

# Model wavefunctions for Chiral Topological Order Interfaces

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**Topological phases of matter [TOPMAT]:  
from the quantum Hall effect to spin liquids**

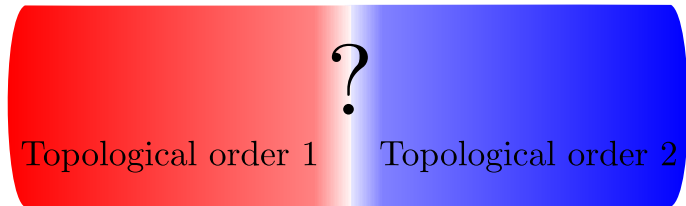
# Acknowledgements

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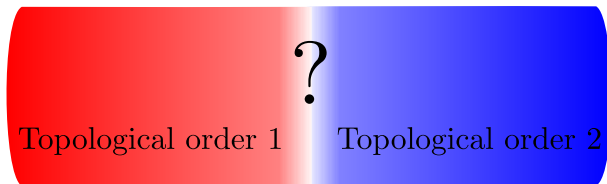
**V. Crépel et al., arXiv :1806.06858**  
**V. Crépel et al., PRB 97, 165136 (2018)**

# Motivations

What's going on at the interface between two topologically ordered phases?



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PRL **103**, 076803 (2009)

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## Non-Abelian Anyons: When Ising Meets Fibonacci

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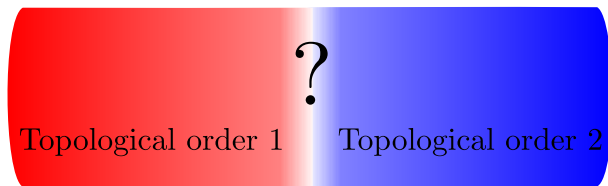
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(Received 20 October 2008; published 13 August 2009)

We consider an interface between two non-Abelian quantum Hall states: the Moore-Read state, supporting Ising anyons, and the  $k = 2$  non-Abelian spin-singlet state, supporting Fibonacci anyons. It is shown that the interface supports neutral excitations described by a  $(1 + 1)$ -dimensional conformal field theory with a central charge  $c = 7/10$ . We discuss effects of the mismatch of the quantum statistical properties of the quasiholes between the two sides, as reflected by the interface theory.

# Motivations

What's going on at the interface between two topologically ordered phases?



**Two questions we want to address:**

- Can you build accurate model wavefunctions for the full system (bulk+interface)?
- Can we characterize the interface down to the microscopic level?

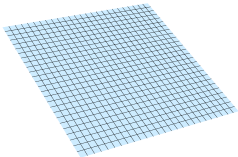
# Outline

- Fractional Quantum Hall (FQH) model wavefunctions
- Matrix Product States (MPS) for the FQH model wavefunctions
- Building a model state for the Laughlin/Halperin interface
- Characterizing the interface

FQH (Abelian) model states

# A single electron in 2D and in a $\perp$ magnetic field $B$ .

**Uniform  $\perp$  magnetic field** : gauge choice



$$H = \frac{1}{2m} \left( \vec{p} - e\vec{A} \right)^2, \quad \vec{A} = \frac{B}{2} \begin{pmatrix} -y \\ x \end{pmatrix}$$

$$H = \frac{1}{2m} \left( -i\hbar \frac{\partial}{\partial x} + \frac{eB}{2} y \right)^2 + \frac{1}{2m} \left( -i\hbar \frac{\partial}{\partial y} - \frac{eB}{2} x \right)^2$$

- **energy scale** cyclotron frequency  $\omega_c = \frac{|eB|}{m}$ ,
- **length scale** : magnetic length  $l_B = \sqrt{\frac{\hbar}{|eB|}}$

$$H = \frac{1}{2} \hbar \omega_c \left[ \left( -i l_B \frac{\partial}{\partial x} + \frac{y}{2 l_B} \right)^2 + \left( -i l_B \frac{\partial}{\partial y} - \frac{x}{2 l_B} \right)^2 \right]$$



# Landau levels

In (dimensionless) complex coordinate  $z = (x + iy)/l_B$ , and setting

$$a = \sqrt{2} \left( \frac{\partial}{\partial \bar{z}} + \frac{z}{2} \right), \quad a^\dagger = -\sqrt{2} \left( \frac{\partial}{\partial z} - \frac{\bar{z}}{2} \right)$$

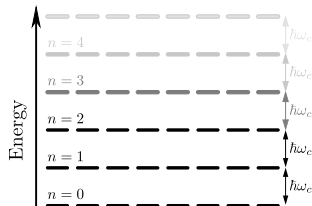
## Familiar form of the Hamiltonian

$$H = \hbar\omega_c \left( a^\dagger a + \frac{1}{2} \right) \quad [a, a^\dagger] = 1$$

