The QCD CEP in the 3 flavoured constituent quark model

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Rab, aug. 30 - sept. 3, 2007

- Motivation for using effective models to describe the QCD CEP
- The model and its parametrization for zero and non-zero μ_B, μ_I, μ_Y chemical potential
- The location and the scaling region of the CEP at $\mu_I = \mu_Y = 0$
- Introduction of the chemical potentials: μ_B, μ_I, μ_Y
- Effects of isospin breaking on the location of the CEP
- Conclusions

Contradicting lattice results, and the role of effective models

The physical point is in the crossover regime. Y. Aoki, *et al.*, Nature 443, 675 (2006) Envisaged chiral critical surfaces from lattice simulations:



The common expectation is that this surface bends towards the physical point. However negative curvature according to O. Philipsen, Ph. de Forcrand, hep-lat/0607017 CEP found at:

 $T_{\text{CEP}} = 160 \pm 3.5 \text{ MeV } \mu_{\text{B,CEP}} = 725 \pm 35 \text{ MeV}$, lattice volume: 12×4^3 and $m_{\pi} \approx 2m_{\pi}^{\text{phys}}$ Z. Fodor, S. D. Katz, JHEP 0203:014,2002 $T_{\text{CEP}} = 162 \pm 2 \text{ MeV } \mu_{\text{B,CEP}} = 360 \pm 40 \text{ MeV}$, lattice volume: 12×4^3 and $m_{\pi} = m_{\pi}^{\text{phys}}$ Z. Fodor, S. D. Katz, JHEP 0404:050,2004

The simulation at finite μ is very difficult. \longrightarrow qualitative difference between lattice results \longrightarrow role of effective models

Relevance of the study of the CEP



the CEP is experimentally accesible $\mu_B, \mu_I \neq 0$ in heavy ion collison experiments μ_B is tunable \rightarrow beam energy, centrality μ_I is tunable \rightarrow different isotopes of an element focusing effect: if CEP exist it cannot be missed

analogy to the CEP of a liquid-gas phase transition which is easy to hit

lattice simulation at μ_B is very difficult

 \implies not all the methods predict/find the CEP

CEP found at: $(T, \mu_B)_{\text{CEP}} = (162 \pm 2, 360 \pm 40) \text{ MeV}, \text{ volume: } 12 \times 4^3 \text{ and } m_{\pi} = m_{\pi}^{\text{phys}}$

Z. Fodor, S. D. Katz, JHEP 0404:050,2004

lattice simulation at μ_I is free of the sign problem

it is important to study the CEP and its μ_I , μ_Y dependence in effective models

Influence of μ_I on the $\mu_B - T$ diagram

Barducci et. al, PLB 564, 217

without $U(1)_A$ breaking \rightarrow generic result for low $T \mu_I$ induces two 1st order transitions \implies 2 critical endpoints



the structure cease to exist in case of a sufficiently strong $U(1)_A$ breaking

Frank et. al, PLB **562**, 221



$SU_L(3) imes SU_R(3)$ symmetric chiral quark model

$$\mathcal{L} = \frac{1}{2} \operatorname{Tr}(\partial_{\mu} M^{\dagger} \partial^{\mu} M + m_0^2 M^{\dagger} M) - f_1 \left(\operatorname{Tr}(M^{\dagger} M) \right)^2 - f_2 \operatorname{Tr}(M^{\dagger} M)^2 - g \left(\det(M) + \det(M^{\dagger}) \right) + \epsilon_0 \sigma_0 + \epsilon_3 \sigma_3 + \epsilon_8 \sigma_8 + \bar{\psi} \left(i \partial \!\!\!/ - g_F M_5 \right) \psi.$$

