

# Lattice QCD: a brief introduction

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## Study of QCD thermodynamics

Theoretical investigations of QCD thermodynamics make use of different methods and tools

### From first principles:

- **Lattice QCD**
- Perturbation theory (large  $T$  and/or  $\mu$ )
- Functional methods (functional renormalization group - FRG, Dyson-Schwinger equations, etc...)

### Models:

- Nambu-Jona-Lasinio (NJL) -type models (Nambu and Jona-Lasinio, *Phys. Rev.* **122** (1961) 345, *Phys. Rev.* **124** (1961) 246)
- Hadron Resonance Gas (HRG) -type models (Hagedorn, *Nuovo Cim. Suppl.* **3** (1965), 147)
- ...

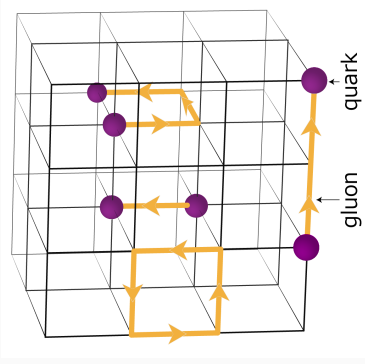
## Lattice formulation of QCD

In the non-perturbative regime where the coupling  $g_s$  is not small, lattice QCD is the major tool of investigation of equilibrium (thermodynamic) properties of QCD

- Essentially, it amounts to calculating path integrals of the like:

$$Z[U, \bar{\psi}, \psi](V, T, \mu) = \int \mathcal{D}U \mathcal{D}\bar{\psi} \mathcal{D}\psi e^{-S_G[U, \bar{\psi}, \psi] - S_F[U, \bar{\psi}, \psi]}$$

- $S_G$  and  $S_F$  are the *Euclidean* gauge (gluonic) and fermionic actions,  $U$  the gauge fields,  $\psi, \bar{\psi}$  the fermion fields.
- The theory is defined on a discretized 3+1d lattice of size  $N_s^3 \times N_\tau$ , with lattice spacing  $a$



2/19

## Lattice formulation of QCD

- Euclidean* actions  $S_G$  and  $S_F$  only depend on coupling  $g$  and fermion masses  $m_f$
- Periodic boundary conditions for bosonic fields  $U$ , anti-periodic for fermionic fields  $\psi, \bar{\psi}$  (Grassmann variables)

- The finiteness of the lattice spacing  $a$  serves as a *regulator* for the theory. At the end one wishes to recover the continuum theory with  $\lim_{a \rightarrow 0} (\lim_{N_\tau \rightarrow \infty})$ : **continuum limit**

- Calculations are done in finite volume. When possible, one wishes to study the

thermodynamic limit  $\lim_{V \rightarrow \infty}$ : a.k.a. **infinite volume limit**

- Simulations at finite temperature have set length in “temporal” direction

$$T^{-1} = \beta = N_\tau a$$

- Summarizing

$$V = (N_s a)^3, \quad T = \frac{1}{N_\tau a}$$

- NOTE:** the integrals **cannot** be calculated by brute force! Even for a small  $10^4$  lattice, integral is 320000-dimensional!  $\Rightarrow$  **Importance sampling**

3/19

## Lattice QCD at $\mu_B = 0$

Thanks to asymptotic freedom (and perturbative calculations) and confinement, we expect a phase transition between *confined* and *deconfined* matter

- At low  $T$ /low  $\mu$ :
- **Confined matter**  $\langle \vec{L} \rangle = 0$
- **Chiral symmetry is spontaneously broken** by a non-zero chiral condensate

$$\langle \bar{\psi}\psi \rangle = \frac{T}{V} \frac{\partial}{\partial m_f} \log \mathcal{Z}(T, V, \mu) \neq 0$$

- At high  $T$ /high  $\mu$ :
- **Deconfined matter**  $\langle \vec{L} \rangle \neq 0$
- **Chiral symmetry is restored**

