

Phase diagram of strong matter at finite chemical potentials

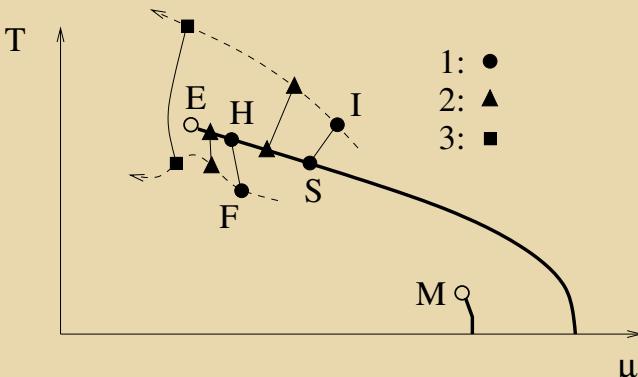
Péter Kovács

KFKI Research Inst. of Part. and Nucl. Phys. of HAS, Theoretical Department

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- Motivation for investigating QCD at non-zero μ_I, μ_Y
- The constituent quark model and its parameterization
- Introduction of the chemical potentials: μ_B, μ_I, μ_Y
- Results at $\mu_I = \mu_Y = 0$
- Results at non-zero μ_I, μ_Y
- Conclusions

Relevance of the isospin and strange chemical potentials



- the CEP is experimentally accessible
- $\mu_B, \mu_I \neq 0$ in heavy ion collision experiments
- μ_B is tunable \rightarrow beam energy, centrality
- μ_I is tunable \rightarrow different isotopes of an element
- In some experiment even μ_Y plays role
- Focusing effect: if CEP exist it cannot be missed

Brookhaven AGS exp. Si+Au collision: at $\mu_B = 540$ MeV $\rightarrow \mu_Y \approx 150$ MeV

CERN SPS exp. Pb+Pb collision:

at $\mu_B = 233 - 266$ MeV $\rightarrow \mu_Y \approx 70 - 80$ MeV, $\mu_I \approx 12 - 13$ MeV

CMB exp. at FAIR will explore QCD phase diagram at high μ_B

Analogy to the QCD CEP \rightarrow liquid-gas phase transition which is easy to hit

lattice simulations at finite chemical potential is very difficult

\Rightarrow not all the methods predict/find the CEP

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

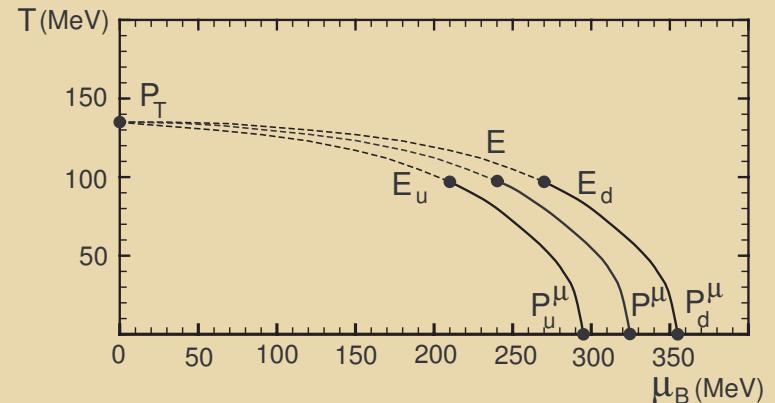
Z. Fodor, S. D. Katz, JHEP 0404:050,2004

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

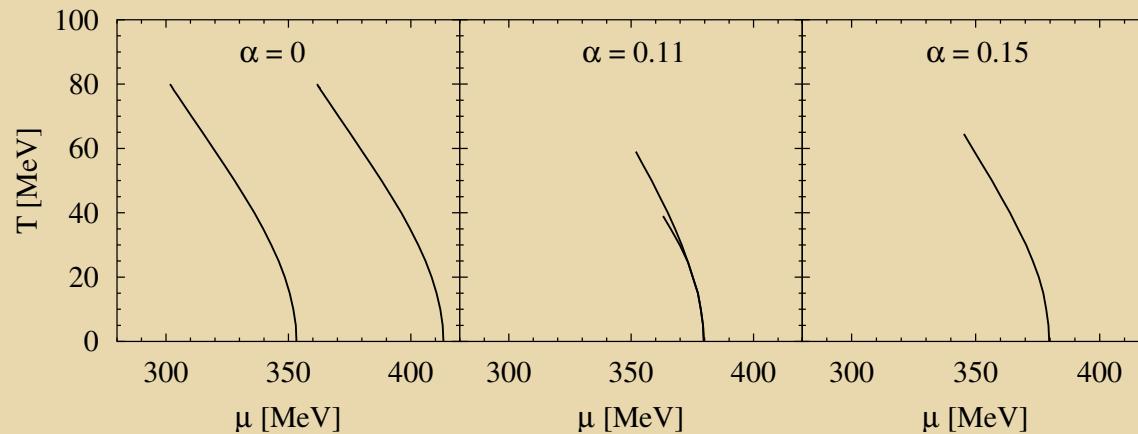
Barducci et. al, PLB 564, 217

without $U(1)_A$ breaking \rightarrow generic result
for low T μ_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



$SU_L(3) \times SU_R(3)$ symmetric chiral quark model

$$\begin{aligned}\mathcal{L} = & \frac{1}{2} \text{Tr}(\partial_\mu M^\dagger \partial^\mu M + m_0^2 M^\dagger M) - f_1 (\text{Tr}(M^\dagger M))^2 - f_2 \text{Tr}(M^\dagger M)^2 \\ & - g (\det(M) + \det(M^\dagger)) + \epsilon_0 \sigma_0 + \epsilon_3 \sigma_3 + \epsilon_8 \sigma_8 + \bar{\psi} (i\partial - g_F M_5) \psi.\end{aligned}$$

