

THOR meets THOR

exploring the interface of fundamental and effective approaches to extreme matter

COST THOR Working Group I & II & GDRI Meeting

COST Action CA15213

Instituto Superior Técnico, Universidade de Lisboa

Monday June 11 - Thursday June 14 2018



Canonical partition functions, virial expansion and the critical point(s) of QCD

Vitaly G.Bornyakov(IHEP&FEFU&ITEP), Vladimir A. Goy(FEFU), Maria Paola Lombardo(INFN) and Atsushi Nakamura (FEFU&RIKEN&Osaka U.)









Searching for the QCD critical point

G. Aarts, S. Hands, L. von Smekal

Series expansions are useful tools to study the phase diagram in the region which is not directly accessible to MonteCarlo simulations due to the sign problem.

Typical parameters for the expansion are the chemical potential μ and the fugacity $e^{\mu/T}$, while the observables to be expanded include the Grand Canonical Partition Function or the associated thermodynamic quantities, derivatives of the GCPF

$$Z(\mu) = \sum_{k=-3n_s^3}^{3n_s^3} < b_i > e^{k\mu T} = \sum_{k=-3n_s^3}^{3n_s^3} Q_k f^k$$

GCFP on the lattice Glasgow-style

I.M. Barbour, N. Behilil, E. Dagotto, F. Karsch, A.Moreo, M.Stone, H. Wyld, Nuclear Physics **B275** (FS17)(1986) 296.

$$\frac{p(\mu/T)}{T^4} = \sum_{n=0}^{\infty} c_n(T) \left(\frac{\mu}{T}\right)^n$$

$$p = (T/V)\ln \mathcal{Z}$$

Pressure on the lattice Bielefeld-Swansea -style

C. R. Allton, S. Ejiri, S. J. Hands, O. Kaczmarek, F. Karsch, E. Laermann, and C. Schmidt, Phys. Rev. **D66**, 074507 (2002); **D68**, 014507 (2003); C. R. Allton, M. Doering, S. Ejiri, S. J. Hands, O. Kaczmarek, F. Karsch, E. Laermann, and K. Redlich, **D71**, 054508 (2005).

Exploring QCD at finite density

G. Aarts, S. Hands, L. von Smekal

Series expansions are useful tools to study the phase diagram in the region which is not directly accessible to MonteCarlo simulations due to the sign problem.

Typical parameters for the expansion are the chemical potential μ and the fugacity $e^{\mu/T}$, while the observables to be expanded include the Grand Canonical Partition Function or the associated thermodynamic quantities, derivatives of the GCPF

Compute $\mathcal{Z} = <$ Det >. This is a polynomial in complex $e^{3\mu/T}$ whose zeros can be computed.

GCFP on the lattice Glasgow-style

I.M. Barbour, N. Behilil, E. Dagotto, F. Karsch, A.Moreo, M.Stone, H. Wyld, Nuclear Physics **B275** (FS17)(1986) 296.

Compute c_n at $\mu = 0$ Study the radius of convergence of the Taylor series in μ/T

Gavai-Gupta et al. 2003-

Pressure

on the lattice Bielefeld-Swansea -style

C. R. Allton, S. Ejiri, S. J. Hands, O. Kaczmarek,
F. Karsch, E. Laermann, and C. Schmidt, Phys. Rev.
D66, 074507 (2002); D68, 014507 (2003); C. R. Allton, M. Doering, S. Ejiri, S. J. Hands, O. Kaczmarek,
F. Karsch, E. Laermann, and K. Redlich, D71, 054508 (2005).

Exploring QCD at finite density

G. Aarts, S. Hands, L. von Smekal

Series expansions are useful tools to study the phase diagram in the region which is not directly accessible to MonteCarlo simulations due to the sign problem.

Typical parameters for the expansion are the chemical potential μ and the fugacity $e^{\mu/T}$, while the observables to be expanded include the Grand Canonical Partition Function or the associated thermodynamic quantities, derivatives of the GCPF

Compute $\mathcal{Z} = \langle \text{Det} \rangle$. This is a polynomial in complex $e^{3\mu/T}$ whose zeros can be computed.

