

Title (even content) has been changed to

Phase structure of many flavor lattice QCD at finite temperature

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in collaboration with

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Introduction

- The nature of the chiral phase transition of Many Flavor QCD [MFQCD] depends on the number of flavors and the masses.

Pisarski and Wilczek, PRD 29, 338 (1984) and many lattice calcs.

- EW baryogenesis in TC models.

Appelquist, Schwetz and Selipsky, PRD52, 4741 (1995);

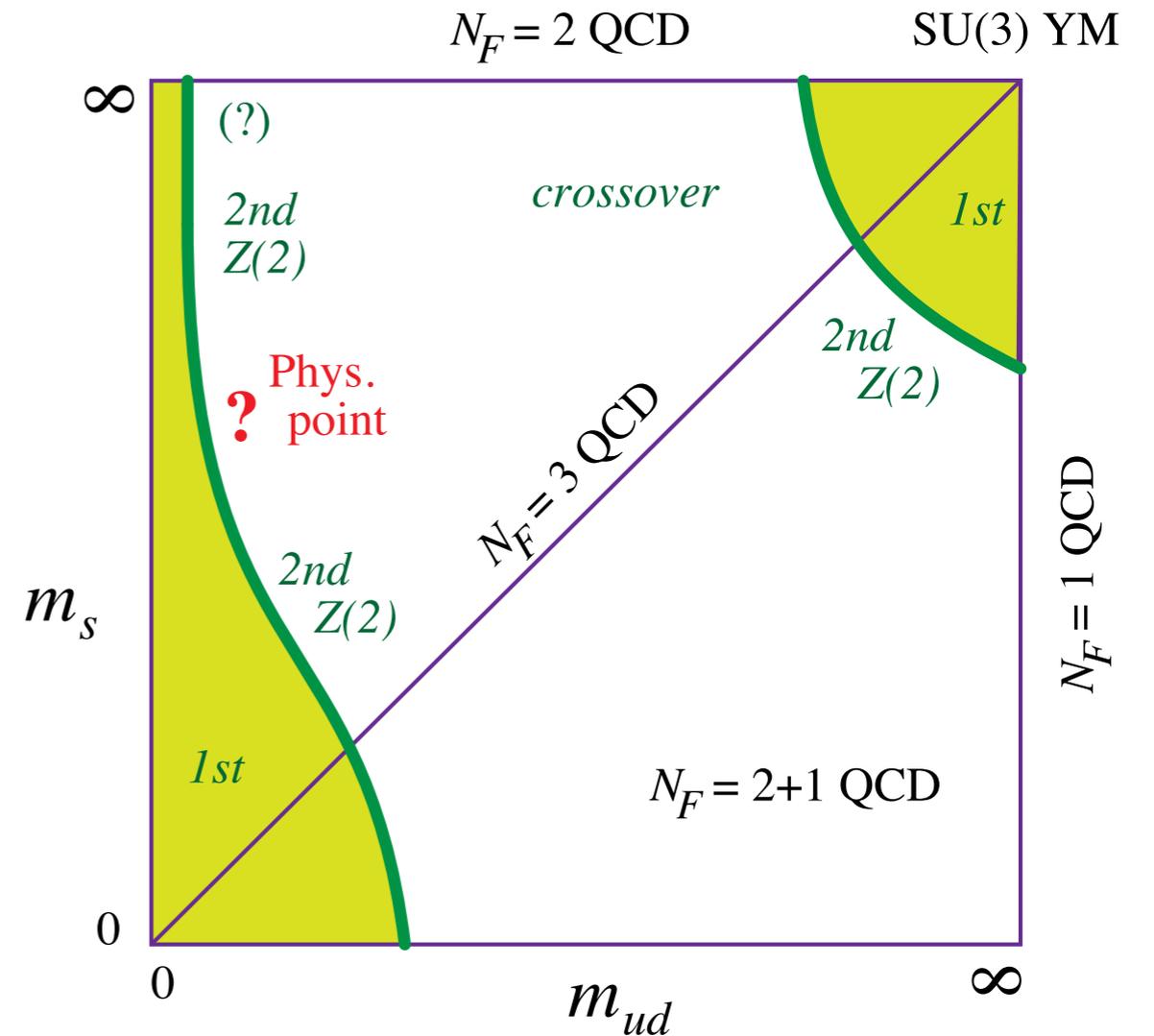
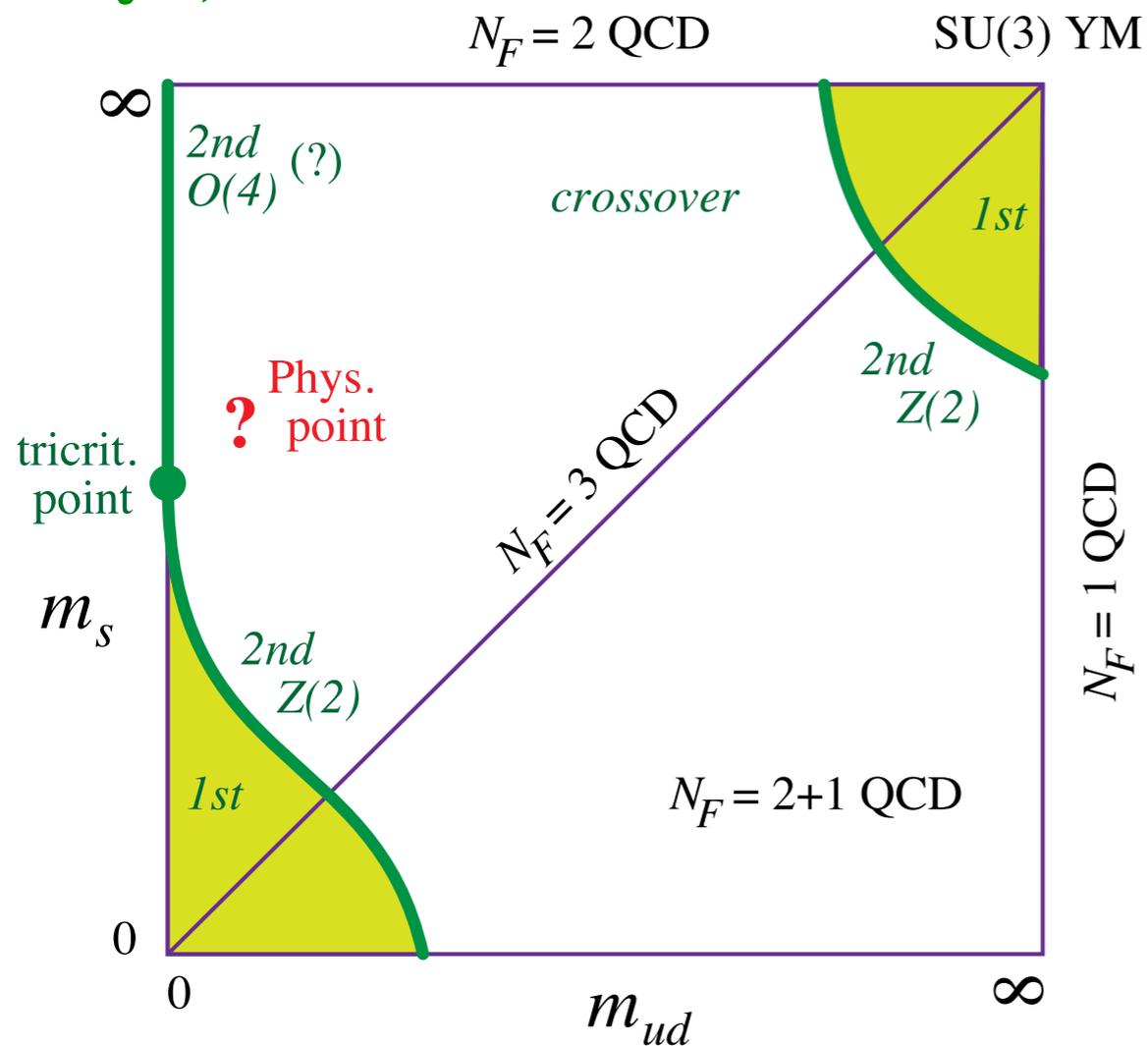
Kikukawa, Kohda and Yasuda, PRD77 (2008) 015014