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

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

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 $(\langle \sigma_0 \rangle, \langle \sigma_8 \rangle), \langle \sigma_3 \rangle \longleftrightarrow (x, y), v_3$

x: non-strange, y: strange

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

Parameterization and thermodynamics at one-loop level

13 unknown parameters:

couplings	m_0^2, f_1, f_2, g, g_F
condensates	x, y, v_3
external fields	$\epsilon_x, \epsilon_y, \epsilon_3$
renormalization scales	l_f, l_b

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

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

fastest apparent convergence $M_\pi^2 = iG^{-1}(p^2=0)|_{\text{1-loop}} \stackrel{!}{=} m_\pi^2|_{\text{tree}} \implies \text{equation for the effective mass:}$

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

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_η from:

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

(5) tree-level 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:

$$\epsilon_x = -m_0^2x + 2gxy + 4f_1xy^2 + 2(2f_1 + f_2)x^3 + \sum_{\alpha,i,j} t_{\alpha_i,j}^x \langle \alpha_i \alpha_j \rangle + \frac{g_F}{2}(\langle \bar{u}u \rangle + \langle \bar{d}d \rangle)$$

(9) EOS for y: $\epsilon_y = -m_0^2y + gx^2 + 4f_1x^2y + 4(f_1 + f_2)y^3 + \sum_{\alpha,i,j} t_{\alpha_i,j}^y \langle \alpha_i \alpha_j \rangle + \frac{g_F}{\sqrt{2}}\langle \bar{s}s \rangle$

Differences in case of isospin breaking

New variable: v_3

Equation for $v_3 \rightarrow$ third EoS:

$$\left\langle \frac{\partial \mathcal{L}}{\partial \sigma_3} \right\rangle = 0 \quad (1)$$

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

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

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

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

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

Introduction of chemical potentials

21 particles:

pseudoscalars	$\pi^0, \eta, \eta', \pi^+, \pi^-, K^+, K^-, K^0, \bar{K}^0$
scalars	$\sigma, a_0^0, f_0, a_0^+, a_0^-, \kappa^+, \kappa^-, \kappa^0, \bar{\kappa}^0$
fermions	m_u, m_d, m_s

Lagrangian is invariant under

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

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:

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

The conserved charges:

$$Q^B = \frac{1}{3}(N_u + N_d + N_s - N_{\bar{u}} - N_{\bar{d}} - N_{\bar{s}}),$$

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

$$Q^Y = \frac{1}{3}(N_u - N_{\bar{u}} + N_d - N_{\bar{d}} - 2N_s + 2N_{\bar{s}}) + N_{\kappa^+} - N_{\kappa^-} + N_{\kappa^0} - N_{\bar{\kappa}^0} + N_{K^+} - N_{K^-} + N_{K^0} - N_{\bar{K}^0}$$

Statistical density matrix of the system:

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

The following chemical potentials can be introduced:

$$\mu_u = -\mu_{\bar{u}} = \frac{1}{3}\mu_B + \frac{1}{2}\mu_I + \frac{1}{3}\mu_Y,$$

$$\mu_d = -\mu_{\bar{d}} = \frac{1}{3}\mu_B - \frac{1}{2}\mu_I + \frac{1}{3}\mu_Y,$$

$$\mu_s = -\mu_{\bar{s}} = \frac{1}{3}\mu_B - \frac{2}{3}\mu_Y,$$

$$\mu_{a_0^+} = \mu_{\pi^+} = -\mu_{a_0^-} = -\mu_{\pi^-} = \mu_I,$$

$$\mu_{\kappa^+} = \mu_{K^+} = -\mu_{\kappa^-} = -\mu_{K^-} = \frac{1}{2}\mu_I + \mu_Y,$$

$$\mu_{\kappa^0} = \mu_{K^0} = -\mu_{\bar{\kappa}^0} = -\mu_{\bar{K}^0} = -\frac{1}{2}\mu_I + \mu_Y$$

Finite temperature propagators of charged fields

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

$$\begin{aligned} K^-(x) &= \int \frac{d^3\mathbf{p}}{(2\pi)^3} \frac{1}{\sqrt{2E_{\mathbf{p}}}} \left(a^+(\mathbf{p}) e^{ip \cdot x} + b(\mathbf{p}) e^{-ip \cdot x} \right) \Big|_{p_0=E_{\mathbf{p}}}, \\ K^+(x) &= \int \frac{d^3\mathbf{p}}{(2\pi)^3} \frac{1}{\sqrt{2E_{\mathbf{p}}}} \left(b^+(\mathbf{p}) e^{ip \cdot x} + a(\mathbf{p}) e^{-ip \cdot x} \right) \Big|_{p_0=E_{\mathbf{p}}} \end{aligned}$$

The two-point functions:

$$\begin{aligned} G_{K^-}(y-x) &:= \langle T K^-(y) K^+(x) \rangle_\beta = \Theta(y_0 - x_0) \langle K^-(y) K^+(x) \rangle_\beta + \Theta(x_0 - y_0) \langle K^+(x) K^-(y) \rangle_\beta, \\ G_{K^+}(y-x) &:= \langle T K^+(y) K^-(x) \rangle_\beta = \Theta(y_0 - x_0) \langle K^+(y) K^-(x) \rangle_\beta + \Theta(x_0 - y_0) \langle K^-(x) K^+(y) \rangle_\beta, \end{aligned}$$

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

$$-i\Sigma_{\pi^+} = \sum_{\substack{f \in (\sigma, \pi) \\ \alpha=0\dots 8}} -\pi^+ \begin{array}{c} \nearrow f_\alpha \\ \searrow \end{array} \pi^+ + \sum_{\delta=0,3,8} \pi^+ \begin{array}{c} \nearrow \sigma_\delta \\ \searrow \end{array} \pi^+ + \pi^+ \begin{array}{c} \nearrow \bar{\kappa}^0 \\ \searrow K^+ \end{array} \pi^+ + \pi^+ \begin{array}{c} \nearrow \kappa^+ \\ \searrow \bar{K}^0 \end{array} \pi^+ + \sum_{\delta=0,3,8} \pi^+ \begin{array}{c} \nearrow a_0^+ \\ \searrow \pi_\delta \end{array} \pi^+ + \pi^+ \begin{array}{c} \nearrow u \\ \searrow \bar{d} \end{array} \pi^+$$

