

# Parameterizing the Geometry of the QGP on an Event-by-Event Basis

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**Abstract.** Ultra-relativistic heavy-ion collisions create a nuclear fireball that serves as a powerful laboratory for probing the frontiers of Quantum Chromodynamics (QCD). In recent years, there has been growing interest in the study of small collision systems—such as proton-proton ( $pp$ ) and proton-nucleus ( $pA$ )—at facilities like RHIC and the LHC. Many of the assumptions underlying the energy loss formalism developed in the Djordjevic-Gyulassy-Levai-Vitev (DGLV) model, break down in these small systems. In this work, we present an extension of the DGLV formalism that specifically accounts for the unique features of small system dynamics. This is achieved by relaxing the large formation time approximation and introducing an additional correction term that accounts for short path lengths in the medium. By relaxing these assumptions, one encounters a more intricate analytic structure for the energy loss, and thus increased computational demands; we address this challenge by developing a novel numerical scheme. Our approach accurately parametrizes the geometry of the quark-gluon plasma (QGP), resulting in a dramatic computational speedup—improving efficiency by up to seven orders of magnitude.

## 1 Introduction

The nuclear modification factor ( $R_{AA}$ ) is a key observable for studying the energy loss of high transverse momentum ( $p_T$ ) particles traversing the Quark-Gluon Plasma (QGP). Rooted in Bjorken’s jet quenching framework [1], the  $R_{AA}$  quantifies the suppression of particle yields in heavy-ion collisions relative to proton-proton systems. Experiments at RHIC observed a suppression in light hadron spectra by a factor of five [2], signaling strong partonic energy loss in the QGP.

More recently, signatures of QGP formation—including quarkonium suppression, strangeness enhancement, and collective flow—have also been identified in small collision systems at RHIC and the LHC [3]. However, small systems pose unique challenges, such as centrality bias, which arises from correlations between soft and hard particles [4].

Azimuthal anisotropies in detected spectra, as quantified by the  $v_n$  flow coefficients, offer deeper insights into the properties of the QGP, including transport properties and the path-length-dependent energy loss of partons [5, 6]. The azimuthal anisotropies can be characterized through a Fourier decomposition of the observed spectra, the Fourier decompositions are given in terms of the  $v_n$  Fourier coefficients and the reaction plane angle  $\psi_n$ . Experimentally, the reaction plane angle is not accessible; as such, observable quantities which couple soft and hard hadrons in a given centrality class serve as approximations to the reaction plane angle have been developed [7, 8].

Various energy loss models have been developed to describe the  $R_{AA}$ , but it has proven to be a challenge on theoretical grounds to simultaneously predict the nuclear modification factor and the  $v_n$  coefficients—a tension known as the  $R_{AA} \otimes v_n$  puzzle. The tension is attributed to the omission of soft-sector fluctuations and event-by-event fluctuations of the initial state [5].

In this work we use the energy loss model developed by Faraday and Horowitz (FH) which has improved the DGLV radiative energy loss formalism by including short path-length corrections and collisional energy loss using Hard Thermal Loop (HTL) kinematics [9]. Their model adopts a static, brick-like medium and parameterizes the scattering center using an effective path length [9, 10, 11, 12, 13].

The primary focus of this manuscript will be to develop a framework that describes the nuclear modification factor while simultaneously describing the anisotropic flow harmonics. We extend upon the work of FH by including event-by-event fluctuations of the bulk geometry and relaxing the brick-like and static simplification of the bulk's geometry.

## 2 Parametrization of Trajectories

The energy loss of a parton in the QGP depends on its path and the medium's geometry. As calculating energy loss for all possible trajectories can be computationally expensive (see section 2.2 for a more in depth discussion), we follow the previous work of [9, 10] by mapping each trajectory to two parameters—which the energy loss can be made to be dependent on—the energy loss can then be averaged over these parameters to capture the global effects of the collision system.