- Towards the study of this possibility, we consider $2(\text{light})+N_f(\text{heavy})\text{-flavor QCD}$, propose an easy way to explore the phase structure and demonstrate the feasibility of the method.

Columbia plot for 2+1 QCD

Brown, Butler, Chen, Christ, Dong, Schaffer, Unger, and Vaccarino (90),
 N.H. Christ, Z. Dong (92) and N.H. Christ(92)

Kanaya, Lattice 2010

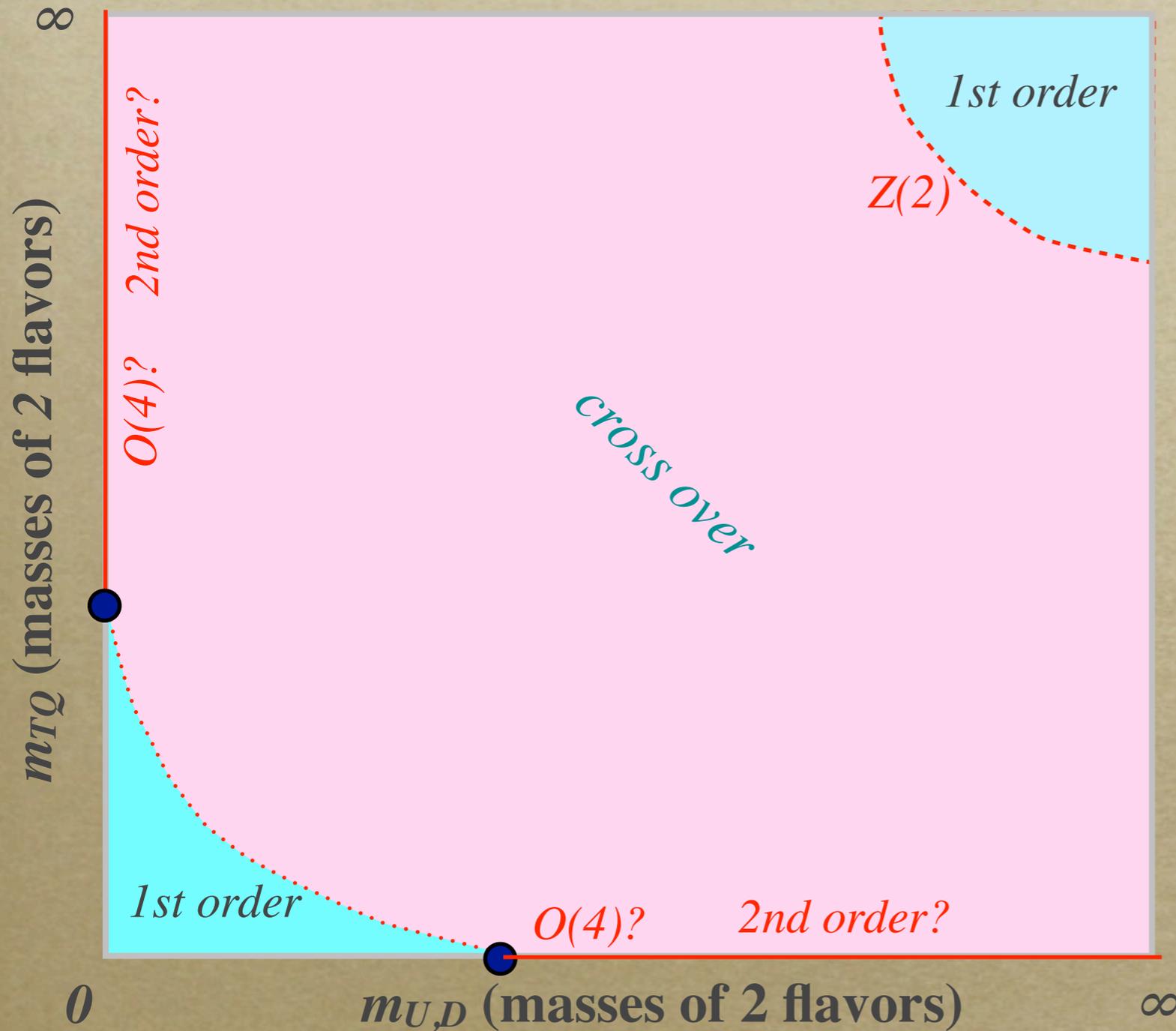


When $N_F \geq 3$ (or 2?), Chiral Phase Transition around the massless limit is 1st order.

Pisarski and Wilczek, PRD 29, 338 (1984); Iwasaki et al. (1997). See also Aoki, Fukaya, Taniguchi (2012)

