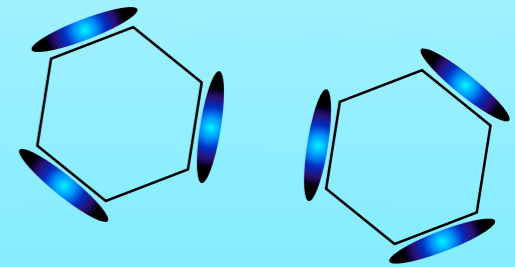


Bulk-Edge correspondence and Fractionalization

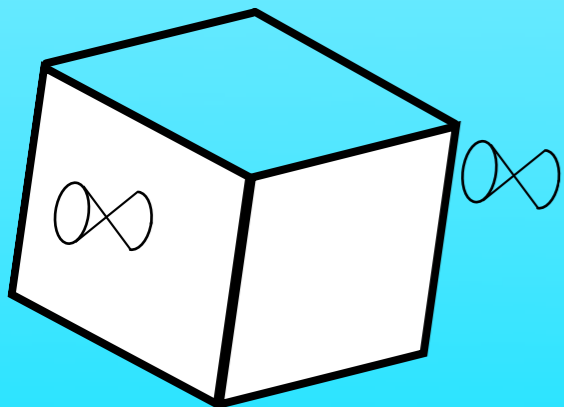
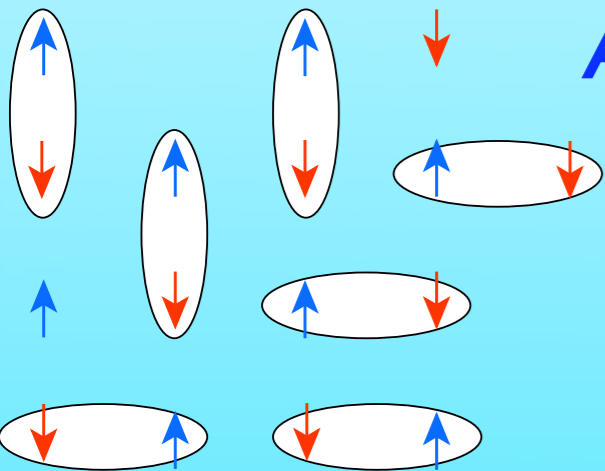
As a topological (spin) insulator
with strong interaction



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

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 ?

How to understand gapped quantum liquids ?

Bulk

classically featureless : need geometrical phase

How to understand gapped quantum liquids ?

Bulk

classically featureless : need geometrical phase

1-st Chern number for QHE TKNN

How to understand gapped quantum liquids ?

Bulk

classically featureless : need geometrical phase

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Edge

low energy localized modes in the gap

How to understand gapped quantum liquids ?

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low energy localized modes in the gap

edge states for QHE

Laughlin, Halperin, YH

How to understand gapped quantum liquids ?

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How to understand gapped quantum liquids ?

Bulk-Edge correspondence

Common property of topological ordered states

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classically featureless : need geometrical phase



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How to understand gapped quantum liquids ?

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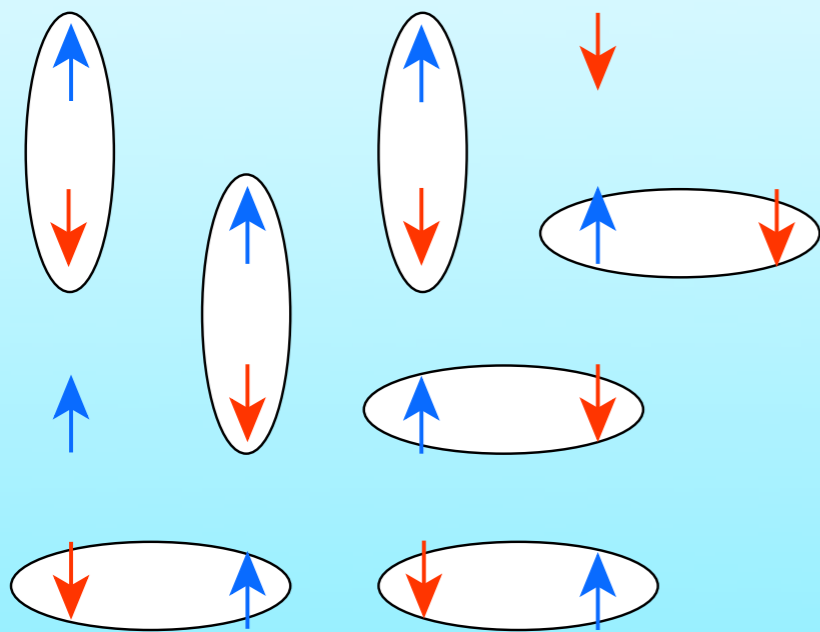
Laughlin, Halperin, YH

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

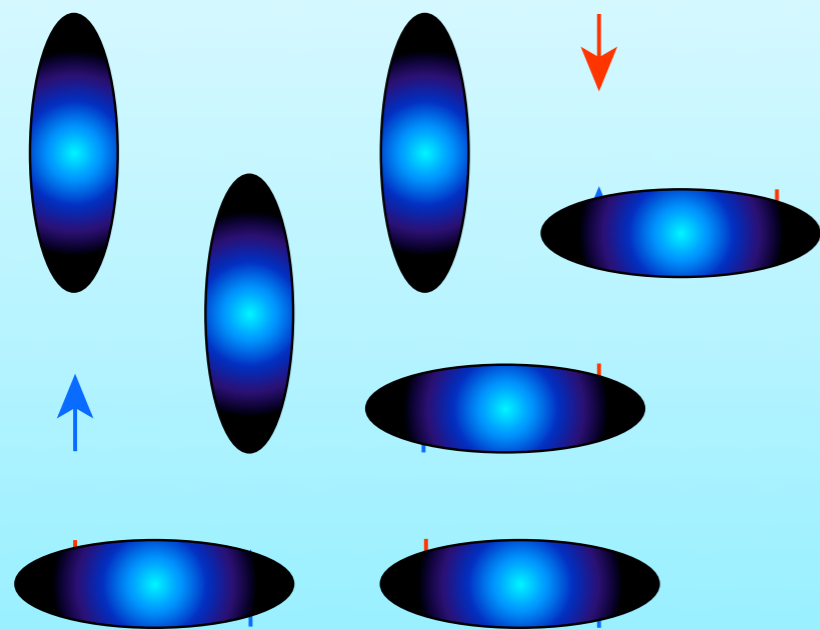


Local Singlet Pairs :
(Basic Objects)

Purely Quantum Objects are basic

Quantum Liquid (Example 1)

★ The *RVB* state by Anderson



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Spins disappear
as a *Singlet pair*

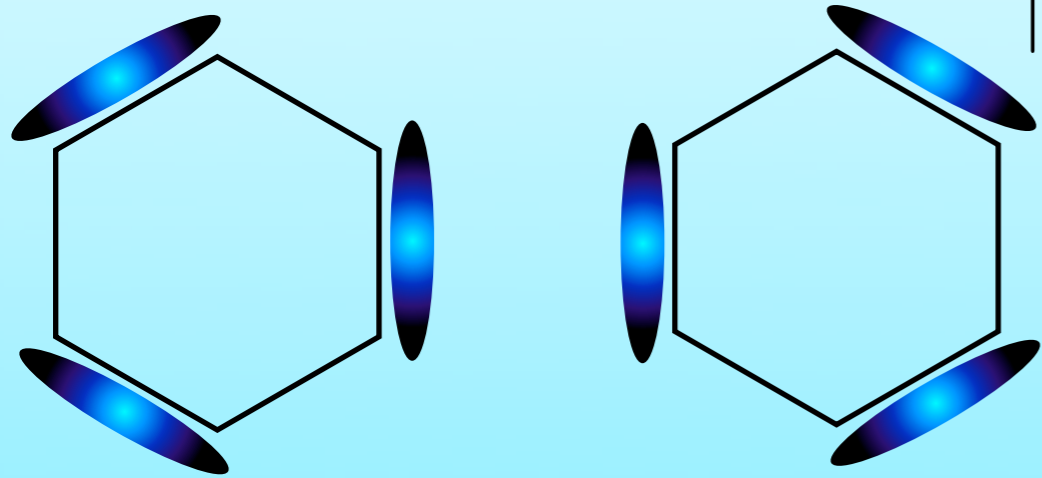


Local Singlet Pairs :
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Purely Quantum Objects are basic

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*

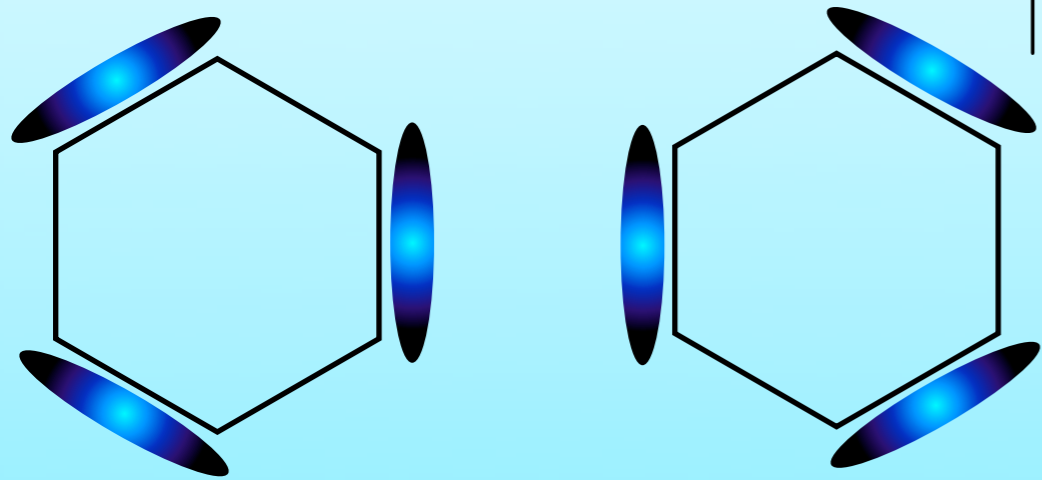
A

Local Covalent Bonds :
(Basic Objects)

