

Determination of **confinement-deconfinement transition** by **Roberge-Weiss periodicity**

Kouji Kashiwa



Collaborator: Akira Ohnishi

- **K.K.**, and A. Ohnishi, Phys. Lett. B750 (2015) 282
- **K.K.**, and A. Ohnishi, Phys. Rev D. 93 (2016) 116002
- **K.K.**, and A. Ohnishi, arXiv:1701.04953

Purpose of this study

To determine

the **confinement-deconfinement transition**

in the system with dynamical quarks



Heavy quark-mass limit

Polyakov-loop describes the confinement-deconfinement transition

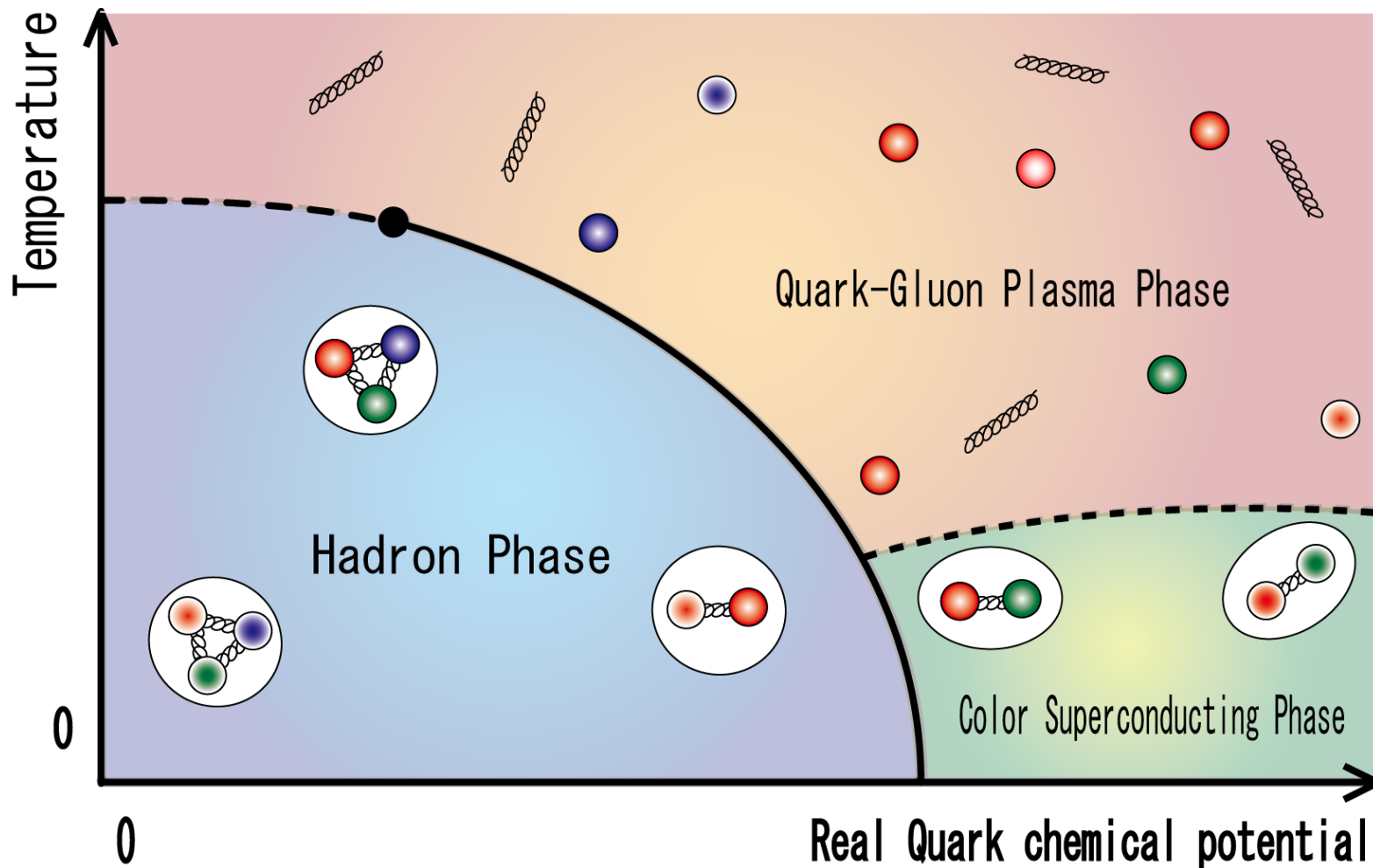
➡ The \mathbb{Z}_3 symmetry relates with the deconfinement transition via the free-energy

We can well determine the deconfinement temperature

Finite quark-mass case

Polyakov-loop is **no longer** the order-parameter !

Schematic QCD phase diagram



Important point

Finite quark-mass case :

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Ordinary phase transition  Spontaneous symmetry breaking

Phase transition described by the **topological order**

X. G. Wen, Int. J. Mod. Phys. B4 (1990) 239.

 Ground-state degeneracy



Question

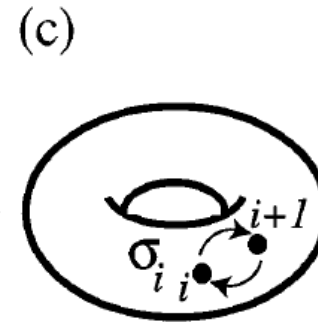
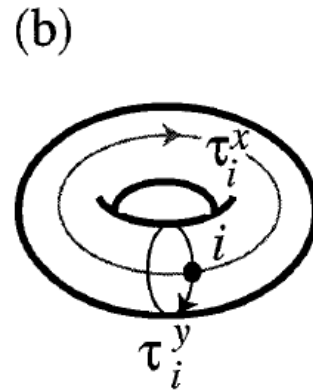
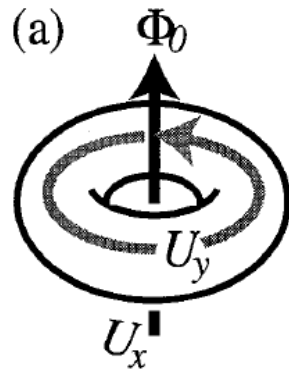
How to see the topological order at $T = 0$?



