

# 有限密度格子QCDの研究

## -高密度への挑戦-

K. Nagata,

Hiroshima Univ., RIISE

S. Motoki (KEK), A. Nakamura (Hiroshima)

in XQCD-J collaboration

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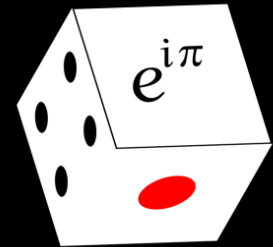


N. E. C. O.

# QCD at finite temperature and density

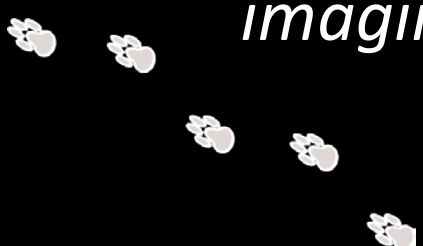


- *nature of matter in our universe*
  - *ordinary* : hadron, nuclei, nuclear matter
  - *dense* : compact stars
  - *hot* : HIC, creation of matter in early universe
- *QCD is non-perturbative.*
- *Finite density LQCD : sign problem*



# To reach at low $T$ & high $\mu$

- *Several methods.*
  - *Most studies have been done for regions near  $T_c$ .*
  - *Each method has its own difficulty*
    - *Taylor : truncation error*
    - *MPR : overlap & sign problems*
    - *Imaginary : uncertainty of functional form*
    - *etc.....*
- *We have been studied those methods,*
  - *Wilson fermion, formula for determinant, real & imaginary*



# Fermion determinant

$$Z(\mu) = \int \mathcal{D}U (\det \Delta(\mu))^{N_f} e^{-S_G}$$


- *effects of chemical potential*
- *sign problem*
- *numerical hot spot*

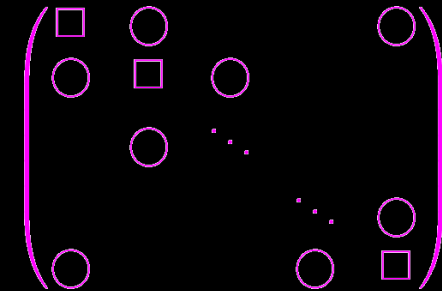


# A formula for determinant

$$Z(\mu) = \int \mathcal{D}U (\det \Delta(\mu))^{N_f} e^{-S_G}$$

- **Quark action as t-t matrix**

- temporal hop accompanies chemical potential
- spatial (diag), temporal (n.n + b.c)



- **Performing temporal determinant by hand**

- rank reduces to  $N/Nt$  (memory & CPU time  $\sim 1/Nt^2$ )
- $\det D$  is an analytic function of  $\mu$ .

- **Reduction formula enables us**

- exact evaluation of determinant (useful for large  $\mu$ )
- for arbitrary values of chemical potential
- suppress CPU time for low  $T$  ( $T=1/a Nt$ )



# Reduction formula

$$\det \Delta(\mu) = C_0 e^{(N_r/2)\mu/T} \det(Q + e^{-\mu/T})$$

- $Q$  : rank =  $N_r = 4 N_c N_s^3$  matrix ( $N_r = N/N_t$ )
  - function of link variables
  - independent of chemical potential

Wilson type : KN&AN,  
PRD82,094027 ('10),

$$\det \Delta(\mu) = C_0 e^{(N_r/2)\mu/T} \prod_{n=1}^{N_r} (\lambda_n + e^{-\mu/T})$$