Extend Columbia plot to $2+N_f$ QCD.

$$N_f = 2 \quad (2+2 < N_f^{\text{crit}})$$



In TC, two flavors must be exact massless.

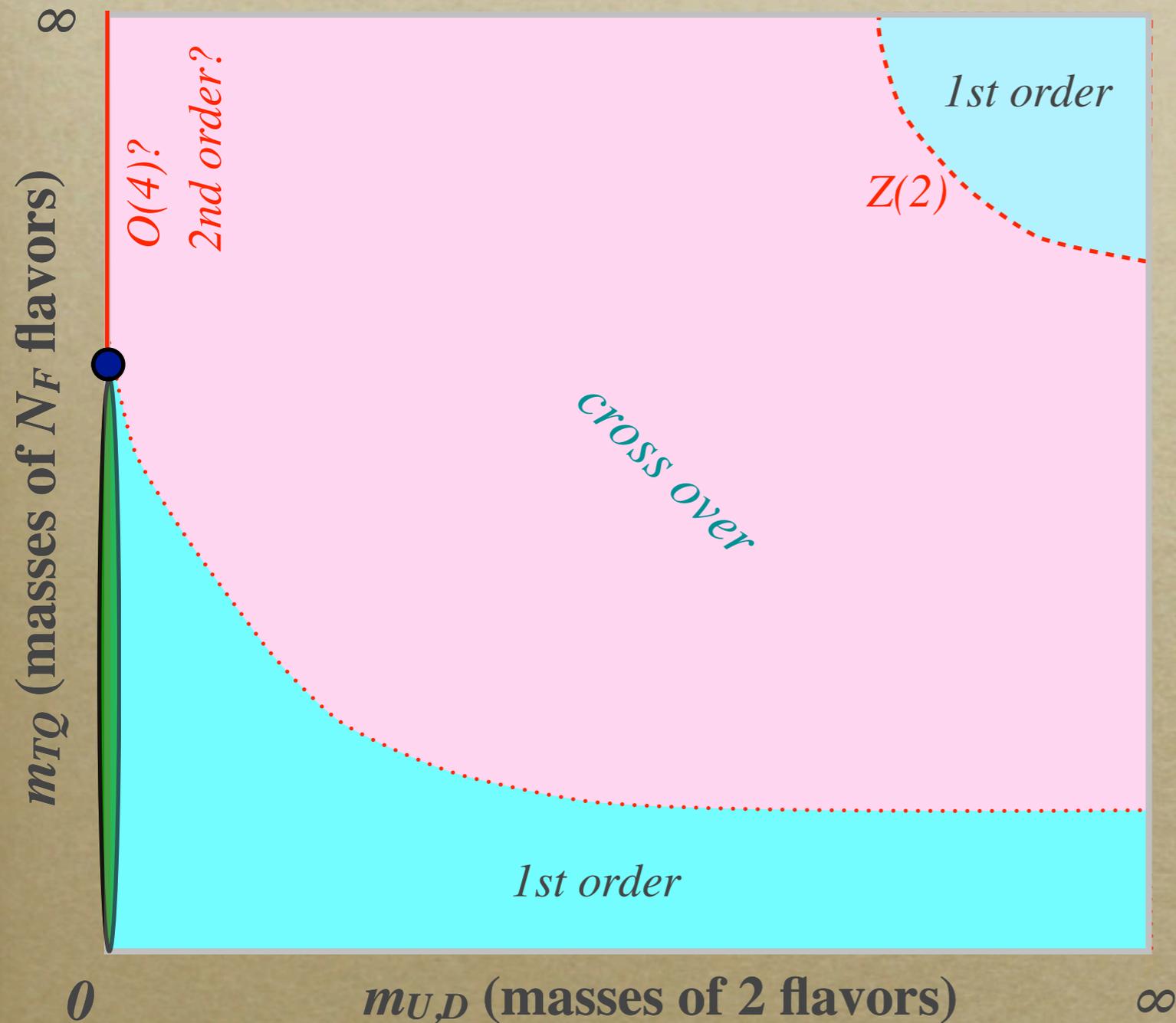
N_f^{crit} separating confining and conformal theories

Symmetric under reflection w.r.t. diagonal line.

Running of g^2 is not slow enough.

Less interesting.

$$3 \equiv N_f (2 + N_f < N_f^{\text{crit}})$$



1st order persists to $m_{U,D} = \infty$ for sufficiently small m_{TQ} .

Walking with $\gamma_m \sim O(1)$ is expected at an appropriate N_f .

