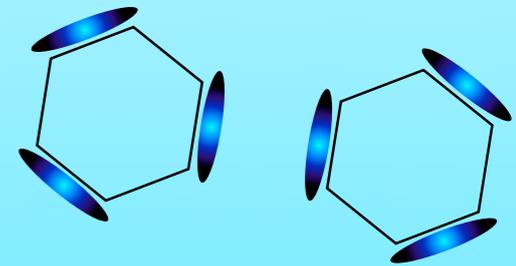


# Bulk-Edge correspondence and Fractionalization

As a topological (spin) insulator  
with strong interaction



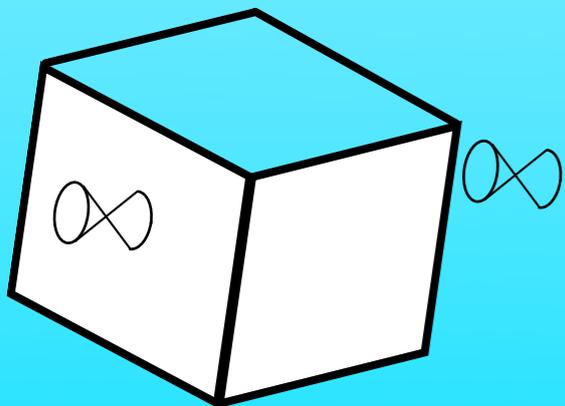
$$i\gamma_C(A_\psi) = \int_C A_\psi$$

Y. Hatsugai

Institute of Physics

Univ. of Tsukuba

JAPAN



# Plan

With time reversal invariance

- ★  $\mathbb{Z}_2$  Berry phase for a topological order parameter
- ★ Fractionalization for the Bulk in 1D & 2D
- ★ Entanglement Entropy to detect edge states
  - ★ (effective) Description by the Edges :
    - ★ Fractionalization at the Edges in 1D
    - deconfined spinons in 2D & 3D ??
- ★ Time Reversal operators with interaction
  - ★ Global to Local : super-selection rule  $\Theta^2 = 1, \text{ or } -1$

Let us consider

Gapped spin liquid as a topological insulator  
with strong interaction

# Quantum Liquids without Symmetry Breaking

## ★ *Quantum Liquids in Low Dimensional Quantum Systems*

★ *Low Dimensionality, Quantum Fluctuations*

★ *No Symmetry Breaking*

Topological Order

★ *No Local Order Parameter*

X.G.Wen

★ *Various Phases & Quantum Phase Transitions*

## ★ *Gapped Quantum Liquids in Condensed Matter*

★ *Integer & Fractional Quantum Hall States*

★ *Dimer Models of Fermions and Spins*

★ *Integer spin chains*

★ *Valence bond solid (VBS) states*

★ *Half filled Kondo Lattice*

*How to understand gapped quantum liquids ?*

# How to understand gapped quantum liquids ?

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**Bulk**

*classically featureless : need geometrical phase*

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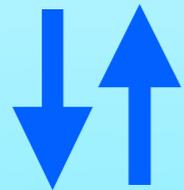
*low energy localized modes in the gap*

*edge states for QHE*

*Laughlin, Halperin, YH*

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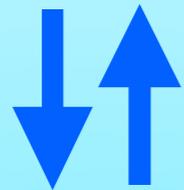
# How to understand gapped quantum liquids ?

## Bulk-Edge correspondence

Common property of topological ordered states

Bulk

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1-st Chern number for QHE TKNN

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*low energy localized modes in the gap*

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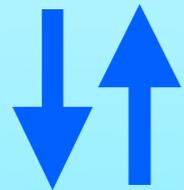
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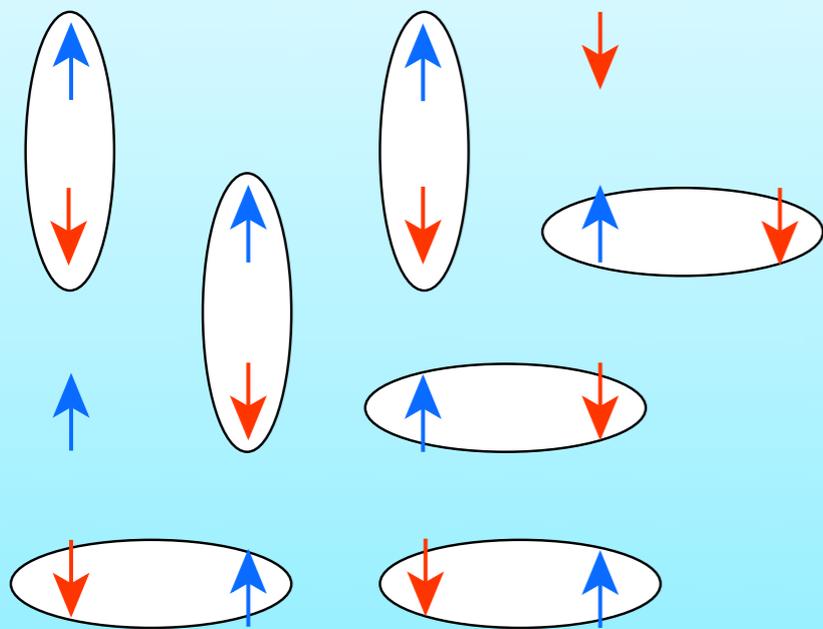
Laughlin, Halperin, YH

## As for quantum spins

- ★  $Z_2$  Berry Phase as a Topological Order Parameter of bulk
- ★ Entanglement Entropy to detect edge states (generic Kennedy triplets)

# Quantum Liquid (Example 1)

## ★ The *RVB* state by Anderson



$$|\text{Singlet Pair}_{12}\rangle = \frac{1}{\sqrt{2}} (|\uparrow_1\downarrow_2\rangle - |\downarrow_1\uparrow_2\rangle)$$

$$|G\rangle = \sum_{J=\text{Dimer Covering}} c_J \otimes_{ij} |\text{Singlet Pair}_{ij}\rangle$$

*small magnets*

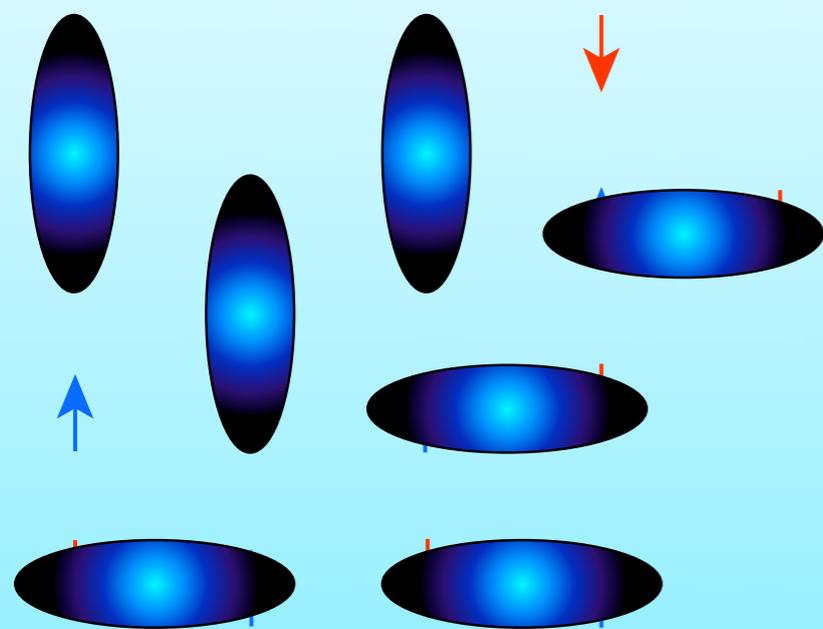


*Local Singlet Pairs :*  
*(Basic Objects)*

*Purely Quantum Objects are basic*

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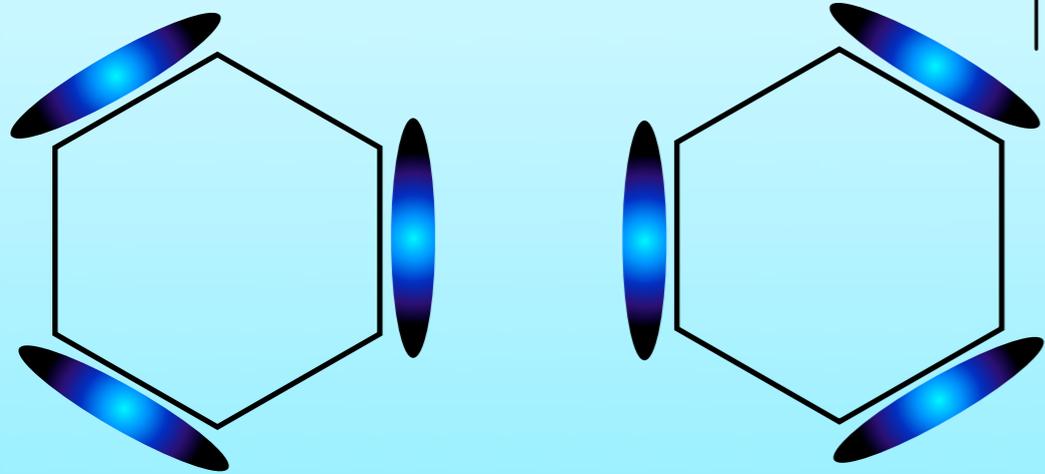


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# Quantum Liquid (Example 2)

## ★ The *RVB* state by Pauling



$$|\text{Bond}_{12}\rangle = \frac{1}{\sqrt{2}}(|1\rangle + |2\rangle) = \frac{1}{\sqrt{2}}(c_1^\dagger + c_2^\dagger)|0\rangle$$

$$|G\rangle = \sum_{J=\text{Dimer Covering}} c_J \otimes_{ij} |\text{Bond}_{ij}\rangle$$

Do Not use the Fermi Sea

localized charge  
at site *A*

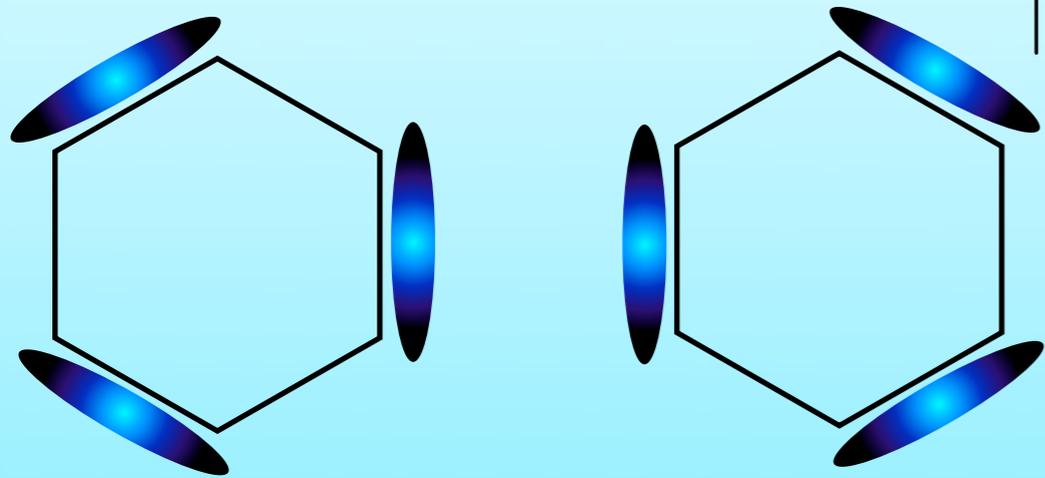
*A*

Local Covalent Bonds :  
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# Quantum Liquid (Example 2)

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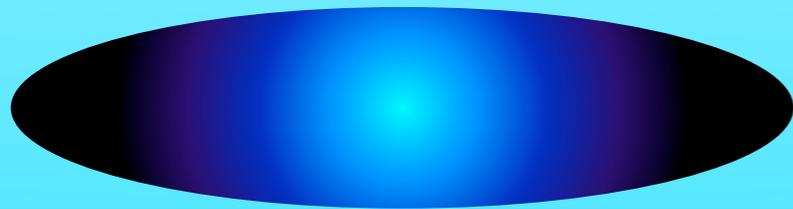
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Do Not use the Fermi Sea