Tree adiabatic operations

M. Sato, M. Kohmoto and Y.-S. Wu, PRL 97 (2006) 010601.

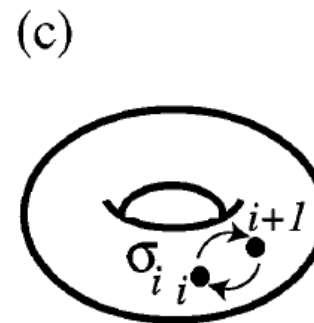
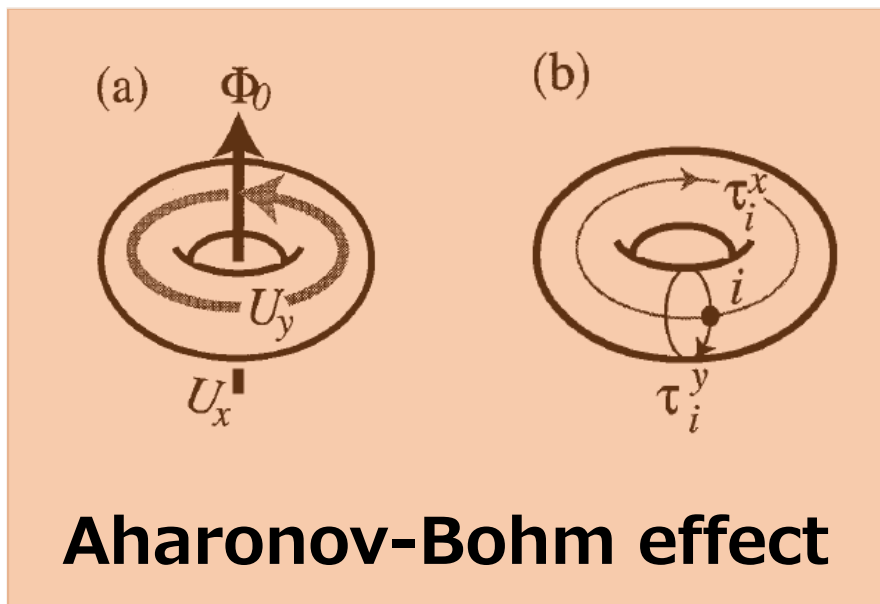
M. Sato, PRD 77 (2008) 045013.



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M. Sato, M. Kohmoto and Y.-S. Wu, PRL 97 (2006) 010601.

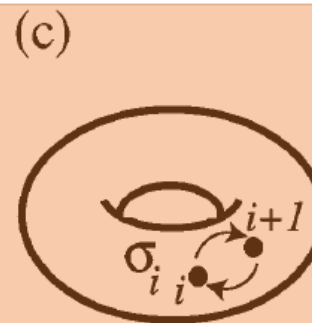
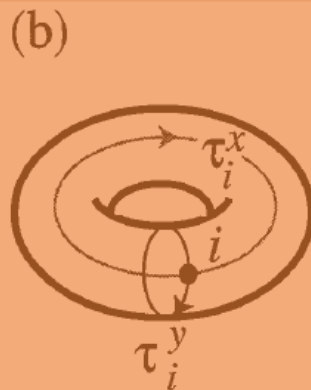
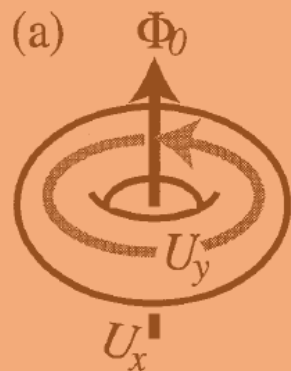
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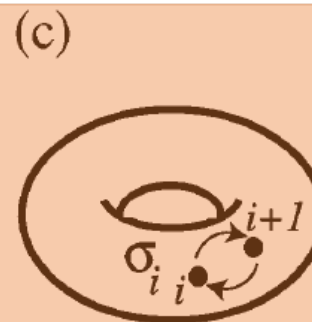
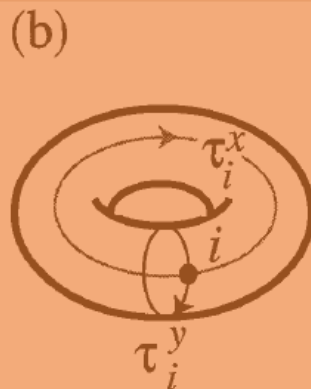
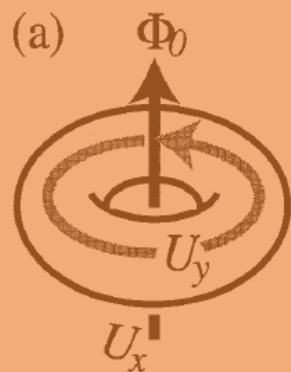
Aharonov-Bohm effect

Braid group

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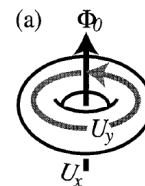
Without fractional charge : Commutable

With fractional charge : **Non-commutable** → Ground-state degeneracy

We wish to **extend** it to **finite temperature QCD!**

However, direct extension of the ground-state degeneracy is difficult...