We model the energy loss in terms of a scattering center density  $\bar{\rho}$ , which is related to the medium density  $\rho$  via

$$\rho = \frac{N_s}{A_\perp} \bar{\rho}, \quad (1)$$

where  $N_s$  is the number of scatterings,  $A_\perp$  the transverse area. This medium density  $\rho$  can be expressed in terms of the medium's temperature through elementary thermodynamic relations,

$$\rho = \frac{4\zeta(3)(4+n_f)}{\pi^2} T^3. \quad (2)$$

To model  $\rho$  along each trajectory, we fit the medium density with power-law profile via

$$\rho_{fit}(z) = \rho_0 \left( \frac{\tau_0}{z} \right)^{1.2} \theta(z_c - z) \theta(z - \tau_0), \quad (3)$$

with the formation time of the medium  $\tau_0 = 0.4$  fm. Then, using eq. (2), the corresponding temperature profile is

$$T_{fit}(z) = \left( \frac{\pi^2 \rho_{fit}(z)}{4\zeta(3)(4+n_f)} \right)^{1/3}. \quad (4)$$

For each path, the cutoff  $z_c$  is determined by the distance at which  $T$  drops below the thermalization temperature, and  $\rho_0$  is fixed by the following condition

$$\int dz T_{Hydro}(\mathbf{x}_0 + z\hat{\phi}, \tau_0) \stackrel{!}{=} \int dz T_{fit}(z), \quad (5)$$

where  $T_{Hydro}$  is a hydrodynamic temperature distribution obtained from the IP-Glasma model [14]. To generalize to an entire event ensemble, we define a probability distribution over  $(\rho_0, z_c)$  for a fixed angle  $\phi$  through the medium via

$$P_{geo}(\rho_0, z_c|\phi) = \int d^2\mathbf{x}_0 n_{coll}(\mathbf{x}_0) \times \delta(\rho_0(\mathbf{x}_0, \phi) - \rho_0) \delta(z_c(\mathbf{x}_0, \phi) - z_c), \quad (6)$$

where we have assumed that the initial parton production scales with  $n_{coll}$ .

### 2.1 Energy Loss

In this work, we model the energy loss of a parton moving through the QGP by taking into account radiative and elastic contributions. Following FH [9, 11, 12, 13] we use the WHDG [10] formalism along with its short path length correction to model the radiative sector of the energy loss, and we model the elastic energy loss by using an effective field theory called Hard Thermal Loops (HTL). Our implementation of the energy loss relaxes the static brick assumption used by FH; the formalism we make use of allows the Debye mass  $\mu$  and gluon mass  $m_g$  to depend on the temperature of the medium at each point of the parton's trajectory through the medium. The radiative and elastic formalism grants us access to the probability distributions  $P_{rad}(x|\rho_0, z_c)$  and  $P_{el}(x|\rho_0, z_c)$  respectively, where  $x$  is the fractional energy loss of the incoming parton; the total energy loss for each event in the azimuthal direction  $\phi$  can then be found via the convolution of  $P_{rad}$  and  $P_{el}$ , and then averaging over the geometry of the medium, i.e.

$$P_{tot}(x|\phi) = \int d\rho_0 dz_c P_{geo}(\rho_0, z_c|\phi) \int d\epsilon P_{el}(\epsilon|\rho_0, z_c) P_{rad}(x - \epsilon|\rho_0, z_c) \quad (7)$$