*Delocalized* charge  
as a *covalent bond*

*Local Covalent Bonds* :  
(Basic Objects)



Purely Quantum Objects are basic

# Quantum Interference for the Classification

## ★ “Classical” Observables

★ Charge density, Spin density, ...  $\mathcal{O} = n_{\uparrow} \pm n_{\downarrow}, \dots$

$$\langle \mathcal{O} \rangle_G = \langle G | \mathcal{O} | G \rangle = \langle G' | \mathcal{O} | G' \rangle = \langle \mathcal{O} \rangle_{G'}$$

$$|G'\rangle = |G\rangle e^{i\phi}$$

## ★ “Quantum” Observables !

★ Quantum Interferences:  $\langle G_1 | G_2 \rangle = \langle G'_1 | G'_2 \rangle e^{i(\phi_1 - \phi_2)}$

★ Probability Amplitude (overlap)  $|G_i\rangle = |G'_i\rangle e^{i\phi_i}$

★ Aharonov-Bohm Effects

★ Phase (Gauge) dependent

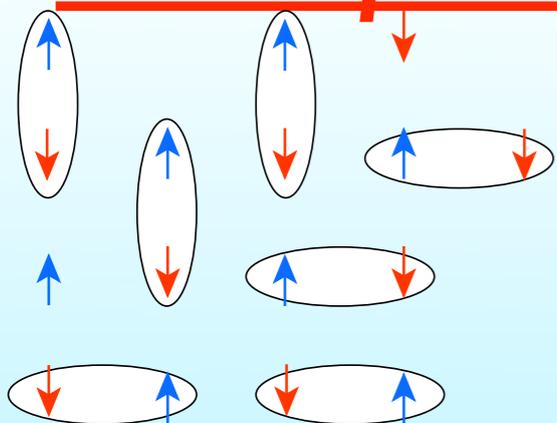
$$\langle G | G + dG \rangle = 1 + \langle G | dG \rangle$$

$$A = \langle G | dG \rangle \text{ :Berry Connection}$$

$$i\gamma = \int A \text{ :Berry Phase}$$

Use Quantum Interferences To Classify Quantum Liquids

# Examples: RVB state by Anderson

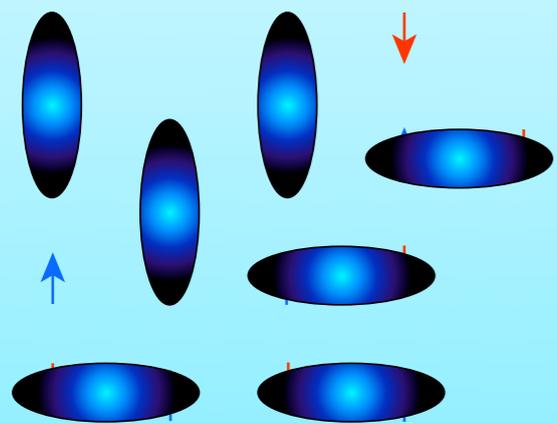


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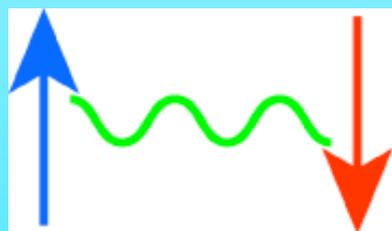
No Long Range Order  
in Spin-Spin Correlation



*Local Singlet Pair is a Basic Object*

## How to Characterize the Local Singlet Pair ?

$$|G\rangle = \frac{1}{\sqrt{2}} (|\uparrow_i\downarrow_j\rangle - |\downarrow_i\uparrow_j\rangle)$$

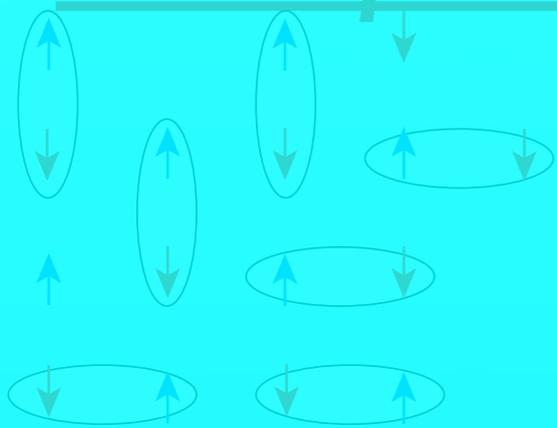


*Use Berry Phase to characterize the Singlet!*

*Singlet does not carries spin but does Berry phase*

$$\gamma_{\text{singlet pair}} = \pi \pmod{2\pi}$$

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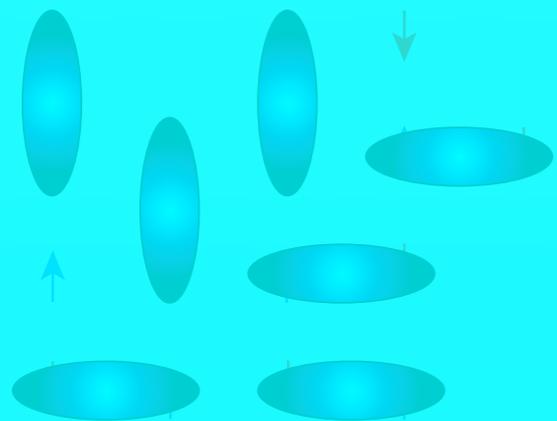


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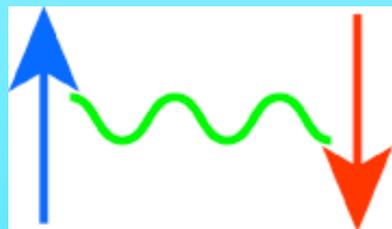
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# $Z_2$ Berry phases for gapped quantum spins

★ generic Heisenberg Models (with *frustration*)

$$H = \sum_{ij} J_{ij} \mathbf{S}_i \cdot \mathbf{S}_j$$

*Time Reversal Invariant*

$$\Theta_N \mathbf{S}_i \Theta_N^{-1} = -\mathbf{S}_i$$

$$[H, \Theta_N] = 0$$

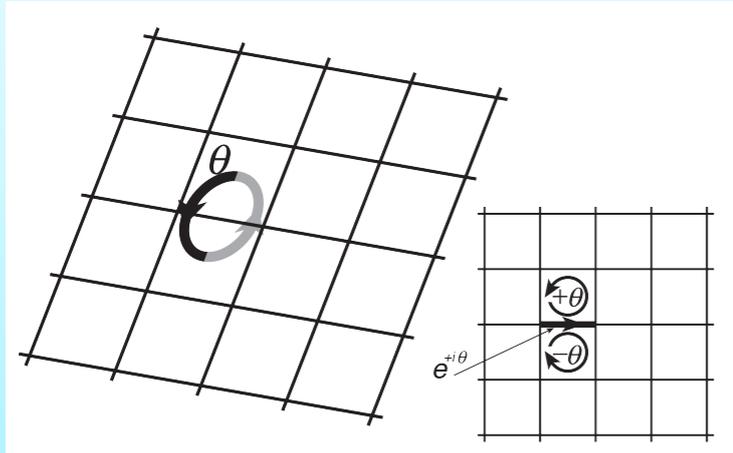
$$\Theta_N = (i\sigma_y^1) \otimes (i\sigma_y^2) \cdots (i\sigma_y^N) K$$

$$\Theta_N^2 = (-1)^N$$

*Mostly N: even*  $\Theta_N^2 = 1$  (probability 1/2 in HgTe)

# $Z_2$ Berry phases for gapped quantum spins

Define a many body hamiltonian by local twist as a parameter



$$H(x = e^{i\theta})$$

$$C = \{x = e^{i\theta} | \theta : 0 \rightarrow 2\pi\} \quad U(1)$$

$$\mathbf{S}_i \cdot \mathbf{S}_j \rightarrow \frac{1}{2} (e^{-i\theta} S_{i+} S_{j-} + e^{+i\theta} S_{i-} S_{j+}) + S_{iz} S_{jz} \quad \text{Only link } \langle ij \rangle$$

Calculate the Berry Phases using the Entire Many Spin Wavefunction *numerically*

$Z_2$  quantization

Require excitation Gap!

$$\gamma_C = \int_C A_\psi = \int_C \langle \psi | d\psi \rangle = \begin{cases} 0 \\ \pi \end{cases} : \text{mod } 2\pi \quad \mathbf{Z}_2$$

Time Reversal ( Anti-Unitary ) Invariance

# Berry Connection and Gauge Transformation

★ **Parameter Dependent Hamiltonian**

$$H(x) \left| \psi(x) \right\rangle = E(x) \left| \psi(x) \right\rangle$$

$$H(x) \left| \psi(x) \right\rangle = E(x) \left| \psi(x) \right\rangle, \quad \langle \psi(x) | \psi(x) \rangle = 1.$$

★ **Berry Connections**  $A_\psi = \langle \psi | d\psi \rangle = \langle \psi | \frac{d}{dx} \psi \rangle dx.$

★ **Berry Phases**  $i\gamma_C(A_\psi) = \int_C A_\psi$  (Abelian)

★ **Phase Ambiguity of the eigen state**

$$\left| \psi(x) \right\rangle = \left| \psi'(x) \right\rangle e^{i\Omega(x)}$$

**Gauge Transformation**

$$A_\psi = A'_\psi + id\Omega = A'_\psi + i \frac{d\Omega}{dx} dx$$

★ **Berry phases are not well-defined without**

$$\gamma_C(A_\psi) = \gamma_C(A_{\psi'}) + \int_C d\Omega$$

**specifying the gauge**  
 $2\pi \times (\text{integer})$  if  $e^{i\Omega}$  is single valued

★ **Well Defined up to mod  $2\pi$**

$$\gamma_C(A_\psi) \equiv \gamma_C(A_{\psi'}) \pmod{2\pi}$$

# Anti-Unitary Operator and Berry Phases

## ★ **Anti-Unitary Operator** (Time Reversal, Particle-Hole)

$$\Theta = KU_{\Theta}, \quad K : \text{Complex conjugate} \\ U_{\Theta} : \text{Unitary} \quad (\text{parameter independent})$$

$$|\Psi\rangle = \sum_J C_J |J\rangle \quad \sum_J C_J^* C_J = \langle\Psi|\Psi\rangle = 1$$

$$|\Psi^{\Theta}\rangle = \Theta|\Psi\rangle = \sum_J C_J^* |J^{\Theta}\rangle, \quad |J^{\Theta}\rangle = \Theta|J\rangle$$

## ★ **Berry Phases and Anti-Unitary Operation**

$$A^{\Psi} = \langle\Psi|d\Psi\rangle = \sum_J C_J^* dC_J \quad \sum_J dC_J^* C_J + \sum_J C_J^* dC_J = 0$$

$$A^{\Theta\Psi} = \langle\Psi^{\Theta}|d\Psi^{\Theta}\rangle = \sum_J C_J dC_J^* = -A^{\Psi}$$

$$\gamma_C(A^{\Theta\Psi}) = -\gamma_C(A^{\Psi})$$

# Anti-Unitary **Invariant State** and $\mathbb{Z}_2$ Berry Phase

★ **Anti-Unitary Symmetry**  $[H(x), \Theta] = 0$

★ **Invariant State**  $\exists \varphi, |\Psi^\Theta\rangle = \Theta|\Psi\rangle = |\Psi\rangle e^{i\varphi}$

★ ex. **Unique Eigen State**  $\simeq |\Psi\rangle$  Gauge Equivalent (Different Gauge)

★ **To be compatible with the ambiguity,**

the Berry Phases have to be **quantized** as

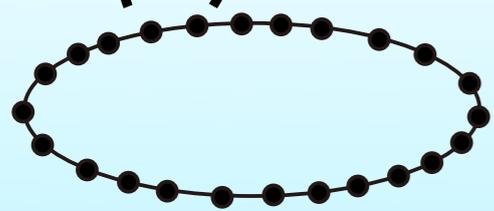
$$\gamma_C(A^\Psi) = \begin{cases} 0 \\ \pi \end{cases} \pmod{2\pi}$$

