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ON THERMODYNAMIC PROPERTIES OF CHROMOPLASMA

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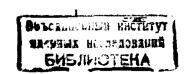
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At present it is well established that there is a transition to the deconfinement phase at high enough temperatures in gauge theories. This transition, predicted in refs. /1.2/, has then been confirmed by the numerical calculations in the framework of the lattice approach (see, for instance, refs. /3~9/). According to these calculations, at high enough temperatures chromoplasma behaves as a gas of almost free gluons and guarks. At the same time at temperatures just above the phase transition temperature  $heta_{ extstyle extsty$ the temperature region close to  $heta_c$  is of great interest for experimentalists as just this temperature interval is probable for obtaining chromoplasma in nucleus-nucleus collisions. At present, we have no unequivocal idea about the excitation modes of chromoplasma. It is not improbable (see review /10/ and refs. therein) that at temperatures close to  $heta_{\mathcal{L}}$  an important role is played by collective excitation modes of chromoplasma which give rise to electric and magnetic masses. These collective excitations can be both phonon /11/ and massive /12/. As a result, qualitative picture may be thought to be the following /12-14/: In the large momentum region the dominating part is played by almost noninteracting gluons whereas at momenta of an order of m, and m (the Debye and magnetic masses, respectively) the contribution comes from collective excitations both massive and massless. These collective excitations are assumed to be singlet with respect to colour (dynamical confinement hypothesis) and electric and magnetic massive modes qive similar contributions /14/.

The role and nature of collective excitation modes of chromoplasma can be elucidated by investigating thermodynamic functions within the lattice approach. The standard Wilson action is

$$S = \beta \sum_{PI_{\underline{a}}} (1 - \frac{1}{2} S_P U_{PI_{\underline{a}}}) + \beta \sum_{PI_{\underline{b}}} (1 - \frac{1}{2} S_P U_{PI_{\underline{b}}}), \qquad (1)$$

where  $U_{PI_{e}}$  and  $U_{PI_{t}}$  are plaquette variables corresponding to spacelike and timelike plaquettes, respectively, and  $\beta=4/g^2_{pare}$ 



The calculations were made on the lattice  $4*12^8$  with periodic boundary conditions in the range of  $\beta$  values :  $2.2 \le \beta \le 2.8$ .

The Monte-Carlo method was used to calculate the contribution of electric and magnetic modes to the internal energy E, pressure P and the so-called interaction measure  $\delta = E - 3P$  in the SU(2)-gauge theory. Temperature  $\theta$  was determined by the known two-loop expression

$$\theta/\Lambda_{L} = \frac{1}{N_{L}a} = \frac{1}{\Lambda_{L}}$$

$$= \frac{1}{N_{L}} \left( \frac{11}{6\pi^{2}\beta} \right)^{5i/42i} \exp\left\{ \frac{3\pi^{2}\beta}{11} \right\} (1 + \Phi(g^{2})), \qquad (2)$$

where a is the lattice spacing and  $N_{\rm L}\!=\!4\,$  is the number of sites along the time axis.

The energy density  $\mathcal{E}$ , pressure P and interaction measure  $\delta$  were determined by the expressions /5/:

$$E/\theta^{4} = 12 \, N_{L}^{4} \left( g^{-2} \left( \langle PI_{s} \rangle - \langle PI_{t} \rangle \right) + \frac{1}{2} \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{t} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{t} \rangle \right) \right) = \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{t} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{t} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{t} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{t} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{t} \rangle \right) - \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{t} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{t} \rangle \right) - \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{t} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle \right) = \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_{sym} \rangle - \langle PI_{sym} \rangle \right) + \frac{1}{2} \left( \langle PI_$$

where  $\langle Pl_s \rangle$  and  $\langle Pl_t \rangle$  are the spacelike and timelike plaquettes, respectively, and  $\langle Pl_{sym} \rangle$  is the plaquette average calculated on the hypercubic lattice. We have used values of coefficients  $C_s$  and  $C_t$  calculated within one-loop aproximation /15/:  $C_s=.1140$ ,  $C_t=-.06758$ . The electric and magnetic contributions to the thermodynamic functions are determined in a standard way through a lattice generalisation of the strength tensor  $F_{\mu\nu}$ . Correspondingly, in formulae (3a) ~ (3c) electric (magnetic) components of the thermodynamic functions are composed of spacelike (timelike) plaquettes and  $\langle Pl_{sym} \rangle$ . It is worth noting that all thermodynamic functions listed above are normalised in such a way that their values at zero temperature are equal to zero by definition.