Purely Quantum Objects are basic

Quantum Liquid (Example 2)

★ The *RVB* state by Pauling



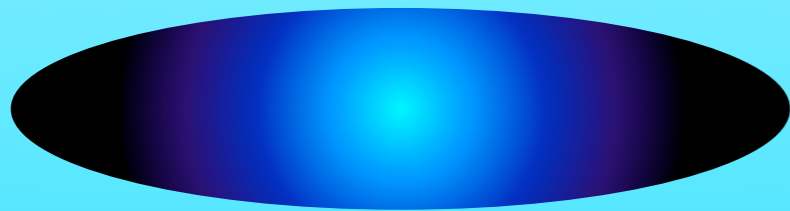
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$$|G\rangle = \sum_{J=\text{Dimer Covering}} c_J \otimes_{ij} |\text{Bond}_{ij}\rangle$$

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

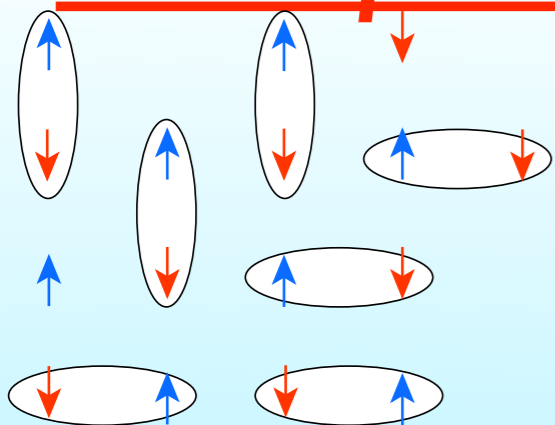
$$\langle G | G + dG \rangle = 1 + \langle G | dG \rangle$$

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

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

Use Quantum Interferences To Classify Quantum Liquids

Examples: RVB state by Anderson

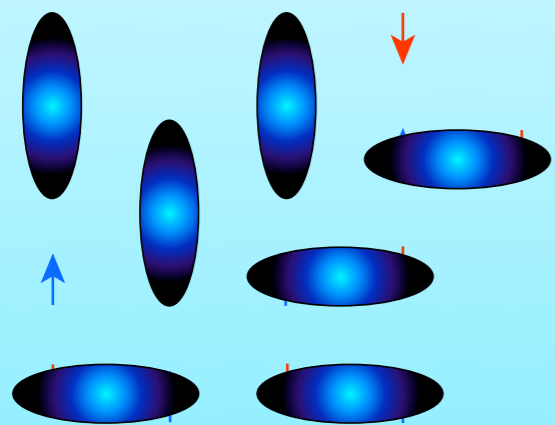


$$|\text{Singlet Pair}_{12}\rangle = \frac{1}{\sqrt{2}} (|\uparrow_1\downarrow_2\rangle - |\downarrow_1\uparrow_2\rangle)$$

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Spins *disappear*
as a *Singlet pair*

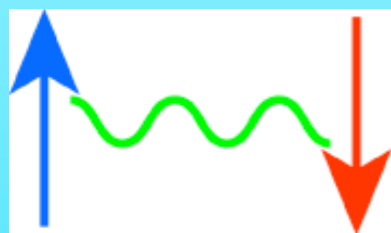
No Long Range Order
in Spin-Spin Correlation



Local Singlet Pair is a Basic Object

How to Characterize the Local Singlet Pair ?

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

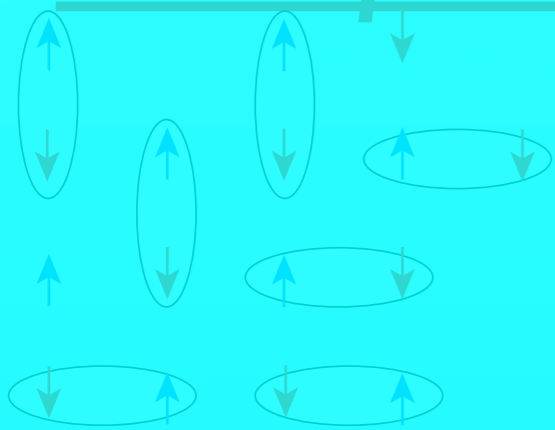


Use Berry Phase to characterize the Singlet!

Singlet does not carries spin but does Berry phase

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

Examples: RVB state by Anderson

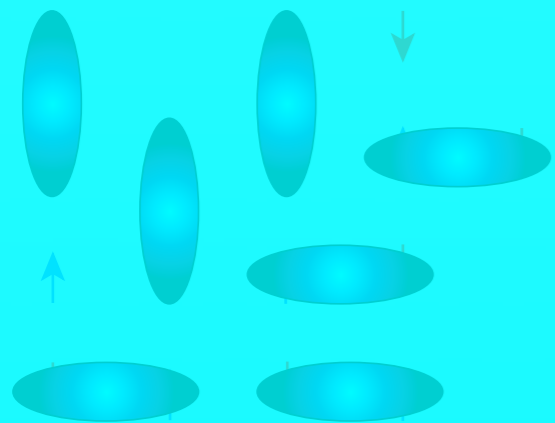


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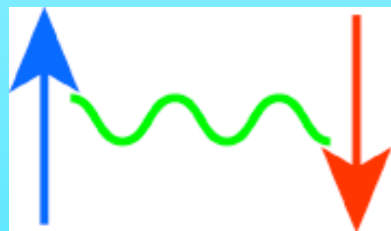
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Use Berry Phase to characterize the Singlet!

Singlet does **not** carries **spin** but does Berry phase

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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

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

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

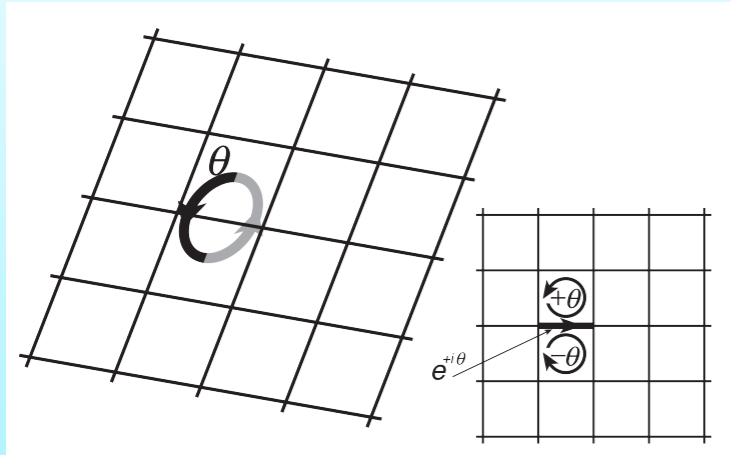
$$\Theta_N = (i\sigma_y^1) \otimes (i\sigma_y^2) \cdots (i\sigma_y^N) K$$

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

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!

$$\gamma_C = \int_C A_\psi = \int_C \langle \psi | d\psi \rangle = \begin{cases} 0 \\ \pi \end{cases} : \text{mod } 2\pi \quad Z_2$$

Time Reversal (Anti-Unitary) Invariance

Berry Connection and Gauge Transformation

★ **Parameter Dependent Hamiltonian**

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

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

★ **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**

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

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 2π**

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

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**

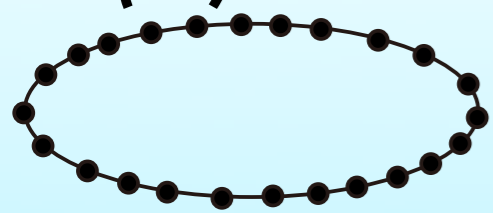
$$\gamma_C(A^\Psi) = \begin{cases} 0 \\ \pi \end{cases} \pmod{2\pi}$$