**$\mathbb{Z}_2$  Berry phase**

$$\gamma_C(A^\Psi) = -\gamma_C(A^{\Theta\Psi}) \equiv -\gamma_C(A^\Psi), \pmod{2\pi}$$

# Numerical Evaluation of the Berry Phases (incl. non-Abelian)

(1) Discretize the periodic parameter space



$$x_0, x_1, \dots, x_N = x_0 \quad \theta_0 = 0, \theta_N = 2\pi$$
$$x_n = e^{i\theta_n} \quad \theta_{n+1} = \theta_n + \Delta\theta_n \quad \forall \Delta\theta_n \rightarrow 0$$

(2) Obtain eigen vectors

$$H(x_n)|\psi_n^i\rangle = E^i(x_n)|\psi_n^i\rangle$$

(3) Define Berry connection in a discretized form

$$A_n = \text{Im} \log \langle \psi_n | \psi_{n+1} \rangle$$

*non-Abelian*  $A_n = \text{Im} \log \det D_n, \{D_n\}_{ij} = \langle \psi_n^i | \psi_{n+1}^j \rangle$

(4) Evaluate the Berry phase

$$\gamma = \sum_{n=0}^{N-1} A_n = \text{Im} \log \langle \psi_0 | \psi_1 \rangle \langle \psi_1 | \psi_2 \rangle \cdots (= \text{Im} \log \det D_1 D_2 \cdots D_n) \quad \text{non-Abelian}$$

**Independent of the choice of the phase**

$$|\psi_n\rangle \rightarrow |\psi_n\rangle' e^{i\Omega_n}$$

**Gauge invariant**

Luscher '82 (Lattice Gauge Theory)

**after the discretization**

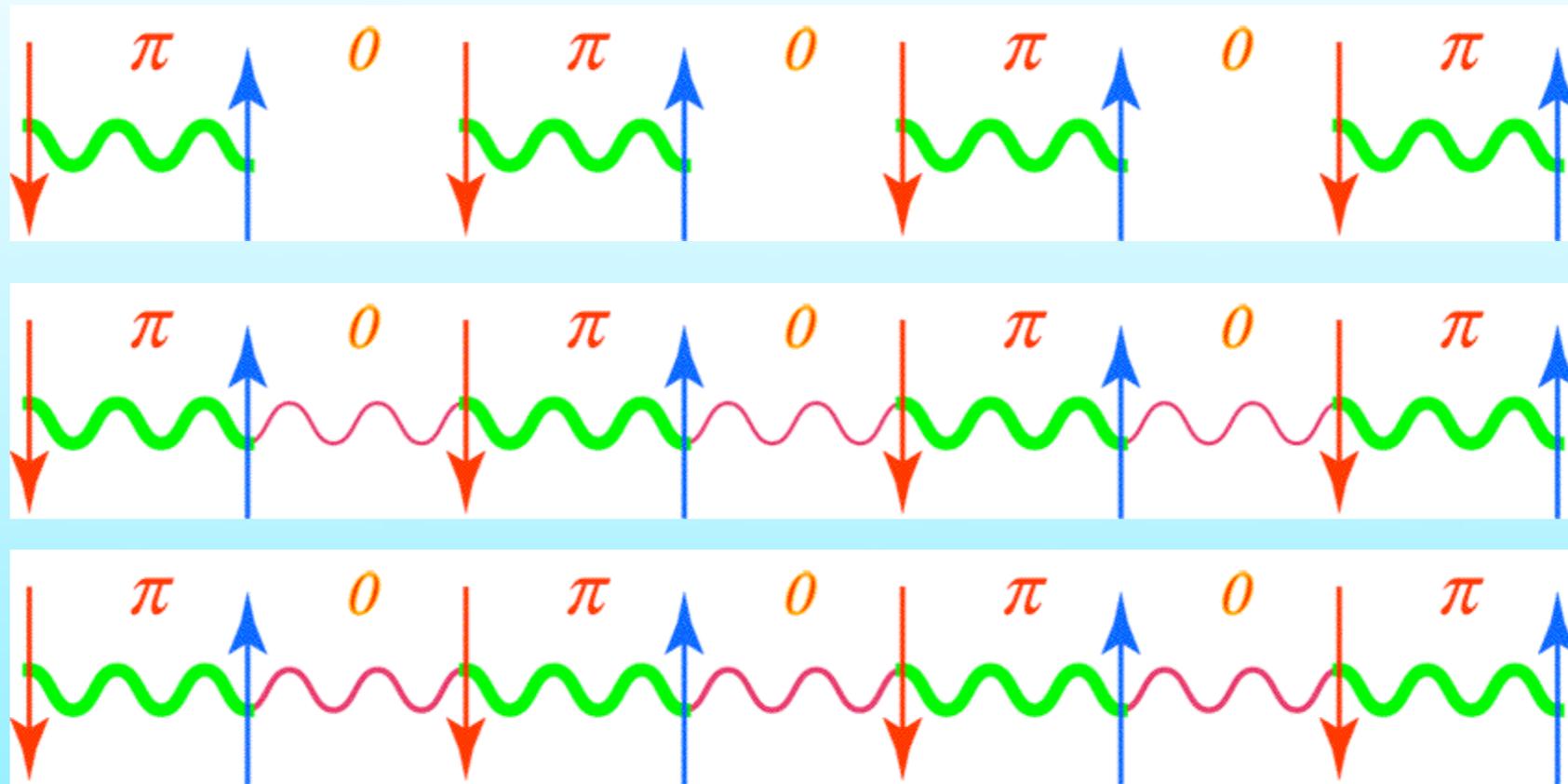
King-Smith & Vanderbilt '93 (polarization in solids)

**Convenient for Numerics**

T. Fukui, H. Suzuki & YH '05 (Chern numbers)

# Adiabatic Continuation & the Quantization

Introduce interaction between singlets



★  $Z_2$ -quantization of the Berry phases **protects** from **continuous change**

Adiabatic Continuation in a **gapped** system



Renormalization Group in a **gapless** system

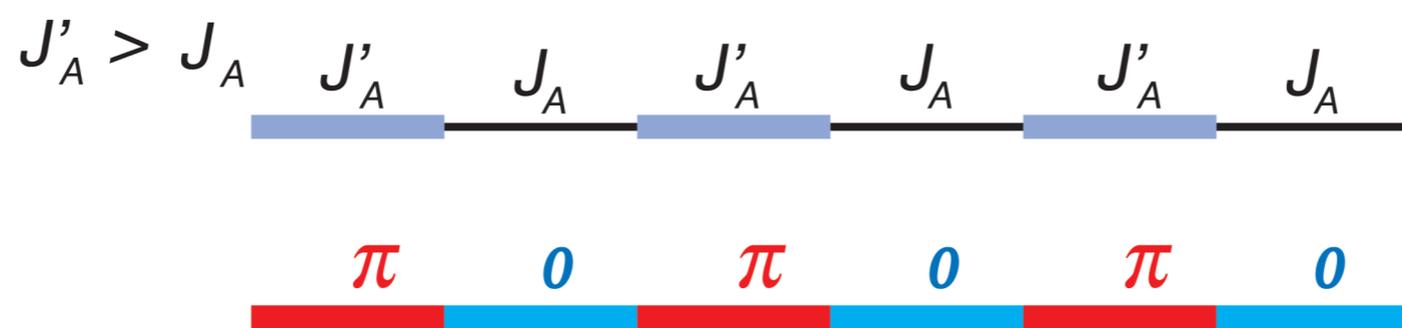
# Local Order Parameters of Singlet Pairs

## ★ 1D AF-AF, AF-F Dimers

Y.H., J. Phys. Soc. Jpn. 75 123601 (2006)

- ★ Strong Coupling Limit of the AF Dimer link is a gapped unique ground state.

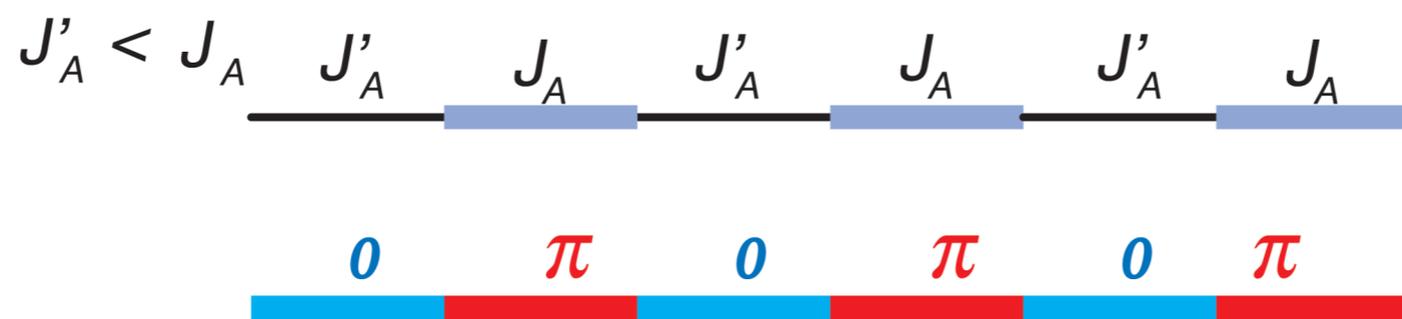
AF-AF



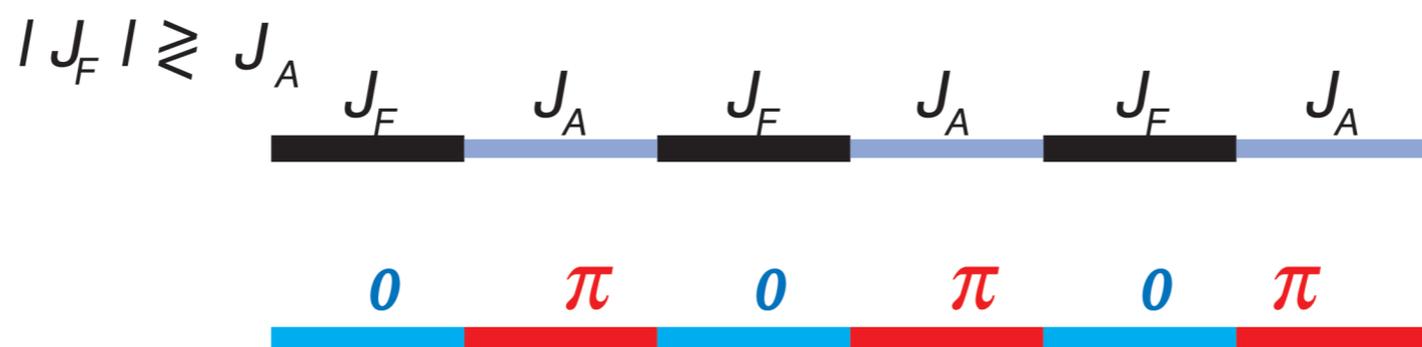
AF-AF case

**Strong** bonds

:  $\pi$  bonds



F-AF



F-AF case

**AF** bonds

:  $\pi$  bonds

Hida

# Local Order Parameters of the Haldane Phase

## ★ Heisenberg Spin Chains with integer $S$

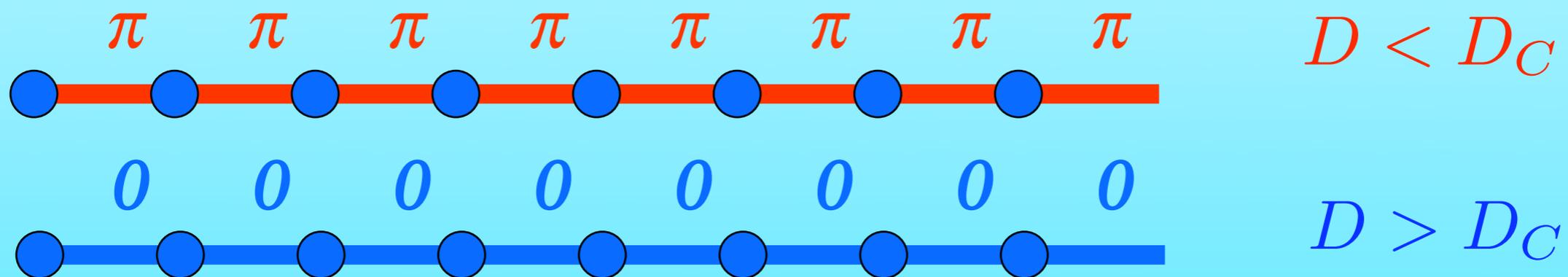
★ No Symmetry Breaking by the Local Order Parameter

★ "String Order": Non-Local Order Parameter!