We consider that the **imaginary chemical potential** is

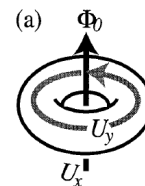


an probe to determine the deconfinement transition

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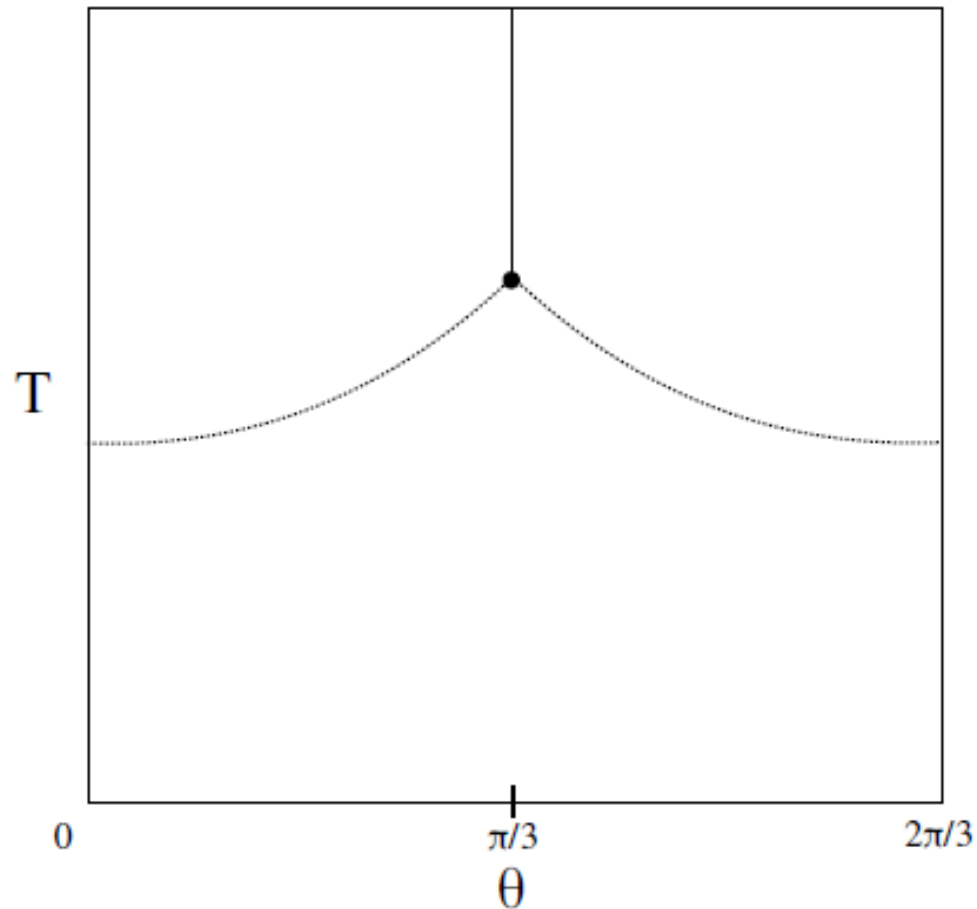
We consider that the **imaginary chemical potential** is



an probe to determine the deconfinement transition

- ➔ There is no sign problem
 - ➔ This region has all information of the region with finite μ_R
 - ➔ There are topological differences between the low and high T regions
 - ➔ We can consider similar special operations
- (Today, I do not explain this point)

Phase diagram at finite imaginary chemical potential



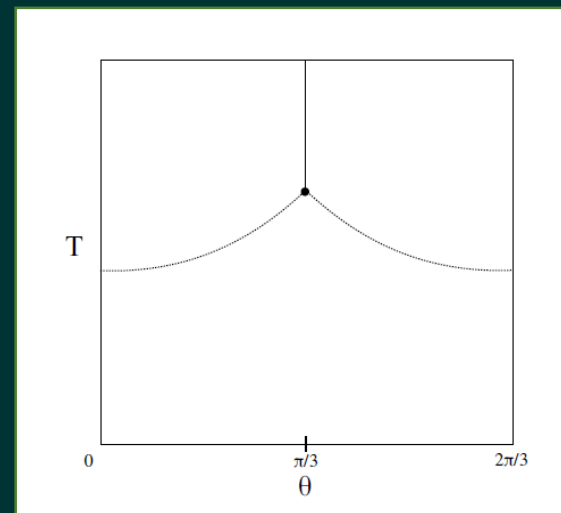
Important point

A. Roberge and N. Weiss,
Nucl. Phys. B275 (1986) 734

Roberge-Weiss periodicity

Special π/N_c periodicity along θ -direction

It appears at low and **high T**



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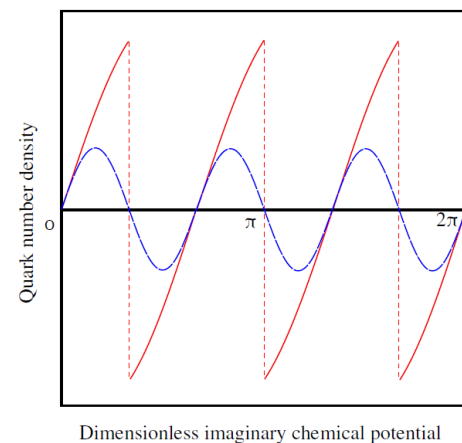
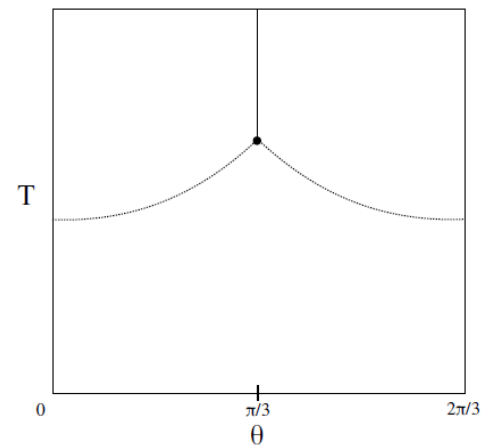
Special π/N_c periodicity along θ -direction

It appears at low and high T

Roberge-Weiss transition

First-order transition along T -direction

It is characterized by the gap
of the quark number density



Question

How to use these properties to determine
the deconfinement transition ?



