

# Validity range of canonical approach to finite density QCD

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for  **$Z_n$  collaboration**

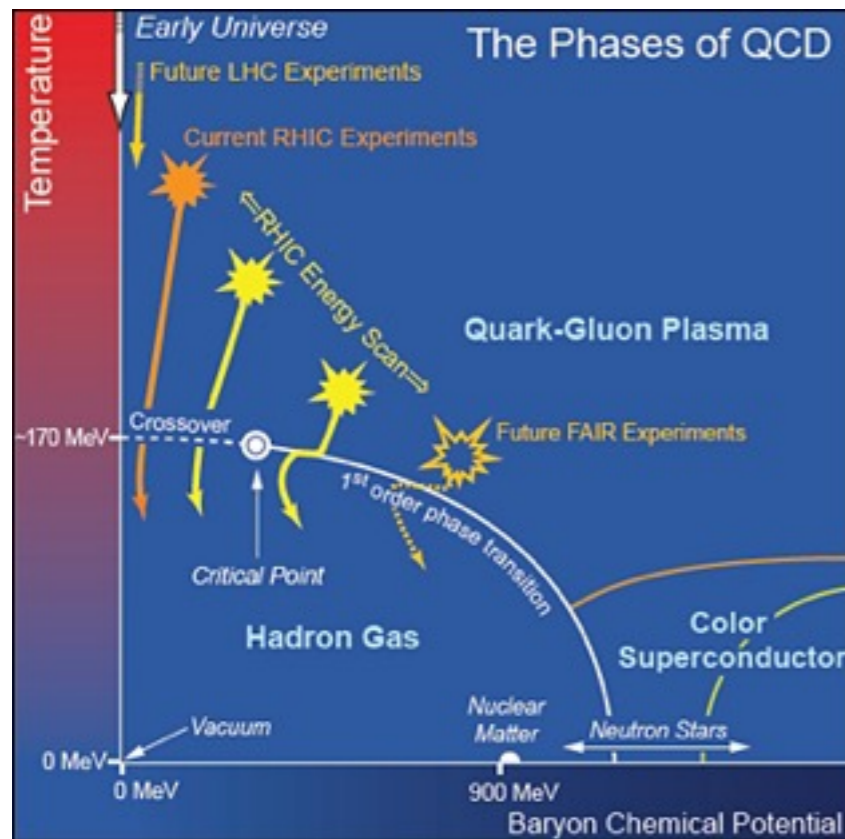


This talk is based on arXiv:1504.06351 .

LATTICE 2015 at Kobe, Japan, July 16, 2015



# Our motivation



We adopt **canonical approach** to explore finite density QCD.

A. Hasenfratz and D. Toussaint, Nucl. Phys. B **371**, 539-549 (1992)

S. Kratochvila and Ph. de Forcrand, Nucl. Phys. B, Proc. Suppl. **140**, 514 (2005)

<http://www.bnl.gov/physics/news/news.asp?a=1870&t=today>

## Purpose of this work

To calculate thermodynamic observables at finite density while comparing with results obtained by Multi parameter reweighing (MPR) method.

# What is the canonical approach ?

Fugacity expansion of grand canonical partition function

$$Z_{GC}(\mu_B, T) = \sum_{n=-\infty}^{\infty} Z_n(T) \left( e^{\mu_B/T} \right)^n$$

$\mu_B$  : baryon chemical potential       $n$  : net baryon number

How to get canonical partition function

$$Z_n(T) = \frac{1}{2\pi} \int_0^{2\pi} d\left(\frac{\mu_I}{T}\right) \underline{Z_{GC}(i\mu_I)} e^{-in\mu_I/T}$$

$\mu_I \in \mathbb{R}$

# How to calculate $Z_n(T)$ ?

$$Z_{GC}(i\mu_I) = \int [dU] \left( \frac{\det \Delta(i\mu_I)}{\det \Delta(\mu_0)} \right) \underbrace{\det \Delta(\mu_0) e^{-S_g}}_{\text{Monte-Carlo weight}} = Z_{GC}(\mu_0) \left\langle \frac{\det \Delta(i\mu_I)}{\det \Delta(\mu_0)} \right\rangle_{\mu_0}$$

