

Title: Topological Order Series

Date: Jun 18, 2015 03:00 PM

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Abstract:

- **$n$ -cluster condition:** No off-particle zeros when  $c = n$  (or the wave function for the  $n$ -clusters is the Laughlin wave function or a  $n$ -electron cluster is “kind of trivial”)  
 $\rightarrow D_{a+b,n} = D_{a,n} + D_{b,n} \rightarrow$

$$S_{a+kn} = S_a + kS_n + \frac{k(k-1)nm}{2} + kma$$

Since  $S_1 = 0$ ,  $(m, S_2, \dots, S_n)$  carries all the information about the pattern of zeros from an  $n$ -cluster symmetric polynomial.

- Additional conditions

$$\Delta_2(a, a) = \text{even}, \quad m > 0, \quad mn = \text{even}, \quad 2S_n = 0 \bmod n.$$

- **A mysterious condition** (the one we want but cannot prove):

$$\Delta_3(a, b, c) = \text{even}$$

## Local dancing rule → global dancing pattern

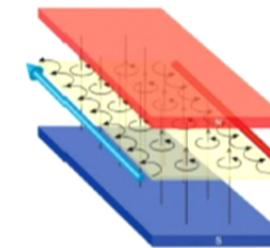
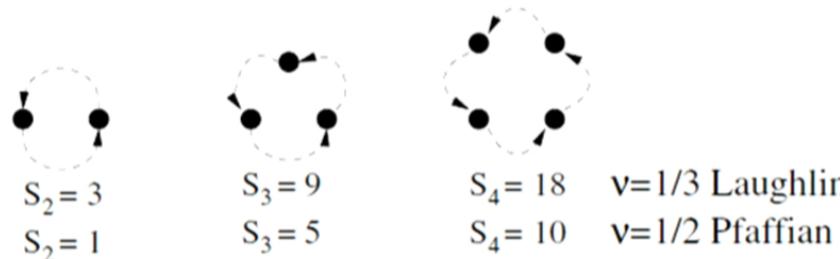
- Local dancing rules of a FQH liquid:
  - (1) every electron dances around clock-wise  
( $\Phi_{\text{FQH}}$  only depends on  $z = x + iy$ )
  - (2) takes exactly three steps to go around each other



(Relative angular momentum  $S_2 = 3$ )    Wen-Wang arXiv:0803.1016

$$\rightarrow \text{Global dancing pattern } \Phi_{\text{FQH}}(\{z_1, \dots, z_N\}) = \prod(z_i - z_j)^3$$

- A systematic theory of FQH state – **Pattern of zeros  $S_a$** :  
 $a$ -electron cluster has a relative angular momentum  $S_a$



- Local dancing rules are enforced by the Hamiltonian to lower energy.
- Only certain sequences  $S_a$  correspond to valid FQH states. Which?
- Different POZ  $S_a$  give rise to different topological properties

## Conditions on pattern of zeros – ground state

- **Concave conditions**

$$\Delta_2(a, b) \equiv S_{a+b} - S_a - S_b = D_{a,b} \geq 0,$$

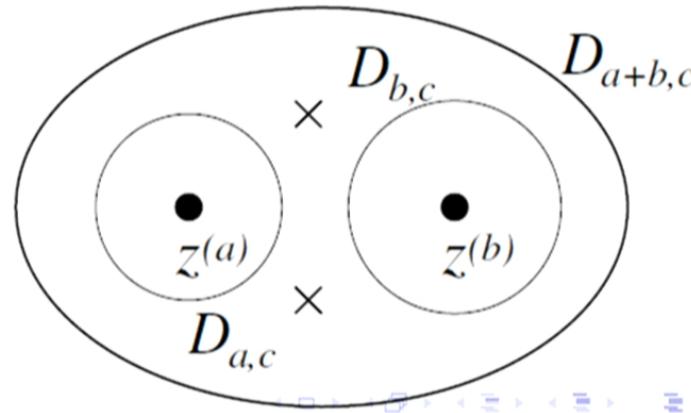
$$\Delta_3(a, b, c) \equiv S_{a+b+c} - S_{a+b} - S_{b+c} - S_{a+c} + S_a + S_b + S_c \geq 0$$

The second one can be shown by considering

$P_{\text{derived}}(z^{(a)}, z^{(b)}, z^{(c)}, \dots)$  as a function of  $z^{(c)}$ , which leads to

$$D_{a+b,c} = D_{a,c} + D_{b,c} + \text{off-particle zeros at "x"} \geq D_{a,c} + D_{b,c}$$

- $D_{a,c}, D_{b,c}$  = the number of on-particle zeros at  $z^{(a)}, z^{(b)}$ .
- As  $z^{(a)} \rightarrow z^{(b)}$ , some off-particle zeros at “x” may also approach  $z^{(b)}$ .
- $D_{a+b,c} \geq D_{a,c} + D_{b,c}$  leads to the second condition.