- *It can be used for many applications*



# Taylor expansion

$$\frac{p(\mu, T)}{T^4} = c_0 + c_2(\mu/T)^2 + c_4(\mu/T)^4 + \dots$$

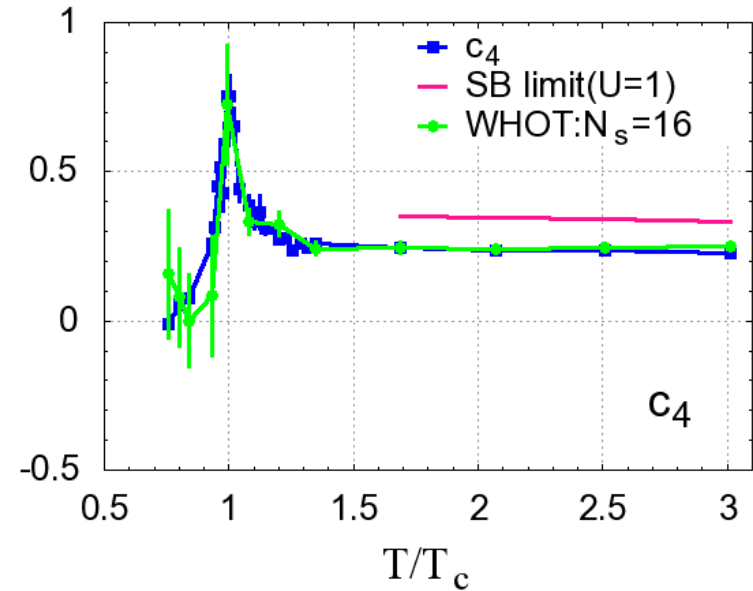
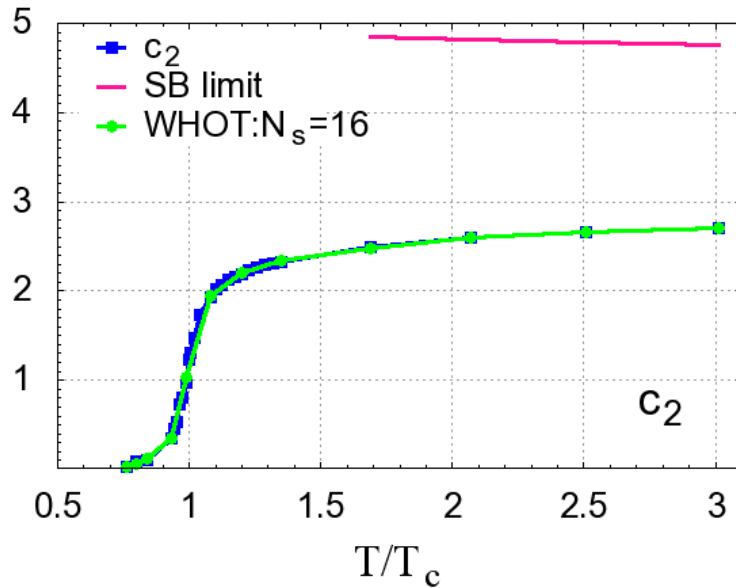
- *Coefficients are calculated at  $\mu=0$ , and functions of  $T$ .*

$$c_n(T) = \frac{1}{n!} T^n \frac{\partial^n}{\partial \mu^n} \frac{p(\mu = 0)}{T^4}$$





# Taylor coefficients



Data of "WHOT" are taken from  
WHOT, Ejiri et al.,  
PRD82, 014508 (2010),  
arXiv:0909.2121

*clover-Wilson + RG-gauge( $N_f=2$ )*

*Volume :  $8^3 \times 4$*

*quark mass :  $m_{ps}/m_V \sim 0.8$*

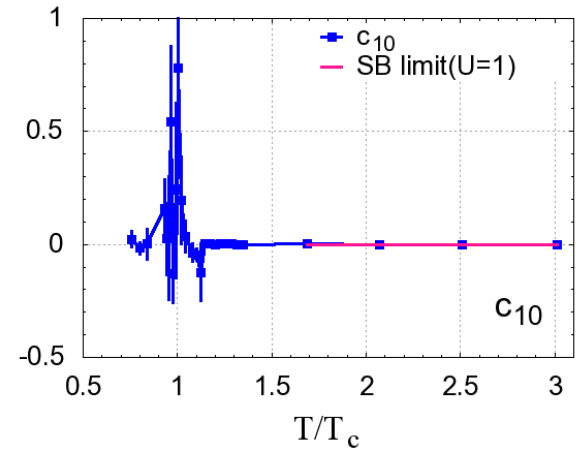
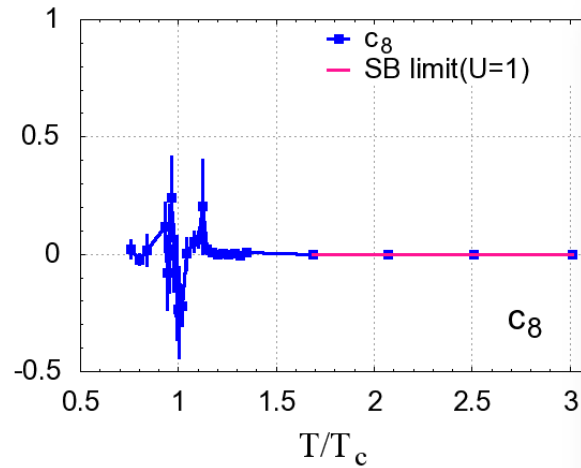
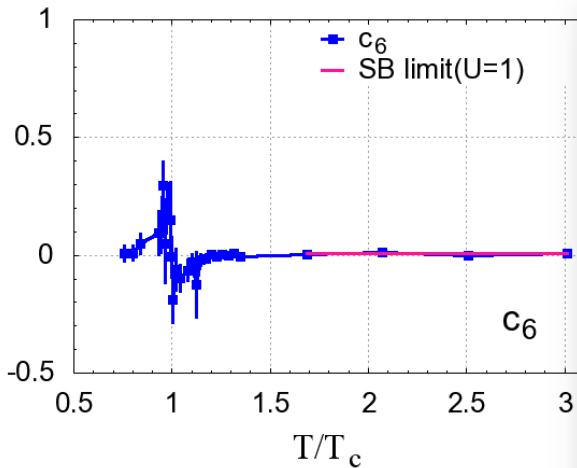
*Configurations : HMC at  $\mu=0$*