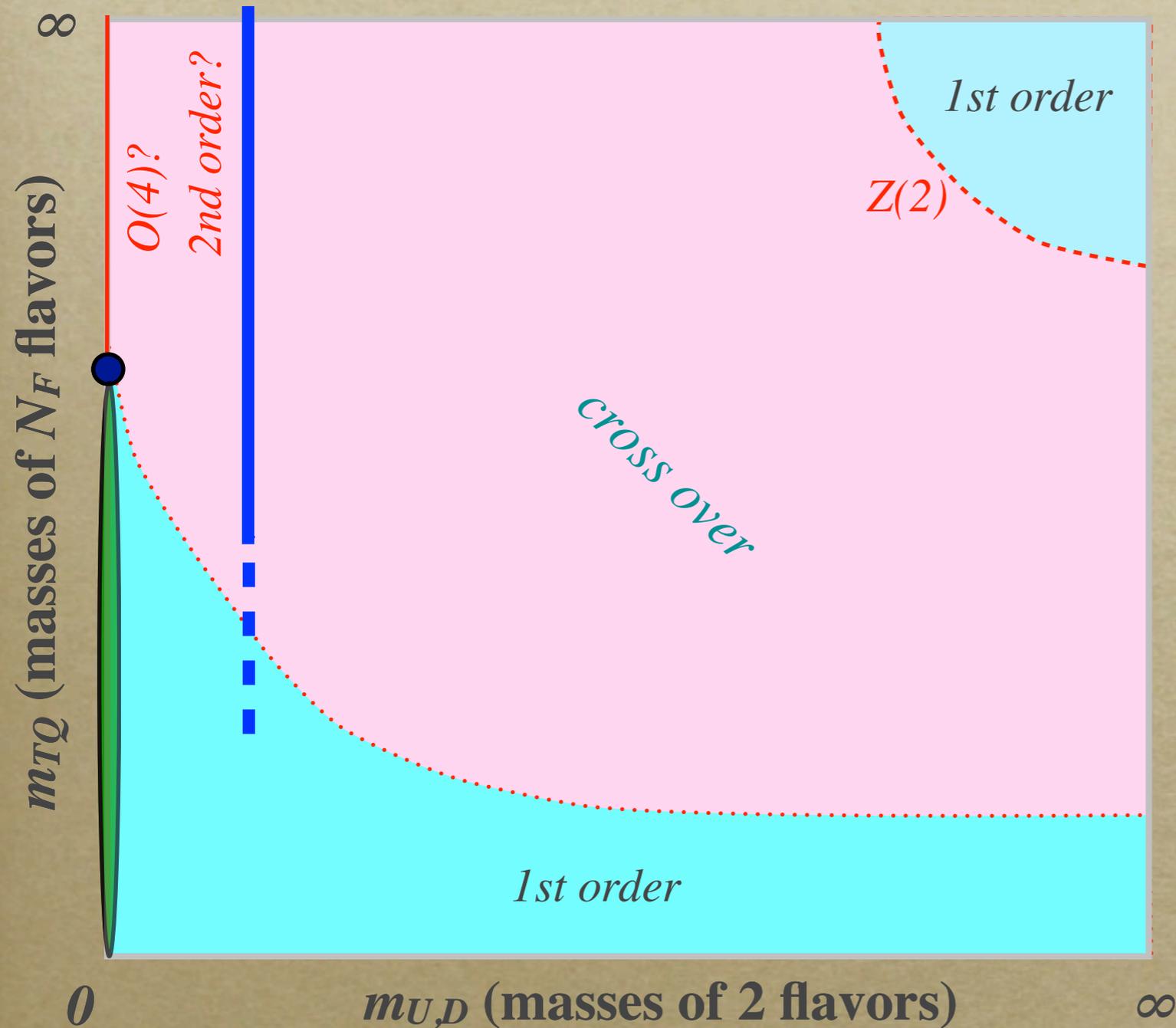
For EW baryogenesis, important to identify the location of the 1st order region and its strength.

Appelquist, Schwetz and Selipsky, PRD52, 4741 (1995).

