(Heavy) dense QCD and nuclear matter from a (3d) effective lattice theory



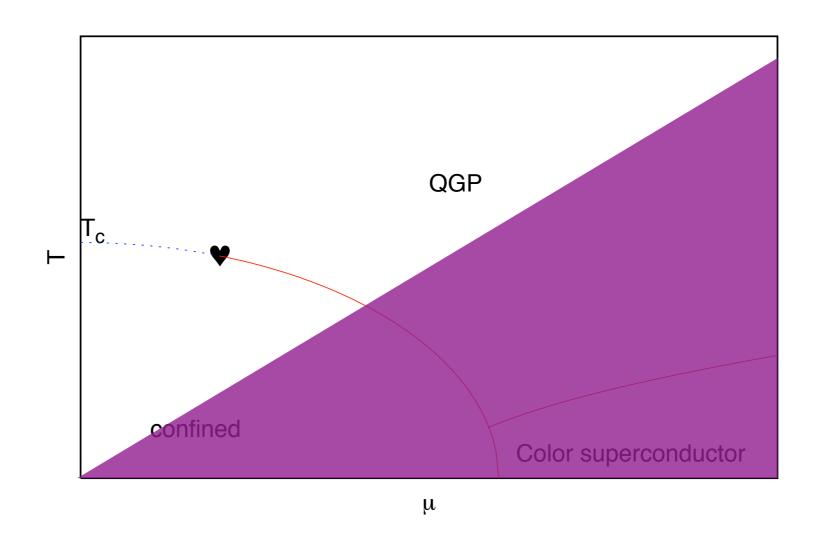
Owe Philipsen



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- 3d effective lattice theories derived by strong coupling methods
- The deconfinement transition in QCD with heavy dynamical quarks JHEP 1201 (2012)
- Cold and dense QCD: transition to nuclear matter JHEP 1409 (2014)
- Cold and dense chiral QCD on (very) coarse lattices PRL 113 (2014)

The (lattice) calculable region of the QCD phase diagram



- Sign problem prohibits direct simulation, circumvented by approximate methods: reweigthing, Taylor expansion, imaginary chem. pot., need $\mu/T \lesssim 1 \ (\mu = \mu_B/3)$
- No critical point in the controllable region
- Complex Langevin: lots of progress, but not in all parameter space, no "guarantees"

The effective lattice theory approach I

- Two-step treatment:
 - I. Calculate effective theory analytically
 - II. Simulate effective theory
- Step I.: split temporal and spatial link integrations:

$$Z = \int DU_0 DU_i \det Q \ e^{S_g[U]} \equiv \int DU_0 e^{-S_{eff}[U_0]} = \int DL \ e^{-S_{eff}[L]}$$

Spatial integration after analytic strong coupling and hopping expansion

- Result: 3d spin model of QCD, infinitely many couplings, ordered parametrically
- Truncation: valid for heavy quarks, sufficiently close to the continuum
- Step II: sign problem milder: Monte Carlo, complex Langevin
- Numerical simulations in 3d without fermion matrix inversion, very cheap!

Starting point: Wilson's lattice Yang-Mills action

Partition function; link variables as degrees of freedom

$$Z = \int \prod_{x,\mu} dU(x;\mu) \exp(-S_{YM}) \equiv \int DU \exp(-S_{YM})$$

Wilson's gauge action

$$S_W = -\frac{\beta}{N} \sum_{p} \operatorname{ReTr}(U_p) = \sum_{p} S_p$$
 $\beta = \frac{2N}{g^2}$

Plaquette:
$$\Box \to 1 + i a^2 g F_{\mu\nu} - \frac{a^4 g^2}{2} F_{\mu\nu} F^{\mu\nu} + O(a^6) + \dots$$
 $U_{\mu}(x) = \mathrm{e}^{-\mathrm{i} a g A_{\mu}(x)}$

$$T=rac{1}{aN_t}$$
 continuum limit $a o 0, N_t o \infty$

Small $\beta(a) \Rightarrow \text{small T}$

■ Leading order graph in case of $N_{\tau} = 4$:



Figure: 4 plaquettes in fundamental representation lead to a 2 Polyakov loop interaction term

Integration of spatial link variables leads to

$$-S_1 = u^{N_\tau} \sum_{\langle ij \rangle} \operatorname{tr} W_i \operatorname{tr} W_j$$

Expansion parameter: $u = a_f(\beta) = \beta/18 + \cdots$

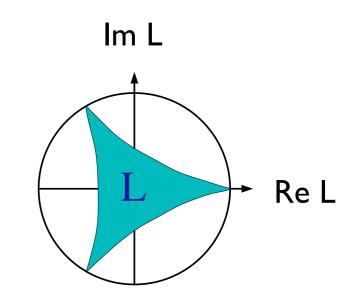
- Possible generalizations: larger distance, higher dimensional representations, larger number of loops involved, . . .
- Here: Decorate LO graph with additional spatial and temporal plaquettes

Effective one-coupling theory for SU(3) YM

(L=TrW)
$$Z = \int [dL] \exp \left[-S_1 + V_{SU(3)}\right]$$

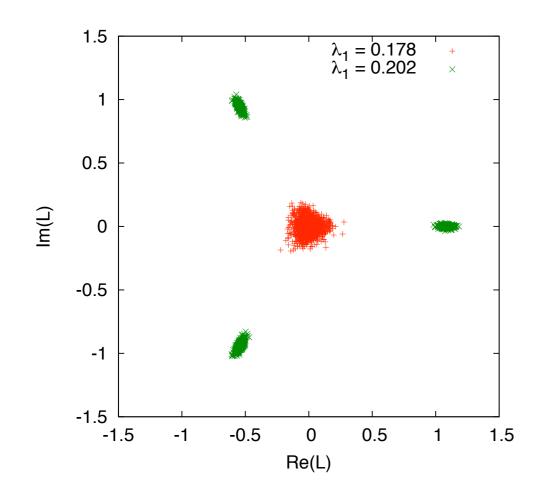
$$= \int [dL] \prod_{\langle ij \rangle} \left[1 + 2\lambda_1 \operatorname{Re}\left(L_i L_j^*\right)\right] *$$