## Primitive solutions for pattern of zeros

The conditions are semi-linear →

if  $(m; S_2, \dots, S_n)$  and  $(m'; S'_2, \dots, S'_n)$  are solutions, then

$(m''; S''_2, \dots, S''_n) = (m; S_2, \dots, S_n) + (m'; S'_2, \dots, S'_n)$  is also a solution  $\sim P''(\{z_i\}) = P(\{z_i\})P'(\{z_i\})$

1-cluster state:  $\nu = 1/m$  Laughlin state

$$P_{1/m} : \quad \mathbf{S} = (m; ), \\ (n_0, \dots, n_{m-1}) = (1, 0, \dots, 0).$$

2-cluster state: Pfaffian state ( $Z_2$  parafermion state)

$$P_{\frac{2}{2}; Z_2} : \quad (m; S_2) = (2; 0), \\ (n_0, \dots, n_{m-1}) = (2, 0)$$

3-cluster state:  $Z_3$  parafermion state

$$P_{\frac{3}{2}; Z_3} : \quad (m; S_2, S_3) = (2; 0, 0), \\ (n_0, \dots, n_{m-1}) = (3, 0)$$



4-cluster state:  $Z_4$  parafermion state

$$P_{\frac{4}{2};Z_4} : (m; S_2, \dots, S_n) = (2; 0, 0, 0), \\ (n_0, \dots, n_{m-1}) = (4, 0),$$

5-cluster states:  $Z_5$  (generalized) parafermion state

$$P_{\frac{5}{2};Z_5} : (m; S_2, \dots, S_n) = (2; 0, 0, 0, 0), \\ (n_0, \dots, n_{m-1}) = (5, 0)$$

$$P_{\frac{5}{8};Z_5^{(2)}} : (m; S_2, \dots, S_n) = (8; 0, 2, 6, 10), \\ (n_0, \dots, n_{m-1}) = (2, 0, 1, 0, 2, 0, 0, 0)$$

6-cluster state:

$$P_{\frac{6}{2};Z_6} : (m; S_2, \dots, S_n) = (2; 0, 0, 0, 0, 0), \\ (n_0, \dots, n_{m-1}) = (6, 0)$$



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7-cluster states:

$$P_{\frac{7}{2};Z_7} : (m; S_2, \dots, S_n) = (2; 0, 0, 0, 0, 0, 0, 0), \\ (n_0, \dots, n_{m-1}) = (7, 0)$$

$$P_{\frac{7}{8};Z_7^{(2)}} : (m; S_2, \dots, S_n) = (8; 0, 0, 2, 6, 10, 14), \\ (n_0, \dots, n_{m-1}) = (3, 0, 1, 0, 3, 0, 0, 0)$$

$$P_{\frac{7}{18};Z_7^{(3)}} : (m; S_2, \dots, S_n) = (18; 0, 4, 10, 18, 30, 42), \\ (n_0, \dots, n_{m-1}) = (2, 0, 0, 0, 0, 1, 0, 0, 0, 2, 0, 0, 0, 0, 0, 0)$$

$$P_{\frac{7}{14};C_7} : (m; S_2, \dots, S_n) = (14; 0, 2, 6, 12, 20, 28), \\ (n_0, \dots, n_{m-1}) = (2, 0, 1, 0, 1, 0, 1, 0, 2, 0, 0, 0, 0, 0, 0)$$

- Also get composite parafermion state  $P = P_{Z_{n_1}} P_{Z_{n_2}}$

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## How good is the pattern-of-zero classification

Not so bad :-) Not so good :-(

- Every symm. poly.  $P$  corresponds to a unique pattern of zero  $\{S_a\}$ .  
Only some patterns of zero correspond to a unique symm. poly.
  - It appears that all the primitive patterns of zero correspond to a unique symm. poly.
  - It is known that
    - some composite patterns of zero  $\rightarrow$  a unique symm. polynomial
    - some composite patterns of zero  $\rightarrow$  several symm. polynomials
- So in general, we need more information than  $\{n; m; S_a\}$  to fully characterize symmetry polynomial of infinite variables**



## Topological properties from pattern of zeros

For those patterns of zeros that uniquely characterize a FQH wave function, we should be able to calculate the topological properties of FQH states from the data  $(n; m; S_2, \dots, S_n)$ .