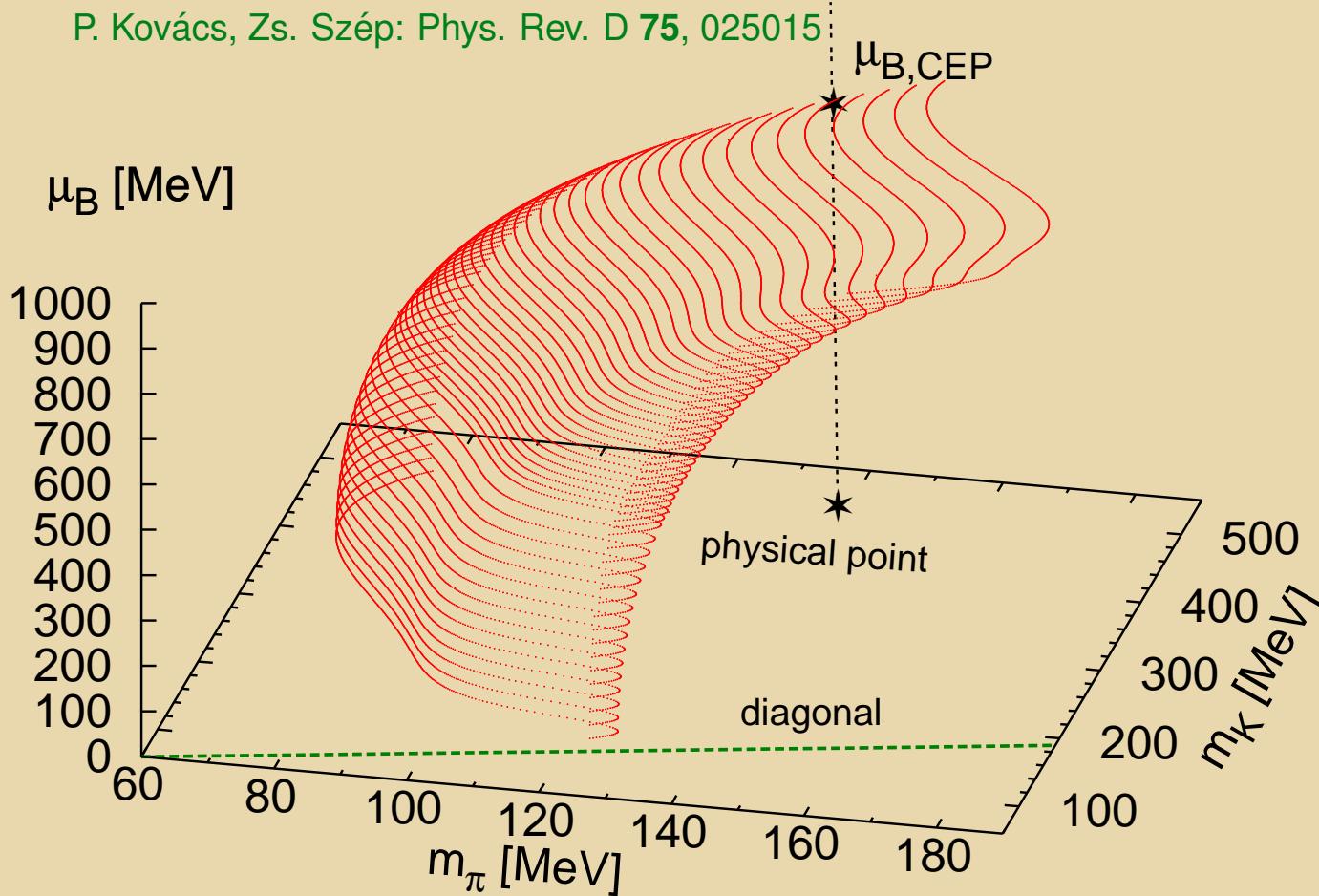
$$-i\Sigma_{0,3,8}^{\gamma\gamma'} = \sum_{\substack{f \in (\sigma, \pi) \\ \alpha=0\dots 8}} -\pi_\gamma \begin{array}{c} \nearrow f_\alpha \\ \searrow \end{array} \pi_\gamma + \sum_{\delta=0,3,8} \pi_\gamma \begin{array}{c} \nearrow a_0^- \\ \searrow \pi^+ \end{array} \pi_{\gamma'} + \pi_\gamma \begin{array}{c} \nearrow a_0^+ \\ \searrow \pi^- \end{array} \pi_{\gamma'} + \pi_\gamma \begin{array}{c} \nearrow \kappa^- \\ \searrow K^+ \end{array} \pi_{\gamma'} + \pi_\gamma \begin{array}{c} \nearrow \kappa^+ \\ \searrow K^- \end{array} \pi_{\gamma'} + \pi_\gamma \begin{array}{c} \nearrow \bar{\kappa}^0 \\ \searrow K^0 \end{array} \pi_{\gamma'}$$

$$+ \pi_\gamma \begin{array}{c} \nearrow \bar{\kappa}^0 \\ \searrow \bar{K}^0 \end{array} \pi_{\gamma'} + \sum_{\delta, \delta'=0,3,8} \pi_\gamma \begin{array}{c} \nearrow \sigma_\delta \\ \searrow \pi_{\delta'} \end{array} \pi_{\gamma'} + \sum_{q=u,d,s} \pi_\gamma \begin{array}{c} \nearrow q \\ \searrow \bar{q} \end{array} \pi_{\gamma'}$$

$$-i\Sigma_{K^+} = \sum_{\substack{f \in (\sigma, \pi) \\ \alpha=0\dots 8}} -K^+ \begin{array}{c} \nearrow f_\alpha \\ \searrow \end{array} K^+ + \sum_{\delta=0,3,8} K^+ \begin{array}{c} \nearrow \kappa^0 \\ \searrow \pi^+ \end{array} K^+ + \sum_{\delta=0,3,8} K^+ \begin{array}{c} \nearrow \sigma_\delta \\ \searrow K^+ \end{array} K^+ + K^+ \begin{array}{c} \nearrow a_0^+ \\ \searrow K^0 \end{array} K^+ + \sum_{\delta=0,3,8} K^+ \begin{array}{c} \nearrow \kappa^+ \\ \searrow \pi_\delta \end{array} K^+ + K^+ \begin{array}{c} \nearrow u \\ \searrow \bar{s} \end{array} K^+$$