$$* \prod_i \sqrt{27 - 18|L_i|^2 + 8\operatorname{Re}L_i^3 - |L_i|^4}$$

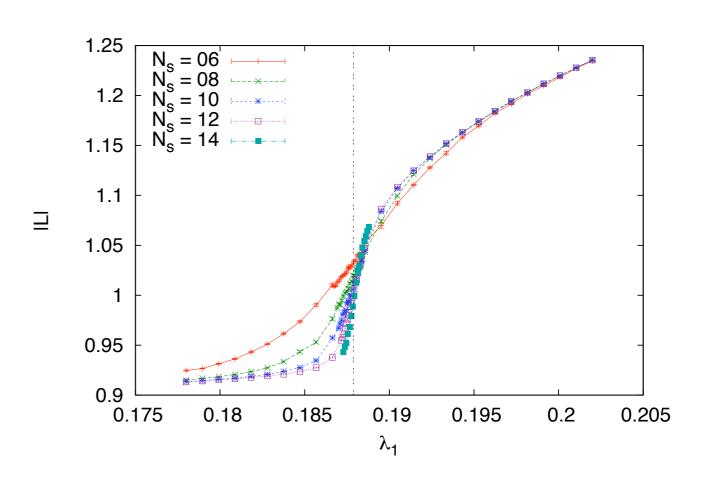


$$\lambda(u, N_{\tau} \ge 5) = u^{N_{\tau}} \exp \left[N_{\tau} \left(4u^4 + 12u^5 - 14u^6 - 36u^7 + \frac{295}{2}u^8 + \frac{1851}{10}u^9 + \frac{1055797}{5120}u^{10} \right) \right]$$

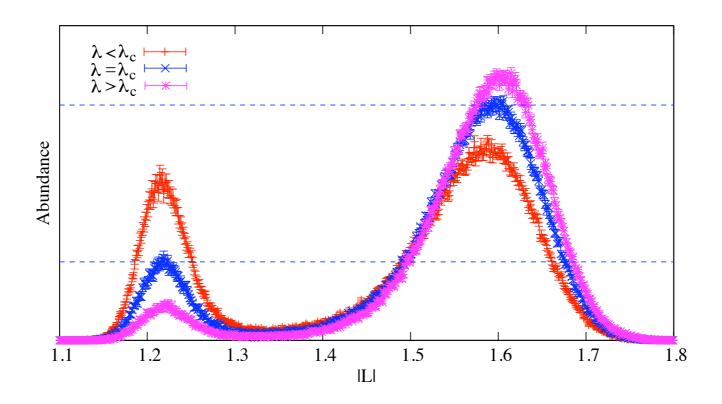
Numerical results for SU(3), one coupling



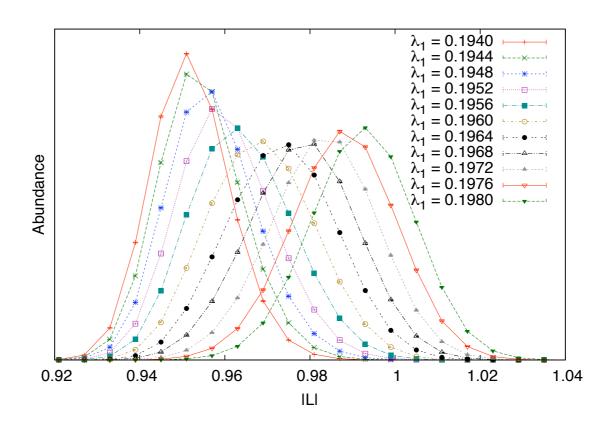
Order-disorder transition = Z(3) breaking



First order phase transition for SU(3):



Second order (3d Ising) phase transition for SU(2):

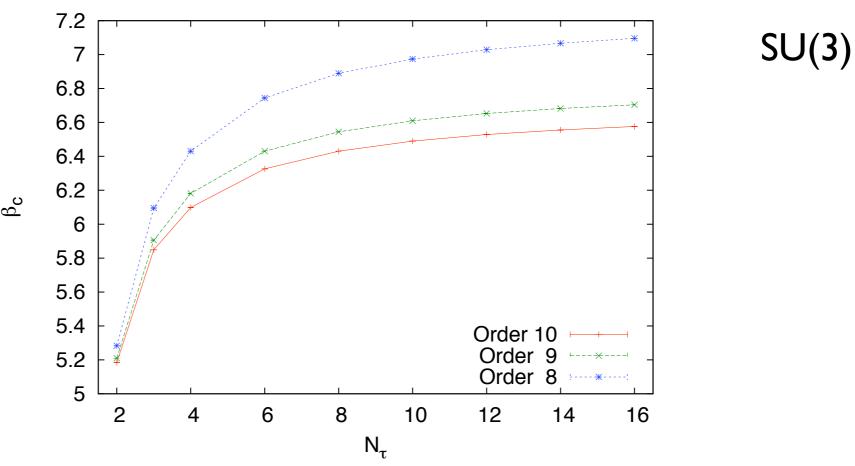


Mapping back to 4d finite T Yang-Mills

Inverting

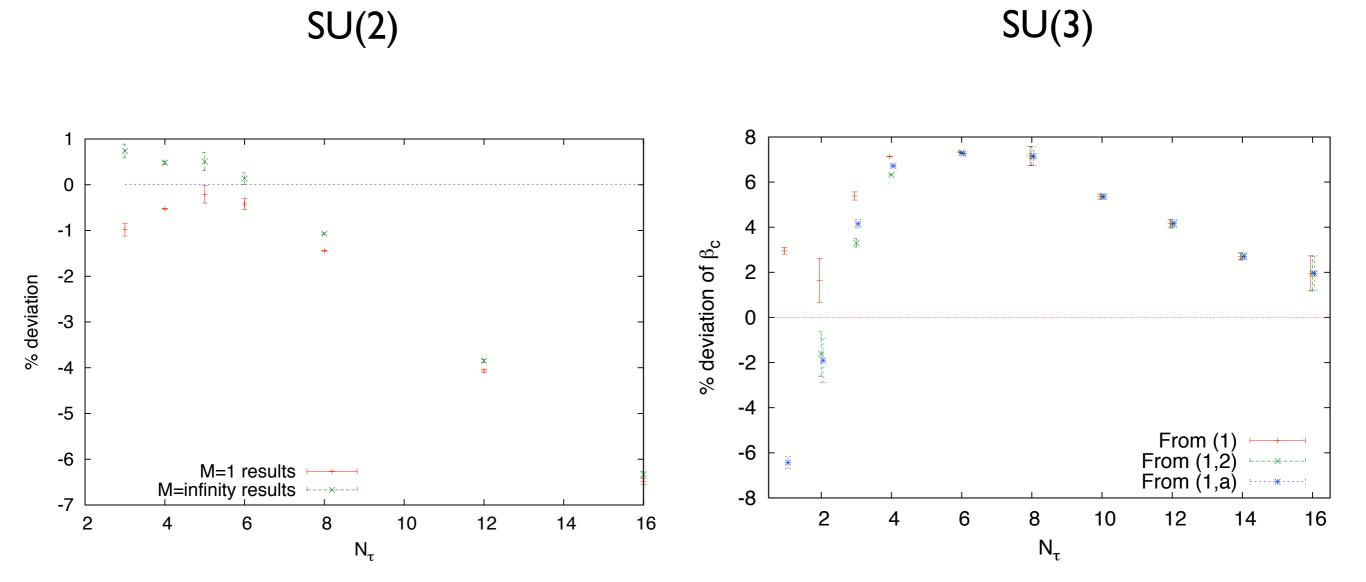
$$\lambda_1(N_{\tau},\beta) \to \beta_c(\lambda_{1,c},N_{\tau})$$

