

# QCD phase diagram with heavy quarks: Universal aspects in one-loop models and extension to 2-loop in Curci-Ferrari

Motivation

generic 1-loop

Curci-Ferrari at  
2-loopVanishing  $\mu=0$ 

Imaginary

 $\mu = i\mu_i$ 

Conclusion

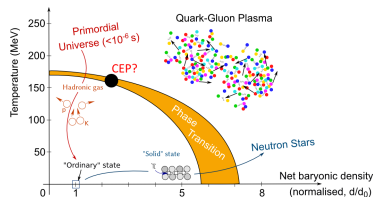
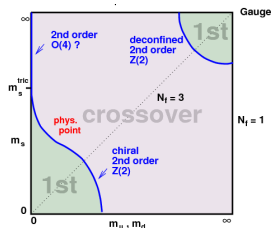
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# Motivation



Several approaches on the market:

- ▶ Lattice QCD [de Forcrand, Philipsen, Rodriguez-Quintero, Mendes, ...]
- ▶ Dyson Schwinger Equations [Alkofer, Fischer, Huber, ...]
- ▶ Functional Renormalization Group [Pawlowski, Mitter, Schaefer...]
- ▶ Variational Approach [Reinhardt, Quandt, ...]
- ▶ Gribov-Zwanziger Action [Dudal, Oliveira, Zwanziger...]
- ▶ Matrix-, QM-, NJL-Model,... [Pisarski, Dumitru, Schaffner-B., Stiele, ...]
- ▶ Curci-Ferrari Model [Reinosa, Serreau, Tissier, Wschebor, ...]

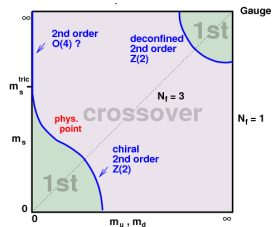
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- ▶ Lattice QCD
- ▶ Dyson Schwinger Equations
- ▶ Functional Renormalization Group
- ▶ Variational Approach
- ▶ Gribov-Zwanziger Action
- ▶ Matrix-, QM-, NJL-Model,...
- ▶ Curci-Ferrari Model

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## Part 1:

- ▶ generic aspects of the heavy quark region
- ▶ common to all approaches at one-loop order

## Part 2:

- ▶ higher order corrections in one particular model
- ▶ Curci-Ferrari at two-loop order

# Polyakov loops & effective potentials

At the YM point, a relevant order parameter for the deconfinement transition is the (anti-)Polyakov loop. It is related to the free energy  $F_q$  necessary to bring a quark into a "bath" of gluons.

$$\ell \equiv \frac{1}{3} \text{tr} \left\langle P \exp \left( ig \int_0^\beta d\tau A_0^a t^a \right) \right\rangle \sim e^{-\beta F_q} \quad \bar{\ell} \sim e^{-\beta F_{\bar{q}}}$$

Hence

$$\ell = 0 \leftrightarrow F_q = \infty \leftrightarrow \text{confinement} \quad \ell \neq 0 \leftrightarrow F_q < \infty \leftrightarrow \text{deconfinement}$$

In all models, for each value of the temperature  $T$ , one then minimizes an effective potential

$$V_{\text{glue}}(\ell, \bar{\ell})$$

to find the physical position of the system. The particular form of this potential is model-dependent.

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# Polyakov loops & effective potentials

Introducing quarks, center symmetry is explicitly broken. For heavy quarks, this breaking is "soft", thus:

$$\ell \approx 0 \leftrightarrow F_q \approx \infty \leftrightarrow \text{confinement} \quad \ell \neq 0 \leftrightarrow F_q < \infty \leftrightarrow \text{deconfinement}$$

Therefore  $\ell, \bar{\ell}$  are still approximately good order parameters.

At leading order, the new effective potential is simply found by adding a quark part at one-loop level:

$$V_{\text{glue}}(\ell, \bar{\ell}) + V_{\text{quark}}(\ell, \bar{\ell}, \mu)$$

→ Let's look at some particular cases in more detail.