It is natural to consider the **free-energy**
since we are interested in the **thermodynamic system**

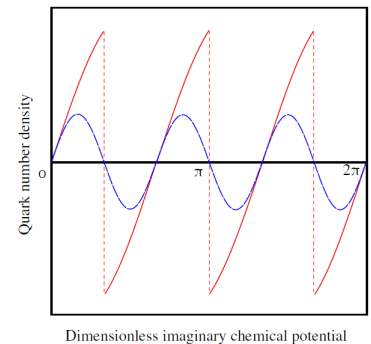
It is natural to consider the **free-energy**
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However ...

The topological order is usually difficult to discuss from thermodynamics
(Bulk thermodynamics is insensitive to the topological change)

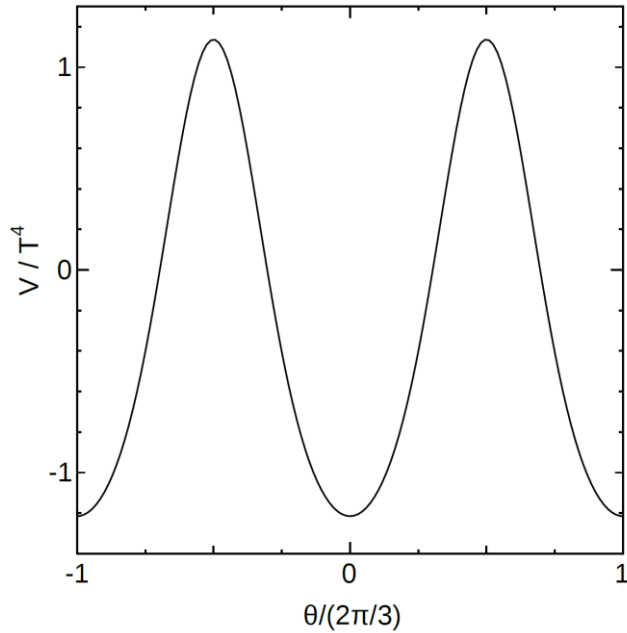
But

We can discuss topological structures at **finite θ**
(The imaginary chemical potential acts as the extra-dimension)



Result 1 : Free-energy degeneracy

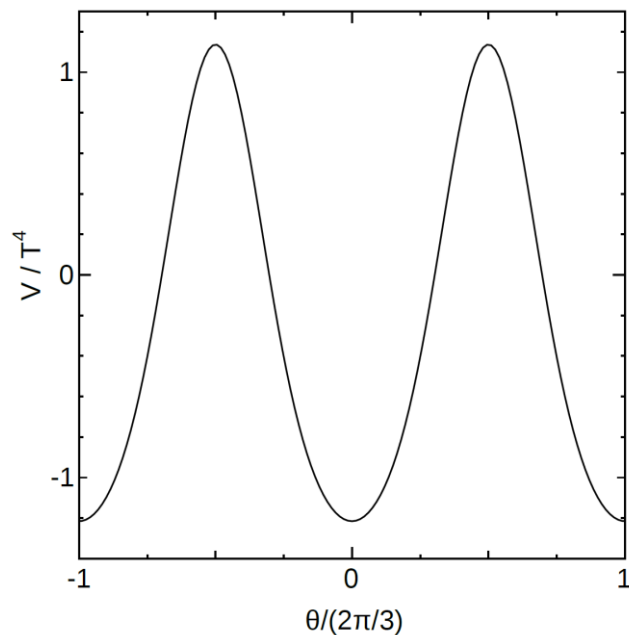
Confined phase



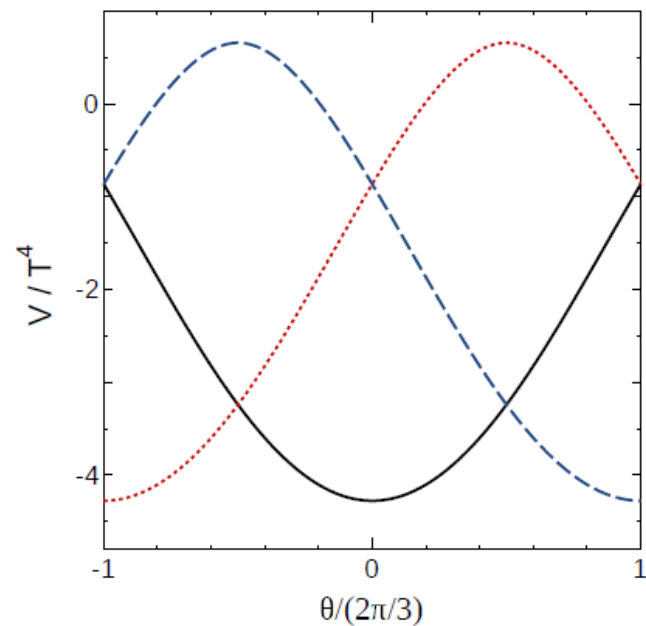
Confined phase : There is no non-trivial degeneracy

