

QCD: finite baryon density ; "sign problem"

Recall the global symmetry of (continuum) QCD :

$$\underbrace{SU_L(N_f) \times SU_R(N_f)}_{\substack{\text{broken explicitly if } M \neq 0, \\ \text{spontaneously if } M \rightarrow 0.}} \times U_V(1) \times U_A(1)$$

↓

broken by anomaly.

Unbroken exact global symmetry!

The Noether theorem: existence of a global symmetry implies existence of a conserved current!

$$U_V(1): \Psi \rightarrow e^{i\alpha v} \Psi, \bar{\Psi} \rightarrow e^{-i\alpha v} \bar{\Psi}.$$

For  $\mathcal{L} = \bar{\Psi} [\gamma_\mu D_\mu + M] \Psi$  (i.e. continuum), the corresponding Noether current is

$$j^\mu \propto \frac{\delta \mathcal{L}_M}{\delta (\bar{\Psi} \Psi)} \frac{\delta \Psi}{\delta \alpha v} = \frac{\delta}{\delta (\bar{\Psi} \Psi)} [\bar{\Psi} i\gamma^\mu \partial_\mu \Psi]. i\Psi = -\bar{\Psi} \gamma^\mu \Psi$$

In Euclidean space:

$$j^\mu = \bar{\Psi} \gamma_\mu \Psi, \text{ where an overall sign has been chosen.}$$

The zeroth component, "charge density":

$$j_0 = \bar{u} \gamma_0 u + \bar{d} \gamma_0 d + \bar{s} \gamma_0 s$$

$\bar{u} \gamma_0 u$  is interpreted as quark number density (quarks minus antiquarks)

$$\frac{2}{3} e \bar{u} \gamma_0 u - 1/3 \text{ electric charge density.}$$

$j_0$  counts the total number of quarks. Since there are three quarks in a baryon, we define the baryon number  $B$  as

$$B = \frac{1}{3} \int d^3 \bar{x} j_0 = \frac{1}{3} \int d^3 \bar{x} \bar{\Psi} \gamma_0 \Psi.$$

The baryon number is exactly conserved within QCD.

(Hadrons: mesons  $\equiv \{B=0\}$  + baryons  $\equiv \{B \neq 0\}$ .)

Systems with a finite baryon number density are of some physical interest:

- inside heavy nuclei : nuclear matter, nuclear liquid
- inside neutron stars : baryon densities much higher still : protons & neutrons overlap, "quark matter"?

It would be important to know the properties ["equation of state";  $p(T, \mu)$ ] of such matter, since they enter the Tolman-Oppenheimer-Volkov equations determining the structure of these objects.

A system with a finite baryon number density is often easier to describe by carrying out a Legendre transform and going into a system with a finite chemical potential:

$$\frac{\partial F(B, T, V)}{\partial B} = \mu_B ; \quad \Omega(\mu_B, T, V) \equiv F - \mu_B B$$

Baryon number (density) can then be recovered by

$$\frac{\partial \Omega(\mu_B, T, V)}{\partial \mu_B} = -B ; \quad F(B, T, V) = \Omega + \mu_B B .$$

Stat mech.:

$$\Omega(\mu_B, T, V) = -V p(\mu_B, T) = -T \ln \text{Tr} \exp \left[ -\beta (\hat{H} - \mu_B \hat{B}) \right] .$$

$\uparrow$   
Hamiltonian       $\uparrow$   
baryon number

Since  $B$  is conserved,  $[\hat{H}, \hat{B}] = 0$ , we can consider  $\mu_B \hat{B}$  simply as a part of the Hamiltonian! Therefore also path integral concept be derived as usual. Denoting  $\mu = 3 \cdot \mu_B$ , and recalling  $\beta \rightarrow \frac{i}{\hbar} \int d\tau$ ,

$$\mathcal{Z} \rightarrow \overline{\Psi} [\gamma^\mu D_\mu - \mu \delta_0 + M] \Psi .$$

How can this system be discretised?

[P. Hasenfratz, F. Karsch, Phys. Lett. B 125 (1983) 308]

Recall (p. 71-72):

$$U_p = e^{i \text{ag}_0 A_p}, \quad A_p = A_p^a T^a$$

$$(p. 93) \Rightarrow S = \sum_{\bar{x}} \alpha^4 \left\{ \bar{\Psi} \left( M + \frac{4r}{a} \right) \Psi - \frac{1}{2a} \sum_p \left[ \bar{\Psi}(\bar{x} + a\hat{p}) U_p^+(\bar{x})(r + \gamma_p) \Psi(\bar{x}) + \bar{\Psi}(\bar{x}) U_p(\bar{x})(r - \gamma_p) \Psi(\bar{x} + a\hat{p}) \right] \right\}$$

In continuum limit:

\* terms with  $r$  are  $\Theta(a)$  (p. 91)