Results at zero μ_I, μ_Y : critical surface and CEP

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

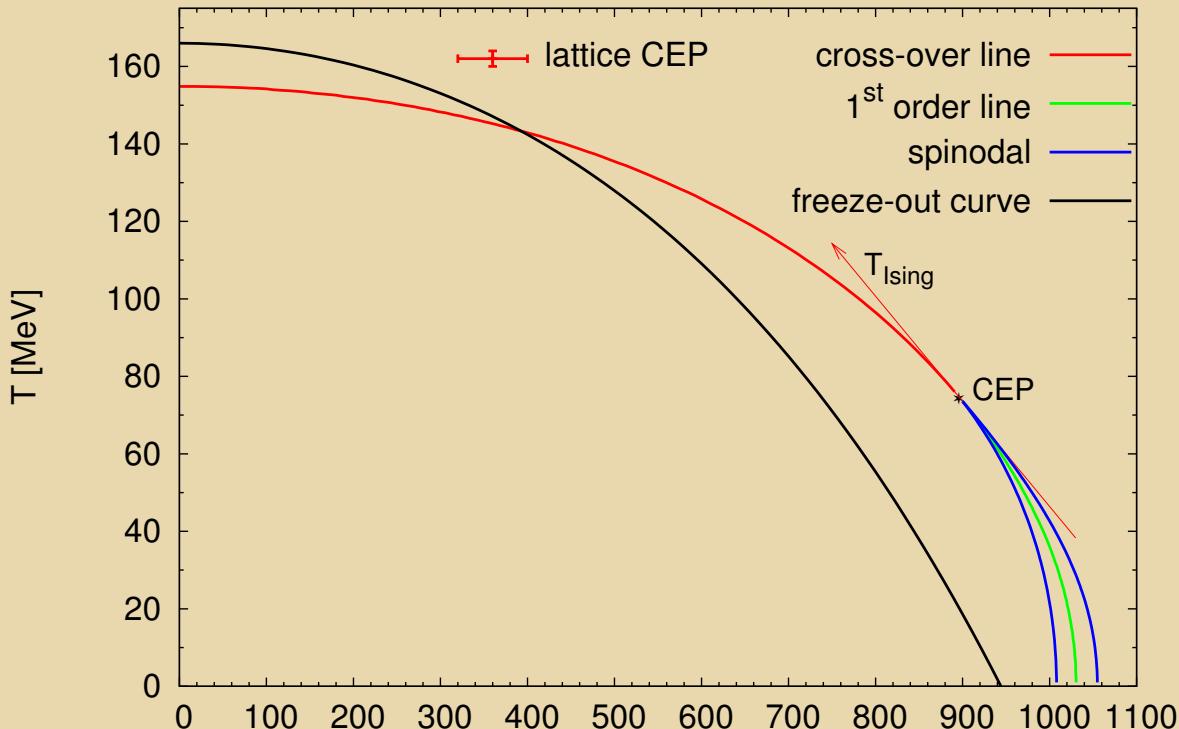


The surface bends towards the physical point \Rightarrow The CEP must exist

The continuation is reliable up to $m_K \approx 500$ MeV and above the diagonal

The CEP at the physical point of the mass plane

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



effective model

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

μ_B [MeV]