Result 1 : Free-energy degeneracy

Confined phase



Deconfined phase

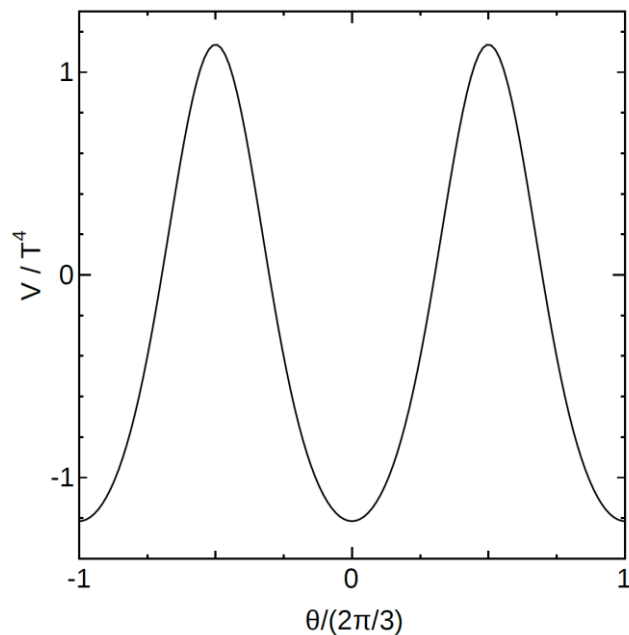


Confined phase : There is no non-trivial degeneracy

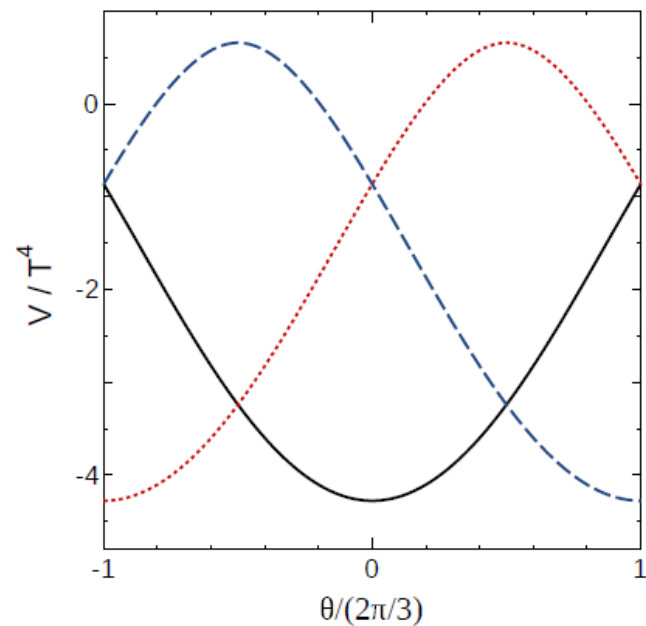
Deconfined phase : There is the **non-trivial degeneracy!**

Result 1 : Free-energy degeneracy

Confined phase



Deconfined phase

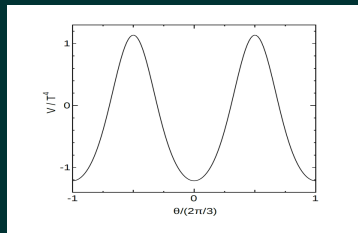


This difference relates to appearance of the quark-gluon dynamics

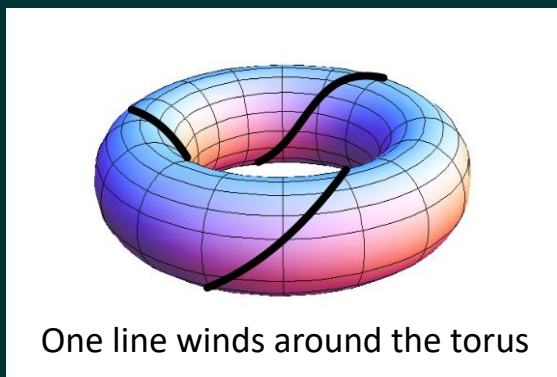
Dominant degree of freedom : Hadrons \rightarrow Quarks

Result 1 : Free-energy degeneracy

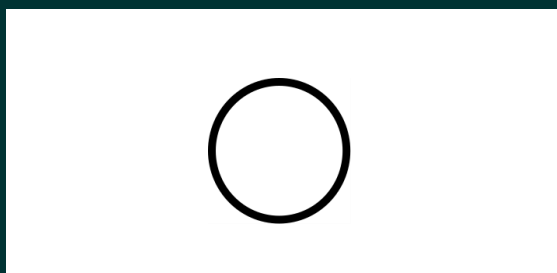
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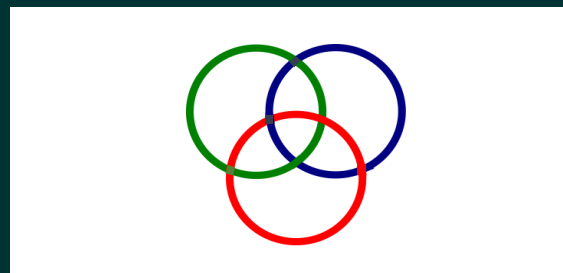
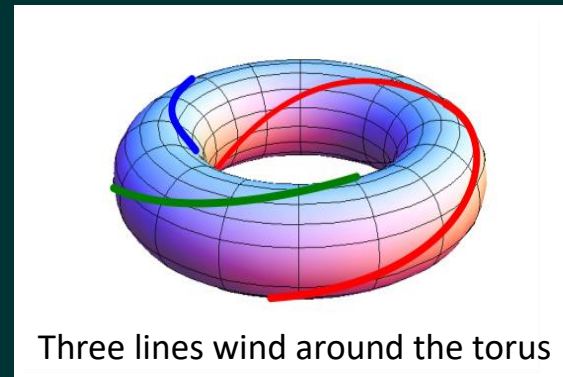
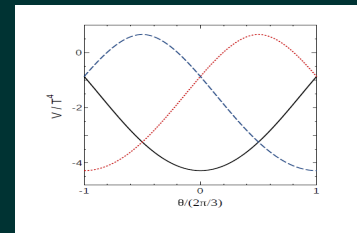
S^1 map