 $M = \frac{1}{\sqrt{2}} \sum_{i=0}^{8} (\sigma_i + i\pi_i) \lambda_i, M_5 = \sum_{i=0}^{8} \frac{1}{2} (\sigma_i + i\gamma_5\pi_i) \lambda_i \quad 3 \times 3 \text{ complex matrices}$

pseudo(scalar) fields: π_i , σ_i , quark field: $\overline{\psi} = (u, d, s)$

Gell-Mann matrices: $\lambda_0 := \sqrt{\frac{2}{3}}\mathbf{1}, \lambda_i : i = 1 \dots 8.$

determinant breaks $U_A(1)$ symmetry explicit symmetry breaking: external fields $\epsilon_0, \epsilon_3, \epsilon_8 \neq 0 \iff m_u \neq m_d \neq 0, m_s \neq 0$

broken symmetry phase: three condensates $(\langle \sigma_0 \rangle, \langle \sigma_8 \rangle), \langle \sigma_3 \rangle \longleftrightarrow (x, y), v_3$

fermion masses:
$$M_u = \frac{g_F}{2}(x+v_3), M_d = \frac{g_F}{2}(x-v_3), M_s = \frac{g_F y}{\sqrt{2}}$$

technical difficulty: mixing in the 0, 3, 8 sector

parameters determined from the T = 0 mass spectrum

Parametrization and thermodynamics at one-loop level

13 unknown parameters:

$$\begin{array}{rl} \mbox{couplings} & m_0^2, f_1, f_2, g, g_F \\ \mbox{condensates} & x, y, v_3 \\ \mbox{external fields} & \epsilon_x, \epsilon_y, \epsilon_3 \\ \mbox{renormalization scales} & l_f, l_b \end{array}$$

resummation using optimized perturbation theory Chiku & Hatsuda, PRD58:076001

change:
$$-m_0^2 \to m^2 \Rightarrow \mathcal{L}_{mass} = \frac{1}{2}m^2 \mathrm{Tr}M^{\dagger}M - \frac{1}{2}\underbrace{(m_0^2 + m^2)\mathrm{Tr}M^{\dagger}M}_{\Delta m^2: \text{ one-loop counterterm}}$$

principle of minimal sensitivity $M_{\pi}^2 = iG^{-1}(p^2=0)|_{1-loon} \stackrel{!}{=} m_{\pi}^2|_{tree} \Longrightarrow$ equation for the effective mass:

$$m^2 = -m_0^2 + \Sigma_\pi(p = 0, m_i(m^2), M_q)$$

From the tree-level pion mass: $m^2 = m_{\pi}^2 - (4f_1 + 2f_2)x^2 - 4f_1y^2 - 2gy$

 \implies introducing into the other tree-level masses \implies self-consistent gap equation for the pion mass

Set of coupled nonlinear equations (for $v_3 = 0$):

(1) gap-equation: $m_{\pi}^2 = -m_0^2 + (4f_1 + 2f_2)x^2 + 4f_1y^2 + 2gy + \text{Re}\Sigma_{\pi}(p=0, m_i(m_{\pi}), M_u)$

(2) pole-mass
$$M_K$$
 from:
 $M_K^2 = -m_0^2 + 2(2f_1 + f_2)(x^2 + y^2) + 2f_2y^2 - \sqrt{2}x(2f_2y - g) + \text{Re}\Sigma_K(p^2 = M_K^2, m_i)$

(3) FAC criterion for M_K : $\Sigma(p^2 = M_K^2) = 0$

(4) pole-mass
$$M_{\eta}$$
 from:

$$Det \begin{pmatrix} p^2 - m_{\eta_{xx}}^2 - \Sigma_{\eta_{xx}}(p^2, m_i) & -m_{\eta_{xy}}^2 - \Sigma_{\eta_{xy}}(p^2, m_i) \\ -m_{\eta_{xy}}^2 - \Sigma_{\eta_{xy}}(p^2, m_i) & p^2 - m_{\eta_{yy}}^2 - \Sigma_{\eta_{yy}}(p^2, m_i) \end{pmatrix} \Big|_{p^2 = M_{\eta}^2, M_{\eta'}} = 0$$