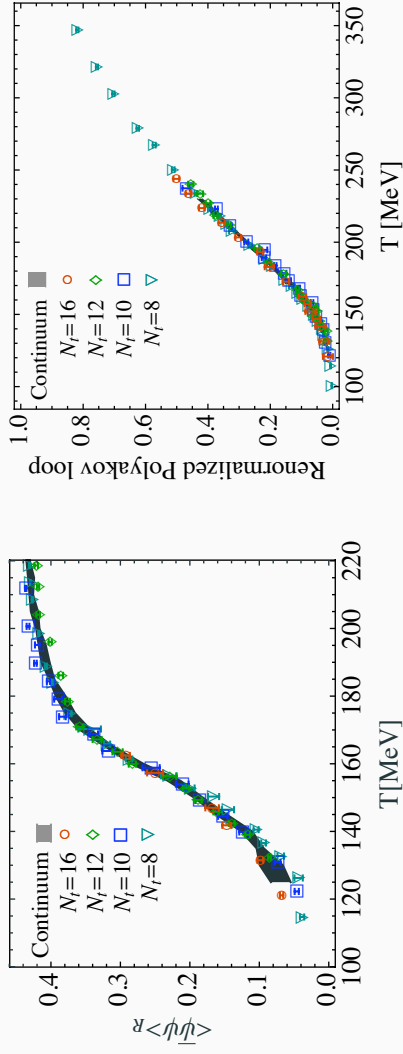
$$\langle \bar{\psi}\psi \rangle = \frac{T}{V} \frac{\partial}{\partial m_f} \log \mathcal{Z}(T, V, \mu) = 0$$

**NOTE:** although deconfinement and chiral transition are two “different” transitions, we have no evidence of them taking place at different temperatures. Unique  $T_C$ !

4/19

## The QCD transition: observables

Both observables are able to distinguish between the two phases:



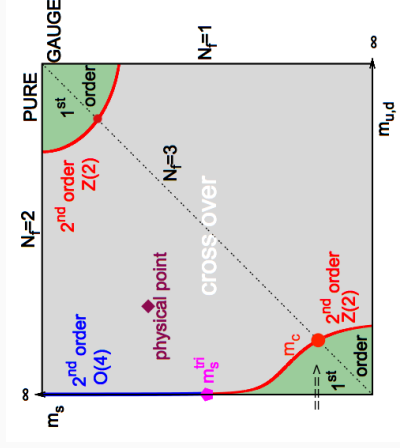
Borsanyi et al. **JHEP 1009:073 (2010)**

**NOTE:** this chiral condensate is renormalized subtracting the zero-temperature value

5/19

## The QCD transition: Columbia plot

As a function of the light (u,d) and strange quark masses, the order of the transition changes



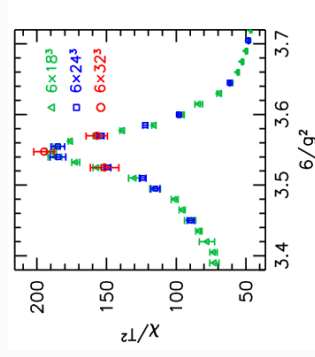
- At the physical point  $m_s/m_{ud} \simeq 27$ , the transition is a smooth crossover!
- In the heavy-quark limit (pure gauge), the transition is first order

6/19

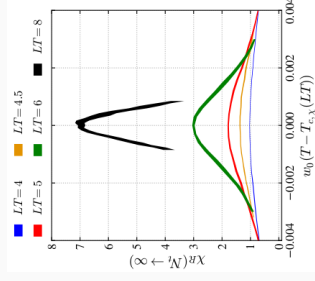
## The QCD transition: crossover vs. first order

On the lattice we study the volume scaling of certain quantities to determine the order of the transition

**Left:** physical masses



**Right:** infinite masses (pure gauge)



- For a crossover (left), the peak height is independent of the volume
- For a first order transition, it scales linearly with the volume

7/19

## Thermodynamic description of QCD

The thermodynamics of QCD is commonly investigated in the grand canonical ensemble

- Grancanonical partition function:

$$\mathcal{Z} = \sum_N Z_N e^{\mu N}$$

where:

- $Z_N$  is the canonical partition function with  $N$  particles
- $\mu$  is the *chemical potential* associated to the particle number  $N$

The chemical potential is the energy associated to a change in the number  $N$  of particles

- In QCD, there are 3 conserved “particle numbers”, which one can see as:

- **Quark numbers:**  $u, d, s$

$$\mu N = \mu_u N_u + \mu_d N_d + \mu_s N_s$$

- **Conserved charges:**  $B, Q, S$

$$\mu N = \mu_B N_B + \mu_Q N_Q + \mu_S N_S$$

8/19

## Thermodynamic description of QCD

The **Equation of State (EoS)** is extremely important since it completely describes the equilibrium properties of QCD matter.