Physical properties that we want to get

- The filling fraction (actually given by  $\nu = n/m$ ).
- Topological degeneracy on torus (and other Riemann surface)
- Number of quasiparticle types
- Quasiparticle charges
- Quasiparticle scaling dimensions
- Quasiparticle fusion algebra
- Quasiparticle statistics (Abelian and non-Abelian)
- The counting of edge excitations (central charge  $c$  and spectrum)

7-cluster states:

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$$P_{\frac{7}{18};Z_7^{(3)}} : (m; S_2, \dots, S_n) = (18; 0, 4, 10, 18, 30, 42), \\ (n_0, \dots, n_{m-1}) = (2, 0, 0, 0, 0, 1, 0, 0, 0, 2, 0, 0, 0, 0, 0, 0)$$

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## The pattern of zeros for quasiparticles

- A quasiparticle  $\gamma$  in a FQH state can also be quantitatively characterized by pattern of zeros  $\{S_{\gamma;a}\}$ :  

- $P_\gamma(\eta; \{z_i\})$  has a quasiparticle at  $z = \eta$
- Let  $z_i = \lambda\xi_i + \eta$ ,  $i = 1, 2, \dots, a$  (bring  $a$  electrons to the quasiparticle)

$$P_\gamma(\eta; \{z_i\}) = \lambda^{S_{\gamma;a}} P_\gamma(z^{(a)} = \eta, z_{a+1}, z_{a+2}, \dots) + O(\lambda^{S_a+1})$$

$S_{\gamma;a}$  is the order of zero of  $\Psi_\gamma(\xi, z_i)$  when we bring  $a$  electrons to  $\xi$ .

## Quasi-holes in the $\nu = 1/m$ Laughlin state and fractional charge

- A hole-like excitation = missing an electron, charge = 1

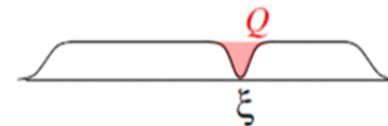
$$\prod_i (\xi - z_i)^m \prod_{i < j} (z_i - z_j)^m e^{-\sum |z_i|^2/4}$$

which can be splitted into  $m$  quasi-hole excitations:

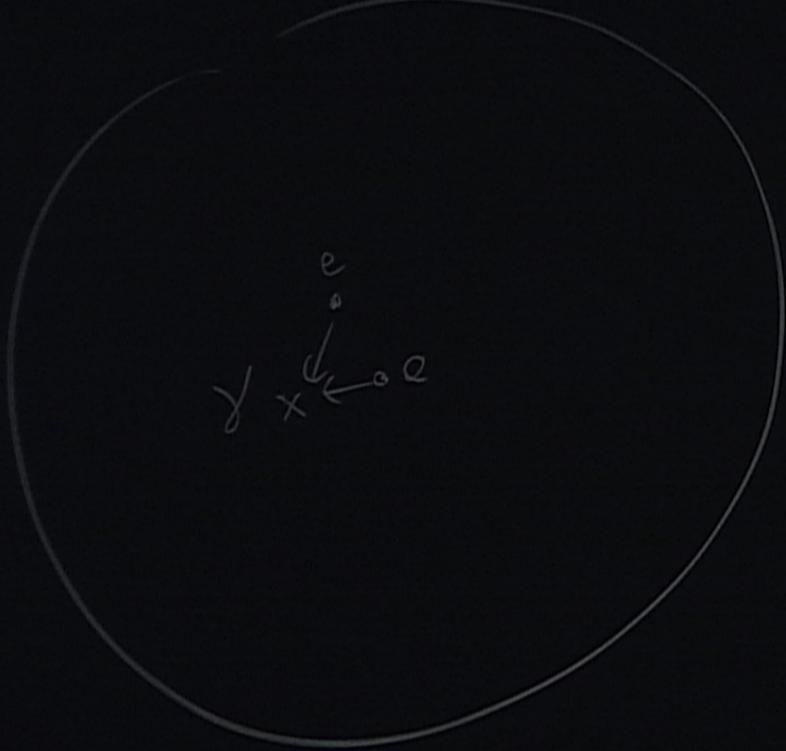
$$\prod_i (\xi_1 - z_i) \cdots \prod_i (\xi_m - z_i) \prod_{i < j} (z_i - z_j)^m e^{-\sum |z_i|^2/4}$$

- A quasi-hole excitation = minimal excitation, charge =  $1/m$

$$\prod_i (\xi - z_i) \prod_{i < j} (z_i - z_j)^m e^{-\sum |z_i|^2/4}$$



- Why the density dip have a small finite size?



$$S_1 = 0 \quad S_2 =$$

$$S_3 = \dots$$

$S_n \rightarrow$  trivial

$$S_{\gamma,1} = 1$$

$$S_{\gamma,2} = m+2$$

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$$P_\gamma(\eta; \{z_i\}) = \lambda^{S_{\gamma;a}} P_\gamma(z^{(a)} = \eta, z_{a+1}, z_{a+2}, \dots) + O(\lambda^{S_a+1})$$

$S_{\gamma;a}$  is the order of zero of  $\Psi_\gamma(\xi, z_i)$  when we bring  $a$  electrons to  $\xi$ .