# Explicit Potentials in various Models

**Gribov-Zwanziger:** [JM, U.Reinosa, J.Serreau (2018)]

$$V_{GZ} = \underbrace{-\frac{d}{2} \frac{\sum_{\kappa} m_{\kappa}^4}{g^2 C_{\text{ad}}} + \frac{d-1}{2} \sum_{\kappa} \int_Q^T \ln \frac{Q_{\kappa}^4 + m_{\kappa}^4}{Q_{\kappa}^2} - \frac{1}{2} \sum_{\kappa} \int_Q^T \ln Q_{\kappa}^2}_{V_{\text{glue}}} - \text{Tr Ln}(\not{\partial} + M)$$

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**Curci-Ferrari:** [U. Reinosa, J. Serreau, M. Tissier (2015)]

$$V_{CF} = \underbrace{\sum_{\kappa} \frac{T}{2\pi^2} \int_0^{\infty} dq q^2 \left\{ 3 \ln [1 - e^{-\beta \varepsilon q + i r \kappa}] - \ln [1 - e^{-\beta q + i r \kappa}] \right\}}_{V_{\text{glue}}} - \text{Tr Ln}(\not{\partial} + M - i g \gamma_0 \vec{A}^k t^k)$$

Curci-Ferrari at  
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**Matrix-Models:** [K.Kashiwa, R.D.Pisarski and V.V.Skokov (2012)]

$$V_M = -\frac{4\pi^2}{3} T^2 T_d^2 \left( c_1 \sum_{i,j=1}^N V_1(q_i - q_j) + c_2 \sum_{i,j=1}^N V_2(q_i - q_j) + \frac{(N^2 - 1)}{60} c_3 \right) - \underbrace{\frac{(N^2 - 1)\pi^2}{45} T^4 + \frac{2\pi^2}{3} T^4 \sum_{i,j=1}^N V_2(q_i - q_j)}_{V_{\text{glue}}} + \ln \det(\gamma^{\mu} \partial_{\mu} + q \delta^{\mu 4} + im)$$

**Lattice:** [M.Fromm, J.Langelage, S.Lottini and O.Philippen (2012)]

$$Z_{\text{eff}} = \int [dU_0] \underbrace{\left( \prod_{\langle ij \rangle} [1 + 2\lambda_1 \text{Re} L_i^* L_j] \right)}_{Z_{\text{glue}}} \left( \prod_{\vec{x}} \det \left[ (1 + h_1 W_{\vec{x}}) (1 + \bar{h}_1 W_{\vec{x}}^{\dagger}) \right]^{2N_f} \right)$$

# Commonalities & Assumptions

- ▶ Potential  $v_{\text{glue}}$  is confining, with a minimum at  $\ell = 0$  at zero temperature
- ▶ Quarks are added at one-loop level, in form of a  $\text{Tr Ln}$

$$V = V_{\text{glue}} - \text{Tr Ln}(\not{\partial} + M)$$

Then in the heavy quark limit, the  $\text{Tr Ln}$  expands and one finds

$$\beta^4 V(\ell, \beta, M) = v_{\text{glue}}(\ell, \beta) - 2N_f f(\beta M) \ell$$

$$f(x) = (3x^2/\pi^2)K_2(x)$$

$K_2(x)$  is the modified Bessel function of the second kind

$\ell$ : Polyakov loop

$\beta$ : inverse temp.

$M$ : deg. quark mass

→ How do we find the 2nd order critical line?

Motivation

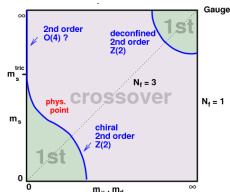
generic 1-loop

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# Determination of the critical line



$$\beta^4 V(\ell, \beta, M) = v_{\text{glue}}(\ell, \beta) - 2N_f f(\beta M) \ell$$

For a **fixed**  $N_f$ :

3 parameters:  $\ell, \beta, \beta M$

3 equations:  $\partial_\ell V = \partial_\ell^2 V = \partial_\ell^3 V = 0$

This yields:

$$\underbrace{\partial_\ell v_{\text{glue}} = 2N_f f(\beta M)}_{\text{determines model-dep. } \beta M},$$

$$\underbrace{\partial_\ell^2 v_{\text{glue}} = \partial_\ell^3 v_{\text{glue}} = 0}_{\text{determines } \ell, \beta, \text{ indep. of } \beta M, N_f}$$

→  $N_f f(\beta M) = N'_f f(\beta M') = f(\beta M_s) + 2f(\beta M_{ud})$  is const. on the critical line!