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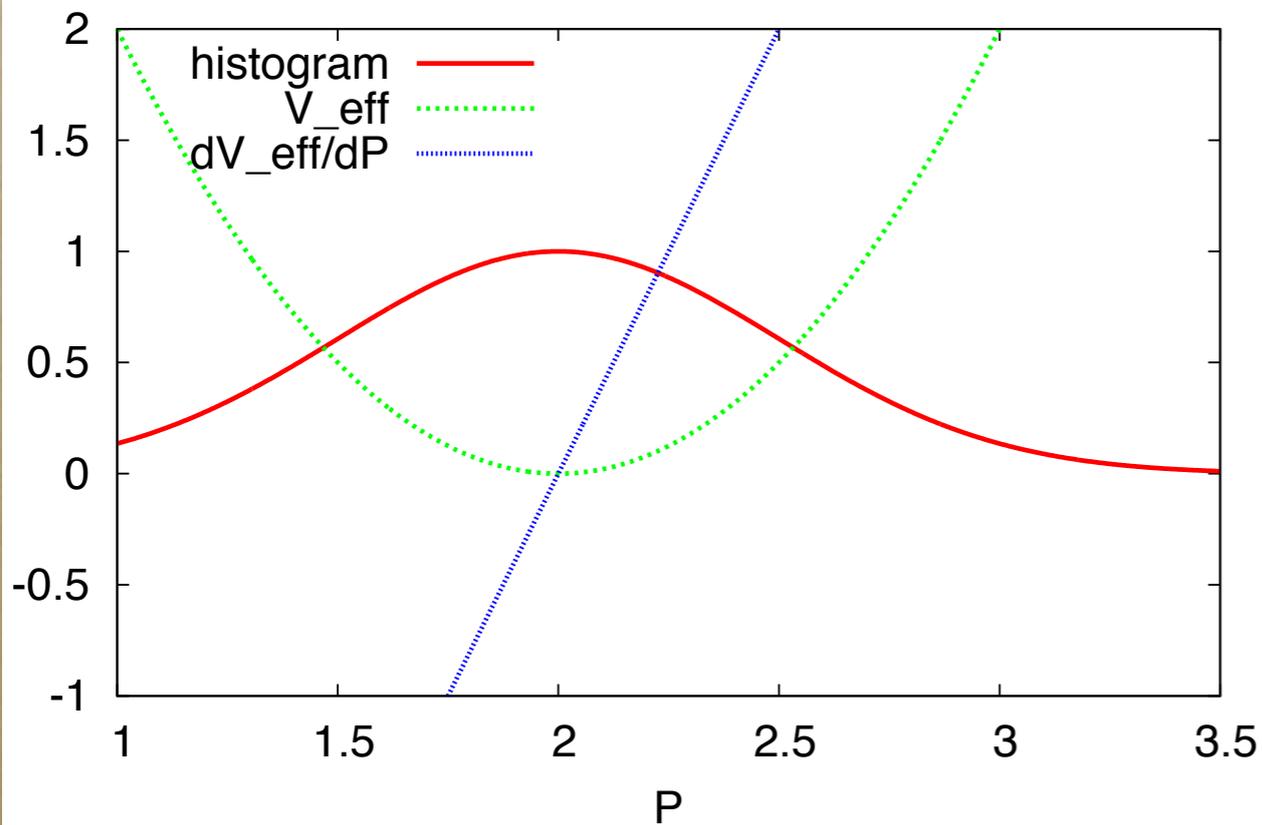
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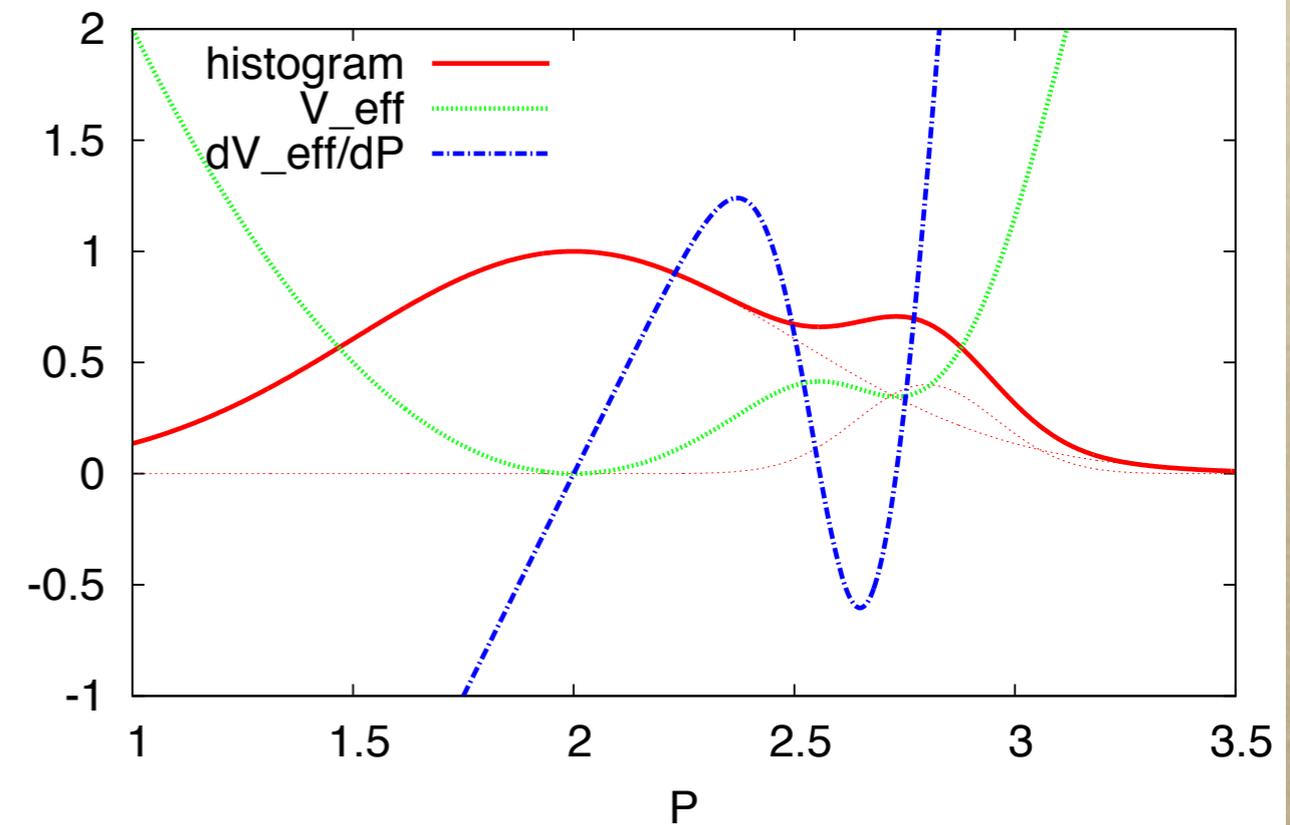
Effective potential through histogram

S. Ejiri, PRD77, 014508 (2008); Saito et al, [WHOT-QCD], PRD84, 054502 (2011)

Example 0



Example 1



$$w(P, \beta) = \langle \delta(P - \hat{P}) \rangle_{\beta}$$

$$V_{\text{eff}}(P, \beta) = -\ln w(P, \beta)$$

$$= -\ln w(P, \beta_0) + 6(\beta - \beta_0)N_{\text{site}}P$$

Double well potential
 = 1st order PT
 \Rightarrow look for $\partial^2 V_{\text{eff}}/\partial P^2 \leq 0$