Issue: reweighting?

Compute c_n at $\mu = 0$ Study the radius of convergence of the Taylor series in μ/T

Issue: higher orders? Alternative Strategy: the Canonical Approach

Beating the Sign problem by Canonical Approach (Hasenfraz-Toussant,1992)

Calculation in pure $\mu_{\rm I}$ regions, where no sign problem.

$$Z_n = \int \frac{d\theta}{2\pi} e^{in\theta} Z(T,\theta \equiv \frac{Im \,\mu}{T})$$

Then calculate Z in real μ regions.

$$Z(\xi, T) = \sum_{n} Z_n(T) \xi^n$$
$$\xi \equiv e^{\mu/T}$$

Recent developments:

A. Nakamura, Mod. Phys. Lett. A 22 (2007) 473; V. Bornyakov, D. Boyda, V. Goy, A. Molochkov, A. Nakamura, A. Nikolaev and V. I. Zakharov, Nucl. Phys. A 956 (2016) 809.

V. G. Bornyakov, D. L. Boyda, V. A. Gov, A. V. Molochkov, A. Nakamura. A. A. Ni kolaev and V. I. Zakharov, Phys. Rev. D **95** (2017) no.9, 094506. V. G. Bornyakov *et al.*, arXiv:1712.02830 [hep-lat].

Alternative Strategy for GCPF: the Canonical Approach

Beating the Sign problem by Canonical Approach (Hasenfraz-Toussant,1992)

Calculation in pure $\mu_{\rm I}$ regions, where no sign problem.

$$Z_n = \int \frac{d\theta}{2\pi} e^{in\theta} Z(T,\theta \equiv \frac{Im \,\mu}{T}) = e^{-\int_0^{\mu_I/T} n^n(x) dx}$$

Then calculate Z in real μ regions.

$$Z(\xi, T) = \sum_{n} Z_n(T) \xi^n$$
$$\xi \equiv e^{\mu/T}$$

Recent developments:

A. Nakamura, Mod. Phys. Lett. A 22 (2007) 473; V. Bornyakov, D. Boyda, V. Goy, A. Molochkov, A. Nakamura, A. Nikolaev and V. I. Zakharov, Nucl. Phys. A 956 (2016) 809.

V. G. Bornyakov, D. L. Boyda, V. A. Gov, A. V. Molochkov, A. Nakamura, A. A. Ni kolaev and V. I. Zakharov, Phys. Rev. D **95** (2017) no.9, 094506. V. G. Bornyakov *et al.*, arXiv:1712.02830 [hep-lat].





Pressure and its derivatives from imaginary μ

- •Strategy: analytically continue from imaginary μ MpL 00 de Forcrand Philipsen 02 d'Elia MpL 02
- Observation: optimal parametrization depends on T



Virial expansion natural

see M.D'Elia plenary review at QM2018

Virial expansion and Cluster Model Expansion



Summary of introduction

Interplay of virial expansion, canonical expansion, radius of convergence at real and imaginary chemical potential

Virial expansion within CME constraints the CEP

Plan

Virial expansion, canonical PF and GCPF in 1dQCD

Check CME in QCD

A simple model : QCD in 1d

One dimensional QCD is a an interesting, exactly solvable model. There is no spontaneous symmetry breaking, but there are baryons. Its partition function is formally the same as the one obtained in 4d QCD at strong coupling, once an explicit mass term is identified with the 4d dynamically generated mass. For any temperature T and varying μ there is a crossover to a baryon rich phase, which turns into a first order transition at zero temperature. The partition function reads

$$\mathcal{Z}(\mu, T) = 2\cosh(\mu/T) + \sinh(4m'/T) / \sinh(m'/T) = \sum_{n=-1,0,1} A_n \xi^n$$

 $A_0 = \sinh(4m'/T) / \sinh(m'/T) \equiv A, A_1 = A_{-1} = 1, \mu$ is understood to be $\mu_B, m' = sh^{-1}m$, with m the quark mass.