# Determination at non-vanishing chemical potential

$$\beta^4 V = v_{\text{glue}}(\ell, \bar{\ell}, \beta) - N_f f(\beta M)(e^{-\beta\mu}\ell + e^{\beta\mu}\bar{\ell})$$

For a fixed  $N_f, \mu$ :

4 parameters:  $\ell, \bar{\ell}, \beta, \beta M$

4 equations:

$$\partial_\ell V = \partial_{\bar{\ell}} V = 0, \quad (1)$$

$$\underbrace{\partial_{\bar{\ell}}^2 V \partial_{\bar{\ell}}^2 V - (\partial_\ell \partial_{\bar{\ell}} V)^2}_{\ell(\beta), \bar{\ell}(\beta) \text{ indep. of } N_f, \mu} = (a\partial_\ell - b\partial_{\bar{\ell}})^3 V = 0 \quad (2)$$

with  $a = \partial_{\bar{\ell}}^2 V|_c$  and  $b = \partial_\ell \partial_{\bar{\ell}} V|_c$ . The first two equations rewrite

$$N_f f(\beta M) = e^{\beta\mu} \partial_\ell v_{\text{glue}} = e^{-\beta\mu} \partial_{\bar{\ell}} v_{\text{glue}} \longrightarrow \underbrace{e^{-2\beta\mu} = \partial_\ell v_{\text{glue}} / \partial_{\bar{\ell}} v_{\text{glue}}}_{\ell(\beta(\mu)), \bar{\ell}(\beta(\mu)) \text{ indep. of } N_f} \quad (3)$$

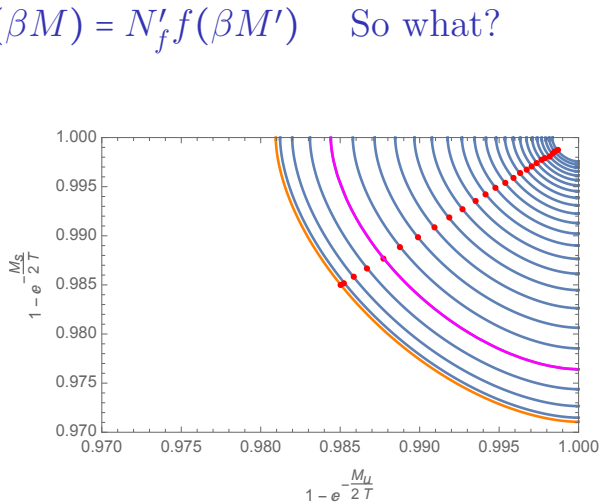
$\longrightarrow \boxed{N_f f(\beta M) = N'_f f(\beta M') = f(\beta M_s) + 2f(\beta M_{ud})}$  for each value of  $\mu$

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**Red:** Model-dep.  $N_f = 3$  input determined from:  $\partial_\ell V = \partial_\ell^2 V = \partial_\ell^3 V = 0$

**Blue:** Model-indep. line from:  $N_f f(\beta M) = f(\beta M_s) + 2f(\beta M_{ud})$

$$N_f f(\beta M) = N'_f f(\beta M') \quad \text{So what?}$$

If expanded in large  $R \equiv \beta M$ , allows for the simple relation

$$R_{N'_f} - R_{N_f} \approx \ln \frac{N'_f}{N_f} \quad \longrightarrow \quad Y_{N_f} \equiv \frac{R_{N_f} - R_1}{R_2 - R_1} \approx \frac{\ln N_f}{\ln 2}$$

- ▶ satisfied both in continuum approaches as well as on the lattice
- ▶ robust against higher order corrections in the large  $\beta M$  expansion
- ▶ independent of chemical potential
- ▶ predict  $R_{N_f}$  for  $N_f > 3$  or  $\notin \mathbb{Z}$

$Y_3$	$\mu = 0$	$\mu = i\pi T/3$
Lattice	1.59	1.59
GZ1	1.58	1.57
GZ2	1.58	1.58
Matrix	1.59	1.56
CF	1.58	1.57

$$Y_3 \approx \frac{\ln 3}{\ln 2} \approx 1.58$$

# Intermediate Summary for One-loop Models

- ▶ Heavy Quark region exhibits generic features among all one-loop models
- ▶  $T_c$  constant along the critical line, whose shape is completely fixed, independently of  $\mu$
- ▶ Flavor dependence of the critical mass is independent of the gluon dynamics, as predicted by the universal quantity  $Y_{N_f}$

Motivation

generic 1-loop

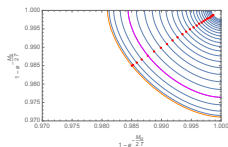
Curci-Ferrari at  
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Conclusion



$$Y_{N_f} \equiv \frac{R_{N_f} - R_1}{R_2 - R_1} \approx \frac{\ln N_f}{\ln 2}$$

Two assumptions were made:

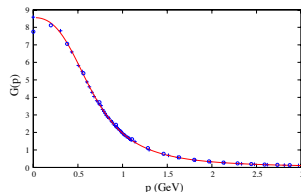
- ▶ large quark mass expansion
- ▶ quarks contribute at one-loop level