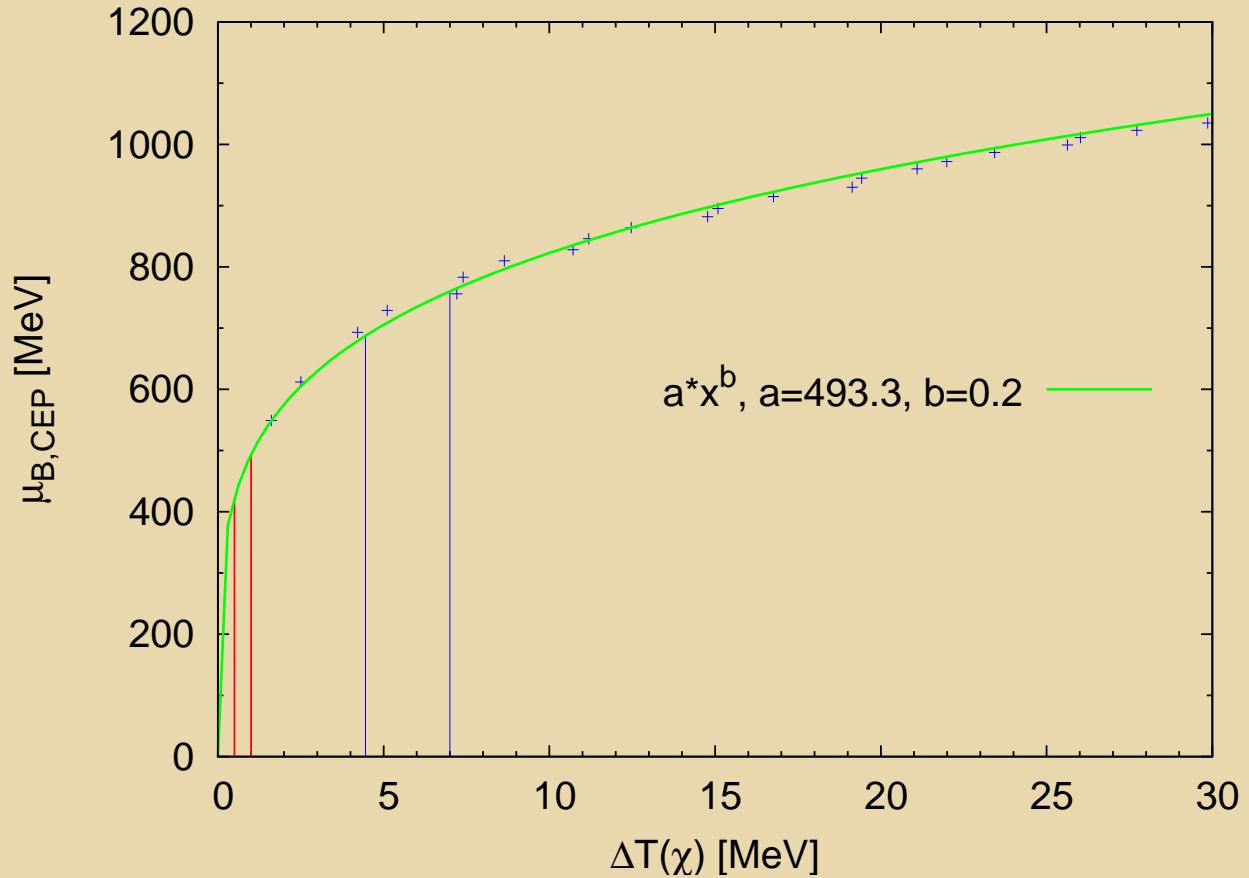
lattice

- $T_c(\mu_B = 0) = 151(3) \text{ MeV}$
 $\Delta T_c(\chi_{\bar{\psi}\psi}) = 28(5) \text{ MeV}$
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

$\mu_{B,CEP} = 725(35)$ MeV →
non-physical quark mass

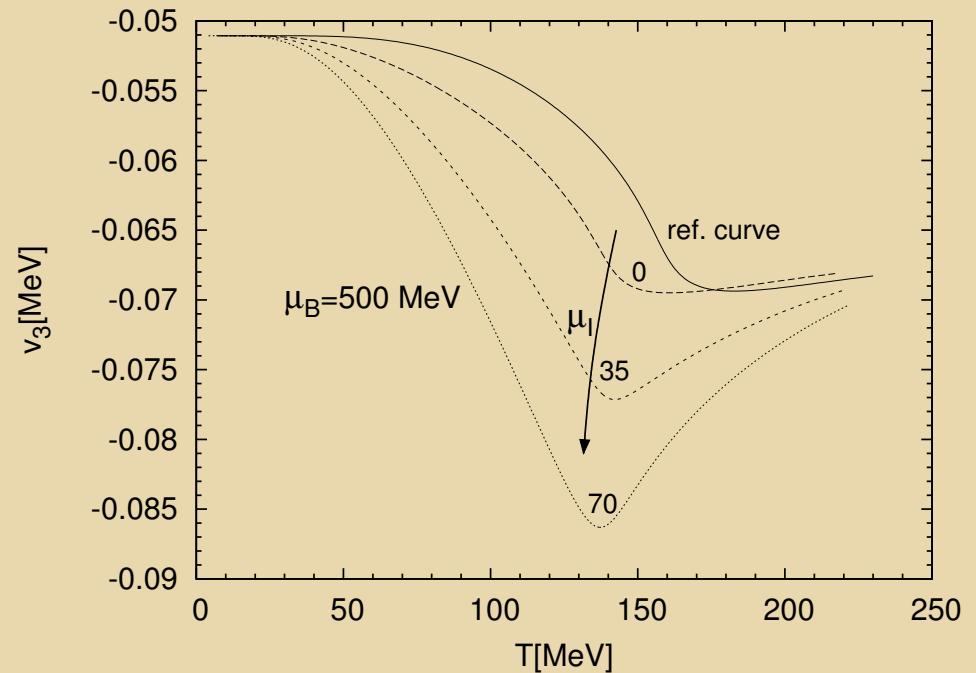
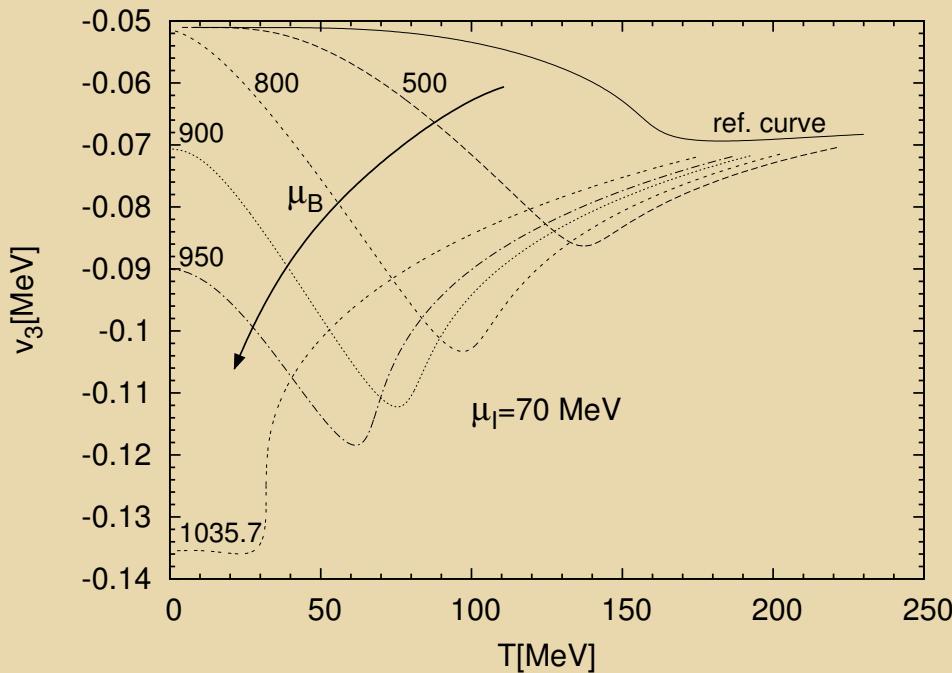
$\mu_{B,CEP} = 360(40)$ MeV →
physical quark mass