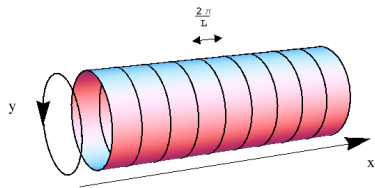
$(n+1)^{\text{th}}$  Landau level :

$$E_n = \hbar\omega_c \left( n + \frac{1}{2} \right)$$

**Discrete** spectrum, large **degeneracy**  
(translation invariance/guiding center).



# Cylinder with perimeter $L$ (we identify $y \equiv y + L$ )



Natural gauge choice :  $\vec{A} = B \begin{pmatrix} 0 \\ x \end{pmatrix}$

$$t_y |\Psi_{k_y}\rangle = k_y |\Psi_{k_y}\rangle, \quad k_y = \frac{2\pi n}{L}$$

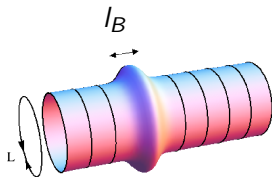
LLL

$$\Psi_{k_y}(x, y) = e^{iyk_y} e^{-\frac{(x - l_B^2 k_y)^2}{2l_B^2}}$$

Momentum  $k_y$  and position  $x$  are locked :

$$x \sim l_B^2 k_y$$

- $[\hat{x}, \hat{y}] = il_B^2$  implies that  $\hbar \hat{x} = l_B^2 \hat{p}_y$ .
- localized in  $\hat{x}$  and delocalized in  $\hat{y}$
- the interorbital distance is  $\frac{2\pi}{L} l_B^2$



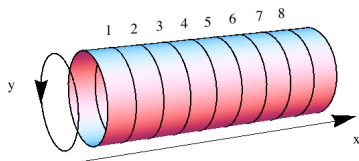
Density profile of the LLL orbital  $\Psi_{k_y}(x, y)$ .

# Projection to the LLL : dimensional reduction

Projection to the LLL :  $x$  and  $y$  no longer commute  $[\hat{x}, \hat{y}] = i l_B^2$ .

**4 dimensional phase space  $\Rightarrow$  2 dimensional phase space**

A **basis** of LLL states



looks like a one-dimensional chain

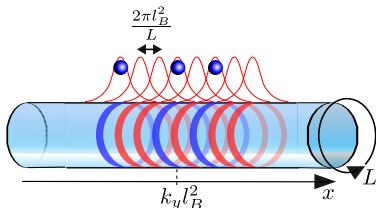
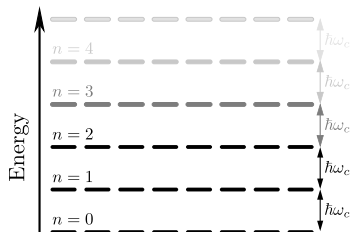


**But !**

Physical short range interactions become long range in this description  
(distance of order  $l_B$  means  $\sim L/l_B$  sites).

# Fractional Quantum Hall effect

## Landau levels (spinless case)



- Partial filling + interaction  $\rightarrow$  FQHE
- Filling factor :  $\nu = \frac{hn}{eB} = \frac{N}{N_\phi}$
- $N$ -body wave function :

$$\Psi = P(z_1, \dots, z_N) e^{-\sum |z_i|^2 / (4l_B^2)}$$

where  $P$  is a polynomial.

- Equivalently in occupation basis

$$|\Psi\rangle = \sum_{\{m_i\}} c_{\{m_i\}} |m_1, \dots, m_{N_{\text{orb}}}\rangle$$

- How to guess  $|\Psi\rangle$  ? ED, DMRG, model wavefunctions, ....

# The mother of all model wavefunctions

The  $\nu = 1/3$  Laughlin state.

**filling fraction  $\nu = 1/3$  + short range model interaction**  
 $\Rightarrow$  **exact ground-state :**

$$\Psi_{\frac{1}{3}}(z_1, \dots, z_N) = \prod_{i < j} (z_i - z_j)^3 e^{-\sum_i |z_i|^2 / 4l_B^2}$$

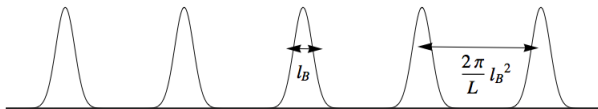
The model interaction is the short range part of Coulomb.

**Extremely high overlap with Coulomb interaction !**  
**(obtained by exact diagonalization)**

First hints of a topological phase :

- excitations with fractional charge  $e/3$
- topology dependent ground state degeneracy :  $3^g$  exact ground states.