It is one of the main inputs to hydro and several other tools for calculations in heavy-ion collisions and higher-density physics.

Thermodynamic quantities follow directly from the grandcanonical partition function  $\mathcal{Z}$  and the relation:

$$-k_B T \ln \mathcal{Z} = U - TS - \mu N$$

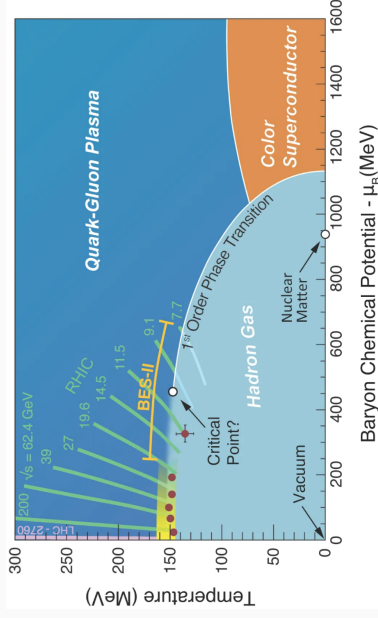
- **Pressure:**  $p = -k_B T \frac{\partial \ln \mathcal{Z}}{\partial V}$
- **Entropy density:**  $s = \left( \frac{\partial p}{\partial T} \right)_{\mu_i}$
- **Charge densities:**  $n_i = \left( \frac{\partial p}{\partial \mu_i} \right)_{T, \mu_{j \neq i}}$
- **Energy density:**  $\epsilon = Ts - p + \sum_i \mu_i n_i$
- **Speed of sound:**  $c_s^2 = \left( \frac{\partial p}{\partial \epsilon} \right)_{s/n_B}$
- More **(Fluctuations, etc...)**

9/19

## The phase diagram of QCD

Different phases of QCD matter (in equilibrium) are depicted in (temperature vs baryo-chemical potential) phase diagram

- **Hadron gas** at low- $T$  and/or low- $\mu_B$
- **Quark Gluon Plasma (QGP)** at large  $T$  and (possibly) at large  $\mu_B$
- **More exotic phases** proposed at low- $T$  and high- $\mu_B$  (color superconductivity, etc...)

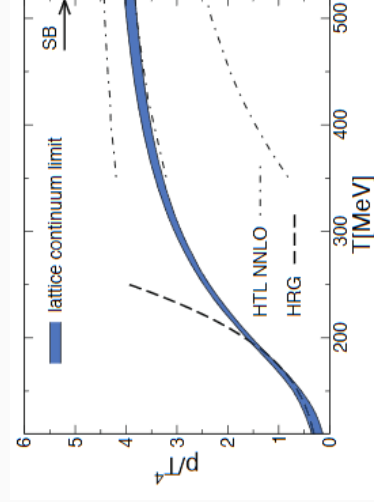


10/19

## Equation of state at $\mu_B = 0$

A combination of methods gives us good understanding of the EoS at  $\mu_B = 0$  at all temperatures

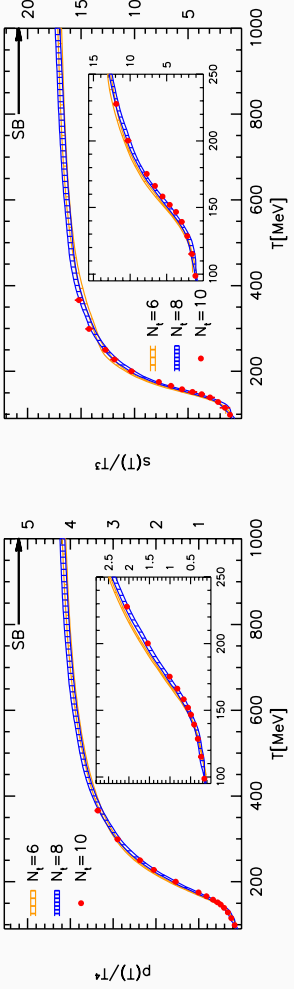
- Perturbative QCD at high temperature  
→ “pure quark-gluon phase”
- HRG model at low temperature  
→ “pure hadron phase”
- **Lattice QCD bridges between regimes and captures the transition**



