Event-by-event pre-equilibrium dynamics — from gluon saturation towards the onset of hydrodynamics

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Based on A. Kurkela, A. Mazeliauskas, J.-F. Paquet, SS, D. Teaney (arXiv:1704.0524 (more in preparation)

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Microscopic early time dynamics & equilibration process

Macroscopic description of equilibration process

Event-by-event simulation of pre-equilibrium dynamics

Conclusions & Outlooks

Early time dynamics & equilibration process



Starting from saturated nuclei before the collisions sequence of processes that eventually leads to the formation of an equilibrated QGP

Early time dynamics ($0 < \tau < 1/Q_s$)

Because of high phase-space density of gluons particle initial particle production and early time dynamics described in terms of classical field theory to leading order $\nabla r^{\mu\nu} - r^{\nu}$

$$D_{\mu}F^{\mu\nu}=J^{\nu}$$

Strong boost invariant classical fields E^{η}, B^{η} created immediately after the collision

Decoherence of classical fields occurs on a time scale $\tau \sim 1/Q_s$ where quasi-particle description starts to become applicable

-> Basis for microscopic initial state calculations (IP-Glasma)

Challenge to understand subsequent equilibration process

- need quantum corrections (beyond NLO)
- effects of plasma instabilities





Equilibration process at weak coupling

Equilibration proceeds as three step process described by "bottom-up" scenario

Baier, Mueller, Schiff, Son PLB502 (2001) 51-58

Phase I: Evolution towards classical-attractor quasi-particle description becomes applicable

Phase II: Mini-jets undergo a radiative break-up cascade

formation of soft thermal bath

Phase III: Quenching of mini-jets in soft thermal bath

isotropization of plasma

Equilibration time determined by the time-scale for a mini-jet Q_s to loose all its energy

Berges,Boguslavski,SS, Venugopalan, PRD 89 (2014) no.7, 074011



Extrapolations from weak-coupling limit to realistic values of α_s (~0.3) at RHIC & LHC energies yield results consistent with phenomenological estimates

Viscous hydrodynamics applicable on time scales ~1 fm/c

Similar to strong coupling picture viscous hydrodynamics becomes applicable when pressure anisotropies are still O(1)



Kurkela, Zhu PRL 115 (2015) 182301

Early time dynamics & equilibration process



description of early-time dynamics can be achieved

Event-by-event pre-equilibrium dynamics

<u>Goal</u>: Obtain event-by-event initial conditions including weakly coupled pre-equilibrium evolution

-> Eliminate uncertainties in extraction of QGP transport properties due to artificial time scale τ_{Hydro} when hydro simulation starts

Challenge: Different degrees of freedom relevant at different times

classical fields, quasi-particles, energy-momentum tensor



Brute force calculation extremely challenging (CYM f(x,p), 3+2+1D EKT)

Ultimately we are only interested in calculation of energy-momentum tensor Exploit memory loss to use macroscopic degrees of freedom for description of pre-equilibrium dynamics

Macroscopic pre-equilibrium evolution

Extract energy-momentum tensor T^{µv}(x) from classical statistical lattice simulation

Evolve $T^{\mu\nu}$ from initial time $\tau_0 \sim 1/Q_s$ to hydro initialization time τ_{Hydro} using eff. kinetic theory description

Causality restricts contributions to $T^{\mu\nu}(x)$ to be localized from causal disc $|x-x_0| < \tau_{Hydro}-\tau_0$ useful to decompose into a local average $T^{\mu\nu}_{BG}(x)$ and fluctuations $\delta T^{\mu\nu}(x)$

Since in practice size of causal disc is small $\tau_{Hydro}-\tau_0 << R_A$ fluctuations $\delta T^{\mu\nu}(x)$ around local average $T^{\mu\nu}_{BG}(x)$ are small and can be treated in a linearized fashion



class.Yang-Mills

Effective kinetic description needs phase-space distribution $f(\tau, p, x)$ Memory loss: Details of initial phase-space distribution become irrelevant as system approaches local equilibrium

Can describe evolution of $T^{\mu\nu}$ in kinetic theory in terms of a representative phase-space distribution

 $f(\tau, p, x) = f_{BG}(Q_s(x)\tau, p/Q_s(x)) + \delta f(\tau, p, x)$

where f_{BG} characterizes typical momentum space distribution, and δf can be chosen to represent local fluctuations of initial energy momentum tensor, e.g. energy density δT^{TT} and momentum flow δT^{TT}

