

QCD phase structure and Heavy Ion Collisions (2)

V. Bornyakov, D. Boyda, V. Goy, H.Iida, A. Molochkov, A.Nakamura, A. Nikolaev, M.Wakayama and V. I. Zakharov

Lattice and Functional Techniques for Exploration of Phase Structure and Transport Properties in Quantum Chromodynamics

Dubna, 10-14 July 2017





I am happy to be in Dubna where 114, 115 and 117 Atoms

are fund.



114 Flerovium





Riken found 113, Nihonium.

On Monday, we got a very useful relation,

$$Z(\xi, T) = \sum_{n} Z_n(T) \xi^n$$

($\xi \equiv e^{\mu/T}$: Fugacity)

We determined Z_n from Experimental data.

Today's Menu,

- 1. To show this is useful to determine Phase Transition Line especially at NICA and J-PARC
- 2. Calculate Lee-Yang Zero from these Z_n .
- 3. Calculate Z_n by Lattice QCD Simulations.



Lower estimation of larger density contribution.

2) Measured multiplicities are not Baryon, but Proton.

For Charge or Strangeness Multiplicity, No problem. I need Baryon Multiplicities ! But there are only Proton Multiplicity data !

What should I do ?

Construct NB dist. from N_p ?

M.Kitazawa and M.Asakawa, Proposal of a Method PhysRevC.86.024904 (arXiv:1205.3292)

NICA White-Paper,

theor0.jinr.ru/twiki/pub/NICA/WebHome/WhitePaper_10.01.pdf

the ratio of the susceptibility can be readily constructed by the high order correlation function for conserved quantities in high-energy nuclear collisions. Baryon numbers are conserved in strong interactions. Simulations have shown that net-proton correlation function is a fair good representation of the net-baryons [16]. Note, at the critical point, the smallness (or disappearance) of the correlation length will lead to a large value of the quark susceptibility ratio. The systematic measurements of the net-charge and net-proton Kurtosis will provide info on the nature of the system. [16] B. Mohanty, private communication, 2009.



OK, for a while, let us use Proton distribution as a Proxi of the Baryon.

Looking for High Density Not too High Energy Regions

NICA $\sqrt{s_{NN}} = 4 - 11 \text{ GeV}$ J-PARC $\sqrt{s_{NN}} = 2 - 6.2 \text{ GeV}$

FAIR $\sqrt{s_{NN}}$ =2.7 -4.9 GeV (CMB@SIS100)









J-PARC

Letter of Intent: http://j-parc.jp/researcher/Hadron/en/pac_1607/pdf/LoI_2016-16.pdf July, 2016



Jets Behavior are model dependent.



Important Observation We construct Z_n from Heavy Ion Collision data at Freeze-out T and μ . Then using

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

for arbitrary values of μ/T at fixed T





Then how RHIC data look like?

i.e., We construct Z_n from RHIC data and calculate the Moments using

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

at arbitrary values of μ/T

on

We construct Z_n on \bigstar and calculate moments

12/36

ata Susceptivility as a function of μ/T







Another Application



This ξ can be complex variables !



Lee-Yang Zeros (1952) Zeros of $Z(\xi)$ in Complex Fugacity Plane. $Z(lpha_k)=0$



But high-order polynomial zero is a famous ill-posed problem !

Read Text book of Numerical Analysis



cut Baum-Kuchen (cBK) Algorithm



50 - 100 number of significant digits

0)

A Coutour is cut into four pieces if there are zeros inside.



Lee-Yang Zeros from Experimental Data(RHIC)



Lee-Yang Zeros: RHIC Experiments



We can calculate Z_n also by Lattice QCD

But Sign Problem on Lattice ?



$$Z_{GC}(\mu, T) = \int \mathcal{D}(\text{Gluon Fields}) \\ \times \det D(\mu) \ e^{-(\text{Gluon Action})} \\ \text{Complex if } \mu \text{ is real.}$$

Our Lattice

- Clover improved Wilson action
- Iwasaki gauge action
- Substitution Lattice 4×16^3 (L \approx 3.2fm, a \approx 0.2fm)
- $m_{\pi}/m_{
 ho} = 0.8$ ($m_{\pi} = 0.7 \text{GeV}$) $T/T_c = 0.84, 0.93, 0.99, 1.08, 1.20, 1.35$

\bigcirc 20 - 40 points Im μ , 1800 - 3800 configurations at each point

- Parameters were taken from S. Ejiri et. al., PRD 82, 014508 (2010)
- Our cluster: Vostok1 (20 GPU K40)

For Pure Imaginary $\mu \quad real$



. M. D'Elia, M. P. Lombardo, Proceedings of the GISELDA Meeting held in Frascati, Italy, 14-18 January 2002, hep-lat/0205022, 22 May 2002 Ph. de Forcrand, O. Philipsen, Nucl. Phys. B642 290 (2002)16, hep-lat/02050,16 May 2002

A.Hasenfratz and Toussant, 1992

$$Z_n = \int \frac{d\theta}{2\pi} e^{i\theta n} Z_{GC}(\theta \equiv \frac{\mathrm{Im}\mu}{T}, T)$$

All information is in Imaginary Chemical Potential regions! Great Idea ! But practically it did not work.

For few years, we must develop several Engineering Methds.