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

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

Temperature dependence of v_3



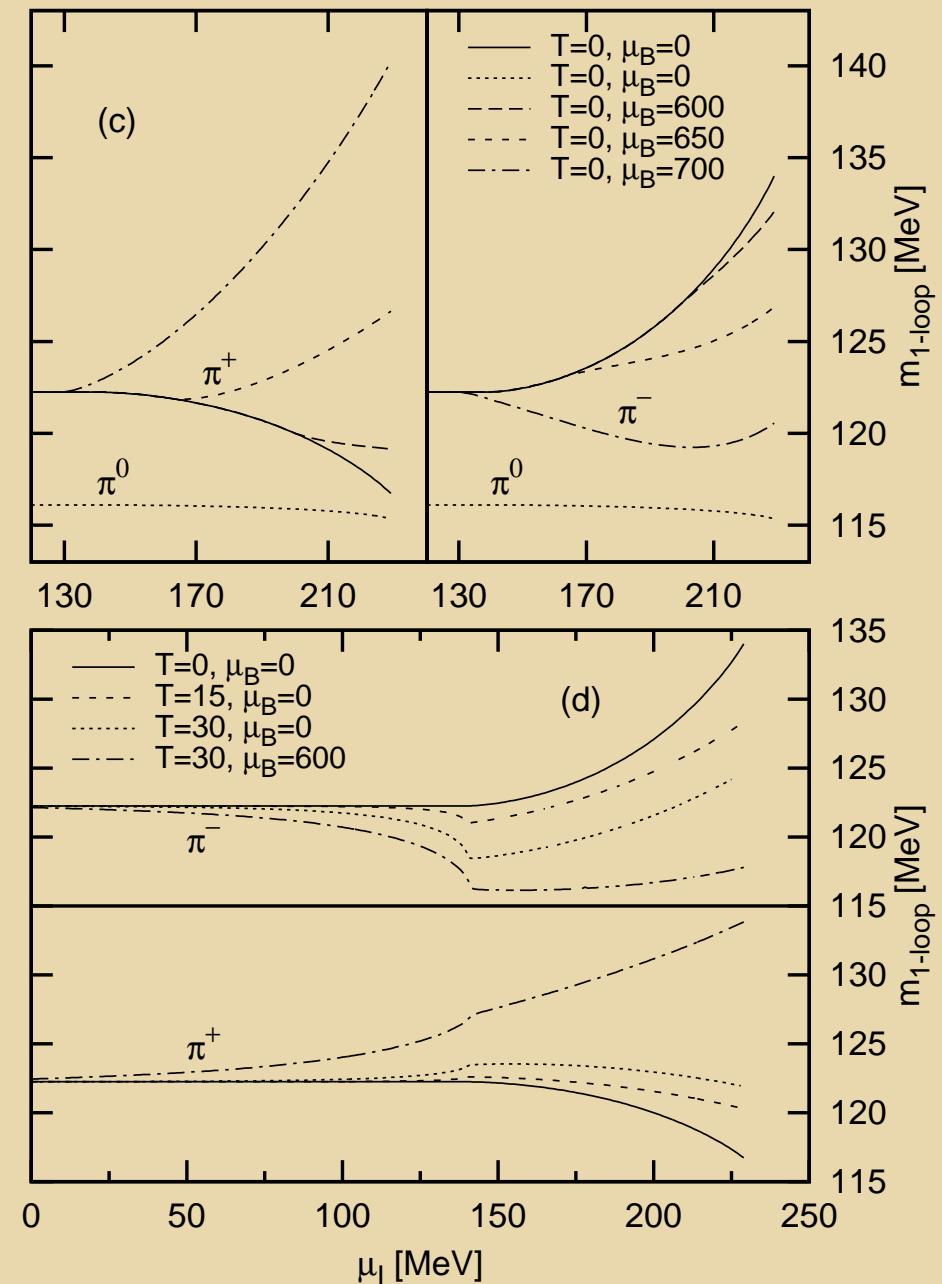
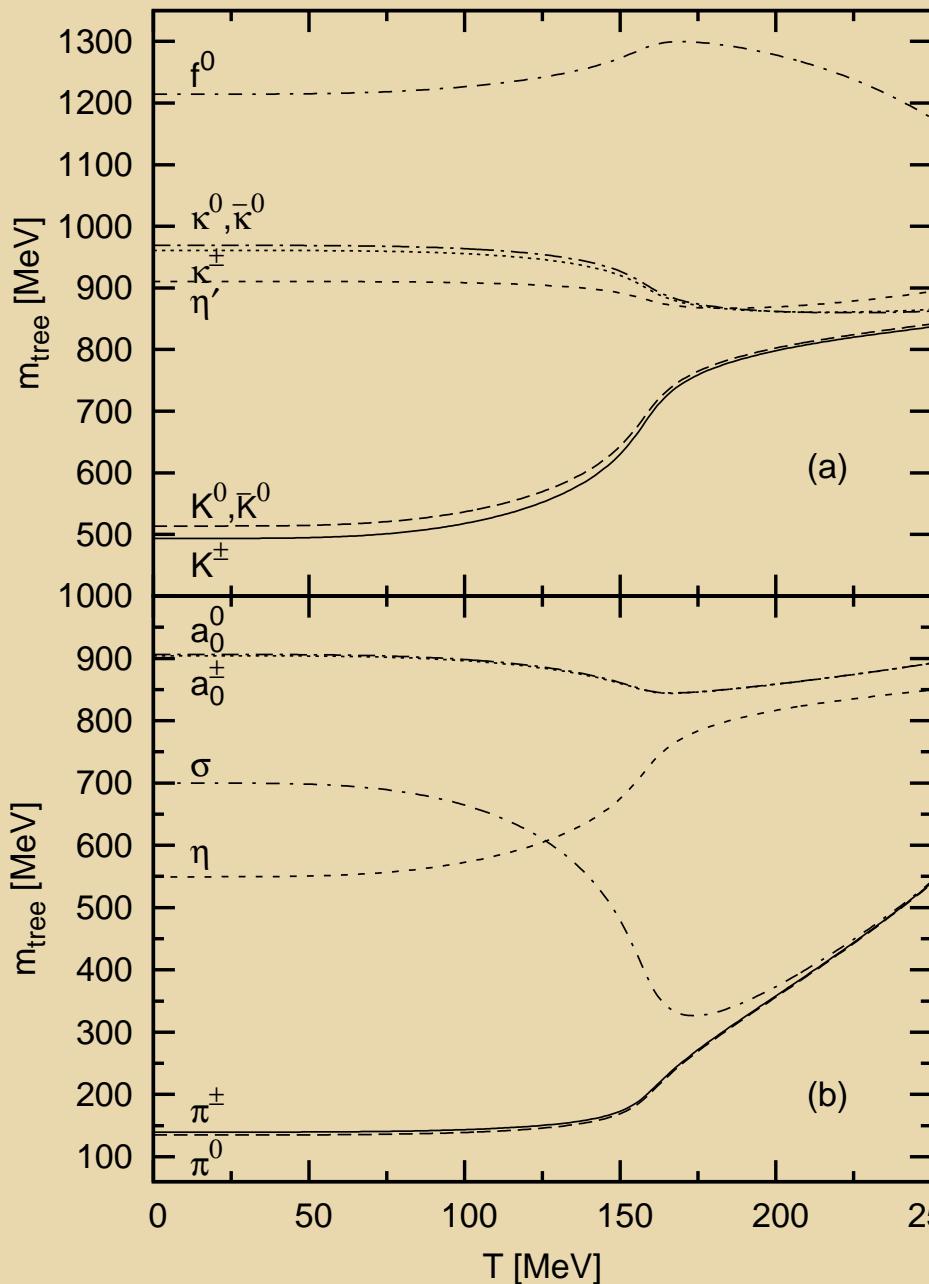
On the left Fig.: μ_B dependence at a given μ_I
 Lowest curve corresponds to a CEP
 v_3 at $T = 0$ significantly depend on μ_B