## Cartoon picture : thin cylinder limit ( $L \ll l_B$ )



Very small cylinder perimeter  $L$  : **LLL orbitals no longer overlap**  
1d problem

Laughlin's Hamiltonian  $\rightarrow$  Haldane's exclusion statistics  
**no more than 1 particle in three orbitals**

At filling fraction  $\nu = 1/3$ , we get three possible states

$$|\Psi_1\rangle = |\cdots 100100100\cdots\rangle$$

$$|\Psi_2\rangle = |\cdots 010010010\cdots\rangle$$

$$|\Psi_3\rangle = |\cdots 001001001\cdots\rangle$$

3-fold degenerate ground state on the cylinder (and torus).

# Metallic boundary : massless edge modes

$$\psi_u = P_u(z_1, \dots, z_N) \prod_{i < j} (z_i - z_j)^3$$

where  $P_u$  is any symmetric, homogeneous polynomial.

Cartoon picture: no more than 1 electron in 3 orbitals.

- dispersion relation :

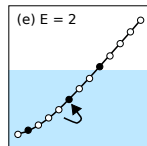
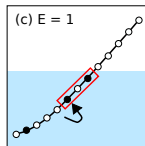
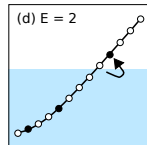
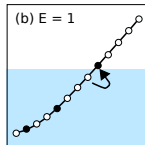
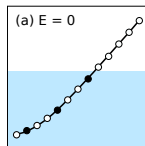
$$E = v_F \Delta P = v_F \frac{2\pi}{L} \Delta N$$

**chiral** and **gapless** edge

- Number of edge states :

- $E = 0$  : 1 state
- $E = 1$  : 1 state
- $E = 2$  : 2 states
- $E = 3$  : 3 states
- $E = 4$  : 5 states
- $E = 5$  : 7 states
- ...

**spectrum of a compact chiral boson ( $R = \sqrt{3}$ ).**



(cartoon picture)

# Metallic boundary : massless edge modes

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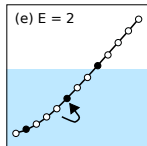
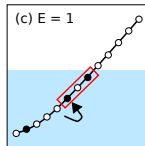
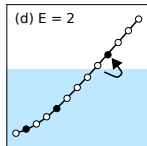
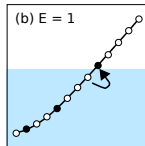
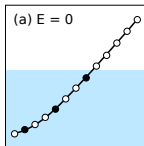
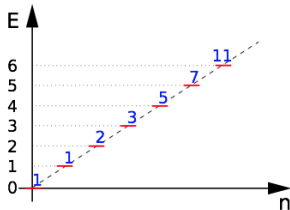
Cartoon picture: no more than 1 electron in 3 orbitals.

- dispersion relation :

$$E = v_F \Delta P = v_F \frac{2\pi}{L} \Delta N$$

**chiral** and **gapless** edge

- Number of edge states :



(cartoon picture)

**spectrum of a compact chiral boson ( $R = \sqrt{3}$ ).**



# Model states and CFT

- **Moore and Read** : A large set of model wavefunctions can be written as a CFT correlator (Laughlin, Moore-Read, Read-Rezayi, Halperin...).

$$\Psi(z_1, \dots, z_N) = \langle \mathcal{O}_{\text{bg}} V(z_1) \cdots V(z_N) \rangle$$

with  $V(z)$  an operator/field in a chiral  $1+1$  CFT and  $\mathcal{O}_{\text{bg}}$  is the background charge.

- **Bulk-edge correspondence** : The CFT used to describe the (gapped) bulk is identical to the CFT that describes the (gapless) edge
- **Laughlin state** :
  - $V(z) =: \exp(i\sqrt{m}\Phi(z))$  :, where  $\Phi(z)$  is a (compact) chiral boson
  - $\langle \Phi(z_1)\Phi(z_2) \rangle = -\log(z_1 - z_2)$
  - $\langle V(z_1) \cdots V(z_N) \rangle = \prod_{i < j} (z_i - z_j)^m$
- **Halperin state** : A two-component (compact chiral) boson.

MPS for the FQH model states

# Entanglement entropy

Cut the system in two parts **A** and **B**  
(the boundary has length  $L$ )

The **entanglement entropy** is

$$S_A = -\text{Tr}(\rho_A \log \rho_A)$$

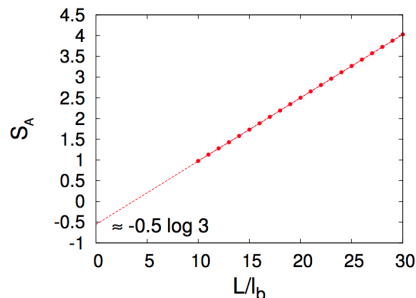
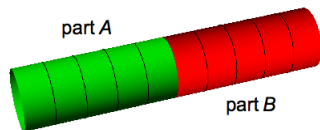
with  $\rho_A$  the reduced density matrix.