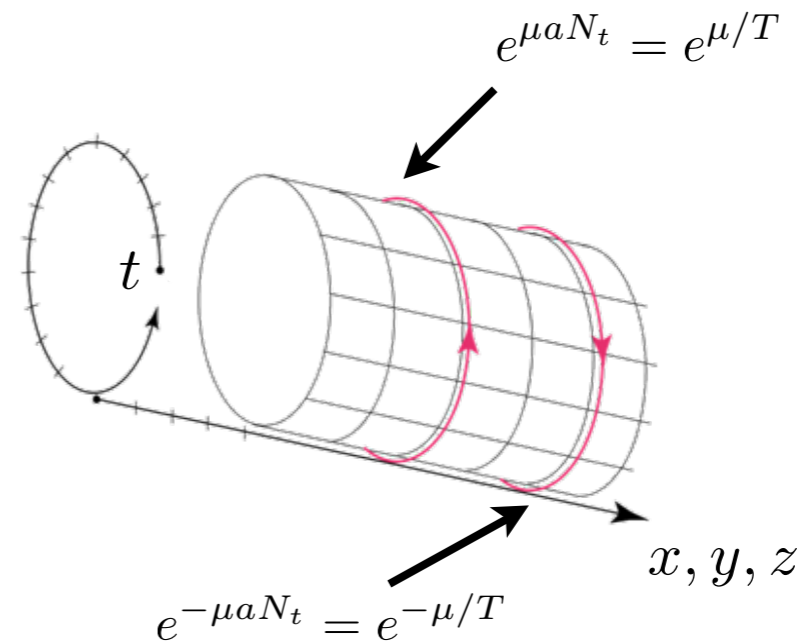
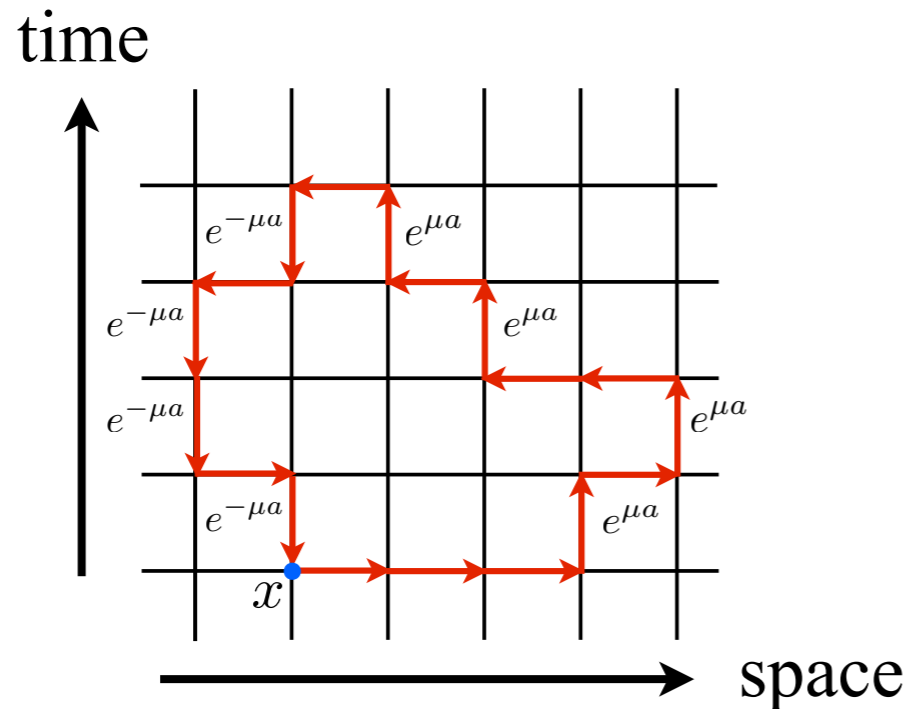


$$Z_n(T) = \underbrace{Z_{GC}(\mu_0)}_{\text{extra constant}} \times \frac{1}{2\pi} \int_0^{2\pi} d\left(\frac{\mu_I}{T}\right) \left\langle \frac{\det \Delta(i\mu_I)}{\det \Delta(\mu_0)} \right\rangle_{\mu_0} e^{-in\mu_I/T}$$

To avoid this constant, we consider  $\frac{Z_n(T)}{Z_0(T)}$ .