Borsányi *et al.*, **PLB 370 (2014) 99-104**

11/19

## Lattice QCD: equation of state at $\mu_B = 0$



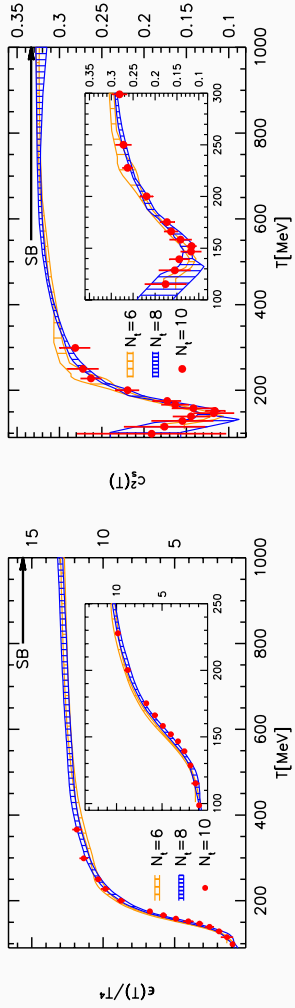
Stefan-Boltzmann limit for free gas (large T):

$$p_{\text{QCD}} = \frac{\pi^2}{45} T^4 (N_c^2 - 1) + \sum_{\text{flavours}} \frac{N_c}{3\pi^2} \left[ \frac{7\pi^4 T^4}{60} + \frac{\mu_f^2 \pi^2 T^2}{2} + \frac{\mu_f^4}{4} \right]$$

Borsányi *et al.*, JHEP 11 (2010) 077

12/19

## Lattice QCD: equation of state at $\mu_B = 0$



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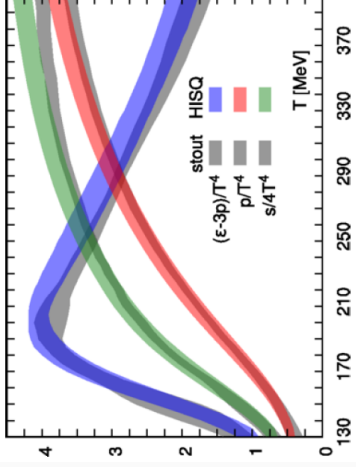
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Borsányi *et al.*, JHEP 11 (2010) 077

13/19

## Lattice QCD: equation of state at $\mu_B = 0$

- Current lattice simulations are performed with realistic setup
- Great agreement between different collaborations
- This means systematics are well under control and results are extremely reliable



Borsányi *et al.*, PLB 370 (2014) 99-104, Bazavov *et al.* PRD 90 (2014) 094503

14/19

## The sign/complex action problem

Euclidean path integrals are calculated with MC methods using importance sampling, interpreting the factor  $\det M[U] e^{-S_G[U]}$  as the Boltzmann weight for the configuration  $U$

$$\begin{aligned} Z(V, T, \mu) &= \int \mathcal{D}U \mathcal{D}\psi \mathcal{D}\bar{\psi} e^{-S_F(U, \psi, \bar{\psi}) - S_G(U)} \\ &= \int \mathcal{D}U \det M(U) e^{-S_G(U)} \end{aligned}$$