For a topological phase :

$$S_A \sim \alpha L - \gamma, \quad \gamma = \log \mathcal{D}$$

where  $\mathcal{D}$  is the quantum dimension.

For  $\nu = 1/3$  Laughlin :  $\mathcal{D} = \sqrt{3}$



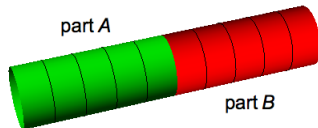
Entanglement entropy of the  $\nu = 1/3$  Laughlin state  
as a function of the cylinder perimeter  $L$   
(N. Regnault)

# Entanglement spectrum

Schmidt decomposition

$$|\Psi\rangle = \sum_{\alpha} \exp(-\xi_{\alpha}/2) |A, \alpha\rangle \otimes |B, \alpha\rangle$$

$$\rho_A = \sum_{\alpha} \exp(-\xi_{\alpha}) |A, \alpha\rangle \langle A, \alpha|$$



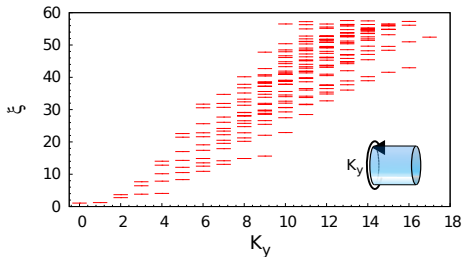
## Entanglement spectrum

Li and Haldane (2008):  
spectrum of  $\xi = -\log \rho_A$   
(plot  $\xi$  vs momentum)

⇒ Reproduces the physical  
edge spectrum !



OES Laughlin  $N=12$ ,  $N_A=6$  on a cylinder  $L=15$



# Matrix Product States

Any state can be written as

$$|\Psi\rangle = \sum_{\{m_i\}} \langle \alpha_L | A^{[m_1]} \dots A^{[m_{N_{\text{orb}}}] } | \alpha_R \rangle | m_1, \dots, m_{N_{\text{orb}}} \rangle$$

- $\{A^{[m]}\}$  is a set of  $\chi \times \chi$  matrices
- $(\alpha_L, \alpha_R)$  encode the boundary conditions for an open system.

The  $A_{\alpha,\beta}^{[m]}$  matrices have two types of indices

- $[m]$  is the physical index ( $m \in \{0, 1\}$  for fermions,  $m \in \mathbb{N}$  for bosons,  $m \in \{\uparrow, \downarrow\}$  for spins ...)
- $(\alpha, \beta)$  are the bond indices (auxiliary space), ranging from  $1, \dots, \chi$ .
- The **bond dimension**  $\chi$  is of the order of  $\exp S_A$   
 $\Rightarrow$  for 2d gapped phases, it grows exponentially with  $L$ .

An exponential improvement over the  $\exp(\text{surface})$  of ED...

## Starting from a model wavefunction given by a CFT correlator

$$\Psi(z_1, \dots, z_N) = \langle u | \mathcal{O}_{\text{b.c.}} V(z_1) \cdots V(z_N) | v \rangle$$

and expanding  $V(z) = \sum_n V_{-n} z^n$ , one finds (up to orbital normalization)

$$c_{(m_1, \dots, m_n)} = \langle u | \mathcal{O}_{\text{b.c.}} \frac{1}{\sqrt{m_n!}} V_{-n}^{m_n} \cdots \frac{1}{\sqrt{m_2!}} V_{-2}^{m_2} \frac{1}{\sqrt{m_1!}} V_{-1}^{m_1} | v \rangle$$

## This is a site/orbital dependent MPS

$$c_{(m_1, \dots, m_n)} = \langle u | \mathcal{O}_{\text{b.c.}} B^{[m_n]}(n) \cdots B^{[m_2]}(2) B^{[m_1]}(1) | v \rangle$$

with matrices at site/orbital  $j$  (including orbital normalization)

$$B^{[m]}(j) = \frac{e^{\left(\frac{2\pi}{L} j\right)^2}}{\sqrt{m!}} (V_{-j})^m$$

# Translation invariant MPS

A relation of the form  $B^{[m]}(j) = U^{-1} B^{[m]}(j-1) U$  yields

$$B^{[m]}(\textcolor{red}{j}) = U^{-\textcolor{red}{j}} B^{[m]}(0) U^{\textcolor{red}{j}}$$

and then

$$B^{[m_n]}(\textcolor{red}{n}) \cdots B^{[m_1]}(\textcolor{red}{1}) = U^{-n} \times B^{[m_n]}(\textcolor{red}{0}) U \cdots B^{[m_1]}(\textcolor{red}{0}) U$$