Re-weighting method

- Start with the partition function with $2+N_f$ -flavor QCD.

$$Z(\beta, m_f, \mu_f) = \int \mathcal{D}U \mathcal{D}\psi \mathcal{D}\bar{\psi} e^{-S_q - S_g} = \int \mathcal{D}U e^{6\beta N_{\text{site}} \hat{P}} \prod_{f=1}^{N_f+2} (\det M(m_f, \mu_f)),$$

$$\prod_{f=1}^{N_f+2} (\det M(m_f, \mu_f)) \rightarrow (\det M(m_1, 0))^2 \left(\frac{\det M(m_1, \mu)}{\det M(m_1, 0)} \right)^2 \prod_{h=1}^{N_f} \left(\frac{\det M(m_h, 0)}{\det M(\infty, 0)} \right)$$

- N_f flavors are heavy and the chemical potential for u and d quarks are small.

$$\ln \left[\frac{\det M(m_1, \mu)}{\det M(m_1, 0)} \right] = \sum_{n=1}^{N_\mu} \frac{1}{n!} \left[\frac{\partial^n (\ln \det M)}{\partial (\mu/T)^n} \right] \left(\frac{\mu}{T} \right)^n$$

M : quark matrix

P : plaquette

Ω : Re[Polyakov loop]

$$\ln \left[\frac{\det M(\kappa_h, 0)}{\det M(0, 0)} \right] = 288 N_{\text{site}} \kappa_h^4 P + 12 N_s^3 (2\kappa_h)^{N_t} \Omega + \dots$$

Re-weighting method

Expectation value in $2+N_f$ -flavor QCD and its μ - and m_h -dependence can be calculated by ensemble average over two-flavor QCD configurations upto truncation errors.

$$V_{\text{eff}}(P, \beta, m_l, m_h, \mu) = -\ln w^{(2+N_f)}(P, \beta, m_l, m_h, \mu)$$

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$$V_{\text{eff}}(P, \beta, m_h, \mu) = V_0(P, \beta_0) - \ln R(P; \beta, m_1, m_h, \mu; \beta_0)$$

Two-flavor part

Always single well

Effect of additional heavy flavors

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$$\ln R(P; \beta, m_h, \mu; \beta_0) = 6(\beta - \beta_0)N_{\text{site}}P$$

$$+ \ln \frac{\left\langle \delta(P - \hat{P}) \left(\frac{\det M(m_1, \mu)}{\det M(m_1, 0)} \right)^2 \prod_{h=1}^{N_f} \left(\frac{\det M(m_h, 0)}{\det M(\infty, 0)} \right) \right\rangle_{\text{two-flavors}, \beta}}{\left\langle \delta(P - \hat{P}) \right\rangle_{\text{two-flavors}, \beta}}$$

Heavy quark mass and N_f dependence

$$\ln R(P; \beta, \kappa_h, 0; \beta_0) = \ln \bar{R}(P; \kappa_h, 0) + (\text{plaquette term}) + O(\kappa_h^{N_t+2})$$

$$\bar{R}(P; \kappa_h, 0) = \frac{\left\langle \delta(P - \hat{P}) \exp[6hN_s^3\Omega] \right\rangle_\beta}{\left\langle \delta(P' - P) \right\rangle_\beta}$$