...points at reasonable convergence



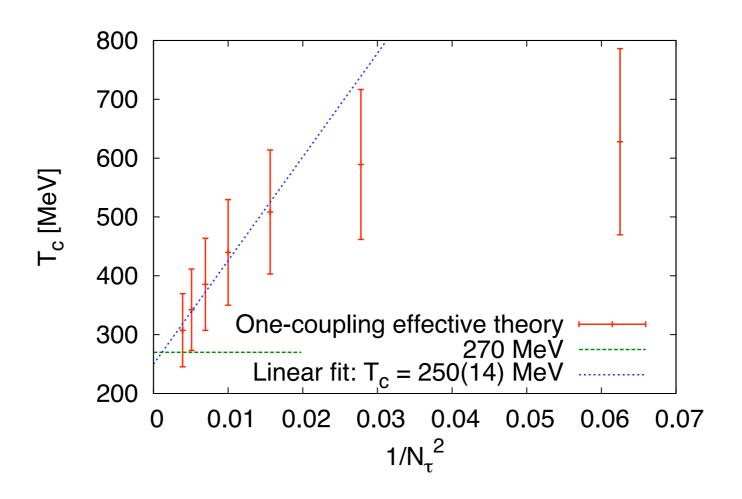
Comparison with 4d Monte Carlo

Relative accuracy for β_c compared to the full theory



Note: influence of additional couplings checked explicitly!

Continuum limit feasible!



- -error bars: difference between last two orders in strong coupling exp.
- -using non-perturbative beta-function (4d T=0 lattice)
- -all data points from one single 3d MC simulation!

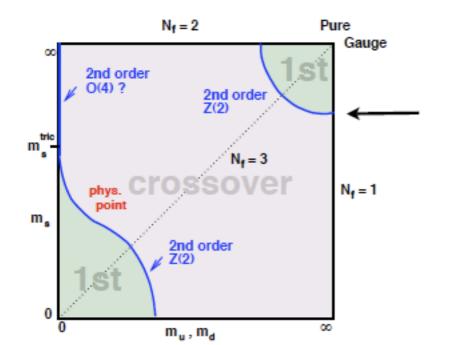
Including heavy, dynamical Wilson fermions

Expand in the *hopping parameter* $\kappa = 1/(2aM + 8)$:

$$-S_{\text{eff}} = \sum_{i} \lambda_{i}(u, \kappa, N_{\tau}) S_{i}^{S} - 2N_{f} \sum_{i} \left[h_{i}(u, \kappa, \mu, N_{\tau}) S_{i}^{A} + \overline{h}_{i}(u, \kappa, \mu, N_{\tau}) S_{i}^{\dagger A} \right]$$

Now, keep only $\lambda_1 S_1^S$ and $h_1 S_1^A + \overline{h}_1 S_1^{\dagger A}$

NLO: $\sim \kappa^2$

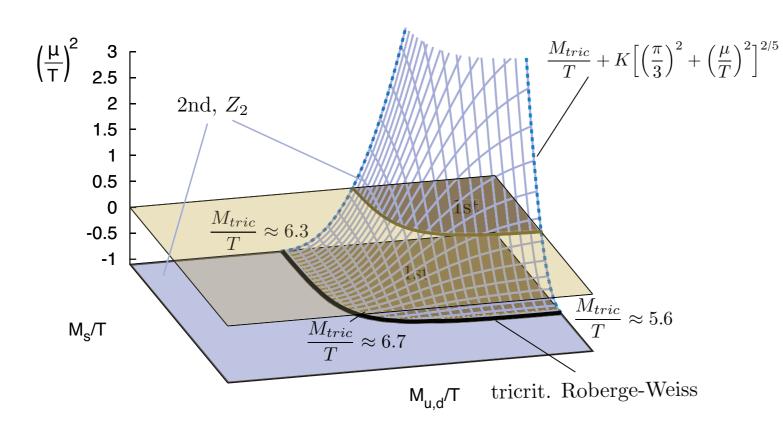


eff. theory 4d MC, WHOT 4d MC, de Forcrand et al

N_f	M_c/T	$\kappa_c(N_{\tau}=4)$	$\kappa_c(4)$, Ref. [23]	$\kappa_c(4)$, Ref. [22]
1	7.22(5)	0.0822(11)	0.0783(4)	~ 0.08
2	7.91(5)	0.0691(9)	0.0658(3)	_
3	8.32(5)	0.0625(9)	0.0595(3)	_

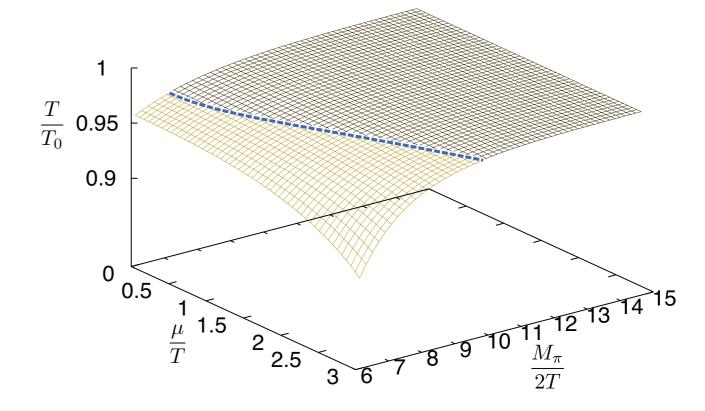
Accuracy ~5%, predictions for Nt=6,8,... available!