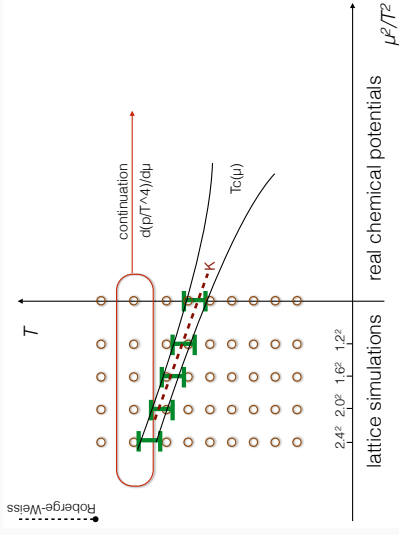
- If there is particle-antiparticle-symmetry ( $\mu = 0$ )  $\det M(U)$  is real
- For real chemical potential ( $\mu^2 > 0$ )  $\rightarrow \det M(U)$  is complex (**complex action problem**) and has wildly oscillating phase (**sign problem**)  
 $\Rightarrow$  It cannot serve as a statistical weight
- For *purely imaginary* chemical potential ( $\mu^2 < 0$ )  $\rightarrow \det M(U)$  is real again, simulations can be made!

15/19



## Quick note: simulations at imaginary chemical potential

- While for real chemical potential ( $\mu^2 > 0$ )  $\det M(U)$  is complex, for **imaginary** chemical potential ( $\mu^2 < 0$ )  $\det M(U)$  is real
- Perform simulations at imaginary chemical potentials:
- **Analytically continue** to  $\mu^2 > 0$

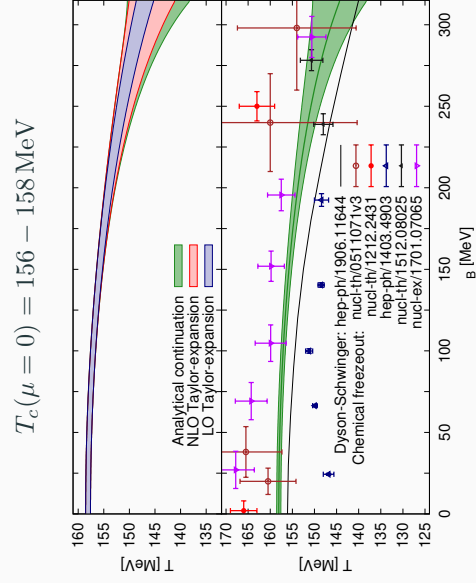


We can do this because the transition at  $\mu = 0$  is a crossover, so the partition function is analytic!

16/19

## The transition at finite chemical potential

Current results for transition temperature give:



17/19

## The width of the transition at finite chemical potential

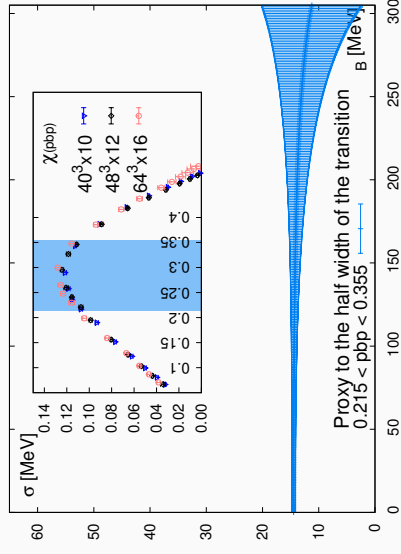
We get at  $\mu_B = 0$  for  $T_c$ :

$$T_c(LT = 4, \mu_B = 0) = 158.0 \pm 0.6 \text{ MeV}$$

while for the width

$$\Delta T(LT = 4, \mu_B = 0) = 15.0 \pm 1.0 \text{ MeV}$$

The width of the transition has a very mild chemical potential dependence



18/19

## Summary

- Lattice QCD is a regularization scheme for QCD that does not rely on a perturbative expansion
- Thermodynamic results show that chiral restoration and deconfinement apparently at same temperature  $T_c \simeq 156 - 158 \text{ MeV}$
- At physical quark masses, transition is a smooth crossover (at infinite quark masses, it is of first order)
- Equation of state of QCD known to high precision, shows liberation of degrees of freedom and approach to Stefan-Boltzmann limit
- Lattice QCD cannot at the moment give *direct* results at non-zero chemical potential (sign problem), but..
- ..thanks to extrapolations, we know the location of the phase transition line in the phase diagram (among other things)

19/19