

Initial conditions for heavy ion collisions with QCD kinetic theory

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We employ the leading order QCD kinetic theory to describe a consistent matching between the initial stage of a heavy ion collision and the subsequent hydrodynamic evolution. We use the linearized kinetic response functions around the non-equilibrium longitudinally expanding background to map initial energy and momentum perturbations to the energy momentum tensor at hydrodynamic initialization time τ_{hydro} . We check that hadronic observables then become rather insensitive to the cross-over time between kinetic theory and viscous hydrodynamics. The universal scaling of kinetic response in units of kinetic relaxation time $\tau_R \sim \eta/(sT)$ allows for a straightforward application of kinetic pre-equilibration event-by-event.

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1. Introduction

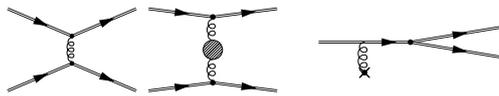
The heavy ion collision is a multi-phase process spanning a wide range of energy scales. Many of experimental signatures are rather successfully reproduced by modern heavy ion models, which describe each stage and energy regime in a suitable theoretical approximation. In particular, the low energy expansion of Quark Gluon Plasma (QGP) is well described by relativistic viscous hydrodynamics [1, 2, 3]. The initial conditions for the hydrodynamic expansion are supplied by initial state models at hydrodynamic initialization time τ_{hydro} . However most of initial state models do not contain the necessary physics to describe the approach to hydrodynamic behaviour and, therefore, create a mismatch of descriptions at the crossover time, which results in the dependence of physical observables on the auxiliary model parameter τ_{hydro} [6]. A consistent theoretical overlap of initial stages and hydrodynamics could be important not only for low energy medium physics, like a build up of radial flow, but also for the high- p_T physics of QGP, e.g. jets quenching and photon emission [7, 8].

In the high energy limit, the equilibration of weakly coupled QCD matter can be described by the leading order QCD kinetic theory [9, 10, 11]. We develop a framework of non-equilibrium kinetic response functions to smoothly map the initial conditions of far-from-equilibrium initial state models to hydrodynamically expanding matter and thus reduce the sensitivity to the hydrodynamic initialization time τ_{hydro} [12, 13].

2. Kinetic theory and universal scaling

We use the leading order kinetic theory known as AMY formalism [9]. The QGP is then modelled by a collection of massless high energy quasi-particles (quarks and gluons) in a bath of soft momentum excitations. At leading order in the coupling constant $\lambda = N_c g^2$ the quasi-particles experience elastic $2 \leftrightarrow 2$ scatterings and inelastic $1 \leftrightarrow 2$ medium induced splittings. In a boost invariant expansion the QGP equilibrates according to the so called "bottom-up" thermalization scenario, which is supported by the first principles classical-statistical Yang-Mills simulations [10, 14, 15]. We would like to emphasize that the same kinetic framework can be also used to describe the energy loss of high- p_T partons, and the electromagnetic emission in QGP [7, 8].

In this work we solve the Boltzman equation for the color averaged gluon distribution function $f(\tau, \mathbf{p}, \mathbf{x})$

$$\partial_\tau f + \frac{\mathbf{p}}{|p|} \cdot \nabla f - \frac{p_z}{\tau} \partial_{p_z} f = - \underbrace{\mathcal{C}_{2 \leftrightarrow 2}[f]}_{\text{diagram}} - \underbrace{\mathcal{C}_{1 \leftrightarrow 2}[f]}_{\text{diagram}}, \quad (2.1)$$


but neglect the effect of quark degrees of freedom¹. The distribution function $f(\tau, \mathbf{p}, \mathbf{x})$ is decomposed into a spatially homogeneous, but anisotropic $\langle p_\perp^2 \rangle \gg \langle p_z^2 \rangle$ background $\bar{f}(\tau, \mathbf{p})$ and linearized fixed wavenumber \mathbf{k} perturbations $\delta f(\tau, \mathbf{p}) e^{i\mathbf{k} \cdot \mathbf{x}}$ in the transverse plane of the collision. The resulting system of coupled equations is then solved [12].

¹At early times the gluon occupation far exceeds that of quarks due to Pauli exclusion principle.