*11 K steps including ( therm. as 3-5K)*

*Eigen values : 400 configs.*



# Taylor coefficients



- *At high  $T$ , almost converge up to  $O((\mu/T)^4)$ .*
- *Near and below  $T_c$ , slow convergence. Higher order terms are non-negligible.*

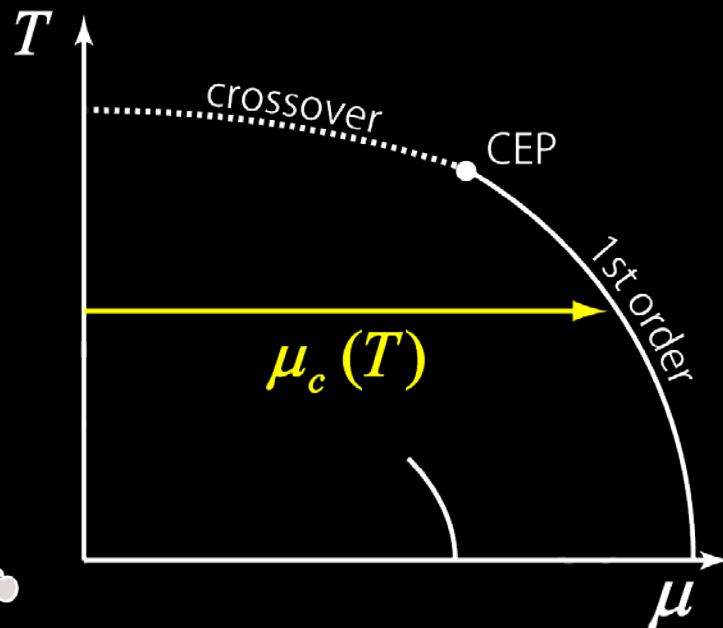
*We are preparing for higher order coefficients and those at low  $T$*



# EoS & Phase boundary

Phase boundary can be obtained from convergence radius  
(in principle)

$$r(T) = \lim_{n \rightarrow \infty} \left| \frac{C_n}{C_{n+2}} \right|$$



$$\mu_c = T \sqrt{r(T)}$$



# Lee-Yang zeros

- Zeros of partition function (Lee-Yang) implies phase transitions

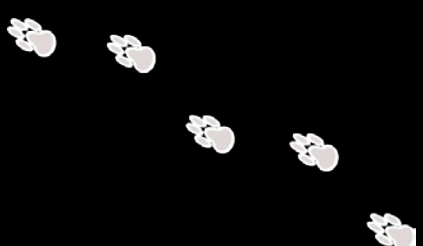
$$Z(\xi) \rightarrow 0$$

$$F = -T \ln Z(\xi)$$

$$\frac{\partial}{\partial \mu} F = -T \frac{1}{Z} \frac{\partial Z}{\partial \mu}$$

$$\xi = \exp(-\mu/T)$$

- LY zeros are distributed on the complex fugacity plane.
- LY zeros approach to  $\text{Re}[\xi]$  AXIS as  $V \rightarrow 0$ .