(5) PCAC: $x = f_{\pi}$

(6) From non-strange quark mass: $g_F = \frac{2M_u}{x}$

(7) From strange quark mass: $y = \frac{\sqrt{2}M_s}{g_F}$

(8) EOS for x:

 $\epsilon_x = -m_0^2 x + 2gxy + 4f_1 xy^2 + 2(2f_1 + f_2)x^3 + \sum_{\alpha,i,j} t_{\alpha_{i,j}}^x \langle \alpha_i \alpha_j \rangle + \frac{g_F}{2} (\langle \bar{u}u \rangle + \langle \bar{d}d \rangle)$

(9) EOS for y: $\epsilon_y = -m_0^2 y + gx^2 + 4f_1 x^2 y + 4(f_1 + f_2) y^3 + \sum_{\alpha,i,j} t^y_{\alpha_{i,j}} \langle \alpha_i \alpha_j \rangle + \frac{g_F}{\sqrt{2}} \langle \bar{s}s \rangle$

Differences in case of isospin breaking

New variable: v_3 Equation for $v_3 \longrightarrow$ third EoS:

$$\langle \frac{\partial \mathcal{L}}{\partial \sigma_3} \rangle = 0 \tag{1}$$

Even if $\epsilon_3 = 0 \iff v_3 = 0$ at T = 0) non zero μ_I will generate v_3 at non zero temperature

Cosequence: charged and neutral particle masses will be different at tree level

If explicit isospin breaking is also introduced another equation is needed:

$$m_{\pi^+,\text{tree}} - m_{\pi^0,\text{tree}} = 4.594 MeV$$
 (2)

This equation will determine v_3 at T = 0 and EoS for v_3 at T = 0 will determine ϵ_3

Deviation from the physical mass spectrum

The remaining two unknown parameters, l_f and l_b are determined trough accurate parametrization

Better parametrization \longleftrightarrow closer to the physical spectrum

$$\begin{split} R = \frac{1}{|T|} \sum_{i \in T} \frac{|m_i^{\text{tree}} - m_i^{\text{phys}}|}{m_i^{\text{phys}}} + \frac{1}{|L|} \sum_{i \in L} \frac{|m_i^{\text{tree}} - m_i^{1-\text{loop}}|}{m_i^{\text{tree}}}, \\ T = \{\eta, \eta', a_0, f_0, \sigma\}, L = \{\eta', a_0, \kappa, f_0\}, |T| = 5, |L| = 4. \end{split}$$

Physical mass spectrum:

$$\begin{array}{ll} m_{\pi} = 138 \; {\rm MeV} & m_{a_0} = 980 \; {\rm MeV} \\ m_K = 495.6 \; {\rm MeV} & m_{\kappa} = 900 \; {\rm MeV} \\ m_{\eta} = 547.8 \; {\rm MeV} & m_{f_0} = 1370 \; {\rm MeV} \\ m_{\eta'} = 958 \; {\rm MeV} & m_{\sigma} = 700 \; {\rm MeV} \end{array}$$

The closer we are to the physical spectrum the smaller R we get.

 \implies We have located the minimum of R.



Star on fig.: $l_b = 520 \text{ MeV}$, $l_f = 1210 \text{ MeV} \longrightarrow \text{corresponds}$ to the minimum of R1-loop masses of σ and f_0 : $m_{\sigma} = 614.2 \text{ MeV}$, $m_{f_0} = 1210.9 \text{ MeV}$

Close to the physical spectrum, the phase transition is of first order / crossover type on the T = 0 / $\mu_B = 0$ axes.