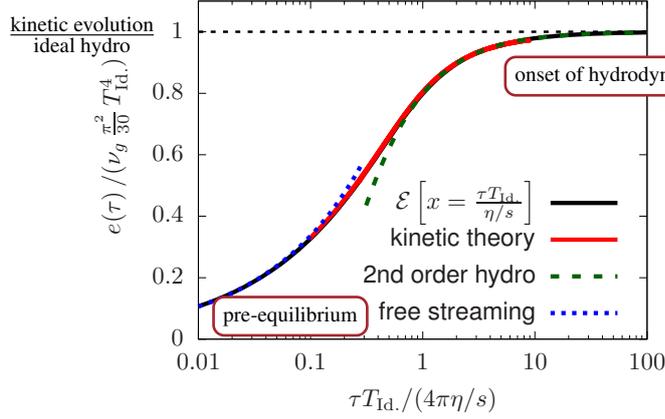


Figure 1: The universal scaling curve, Eq. (2.3), for energy density. Note that the kinetic theory curve corresponds to an overlapping kinetic evolution for a range of couplings $\lambda = 10 - 25$, which correspond to $\eta/s = 0.62 - 0.16$.

The key observation is that, as the system approaches equilibrium, many of the details of the initial distribution function is lost and only the first moments of the distribution function, i.e. energy momentum tensor, carries the necessary information for the subsequent hydrodynamic evolution. Furthermore at late times, the approach to hydrodynamics is approximately universal if measured in units of kinetic relaxation time $\tau_R \sim \eta/(sT)$. For example the energy evolution of a uniform Bjorken expanding background can be described by an asymptotic expansion in second order hydrodynamics [16]

$$\text{scaling in hydro: } e(\tau) = \underbrace{v_g \frac{\pi^2}{30} T_{\text{id.}}^4}_{\text{"ideal" temp.}} \left(\underbrace{1}_{\text{ideal}} - \underbrace{\frac{8}{3} \frac{\eta/s}{\tau T_{\text{id.}}}}_{\text{viscous}} + \underbrace{C_2 \left(\frac{\eta/s}{\tau T_{\text{id.}}} \right)^2}_{\text{2nd order hydro}} \right) \quad (2.2)$$

where $T_{\text{id.}}$ characterises the asymptotic temperature scale and C_2 is a second order transport coefficient with only weak dependence on the coupling constant λ . In practice, we find that for a range of moderate values of λ , the entire kinetic pre-equilibrium exhibits scaling and we define a scaling function $\mathcal{E}(x)$

$$\text{generalized scaling: } e(\tau) = v_g \frac{\pi^2}{30} T_{\text{id.}}^4(\tau) \times \underbrace{\mathcal{E} \left[x = \frac{\tau T_{\text{id.}}(\tau)}{\eta/s} \right]}_{\text{non-equilibrium evolution}}, \quad (2.3)$$

which smoothly interpolates between the early time free-streaming like behavior and the late time hydrodynamics, as shown in Fig. 1. For homogeneous boost invariant background, the energy density function $e(\tau)$ completely determines the entire background energy momentum tensor $\bar{T}^{\mu\nu} = \text{diag}(e, P_T, P_T, \frac{1}{\tau^2} P_L)$, where $P_L = -\partial_\tau(\tau e(\tau))$ and $e + 2P_T + P_L = 0$.

3. Linear response functions

For central heavy ion collisions, e.g. PbPb, the physical description of initial stages $\tau \lesssim 2$ fm can be considerably simplified by using the approximate boost invariance symmetry and the fact

that the transverse size of the collision is much larger than the duration of the pre-equilibrium stage, i.e. $R_{\text{Pb}} \gg \Delta\tau$. Then the energy momentum tensor at a given point in the transverse plane \mathbf{x}_0 is completely determined by a small neighbourhood of points which lie in the causal past of the said point, i.e. $|\mathbf{x} - \mathbf{x}_0| < \Delta\tau$. Furthermore, the energy momentum tensor $T^{\mu\nu}$ in the causal neighbourhood can be decomposed into a homogeneous background $\bar{T}^{\mu\nu}$ and perturbations $\delta T^{\mu\nu}$. At linear order the evolution of such perturbations is determined by linear response functions, which, in Fourier space, is a linear map between initial energy and momentum perturbations and the final time energy momentum tensor²

$$\underbrace{\frac{\delta T^{\mu\nu}(\tau, \mathbf{k})}{e(\tau)}}_{\text{output for hydro}} = \underbrace{\tilde{G}_{(\tau,i)}^{\mu\nu}(\mathbf{k}, \tau, \tau_0)}_{\text{non-equilibrium kinetic response}} \times \underbrace{\left[\frac{\delta T^{\tau\tau}(\tau_0, \mathbf{k})}{e(\tau_0)} \text{ or } \frac{\delta T^{\tau i}(\tau_0, \mathbf{k})}{e(\tau_0)} \right]}_{\text{initial perturbations}} \quad (3.1)$$

Such \mathbf{k} -space Green functions are obtained from the moments of the distribution function of the linearised perturbations $\delta f(\tau, \mathbf{p})e^{i\mathbf{k}\mathbf{x}}$, which are evolved according to the coupled system of equations discussed in Sec. 2. Notably, the linear response functions become approximately invariant if written in terms of scaling variables $\tau T_{\text{Id.}}/(\eta/s)$ and $|\mathbf{k}|(\tau - \tau_0)$. This allows performing a computational intensive kinetic theory evolution for a particular value of λ once, and then applying the scalable response functions to different values of effective η/s , time τ or temperature scale $T_{\text{Id.}}$.