1) Integration method 2) Multi-Precision Calculations



Integration Method

$$n_B = \frac{1}{3V} T \frac{\partial}{\partial \mu} \log Z_G$$
$$= \frac{N_f}{3N_s^3 N_t} \int \mathcal{D}U e^{-S_G} \operatorname{Tr} \Delta^{-1} \frac{\partial \Delta}{\partial \mu} \det \Delta$$

(For pure imaginary μ , n_B is also imaginary)

Then, for fixed T

$$Z(\theta \equiv \frac{\mu}{T}) = \exp(V \int_0^\theta n_B d\theta')$$

$$Z_k = \frac{3}{2\pi} \int_{-\pi/3}^{+\pi/3} d\theta \exp\left(i\,k\theta + \int_0^\theta n_B d\theta'\right)$$

We map Information in Pure Imaginary Chemical Potential to Real ones.

See also, D'Elia, Gagliardi and Sanfilippo, Phys. Rev. D 95, 094503 (2017)

Roberge-Weiss

 $\xi \equiv e^{\mu/2}$

We measure the number density at many pure imaginary chemical potential $n_B(\mu_I)$.

We construct Grand Partition Function Z_G , by integrating $\,n_B(\mu_I)\,$

 ${\ensuremath{{}^{\hspace*{-.5mm} \ensuremath{}^{\hspace*{-.5mm} \ensuremath{}^{\hspace*$

F Then we can calculate Real μ regions by

$$Z(\xi, T) = \sum Z_n(T) \,\xi^n$$

n







References

A.Li et al.(Kentucky), Phys.Rev.D82:054502,2010, arXiv:1005.4158

A.Suzuki et al.(Zn Collaboration), Lattice 2016 Proceedings,

V.Goy et al.(Vladivostok), Prog Theor Exp Phys (2017) (3): 031D01,arXiv:1611.08093 Where comes the phase of \mathcal{Z}_{n} ?

A.Li et al.(Kentucky), Phys.Rev.D82:054502,2010, arXiv:1005.4158

$$Z = \int \mathcal{D}U \left(\det \Delta(\mu)\right)^{N_f} e^{-S_G} = e^{\log(1-\kappa Q)}$$

$$\det \Delta(\mu) = \det(1-\kappa Q(\mu))$$

$$= \exp\left(A_0 + \sum_{n>0} [e^{in\phi}W_n + e^{-in\phi}W_n^{\dagger}]\right)$$

$$= \exp\left(A_0 + \sum_n A_n \cos(n\phi + \delta_n)\right)$$

$$A_n \equiv 2|W_n| \qquad \text{We use } W_{-n} = W_n$$

$$\delta_n \equiv \arg(W_n)$$

Then,

$$Z_n \propto \int_0^{2\pi} \frac{d\phi}{2\pi} e^{-ik\phi} e^{A_0 + A_1 \cos(\phi + \delta_1) + A_1 \cos(2\phi + \delta_2) \cdots}$$

In the lowest order,

$$\int_{0}^{2\pi} \frac{d\phi}{2\pi} e^{-ik\phi} e^{A_{0}+A_{1}\cos(\phi+\delta_{1})} = e^{A_{0}} \int_{\delta_{1}}^{2\pi+\delta_{1}} \frac{d\phi'}{2\pi} e^{-ik(\phi'-\delta_{1})} e^{A_{1}\cos\phi'}$$

$$= e^{A_{0}+ik\delta_{1}} \int_{\delta_{1}}^{2\pi+\delta_{1}} \frac{d\phi'}{2\pi} e^{-ik\phi'} e^{A_{1}\cos\phi'}$$

$$= e^{A_{0}+ik\delta_{1}} \int_{0}^{2\pi} \frac{d\phi'}{2\pi} e^{-ik\phi'} e^{A_{1}\cos\phi'}$$

$$= e^{A_{0}+ik\delta_{1}} I_{k}(A_{1})$$

 $\propto z_k$

where we use

$$I_n(z) = \frac{(-1)^n}{2\pi} \int_0^{2\pi} e^{z \cos t} e^{-int} dt$$

A Remark of Function Form of $n_B(\mu_I)$

Preliminary

n_B/T³



 $n_B(\mu_I)$ is well approximated by sine function at *T*<*Tc*.

Takahashi et al. Phy. Rev. D 91 (1) (2015) 014501. Bornyakov et al., Phys.Rev. D95, 094506 (2017)

A Remark of Function Form of $n_B(\mu_I)$

Preliminary

n_B/T³



 $n_B(\mu_I)$ is well approximated by sine function at *T*<*Tc*.

Takahashi et al. Phy. Rev. D 91 (1) (2015) 014501. Bornyakov et al., Phys.Rev. D95, 094506 (2017) In general,

$$n_B/T^3 = \sum_k f_{3k} \sin(k\theta_I)$$

f3 = 0.0871(3), f6 = -0.00032(27) (\chi_2/\dof = 0.93)

Lowest order,



This is Skellam Model, which is used in Heavy Ion Collisions to describe the gross structure.

Skellam is the difference of two independent Poisson Distributions. Structures of the dynamics.

Concluding Remarks

We have developed the Canonical Approach for revealing QCD Phase Structure. The canonical partition functions Z_n drop very rapidly as *n* goes large, and we need multiprecision calculations.

The phase of \mathcal{Z}_n fluctuates rapidly as *n* goes large in the confinement phase. No such problem in the deconfinement phase. i.e., we plan to predict thermo-dynamical quantities

for LHC.

☆Our Quark masses are heavy. Now it is time to go towards Physical Parameters. Coming Soon !