This is a **translation invariant MPS**, with matrices

$$\textcolor{red}{A}^{[m]} = \textcolor{red}{B}^{[m]}(0) U$$

# Translation invariant MPS on the cylinder

## Site independant MPS

$$B^{[m]}(\textcolor{red}{j}) = \frac{e^{(\frac{2\pi}{L}\textcolor{red}{j})^2}}{\sqrt{m!}} (V_{-j})^m \quad \Rightarrow \quad A^{[m]} = \frac{1}{\sqrt{m!}} (V_0)^m U$$

where  $U$  is the operator is (Zaletel and Mong (2012))

$$\textcolor{red}{U} = e^{-\frac{2\pi}{L}H - i\sqrt{\nu}\varphi_0}$$

where

- $\varphi_0$  is the bosonic zero mode ( $e^{-i\sqrt{\nu}\varphi_0}$  shifts the electric charge by  $\nu$ )
- $H$  is the CFT cylinder Hamiltonian :  $H = \frac{2\pi}{L}L_0$
- $V_0$  is the zero mode of  $V(z)$

**auxiliary space = CFT Hilbert space**

infinite bond dimension :/

Extension to spinfull FQH : **V. Crépel et al., PRB 97, 165136 (2018)**



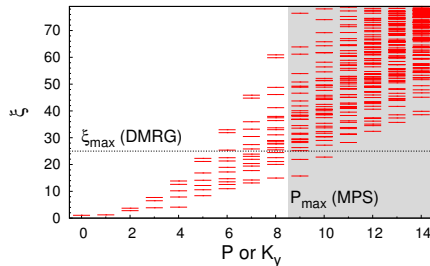
# Truncation of the auxiliary CFT basis

$\chi$  is infinite  $\rightarrow$  a truncation scheme is required.

- The natural cut-off is the total conformal dimension  $\rightarrow P_{\max}$ .
- Truncation over the momentum in the OES.
- In finite size, the truncated MPS becomes exact for  $P_{\max}$  large enough.

- DMRG : cut-off in  $\xi$  (remove the smallest weight of  $\rho_A$ ).
- MPS : cut-off in momentum.

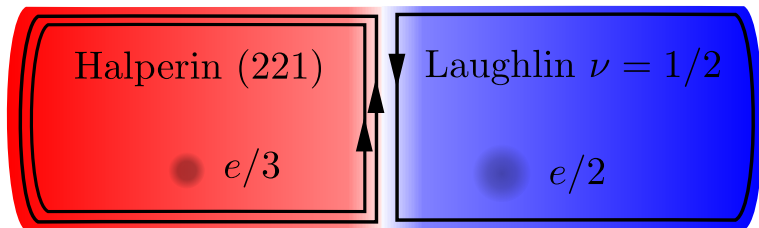
Morally equivalent as long as the ES mimics the chiral edge mode spectrum (linear dispersion).



Building a model state for the Halperin/Laughlin interface

# Let's look at the following interface

We consider the (bosonic) case interfacing the Halperin (221) and the Laughlin  $\nu = 1/2$  phases.



## Halperin (221)

- Spinful,  $SU(2)$  symmetric,  $\nu_{\uparrow} = \nu_{\downarrow} = 1/3$ .
- $e/3$  excitations.
- Two  $U(1)$  chiral edge modes (charge and spin).

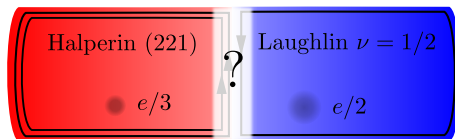
## Laughlin $\nu = 1/2$

- Spinless,  $\nu_{\uparrow} = 0, \nu_{\downarrow} = 1/2$ .
- $e/2$  excitations.
- One  $U(1)$  chiral edge mode (charge).

# A microscopic model

$$\mathcal{H}_{\text{int}} = \int d^2\vec{r} \left( \sum_{\sigma, \sigma'=\uparrow, \downarrow} : \rho_{\sigma}(\vec{r}) \rho_{\sigma'}(\vec{r}) : \right) + \mu_{\uparrow}(\vec{r}) \rho_{\uparrow}(\vec{r})$$

- Use the chemical potential  $\mu_{\uparrow}(\vec{r})$  to polarize half of the system.
- Laughlin  $\nu = 1/2$  is the densest **polarized** zero energy state.
- Halperin (221) is the densest **unpolarized** zero energy state.
- The two quantum liquids are sewed together by the interaction.