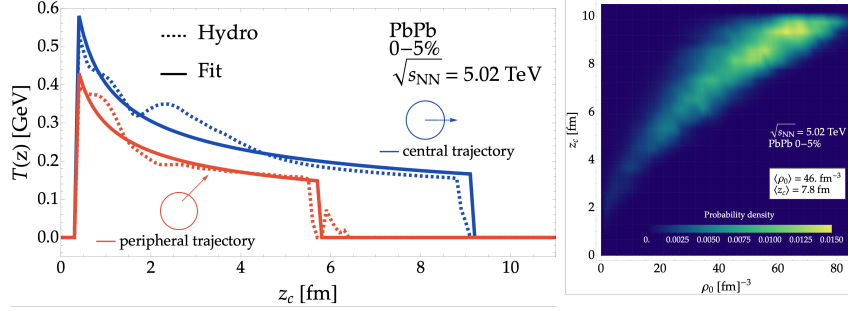


Figure 1: (Left): Comparison of temperature distributions obtained from hydrodynamic simulations (dotted curves) and the corresponding fitted temperature distribution (solid curves) for two paths through the medium. The red curves have initial positions of  $\mathbf{x}_0 = (1.5, 4.5)$  fm and angles  $\phi = \pi/4$ , and are representative of characteristic *peripheral* trajectories; the blue curves have initial positions of  $\mathbf{x}_0 = (0, 0)$  fm and angles  $\phi = 0$ , and are representative of characteristic *central* trajectories. The hydrodynamic temperature distribution comes from a PbPb 0–5% collision system at  $\sqrt{s_{NN}} = 5.02$  TeV. For the peripheral trajectory,  $z_c = 5.7$  fm and  $\rho_0 = 30.5$  fm $^{-3}$ ; for the central trajectory,  $z_c = 9.1$  fm and  $\rho_0 = 74.9$  fm $^{-3}$ . (Right): Correlation between  $\rho_0$  and  $z_c$  parameters for 0–5% PbPb collision system at  $\sqrt{s_{NN}} = 5.02$  TeV. The first moments for the  $P_{geo}(\rho_0, z_c)$  distribution (with the angular dependency integrated out) are found for this collision system to be  $\langle \rho_0 \rangle = 46.0$  fm $^{-3}$  and  $\langle z_c \rangle = 7.8$  fm.

## 2.2 Numerical challenges of the energy loss

Conceptually, the most simple method of calculating the energy loss on a path-by-path basis would be to compute the energy loss using the *true* hydrodynamic temperature profile and then using eq. (2) to find the associated medium density. To capture the global effects, one would then simply average over all the different paths through the medium. However, the total number of paths one would need to consider to resolve the QGP would typically involve  $\sim 20$  angles and a grid of size  $\sim 15 \times 15$  (fm) $^2$  with intervals of  $\sim 0.05$  (fm). To capture the event-by-event fluctuations, one would need to calculate the energy loss another  $\sim 10^3$  times, one for each event.

This computational expensiveness is why we introduce the  $\rho_0$  and  $z_c$  parameters, as these parameters allow us to calculate the energy loss as a function of the two parameters and then use the  $P_{geo}(\rho_0, z_c|\phi)$  distribution (which is computationally inexpensive) to capture the global effects. The entire phase space of possible paths requires a grid of  $\rho_0$  and  $z_c$  parameters such that  $\rho_0 \in [1, 100]$  fm $^{-3}$  and  $z_c \in [0.4, 10.4]$  fm with intervals of 10 fm $^{-3}$  and 1 fm respectively. Thus, calculating the energy loss by using the fitted parameter method proves to be seven orders of magnitude quicker than more direct method describe in the preceding paragraph.

## 3 Observables

### 3.1 Nuclear Modification Factor

The nuclear modification factor of a parton  $q$  in an  $A + A$  collision system is defined as

$$R_{AA}^q(p_T, \phi) \equiv \frac{d^2 N_{AA,f}^q / dp_T d\phi}{N_{coll} d^2 N_{pp,f}^q / dp_T d\phi} \quad (8)$$

$$= \int \frac{dx}{1-x} P_{tot}(x|\phi) \frac{f_{AA}^q\left(\frac{p_T}{1-x}, \phi\right)}{f_{pp}^q(p_T, \phi)} \quad (9)$$

where  $d^2 N_{AA,f}^q / dp_T d\phi$  and  $d^2 N_{pp,f}^q / dp_T d\phi$  are the final-state parton spectra in  $A + A$  and  $p + p$  collisions respectively;  $N_{coll}$  is the average number of binary collisions, typically calculated using the Glauber model [15]. The second equality (eq. (9)) grants us access to the  $R_{AA}$  on theoretical grounds, and is true if one assumes the following: (1) The partonic spectrum of the initial state in a  $p + p$  collision, scales like the partonic spectrum of the initial state in an  $A + A$  collision, weighted by  $1/N_{coll}$ . (2) All modifications to the  $A + A$  differential spectrum arises from the energy loss through interactions with the medium. (3) The proton-proton spectrum is unmodified by the medium. Note that we have defined the notational device  $f_{pp/AA}^q(p_T, \phi)$  as:

$$f_{pp}^q(p_T, \phi) \equiv \frac{d^2 N_{pp,i}^q}{dp_T d\phi} \quad \& \quad f_{AA}^q(p_T, \phi) \equiv \frac{1}{N_{coll}} \frac{d^2 N_{AA,i}^q}{dp_T d\phi} \quad (10)$$

where in eq. (10), the  $i$  subscript is used to denote the spectrum of the (theoretically accessible but experimentally inaccessible) initial state.