\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

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

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

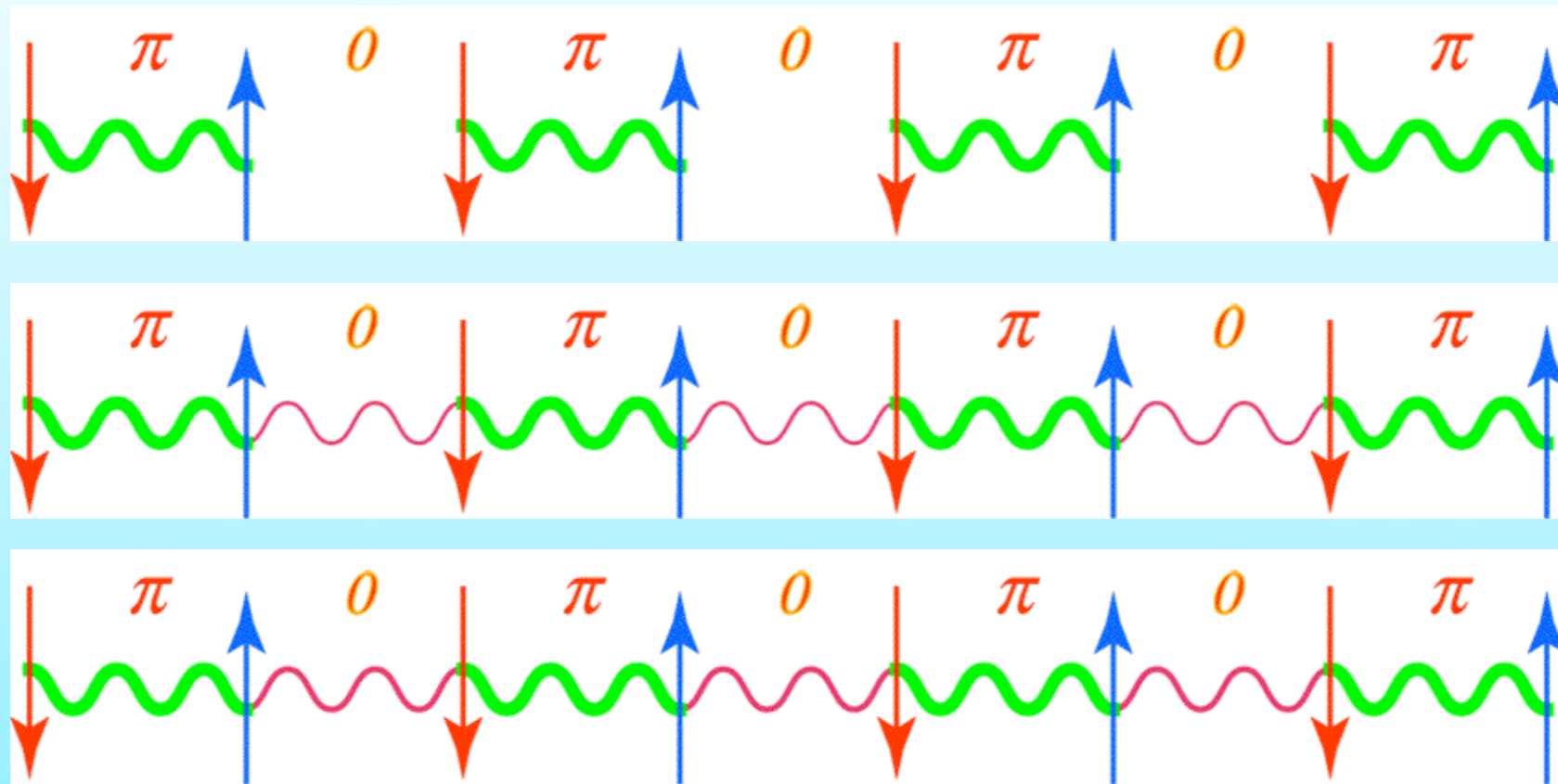
King-Smith & Vanderbilt '93 (polarization in solids)

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

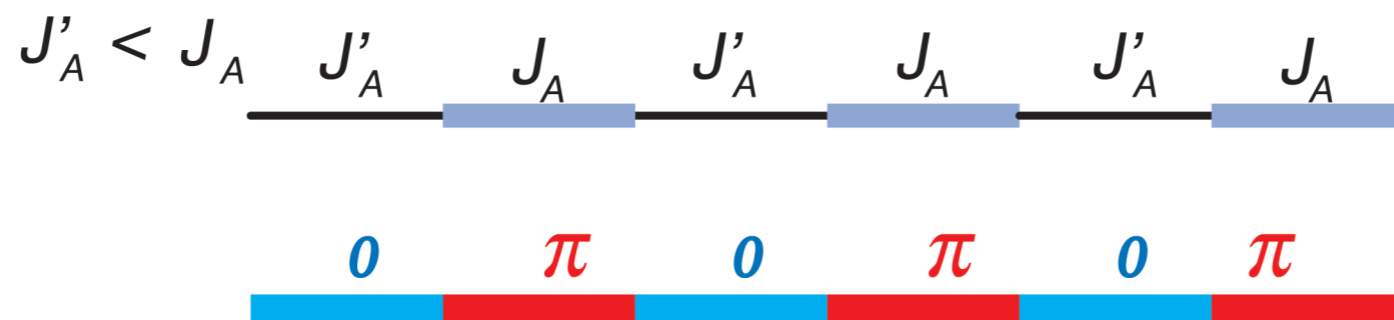
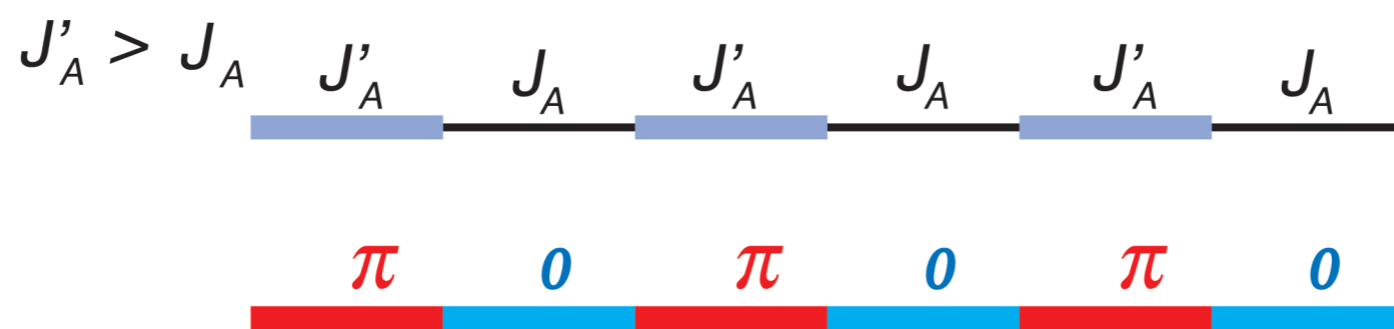
Local Order Parameters of Singlet Pairs

★ 1D AF-AF, AF-F Dimers

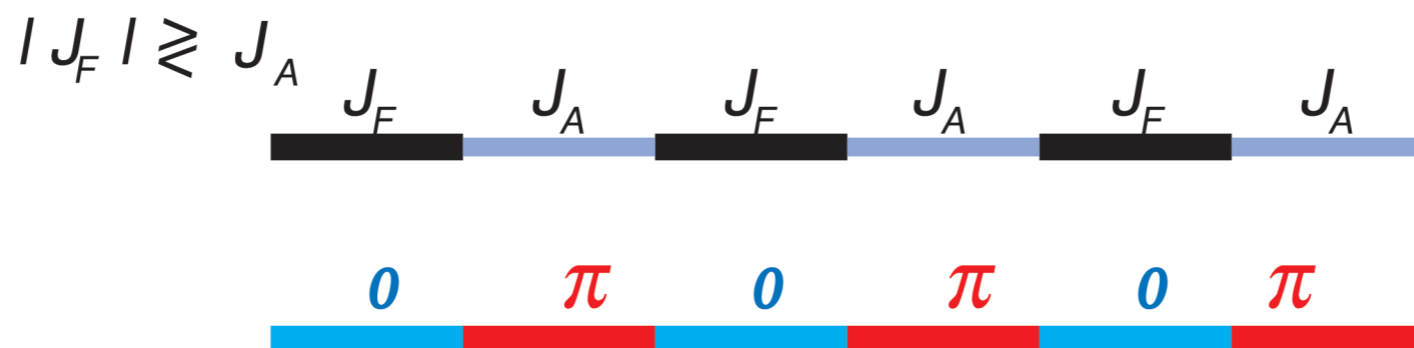
Y.H., J. Phys. Soc. Jpn. 75 123601 (2006)

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

AF-AF



F-AF



AF-AF case

Strong bonds

: π bonds

F-AF case

AF bonds

: π bonds

Local Order Parameters of the Haldane Phase

★ Heisenberg Spin Chains with integer S

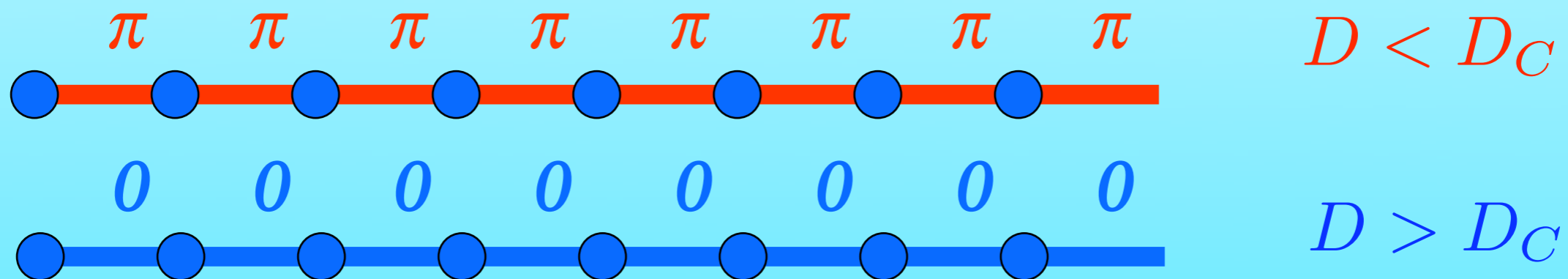
★ No Symmetry Breaking by the Local Order Parameter

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

$$S=1 \quad (\mathbf{S}_i)^2 = S(S+1), \quad S = 1$$

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

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

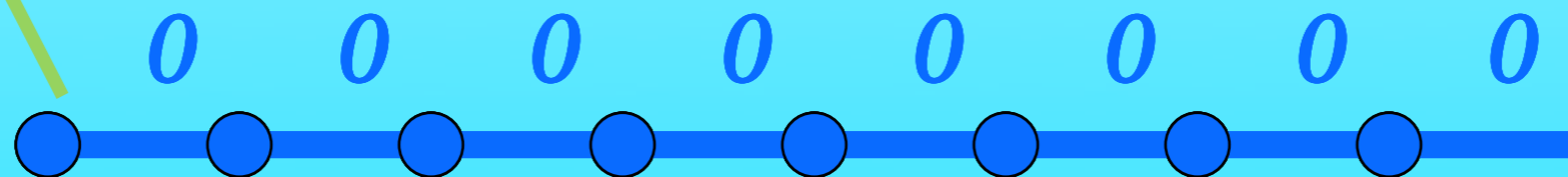
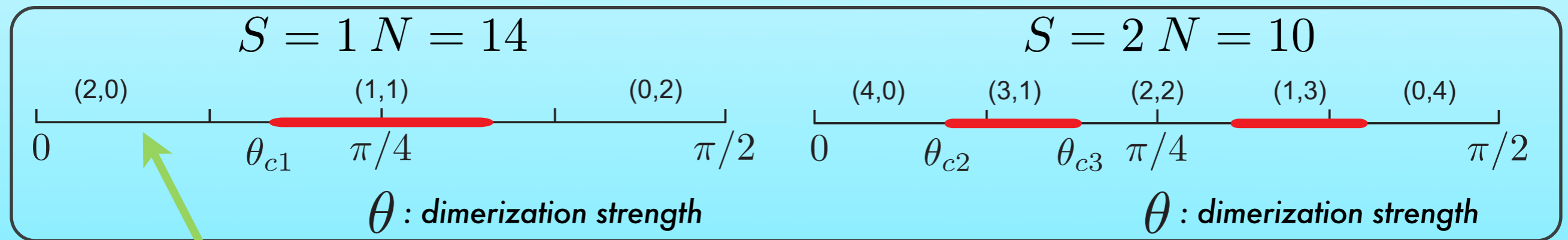
Topological Classification of Gapped Spin Chains