**The sequence of integers  $\{S_{\gamma;a}\}$  characterizes the quasiparticle  $\gamma$ .**

- $\{S_a\}$  correspond to the trivial quasiparticle  $\gamma = 0$ :  $\{S_{0;a}\} = \{S_a\}$

## Conditions on $S_{\gamma;a}$

- Concave condition

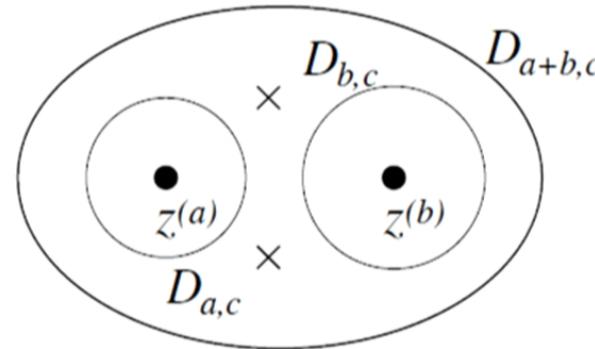
$$S_{\gamma;a+b} - S_{\gamma;a} - S_b \geq 0,$$

$$S_{\gamma;a+b+c} - S_{\gamma;a+b} - S_{\gamma;a+c} - S_{b+c} + S_{\gamma;a} + S_b + S_c \geq 0$$

- Taking  $b = c = 1$ , we obtain

$$S_{\gamma;a+2} - 2S_{\gamma;a+1} + S_{\gamma;a} = I_{\gamma;a+2} - I_{\gamma;a+1} \geq S_2, \quad a \geq 0.$$

$I_{\gamma;a}$  increases with  $a$ :  $I_{\gamma;a+1} - I_{\gamma;a} \geq S_2 \geq 0$ .



$z^{(a)}$  is a quasiparticle  $\gamma$  fused with  $a$  electrons.

For the  $\nu = 1$  Pfaffian state ( $n = 2$  and  $m = 2$ )

$$S_1, S_2, \dots : 0, 0, 2, 4, 8, 12, 18, 24, \dots$$

$$n_0 n_1 n_2 \dots : 2020202020202020202 \dots$$

- Quasiparticle solutions ( $S_{\gamma;a} \rightarrow l_{\gamma;a} \rightarrow n_{\gamma;l}$ ):

$$n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 2020202020202020202 \dots \quad Q_\gamma = 0$$

$$n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 0202020202020202020 \dots \quad Q_\gamma = 1$$

$$n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 11111111111111111111 \dots \quad Q_\gamma = 1/2$$

- All other quasiparticle solutions can be obtained from the above three by removing extra electrons  $\rightarrow$  only 3 quasiparticle types.
- Ground state degeneracy on torus = number of quasiparticle types
- Charge of quasiparticles

$$Q_\gamma = \frac{1}{m} \sum_{a=1}^n (l_{\gamma;a} - l_a)$$

the average increase of  $l_{\gamma;a}$  from the ground state values  $l_a$ .



# Quasi-holes in the $\nu = 1$ Pfaffian state: What is non-Abelian statistics?

- Ground state:  $z_1 \approx z_2$ , no zero;  $z_1 \approx z_2 \approx z_3$ , second-order zero;

$$\Psi_{\text{Pf}} = \mathcal{A} \left( \frac{1}{z_1 - z_2} \frac{1}{z_3 - z_4} \cdots \frac{1}{z_{N-1} - z_N} \right) = \text{Pf} \left( \frac{1}{z_i - z_j} \right)$$

- A charge-1 quasi-hole state

$$\begin{aligned} \Psi_{\text{charge-1}} &= \prod (\xi - z_i) \mathcal{A} \left( \frac{1}{z_1 - z_2} \frac{1}{z_3 - z_4} \cdots \frac{1}{z_{N-1} - z_N} \right) \\ &= \mathcal{A} \left( \frac{(\xi - z_1)(\xi - z_2)}{z_1 - z_2} \frac{(\xi - z_3)(\xi - z_4)}{z_3 - z_4} \cdots \right) = \text{Pf} \left( \frac{(\xi - z_i)(\xi - z_j)}{z_i - z_j} \right) \end{aligned}$$

- A state with two charge-1/2 quasi-holes

$$\begin{aligned} \Psi_{(\xi)(\xi')} &= \mathcal{A} \left( \frac{(\xi - z_1)(\xi' - z_2) + (1 \leftrightarrow 2)}{z_1 - z_2} \frac{(\xi - z_3)(\xi' - z_4) + (3 \leftrightarrow 4)}{z_3 - z_4} \cdots \right) \\ &= \text{Pf} \left( \frac{(\xi - z_i)(\xi' - z_j) + (\xi - z_j)(\xi' - z_i)}{z_i - z_j} \right) \end{aligned}$$