### 3.2 Azimuthal Anisotropies

The observed azimuthal anisotropy in high- $p_T$  hadrons allows for the study of the energy loss and the path length dependence of hard partons moving through the QGP [5]. In this work, an analogue to the  $v_n\{SP\}$  coefficients observed experimentally is developed for the first time in the field from a theoretical point of view (see section 3.4). This is done so that the theory predictions made in our study are more comparable to what is being observed experimentally. As part of the framework developed here, we incorporate event-by-event fluctuations of the initial state into the calculation of the  $v_n$  and  $v_n\{SP\}$  coefficients. We demonstrate that accounting for these fluctuations is crucial for addressing the  $R_{AA} \otimes v_2$  puzzle (see fig. 2).

### 3.3 Azimuthal Anisotropies from Fourier Expansions

The azimuthal anisotropy in the distributions of the observed final state hadrons can be characterized by a Fourier expansion [16]

$$\frac{2\pi}{N_{AA,f}^{h,k}} \frac{dN_{AA,f}^{h,k}}{d\phi} = 1 + 2 \sum_{n=1}^{\infty} v_n^k \cos(n[\phi - \psi_n^k]) \quad (11)$$

where the  $v_n^k$  coefficients are the uniquely determined Fourier coefficients and the  $k$  index specifies a particular event. Normalizing the angular hadronic  $R_{AA}^{h,k}(p_T, \phi)$  by the  $R_{AA}^{h,k}(p_T)$  averaged over all angles, one finds the following expression

$$\frac{R_{AA}^{h,k}(p_T, \phi)}{R_{AA}^{h,k}(p_T)} = 1 + 2 \sum_{n=1}^{\infty} v_n^k \cos(n[\phi - \psi_n^k]). \quad (12)$$

### 3.4 Scalar Product $v_n$

The scalar product (SP) method of determining the reaction plane experimentally is achieved through the use of  $\vec{Q}_n$  flow vectors. The  $\vec{Q}_n$  vectors provide an approximation to the reaction plane [16] and are determined from the final state particles via

$$\mathcal{Q}_n \equiv Q_n e^{in\Psi_n} = \sum_j e^{in\phi_j}, \quad (13)$$

One may then take the real and imaginary parts of  $\mathcal{Q}_n$  to be the  $x$  and  $y$  component of the  $\vec{Q}_n$  vector respectively. To access the  $\vec{Q}_n$  in our energy loss formalism, we make use of the IP-Glasma model which is then evolved with the MUSIC viscous relativistic (2 + 1) D hydrodynamics code, followed by UrQMD microscopic hadronic transport [14]. The  $j$  index in eq. (13) is summed over final state particles from all oversampled UrQMD simulations and the azimuthal angle is  $\phi_j = \arctan 2(p_j^y, p_j^x)$ . The  $v_n\{SP\}$  coefficients are defined here as

$$v_n\{SP\} \equiv \frac{\langle \vec{u}_n \cdot \vec{Q}_n \rangle}{\sqrt{\langle Q_n^2 \rangle}} \quad (14)$$

$$= \sum_{k=1}^{N_e} \frac{\int_0^{2\pi} d\phi R_{AA}^{h,k}(p_T, \phi) \vec{Q}_n^k \cdot \vec{u}_n^k(\phi)}{2\pi \sqrt{\langle Q^2 \rangle} N_e R_{AA}^{h,k}(p_T)} \quad (15)$$

in line with what is commonly done experimentally [7, 8]. In eq. (14) the vector  $\vec{u}_n$  is defined as  $\vec{u}_n \equiv (\cos n\phi, \sin n\phi)$ , where  $\phi$  is the azimuthal direction of the hadronic candidate. The quantity  $\langle \vec{u}_n \cdot \vec{Q}_n \rangle$  in eq. (14), is an average over all events and over different  $\vec{u}_n$  vectors associated with the observed hadrons;  $\langle Q_n^2 \rangle$  is only averaged over the events. The second equality (eq. (15)) provides us access to the  $v_n\{SP\}$  coefficient on theoretical grounds; eq. (15) follows if one assumes that the distribution of initial hard jet production is proportional to the number of binary collision density  $n_{coll}^k(\vec{x})$ .

