

A REAL-TIME LATTICE SIMULATION OF THE THERMALIZATION OF A GLUON PLASMA: FIRST RESULTS*

MAXIMILIAN ATTEMS

Departament de Física Quàntica i Astrofísica
and

Institut de Ciències del Cosmos (ICC)

Universitat de Barcelona, Martí i Franquès 1, 08028 Barcelona, Spain

OWE PHILIPSEN, CHRISTIAN SCHÄFER, BJÖRN WAGENBACH
SAVVAS ZAFEIROPOULOS

Institut für Theoretische Physik, Johann Wolfgang Goethe-Universität
Max-von-Laue-Str. 1, 60438 Frankfurt am Main, Germany

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To achieve an understanding of the thermalization of a quark–gluon plasma, starting from QCD without using model assumptions, is a formidable task. We study the early stage dynamics of a relativistic heavy-ion collision in the framework of real-time simulations of the classical Yang–Mills theory in a static box with the color glass condensate as an initial condition. Our study generalizes a previous one by Fukushima and Gelis from SU(2) to the realistic case of SU(3). We calculate the chromo-electric and chromo-magnetic energy densities as well as the ratio of longitudinal and transverse pressure as a function of time as probes for thermalization. Our preliminary results on coarse lattices show the occurrence of Weibel instabilities prior to thermalization.

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1. Introduction

Experimental results from heavy-ion collisions at the Large Hadron Collider (LHC) and the Relativistic Heavy Ion Collider (RHIC) give strong indications for the occurrence of a quark–gluon plasma (QGP). The successful description of the QGP, even at early times, by relativistic viscous

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hydrodynamics suggests a rather fast thermalization within a few fm [1, 2]. Starting from the highly anisotropic scenario of two colliding ultrarelativistic nuclei, it is still an open question which processes are responsible for this fast thermalization. One candidate are color instabilities [3, 4] such as the chromo-Weibel instability [5–8], which is the analogue of the Weibel instability in electromagnetic plasmas [9].

In the limit of weak coupling, which is assumed for the present energy scales at the LHC and RHIC, the two nuclei can be described in the Color Glass Condensate (CGC) framework [10–12]. It provides an expansion of inclusive quantities, such as the expectation value of the energy-momentum tensor $T^{\mu\nu}$, in powers of the strong coupling α_s , where the leading order is obtained by solving the Yang–Mills equations. We want to approach this problem from first principles by applying real-time lattice QCD techniques, approximating QCD by Yang–Mills theory in the classical limit. The classical approximation is justified by the high occupation of gluons in the non-equilibrium medium produced after the heavy-ion collision. In contrast to other works using Bjorken (expanding) coordinates [13], our exploratory studies focus on finite, static box simulations.

2. Color glass condensate

The classical Yang–Mills equations for the gauge fields and their conjugate momenta can be solved using the framework of lattice gauge theory [14], where the continuum gauge fields $A_\mu(x)$ are represented in terms of lattice gauge fields $U_\mu(x) = \exp(igaA_\mu(x))$. We follow the McLerran–Venugopalan model [15–17] to construct the initial conditions for the lattice gauge fields U and the chromo-electric fields E within the CGC framework.

2.1. Lattice gauge fields

We set the longitudinal component of the collective gauge field $U_z = 1$ and construct the transverse components by merging the color fields from the two colliding nuclei, denoted by $U^{(1)}$ and $U^{(2)}$, by solving [18]

$$\text{Tr} \left\{ T^a \left[\left(U^{(1)} + U^{(2)} \right) \left(1 + U^\dagger \right) - \text{h.c.} \right] \right\} = 0 \quad (2.1)$$

for given $U^{(1)}$ and $U^{(2)}$, with T^a — the generators of SU(3).