For the  $\nu = 1$  Pfaffian state ( $n = 2$  and  $m = 2$ )

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- Quasiparticle solutions ( $S_{\gamma;a} \rightarrow l_{\gamma;a} \rightarrow n_{\gamma;l}$ ):

$$n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 20202020202020202 \dots \quad Q_\gamma = 0$$

$$n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 02020202020202020 \dots \quad Q_\gamma = 1$$

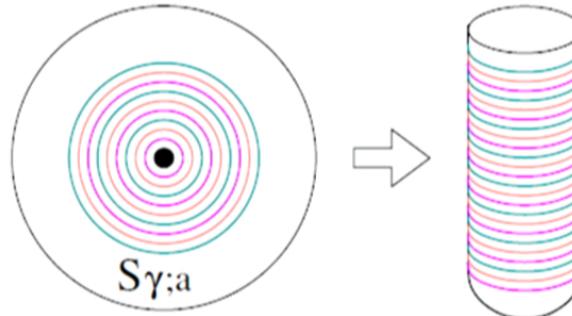
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## Number of quasiparticles types = degeneracy on torus

Consider a quasiparticle  $\gamma$  on the disk at  $z = 0$ , whose type is determined by  $n_{\gamma,I}$  in large  $I$  limit.

- On the disk, the wave function has the POZ  $S_a$  everywhere with  $z \neq 0 \rightarrow$  zero energy.
- Deform the disk into a torus
- On the torus, the wave function has the POZ  $S_a$  everywhere  $\rightarrow$  zero energy.
- The number of the zero-energy patterns (wave functions) is the same as the numbers of quasiparticle types.



For the  $\nu = 1$  Pfaffian state ( $n = 2$  and  $m = 2$ )

$$S_1, S_2, \dots : 0, 0, 2, 4, 8, 12, 18, 24, \dots$$

$$l_1 l_2 l_3 \dots : 0, 0, 2, 2, 4, 4, 6, 6, 8, 8, \dots$$

$$n_0 n_1 n_2 \dots : 2020202020202020202 \dots ,$$

- GPE rule: (1) Three particles occupy  $\geq l_3 + 1 = 3$  orbitals or more ( $c = 1$ ). (2) The spread of two particles plus the spread of next two particles  $\geq (l_2 - l_1) + (l_3 - l_2) = 2$  ( $c = 2$ ). (3) ...
- the quasiparticle occupation patterns

$$n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 2020202020202020202 \dots \quad Q_\gamma = 0$$

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are the only **close-packed** occupation patterns satisfying the above generalized Pauli exclusion rules.

- **close-packed:**

(1) add any electrons  $\rightarrow$  violate the GPE rules.

(2) shift any electrons  $\rightarrow$  violate the GPE rules.



$$\ell=0 \quad 1 \quad 2 \quad 3 \quad 4$$



## Conditions on $S_{\gamma;a}$

- Concave condition

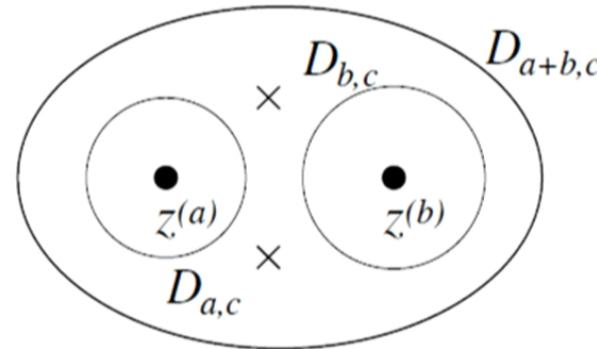
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$z^{(a)}$  is a quasiparticle  $\gamma$  fused with  $a$  electrons.

## Pattern of zeros and generalized Pauli exclusion rule

- $l_{\gamma;a+1} - l_{\gamma;a} \geq S_2$  → the spacing between any two occupied orbitals by different electrons is no less than  $S_2$   
→ a generalized Pauli exclusion rule

### More general Pauli exclusion (GPE) rule

In terms of  $l_{\gamma;a} = S_{\gamma;a} - S_{\gamma;a-1}$ , the concave condition for quasiparticles becomes

$$\sum_{k=1}^b l_{\gamma;a+k} \geq S_b = \sum_{k=1}^b l_{a+k}, \quad \rightarrow \quad l_{\gamma;a} \geq l_a,$$