# Curci-Ferrari and gluon mass term

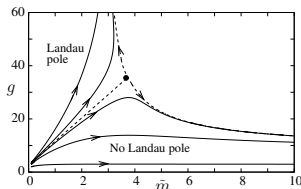
$$S = \int_x \left\{ \frac{1}{4} (F_{\mu\nu}^a)^2 + \bar{\psi} (\not{D} + M + \mu\gamma_0) \psi \right\} + S_{FP} + \int_x \left\{ \frac{1}{2} m^2 (A_\mu^a)^2 \right\}$$

This gluon mass term can be motivated in several ways

- ▶ phenomenologically from lattice data of the Landau gauge gluon propagator saturating in the IR
- ▶ Residual ambiguity after non-complete gauge-fixing in Fadeev-Popov procedure due to presence of Gribov copies



one-loop gluon propagator against lattice data,  
from [Tissier, Wschebor (2011)]



YM one-loop RG flow,  
from [Serreau, Tissier (2012)]

[Bogolubsky et al. (2009), Dudal, Oliveira,  
Vandersickel (2010) ]

# Landau-DeWitt gauge [Braun, Pawłowski, Gies (2010)]

$$A_\mu^a = \bar{A}_\mu^a + a_\mu^a$$

In practice, at each temperature, the background field  $\bar{A}_\mu^a$  is chosen such that the expectation value  $\langle a_\mu^a \rangle$  vanishes in the limit of vanishing sources.

This corresponds to finding the **absolute minimum of  $\tilde{\Gamma}[\bar{A}] \equiv \Gamma[\bar{A}, \langle a \rangle = 0]$** , where  $\Gamma[\bar{A}, \langle a \rangle]$  is the effective action for  $\langle a \rangle$  in the presence of  $\bar{A}$ .

Seek the minima in the subspace of configurations  $\bar{A}$  that respect the symmetries of the system at finite temperature.

→ One restricts to temporal and homogenous backgrounds:

$$\bar{A}_\mu(\tau, \mathbf{x}) = \bar{A}_0 \delta_{\mu 0}$$

→ functional  $\tilde{\Gamma}[\bar{A}]$  reduces to an effective potential  $V(\bar{A}_0)$  for the constant matrix field  $\bar{A}_0$ .

One can always rotate this matrix  $\bar{A}_0$  into the Cartan subalgebra:

$$\beta g \bar{A}_0 = r_3 \frac{\lambda_3}{2} + r_8 \frac{\lambda_8}{2}$$

	$r_3$	$r_8$
$\mu = 0$	$\mathbb{R}$	$0$
$\mu \in i\mathbb{R}$	$\mathbb{R}$	$\mathbb{R}$
$\mu \in \mathbb{R}$	$\mathbb{R}$	$i\mathbb{R}$

Then  $V(\bar{A}_0)$  reduces to a function of 2 components  **$V(r_3, r_8)$** .

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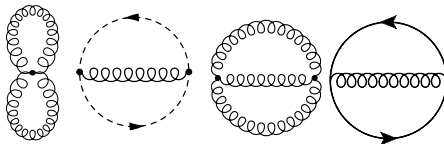
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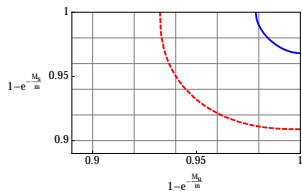
Conclusion

# Two-loop Expansion

$$\begin{aligned} V(r_3, r_8) = & -\text{Tr Ln}(\not{\partial} + M + \mu\gamma_0 - ig\gamma_0\bar{A}^k t^k) \\ & + \frac{3}{2}\text{Tr Ln}(\bar{D}^2 + m^2) - \frac{1}{2}\text{Tr Ln}(\bar{D}^2) \\ & + \end{aligned}$$



# Vanishing chemical potential



$$R_{N_f} \equiv \frac{M_c(N_f)}{T_c(N_f)}$$

$$\mathcal{O}(1): M_{\text{bare}} = M_{\text{ren.}}$$

$$\mathcal{O}(g^2): M_{\text{bare}} = Z_M M_{\text{ren.}} + C_M$$

→ hard to compare between different approaches!