**$S=1$**   $(S_i)^2 = S(S+1), S=1$

$$H = J \sum_{\langle ij \rangle} \mathbf{S}_i \cdot \mathbf{S}_j + D \sum_i (S_i^z)^2$$

Y.H., J. Phys. Soc. Jpn. 75 123601 (2006)



Describe the Quantum Phase Transition locally

c.f.  $S=1/2$ , 1D dimers, 2D with Frustrations, Ladders  
 $t$ -J with Spin gap

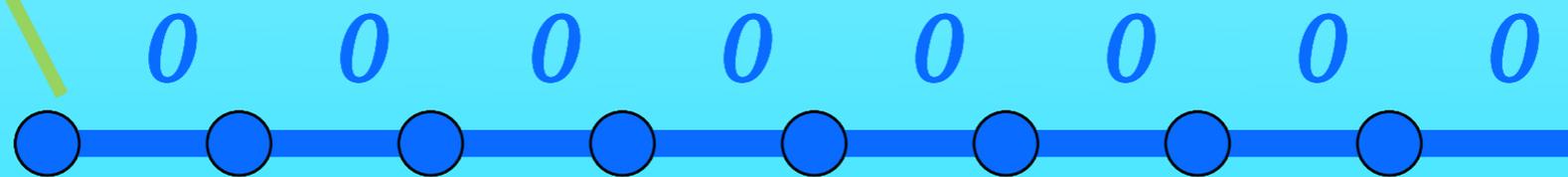
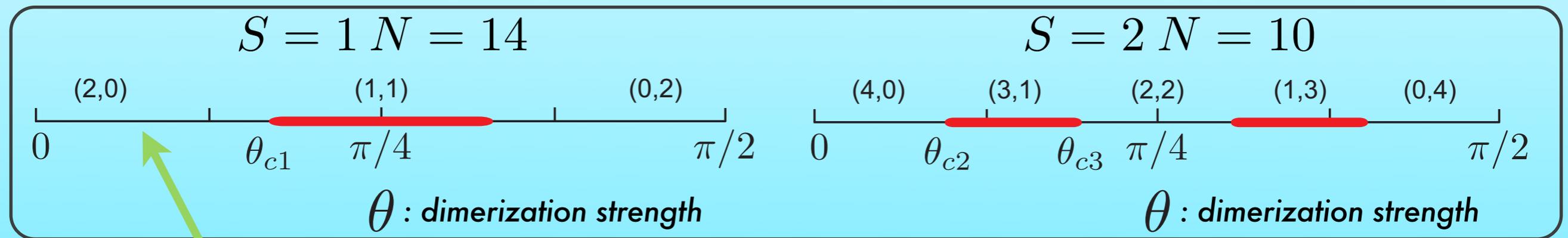
# Topological Classification of Gapped Spin Chains

T.Hirano, H.Katsura & YH, Phys.Rev.B77 094431'08

- ♦  $S=1,2$  dimerized Heisenberg model

$$H = \sum_{i=1}^{N/2} (J_1 \mathbf{S}_{2i} \cdot \mathbf{S}_{2i+1} + J_2 \mathbf{S}_{2i+1} \cdot \mathbf{S}_{2i+2}) \quad J_1 = \cos \theta, J_2 = \sin \theta$$

## $Z_2$ Berry phase



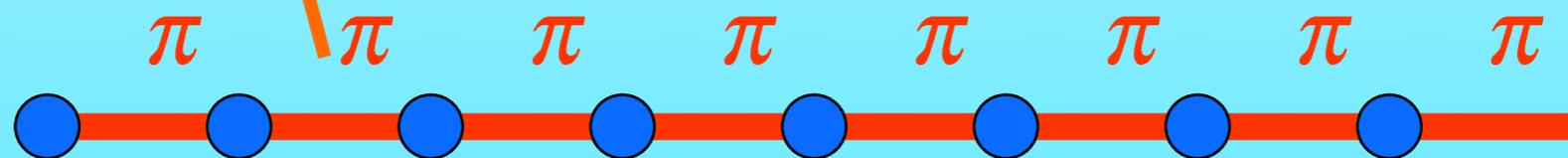
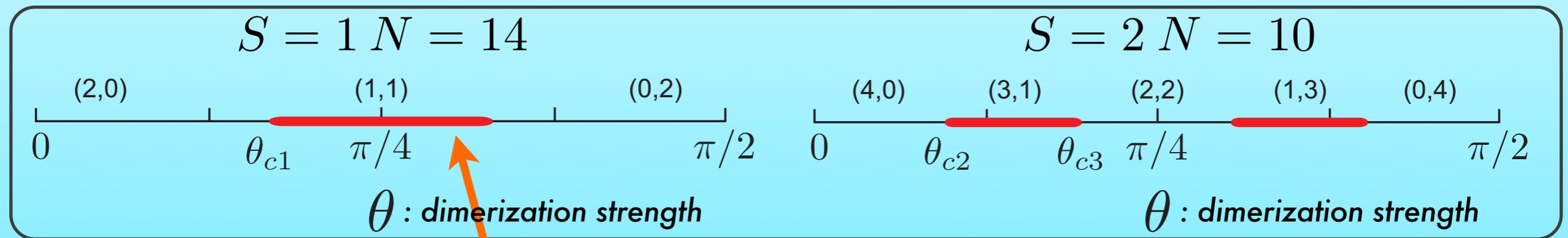
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## Z<sub>2</sub>Berry phase



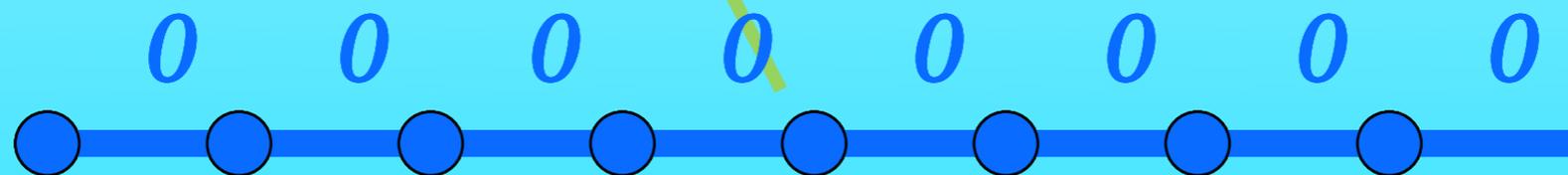
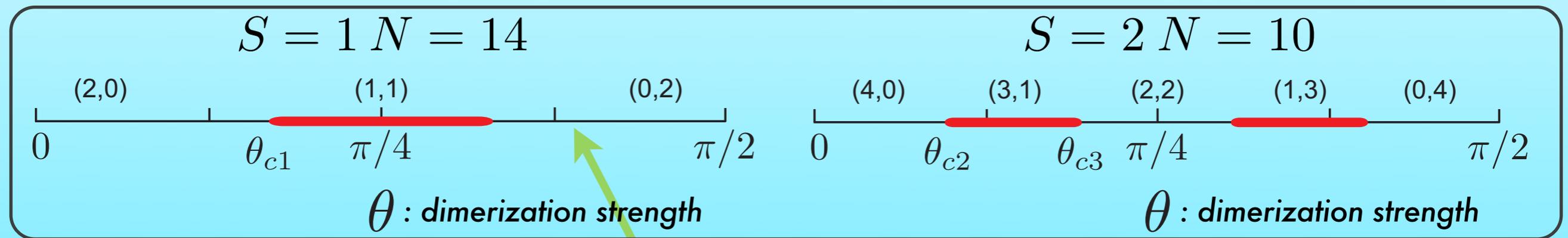
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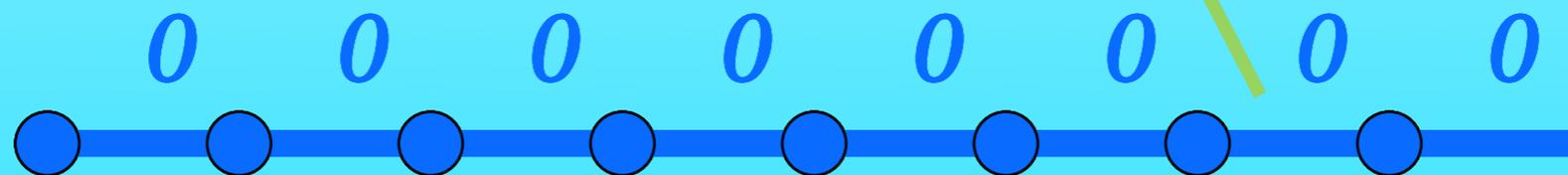
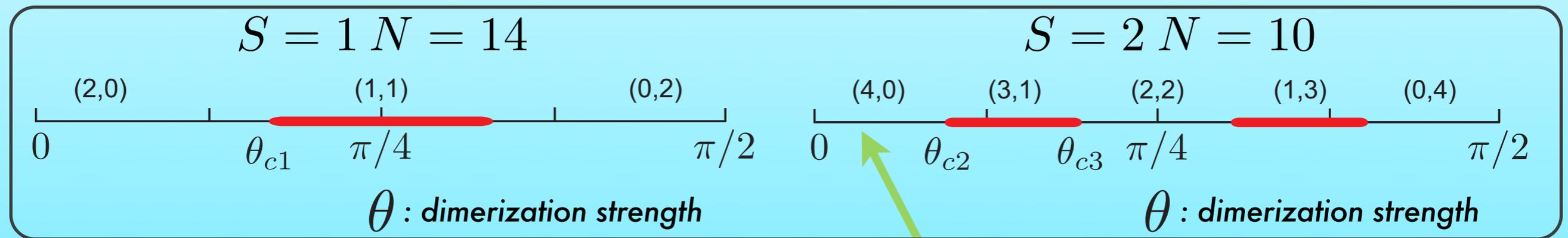
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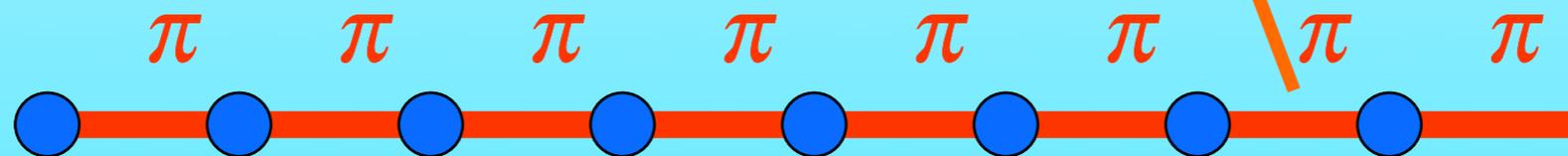
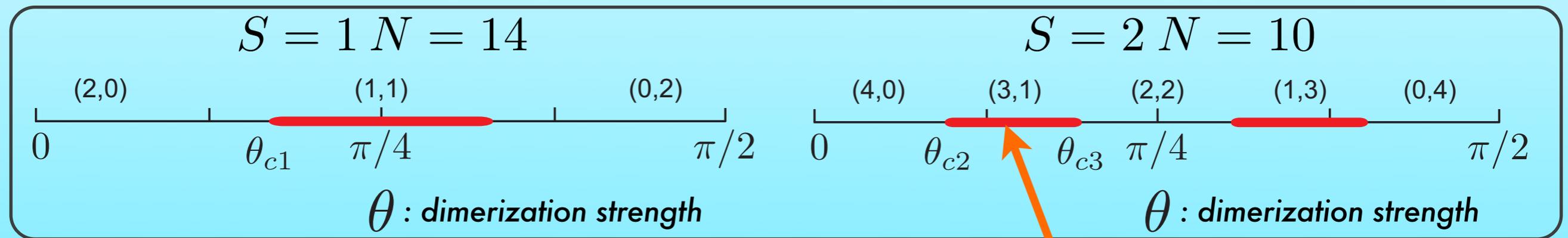
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$$H = \sum_{i=1}^{N/2} (J_1 \mathbf{S}_{2i} \cdot \mathbf{S}_{2i+1} + J_2 \mathbf{S}_{2i+1} \cdot \mathbf{S}_{2i+2}) \quad J_1 = \cos \theta, J_2 = \sin \theta$$