We need to evaluate  $\left\langle \frac{\det \Delta(i\mu_I)}{\det \Delta(\mu_0)} \right\rangle_{\mu_0}$  at various pure imaginary chemical potentials for discrete Fourier transformation (DFT).

# Strategy for calculation of Wilson fermion determinant



$$\log \det \Delta(\mu, U) = \text{Tr} \log \left( 1 - \kappa Q(\mu, U) \right) \sim \sum_{n=1}^N \frac{\kappa^n}{n} \text{Tr} Q^n(\mu, U)$$

$$= \dots + W_{-1}(U) e^{-\mu/T} + W_0(U) + W_1(U) e^{\mu/T} + \dots \equiv \sum_{n=-N/N_t}^{N/N_t} W_n(U) e^{n\mu/T}$$

- ① Generate gauge configuration  $U$  at  $\mu_0 = 0$ .
- ② Calculate a set of  $W_n$  using hopping parameter expansion (HPE).

Note : Sufficiently large  $N$  for HPE is needed .

$\text{Tr} Q^n(\mu, U)$  is evaluated through noise method.

- ③ Calculate  $\det \Delta(i\mu_I, U)$  at desired  $\mu = i\mu_I$  using  $W_n$ .

# Our strategy for discrete Fourier transformation

$$\begin{aligned} \frac{Z_n(T)}{Z_{GC}(\mu_0)} &= \frac{1}{2\pi} \int_0^{2\pi} d\left(\frac{\mu_I}{T}\right) \left\langle \frac{\det \Delta(i\mu_I)}{\det \Delta(\mu_0)} \right\rangle_{\mu_0} e^{-in\mu_I/T} \\ &\sim (\text{const.}) \times \sum_{k=0}^{M-1} Z_{GC} \left( i\frac{\mu_I}{T} = i\frac{2\pi k}{M} \right) e^{-i\frac{2\pi k}{M}n} \end{aligned}$$

In case of large  $n$ , cancellation of significant digits is not negligible in DFT because of oscillatory integral.



We use multiple precision computation for DFT.

# Numerical set up

## Action

Fermion part : 2-flavor clover improved Wilson fermion action

Gauge part : Iwasaki gauge action

## Parameters

(S. Ejiri et al, Phys. Rev. D82:014508, 2010)

$$N_s^3 \times N_t = 8^3 \times 4, \quad m_{\text{ps}}/m_v = 0.8$$

$$T/T_c = 1.35(7), 1.20(6), 1.08(5), 0.99(5), 0.93(5), 0.84(4)$$

Statistics : 400 configurations

## Observables

$$\frac{\delta p(\mu_B/T)}{T^4} \equiv \frac{p(\mu_B/T) - p(0)}{T^4}, \quad \frac{n_B(\mu_B/T)}{T^3}, \quad \frac{\chi(\mu_B/T)}{T^2}$$

# Truncation error estimation for fugacity expansion

$$Z_{GC}(\mu_B) = \sum_{n=-\infty}^{\infty} Z_n e^{n\mu_B/T} \sim \sum_{n=-N_{\max}}^{N_{\max}} Z_n e^{n\mu_B/T}$$

In numerical calculation, the fugacity expansion is finite series.

Truncation error cannot be negligible at large chemical potential.