Two dim. map



Deconfined phase



Based on the topological difference,
we can construct the **quantum order-parameter**

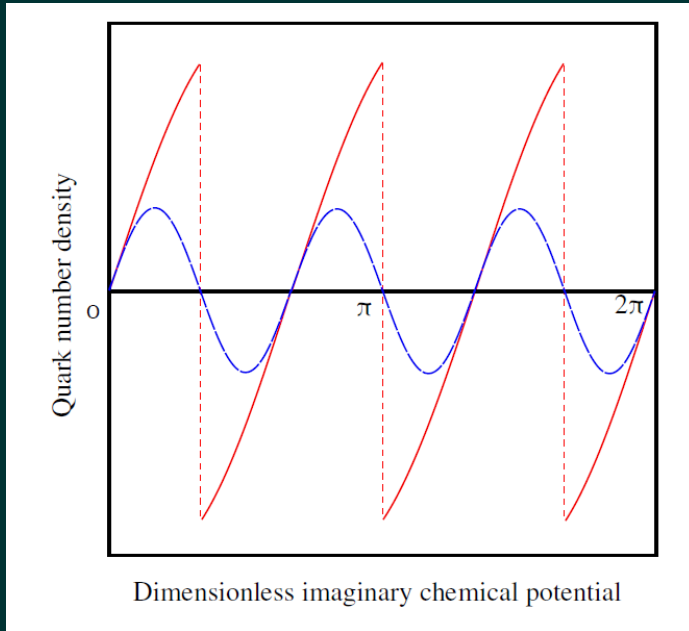
Quark number holonomy

$$\Psi = \left[\oint_0^{2\pi} \left\{ \text{Im} \left(\frac{d\tilde{n}_q}{d\theta} \Big|_T \right) \right\} d\theta \right]$$

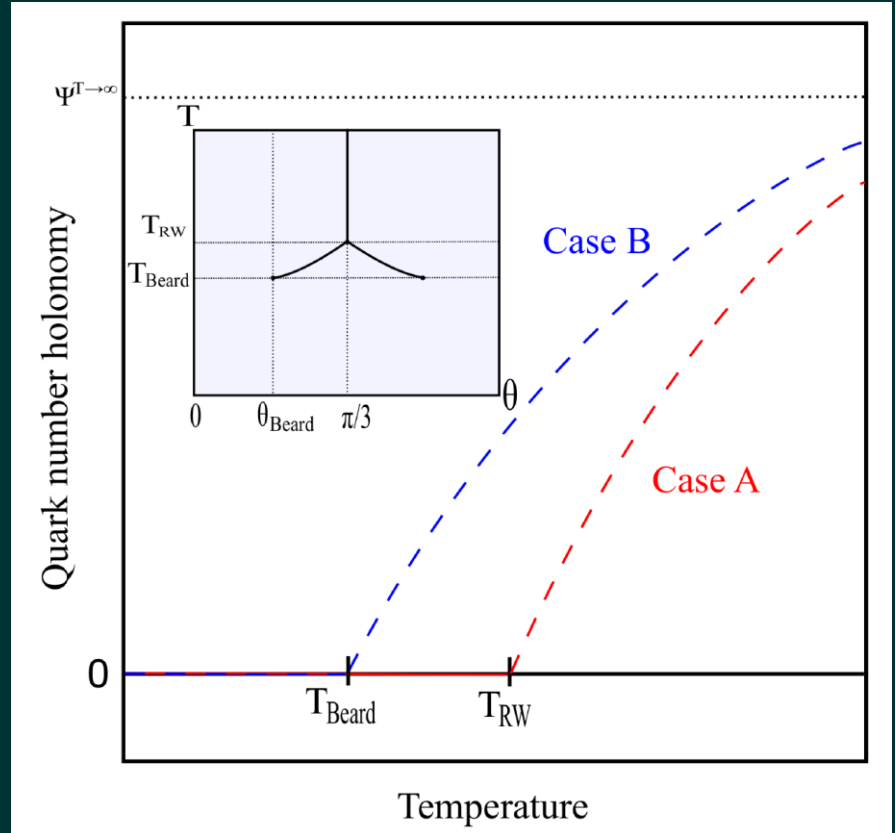
$$\tilde{n}_q \equiv C n_q$$

This quantity basically counts the number of gap in
the quark number density along θ -axis

Important point



We can well determine
the deconfinement transition



Important point

- 2+1 flavor lattice QCD

C. Bonati, M D'Elia, M. Mariti, M. Mesiti and F. Negro,
Phys. Rev. D 93 (2016) 074504

$$T_{RW} = 208(5) \text{ [MeV]}$$

Quark number holonomy becomes nonzero above this temperature

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Quark number holonomy becomes nonzero above this temperature