Energy perturbations:

$$\delta f_s(\tau_0, p, x) \propto \frac{\delta T^{\tau \tau}(x)}{T_{BG}^{\tau \tau}(x)} \times \frac{\partial}{\partial Q_s(x)} f_{BG}$$

local amplitude

 $\frac{\partial}{\partial Q_s(x)} f_{BG}\Big(\tau_0, p/Q_s(x)\Big)$

representative form of phase-space distribution

Macroscopic pre-equilibrium evolution

Energy-momentum tensor on the hydro surface can be reconstructed directly from initial conditions according to



$$T^{\mu\nu}(\tau,x) = T^{\mu\nu}_{BG}\Big(Q_s(x)\tau\Big) + \int_{Disc} G^{\mu\nu}_{\alpha\beta}\Big(\tau,\tau_0,x,x_0,Q_s(x)\Big) \,\,\delta T^{\alpha\beta}(\tau_0,x_0)$$

non-equilibrium evolution of (local) average background non-equilibrium Greens function of energy-momentum tensor

Effective kinetic theory simulations only need to be performed once to compute background evolution and Greens functions



Background evolution and Greens functions still depend on variety of variables, $Q_s(x)$, α_s , ...

-> Identify appropriate scaling variables to reduce complexity

Since ultimately evolution will match onto visc. hydrodynamics, check wether hydrodynamics admits scaling solution

1st order hydro:
$$T^{\tau\tau}(\tau) = T^{\tau\tau}_{Ideal}(\tau) \left(1 - \frac{8}{3} \frac{\eta/s}{T_{eff}\tau} + \dots\right)$$

where $T_{Ideal}^{\tau\tau}(\tau)$ is the Bjorken energy density and $T_{eff} = \tau^{-1/3} \lim_{\tau \to \infty} T(\tau) \tau^{1/3}$

Natural candidate for scaling variable is $x_s = T_{eff} \tau / (\eta / s)$

Background — Scaling variables

Numerical simulation of background evolution in effective kinetic theory

$$\left(\partial_{\tau} - \frac{p_z}{\tau}\right) f(\tau, |\mathbf{p}_{\perp}|, p_z) = \mathcal{C}[f]$$





Non-equilibrium evolution of background $T^{\mu\nu}$ is a unique function of x_s

Scaling property extends beyond hydrodynamic regime in the relevant range of (large) couplings

Greens functions

Greens functions describe evolution of energy/momentum perturbations on top of a (locally) homogenous boost-invariant background

-> Description of perturbations in Fourier space

Decomposition in a complete basis of tensors leaves a total of 10 independent functions, e.g. for energy perturbations

energy response momentum response $\tilde{G}_{\tau\tau}^{\tau\tau}(\tau, \tau_0, \mathbf{k}) = \tilde{G}_s^s(\tau, \tau_0, |\mathbf{k}|)$, $\tilde{G}_{\tau\tau}^{\tau i}(\tau, \tau_0, \mathbf{k}) = \frac{\mathbf{k}^i}{|\mathbf{k}|} \tilde{G}_s^v(\tau, \tau_0, |\mathbf{k}|)$, shear stress response

$$\tilde{G}_{\tau\tau}^{ij}(\tau,\tau_0,\mathbf{k}) = \tilde{G}_s^{t,\delta}(\tau,\tau_0,|\mathbf{k}|) \ \delta^{ij} + \tilde{G}_s^{t,k}(\tau,\tau_0,|\mathbf{k}|) \ \frac{\mathbf{k}^i \mathbf{k}^j}{|\mathbf{k}|^2} \ ,$$

Numerically computed in eff. kinetic theory by solving linearized Boltzmann equation on top of non-equilibrium background

$$\left(\partial_{\tau} + \frac{i\mathbf{p}_{\perp}\mathbf{k}_{\perp}}{p} - \frac{p_{z}}{\tau}\right)\,\delta\widetilde{f}(\tau, |\mathbf{p}_{\perp}|, p_{z}; \mathbf{k}_{\perp}) = \delta\mathcal{C}[f, \widetilde{\delta f}]$$

and computing appropriate moments of δf

Greens functions

Free-streaming:

Energy-momentum perturbations propagate as a concentric wave traveling at the speed of light

energy/momentum response:

$$G_s^{s/v}(\tau,\tau_0,\mathbf{x}-\mathbf{x}_0) = \frac{1}{2\pi(\tau-\tau_0)}\delta\Big(|\mathbf{x}-\mathbf{x}_0| - (\tau-\tau_0)\Big)$$