## $Z_2$ Berry phase



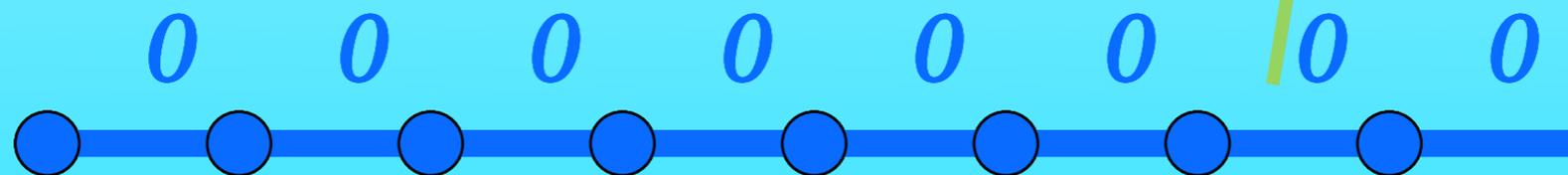
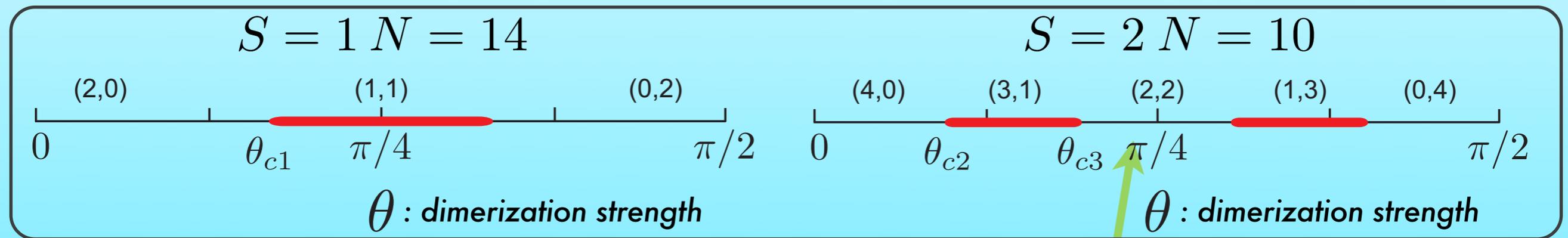
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T.Hirano, H.Katsura & YH, Phys.Rev.B77 094431'08

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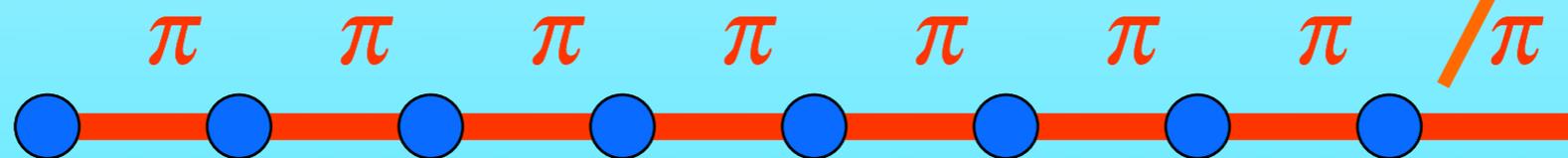
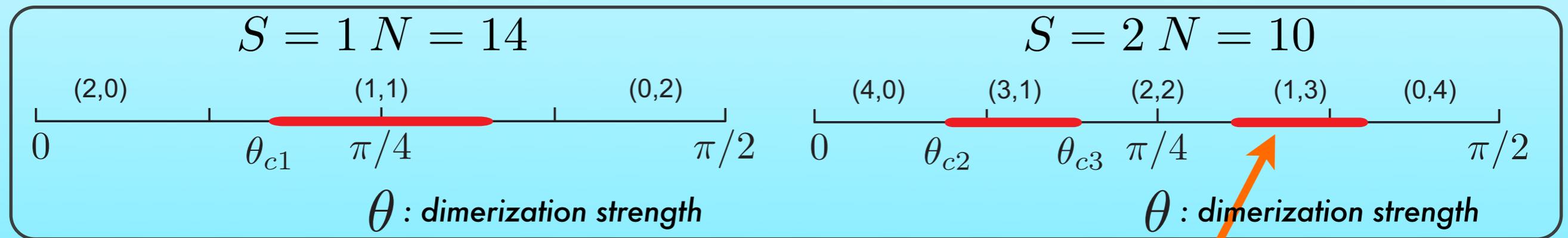
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## Z<sub>2</sub>Berry phase



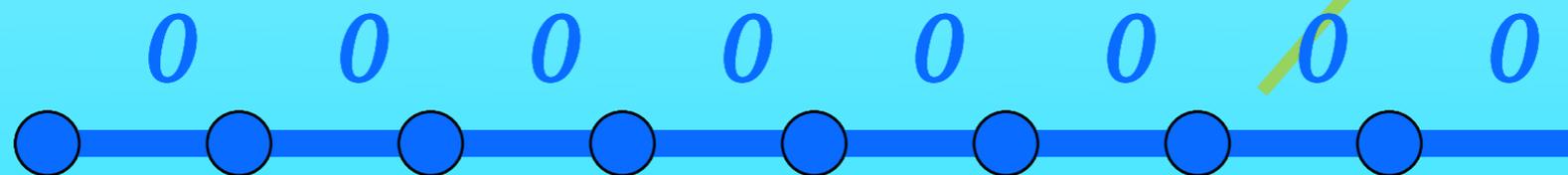
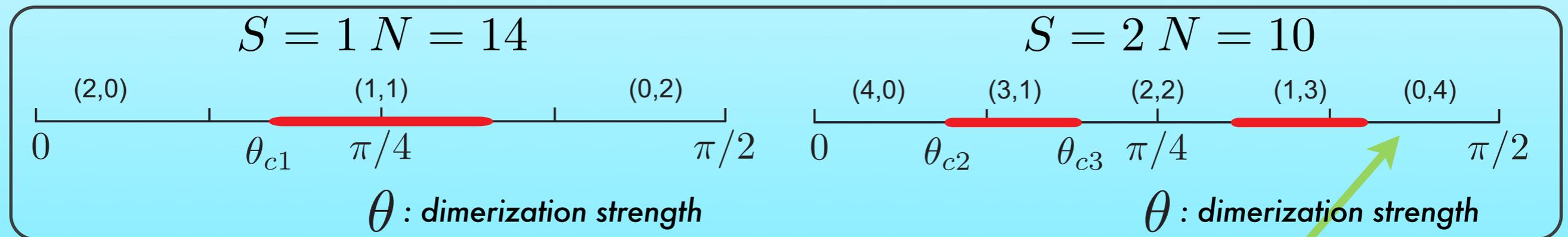
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## Z<sub>2</sub>Berry phase



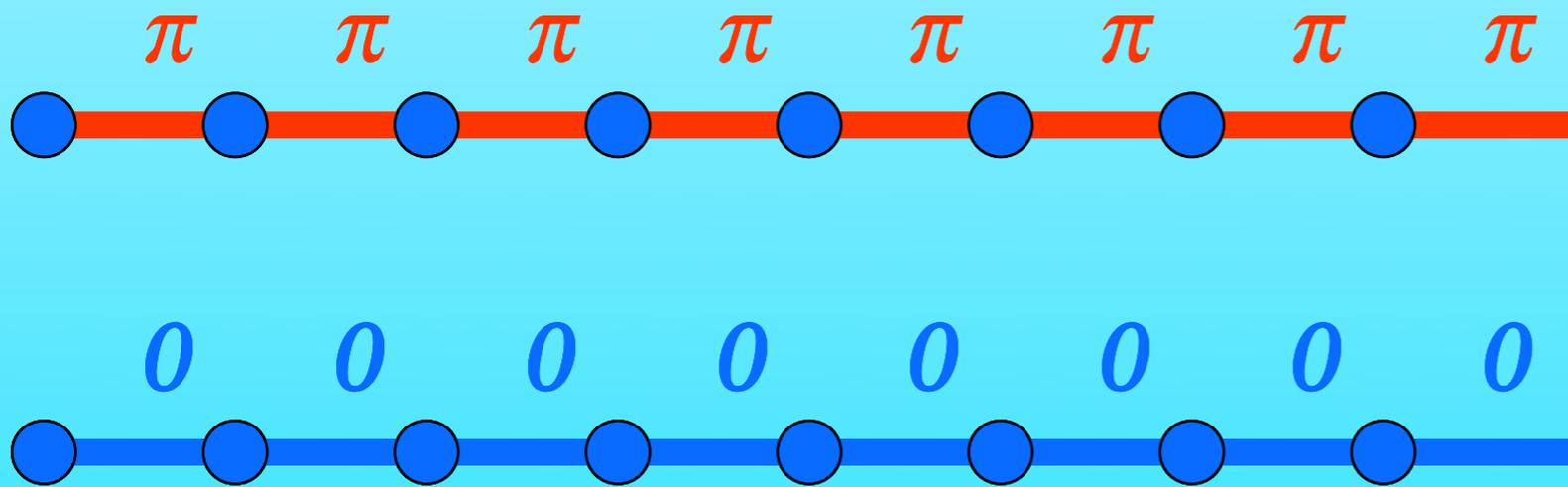
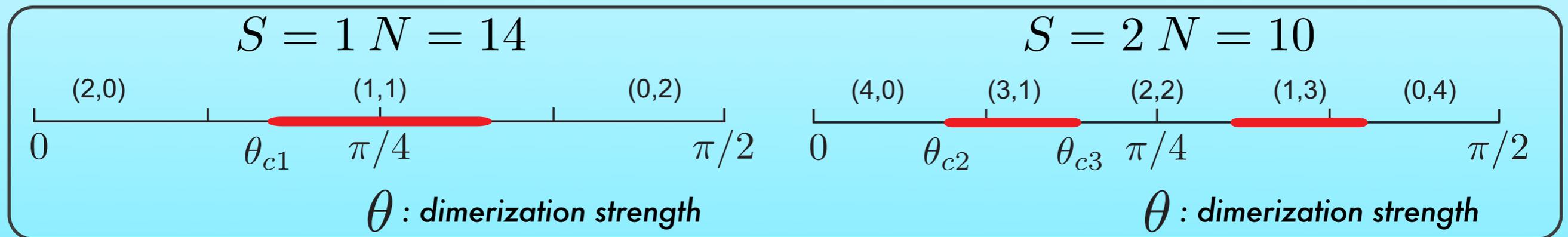
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T.Hirano, H.Katsura & YH, Phys.Rev.B77 094431'08