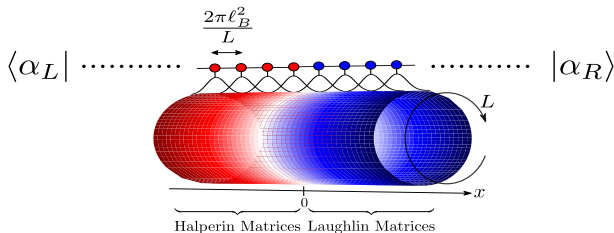
What shall we observe at the interface ? **A single gapless mode described by a free chiral boson (Haldane, PRL 94).**

# MPS and variational Ansatz

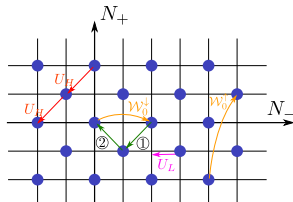
- We know the exact MPS for Halperin  $B^{[n]}$  and Laughlin  $A^{[n]}$ .



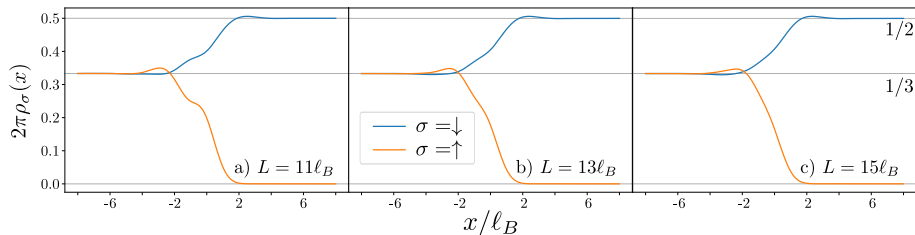
- Brutal gluing :  $\langle \alpha_L | \dots B^{[m-2]} B^{[m-1]} A^{[m_0]} A^{[m_1]} \dots | \alpha_R \rangle$



- Does  $B^{[m-1]} A^{[m_0]}$  make any sense?
- Yes : conformal embedding !**
- A careful choice for the electron, bg charge operators.

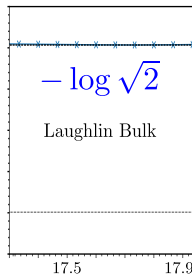
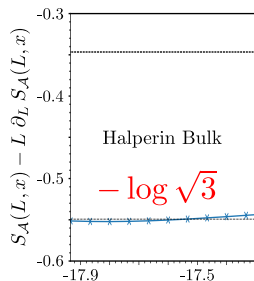


# Density



- Translation invariance along the cylinder perimeter.
- We recover the spin up and down densities in the bulk both on the Halperin and Laughlin side.
- Finite size effects (with respect to  $L$ ) quickly vanish.
- Width of interface  $\simeq 5/\ell_B$

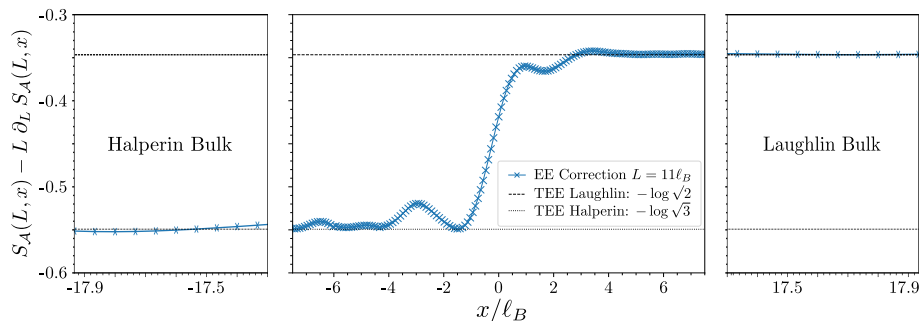
# Topological Entanglement Entropy



$x/\ell_B$

- We extract the TEE  $\gamma$  from the derivative  $S_A - L \partial_L S_A = -\gamma$ .
- Good agreement with the predicted values deep in the bulks (-0.549 and -0.347).

# Topological Entanglement Entropy

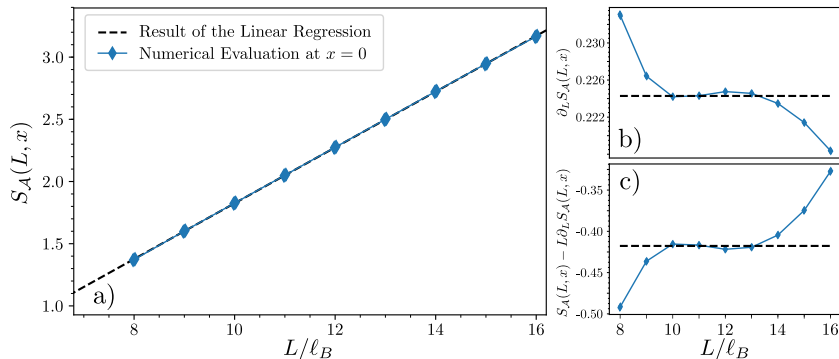


- Up to a small oscillations (finite size effects more important for subleading terms), a rather smooth transition between the two bulk TEE.
- No sign of the gapless mode (as recently predicted by Santos et al. arXiv:1803.04418).