Zeros of the GCPF and radius of convergence in QCD 1d



 $4 < A < \infty \rightarrow y_1 = 1/y_2$ are two negative real roots.

$$\Re e^{\mu/T} = (-A + / - \sqrt{A^2 - 4})/2 \equiv y_{1/2}$$

$$\mu = \pm T(\log|y_1|) + i(2k+1)T\pi.$$



Partial sums reveal the radius of convergence

$$n^{n}(\mu_{I}) = \sum_{1}^{n} a_{k} \sin(k\mu_{I}/T)$$

$$\downarrow \qquad \text{Analytic continuation}$$

$$n^{n}(\mu) = \sum_{1}^{n} a_{k} \sinh(\mu/Tk)$$



The GCPF from the particle number: inherits the same radius of convergence



The GCPF from the particle number: inherits the same radius of convergence



The GCPF in QCD 1d beyond the radius of convergence



The convergence of the CPF to their exact result: an insufficient approximation produces a *negative* Z2

CPF, Virial Expansion, radius of convergence and GCPF QCD 1d



We use gauge field configurations generated in $N_f = 2$ lattice QCD with clover improved Wilson quarks and Iwasaki improved gauge field action. More details about the configurations can be found in Ref.[5]

The simulations were made on $16^3 \times 4$ lattices at temperatures $T/T_c = 1.35, 1.20, 1.08$, and 1.035 in the deconfinement phase and 0.99, 0.93, 0.84 in the confinement phase along the line of constant physics with $m_{\pi}/m_{\rho} = 0.8$. The parameters of the action, including c_{SW} value were borrowed from the WHOT-QCD collaboration paper[6]. We compute the number density on samples of 1800 or 3800 configurations using every 10-th trajectory produced with Hybrid Monte Carlo algorithm.



We measure the quark number density n_q at imaginary quark chemical potential $\mu_q = i\mu_{qI}$. It is defined as

$$\frac{n_q}{T^3} = \frac{1}{VT^2} \frac{\partial}{\partial \mu_q} \ln Z_{GC} = \frac{N_f N_t^3}{N_s^3 Z_{GC}} \int \mathcal{D}U e^{-S_G} (\det \Delta(\mu_q))^{N_f} \operatorname{tr} \left[\Delta^{-1} \frac{\partial \Delta}{\partial \mu_q / \mathrm{T}} \right]$$

We then fitted the number density to functions

$$n_{qI}(\theta)/T^3 = \sum_{k=1}^{k_{max}} a_k \sin(3k\theta)$$

for $T < T_{RW}$ and

$$n_{qI}(\theta)/T^3 = \sum_{n=1}^{2} c_{2n-1}(\theta^{2n-1})$$

for $T > T_{RW}$ ($\theta = \mu_{qI}/T$). In this case we used Fourier transform to compute the virial coefficients a_k .



Comparison of a_k (circles) with predictions of CEM (solid curves). There is a good agreement for $T < T_{RW}$. There is no agreement for $T > T_{RW}$.



Ratio $|a_k/a_{k+1}|$ for four temperature values. Curves show predictions of CEM. There is good agreement for $T/T_c = 1.08$. Note, that for this temperature already a_2/a_3 can be taken as a good estimate of the $k \to \infty$ limit.



Radius of convergence of the virial expansion $(\mu_B/T)_{conv}$ computed using CEM and estimated from the ratio a_2/a_3 .

Summary

Results at imaginary chemical potential offer a simple way to compute the virial expansion.

The series may converge at imaginary chemical potential, but not at real chemical potential.

However, the canonical partition functions computed at imaginary chempot allow the reconstruction of the GCPF everywhere.

In the case of 1d QCD, an insufficient approximation may produce negative canonical partition functions.

Negative partition functions have been observed in real QCD and it remains to be seen whether the origin is the same as in QCD1d.

Virial coefficients can be computed within different phenomenological frameworks. The CEM parametrization puts limit on CEP..

Preliminary numerical results suggest discrepancies with CEM in QCD