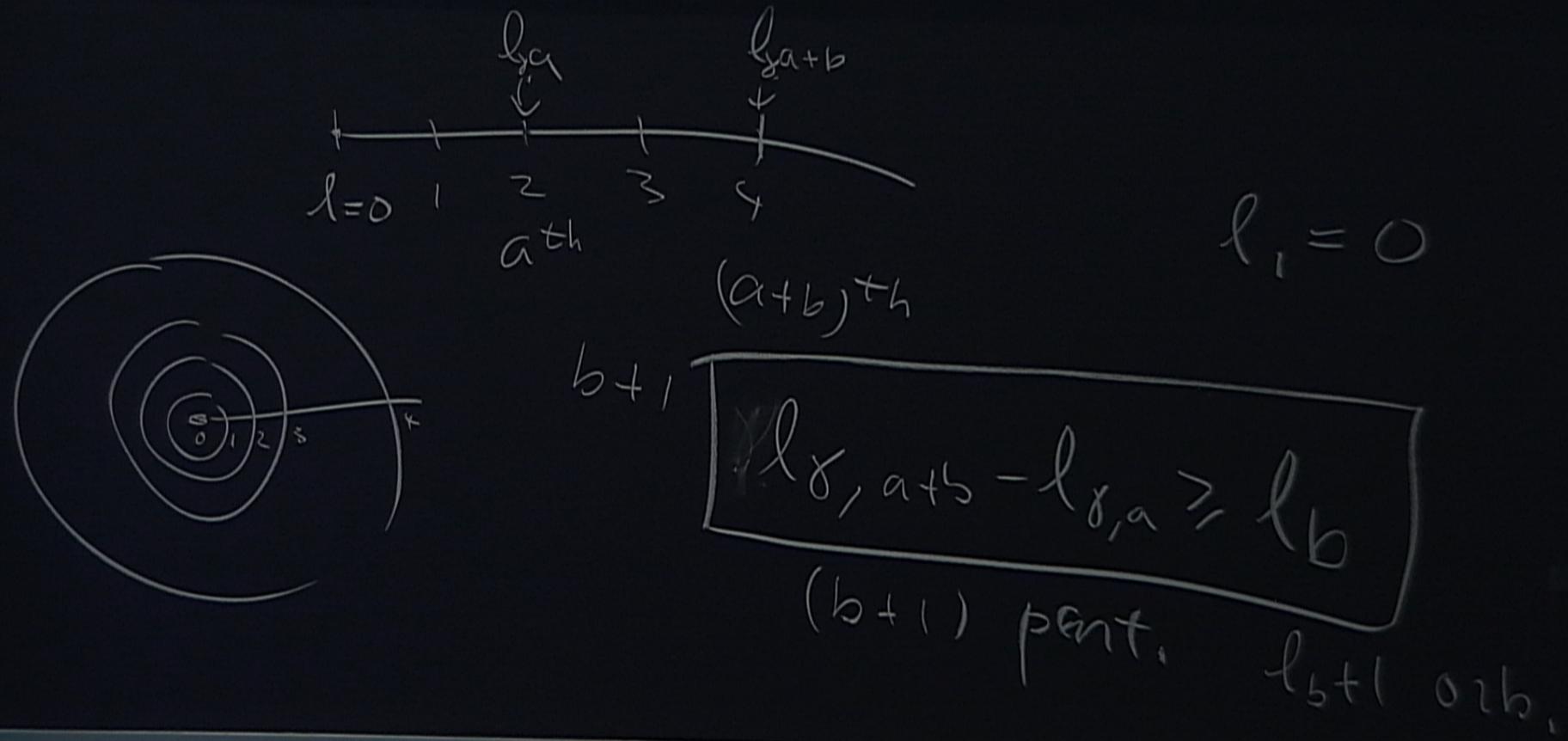
$$\sum_{k=1}^c (l_{\gamma;a+b+k} - l_{\gamma;a+k}) \geq S_{b+c} - S_b - S_c = \sum_{k=1}^c (l_{b+k} - l_k)$$

for any  $a, b, c \in \mathbb{Z}_+$ . Note that  $l_{\gamma;a+b} - l_{\gamma;a}$  is the spread of  $b + 1$  electrons. Setting  $c = 1 \rightarrow$  **The spread of  $b$  electrons  $\geq l_b$ .**

→ **The number of orbitals occupied by  $b$  electrons  $\geq l_b + 1$ .**

( $l_b$  = the spread of the first  $b$  electrons in the ground state.)