The surface of 2^{nd} order phase transition in the $m_{u,d} - m_s - \mu_B$ space

Away from the physical point we re-parametrized the model using CHPT

 \longrightarrow for mesons in the large N_c limit for f_{π} , m_{η} : P. Herrera-Siklódy *et al.*, PLB 419 (1998) 326

$$f_{\pi} = f\left(1 + 4L_5 \frac{m_{\pi}^2}{f^2}\right)$$
$$m_{\eta}^2 = \frac{4m_K^2 - m_{\pi}^2}{3} + \frac{32}{3}(2L_8 - L_5) \frac{\left(m_K^2 - m_{\pi}^2\right)^2}{f^2},$$

 \longrightarrow for baryons ($B \in \{N, \Sigma, \Lambda, \Xi, \}$):

V. Bernard et al., Int. J. Mod. Phys. E4, 193 (1995)

$$M_B = M_0 - 2b_0(m_{\pi,2}^2 + m_{K,2}^2) + b_D \gamma_B^D(m_\pi, m_K) + b_F \gamma_B^F(m_\pi, m_K)
onumber \ - rac{1}{24\pi f^2} \left[lpha_B^\pi m_\pi^3 + lpha_B^K m_K^3 + lpha_B^\eta m_\eta^3
ight]$$

The constituent kvark masses:

$$egin{array}{rcl} m_u &=& rac{M_N(m_\pi,m_K)}{3} \ m_s &=& rac{M_\Lambda(m_\pi,m_K)+M_\Sigma(m_\pi,m_K)}{2}-2M_u(m_\pi,m_K) \end{array}$$



The surface bends towards the physical point \implies The CEP must exist The continuation is reliable up to $m_K \approx 400$ MeV and above the diagonal

The CEP at the physical point of the mass plane

P. Kovács, Zs. Szép: Phys. Rev. D 75, 025015



- $T_c(\mu_B = 0) = 154.84 \text{ MeV}$ $\Delta T_c(x\chi) = 15.5 \text{ MeV}$
- $T_{CEP} = 74.83 \text{ MeV}$ $\mu_{B,CEP} = 895.38 \text{ MeV}$
- $T_c \frac{d^2 T_c}{d\mu_B^2}\Big|_{\mu_B=0} = -0.09$

- $T_c(\mu_B = 0) = 151(3) \text{ MeV}$ $\Delta T_c(\chi_{\bar{\psi}\psi}) = 28(5) \text{ MeV}$ Y. Aoki,*et al.*, PLB **643**, 46 (2006)
- $T_{CEP} = 162(2) \text{ MeV}$ $\mu_{B,CEP} = 360(40) \text{ MeV}$
- -0.058(2)
 - Z. Fodor, et al., JHEP 0404 (2004) 050

Dependence of the $\mu_{B,CEP}$ on the width of the susceptibility



Preliminary lattice estimation by S. Katz: $\Delta T_c(\chi_{\bar{\psi}\psi}) \approx 0.5 - 1 \text{ MeV}$ $\Delta T_c(\chi_{\bar{\psi}\psi}) \approx 2 - 4 \text{ MeV}$

Since $\Delta T_c(\chi_{\bar{\psi}\psi}) \approx 28$ MeV at the physical point \longrightarrow higher $\mu_{B,CEP}$ expected

The critical region of the CEP



For the asymptotically parallel path we get $\gamma = 1.01$, which corresponds to the mean-field Ising exponent.

 \longrightarrow This path is the tangent line of the phase boundary curve at the CEP in the $\mu_B - T$ plane.

Introduction of chemical potentials

21 particles:

Lagrangian is invariant under

$$M \rightarrow e^{-i\alpha_G G} M e^{i\alpha_G G} = M - i\alpha_G [G, M] + \mathcal{O}(\alpha_G^2),$$

$$\psi \rightarrow e^{-i\alpha_G G} \psi = \psi - i\alpha_G \psi + \mathcal{O}(\alpha_G^2),$$

where G can be $B = \sqrt{\frac{3}{2}}\lambda_0$, $I = \frac{1}{2}\lambda_3$ and $Y = \frac{1}{\sqrt{3}}\lambda_8$