 $X-X_0 = T-T_0$



(c.f. Vredevoogd, Pratt PRC79 (2009) 044915, Keegan, Kurkela, Mazeliauskas, Teaney JHEP 1608 (2016) 171)

energy response:
$$\tilde{G}_{s}^{s}(\tau, \tau_{0}, k) = \tilde{G}_{s}^{s}(\tau, \tau_{0}, k = 0) \left(1 - \frac{1}{2}k^{2}(\tau - \tau_{0})^{2} \tilde{s}_{s}^{(2)} + ...\right),$$

momentum response: $\tilde{G}_{s}^{v}(\tau, \tau_{0}, k) = \tilde{G}_{s}^{s}(\tau, \tau_{0}, k = 0) \left(-ik(\tau - \tau_{0}) \tilde{s}_{v}^{(1)} + ...\right),$

shear response: determined by hydrodynamic constitutive relations

$$\begin{split} \tilde{G}_s^s(\tau,\tau_0,k=0) &= \left(\frac{T^{\tau\tau}(\tau_0)}{T^{\tau\tau}(\tau)}\right) \left(\frac{3T^{\tau\tau}(\tau) - T^{\eta}_{\eta}(\tau)}{3T^{\tau\tau}(\tau_0) - T^{\eta}_{\eta}(\tau_0)}\right) \qquad \tilde{s}_v^{(1)} = \frac{1}{2} \qquad \tilde{s}_s^{(2)} = \frac{1}{2} \left(1 + \frac{2}{3x_s}\right) \\ \text{background evolution} \qquad \text{``long wave-length constants''} \end{split}$$

Greens functions — Scaling variables



Non-equilibrium Greens functions show universal scaling in $x_s = T_{eff} \tau / (\eta / s)$ and $k(\tau - \tau_0)$ beyond hydro limit

Satisfy hydrodynamic constitutive relations for sufficiently large times $x_s >> 1$ and long wave-length k (τ - τ_0) << 1

Scaling properties ensure that pre-equilibrium evolution of energy momentum tensor can be expressed in terms of

Background:
$$T_{BG}^{\mu\nu}(x_s)$$
 Greens-functions: $G_{\alpha\beta}^{\mu\nu}\left(x_s, \frac{x-x_0}{\tau-\tau_0}\right)$

which are computed once and for all in numerical kinetic theory simulation

Since dependence of coupling constant α_s has been re-expressed in terms of physical parameter η/s , can now perform event-by-event simulations for variety of physical parameters

- 1) Evolve IP-Glasma initial conditions to early time $\tau_0 = 0.2$ fm/c 2) Macroscopic pre-equilibrium evolution to hydro initialization time τ_{Hydro}
- 3) Hydrodynamic evolution from τ_{Hydro} ($\eta/s = 2/(4\pi)$ | conformal EoS)

Energy density in the central region of Pb+Pb collision



Consistent description of pre-equilibrium dynamics ensures smooth transition to hydrodynamics at times τ>τ_{Hydro}

Energy density profile in Pb+Pb collision



No sensitivity to switching time τ_{Hydro} from pre-equilibrium to hydro

Energy density profile in Pb+Pb collision



Pre-equilibrium flow profile in Pb+Pb collision:



No sensitivity to switching time τ_{Hydro} from pre-equilibrium to hydro

Pre-equilibrium flow profile in Pb+Pb collision:



Universality of long wave-length response leads to very similar results for free-streaming and effective kinetic theory

(c.f. Vredevoogd, Pratt PRC79 (2009) 044915, Keegan, Kurkela, Mazeliauskas, Teaney JHEP 1608 (2016) 171)

Hadronic observables in single Pb+Pb event:



Hydro initialization time τ_{Hydro} [fm/c]

Very little to no sensitivity to switching time τ_{Hydro} from pre-equilibrium to hydro for dN/dy, <p τ >, <v $_2$ >, ...

Significant progress in understanding early time dynamics of heavy-ion collisions from weak-coupling perspective

Development of macroscopic description which enables event-by-event description and can be used in phenomenological modeling of heavy-ion collisions

Description in macroscopic framework is completely general and can be used beyond weak coupling limit

-> Direct comparisons with strong coupling limit should be possible

Several interesting directions beyond bulk phenomenology are starting to be explored

topological transitions, quark production & anomalous transport, photon production