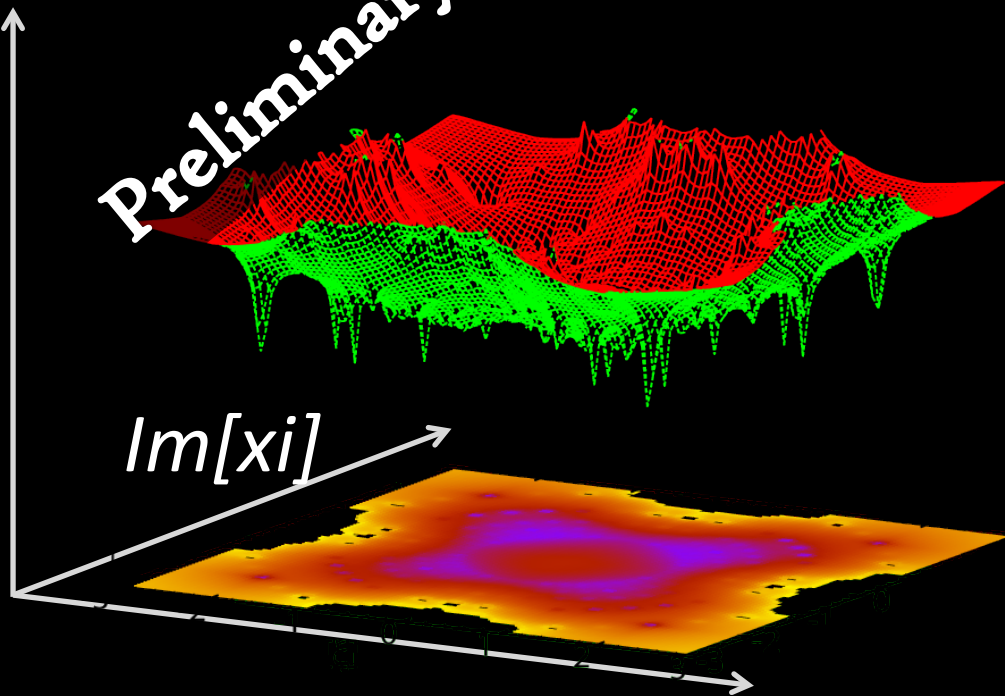
# Lee-Yang zeros

$$Z(\xi) = \left\langle \frac{\det^2 \Delta(\mu)}{\det^2 \Delta(0)} \right\rangle_0$$

$\ln Z$

Preliminary

$$F = -T \ln Z(\xi)$$



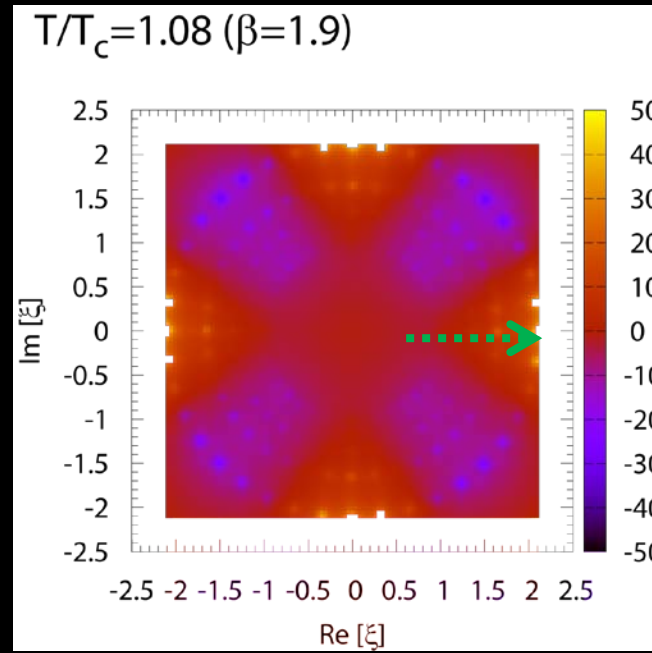
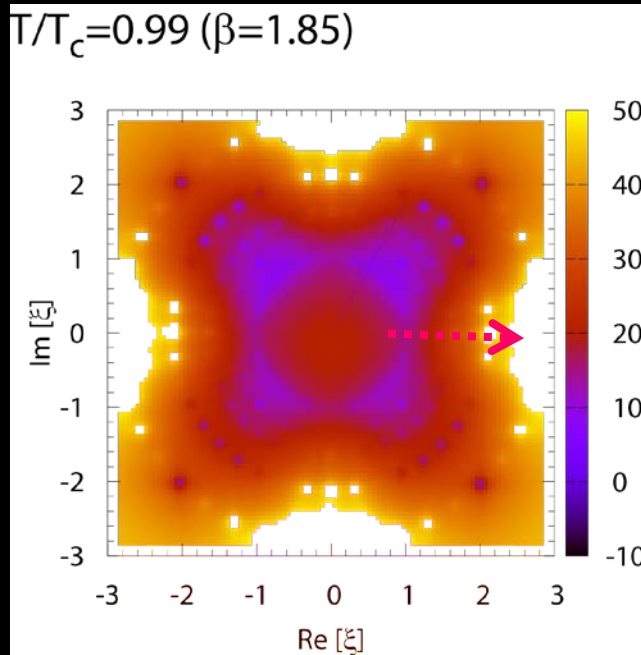
Preliminary data :  
config : clover-Wilson  
Measurement : w/o clover-term  
statistics 100