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## $Z_2$ Berry phase



Topological Quantum Phase Transitions with **translation** invariance

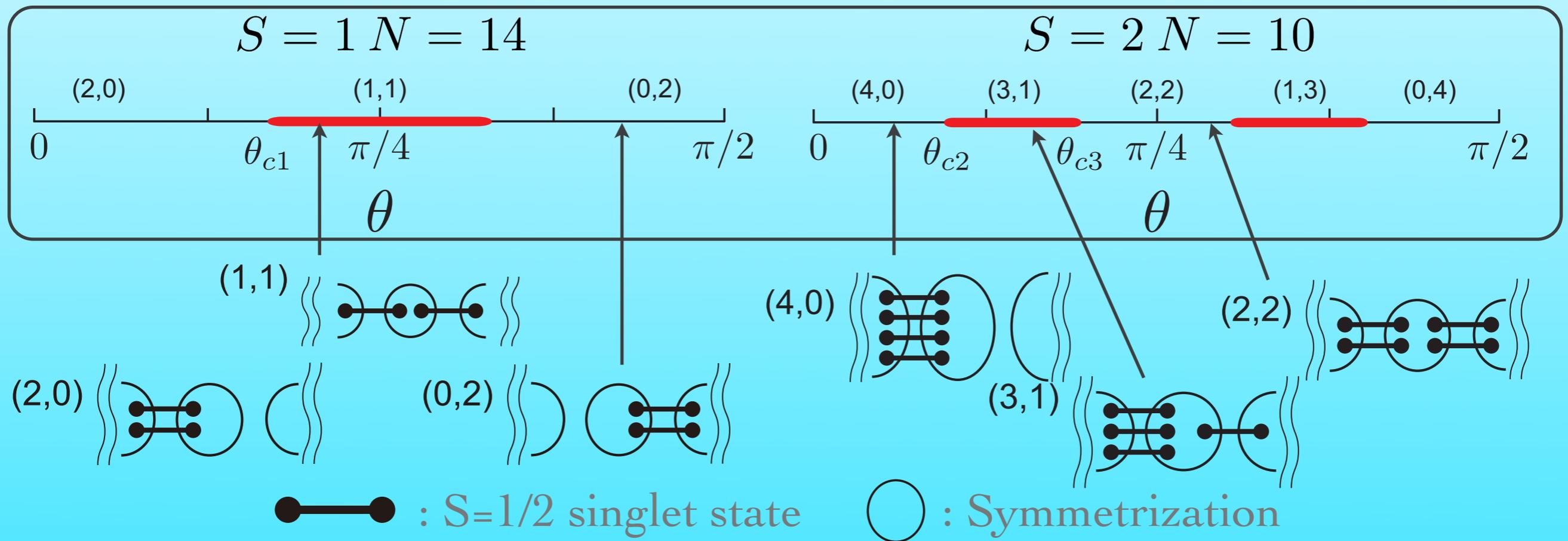
# Topological Classification of Gapped Spin Chains

T.Hirano, H.Katsura & YH, Phys.Rev.B77 094431'08

- ◆ S=1,2 dimerized Heisenberg model

$$H = \sum_{i=1}^{N/2} (J_1 \mathbf{S}_{2i} \cdot \mathbf{S}_{2i+1} + J_2 \mathbf{S}_{2i+1} \cdot \mathbf{S}_{2i+2}) \quad J_1 = \cos \theta, J_2 = \sin \theta$$

## Berry phase



**Reconstruction of valence bonds!**

# Topological Classification of Gapped Spin Chains (cont.)

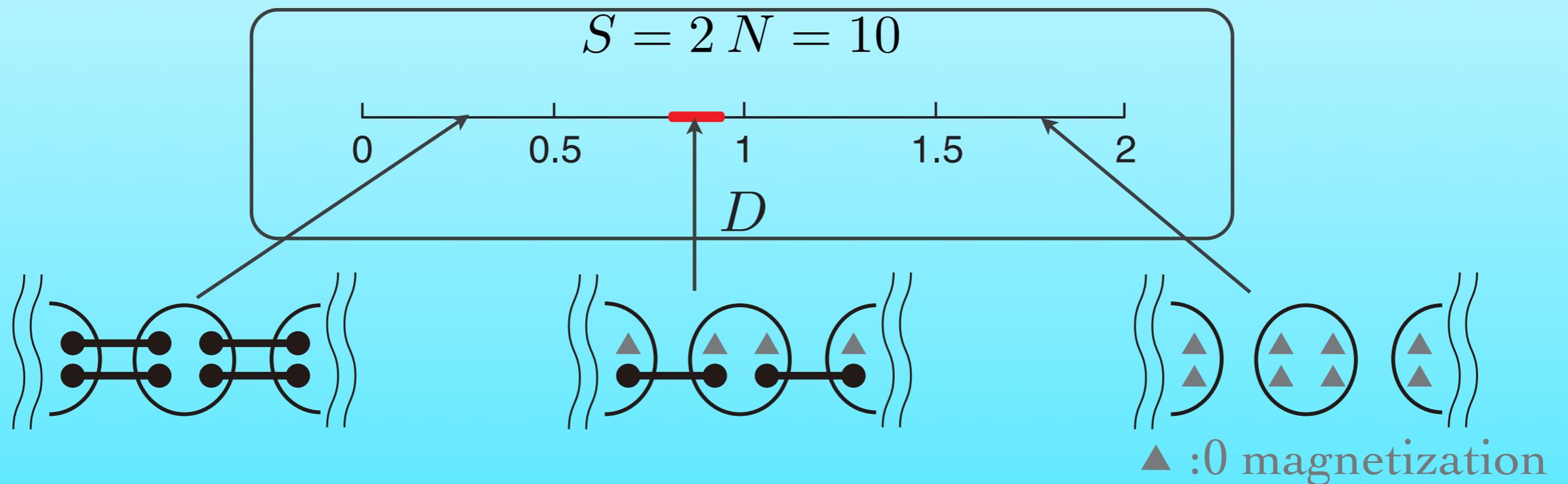
T.Hirano, H.Katsura & YH, Phys.Rev.B77 094431'08

- ♦ S=2 Heisenberg model with D-term

$$H = \sum_i^N \left[ J \mathbf{S}_i \cdot \mathbf{S}_{i+1} + D (S_i^z)^2 \right]$$

## Berry phase

Red line denotes the non trivial Berry phase



Reconstruction of valence bonds!

# Topological Classification of Generic AKLT (VBS) models

T.Hirano, H.Katsura & YH, Phys.Rev.B77 094431'08

Twist the link of the generic AKLT model

$$H(\{\phi_{i,i+1}\}) = \sum_{i=1}^N \sum_{J=B_{i,i+1}+1}^{2B_{i,i+1}} A_J P_{i,i+1}^J[\phi_{i,i+1}]$$

$$|\{\phi_{i,j}\}\rangle = \prod_{\langle ij \rangle} \left( e^{i\phi_{ij}/2} a_i^\dagger b_j^\dagger - e^{-i\phi_{ij}/2} b_i^\dagger a_j^\dagger \right)^{B_{ij}} |\text{vac}\rangle$$

Berry phase on a link (ij)

$$\gamma_{ij} = B_{ij} \pi \text{ mod } 2\pi$$

$S=1/2$

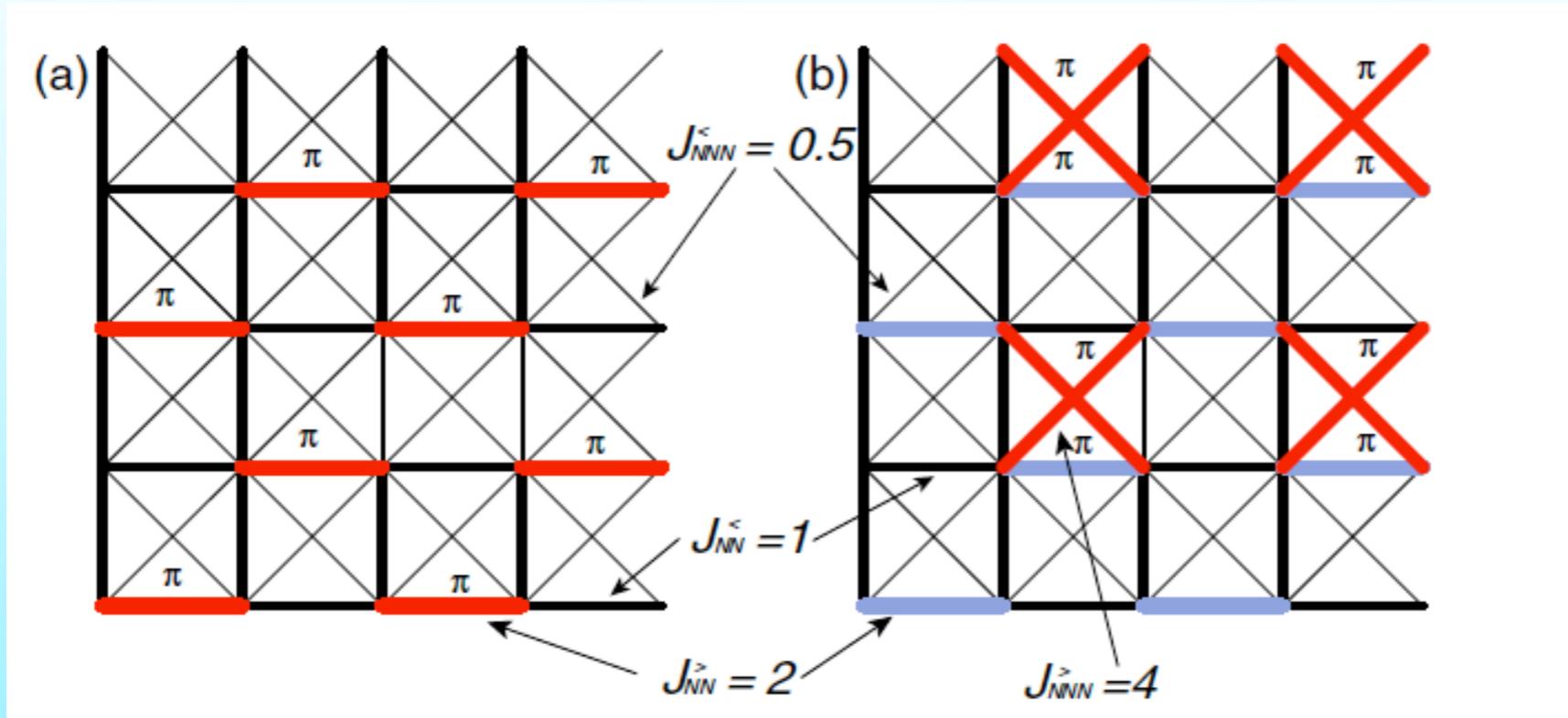
The Berry phase counts the number of the valence bonds!

$S=1/2$  objects are fundamental in  $S=1$  &  $2$  spin chains

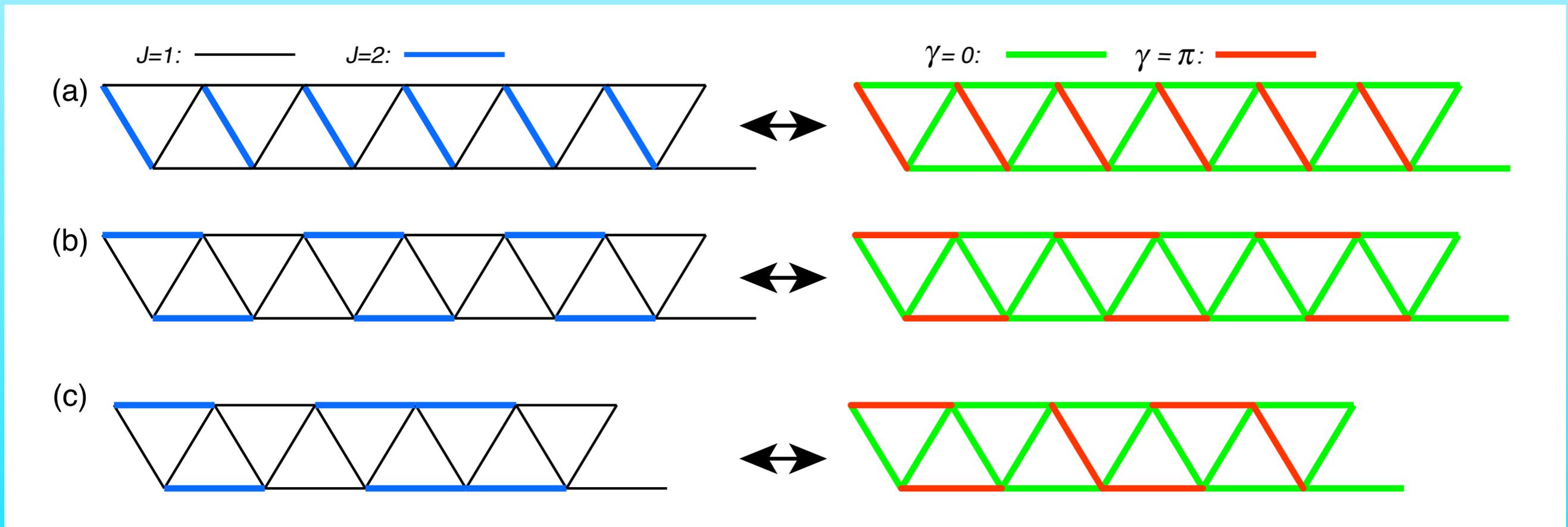
**FRACTIONALIZATION**

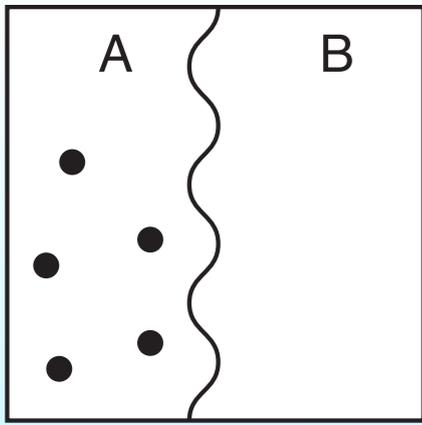
Contribute to the  
Entanglement Entropy  
as of Edge states