For the  $\nu = 1$  Pfaffian state ( $n = 2$  and  $m = 2$ )

$$S_1, S_2, \dots : 0, 0, 2, 4, 8, 12, 18, 24, \dots$$

$$l_1 l_2 l_3 \dots : 0, 0, 2, 2, 4, 4, 6, 6, 8, 8, \dots$$

$$n_0 n_1 n_2 \dots : 2020202020202020202 \dots ,$$

- GPE rule: (1) Three particles occupy  $\geq l_3 + 1 = 3$  orbitals or more ( $c = 1$ ). (2) The spread of two particles plus the spread of next two particles  $\geq (l_2 - l_1) + (l_3 - l_2) = 2$  ( $c = 2$ ). (3) ...
- the quasiparticle occupation patterns

$$n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 2020202020202020202 \dots \quad Q_\gamma = 0$$

$$n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 0202020202020202020 \dots \quad Q_\gamma = 1$$

$$n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 11111111111111111111 \dots \quad Q_\gamma = 1/2$$

are the only **close-packed** occupation patterns satisfying the above generalized Pauli exclusion rules.

- **close-packed:**

(1) add any electrons  $\rightarrow$  violate the GPE rules.

(2) shift any electrons  $\rightarrow$  violate the GPE rules.

7-cluster states:

$$P_{\frac{7}{2};Z_7} : (m; S_2, \dots, S_n) = (2; 0, 0, 0, 0, 0, 0, 0), \\ (n_0, \dots, n_{m-1}) = (7, 0)$$

$$P_{\frac{7}{8};Z_7^{(2)}} : (m; S_2, \dots, S_n) = (8; 0, 0, 2, 6, 10, 14), \\ (n_0, \dots, n_{m-1}) = (3, 0, 1, 0, 3, 0, 0, 0)$$

$$P_{\frac{7}{18};Z_7^{(3)}} : (m; S_2, \dots, S_n) = (18; 0, 4, 10, 18, 30, 42), \\ (n_0, \dots, n_{m-1}) = (2, 0, 0, 0, 0, 1, 0, 0, 0, 2, 0, 0, 0, 0, 0, 0)$$

$$P_{\frac{7}{14};C_7} : (m; S_2, \dots, S_n) = (14; 0, 2, 6, 12, 20, 28), \\ (n_0, \dots, n_{m-1}) = (2, 0, 1, 0, 1, 0, 1, 0, 2, 0, 0, 0, 0, 0, 0)$$

- Also get composite parafermion state  $P = P_{Z_{n_1}} P_{Z_{n_2}}$

## Pattern of zeros and generalized Pauli exclusion rule

- $l_{\gamma;a+1} - l_{\gamma;a} \geq S_2$  → the spacing between any two occupied orbitals by different electrons is no less than  $S_2$   
→ a generalized Pauli exclusion rule

### More general Pauli exclusion (GPE) rule

In terms of  $l_{\gamma;a} = S_{\gamma;a} - S_{\gamma;a-1}$ , the concave condition for quasiparticles becomes

$$\sum_{k=1}^b l_{\gamma;a+k} \geq S_b = \sum_{k=1}^b l_{a+k}, \quad \rightarrow \quad l_{\gamma;a} \geq l_a,$$

$$\sum_{k=1}^c (l_{\gamma;a+b+k} - l_{\gamma;a+k}) \geq S_{b+c} - S_b - S_c = \sum_{k=1}^c (l_{b+k} - l_k)$$

for any  $a, b, c \in \mathbb{Z}_+$ . Note that  $l_{\gamma;a+b} - l_{\gamma;a}$  is the spread of  $b + 1$  electrons. Setting  $c = 1 \rightarrow$  **The spread of  $b$  electrons  $\geq l_b$ .**

→ **The number of orbitals occupied by  $b$  electrons  $\geq l_b + 1$ .**

( $l_b$  = the spread of the first  $b$  electrons in the ground state.)



## Number of quasiparticle types from pattern of zeros

For the parafermion states  $P_{\nu=\frac{n}{2};Z_n}$  ( $m = 2$ ),

$P_{\frac{2}{2};Z_2}$	$P_{\frac{3}{2};Z_3}$	$P_{\frac{4}{2};Z_4}$	$P_{\frac{5}{2};Z_5}$	$P_{\frac{6}{2};Z_6}$	$P_{\frac{7}{2};Z_7}$	$P_{\frac{8}{2};Z_8}$	$P_{\frac{9}{2};Z_9}$	$P_{\frac{10}{2};Z_{10}}$
3	4	5	6	7	8	9	10	11

For the parafermion states  $P_{\nu=\frac{n}{2+2n};Z_n}$  ( $m = 2 + 2n$ )

$P_{\frac{2}{6};Z_2}$	$P_{\frac{3}{8};Z_3}$	$P_{\frac{4}{10};Z_4}$	$P_{\frac{5}{12};Z_5}$	$P_{\frac{6}{14};Z_6}$	$P_{\frac{7}{16};Z_7}$	$P_{\frac{8}{18};Z_8}$	$P_{\frac{9}{20};Z_9}$	$P_{\frac{10}{22};Z_{10}}$
9	16	25	36	49	64	81	100	121

