, 6, 13, ... at $\Delta L = 0, 1, 2, 3, \dots$

$\Delta L = 0$	$\Delta L = 1$	$\Delta L = 2$	$\Delta L = 3$
101011010110101 0000	101011010110100 1000	101011010110100 0100	101011010110100 0010
	101011010110011 0000	101011010110010 1000	101011010110010 0100
	101011010101101 0000	101011010101100 1000	10101101010100 0100
		101011010101011 0000	101011010110001 1000
		101011001110011 0000	101011010101010 1000
		101010110101101 0000	101011001110010 1000
			101011010011100 1000
			101010110101100 1000
			101011010100111 0000
			101011001101011 0000
			100111001110011 0000
			101010110101011 0000
			011010110101101 0000

TABLE IV: In this table, we analyze the counting rule of the edge excitations in the ... 1010110101100|00 ... sector, which has multiplicities 1, 2, 5, 10, ... at $\Delta L = 0, 1, 2, \dots$

$\Delta L = 0$	$\Delta L = 1$	$\Delta L = 2$	$\Delta L = 3$
1010110100 0000	1010110010 0000	1010110001 0000	1010110000 1000
	1010101100 0000	1010101010 0000	1010101001 0000
		1001110010 0000	1001110001 0000
		1010011100 0000	1010100110 0000
		0110101100 0000	1010011010 0000
			1001101010 0000
			0110101010 0000
			0110011100 0000
			011010110101100 0000
			100111001110010 0000

III. Results for finite-layer thickness

In a two dimensional system, the ideal Coulomb interaction between electrons has $V(q) = 1/q$. The finite thickness in the normal direction of an experimental quantum Hall system modifies the short-distance part of the ideal 2D interaction, yielding

Entanglement Spectrum: 12/5 FQHE as Read-Rezayi state

Detected by Entanglement Spectrum:

111001110011100.....

1: 1,1,3,6,12....

111001110011.....

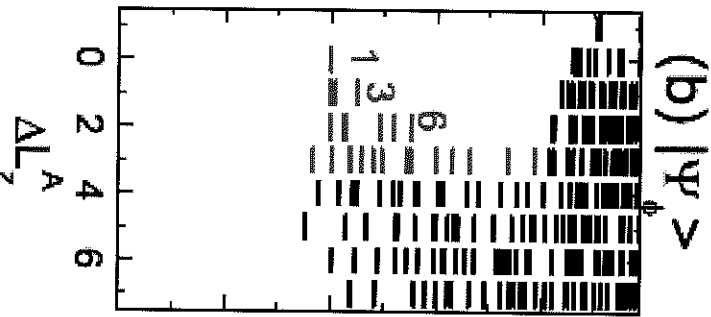
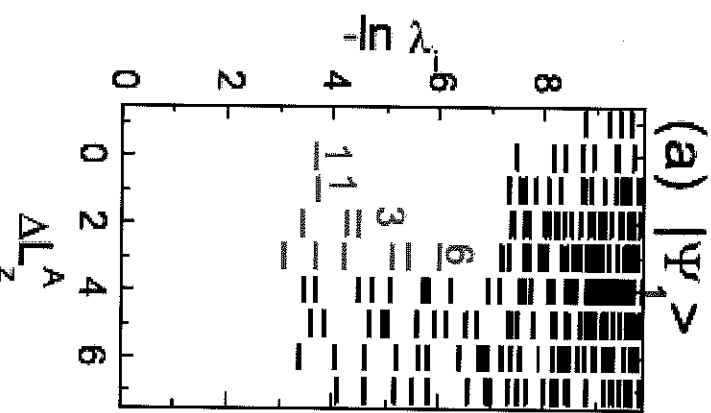
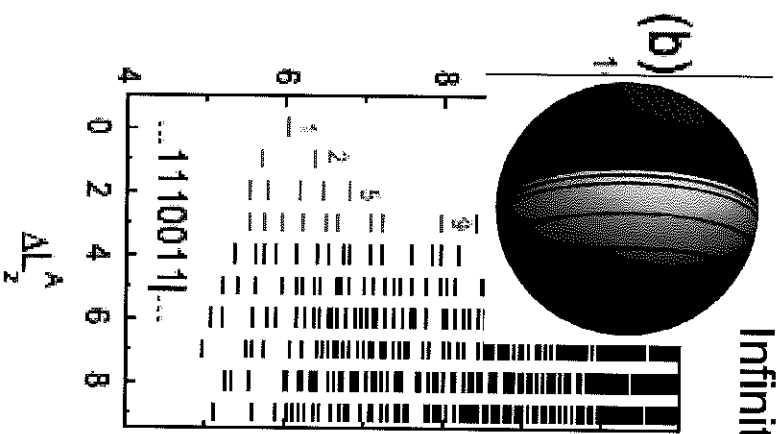
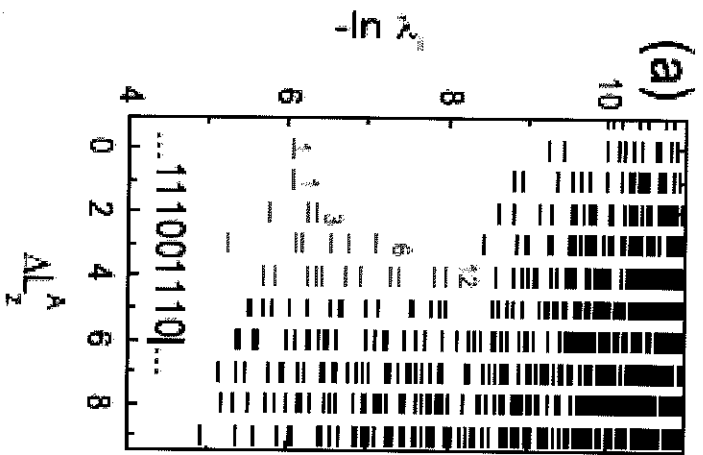
: 1,2,5,9,....

