



Contribution ID: 183

Type: **Contributed talk**

Efficiently simulating quarkonium's master equation beyond the dipole approximation

Wednesday, 10 July 2024 17:35 (20 minutes)

QTRAJ is a computer code that simulates the propagation of quarkonium in the quark-gluon plasma (QGP) based on the quantum trajectories' algorithm. This algorithm solves a master equation in which the quarkonium is treated as an open quantum system (OQS). The specific master equation is obtained through the potential non-relativistic QCD (pNRQCD) approach, but so far has been restricted to the regime $rT \ll 1$, where r is the size of the color dipole and T is the temperature. This limit is accurate for $\Upsilon(1S)$ but the applicability to other quarkonium states is dubious.

A major advantage of this approach is that it turns a 3D spatial evolution into a 1D Schrödinger equation with a non-hermitian Hamiltonian, drastically reducing the computational cost.

We generalize the code by extending to the regime $rT \sim 1$ in the one-gluon exchange approximation. This is done by implementing new jump operators between the resonances and expanding them in plane waves, giving rise to a variation of the algorithm present in QTRAJ 1.0. We will be showing a review of this approach comparing the $rT \ll 1$ and $rT \sim 1$ cases, and we will discuss prospect for phenomenological application to excited states.

session

F. Heavy Flavor and Quarkonia

Primary authors: Dr BERAUDO, Andrea (INFN - Torino); MARTÍNEZ VERA, Jorge Manuel (Università degli Studi di Torino & Universitat de Barcelona); Dr ESCOBEDO ESPINOSA, Miguel Ángel (ICCUB); Dr PAROTTO, Paolo (UniTo)

Co-author: Prof. STRICKLAND, Michael (Kent State University)

Presenter: MARTÍNEZ VERA, Jorge Manuel (Università degli Studi di Torino & Universitat de Barcelona)

Session Classification: F. Heavy Flavor and Quarkonia