## Our analysis

Choose baryon number  $N_{\max}$  and calculate observable  $\langle O^{(1)} \rangle, \langle O^{(2)} \rangle$ .

$$\langle O^{(1)}(\mu_B) \rangle = \sum_{n=-N_{\max}}^{N_{\max}} O_n Z_n e^{n\mu_B/T}$$

$$\langle O^{(2)}(\mu_B) \rangle = \sum_{n=-N_{\max}+1}^{N_{\max}-1} O_n Z_n e^{n\mu_B/T}$$

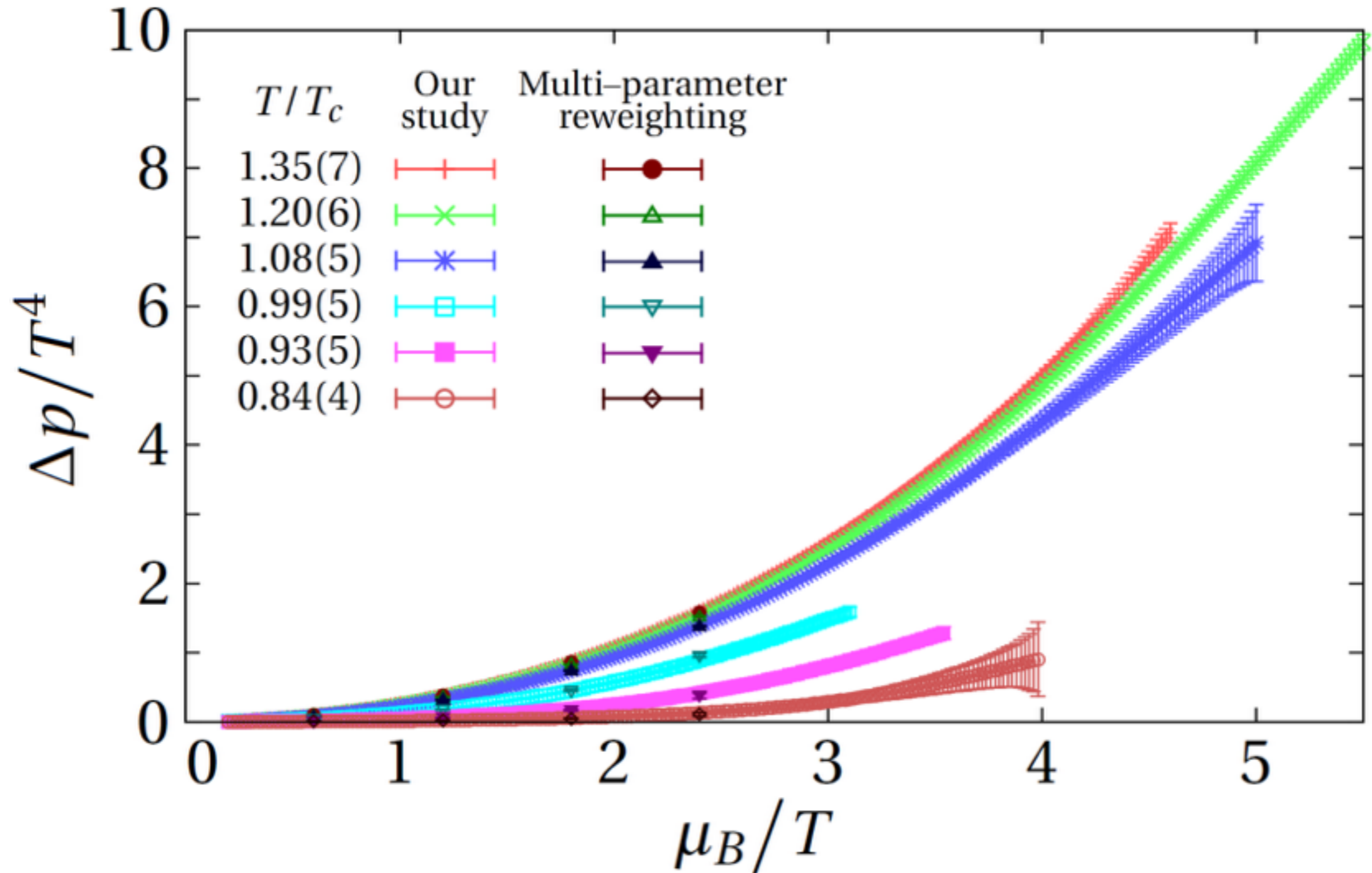
Consider chemical potential region where

$$1 - \frac{\langle O^{(1)}(\mu_B) \rangle}{\langle O^{(2)}(\mu_B) \rangle} < 10^{-3}$$