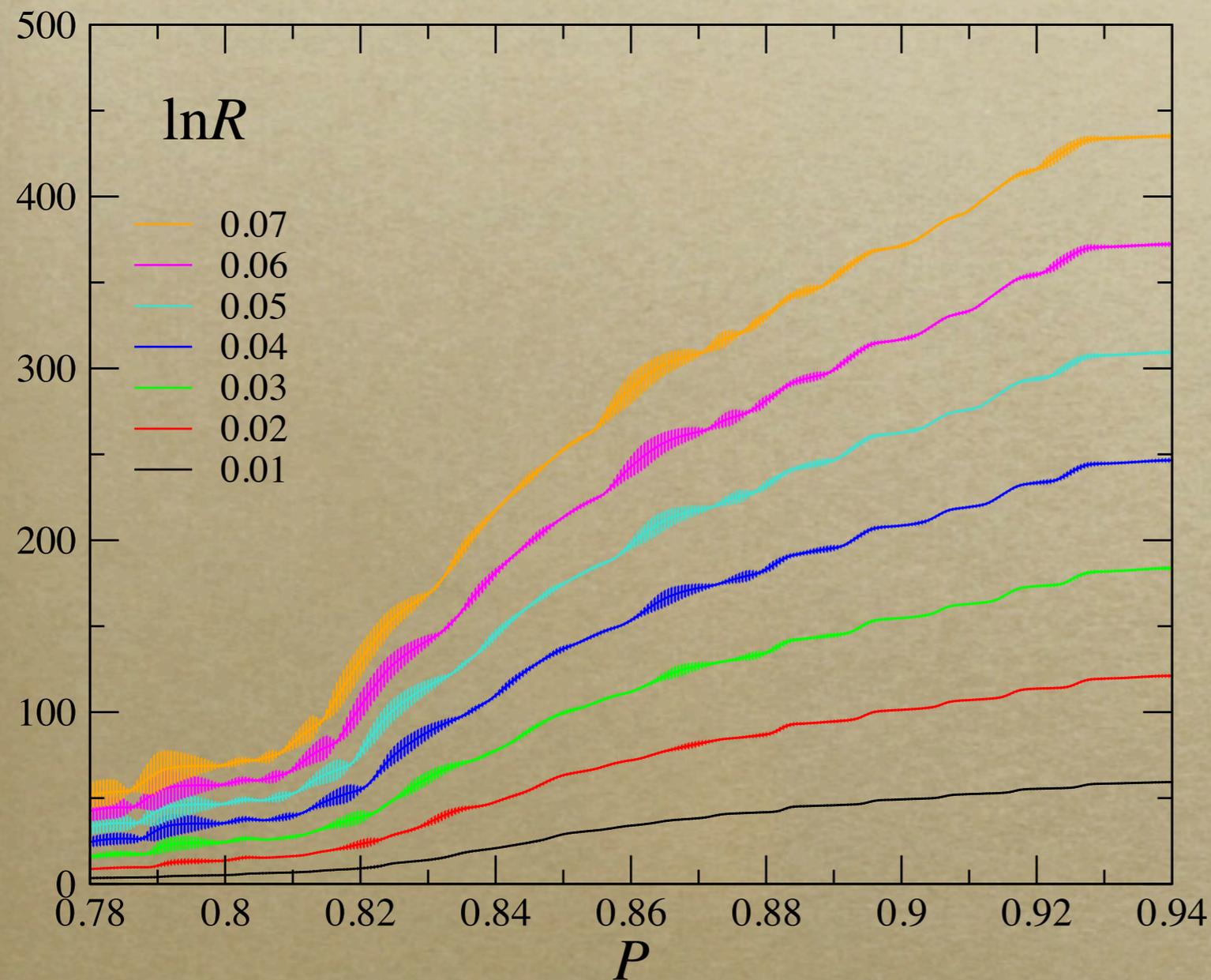
where

$$h = 2N_f(2\kappa_h)^{N_t} \quad \text{for the Wilson quark action}$$

$$h = N_f/(4 \times (2m_h)^{N_t}) \quad \text{for the staggered quark action}$$

Heavy quark dependence and N_f -dependence can be parameterized in a single parameter h .

Result: $\ln R$



Two-flavor QCD configurations:
p4-improved staggered quark &
the standard plaquette gauge

$a m_1 = 0.1$,

10,000-40,000 traajs.

$V=16^3 \times 4$,

16 β in [3.52, 4.00],

$T/T_c = [0.76, 1.98]$,

$M_{PS}/M_V \sim 0.7$

[C.R. Allton, et al., PRD71,054508 (2005)]

Calculated with $h = [0.01, 0.07]$

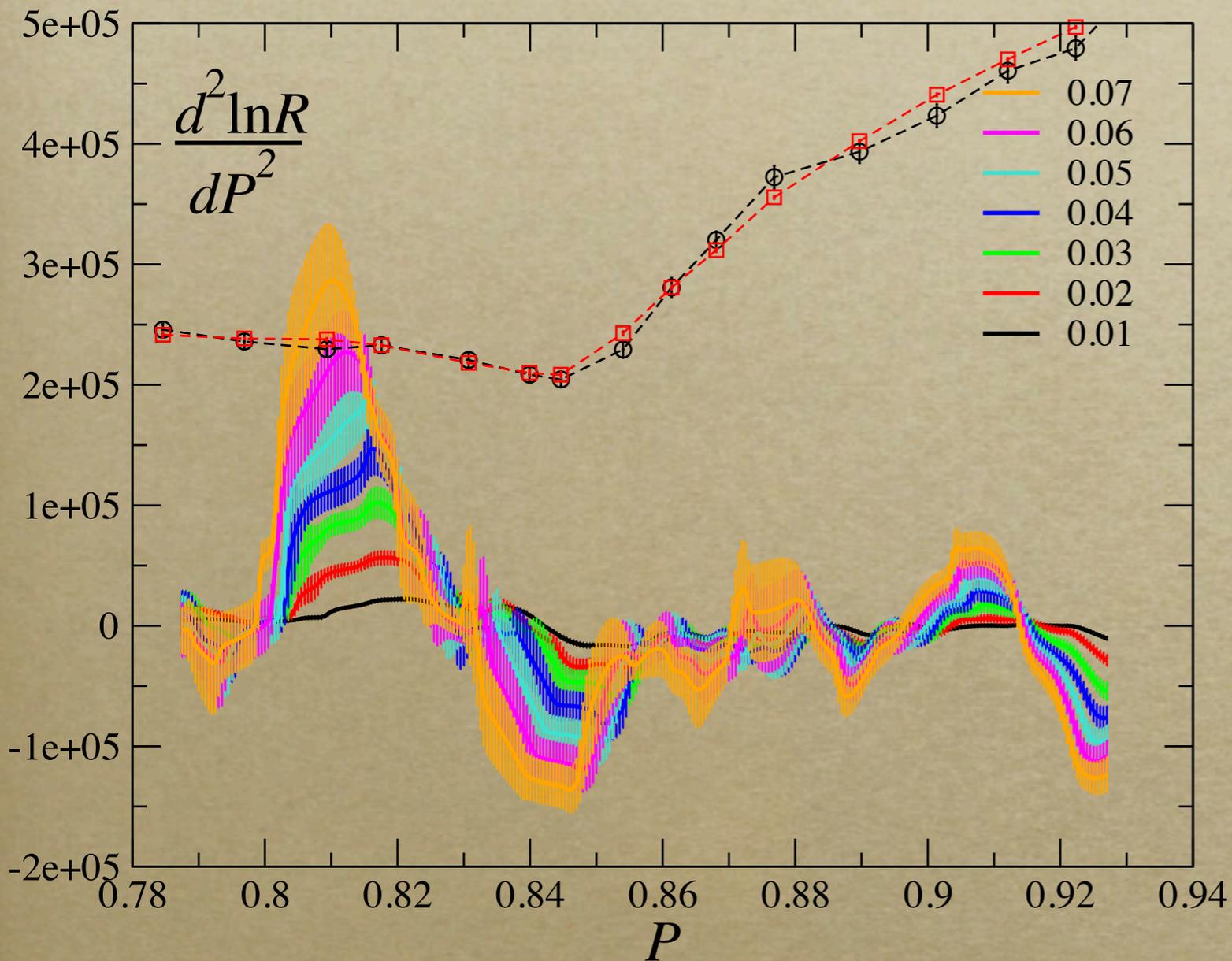
► $\ln R$ increases with h .