Our values for  $\mathcal{E}(\theta)$ ,  $P(\theta)$  and  $\delta(\theta)$  and their electric and magnetic components are shown in figs.1-3<sup>1</sup>. The dashed lines in these figures denote the behaviour of the thermodynamic quantities  $\mathcal{E}/\theta^4$ ,  $P/\theta^4$  and  $\delta/\theta^4$  calculated by the perturbation theory within the one-loop approximation /16/. It is seen from these figures that at large enough temperatures the behaviour of energy  $\mathcal{E}$  and pressure P is in the agreement with the Stefan-Boltzman law (taking into account one-loop corrections). At temperatures close to the critical one in the deconfinement phase the values of  $P/\theta^4$  are strongly suppressed in comparison with the values of  $\mathcal{E}/\theta^4$ . This is quite natural since for the second order phase transition in the vicinity of the critical temperature the thermodynamic functions should have the following behaviour:

 $E \simeq {\rm const} \ | \ \theta - \theta_c \ |^{4-\alpha}$ ;  $P \simeq {\rm const} \ | \ \theta - \theta_c \ |^{2-\alpha}$ , where  $\alpha$  is the relevant critical exponent for specific heat (the notion of universality classes /17,18/ suggested us that  $\alpha \simeq 0.12$  as in the 3D-1sing model).

It is possible to reach some understanding of the behaviour of the chromoelectric part of energy using the Hamiltonian approach to SU(2)-gluodynamics in the strong coupling limit. In this regime the partition function is of the form

 $<sup>^4 \</sup>rm We$  are indebted to prof. H. Satz for providing us with the values of  $\langle PI_{\rm SVm} \rangle$  on the lattice 12.

$$Z = \int \prod_{i} d\gamma_{i} \prod_{l \mid r \mid k \in D} D(\gamma_{k}, \gamma_{j})$$
 (4a)

$$D(\gamma_{i},\gamma_{j}) = 1 + \sum_{k} \exp(-\frac{g^{2}}{2 \cdot a \cdot \theta} \cdot k(k+1)) \cdot X_{k}(\gamma_{i}) \cdot X_{k}(\gamma_{j}), \qquad (4b)$$

where  $\gamma_i$  is the SU(2) group determined in site i,  $\theta$  is the temperature,  $a=a_n$  is the lattice spacing in spatial directions and  $X_k(\gamma)$  is the character for kth representation of SU(2) group ( k=1/2, 1, 3/2...). At temperatures close to the critical one the expression for the partition function can be exposed in the form /1,2/

$$Z = \int \prod_{i} d\gamma_{i} \exp(-S_{eff})$$
 (5)

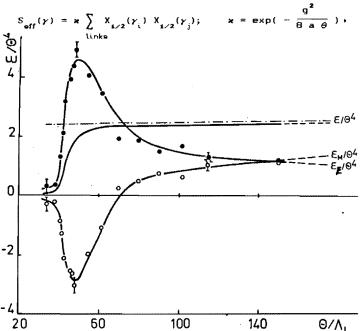


Fig.1. The solid line without points represents qualitatively—the behaviour of  $E/\theta^4$ , full circles and open circles correspond to  $E_{\underline{z}}/\theta^4$  and  $E_{\underline{z}}/\theta^4$ , respectively. The dash-dotted line denotes the perturbative contribution. Some points represent a typical statistical error.

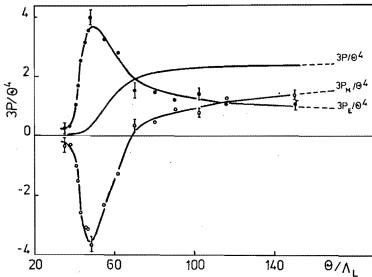


Fig.2. The solid line without points represents qualitatively—the behaviour of  $3P/\theta^4$ ,—full circles—and open circles—correspond to  $3P_{\rm g}/\theta^4$  and  $3P_{\rm m}/\theta^4$ , respectively. Some points represent—a typical statistical error.

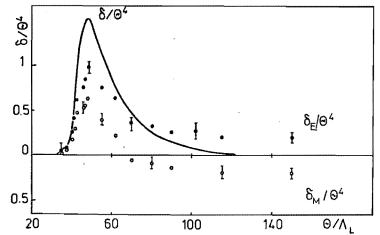


Fig.3. The quantities  $\delta/\theta^4$ ,  $\delta_{\rm E}/\theta^4$  and  $\delta_{\rm M}/\theta^4$  with the same notation as in figs.1,2.