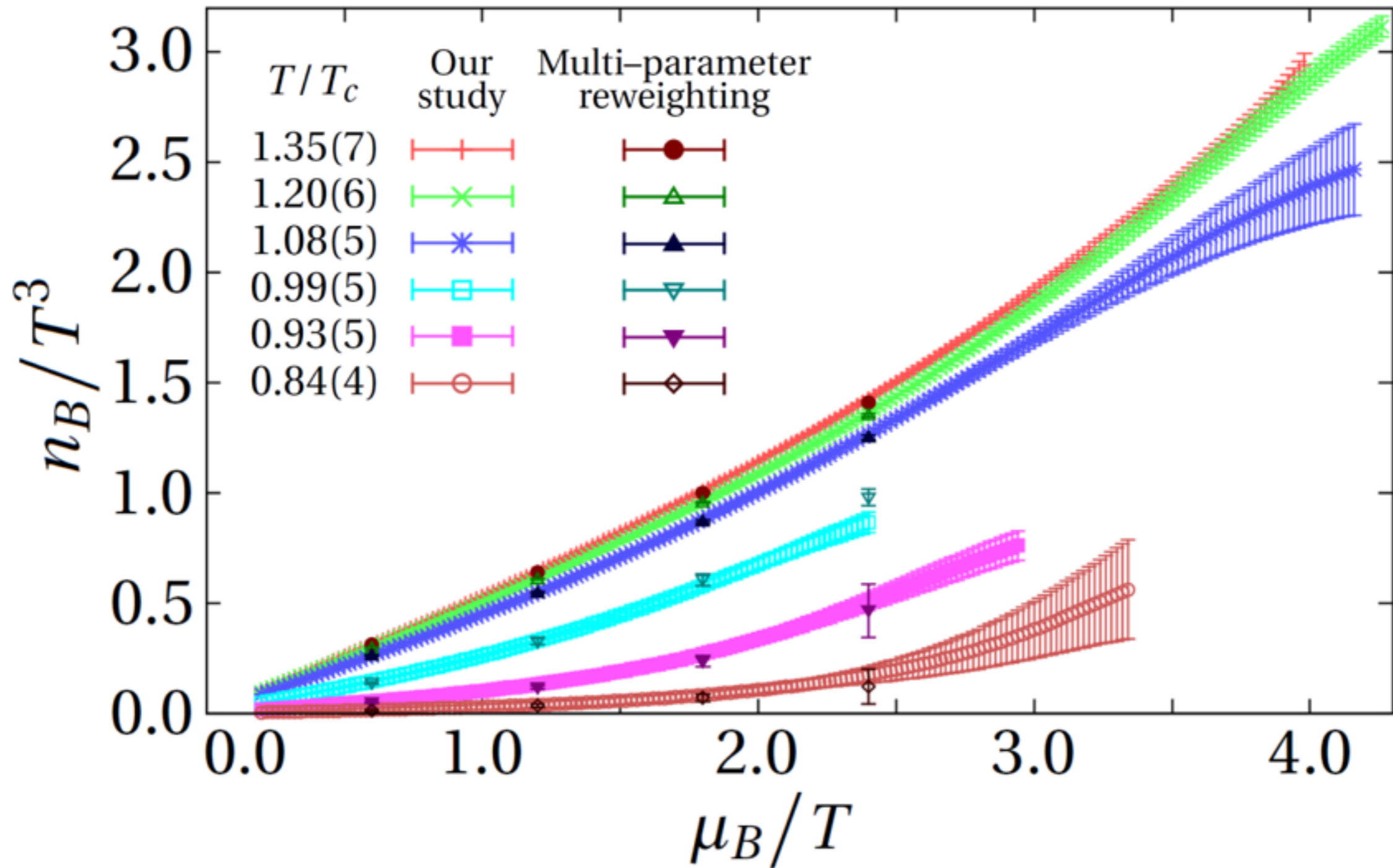
is fulfilled in our work.