- Recent 2+1 flavor effective model

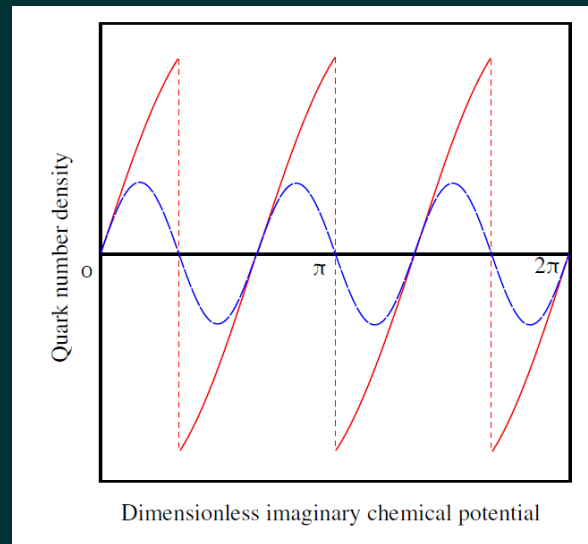
A. Miyahara, Y. Torigoe, H. Kouno, M. Yahiro,
Phys. Rev. D 94 (2016) 016003

$$T_d = \mathbf{215} \text{ [MeV]}$$

Using quantities to determine the deconfinement transition are different,
but there is good agreement. (accidental?)

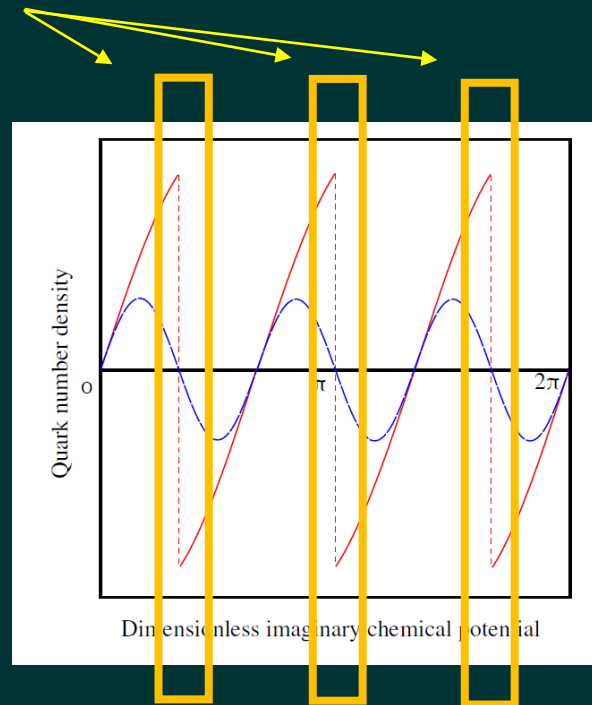
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Why the $\theta = \pi\mathbf{k}/N_c$ can detect the topological differences?



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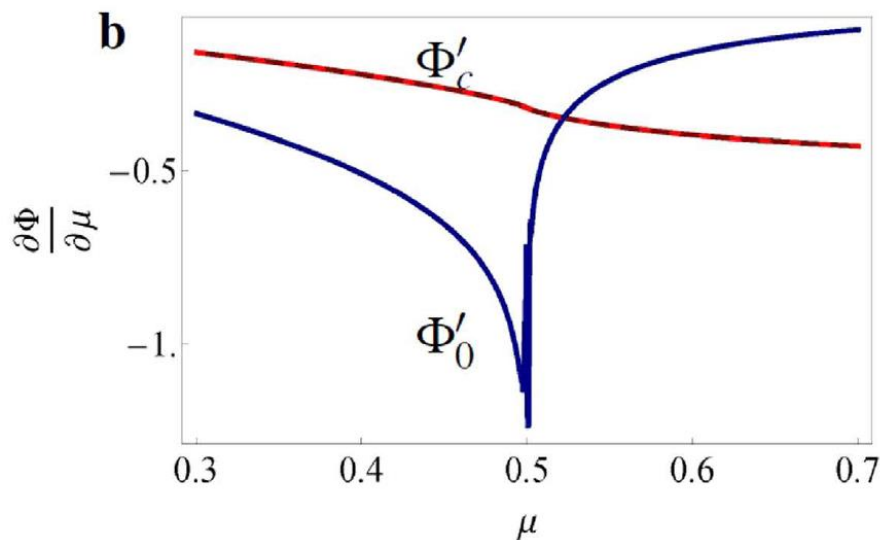


The topological order can be clarify from **thermodynamics (at edges)**

S. Kempkes, A. Quelle, and C. M. Smith, Scientific Reports 6 (2016)

Example : Kitaev chain model

$$\mathcal{H} = - \sum_{j=1}^N \left[\kappa (a_j^\dagger a_{j+1} + a_{j+1}^\dagger a_j) + \mu a_j^\dagger a_j - \Delta a_j a_{j+1} - \Delta^* a_{j+1}^\dagger a_j^\dagger \right]$$

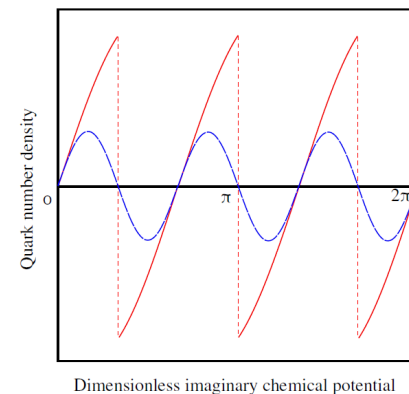


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S. Kempkes, A. Quelle, and C. M. Smith, Scientific Reports 6 (2016)

Possible analogy

We treat the **imaginary chemical potential** as the extra-dimension (parameter) and thus the $\theta = \pi k/N_c$ points acts as edges of the extended system



Therefore, thermodynamics (gaps in the quark number density) can describe the deconfinement phase transition

What happen at finite **real chemical potential**?