To get the color fields $U^{(1)}$ and $U^{(2)}$, one solves the Poisson equation $\Delta_L A_m^a(x) = -\rho_m^a(x)$ for nuclear color sources with the following Gaussian distribution

$$\left\langle \rho_m^a(x) \rho_n^b(y) \right\rangle = g^4 \mu^2 a^2 \delta_{mn} \delta^{ab} \delta(x - y). \quad (2.2)$$

The color fields are then constructed from the obtained “color potentials” $\Lambda_m = T^a \Lambda_m^a$ via $U_i^m(x) = \exp(i\Lambda_m(x)) \exp(i\Lambda_m(x + \hat{i}))$, with $m = 1, 2$ and $i = x, y$.

2.2. Chromo-electric fields

The transverse chromo-electric fields are set to zero and the longitudinal part is constructed out of the gauge fields U via

$$E_z^a(x) = -\frac{i}{2g} \text{Tr} \left(T^a \sum_{i=x,y} \left\{ [U_i(x) - 1] [U_i^{(2)\dagger}(x) - U_i^{(1)\dagger}(x)] - \text{h.c.} \right. \right. \\ \left. \left. + [U_i^\dagger(x - \hat{i}) - 1] [U_i^{(2)\dagger}(x - \hat{i}) - U_i^{(1)\dagger}(x - \hat{i})] - \text{h.c.} \right\} \right). \quad (2.3)$$

2.3. Fluctuations

In order to reach isotropization, fluctuations need to be added to the initial fields [19,20]. This is done by a small random variation of the idealized δ -function-like distribution of the initial chromo-electric fields. To this end, we add a fluctuation term $\delta E \sim \Delta$. We will discuss the effect of the seed Δ in Sec. 4.

3. Observables

To study the thermalization time, we observe the ratio of the longitudinal to the transverse pressure P_L/P_T [21]. The pressure components are related to the spatial diagonal elements of the energy-momentum tensor T^{ii} [20] and the chromo-electric and chromo-magnetic longitudinal and transverse energy densities as:

$$P_L(t) = -T_z^z(t) = \epsilon_{E_T}(t) + \epsilon_{B_T}(t) - \epsilon_{E_L}(t) - \epsilon_{B_L}(t), \quad (3.1)$$

$$P_T(t) = -\frac{1}{2} [T_x^x(t) + T_y^y(t)] = \epsilon_{E_L}(t) + \epsilon_{B_L}(t). \quad (3.2)$$

The lattice expressions for the components of the chromo-electric and chromo-magnetic energy densities $H_i^{E/B}$, with $i = x, y, z$, are:

$$H_i^B(t, x) = \frac{2N_c}{g^2} \sum_{\substack{j < k \\ j, k \neq i}} \left[1 - \frac{1}{N_c} \text{Re Tr } U_{jk}(x) \right], \quad (3.3)$$

$$H_i^E(t, x) = \frac{1}{g^2} \text{Re Tr } [E_i(x)E_i(x)] \quad (3.4)$$

from which we get the energy densities $\epsilon_{B/E_{L/T}}$ by averaging over the transverse lattice plane (x, y) or the longitudinal axis (z) , respectively.

4. Numerical results

Our results are based on calculations on a 40^3 -lattice averaged over 30 independent measurements. We set the coupling constant $g = 2$ and fix $g^2\mu L = 120$ in accordance with the literature and conforming to the expectations of RHIC physics [13].

The value for the seed Δ , which controls the fluctuations, is purely artificial. The higher it is, the faster the system thermalizes, but there is an upper bound: Since Δ only enters the electric part of the initial energy density, we have to monitor its influence on the initial total energy density, which should not significantly change, otherwise the classical approximation is no longer tenable. In our setup, we chose $\Delta = 0.025$, which leads to a change of the initial value of the total energy density, which stays constant during the simulation, by less than 10% (*cf.* figure 1 (left)).