$$\xi = \exp(-\mu/T)$$

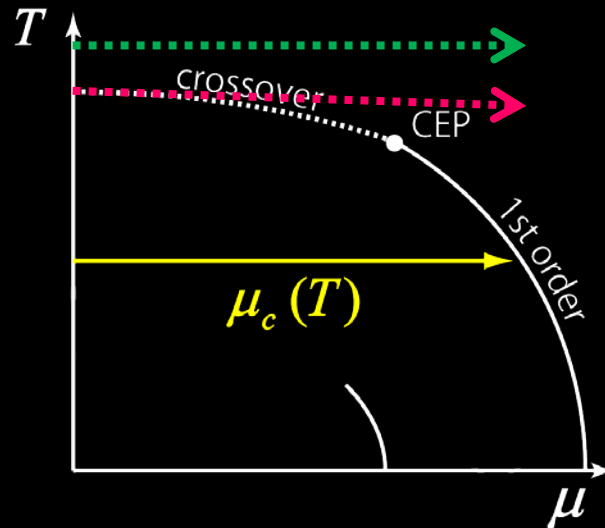
$Re[xi]$



# Lee-Yang zeros



$Im[xi]$   
 $Re[xi]$



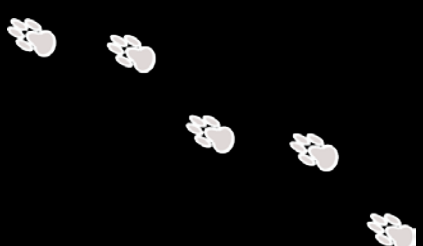
*quark : inside unit circle*  
*anti-quark : outside unit circle*



# *Imaginary chemical potential*

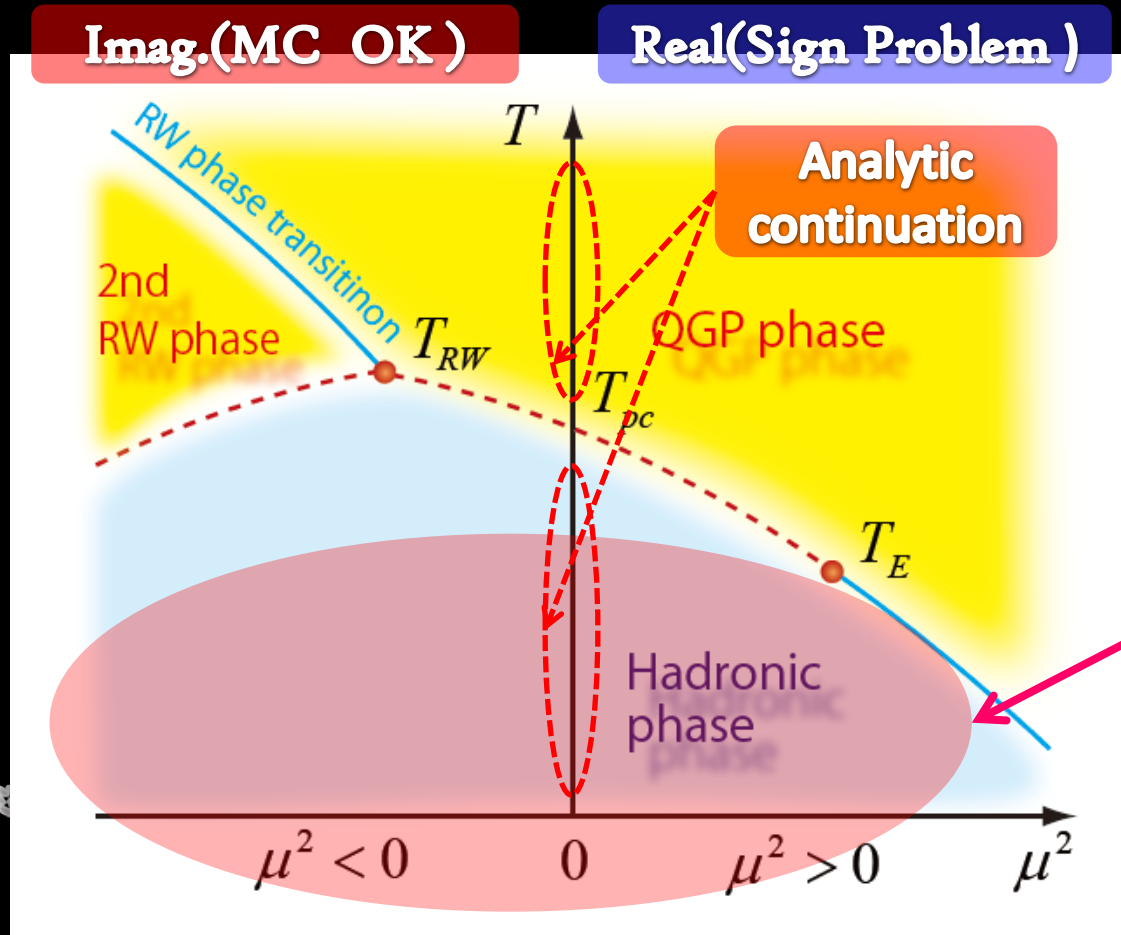
$$Z(\mu) = \int \mathcal{D}U (\det \Delta(\mu))^{N_f} e^{-S_G}$$