In our determination, we should consider

the **complex chemical potential**

and thus it is very difficult to discuss it

In this study,

we employ **Polyakov-loop extended Nambu—Jona-Lasinio** model

● **PNJL Lagrangian density**

$$\mathcal{L} = \bar{q}(\not{D} + m_0) - G[(\bar{q}q)^2 + (\bar{q}i\gamma_5\vec{\tau}q)^2] + \mathcal{V}_g(\Phi, \bar{\Phi})$$

(Good point)

It can describe the chiral phase transition and approximately treat the deconfinement transition via the Polyakov-loop

It can reproduce the RW periodicity and transition

(Bad point)

Unfortunately,
this model still has the model sign problem at finite real chemical potential

So, we consider the isospin chemical potential

- **Sign problem free**

$$\tau_2 \gamma_5 D \gamma_5 \tau_2 = D^\dagger \quad \rightarrow \quad \det(D) \geq 0$$

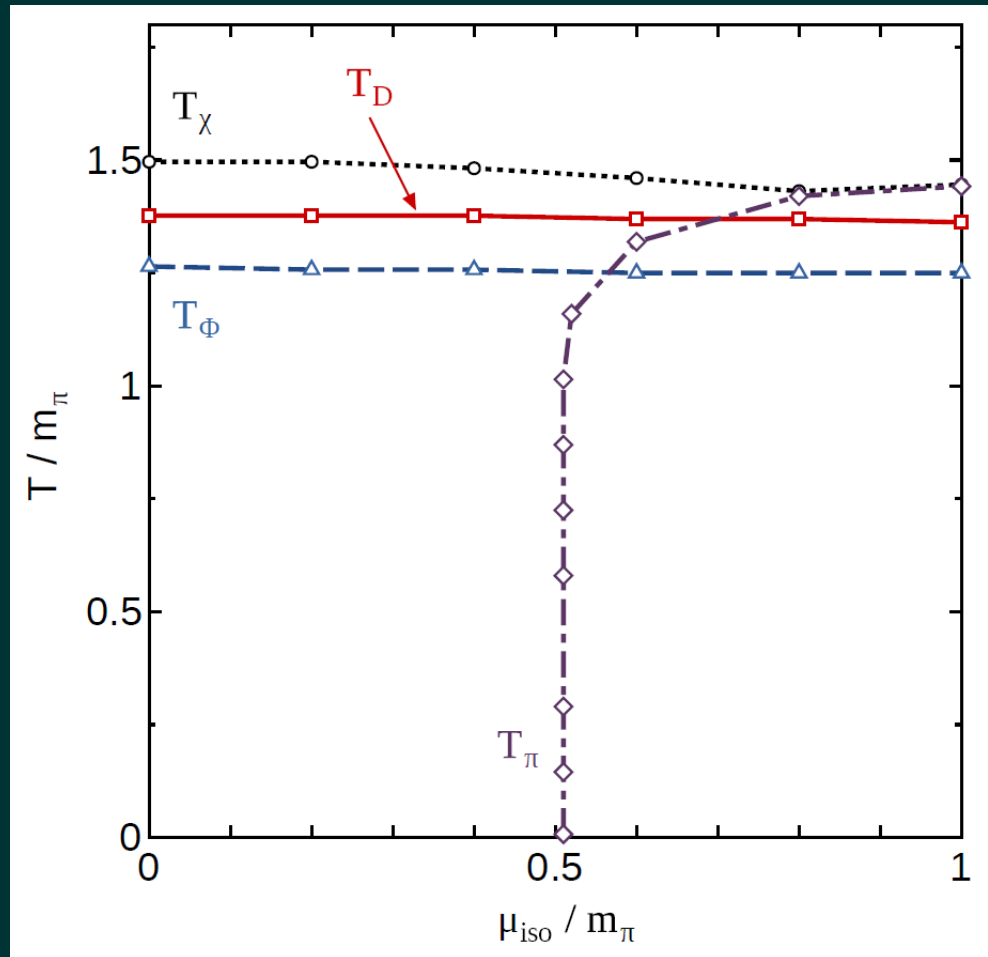
- **Orbifold equivalence**

Outside of the pion condensed region,

**the phase diagram at finite real μ and
the phase diagram at finite isospin μ**

are identical to each other in the large N_c limit

Phase diagram from PNJL model



How to investigate the deconfinement transition at finite real chemical potential?

We need better ways to handle the **sign problem**

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- Lefschetz-thimble path-integral method ?
- Complex Langevin method ?

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We need better ways to handle the **sign problem**

- Lefschetz-thimble path-integral method ?
- Complex Langevin method ?
- **Path optimization method !**

Yuto Mori, K.K., Akira Ohnishi, in preparation

Unfortunately, there is no talk about it ...

If you have interesting on it,
please check our forthcoming paper

We investigate the deconfinement transition from **topological viewpoints**

1. To discuss the deconfinement transition at finite temperature, we use the **nontrivial free-energy degeneracy**
2. We determine the **new order-parameter** of deconfinement transition

$$\Psi = \left[\oint_0^{2\pi} \left\{ \text{Im} \left(\frac{d\tilde{n}_q}{d\theta} \Big|_T \right) \right\} d\theta \right]$$

3. The **density-dependence** of the deconfinement transition is shown by introducing the isospin chemical potential to the PNJL model

What happen in the spatial topology ?

In discussions of the topological order, spatial topology is important

1. Spatial Polyakov-loop with the spatial imaginary chemical potential
2. Entanglement entropy
3. Uhlmann phase

It is an extended quantity of the Berry phase to quantum mixed states

In these calculations, we need heavy numerical computations