On the right Fig.: μ_I dependence at a given μ_B

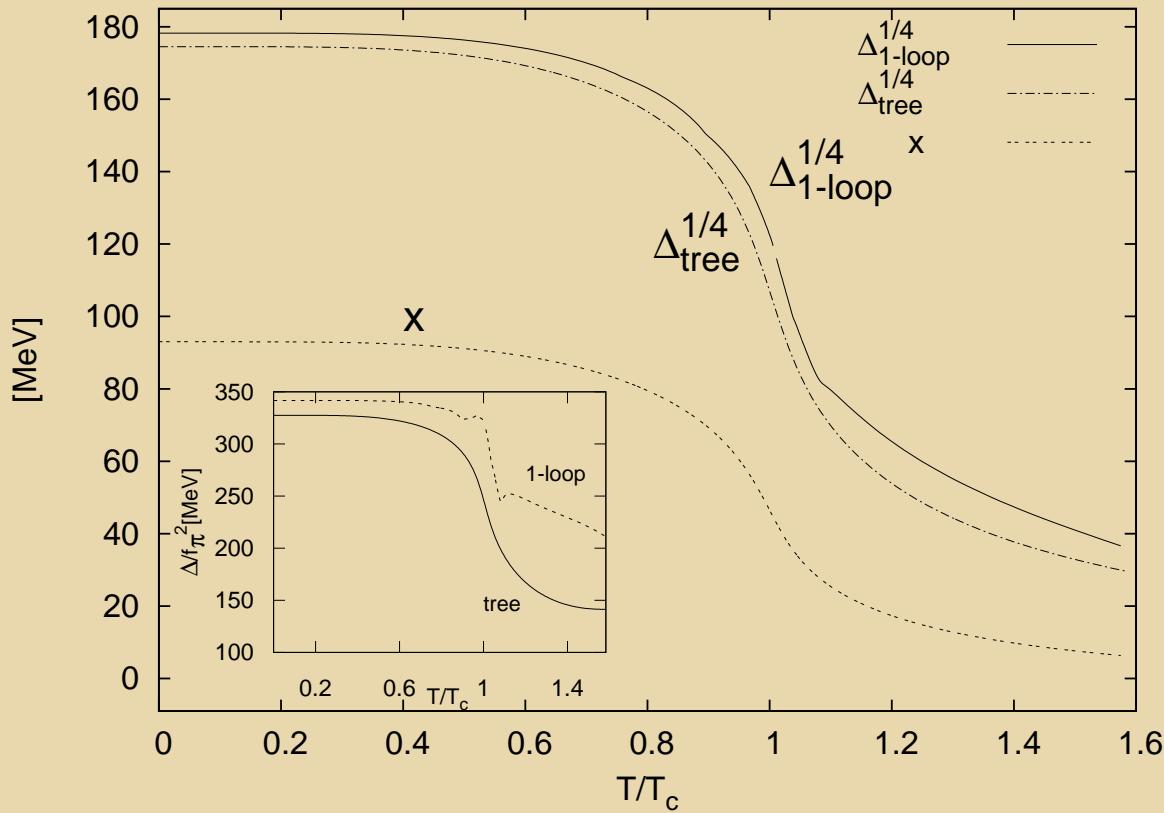
Increasing of either μ_B or μ_I \rightarrow influence of v_3 becomes stronger
 CEP at $\mu_I = 0$: $T_{\text{CEP}} = 63.08$ MeV, $\mu_{B,\text{CEP}} = 960.8$ MeV \rightarrow large diff. to case $v_3 = 0$