- *Sign problem is absent for pure imag.  $\mu$* 
  - *Generating configs. at imaginary  $\mu$*
  - *approach to real  $\mu$* 
    - *analytic continuation*
    - *canonical approach*



# Imaginary chemical potential

- MC simulations in  $\mu^2 < 0$  region and analytic continuation
- This approach may be useful for the study of low- $T$  region.



For properties of this phase diagram e.g. KN&AN, PRD83,114507, arXiv:1104.2142.

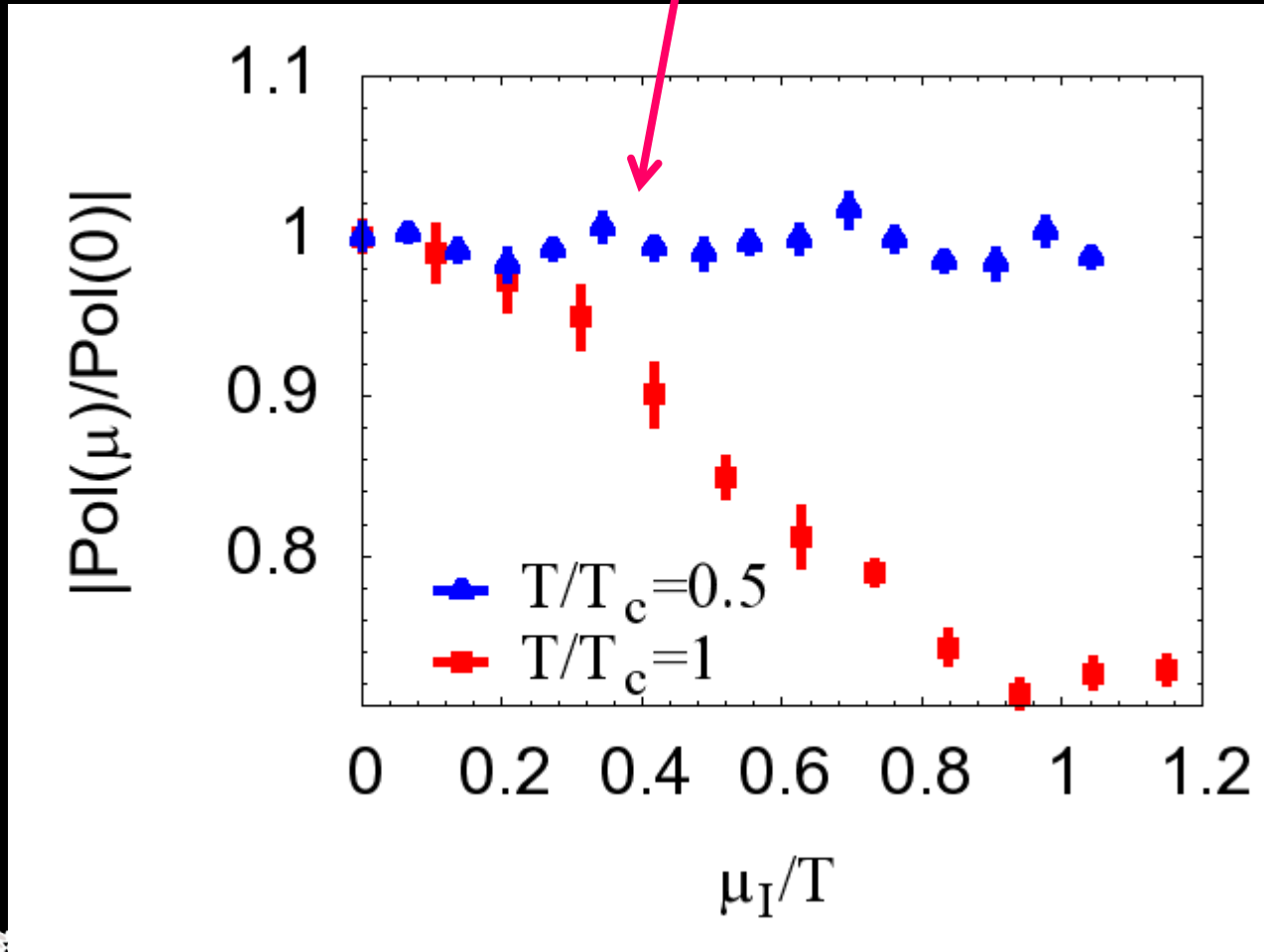
- wide regions
- RW periodicity





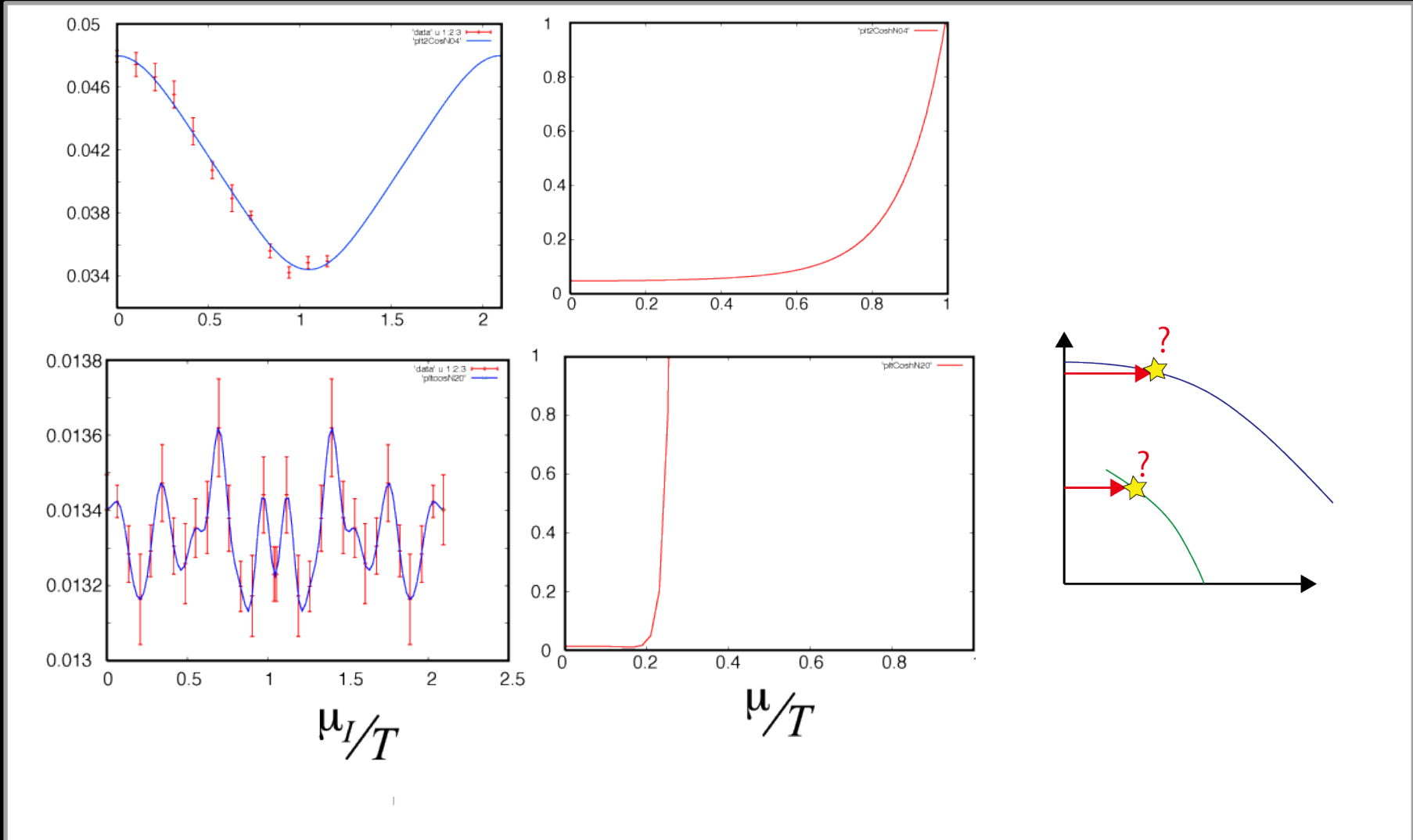
# Polyakov loop at low $T$

We found the effect of chemical potential at  $T/T_c=0.5$ .