# 2D, Ladders ( $S=1/2$ ), $t$ - $J$ (spin gapped)



Y.H., J. Phys. Soc. Jpn. 75 123601 (2006), J. Phys. Cond. Matt.19, 145209 (2007)





- Entanglement Entropy to detect edge states*
- direct calculation of spectrum with boundaries*

# Entanglement Entropy

## ★ Mixed State From Entanglement

Vidal, Latorre, Rico, Kitaev '02

### ★ Direct Product State

$$|\Psi_{AB}\rangle = |\Psi_A\rangle \otimes |\Psi_B\rangle$$

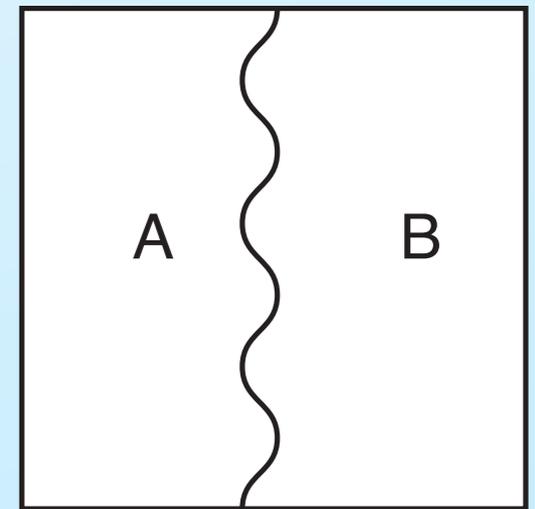
System =  $A \oplus B$

$$\text{State} = \sum \Psi_A \otimes \Psi_B$$

### ★ Entangled State

## ★ Partial Trace

$$|\Psi_{AB}\rangle = \frac{1}{\sqrt{D}} \sum_j |\Psi_A^j\rangle \otimes |\Psi_B^j\rangle$$



$$\rho_{AB} = |\Psi_{AB}\rangle\langle\Psi_{AB}|$$

Pure State

$D = 1$

$$\rho_A = \text{Tr}_B \rho_{AB}$$

$$= \frac{1}{D} \sum_j |\Psi_A^j\rangle\langle\Psi_A^j|$$

Mixed State

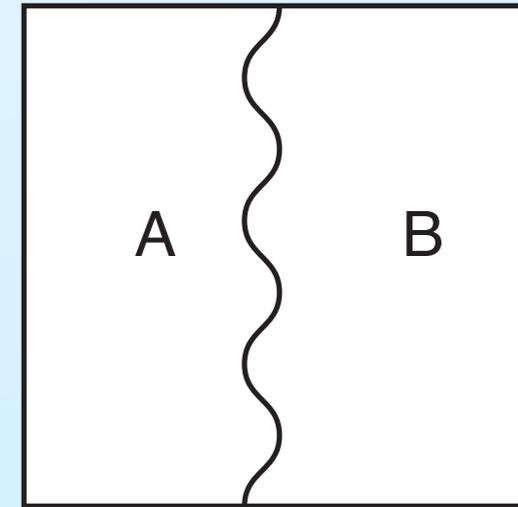
$$\rho_{AB} = \frac{1}{D} \sum_{jk} |\Psi_A^j\rangle\langle\Psi_A^k| \otimes |\Psi_B^j\rangle\langle\Psi_B^k|$$

## ★ How much the State is Entangled between A & B?

Entanglement Entropy :

$$S_A = -\langle \log \rho_A \rangle = \log D$$

# *E.E. & Edge states (Gapped)* *(of spins, fermions...)*



★ *Partial Trace induces effective edge states*

★ Requirement: Finite Energy Gap for the Bulk

★ *The effective edge states contribute to the E.E.*

★ Let us assume that the edge states has degrees of freedom  $D_E$

*Entanglement Entropy > (# edge states) Log  $D_E$*

S. Ryu & YH, *Phys. Rev. B* 73, 245115 (2006)  
(Fermions)

# EE of the Generic VBS States ( $S=1,2,3,\dots$ )

H. Katsura, T.Hirano & YH, Phys. Rev. B76, 012401 (2007)

T.Hirano & YH, J. Phys. Soc. Jpn. 76, 113601 (2007)

$$H_{VBS} = \sum_{i=1}^N \vec{S}_i \cdot \vec{S}_{i+1} + \alpha H_{\text{extra}}^S, \quad \vec{S}_i^2 = S(S+1)$$

$$H_{\text{extra}}^{S=1} = \sum_i \frac{1}{3} (\vec{S}_i \cdot \vec{S}_{i+1})^2$$

$$H_{\text{extra}}^{S=2} = \sum_i \left( \frac{2}{9} (\vec{S}_i \cdot \vec{S}_{i+1})^2 + \frac{1}{63} (\vec{S}_i \cdot \vec{S}_{i+1})^3 \right)$$

$$|\text{VBS}\rangle = \prod_{j=0}^L (a_j^\dagger b_{j+1}^\dagger - b_j^\dagger a_{j+1}^\dagger)^S |\text{vac}\rangle$$

$$\mathcal{S}_L = -\langle \log \rho \rangle_\rho \rightarrow 2 \log(S+1), \quad (L \rightarrow \infty)$$

Boundary Spins:  $S/2$



S	EE	Effective Boundary spins	Degrees of Freedom
1	2 Log 2	$S_{\text{eff}}=1/2$	$2^2=4$
2	2 Log 3	$S_{\text{eff}}=1$	$3^2=9$
S	2 Log (S+1)	$S_{\text{eff}}=S/2$	$(S+1)^2$

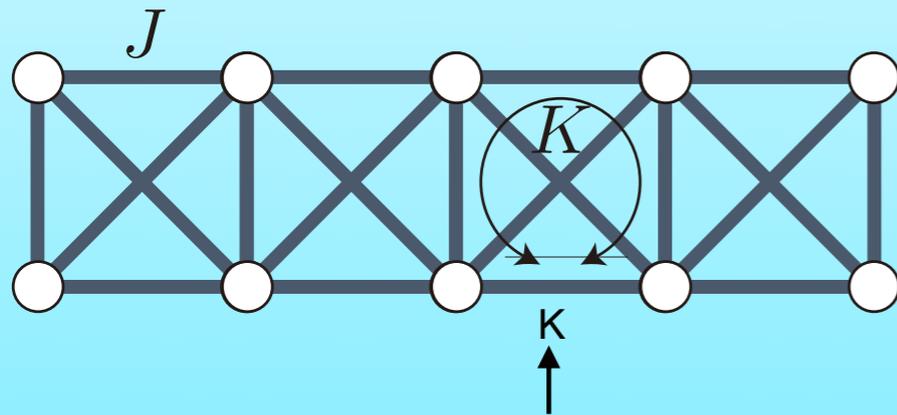
★ **Fractionalization** : Emergent as edge states

(Quantum Resources for qbits)

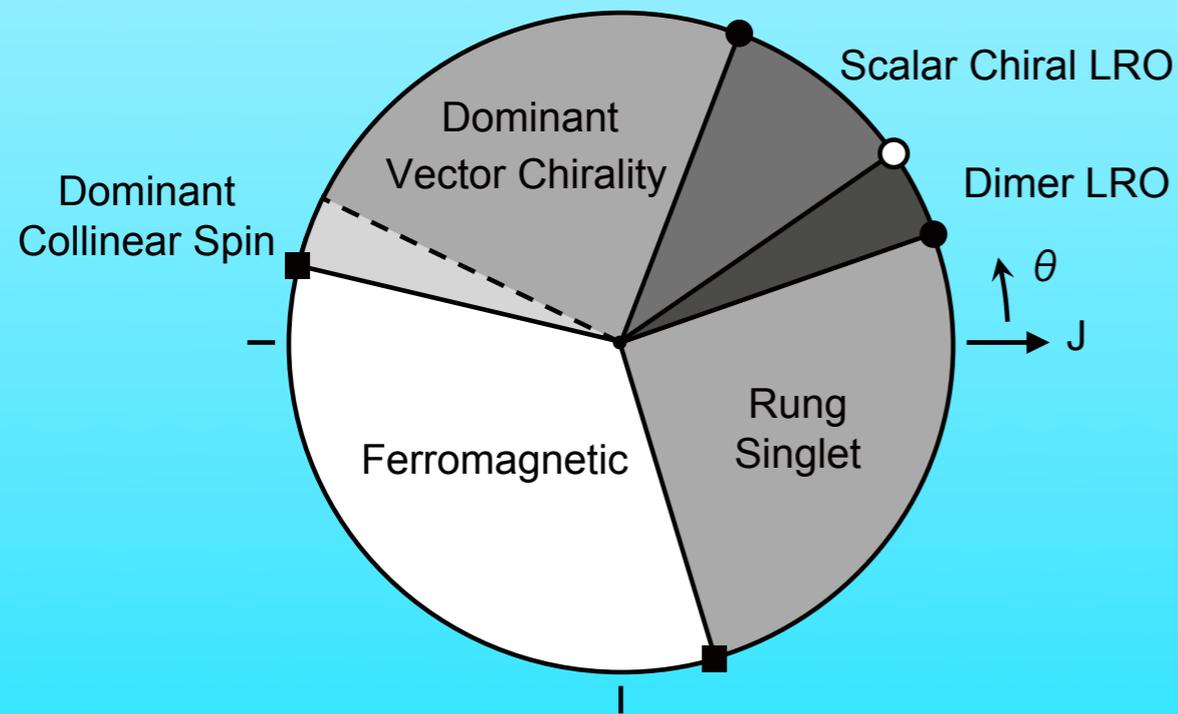
# Another Models

## Spin ladder model with four-spin cyclic exchange

$$\mathcal{H} = \sum_i \{ J_r \mathbf{S}_{1,i} \cdot \mathbf{S}_{2,i} + J_l (\mathbf{S}_{1,i} \cdot \mathbf{S}_{1,i+1} + \mathbf{S}_{2,i} \cdot \mathbf{S}_{2,i+1}) + K (P_i + P_i^{-1}) \}$$



$$(P_i + P_i^{-1}) = \mathbf{S}_{1,i} \cdot \mathbf{S}_{2,i} + \mathbf{S}_{1,i+1} \cdot \mathbf{S}_{2,i+1} + \mathbf{S}_{1,i} \cdot \mathbf{S}_{1,i+1} + \mathbf{S}_{2,i} \cdot \mathbf{S}_{2,i+1} + \mathbf{S}_{1,i} \cdot \mathbf{S}_{2,i+1} + \mathbf{S}_{2,i} \cdot \mathbf{S}_{1,i+1} + 4(\mathbf{S}_{1,i} \cdot \mathbf{S}_{2,i})(\mathbf{S}_{1,i+1} \cdot \mathbf{S}_{2,i+1}) + 4(\mathbf{S}_{1,i} \cdot \mathbf{S}_{1,i+1})(\mathbf{S}_{2,i} \cdot \mathbf{S}_{2,i+1}) - 4(\mathbf{S}_{1,i} \cdot \mathbf{S}_{2,i+1})(\mathbf{S}_{2,i} \cdot \mathbf{S}_{1,i+1}).$$