For the generalized parafermion states  $P_{\nu=\frac{n}{m};Z_n^{(k)}}$

$P_{\frac{5}{8};Z_5^{(2)}}$	$P_{\frac{5}{18};Z_5^{(2)}}$	$P_{\frac{7}{8};Z_7^{(2)}}$	$P_{\frac{7}{22};Z_7^{(2)}}$	$P_{\frac{7}{18};Z_7^{(3)}}$	$P_{\frac{7}{32};Z_7^{(3)}}$	$P_{\frac{8}{18};Z_8^{(3)}}$	$P_{\frac{9}{8};Z_9^{(2)}}$
24	54	32	88	72	128	81	40

where  $k$  and  $n$  are coprime.

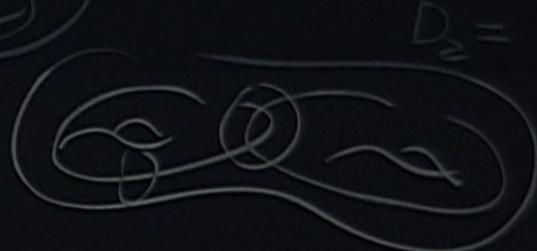


$$\ell = 0 \quad 1 \quad 2 \quad 3 \quad \ell_{a+b}$$

$a^{th}$



$$D_1 = 3$$



$$D_2 = ?$$

$$\ell_1 = 0$$

$$b+1 \boxed{\ell_{\delta, a+b} - \ell_{\delta, a} \geq \ell_b}$$

$(b+1) \text{ pent. } \ell_{b+1} \text{ orb.}$

For the  $\nu = 1$  Pfaffian state ( $n = 2$  and  $m = 2$ )

$$S_1, S_2, \dots : 0, 0, 2, 4, 8, 12, 18, 24, \dots$$

$$l_1 l_2 l_3 \dots : 0, 0, 2, 2, 4, 4, 6, 6, 8, 8, \dots$$

$$n_0 n_1 n_2 \dots : 2020202020202020202 \dots ,$$

- GPE rule: (1) Three particles occupy  $\geq l_3 + 1 = 3$  orbitals or more ( $c = 1$ ). (2) The spread of two particles plus the spread of next two particles  $\geq (l_2 - l_1) + (l_3 - l_2) = 2$  ( $c = 2$ ). (3) ...
- the quasiparticle occupation patterns

$$n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 2020202020202020202 \dots \quad Q_\gamma = 0$$

$$n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 0202020202020202020 \dots \quad Q_\gamma = 1$$

$$n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 11111111111111111111 \dots \quad Q_\gamma = 1/2$$

are the only **close-packed** occupation patterns satisfying the above generalized Pauli exclusion rules.

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(1) add any electrons  $\rightarrow$  violate the GPE rules.

(2) shift any electrons  $\rightarrow$  violate the GPE rules.

## GPE rule and edge spectrum for the Pfaffian state

First kind of edge:

$$M_0 : n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 20202020202|00000000 \dots$$

$$M_0 + 1 : n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 20202020201|10000000 \dots$$

$$M_0 + 1 : n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 20202020112|00000000 \dots \text{ not allowed}$$

$$M_0 + 2 : n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 20202020201|01000000 \dots$$

$$M_0 + 2 : n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 20202020200|20000000 \dots$$

$$M_0 + 2 : n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 20202020111|10000000 \dots$$

$$M_0 + 3 : n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 20202020201|00100000 \dots$$

$$M_0 + 3 : n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 20202020200|11000000 \dots$$

$$M_0 + 3 : n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 20202020111|01000000 \dots$$

$$M_0 + 3 : n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 20202020110|20000000 \dots$$

$$M_0 + 3 : n_{\gamma;0} n_{\gamma;1} n_{\gamma;2} \dots : 20202011111|10000000 \dots$$

$M$	$M_0$	$M_0 + 1$	$M_0 + 2$	$M_0 + 3$	$M_0 + 4$									
# of states $D_n$	1	1	3	5	10									

## Edge states = zero-energy states of ideal Hamiltonian

- For ideal Hamiltonian  $V_{1/3}(z_1, z_2) = -\partial_{z_1}\delta(z_1 - z_2)\partial_{z_1}$ , the  $N$  electron state  $P_{1/3} = \prod_{i < j}(z_i - z_j)^3$ , is the zero-energy state with minimal angular momentum (the order of  $z_i$ 's)  $M_0 = N(N - 1)$ .
- Other zero-energy state has higher angular momenta. Those zero energy states are the so called **edge states**:

The energy spectrum of 100 lowest levels of the ideal Hamiltonian with 6 electrons