i.e., we have developed the three-dimensional "spin" model of the Ising type , with the difference that the integration in (5) is performed over the Haar measures  $\mathrm{d}\gamma_1$ . The role of spins is played by representations  $\frac{1}{2}\,\mathrm{X}_{1/2}(\gamma_1)=s_1$ . (A consistent procedure of reducing (4) to partition function of the "spin" system is demonstrated in ref./19/). A model of this type is known to predict a second order phase transition. For a qualitative investigation of the behaviour of thermodynamic functions we use the mean-field approximation. Using the standard technique of the mean-field approximation (see, for instance, ref./20/) we get the equation for the mean "magnetization" M = <  $s_i$  > :

 $M=1_2(24\times M)/1_1(24\times M), \qquad (6)$  where  $1_2$  and  $I_1$  are the modified Bessel functions. This equation at  $\varkappa \leq \varkappa_c = 1/6$  has only zero solutions for mean magnetization M whereas at  $\varkappa \geq \varkappa_c$  the magnetization changes continuously from zero at  $\varkappa = \varkappa_c$  (i.e. at

 $x = x_{C}$  the second order phase transition occurs.

The effective action  $W(M) = -\ln Z$  in this approximation is

$$\frac{1}{N_{g}} W(M) = 12 \times M^{2} - \ln \left[ I_{1}(24 \times M)/12 \times M \right], \tag{7}$$

where  $N_s$  is the number of sites in the three-dimensional space. Using (7) we get for the internal energy

$$E/\theta^4 = \frac{1}{V\theta^2} \frac{\partial W}{\partial \theta} = \frac{1}{(a\theta)^4} \frac{9g^2}{2} \times M^2,$$
 (B)

where  $V = N_a^a$  is the three-dimensional volume.

With relations (6)-(8) one can easily obtain the behaviour for the internal energy which coincides qualitatively with the one depicted in fig.1 for  $E_{\rm g}/\theta^4$ . This fact is hardly surprising in the light of the universality hypothesis mentioned above /17,18/. (Similar calculations for P and  $\delta$  are more ambiguous because of the presence of derivative  ${\rm dg}^{-2}/{\rm da}$  which is unknown in the strong-coupling limit).

The problem of interpretation like this for the chromomagnetic contributions remains open. At temperatures just above the critical one the magnetic components  $E_{\mathbf{M}}$  and  $P_{\mathbf{M}}$  (both of them

being the difference between the corresponding nonzero temperature and zero temperature values) take negative values (fig.1,2)  $/21/^2$ . This means that the contribution of the chromomagnetic excitation modes differs drastically from that of the chromoelectric excitation modes (at least for SU(2) group), and extra excitation mechanisms of chromoplasma are to be searched for. We assume that the temperature dependence of magnetic components of energy and pressure presented here is maybe an indication of an important role played by unstable modes /23,24/ in the QCD vacuum.

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<sup>&</sup>lt;sup>2</sup>It is worth noting that nonperturbative corrections /22/ changing appreciably the values of coefficients  $C_{\rm g}$  and  $C_{\rm t}$  outside the region of asymptotic scaling do not influence this conclusion.

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О термодинамических свойствах хромоплазмы

Вычислялись вклады электрических и магнитных мод во внутреннюю энергию в SU(2)-калибровочной теории на решетке  $4\cdot 12^3$ . Показано, что поведение хромоэлектрической части энергии  $E_{\rm E}$  вполне объясняется в рамках модели изинговского типа, в полном соответствии с гипотезой универсальности. В то же время, поведение магнитной части внутренней энергии  $E_{\rm M}$  кардинально отличается от поведения  $E_{\rm E}$ . Характер температурной зависимости  $E_{\rm M}$  указывает на существенно несимметричный характер электрических и магнитных мод возбуждения хромоплазмы, и, быть может, проливает свет на роль нестабильных мод в калибровочных теориях.

Работа выполнена в Лаборатории теоретической физики ОИЯИ.

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On Thermodynamic Properties of Chromoplasma

The contributions of electric and magnetic modes to some thermodynamic functions in SU(2)-gauge theory on the lattice  $4\cdot 12^3$  were calculated. It is shown that the behaviour of the chromoelectric part of energy  $E_{\rm E}$  can be interpreted within the Ising-type model in agreement with the universality hypothesis. At the same time the behaviour of the magnetic parts of the internal energy and pressure (i.e.  $E_{\rm M}$  and  $P_{\rm M}$ ) differs drastically from that of  $E_{\rm E}$  and  $P_{\rm E}$ . The character of the temperature dependence of  $E_{\rm M}$  and  $P_{\rm M}$  exhibited here testifies to the presence of a highly nonidentical properties of electric and magnetic modes of chromoplasma and maybe sheds light on the role of unstable modes in gauge theories.

The investigation has been performed at the Laboratory of Theroetical Physics, JINR.