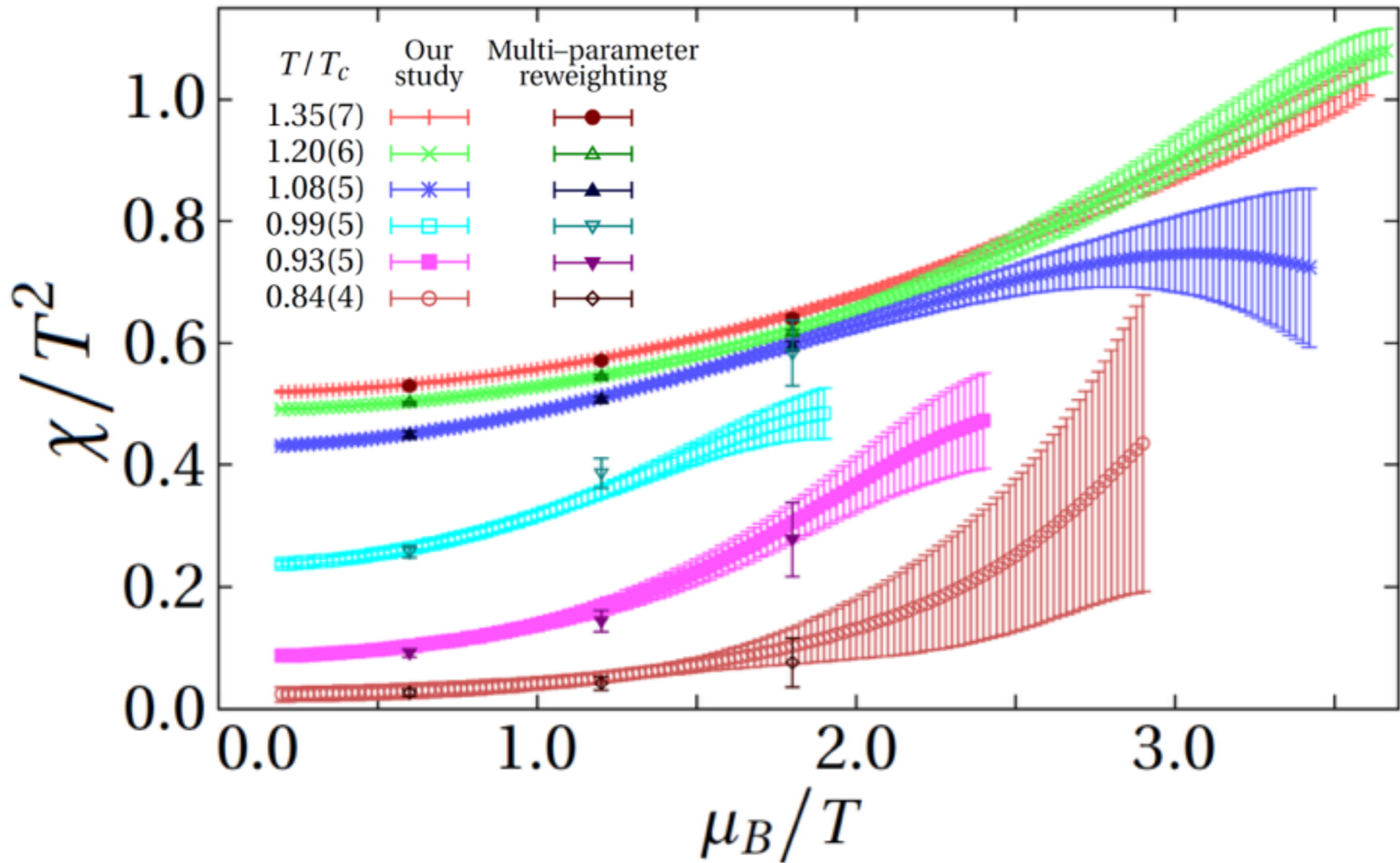
# Pressure



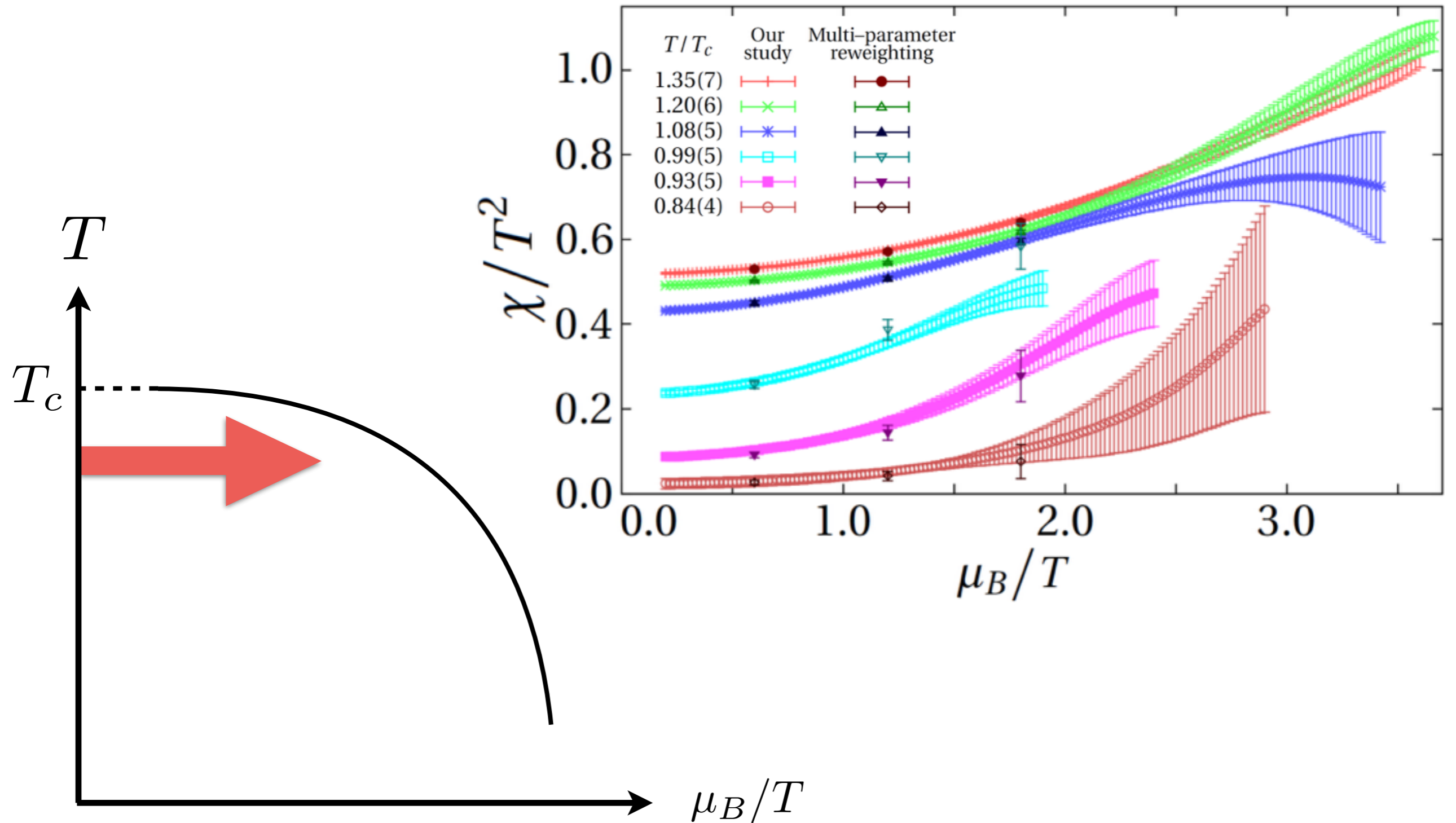
# Baryon number density



# Baryon susceptibility



# Baryon susceptibility



# Summary

We checked that the canonical approach can provide consistent results with multi-parameter reweighting method in low chemical potential region.

In many cases, the canonical approach works not only up to  $\mu_B/T = 3$  but also for  $\mu_B/T > 3$ .

There is a possibility that the **canonical approach is a candidate to overcome the “sign problem”**.

