QCD at nonzero chemical potential and the sign problem

INT lectures 2012

II: standard approaches

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Where are we?

complex weight:

- straightforward importance sampling not possible
- overlap problem

various possibilities:

- preserve overlap as best as possible
- use approximate methods at small μ
- do something radical:
 - rewrite partition function in other dof
 - explore field space in a different way

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discuss first two approaches

Reminder: physics goal

determine

- phase boundary between confined and deconfined phase at small μ
- critical endpoint (if it exists)



"standard conjectured" phase diagram

Reminder: physics goal

phase boundary at small μ :

determine curvature of the phase boundary

$$\frac{T_c(\mu)}{T_c(0)} = 1 + \# \left(\frac{\mu}{T_c(0)}\right)^2 + \# \left(\frac{\mu}{T_c(0)}\right)^4 + \dots$$

(if crossover: this may depend on observable)

determine critical endpoint

- from this expansion
- directly

experimental search for critical endpoint is planned at FAIR (GSI, Darmstadt, Germany) in coming years

general strategy: $Z_w = \int DU w(U)$ $w(U) \in \mathbf{C}$ observable: $\langle O \rangle_w = \frac{\int DU O(U) w(U)}{\int DU w(U)}$

introduce new weight r(U) (*r* for 'reweighting' or 'real'), chosen at will

$$\langle O \rangle_w = \frac{\int DU O(U) \frac{w(U)}{r(U)} r(U)}{\int DU \frac{w(U)}{r(U)} r(U)} = \frac{\langle O \frac{w}{r} \rangle_r}{\langle \frac{w}{r} \rangle_r}$$

reweighting factor, average sign:

$$\left\langle \frac{w}{r} \right\rangle_r = \frac{Z_w}{Z_r} = e^{-\Omega \Delta f} \qquad \Delta f = f_w - f_r \ge 0$$

choose weight r to adapt to problem:

Solution Glasgow reweighting: fix β (or T)



doomed to fail ...

choose weight r to adapt to problem:

- Fodor-Katz reweighting or multi-parameter/overlap preserving reweighting
- \Rightarrow adapt β as well



$$\frac{w}{r} \sim \frac{\det M(\mu)}{\det M(0)} e^{-\Delta\beta S_{\rm YM}}$$

stay on pseudo-critical line $T_c(\mu)$

improved (ensured?) overlap: sample from both phases

Fodor-Katz reweighting: multi-parameter/overlap preserving





never repeated

breakdown of method, (un)expected role of pions?

Splittorff 07

• $Z(\mu)$ is even in μ (charge conjugation invariance) • $\langle n(\mu) \rangle \sim \frac{\partial}{\partial \mu} \ln Z$ is odd in μ

 \Rightarrow Taylor series around $\mu = 0$

Bielefeld-Swansea, Gavai-Gupta 02/05 MILC, hotOCD 10

grand-canonical ensemble $p = \frac{T}{V} \ln Z$

$$\Delta p(\mu) = p(\mu) - p(0) = \frac{\mu^2}{2!} \frac{\partial^2 p}{\partial \mu^2} \Big|_{\mu=0} + \frac{\mu^4}{4!} \frac{\partial^4 p}{\partial \mu^4} \Big|_{\mu=0} + \dots$$



determine coefficients c_{2n}

explicit expressions:

$$Z = \int DU \left(\det M\right)^{N_f} e^{-S_{\rm YM}} = \int DU e^{-S_{\rm YM} + N_f \ln \det M(\mu)}$$

straightforward:

$$\frac{\partial \ln Z}{\partial \mu} = \left\langle N_f \frac{\partial}{\partial \mu} \ln \det M \right\rangle$$
$$\frac{\partial^2 \ln Z}{\partial \mu^2} = \left\langle N_f \frac{\partial^2}{\partial \mu^2} \ln \det M \right\rangle + \left\langle \left(N_f \frac{\partial}{\partial \mu} \ln \det M \right)^2 \right\rangle$$
$$- \left\langle N_f \frac{\partial}{\partial \mu} \ln \det M \right\rangle^2$$

explicit expressions:

 $\ln \det M = \mathrm{Tr} \, \ln M$

$$\frac{\partial}{\partial \mu} \ln \det M = \operatorname{Tr} M^{-1} \frac{\partial M}{\partial \mu}$$
$$\frac{\partial^2}{\partial \mu^2} \ln \det M = \operatorname{Tr} M^{-1} \frac{\partial^2 M}{\partial \mu^2} - \operatorname{Tr} M^{-1} \frac{\partial M}{\partial \mu} M^{-1} \frac{\partial M}{\partial \mu}$$

etc.

straightforward to work out to higher order, but:

- Inumber of terms increases rapidly, $c_n \sim 6^n$ terms
- huge cancelations required: p is intensive, c_n are finite, but individual contributions may scale differently (generalized susceptibilities)

current standard:

- most groups: $\#\mu^2 + \#\mu^4 + \#\mu^6$
- $... + \#\mu^8$

Gavai-Gupta 08

• coarse lattices: $N_{\tau} = 4, 6$

only continuum extrapolated result:

equation of state to $\mathcal{O}(\mu^2)$

Borsanyi, Fodor & Katz et al 12



Method III: imaginary μ

recall: $D^{\dagger}(\mu) = \gamma_5 D(-\mu^*) \gamma_5$

- if $\mu = i\mu_{I}$, det $D(i\mu_{I})$ is real: perform ordinary simulations
- analytical continuation to real μ : $+\mu_{\rm I}^2 \rightarrow -\mu^2$
- determine phase boundary at $\mu^2 < 0$
- fit $T_c(-\mu^2)$
- obtain phase boundary
 at $\mu^2 > 0$

de Forcrand & Philipsen 02-now d'Elia & Lombardo 02 d'Elia et al 02-now



Method III: imaginary μ

- in fact: much richer than just analytical continuation
- \checkmark intricate phase structure at imaginary μ
- see below

other methods (not discussed here):

- canonical ensemble
- histograms

Summary

agreement between methods at small $\mu/T \lesssim 1$

phase boundary:



sign problem under control, fixed $N_{\tau} = 4$

indications for existence of critical endpoint?



- imaginary chemical potential: not obvious (see below)
- Taylor series: number of terms is really small estimate radius of convergence?

crossover at small μ :

- transition temperature not uniquely defined
- depends on observable
- no non-analyticity

study transition using

- Polyakov loop susceptibility
- chiral condensate
- strange quark number susceptibility

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do not have to agree: crossover region loosely defined

two scenarios

- if crossover region shrinks with increasing μ : CEP
- if crossover region extends with increasing μ : no CEP



study crossover region using

- \checkmark chiral condensate $\langle \bar{\psi}\psi \rangle$
- strange quark number susceptibility χ_s

using Taylor series expansion

Endrodi et al 11



conclusion: no indication for scenario I

Summary

standard approaches ...

- ... can be used for some questions
- are limited in applicability
- do not solve the sign problem ...