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



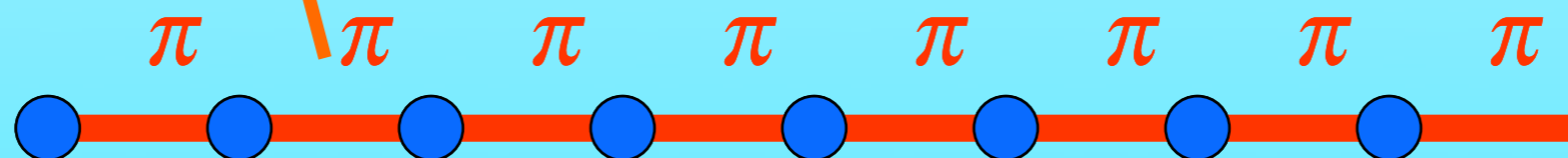
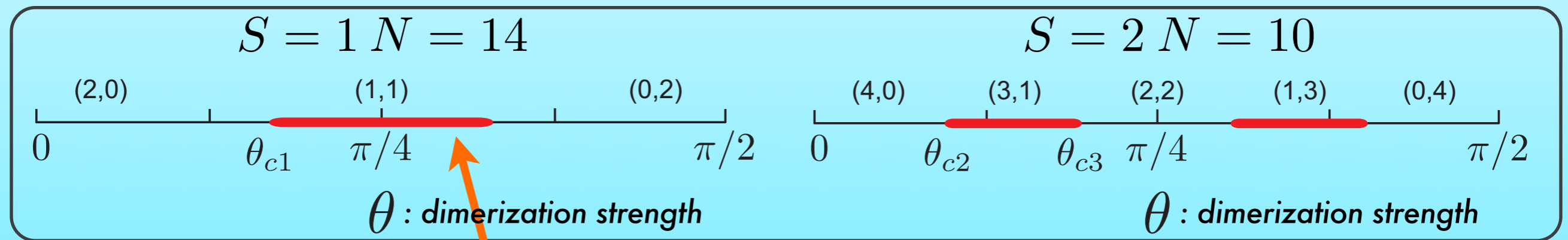
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Z₂Berry phase



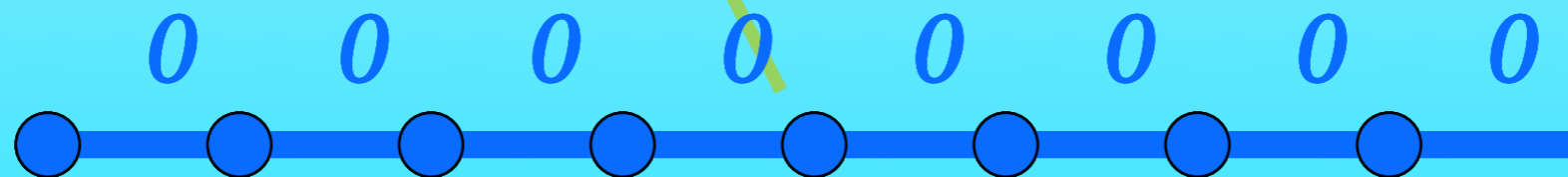
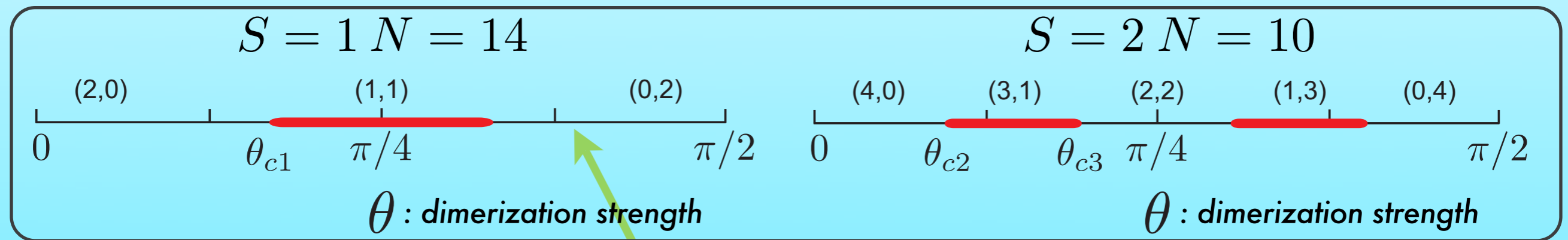
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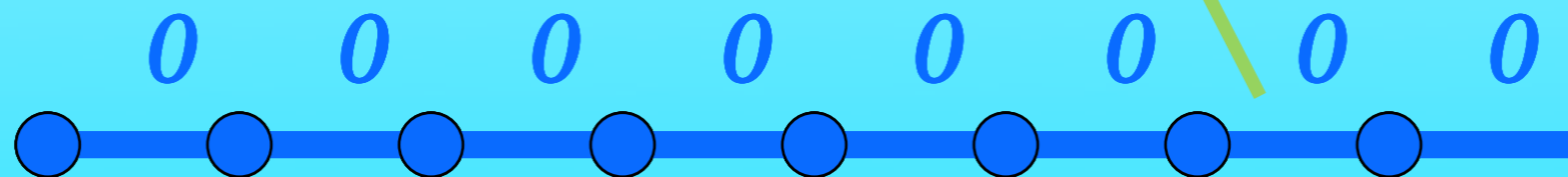
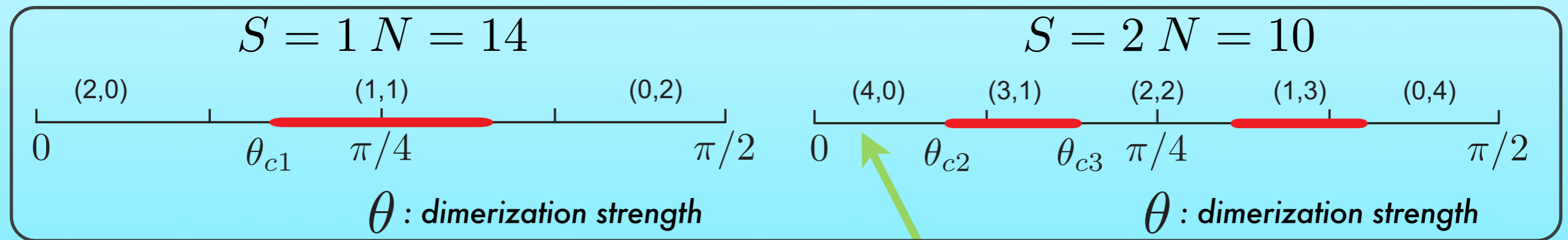
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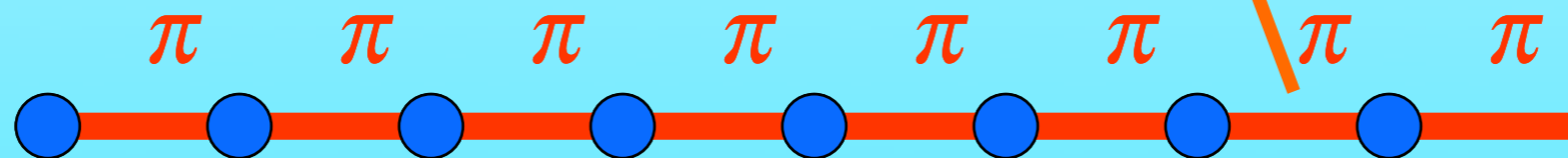
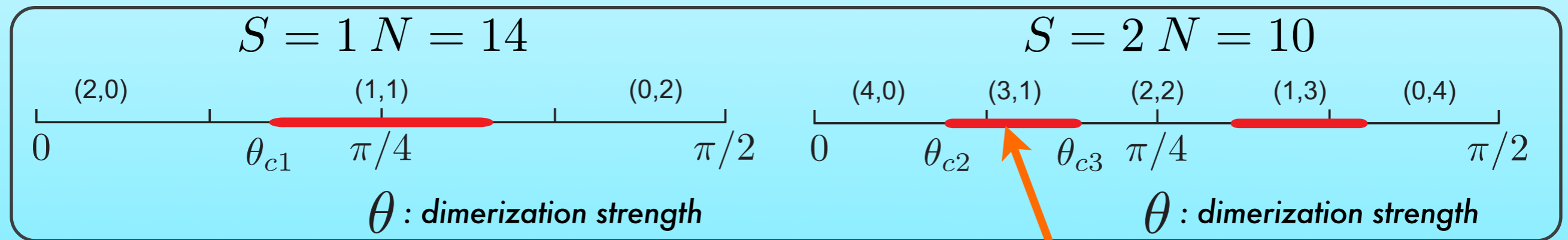
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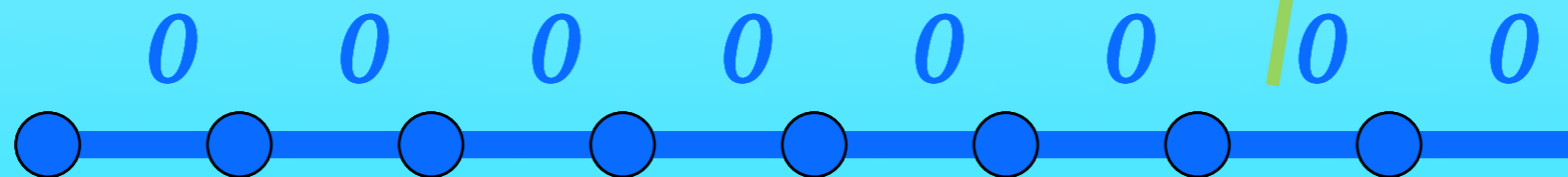
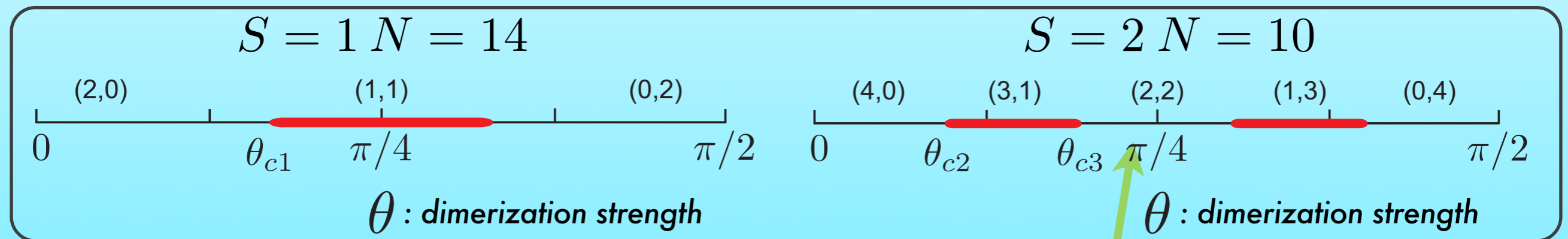
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Z_2 Berry phase