The fully calculated deconfinement transition

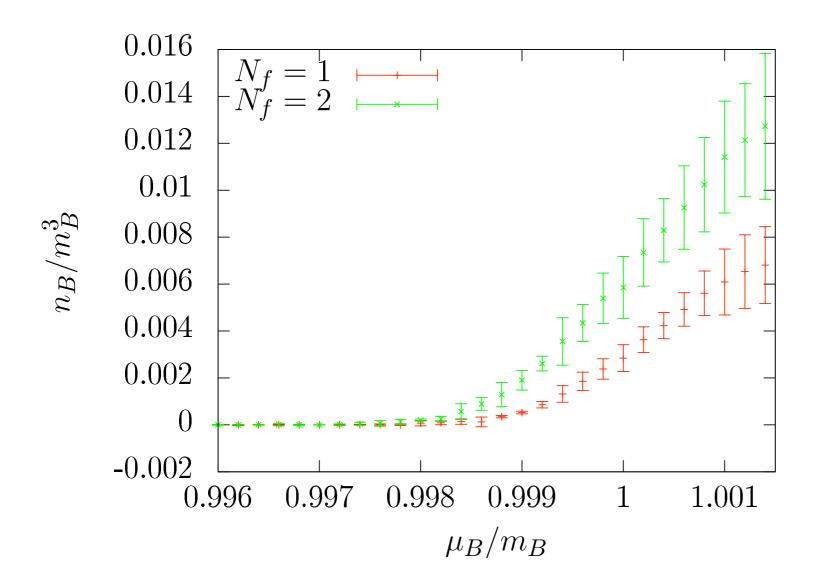


deconfinement critical surface

phase diagram for Nf=2, Nt=6



The equation of state for nuclear matter



$$S_{eff} \sim \kappa^n u^m, \quad n+m=4$$

$$m_{\pi} = 20 \text{ GeV}, T = 10 \text{ MeV}$$

$$\frac{\mu}{T} \sim 4000$$

Effect of binding between baryons:

Binding energy per nucleon:

Transition is smooth crossover:

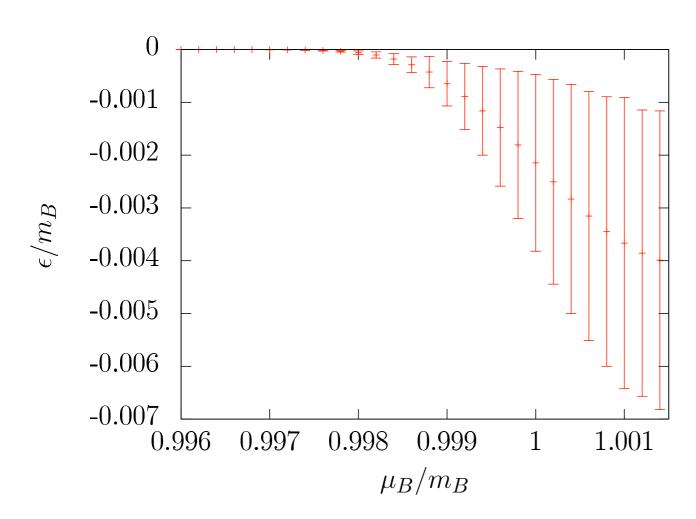
$$\mu_c < m_B$$

$$\epsilon = \frac{\mu_c - m_B}{m_B} \sim 10^{-3}$$

$$T > T_c \sim \epsilon \, m_B$$

Binding energy per nucleon

$$\epsilon = \frac{e - n_B m_B}{n_B m_B} = \frac{e}{n_B m_B} - 1$$



Binding energy per nucleon

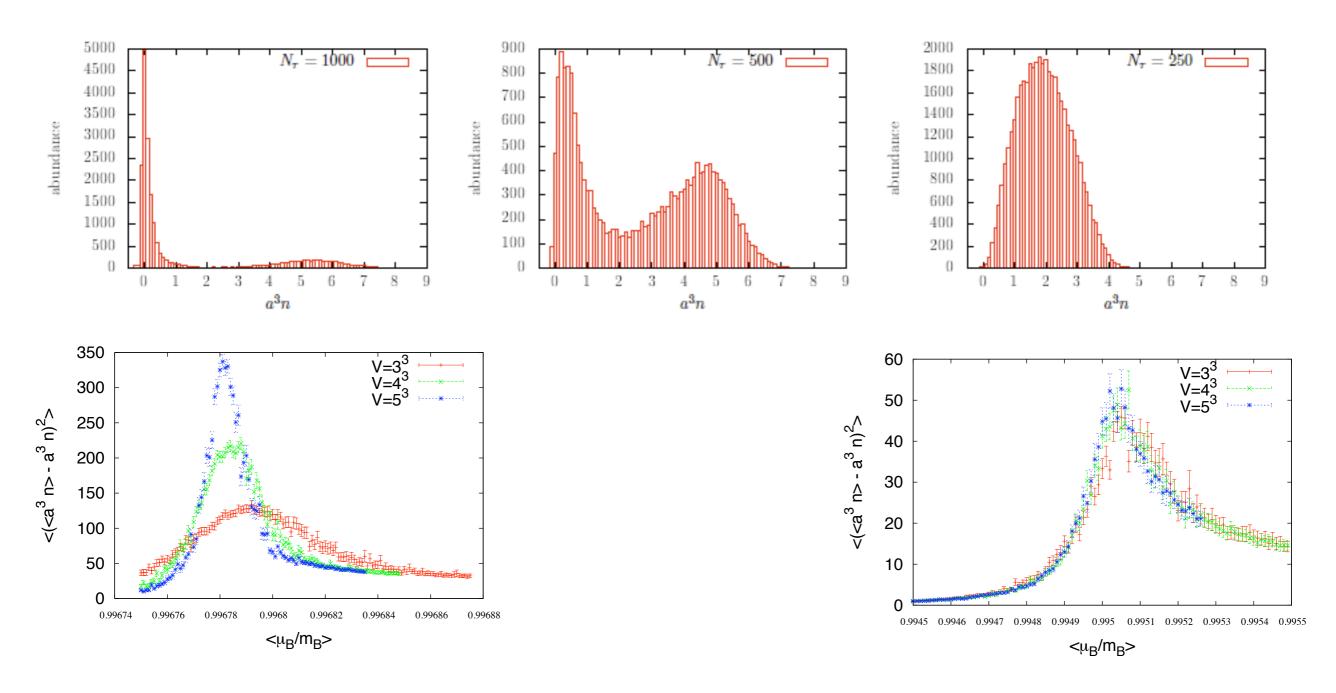
$$\epsilon = \frac{e - n_B m_B}{n_B m_B} = \frac{e}{n_B m_B} - 1$$

$$\begin{array}{c} 0 \\ -0.001 \\ -0.002 \\ -0.003 \\ \hline -0.004 \\ -0.005 \\ -0.006 \\ -0.007 \\ 0.996 \ \ 0.997 \ \ 0.998 \ \ 0.999 \ \ 1 \ \ 1.001 \\ \hline \mu_B/m_B \end{array}$$
 ... to be continued...