# Fit & analytic continuation

RW periodicity  $\rightarrow$  Fourier series (cos)  $\rightarrow$  cosh



# Consistency

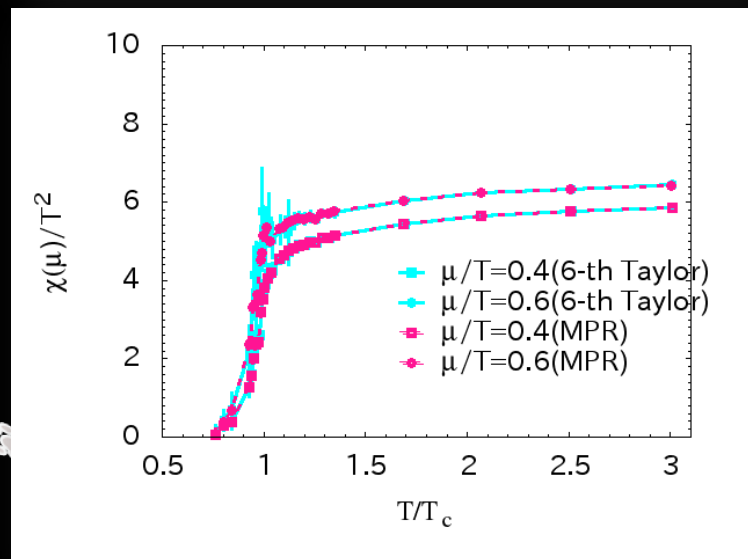
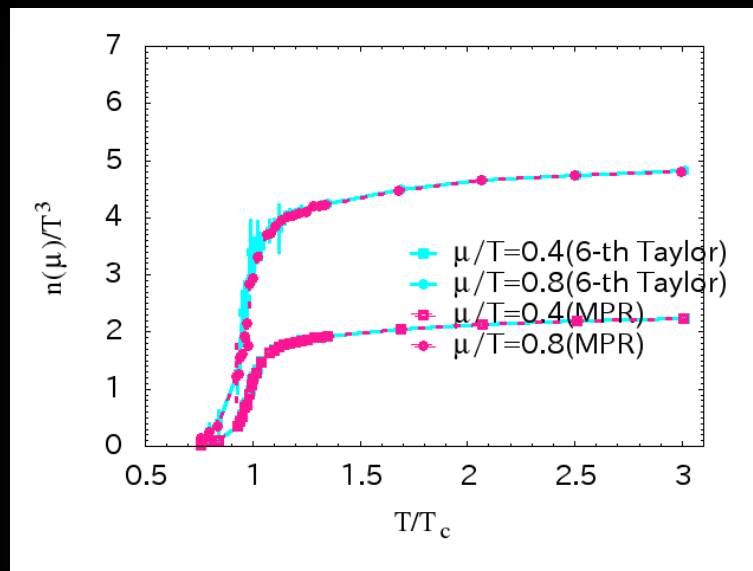
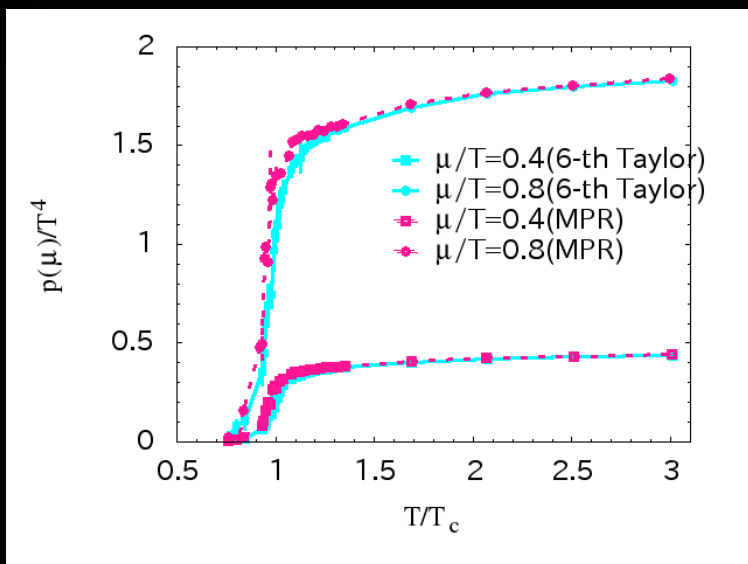
- *Each approach has its own difficulty, consistency check is helpful*

$$Z(\mu) = \int \mathcal{D}U (\det \Delta(\mu))^{N_f} e^{-S_G}$$

- *Taylor : truncation error*
- *MPR : overlap & sign problems*
- *Imaginary : uncertainty of functional form*
- *They would be consistent if calculations are done in enough precision.*



# Consistency



- *Taylor and MPR are consistent*
  - *consistent up to  $\mu/T = 0.6$*
  - *Taylor : truncation error*
  - *MPR : overlap & sign problem*



# Summary

- *We are studying several ideas towards low  $T$  & high  $\mu$* 
  - *Taylor, Lee-Yang, Imaginary etc*
  - *phase boundary, EoS ...*
  
- *Further investigations are in progress*
  - *larger volume, increasing statistics, low  $T$  ( $T=0.5T_c$ - $T_c$ )*