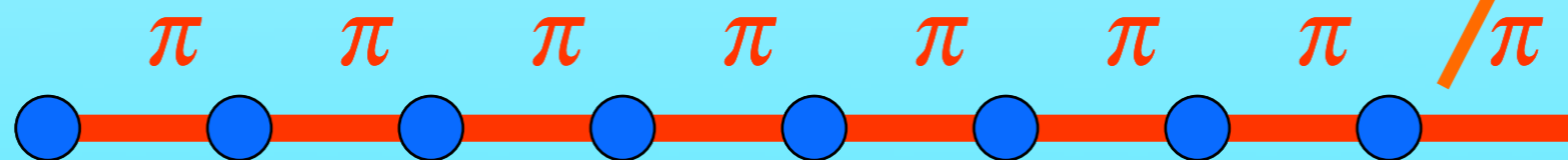
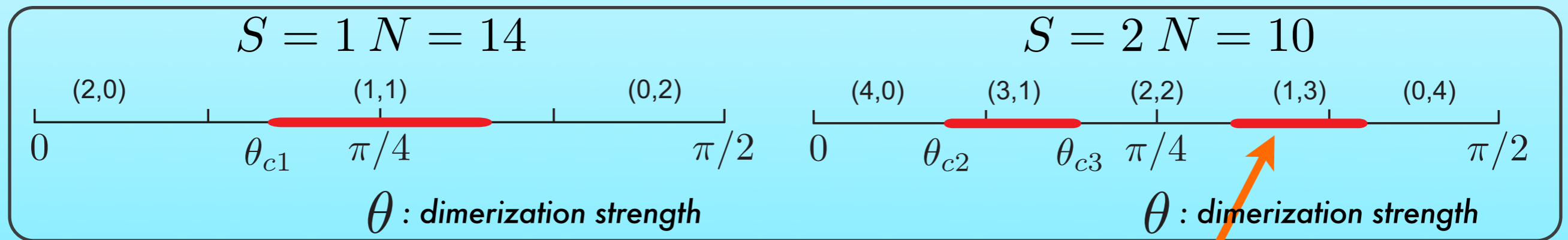
Topological Classification of Gapped Spin Chains

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

- ♦ S=1,2 dimerized Heisenberg model

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

Z₂Berry phase



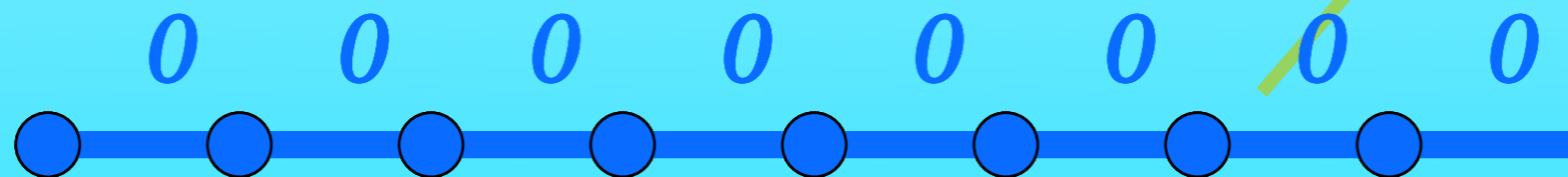
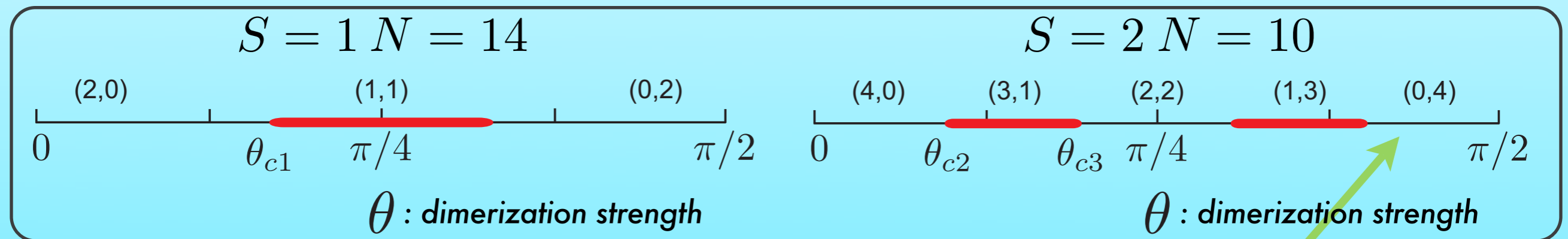
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T.Hirano, H.Katsura & YH, Phys.Rev.B77 094431'08

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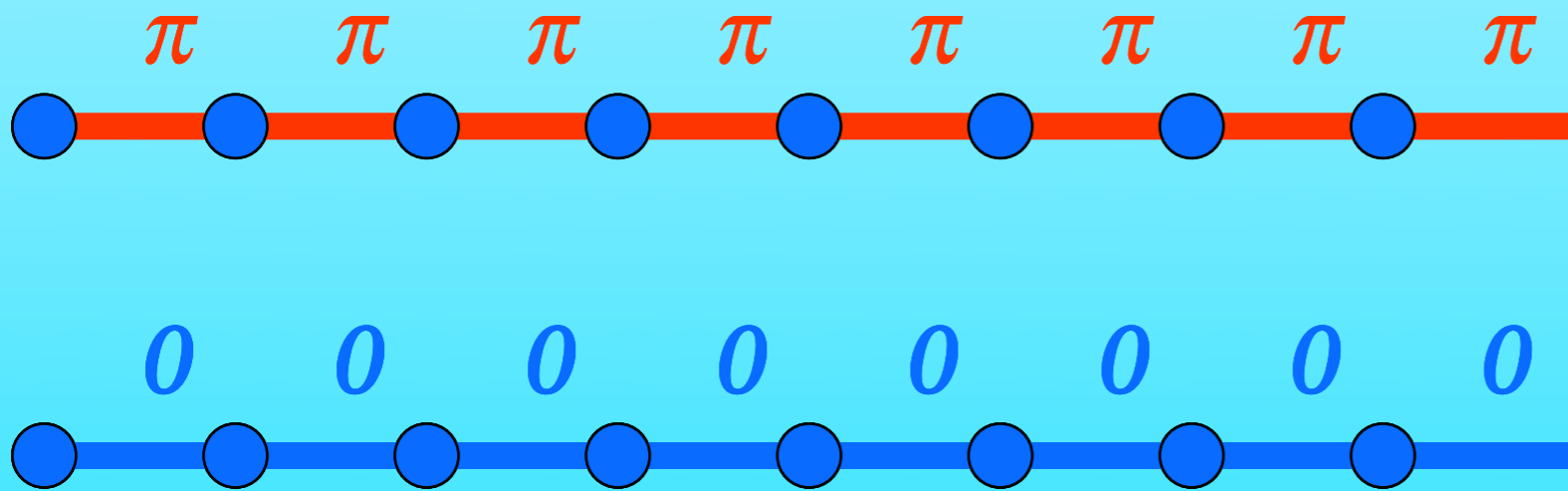
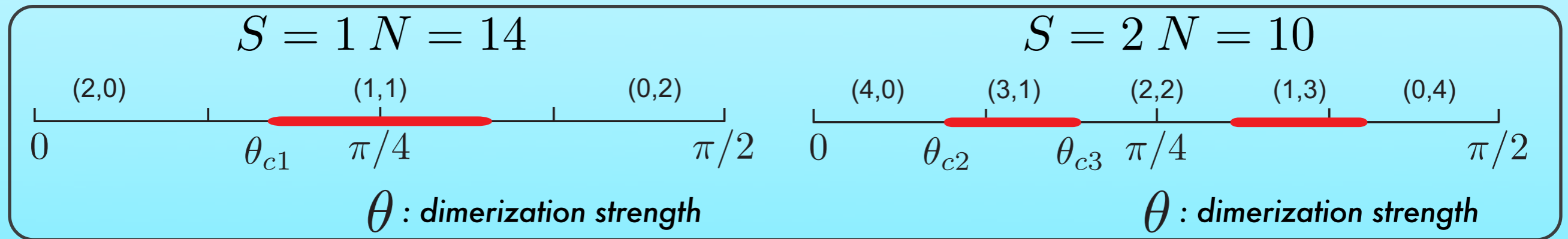
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Z₂Berry phase