Minimum: access to nucl. binding energy, nucl. saturation density!

 $\epsilon \sim 10^{-3}$ consistent with the location of the onset transition

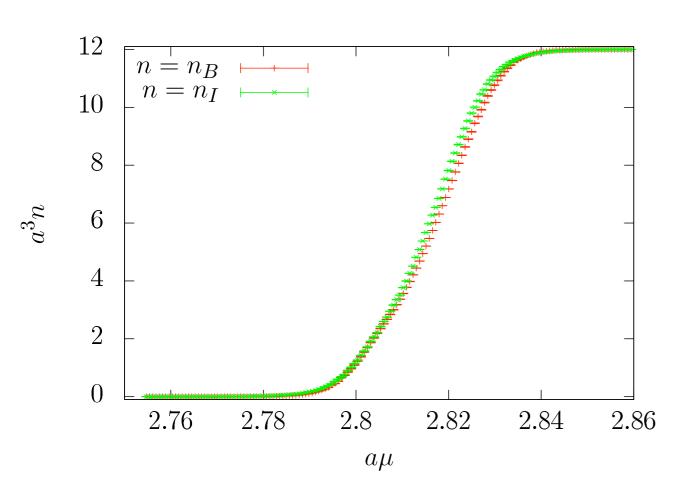
Lighter quarks: first order + endpoint!

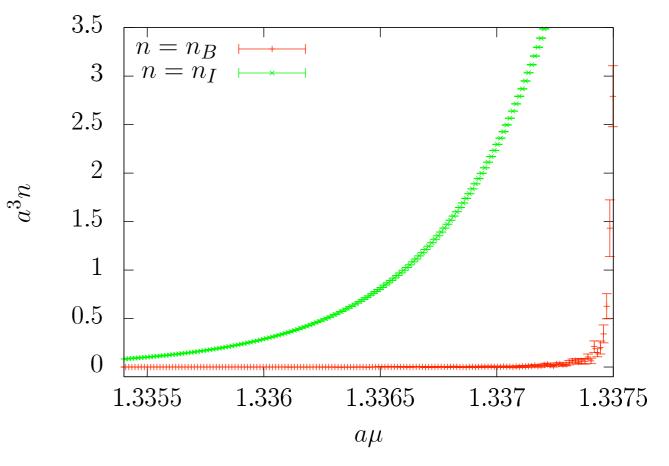


- $O(k^4)$:Stretching the hopping series, $\kappa = 0.12, \beta = 5.6$
- Coexistence of vacuum and finite density phase: 1st order
- If the temperature $T = \frac{1}{aN_{\tau}}$ or the quark mass is raised this changes to a crossover nuclear liquid gas transition!!!

attn: no convergence yet!

Finite isospin vs baryon chemical potential





nearly static quarks

$$\frac{m_{\pi}}{2} \approx \frac{m_B}{3}$$

lighter quarks

$$\frac{m_{\pi}}{2} < \frac{m_B}{3}$$

onset at smaller chemical potential

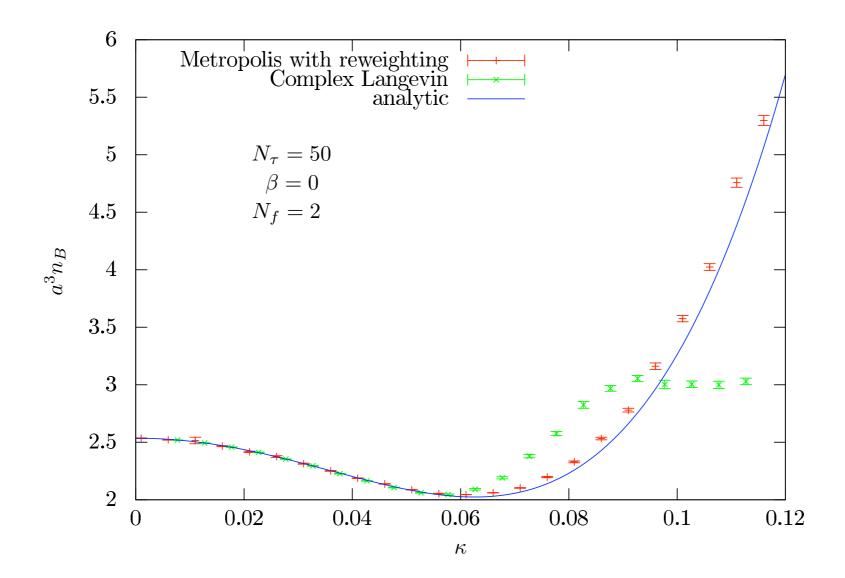
The problem with complex Langevin

No non-analytic phase transition seen yet

Aarts, Seiler, Sexty, Stamatescu

Does not work in all parameter regions!

In effective theory:



The effective lattice theory approach II

Two-step treatment:

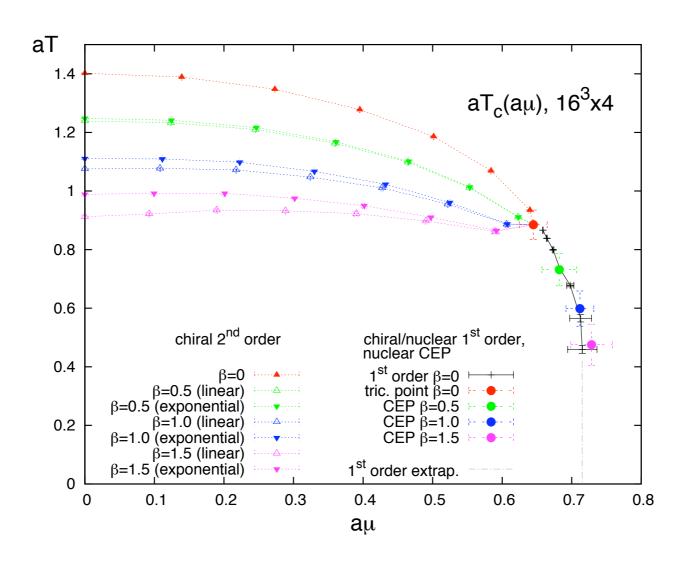
de Forcrand, Langelage, O.P., Unger Phys.Rev.Lett. 113 (2014) 152002