On cylinder, we also find another ground state: 10101 10101 10101

Entanglement spectrum:

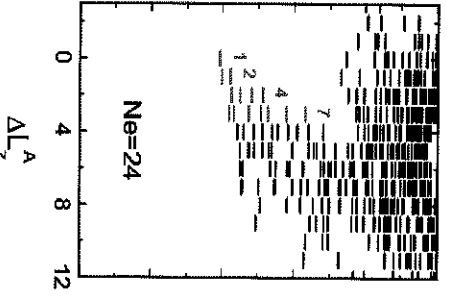
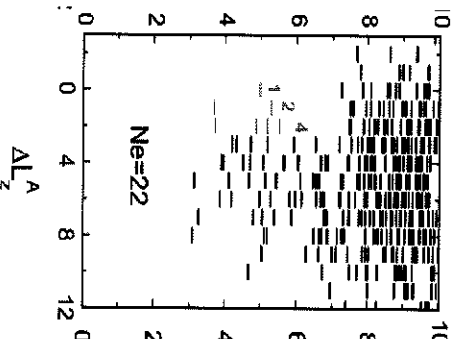
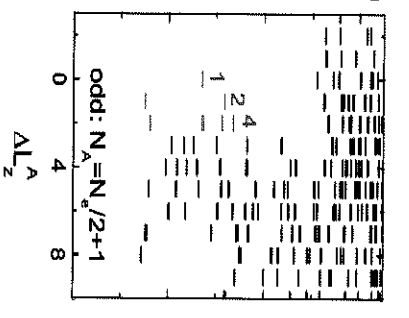
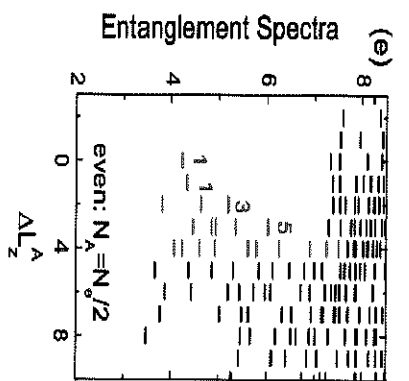
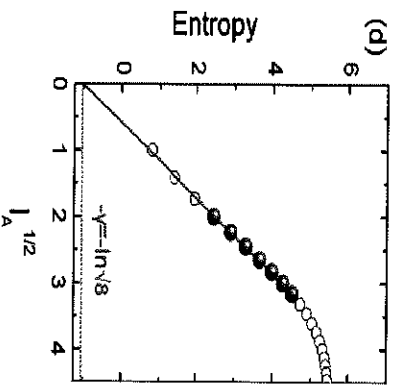
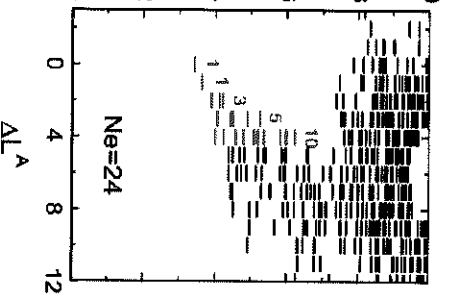
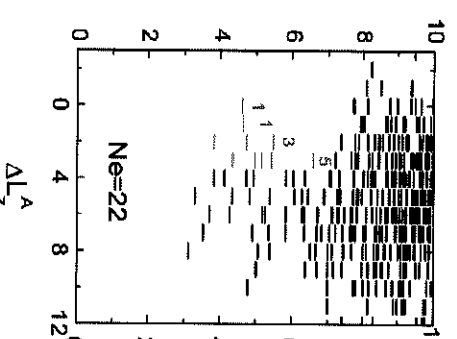
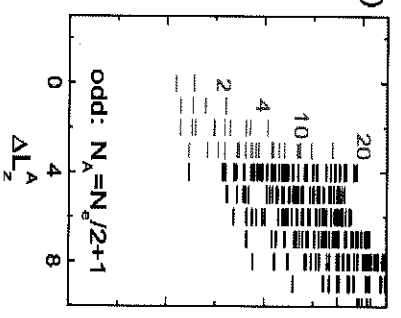
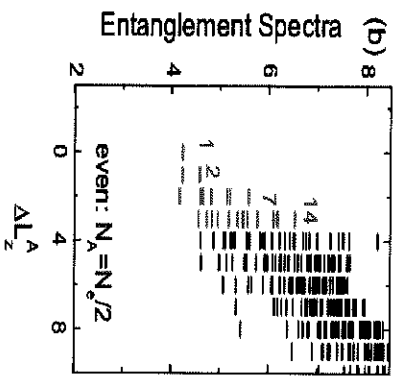
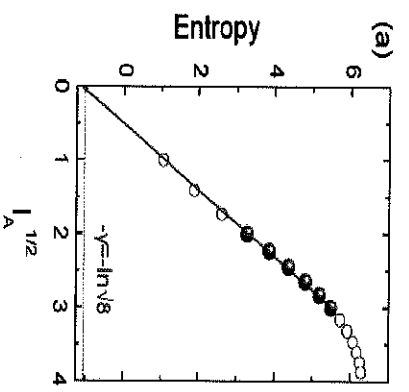
1,1,3,6 ... and 1,3,6, (13)

Infinite Cylinder $L_y=24$



Entanglement spectrum and entropy, Moore-Read vs. Halperin 331

Halperin 331 state



Ne=22

Moore-Read state

Better spectrum for largest system (Ne=24)