Topological Quantum Phase Transitions with **translation** invariance

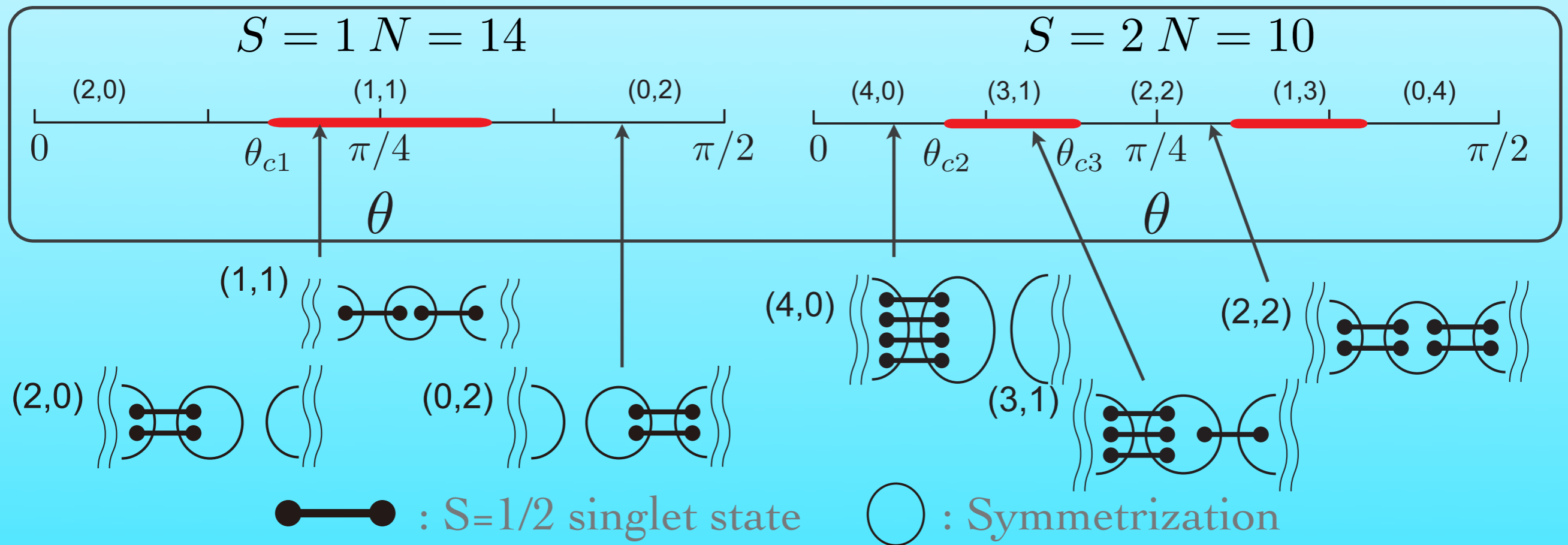
Topological Classification of Gapped Spin Chains

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.)

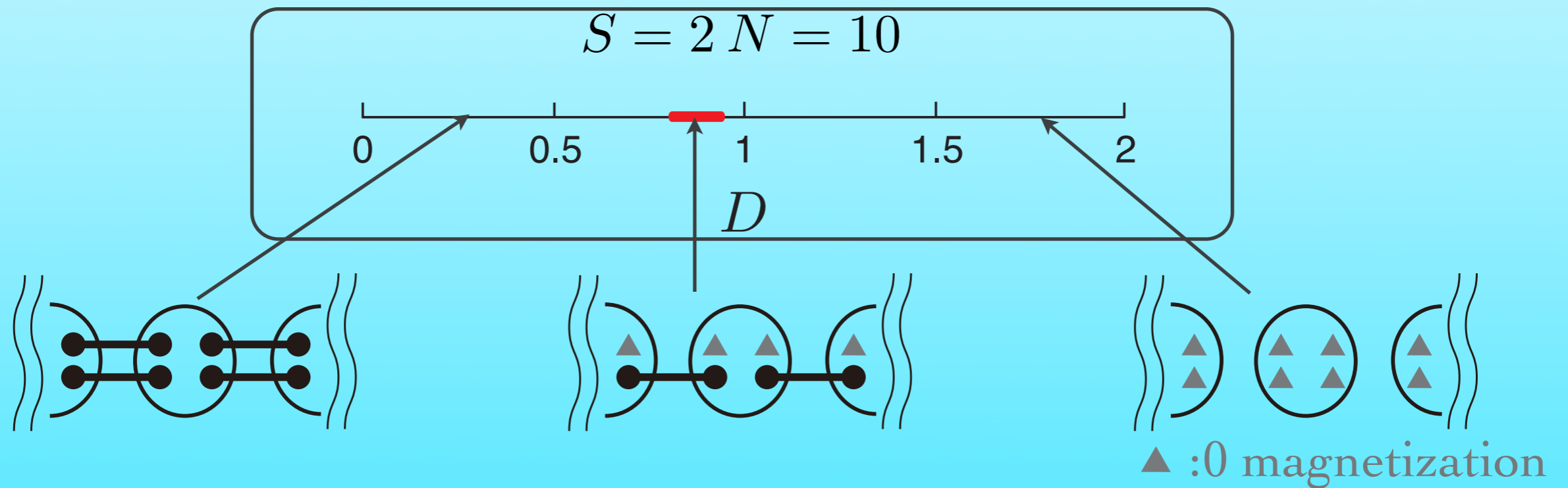
T.Hirano, H.Katsura & YH, Phys.Rev.B77 094431'08

- ♦ 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

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

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

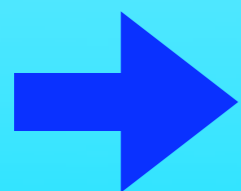
Berry phase on a link (ij)

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

$S=1/2$

The Berry phase counts the number of the valence bonds!

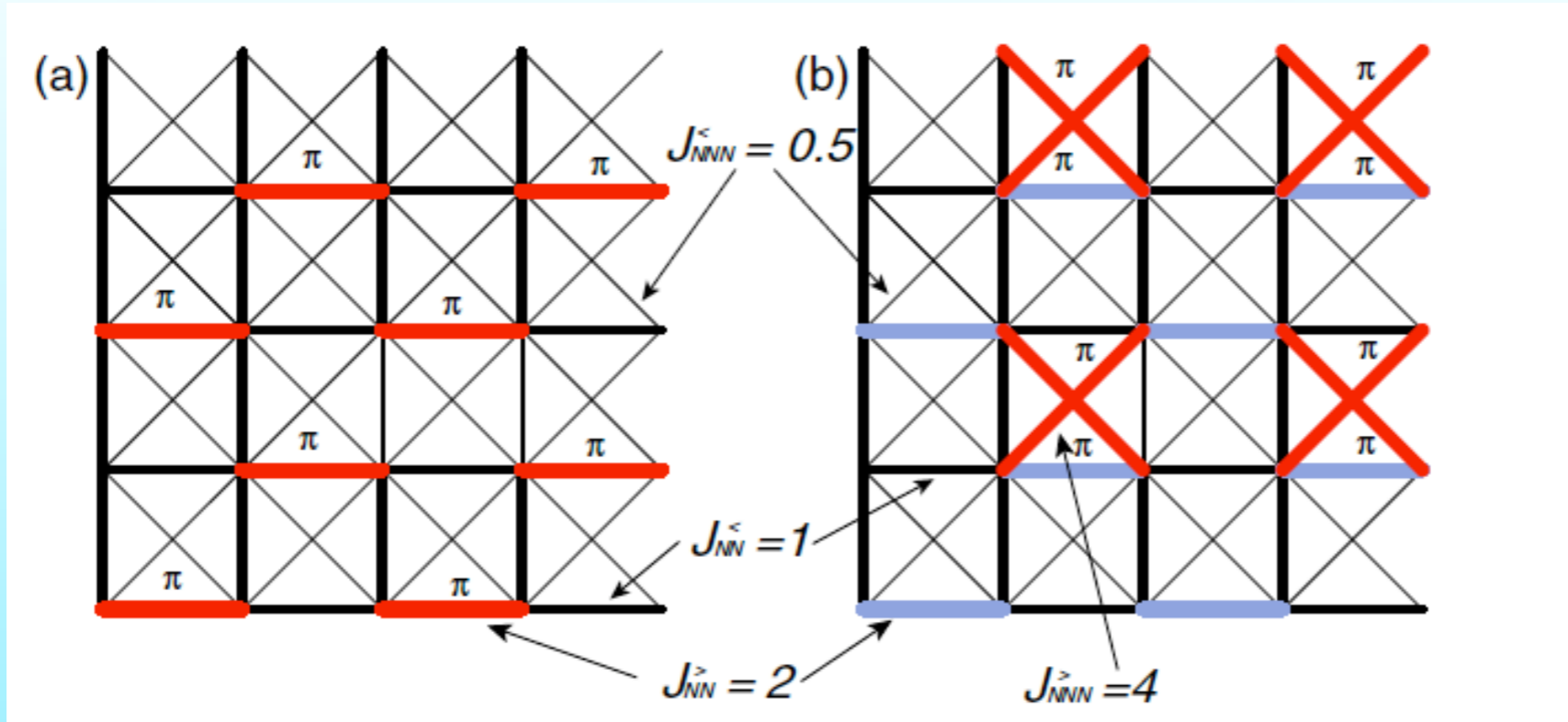
$S=1/2$ objects are fundamental in $S=1$ & 2 spin chains