- I. Calculate effective theory analytically
- II. Simulate effective theory
- Step I.: integrate over gauge links in strong coupling expansion, leave fermions

$$\begin{split} Z_{\rm QCD} &= \int d\psi d\bar{\psi} dU e^{S_F + S_G} = \int d\psi d\bar{\psi} Z_F \left\langle e^{S_G} \right\rangle_{Z_F} \\ \left\langle e^{S_G} \right\rangle_{Z_F} &\simeq 1 + \left\langle S_G \right\rangle_{Z_F} = 1 + \frac{\beta}{2N_c} \sum_P \left\langle \text{tr}[U_P + U_P^\dagger] \right\rangle_{Z_F} \\ \end{split} \qquad Z_F(\psi, \bar{\psi}) = \int dU e^{S_F} \left\langle e^{S_G} \right\rangle_{Z_F} \\ &= 1 + \left\langle S_G \right\rangle_{Z_F} = 1 + \frac{\beta}{2N_c} \sum_P \left\langle e^{S_G} \right\rangle_{Z_F} \\ \end{split}$$

- Result: 4d "polymer" model of QCD (hadronic degrees of freedom!)
 Valid for all quark masses (also m=0!), at strong coupling (very coarse lattices)
- Step II: sign problem milder: Monte Carlo with worm algorithm
- Numerical simulations without fermion matrix inversion, very cheap!

The QCD Phase diagram at strong coupling



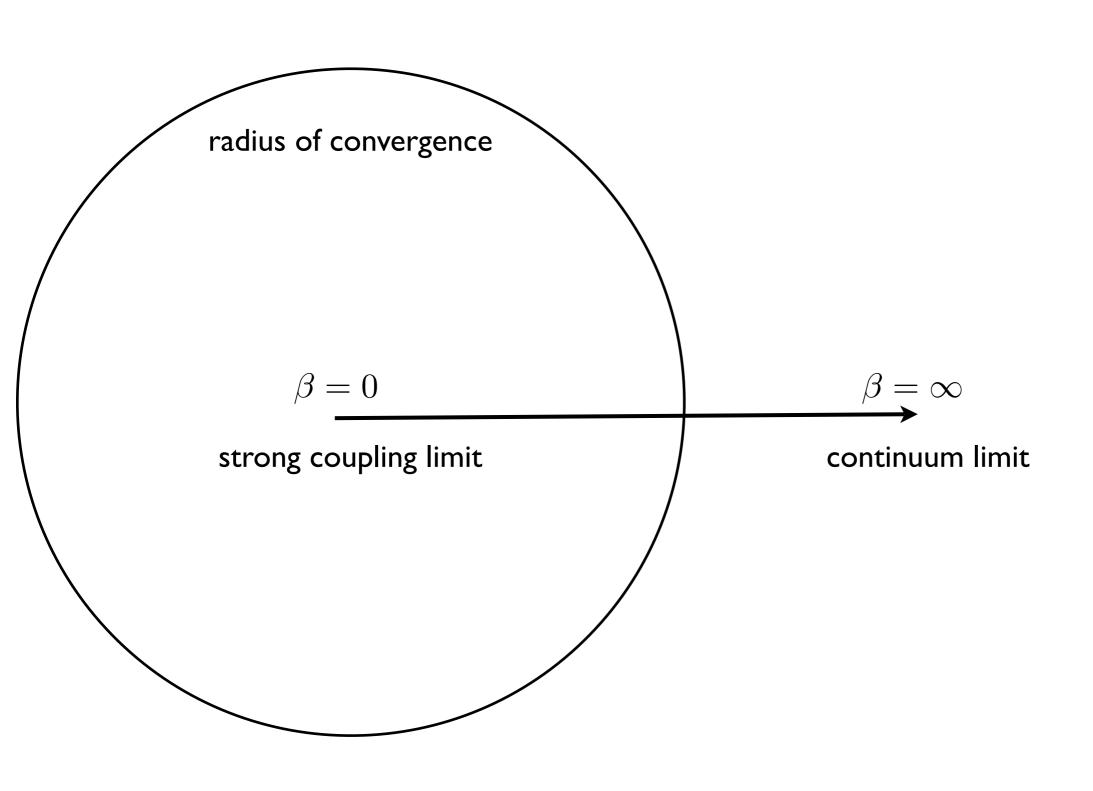
- LO gauge correction included, simulation by worm algorithm
- Chiral phase transition with 2nd order and 1st order line meeting in tricritical point
- Nuclear liquid gas transition on top of first order chiral one at strong coupling