# Area law at the transition

Does we still satisfy the area law at the interface?

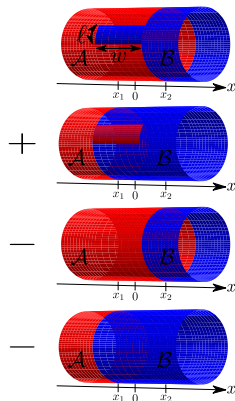


Yes (but hard to spot any deviation with such a limited range).

## Characterizing the interface

# Extracting $c$ : Levin-Wen cut

$$\underbrace{\alpha(x_1)\ell + \alpha(x_2)(L - \ell) + 2 \int_{x_1}^{x_2} \alpha(u) du}_{\text{Area Law}} + \underbrace{\frac{c}{6} \log \left[ \sin \left( \frac{\pi \ell}{L} \right) \right]}_{\text{Critical Mode}} + K(w)$$



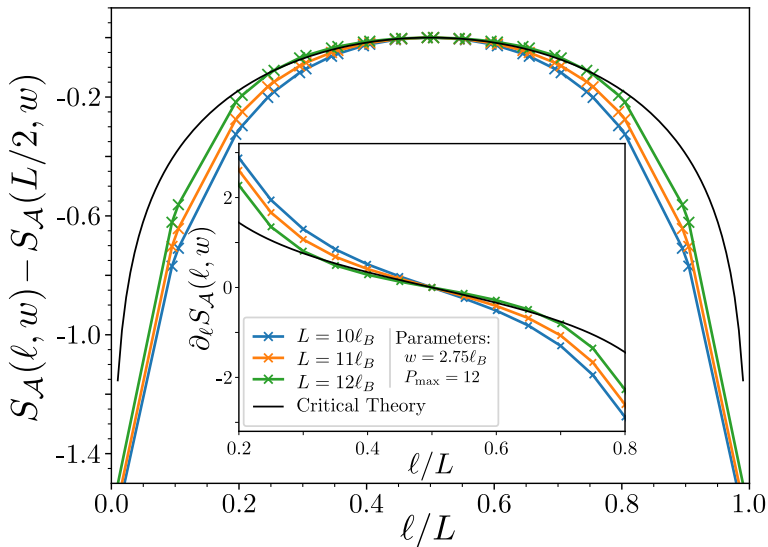
- $K(w)$  contains corrections to the area law, corner contributions,...
- Using a Levin-Wen cut to focus on the critical contribution.

•

$$S_A(\ell, w) = 2 \frac{c}{6} \log \left[ \sin \left( \frac{\pi \ell}{L} \right) \right] + f(w)$$

- To get rid of  $f(w)$  (including the TEE), we compute  $S_A(\ell, w) - S_A(L/2, w)$

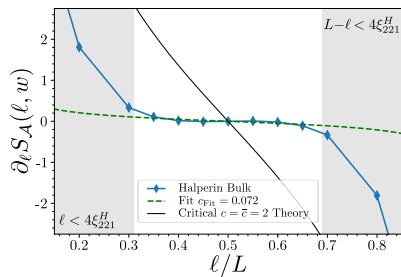
# Extracting $c$ : Levin-Wen cut



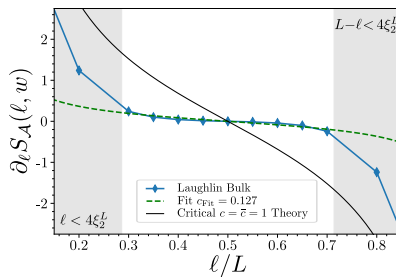
**Fitted central charge  $c = 0.987(1)$ .**