**Thank you for your attention.**

$Z_N$





old slides



# Is canonical approach useful ?

*Absolutely, yes !!*

③ We can get thermodynamic observables directly.

$$\frac{p(\mu, T)}{T^4} = \frac{1}{V_s T^3} \log Z_{GC}(\mu, T) = \frac{1}{V_s T^3} \log \left( \sum_n Z_n(T) e^{n\mu/T} \right)$$

$$\frac{n(\mu, T)}{T^3} = \frac{\partial}{\partial(\mu/T)} \frac{p(\mu, T)}{T^4} \quad , \quad \frac{\chi(\mu, T)}{T^2} = \frac{\partial^2}{\partial(\mu/T)^2} \frac{p(\mu, T)}{T^4}$$

# Hopping parameter expansion

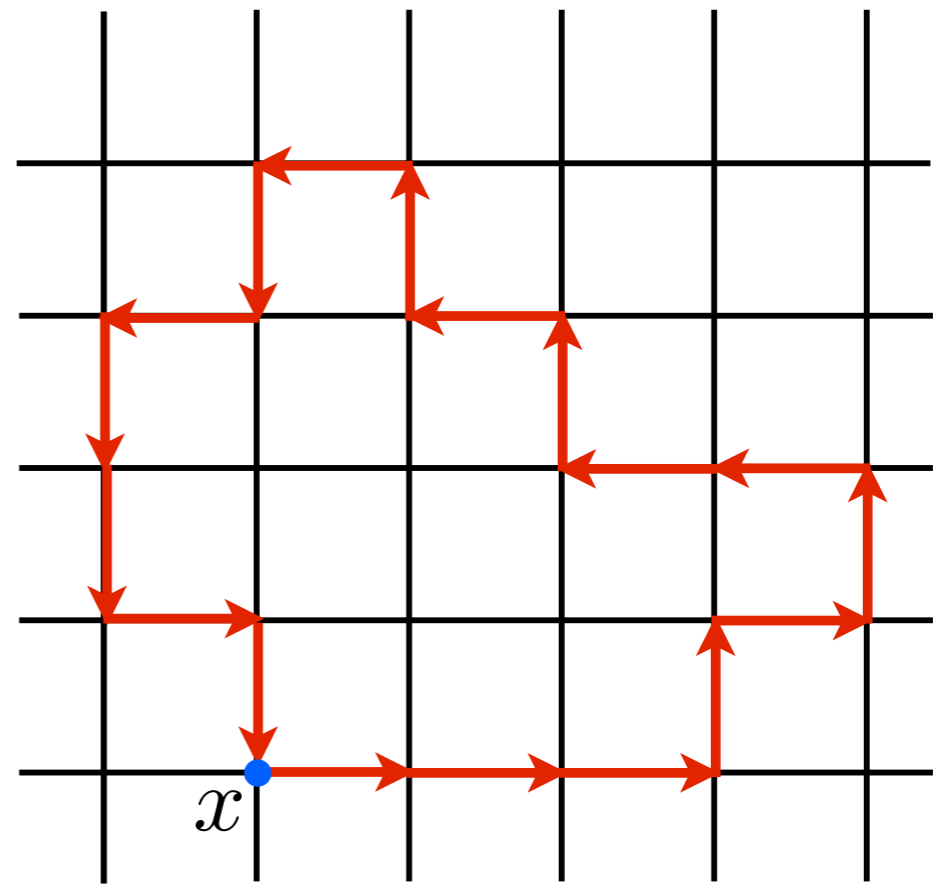
Wilson fermion matrix :  $\Delta(\mu) = 1 - \kappa Q(\mu)$

**Key identity** :  $\det \Delta = e^{\text{Tr} \log \Delta} = e^{\text{Tr} \log(1 - \kappa Q)}$

$$\text{Tr} \log(1 - \kappa Q(\mu)) = \sum_{n=1}^{\infty} \frac{\kappa^n}{n} \boxed{\text{Tr} Q^n(\mu)}$$

$$\sum_{x, \alpha, a} \langle x, \alpha, a | Q^n(\mu) | x, \alpha, a \rangle$$

$x$  : space-time    $\alpha$  : spinor    $a$  : color



Non-zero contribution comes from **closed loops**.

# Cancelled significant digits in F.T.