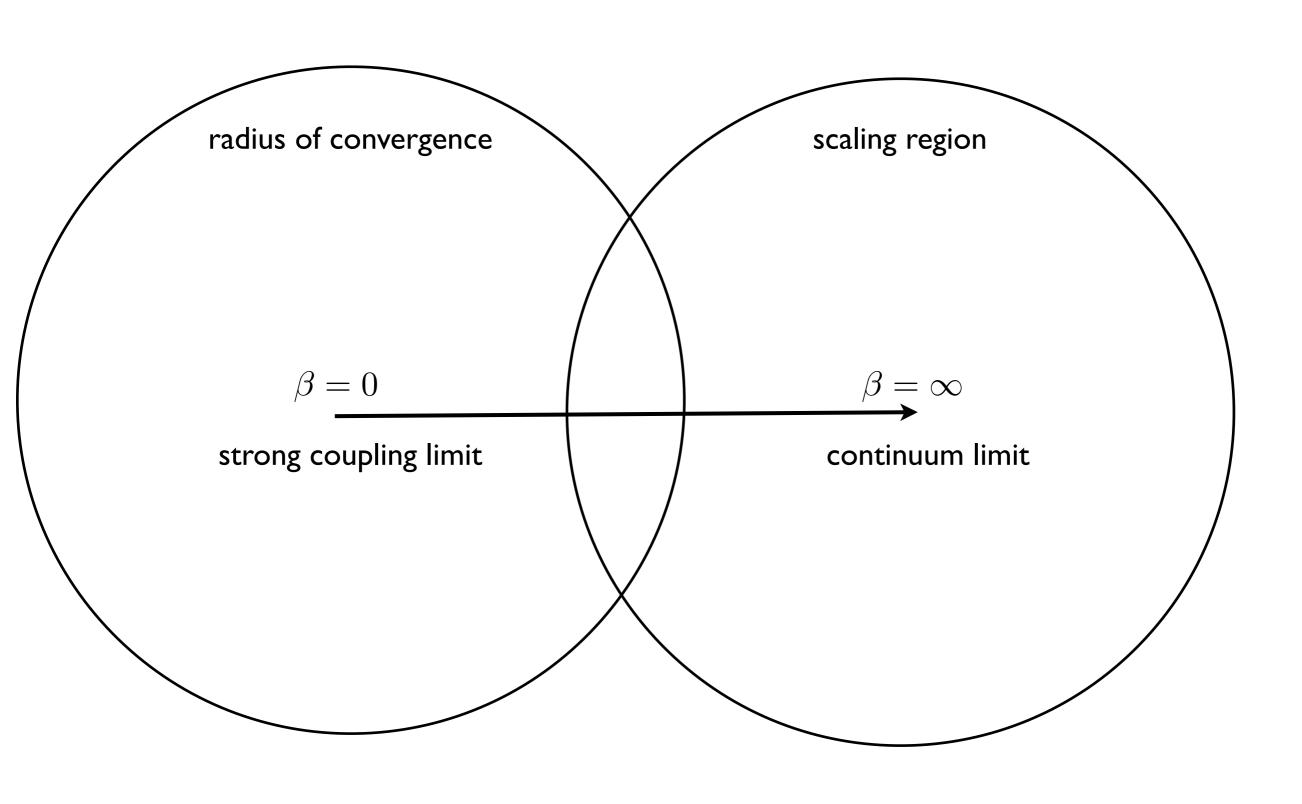
Conclusions

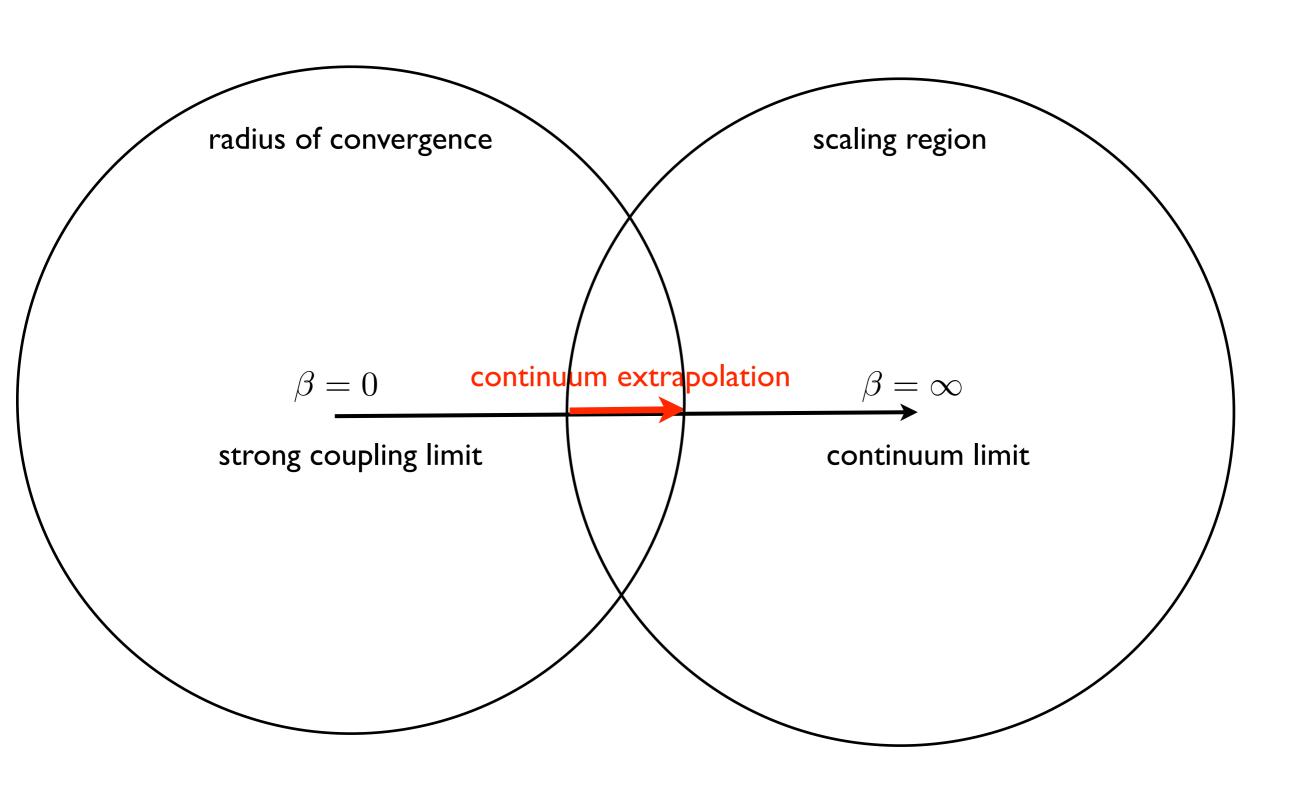
- Two-step treatment of QCD phase transitions:
 - I. Derivation of effective action by strong coupling expansion
 - II. Simulation of effective theory
- Finite T transition + nuclear liquid gas transition for:

Heavy QCD on not so coarse or chiral QCD on very coarse lattices

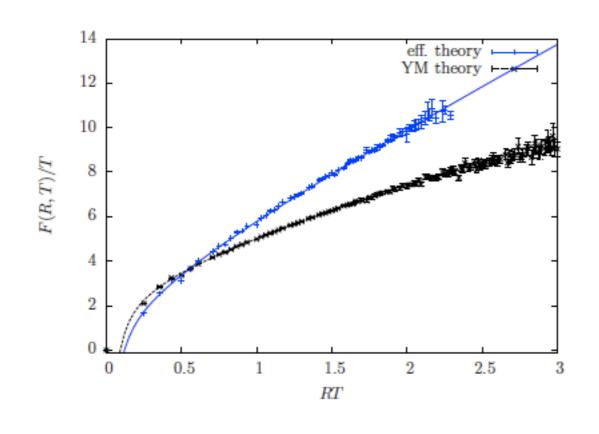
$$\beta = 0 \qquad \qquad \beta = \infty$$
 strong coupling limit continuum limit





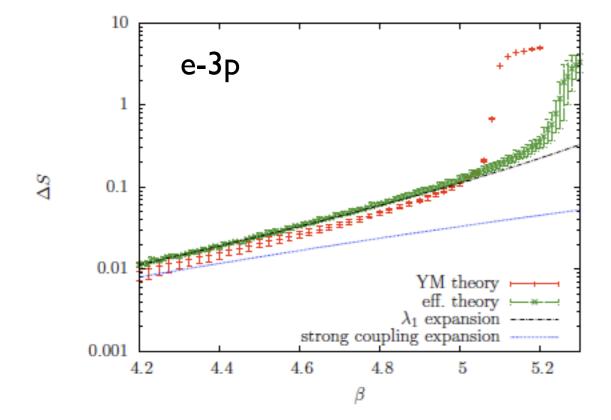


What does and does not work?