4. Results

We apply the kinetic pre-equilibration stage to a central PbPb event with a centre-of-mass energy $\sqrt{s_{\text{NN}}} = 2.76$ TeV and with IP-Glasma initial conditions. IP-Glasma is a successful initial state model motivated by Color Glass Condensate picture and includes a 2D Yang-Mills evolution of initial gluonic fields [17, 18]. However, the classical Yang-Mills evolution does not lead to hydrodynamization, which is achieved by first passing the initial energy and momentum densities to a kinetic pre-equilibrium stage and then smoothly matching to viscous hydrodynamics at a later time τ_{hydro} . We check by varying the crossover time τ_{hydro} that the integrated and differential hydrodynamic fields, e.g. entropy density and radial velocity, are only weakly dependent on the cross-over time³. Finally, we extract the hadronic observables from the constant temperature freeze-out surface. For simplicity we only consider thermal spectra at the moment of freeze-out, i.e. neglecting resonance decays. In Fig. 2 we present the thermal pion multiplicity dN_{π}/dy , mean pion transverse momentum $\langle p_T^{\pi} \rangle$ and the elliptical flow v_2^{π} for our event. We observe that the dependence of hadronic observables on the duration of kinetic pre-equilibrium stage is weak, therefore the hydrodynamic simulations of heavy ion collisions with kinetic pre-equilibrium stage could perhaps yield more reliable extractions of transport properties of QGP.

In summary, we presented a well motivated pre-equilibrium model based on a QCD kinetic theory with a smooth and consistent approach to hydrodynamics. Thanks to an approximate scaling

²In general one can also consider response to the spatial perturbations of energy momentum tensor, but here we focus only on the perturbations sourced by initial fluctuations of conserved charges of energy and momentum.

³One of the remaining shortfalls of the matching is the difference between the conformal equation of state of massless kinetic theory and a non-conformal lattice equation of state used in modern hydrodynamic codes.

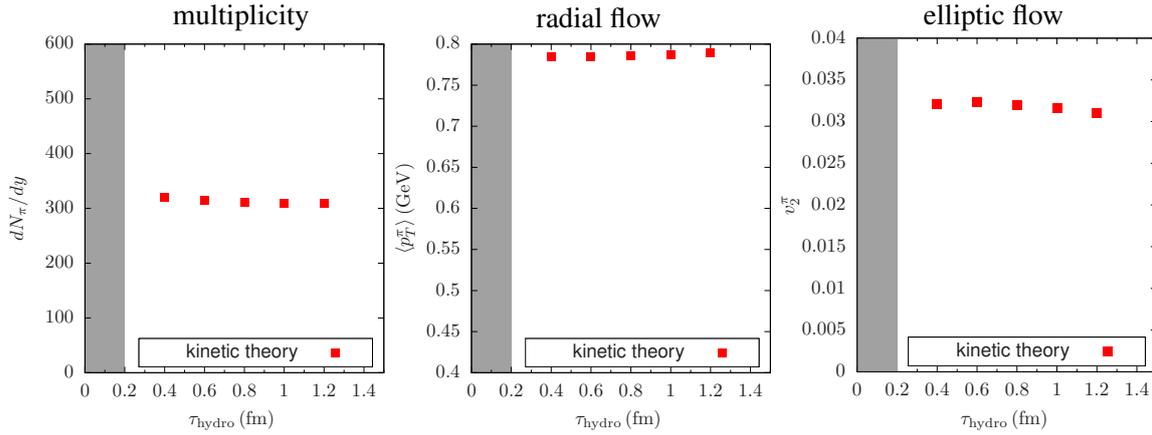


Figure 2: The final hadronic observables as a function of hydrodynamic initialization time τ_{hydro} with kinetic pre-equilibrium stage and IP-Glasma initial conditions for a $\sqrt{s} = 2.76$ TeV central PbPb event. Different panels correspond to thermal pion multiplicity dN_π/dy , the mean transverse momentum $\langle p_T^\pi \rangle$ and the elliptic flow v_2^π .

with $\eta/(sT)$, the tabulated linear response functions can be straightforwardly applied to event-by-event initial conditions of heavy ion collisions and thus reduce the dependence on the auxiliary parameters of a hydrodynamic model of heavy ion collisions.

Acknowledgments

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