Reason: x and v_3 related \rightarrow common transition point

Tree-level and 1-loop pole masses



Estimation of the topological susceptibility



$$\Delta = \frac{1}{6}(m_\eta^2 + m_{\eta'}^2 - 2m_K^2)f_\pi^2$$

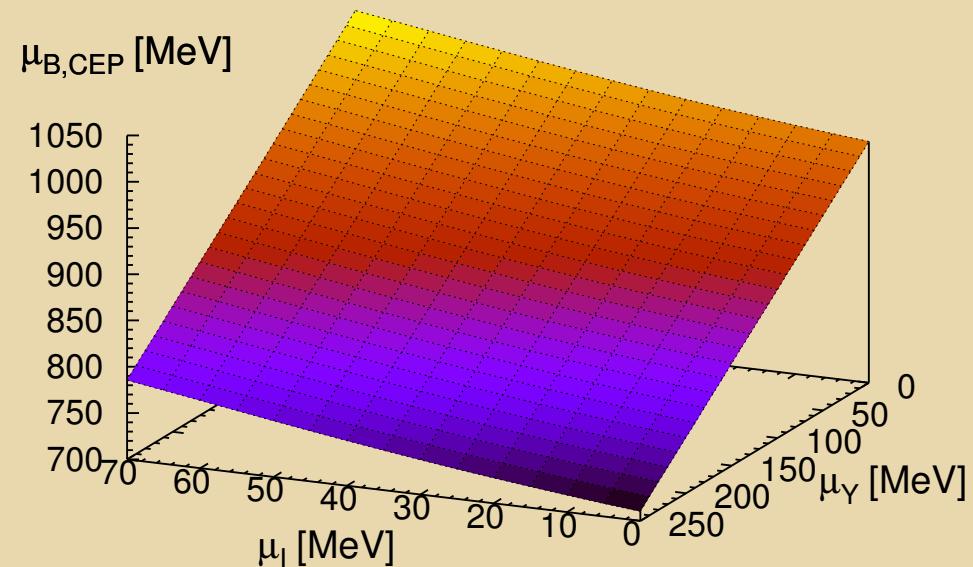
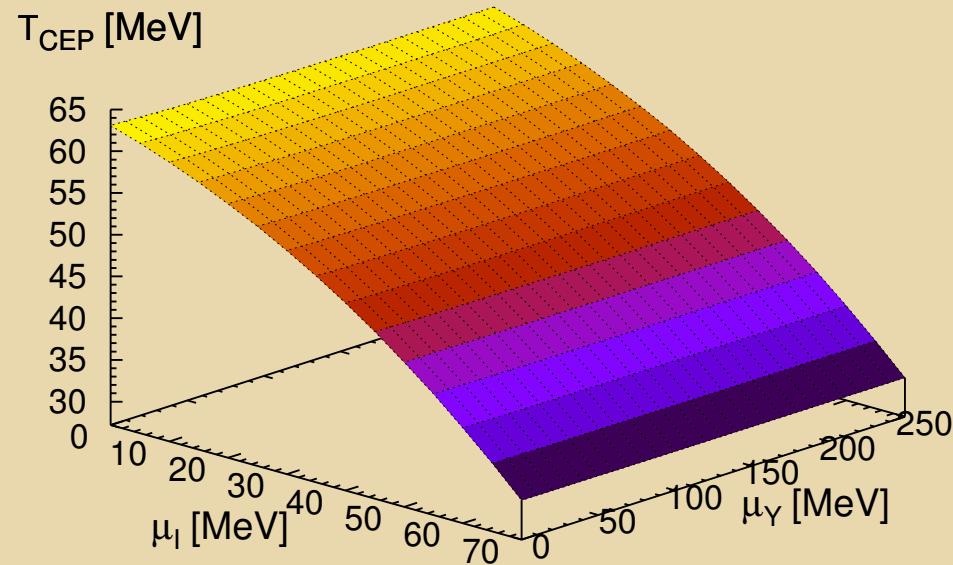
This is an estimation of $\chi_T(T)$ the topological susceptibility trough Witten-Veneziano formula:

$$\frac{2N_f}{f_\pi^2}\chi_T = m_\eta^2 + m_{\eta'}^2 - 2m_K^2$$

→ Connection with the $U_A(1)$ anomaly

Doesn't mean the restoration of $U_A(1) \rightarrow \chi_T(T)$ is dominated by chiral restoration

Dependence of the CEP on μ_I, μ_Y



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

$\mu_{B,\text{CEP}}$ has an almost linear dependence on both chemical potentials μ_I, μ_Y

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

Simple explanation of the influence of μ_I, μ_Y

Approximation: Ideal quantum gas of the quasi-particle degrees of freedom

In the first order region → generalized Clausius-Clapeyron equations

Partition function:

$$\ln Z = V \sum_i \gamma_i (2s_i + 1) \int \frac{d^3 p}{(2\pi)^3} \left[\beta \omega_i + \ln(1 + \alpha_i e^{-\beta(\omega_i - \mu_i)}) + \ln(1 + \alpha_i e^{-\beta(\omega_i + \mu_i)}) \right]$$

Gibbs-Duhem relation: $dp = s dT + n_B d\mu_B + n_I d\mu_I + n_Y d\mu_Y$

In case of $dp|_{phase1} = dp|_{phase2}$:

$$\begin{aligned} \left. \frac{dT}{d\mu_B} \right|_{\mu_Y, \mu_I} &= -\frac{\Delta n_B}{\Delta s}, & \left. \frac{dT}{d\mu_Y} \right|_{\mu_B, \mu_I} &= -\frac{\Delta n_Y}{\Delta s}, & \left. \frac{dT}{d\mu_I} \right|_{\mu_B, \mu_Y} &= -\frac{\Delta n_I}{\Delta s}, \\ \left. \frac{d\mu_B}{d\mu_Y} \right|_{T, \mu_I} &= -\frac{\Delta n_Y}{\Delta n_B}, & \left. \frac{d\mu_B}{d\mu_I} \right|_{T, \mu_Y} &= -\frac{\Delta n_I}{\Delta n_B} \end{aligned}$$

Obtained relations:

$$\Delta n_B, \Delta n_Y, \Delta s > 0, \Delta n_I < 0$$

$$\Delta n_B \approx \Delta n_Y, \Delta s > \Delta n_B, \Delta s > |\Delta n_I|$$

Conclusions and outlook

- The 2nd 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 is robustly predicted and at the physical point of the mass-plane was located at: $T_{CEP} = 74.83 \text{ MeV}$ $\mu_{B,CEP} = 895.38 \text{ MeV}$.
- The dependence of the μ_B on the width of the susceptibility was investigated.
- μ_I dependence of different pole masses were obtained.
- The temperature dependence of the topological susceptibility was estimated.
- 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 $T = 0$).
- A simple ideal gas model was established, which could explain the shifts of the CEP due to μ_I, μ_Y
- **Possible continuation:** Pion condensate, inclusion of the Polyakov-loop