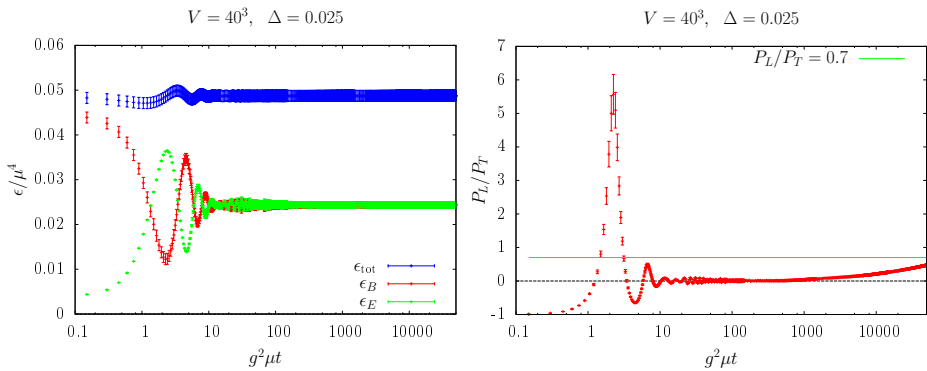


Fig. 1. (Left) Magnetic energy density ϵ_B , electric energy density ϵ_E and total energy density ϵ_{tot} . (Right) Pressure ratio P_L/P_T and hydrodynamization line at $P_L/P_T = 0.7$.

Evidence for thermalization is provided by the pressure ratio P_L/P_T approaching 1. We are especially interested in the so-called hydrodynamization time, which is the time where hydrodynamic models should start to be applicable. Figure 1 (right) shows the value of the pressure ratio, which slowly approaches the hydrodynamization level, which is commonly set to $P_L/P_T \approx 0.7$.

Figure 2 shows the energy density H_x^E (*cf.* (3.4)) in the (y, z) -plane, where we averaged over the x -direction. Since we are only interested in the qualitative structure, we performed only one measurement here. This explains the vertical lines, which are a result of the initial fluctuations.

At $g^2\mu t = 1666$, two filaments start emerging, which are clearly visible at $g^2\mu t = 3333$. Later, at $g^2\mu t = 23333$, the filamentous structure is almost completely dissolved, which suggests an imminent thermalization. This characteristic filamentation is a strong indication for the presence of a chromo-Weibel instability.

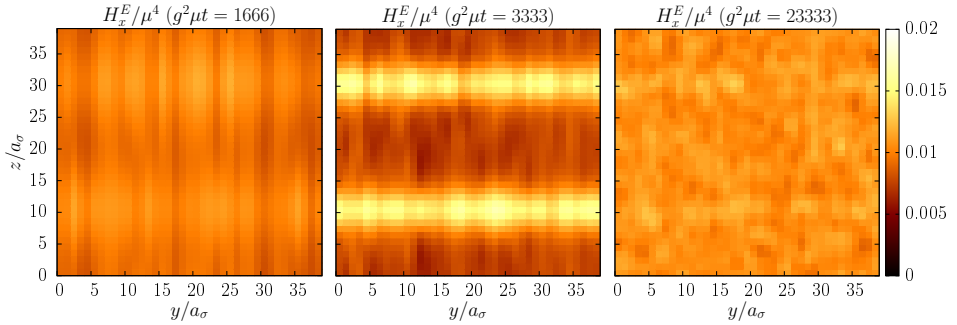


Fig. 2. The energy density H_x^E averaged over the x -direction is plotted in the (y, z) -plane for different times. One can easily see the arising filamentation.

5. Conclusion

We have provided evidence of the thermalization (hydrodynamization) of a color glass condensate such as the one created at relativistic heavy-ion collisions within the classical SU(3) Yang–Mills theory. To this end, we studied observables such as the energy density and its components, and the longitudinal to transverse pressure ratio. By looking at components in a transverse–longitudinal plane, we were able to visualize the occurrence of chromo-Weibel instabilities represented by a clear filamentation of the energy density. The next steps will be an analysis of finite volume and discretization effects as well as studies of an expanding system.

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