However,  $Z_M, C_M$  are independent of  $N_f$  at  $\mathcal{O}(g^2)$ , and observing

$$\frac{T_c(N_f = 3) - T_c(N_f = 1)}{T_c(N_f = 1)} \approx 0.2\%$$

allows for:

$$\overbrace{R_{N_f}'/R_{N_f} \approx M_c(N_f')/M_c(N_f)}^{\text{if } C_M=0} \qquad \overbrace{Y_{N_f} \equiv \frac{R_{N_f} - R_1}{R_2 - R_1}}^{\text{if } C_M \neq 0}$$

is scheme indep. & comparable to other approaches up to higher order corrections.



# Vanishing chemical potential

$\mu = 0$	$R_2/R_1$	$R_3/R_1$	$Y_3$
Matrix [1]	1.10	1.16	1.59
GZ1 [2]	1.12	1.19	1.58
GZ2 [2]	1.08	1.13	1.58
CF 1-loop [3]	1.13	1.20	1.58
CF 2-loop [2]	1.12	1.18	1.57
Lattice [4]	1.10	1.15	1.59
DSE [5]	1.29	1.43	1.51

→ The  $Y_3$  values are still satisfied to very good approximation which underlines its importance as a universal quantity

→ The overall good agreement seems to suggest that the underlying dynamics is well-described within (Curci-Ferrari) perturbation theory.

[1] Kashiwa, Pisarski, Skokov (2012) [2] JM, Reinoso, Serreau (2017+18)

[3] Reinoso, Serreau, Tissier (2015) [4] Fromm, Langelage, Lottini, Philipsen (2012)

[5] Fischer, Luecker, Pawłowski (2015)

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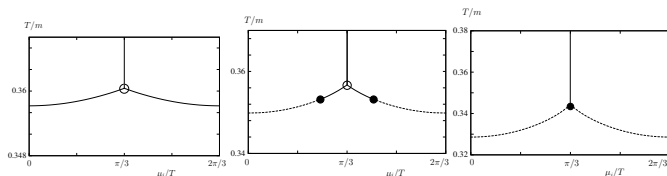
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# Imaginary chemical potential $\mu = i\mu_i$

Heavy Quark  
QCD

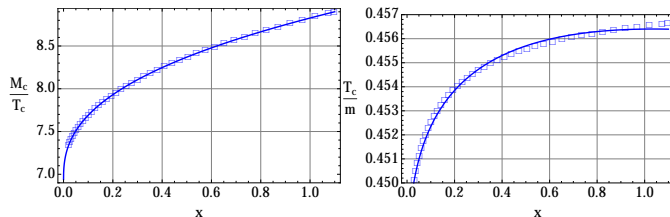
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The vicinity of the tricritical point is approximately described by the mean field scaling behavior

$$\frac{M_c(\mu_i)}{T_c(\mu_i)} = \frac{M_{\text{tric.}}}{T_{\text{tric.}}} + K \left[ \left( \frac{\pi}{3} \right)^2 - \left( \frac{\mu_i}{T_c} \right)^2 \right]^{\frac{2}{5}}$$

[de Forcrand, Philipsen (2010); Fischer, Luecker, Pawłowski (2015)]



$$x \equiv \left( \frac{\pi}{3} \right)^2 + \left( \frac{\mu_i}{T_c} \right)^2 = \left( \frac{\pi}{3} \right)^2 - \left( \frac{\mu_i}{T_c} \right)^2$$

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# Imaginary chemical potential $\mu = i\mu_i$

$\mu = i\pi T/3$	$R_2/R_1$	$R_3/R_1$	$Y_3$
Matrix [1]	1.18	1.28	1.56
GZ1 [2]	1.18	1.28	1.57
GZ2 [2]	1.11	1.17	1.58
CF 1-loop [3]	1.19	1.30	1.57
CF 2-loop [2]	1.17	1.27	1.57
Lattice [4]	1.12	1.20	1.59
DSE [5]	2.07	2.70	1.59

→ The  $Y_3$  values are in overall very good agreement between all cases, one loop models and higher order ones.

[1] Kashiwa, Pisarski, Skokov (2012) [2] JM, Reinos, Serreau (2017+18)

[3] Reinos, Serreau, Tissier (2015) [4] Fromm, Langelage, Lottini, Philipsen (2012)

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## One-loop:

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- ▶  $T_c$  constant along the critical line, whose shape is completely fixed, independently of  $\mu$
- ▶ Flavor dependence of the critical mass is independent of the gluon dynamics, as predicted by the universal quantity  $Y_{N_f}$

## Higher order:

- ▶ updated  $Y_3$  values still agree with one-loop predictions
- ▶ suggests that the perturbative description of the phase diagram within the CF model is robust

## Outlook:

- ▶ Can we describe the chiral transition in the lower left part of the Columbia plot?

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