We set parameters as

$$\begin{cases} J = J_r = J_l = \cos \theta \\ K = \sin \theta \end{cases}$$

Self dual at the point of  $J = 2K$

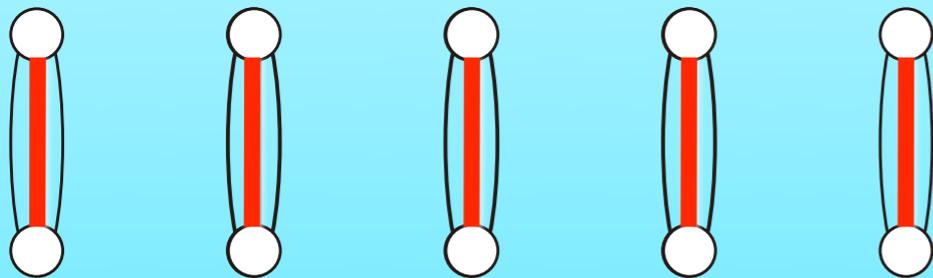
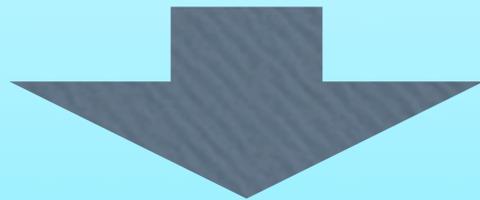
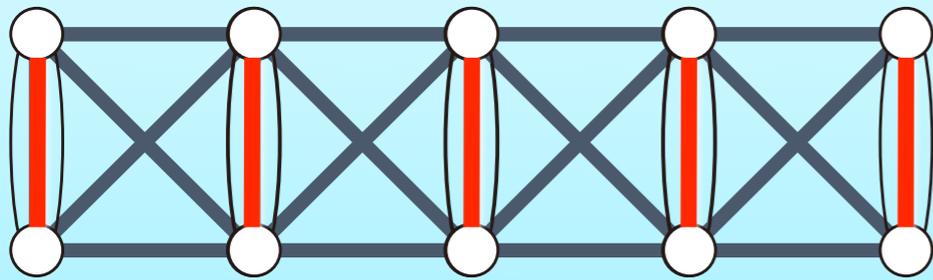
T. Hikihara, T. Momoi and X. Hu (2003)

# Adiabatic deformation

*I. Maruyama, T. Hirano, YH, arXiv:0806.4416*

## Rung singlet phase

$$\theta = 6$$

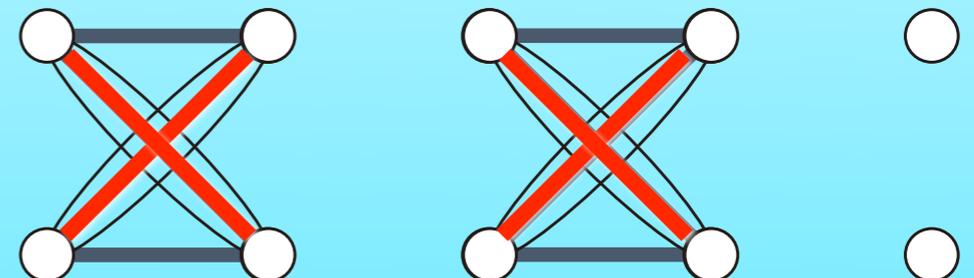
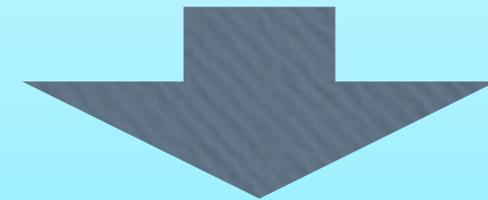
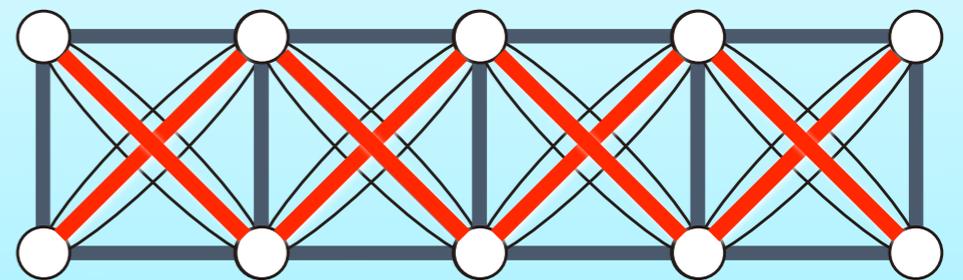


$$H_r = \sum_{i=1} \mathbf{S}_{1,i} \cdot \mathbf{S}_{2,i}$$

Rung singlets

## Vector chirality phase

$$\theta = 2.6$$



$$H_{ps} = \sum_{i \in \text{odd}} (\mathbf{S}_{1,i} \times \mathbf{S}_{2,i}) \cdot (\mathbf{S}_{1,i+1} \times \mathbf{S}_{2,i+1})$$

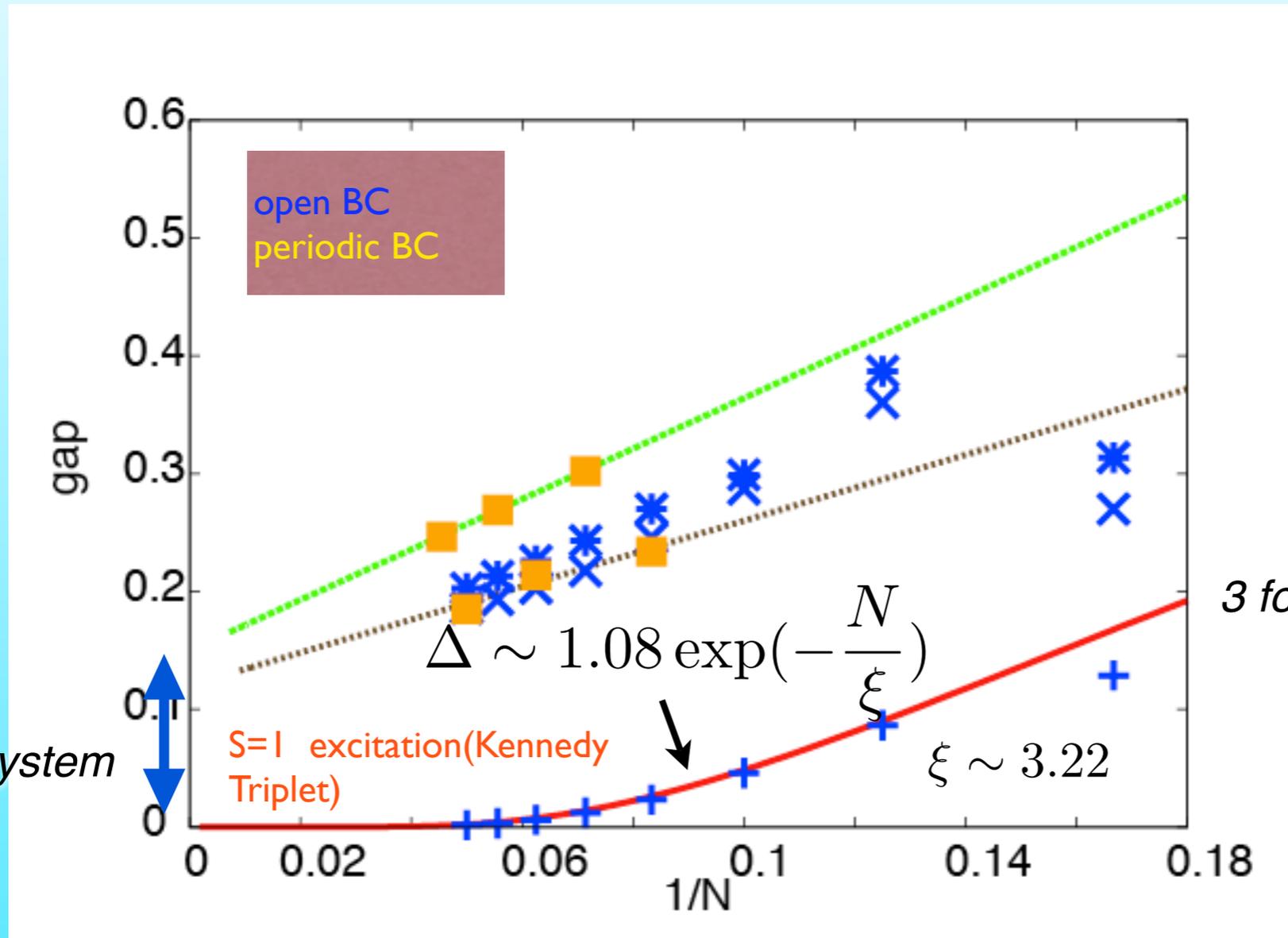
Plaquette singlet (PS)

Berry phase remains the same

Topologically equivalence

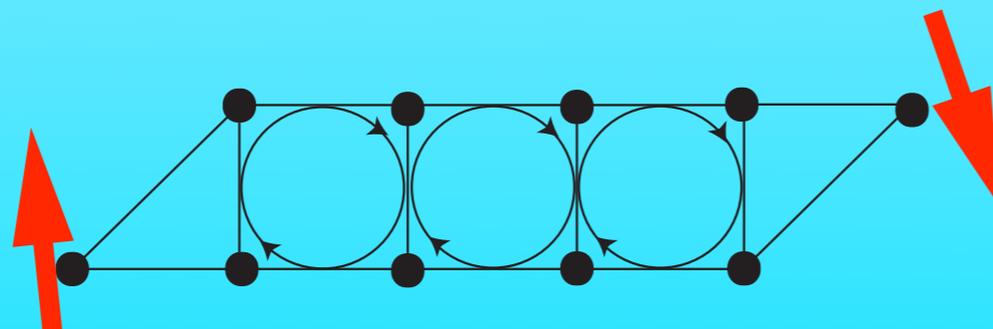
# Energy spectrum with boundaries (diagonal)

*M. Arikawa, S. Tanaya, I. Maruyama, YH, unpublished*



*Ex. for Haldane spin chain*

*Kennedy '90*



*Interaction between effective boundary spins*

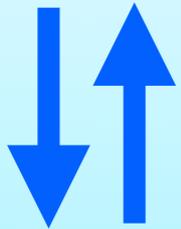
$$H_{eff} = \Delta \mathbf{S}_R \cdot \mathbf{S}_L$$

# Bulk-Edge correspondence for spins

$$\Theta_N^2 = 1 \quad \Theta_N = (i\sigma_y^1) \otimes (i\sigma_y^2) \cdots (i\sigma_y^N) K$$

$K$  : complex conjugate

**Bulk:  $Z_2$  Berry phases**



**Edge: Entanglement Entropy  
& low energy states in the gap**

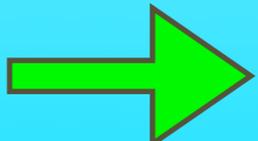
$S = 1/2$  is always fundamental ( electron spin )



$$\Theta_{edges} = \Theta_L \otimes \Theta_R$$

$$\Theta_R^2 = -1$$

$$\Theta_L^2 = -1$$

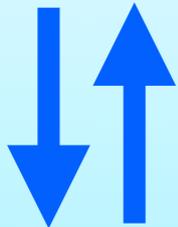
**Global TR**  $\Theta_N$   **Local (edge) TR**  $\Theta_L, \Theta_R$

# Bulk-Edge correspondence for spins

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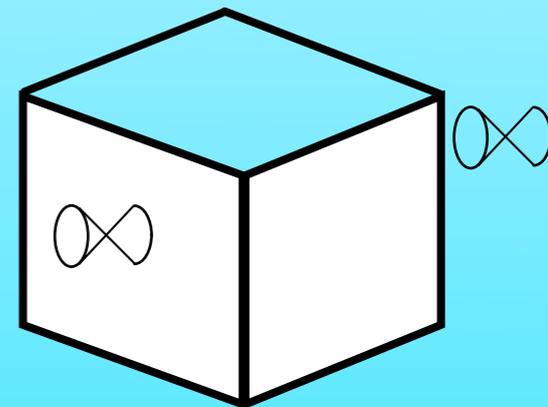
**Bulk:  $Z_2$  Berry phases**



**Edge: Entanglement Entropy  
& low energy states in the gap**

$S = 1/2$  is always fundamental ( electron spin )

**3D ?**



$$\Theta_{edges} = \Theta_L \otimes \Theta_R \quad \Theta_R^2 = -1 \quad \Theta_L^2 = -1$$

**Global TR**  $\Theta_N$   $\longrightarrow$  **Local (edge) TR**  $\Theta_L, \Theta_R$

***Thank you***