The conserved Noether currents:

$$J^G_{\mu} = -\frac{\delta L}{\delta(\partial^{\mu}M)_{ij}} i[G,M]_{j,i} - \frac{\delta L}{\delta(\partial^{\mu}M^+)_{ij}} i[G,M^+]_{j,i} - \frac{\delta L}{\delta(\partial^{\mu}\psi_i)} iG_{ij}\psi_j$$

The conserved charges:

$$\begin{aligned} Q^B &= \frac{1}{3} (N_u + N_d + N_s - N_{\bar{u}} - N_{\bar{d}} - N_{\bar{s}}), \\ Q^I &= \frac{1}{2} (N_u - N_{\bar{u}} - N_d + N_{\bar{d}} + N_{\kappa^+} - N_{\kappa^-} + N_{\bar{\kappa}^0} - N_{\kappa^0} + N_{K^+} - N_{K^-} + N_{\bar{K}^0} - N_{K^0}) \\ &+ N_{a_0^+} - N_{a_0^-} + N_{\pi^+} - N_{\pi^-}, \\ Q^Y &= \frac{1}{3} (N_u - N_{\bar{u}} + N_d - N_{\bar{d}} - 2N_s + 2N_{\bar{s}}) + N_{\kappa^+} - N_{\kappa^-} + N_{\kappa^0} - N_{\bar{\kappa}^0} + N_{K^+} - N_{K^-} + N_{K^0} - N_{\bar{K}^0}) \end{aligned}$$

Statistical density matrix of the system:

$$\rho = \exp[-\beta(H - \mu_i N_i)]$$

The following chemical potencials can be introduced:

$$\begin{split} \mu_{u} &= -\mu_{\bar{u}} = \frac{1}{3}\mu_{B} + \frac{1}{2}\mu_{I} + \frac{1}{3}\mu_{Y}, \\ \mu_{d} &= -\mu_{\bar{d}} = \frac{1}{3}\mu_{B} - \frac{1}{2}\mu_{I} + \frac{1}{3}\mu_{Y}, \\ \mu_{s} &= -\mu_{\bar{s}} = \frac{1}{3}\mu_{B} - \frac{2}{3}\mu_{Y}, \\ \mu_{a_{0}^{+}} &= \mu_{\pi^{+}} = -\mu_{a_{0}^{-}} = -\mu_{\pi^{-}} = \mu_{I}, \\ \mu_{\kappa^{+}} &= \mu_{K^{+}} = -\mu_{\kappa^{-}} = -\mu_{K^{-}} = \frac{1}{2}\mu_{I} + \mu_{Y}, \\ \mu_{\kappa^{0}} &= \mu_{K^{0}} = -\mu_{\bar{\kappa}^{0}} = -\mu_{\bar{K}^{0}} = -\frac{1}{2}\mu_{I} + \mu_{Y} \end{split}$$

Finite temperature propagators of charged fields

For example the K^-, K^+ field operators:

$$K^{-}(x) = \int \frac{d^{3}\mathbf{p}}{(2\pi)^{3}} \frac{1}{\sqrt{2E_{\mathbf{p}}}} \left(a^{+}(\mathbf{p})e^{ip\cdot x} + b(\mathbf{p})e^{-ip\cdot x} \right) \Big|_{p_{0}=E_{\mathbf{p}}},$$

$$K^{+}(x) = \int \frac{d^{3}\mathbf{p}}{(2\pi)^{3}} \frac{1}{\sqrt{2E_{\mathbf{p}}}} \left(b^{+}(\mathbf{p})e^{ip\cdot x} + a(\mathbf{p})e^{-ip\cdot x} \right) \Big|_{p_{0}=E_{\mathbf{p}}},$$

The two-point functions:

 $\begin{aligned} G_{K^{-}}(y-x) &:= \langle TK^{-}(y)K^{+}(x)\rangle_{\beta} &= \Theta(y_{0}-x_{0})\langle K^{-}(y)K^{+}(x)\rangle_{\beta} + \Theta(x_{0}-y_{0})\langle K^{+}(x)K^{-}(y)\rangle_{\beta}, \\ G_{K^{+}}(y-x) &:= \langle TK^{+}(y)K^{-}(x)\rangle_{\beta} &= \Theta(y_{0}-x_{0})\langle K^{+}(y)K^{-}(x)\rangle_{\beta} + \Theta(x_{0}-y_{0})\langle K^{-}(x)K^{+}(y)\rangle_{\beta}, \end{aligned}$

In momentum space the finite temperature propagators:

$$\begin{aligned} G_{K^{-}}(k) &= \frac{i}{2E_{\mathbf{k}}} \left[\frac{1 + n_{K^{-}}(E_{\mathbf{k}})}{k_{0} - E_{\mathbf{k}} + i\epsilon} - \frac{n_{K^{-}}(E_{\mathbf{k}})}{k_{0} - E_{\mathbf{k}} - i\epsilon} - \frac{1 + n_{K^{+}}(E_{\mathbf{k}})}{k_{0} + E_{\mathbf{k}} - i\epsilon} + \frac{n_{K^{+}}(E_{\mathbf{k}})}{k_{0} + E_{\mathbf{k}} + i\epsilon} \right] \\ G_{K^{+}}(k) &= \frac{i}{2E_{\mathbf{k}}} \left[\frac{1 + n_{K^{+}}(E_{\mathbf{k}})}{k_{0} - E_{\mathbf{k}} + i\epsilon} - \frac{n_{K^{+}}(E_{\mathbf{k}})}{k_{0} - E_{\mathbf{k}} - i\epsilon} - \frac{1 + n_{K^{-}}(E_{\mathbf{k}})}{k_{0} + E_{\mathbf{k}} - i\epsilon} + \frac{n_{K^{-}}(E_{\mathbf{k}})}{k_{0} + E_{\mathbf{k}} - i\epsilon} \right] \end{aligned}$$

Self-energies









Temperature dependence of v_3



On the left Fig.: μ_B dpendence at a given μ_I Lowest curve correspond to a CEP v_3 at T = 0 significantly depend on μ_B

On the left Fig.: μ_I dpendence at a given μ_B

Increasing of either μ_B or $\mu_I \longrightarrow$ influence of v_3 becomes stronger CEP at $\mu_I = 0$: $T_{CEP} = 63.08$ MeV, $\mu_{B,CEP} = 960.8$ MeV \rightarrow large diff. to case $v_3 = 0$

Reason: x and v_3 related \longrightarrow common transition point

Dependence of the CEP on μ_I, μ_Y



 T_{CEP} is almost independent of μ_Y , but significantly depend on μ_I

 $\mu_{B,\text{CEP}}$ has an almost linear dependence on both other chenical potential

As μ_Y is increased the phase transition at T = 0 becomes stronger

Conclusions and outlook

- The best parametrization of the model gives first order / crossover type phase transition at $T = 0 / \mu_B = 0$ as a function of μ_B / T of the physical point.
- The 2^{nd} order surface was determined in the $m_{\pi} m_K \mu_B$ space using ChPT to obtain the m_{π}, m_K dependence of the couplings and of the constituent quark masses.
- The CEP was located at the physical point: $T_{CEP} = 74.83 \text{ MeV } \mu_{B,CEP} = 895.38 \text{ MeV}.$
- The dependence of the μ_B on the width of the susceptibility was investigated.
- The scaling properities were studied and the Ising temperature direction was found at the CEP.
- Effects of isospin and hyper chemical potential on the CEP was investigated. $T_{CEP} = 63.08 \text{ MeV } \mu_{B,CEP} = 960.8 \text{ MeV}$ at $\mu_I = 0$ ($v_3 \neq 0$ at T = 0).
- In progress: μ_I dependence of different pole masses and the study of pion condensation.