Correlation functions and spectrum: NO

couplings over large distances needed



Thermodynamics and critical coupling: YES

partition function needed, ultra-local!

Subleading couplings

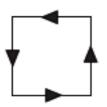
Subleading contributions for next-to-nearest neighbours:

$$\lambda_2 S_2 \propto u^{2N_\tau+2} \sum_{[kl]}' 2 \operatorname{Re}(L_k L_l^*) \text{ distance } = \sqrt{2}$$
 $\lambda_3 S_3 \propto u^{2N_\tau+6} \sum_{\{mn\}}'' 2 \operatorname{Re}(L_m L_n^*) \text{ distance } = 2$

as well as terms from loops in the adjoint representation:

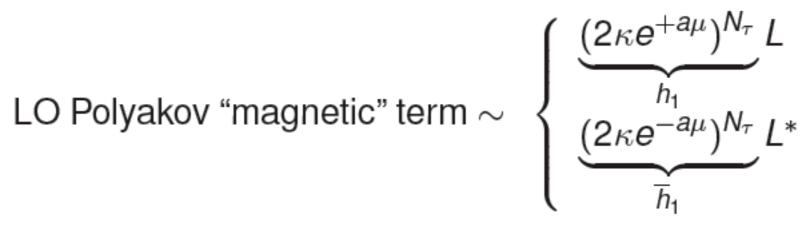
$$\lambda_a S_a \propto u^{2N_\tau} \sum_{\langle ij \rangle} \text{Tr}^{(a)} W_i \text{Tr}^{(a)} W_j$$
 ; $\text{Tr}^{(a)} W = |L|^2 - 1$

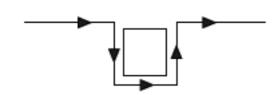
Links along imaginary time gain $\exp(\pm \mu a)$



reabsorbed in gauge part: $\begin{cases} \beta \to \beta + \mathcal{O}(\kappa^4) \\ u(\beta) \to u(\beta, \kappa) \end{cases}$

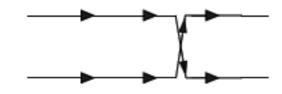






higher corrections to the above:

$$h_1 = (2\kappa e^{a\mu})^{N_\tau} \left[1 + \mathcal{O}(k^2) f(u) + \dots \right]$$



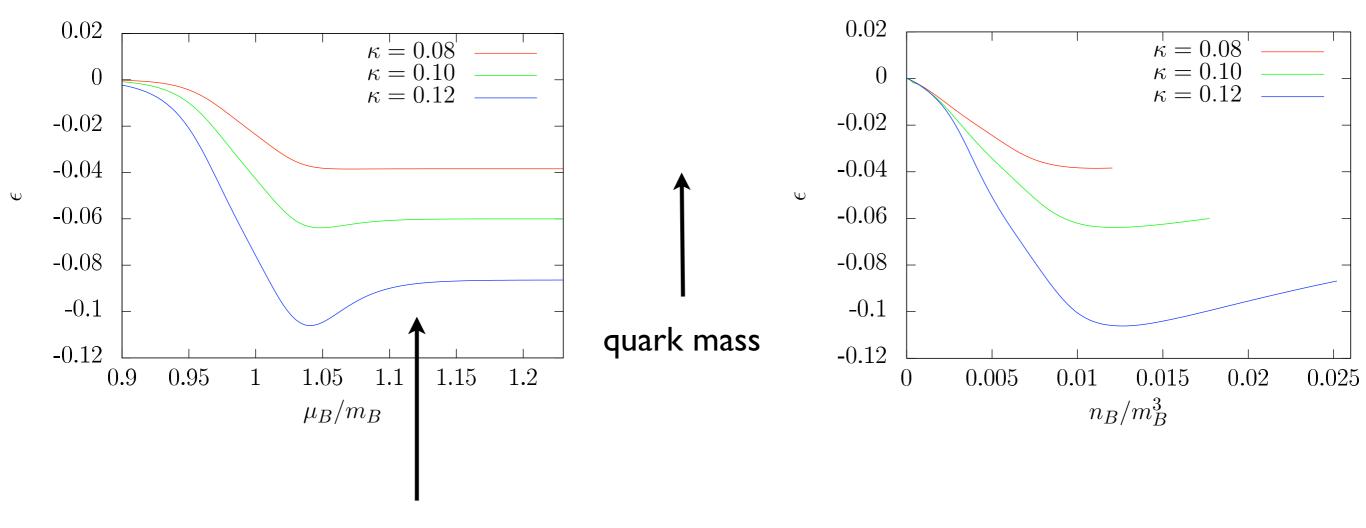
other (suppressed) terms, such as $h_2(L_x L_{x+\hat{i}})$,

$$h_2 \sim (2\kappa e^{a\mu})^{2N_\tau} \kappa^2 \dots$$

Quark mass dependence of the binding energy:

Expect short range nucl. potential for heavy pions, $V \sim \frac{e^{-m_{\pi}r}}{r}$

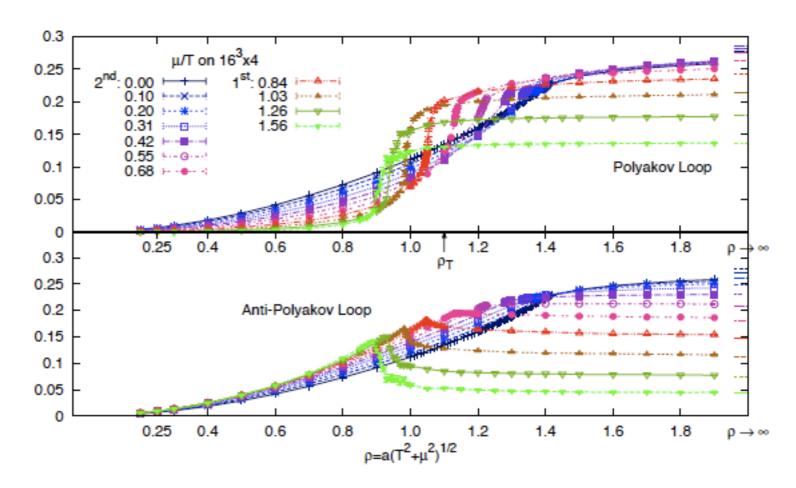
Analytic solution, finite lattice spacing:



lattice saturation

Two transitions seen:

Chiral/deconfinement



Nuclear onset