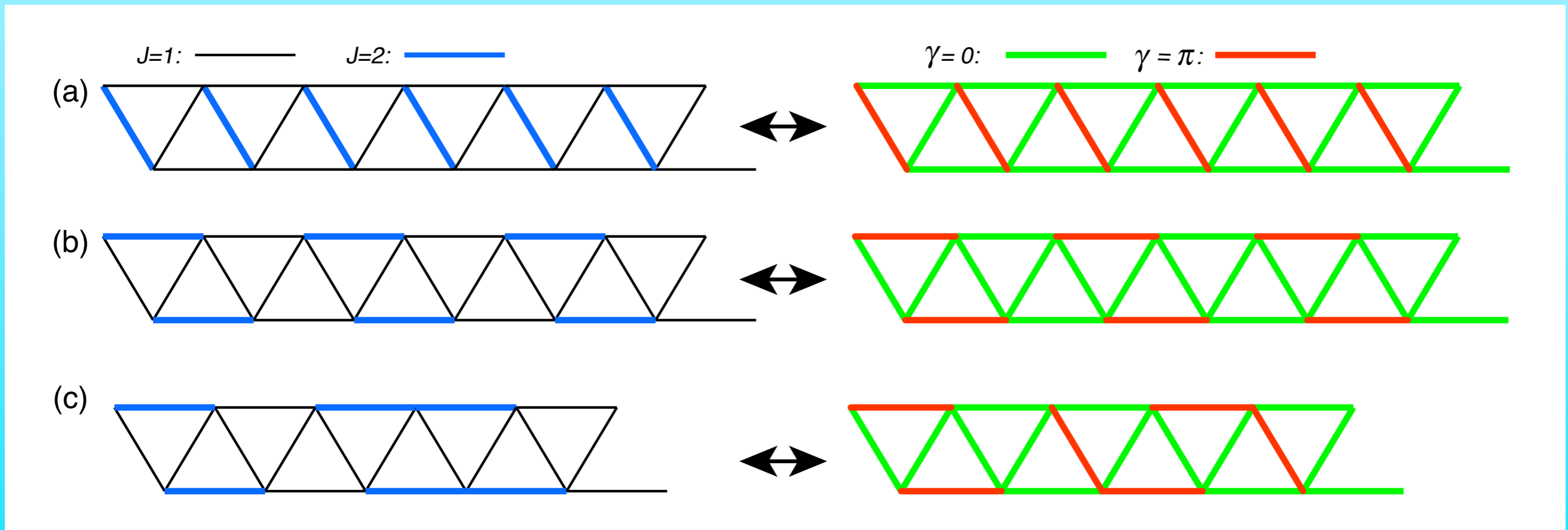
FRACTIONALIZATION

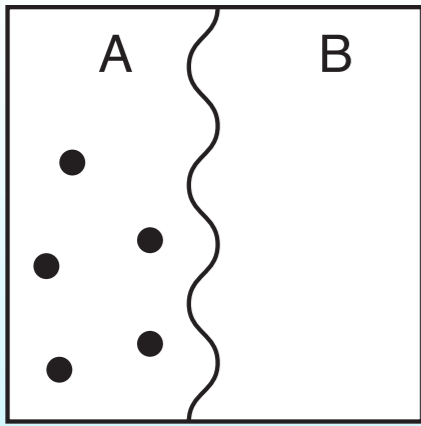
Contribute to the
Entanglement Entropy
as of Edge states

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

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

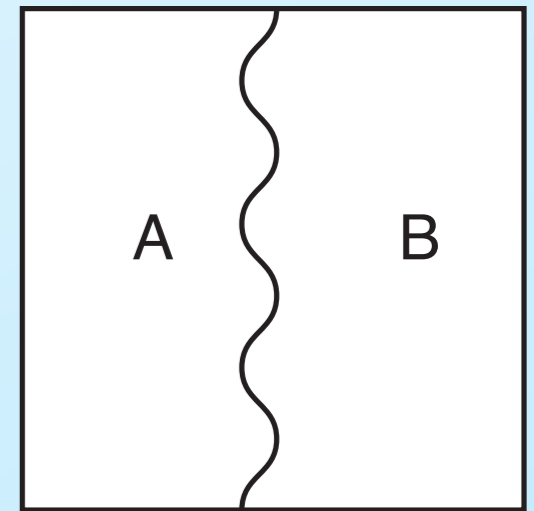
System = $A \oplus B$

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

★ 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

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

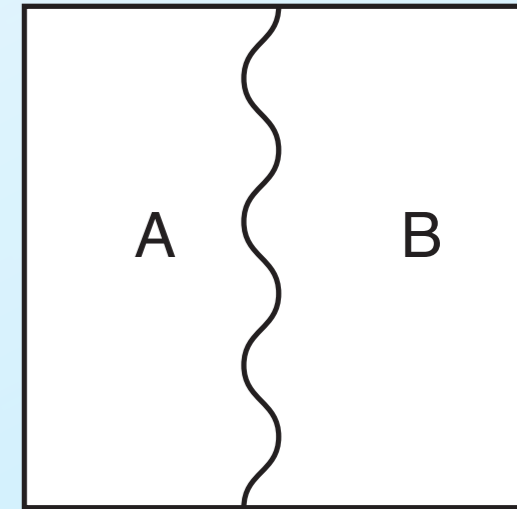
★ How much the State is Entangled between A & B?

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)	$S_{\text{eff}}=S/2$	$(S+1)^2$

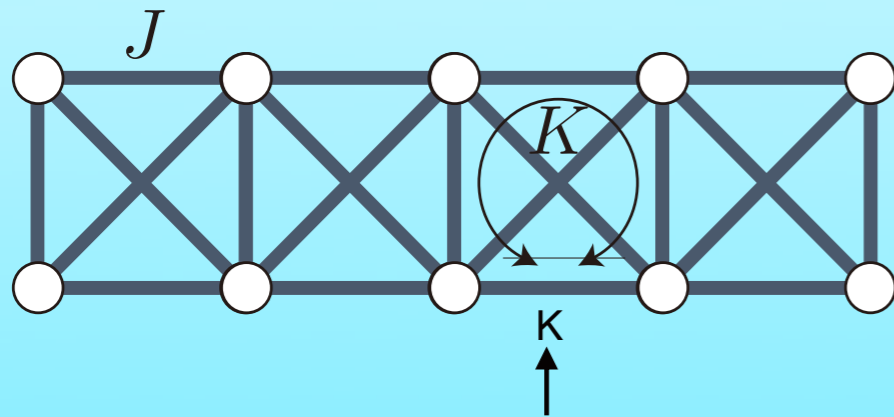
★ **Fractionalization** : Emergent as edge states

(Quantum Resources for qbits)

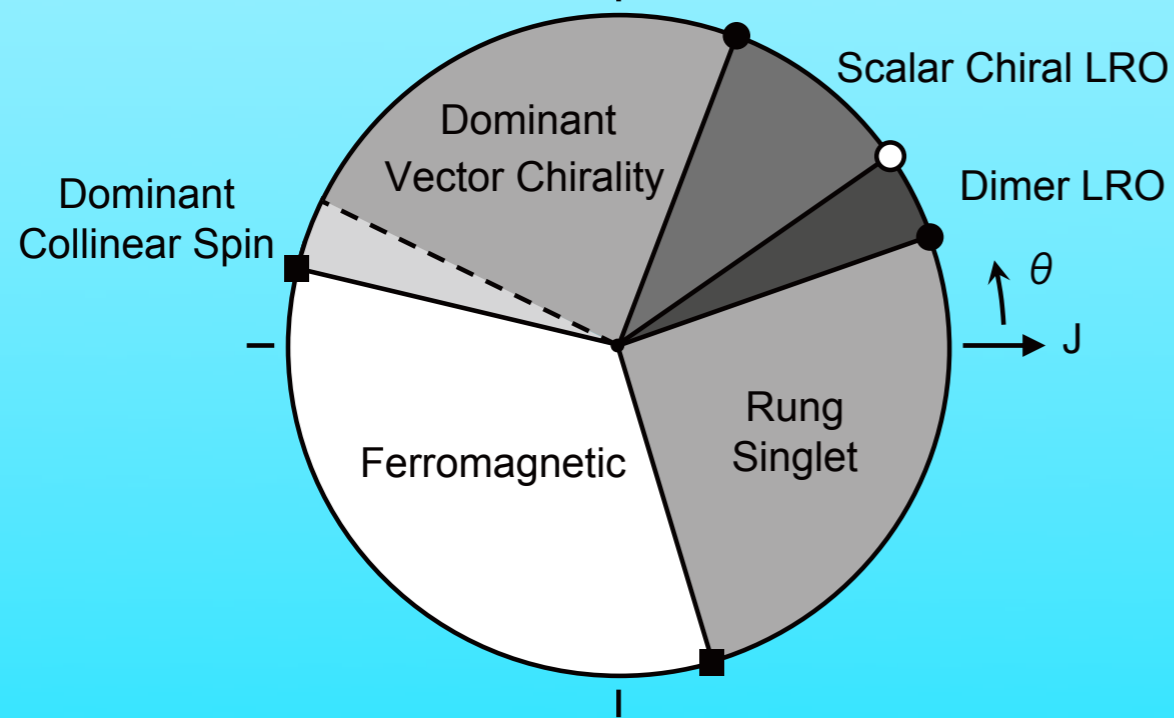
Another Models

Spin ladder model with four-spin cyclic exchange

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



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



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)