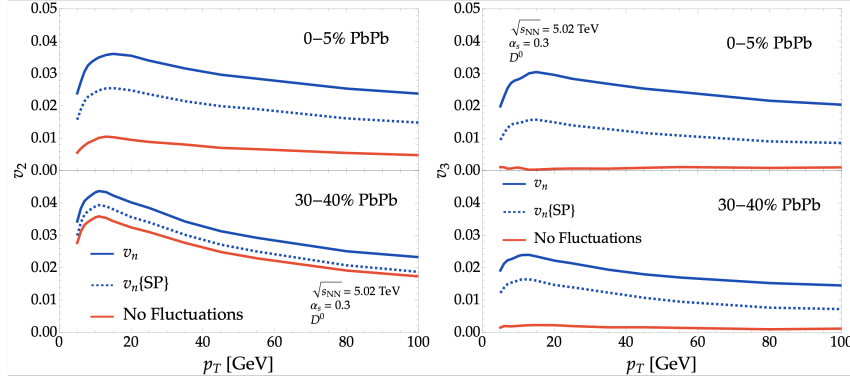


Figure 2: (Left): Comparison of Fourier, Scalar Product  $v_2$  and the Fourier  $v_2$  without fluctuations as a function of  $p_T$  for  $D^0$  mesons. All curves are calculated with the strong coupling fixed at  $\alpha_s = 0.3$ . The top (bottom) panel shows the results where the geometric average is calculated over a PbPb 0-5% (30-40%) centrality collision system. (Right): The same as (Left) but for  $v_3$  instead of  $v_2$ .

#### 4 Results

In this section we present our model's predictions for the  $R_{AA}$  and  $v_2$  coefficients. All comparisons to data are made with the strong coupling  $\alpha_s$  varied between 0.25 and 0.3. Note that we only present results for large system data (PbPb) as calculations for small systems are yet to be complete.

In fig. 3 we show the prediction of the model for charged particles in PbPb collisions at 0-5% and 30-40% centrality for both the  $R_{AA}$  and the  $v_2$ . The  $R_{AA}$  data is from the CMS [17] and ATLAS [18] experiments respectively; the  $v_2$  data comes from the same collaborations [7, 8].

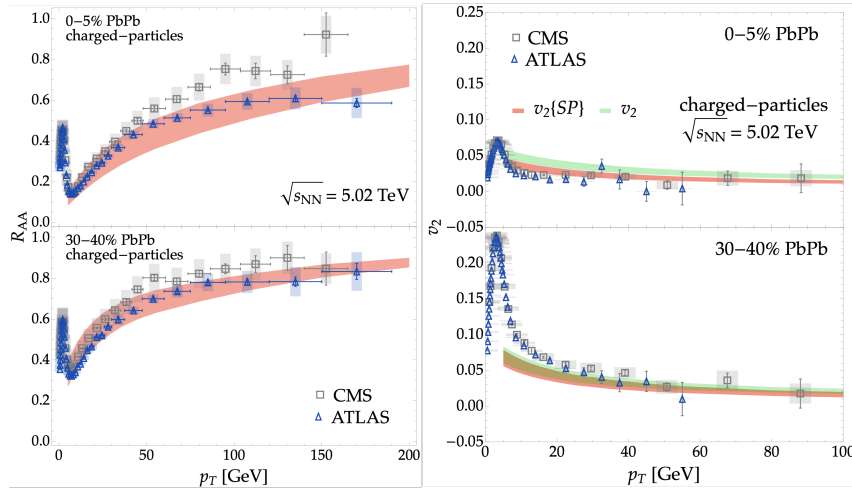


Figure 3: (Left): Comparison of  $R_{AA}$  predictions to CMS [17] and ATLAS [18] data for charged particles in 0-5% (top) and 30-40% (bottom) PbPb. (Right):  $v_2$  predictions for charged particles vs CMS/ATLAS in 0-5% (top) and 30-40% (bottom) PbPb [7, 8].

#### 5 Conclusion

In this manuscript, we presented a modification to the formalism developed by FH in [9, 11, 12, 13], in which the static brick assumption has been relaxed. The model we present here takes into account the temperature profile of the medium as a function of the parton's path through said medium; the dynamic nature of the temperature on a path-by-path basis is captured by parametrizing a power law dependency (eq. (3)) of the medium density through two fitted parameters. The energy loss of a parton moving through the medium can be calculated in terms of these

two fitted parameters, this leads to a dramatic numerical speed up—which is vitally important for enough data to be generated to capture the effects of the event-by-event fluctuations.

For high- $p_T$  charged particles, the model is in good agreement with  $v_2$  data for PbPb collisions systems and shows the correct centrality dependence. The same conclusions can be drawn for the models predictive capabilities when comparing to the  $R_{AA}$  of high- $p_T$  charged particles in PbPb collision systems.

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