\* the rest (apart from the mass term):

$$- \frac{1}{2a} \sum_p \left[ \bar{\Psi}(\bar{x} + a\hat{p}) U_p^+(\bar{x}) \gamma_p \Psi(\bar{x}) - \bar{\Psi}(\bar{x}) U_p(\bar{x}) \gamma_p \Psi(\bar{x} + a\hat{p}) \right]$$

$$\approx - \frac{1}{2a} \sum_p \left[ \bar{\Psi}(\bar{x} + a\hat{p}) \gamma_p \Psi(\bar{x}) - \bar{\Psi}(\bar{x}) \gamma_p \Psi(\bar{x} + a\hat{p}) + \bar{\Psi}(\bar{x}) (-i \text{ag}_0 A_p) \gamma_p \Psi(\bar{x}) - \bar{\Psi}(\bar{x}) (i \text{ag}_0 A_p) \gamma_p \Psi(\bar{x}) \right]$$

$$\text{Summing over } \bar{x} \Rightarrow = \sum_p \bar{\Psi}(\bar{x}) \gamma_p \frac{\Psi(\bar{x} + a\hat{p}) - \Psi(\bar{x} - a\hat{p})}{2a} + \bar{\Psi}(\bar{x}) \gamma_p i \text{ag}_0 A_p \Psi(\bar{x})$$

$$\approx \bar{\Psi}(\bar{x}) \gamma_p [ \partial_p + i \text{ag}_0 A_p ] \Psi(\bar{x}).$$

Now we note that the addition of a chemical potential corresponds to

$$i \text{ag}_0 A_0 \rightarrow i \text{ag}_0 A_0 - \mu \cdot \mathbf{1}_{N_c \times N_c}$$

Thus, on the lattice,

$$U_0 \rightarrow e^{-a\mu} U_0$$

$$U_0^+ \rightarrow e^{+a\mu} U_0$$

## The "sign problem"

Let us look at the Dirac operator again.

For simplicity, we work in the continuum.

$$\not{D} + M - \mu \gamma_0 = \underbrace{\gamma_\mu [\not{p} + i g_0 A_\mu^a T^a]}_{\equiv D_0} + M \underbrace{- \mu \gamma_0}_{\equiv D_\mu}; \text{ indices suppressed; } \gamma^\mu = \gamma_\mu.$$

Let us assume  $M$  to be degenerate for simplicity,  $M = \text{diag}(m, m, m)$ .

Then it is enough to consider one flavour only.

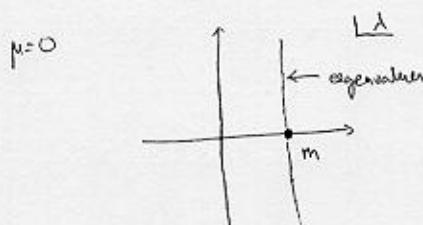
What can we say about the eigenvalues?

$$(1) \quad (D_0 + D_\mu)v = \lambda v \Rightarrow (D_0 + m + D_\mu)v = (\lambda + m)v \Rightarrow m \text{ plays a trivial role.}$$

$$(2) \quad (D_0 + D_\mu)v = \lambda v \Rightarrow (D_0 + D_\mu)\gamma_5 v = -\gamma_5(D_0 + D_\mu)v \\ = -\lambda \gamma_5 v \Rightarrow \text{eigenvalues come in pairs, } \lambda \& -\lambda, \text{ for } m=0.$$

$$(3) \quad D_0^+ = -D_0 \Rightarrow \lambda^* = (v^+ D_0 v)^+ = v^+ D_0^+ v \\ = -v^+ D_0 v = -\lambda$$

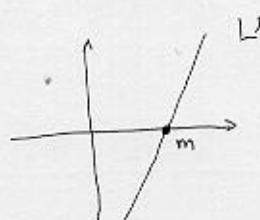
$\Rightarrow$  for  $\mu=0, m=0$ , they are purely imaginary!



$$\Rightarrow \det[D_0 + m] = (m + \lambda_1)(m + \lambda_1^*)(m + \lambda_2)(m + \lambda_2^*) \dots \in \mathbb{R}, \underline{> 0}.$$

$$(4) \quad D_\mu^+ = D_\mu$$

$\Rightarrow$  for  $\mu \neq 0$ , eigenvalues no longer purely imaginary for  $m=0$ .



$$\Rightarrow \det[D_0 + m + D_\mu] \in \mathbb{C}$$

$\Rightarrow$  weight  $\det[D_0 + m + D_\mu] \exp(-S^{(\text{gaugy})}) \in \mathbb{C}$ ; can be positive or negative or oscillatory

Summary

- For applications in astrophysics, it would be interesting to compute the pressure of QCD as a function of the baryon chemical potential (even at  $T=0$ ).
- The baryon chemical potential can easily be included in the QCD action.
- The inclusion of the chemical potential makes, however, the weight of the functional integral in general complex. This turns out to be a major problem, since importance sampling ceases to work.  
⇒ some fresh ideas needed !