► Rapid increase @ $P \sim 0.81$

⇒ large curvature

Result:

$$\partial^2 V_{\text{eff}}/\partial P^2 = \partial^2 V_0/\partial P^2 - \partial^2(\ln R)/\partial P^2$$



1st term is calculate in two different ways (black and red). Maximum of the 2nd term exceeds 1st term at $P \sim 0.81$ for $h \gtrsim 0.06$.

There $\partial^2 V_{\text{eff}}/\partial P^2$ is negative. \Rightarrow **1st order phase transition**

$$h_c = 0.0614(69)$$

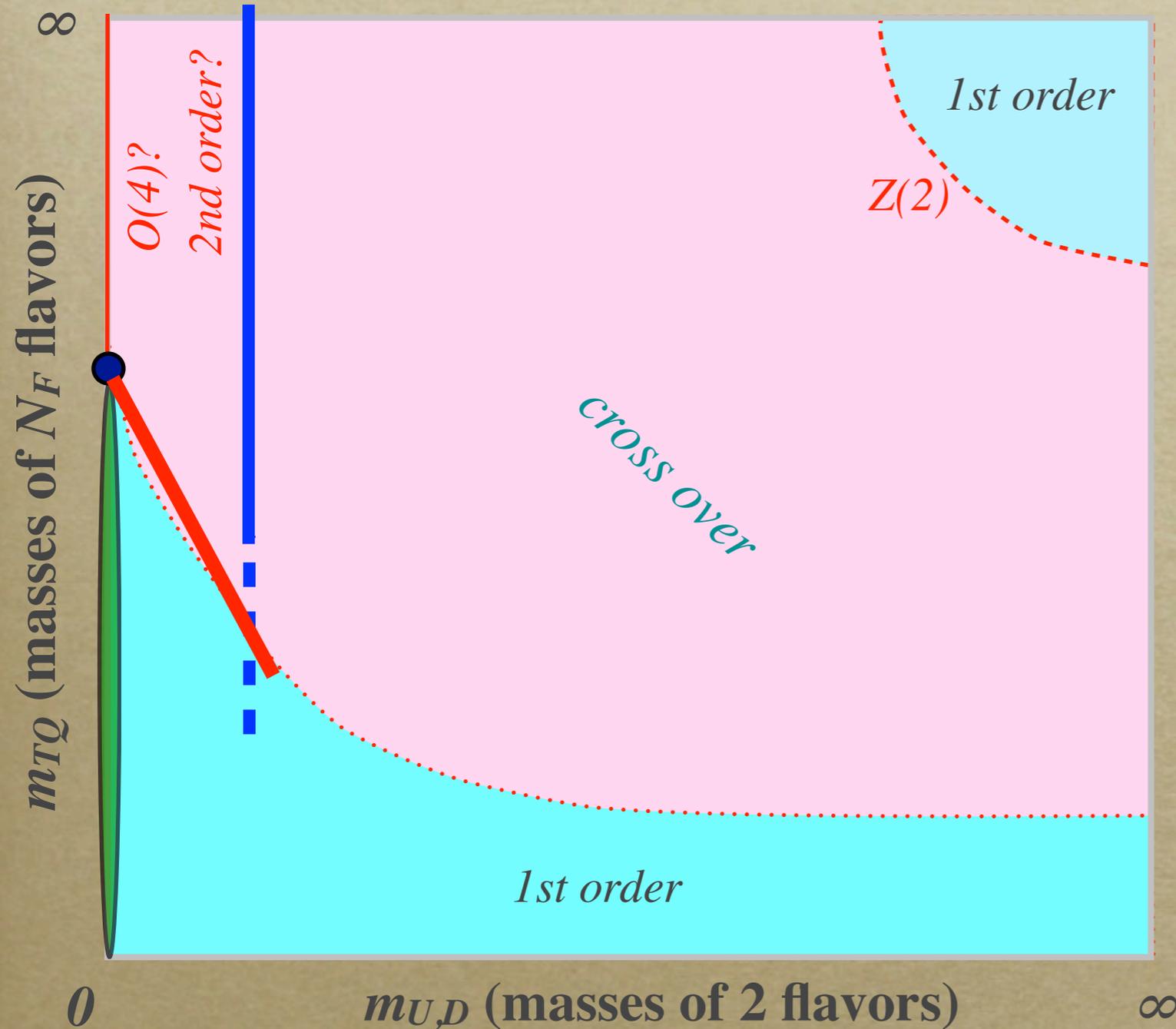
Comments

- ▶ $h_c = 0.0614(69) = 2 N_f (2K_{hc})^{N_f}$
- ▶ Critical kappa K_{hc} becomes small as N_f increases.
- ▶ For example, for $N_f=10$, $K_{hc} \sim 0.118$.
- ▶ Convergence of the HPE is studied in quenched theory in [Ejiri et al. in preparation], and found to be (LO)~(NLO) @ $K_h \sim 0.18$.
- ▶ Study of phase structure at finite density can be done in the same footing. [S. Ejiri, PRD77, 014508 (2008)]
- ▶ We found that a finite μ also makes K_{hc} small.

Summary and outlooks

- ✓ In general, QCD with many flavors is computationally demanding.
- ✓ We proposed an easy way to explore the phase structure of MFQ, and gave the critical mass.
- ✓ Future works:
 - Quantify the strength of 1st order transition
 - Check the universal power behavior at the tri-critical point

Structure around the tri-critical point



Meanfield approximation predicts

$$m_{lc} \sim (m_h - m_{hc})^{5/2}.$$

The power (5/2) is independent of N_f .