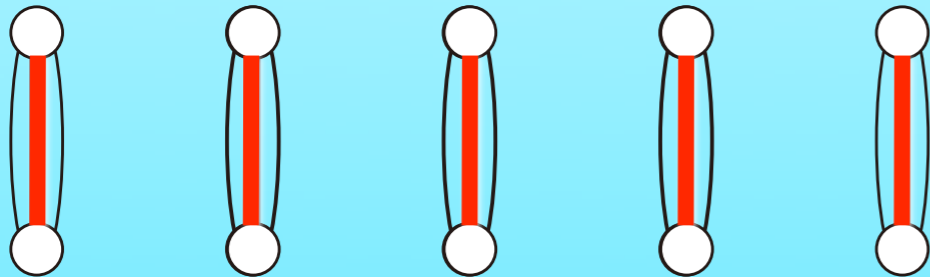
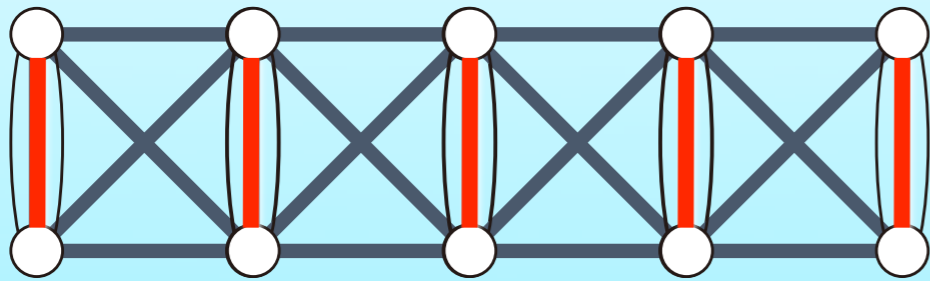
A. Lauchli, G. Schmid and M. Troyer (2003)

Adiabatic deformation

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

Rung singlet phase

$$\theta = 6$$

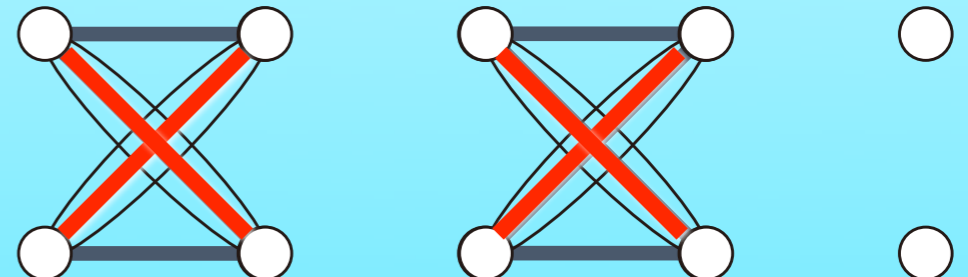
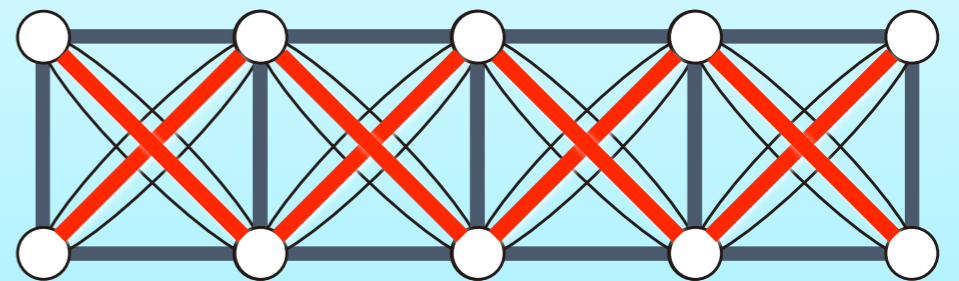


$$H_r = \sum_{i=1} \mathbf{S}_{1,i} \cdot \mathbf{S}_{2,i}$$

Rung singlets

Vector chirality phase

$$\theta = 2.6$$



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

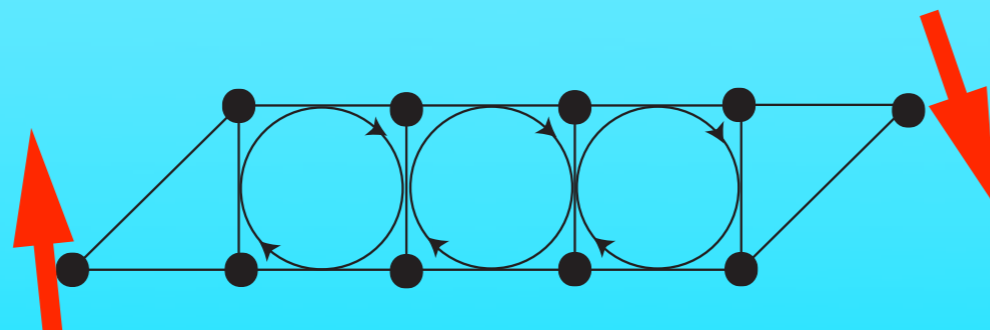
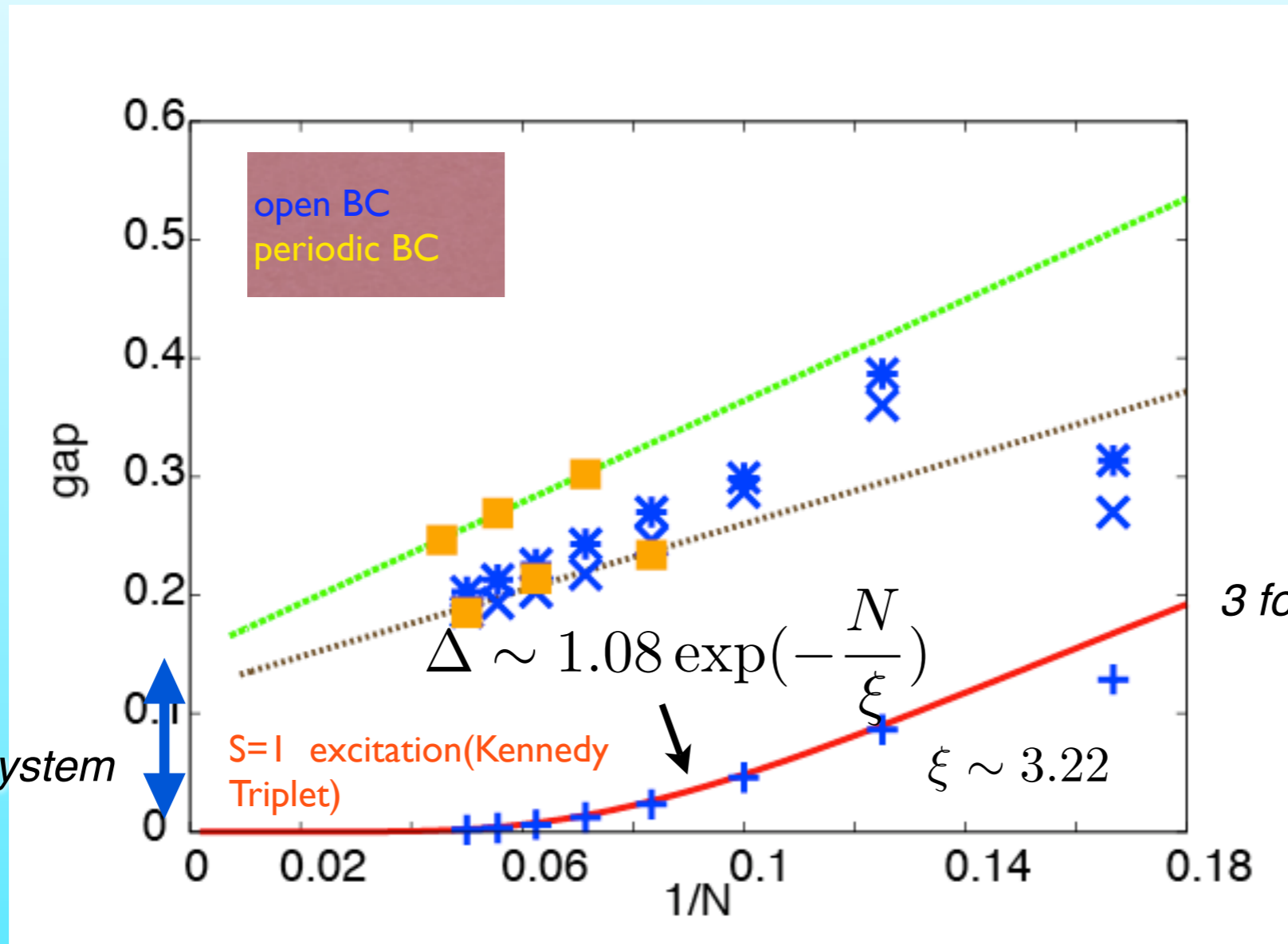
Plaquette singlet (PS)

Berry phase remains the same

Topologically equivalence

Energy spectrum with boundaries (diagonal)

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



Interaction between effective boundary spins

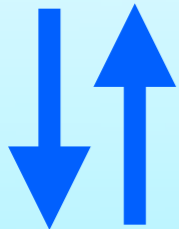
$$H_{eff} = \Delta \mathbf{S}_R \cdot \mathbf{S}_L$$

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)



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

$$\Theta_R^2 = -1$$

$$\Theta_L^2 = -1$$

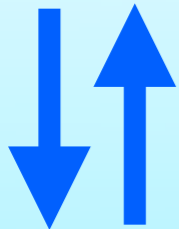
Global TR Θ_N **Local (edge) TR** Θ_L, Θ_R

Bulk-Edge correspondence for spins

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K : complex conjugate

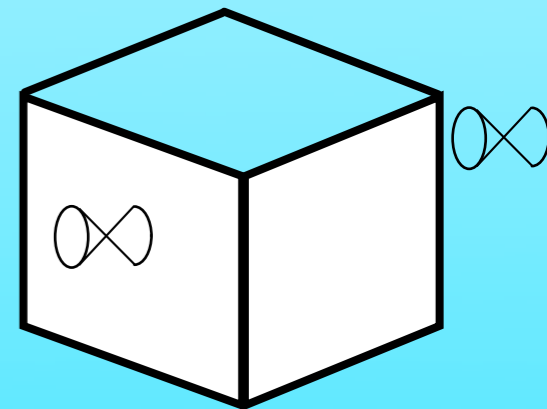
Bulk: Z_2 Berry phases



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

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

3D ?



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

Global TR Θ_N **Local (edge) TR** Θ_L, Θ_R

Thank you