# What about the bulk?

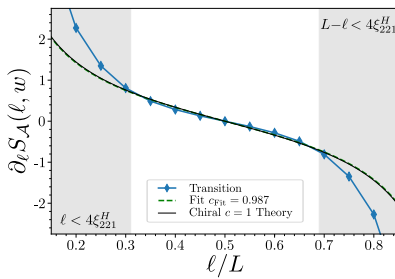
## Halperin



## Laughlin

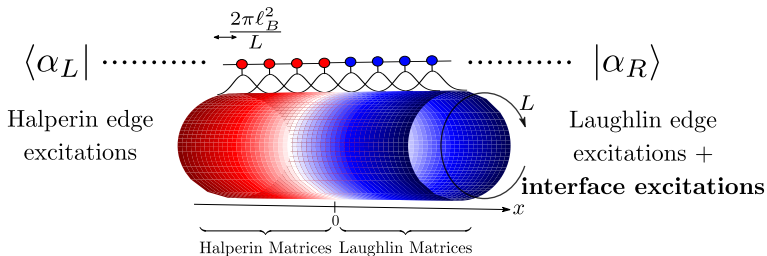


## Interface

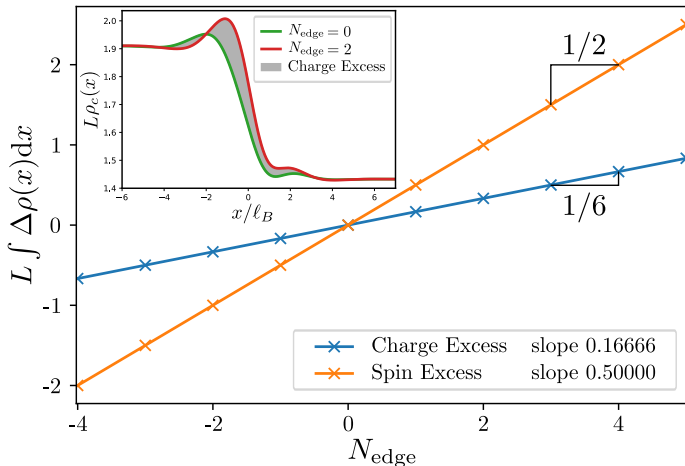


# Compactification radius, fractional charge

- Central charge is only part of the information.
- Mutual information  $\rightarrow$  full partition function of the CFT but hard to evaluate.
- Compactification radius  $\leftrightarrow$  charge of the elementary edge excitation.
- Directly measure the charge along the edge.
- Play with the MPS boundary conditions.



# Compactification radius, fractional charge

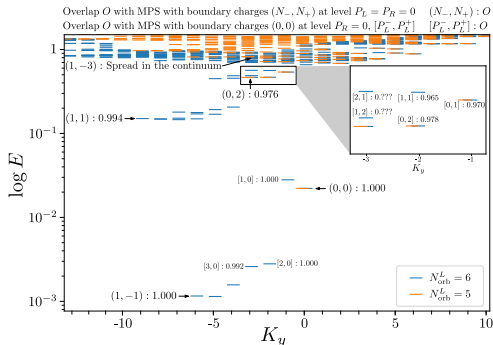


Excitations with a  $e/6$  charge ( $e/6 = e/2 - e/3$ ).

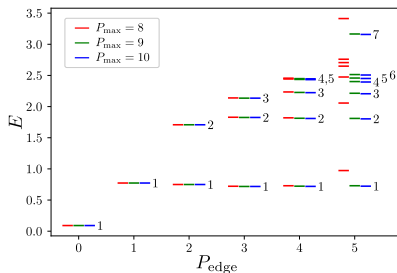
**Edge = compact boson with radius  $R = \sqrt{6}$ .**

# Is it a good variational wavefunction?

- It has all the features that we expect but *does it capture the microscopic model low energy properties?*
- Overlap with ED :  $4 \uparrow + 9 \downarrow$  particles, 21 orbitals  $\rightarrow 0.998$  (Hilbert space dim  $\simeq 2.2 \times 10^8$ ).



ED with 15 orbitals.

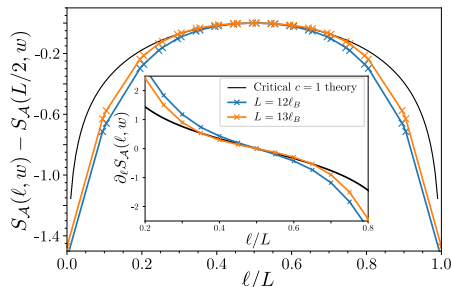
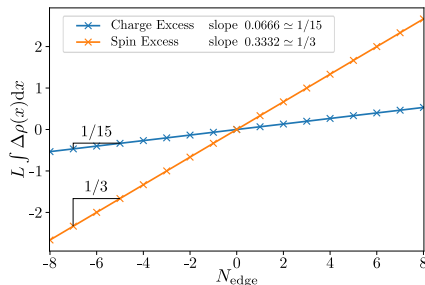


MPO  $L = 12$ .



# Fermions : Laughlin $\nu = 1/3$ / Halperin (332)

- No conceptual difference with the bosonic example.
- Transition from Laughlin  $\nu = 1/3$  to Halperin (332) at  $\nu = 2/5$ .
- Experimental relevance : graphene using the valley degree of freedom (spontaneous polarization at  $\nu = 1/3$ ).



# Conclusion

- A variational ansatz to describe the interface between the Halperin and the Laughlin liquids.
- Microscopic characterization of the interface gapless mode ( $c$  and  $R$ ).
- This scheme can be extended to
  - any case where a MPS/PEPS/TN description is known on both sides.
  - other sewing approach (e.g. superconductor).
  - a more generic approach ?

