

# **pNRQCD for the study of nonequilibrium evolution of quarkonium in a medium**

**M. Escobedo**

**pNRQCD and open quantum system,  
singlet and octet density master evolution  
equations, Lindblad equations**

Results obtained in collaboration with :

M. Escobedo, A. Vairo, J. Soto (van der Griend) N.B. (master equations);

**based on**

- (1) N. Brambilla, M.A. Escobedo, A. Vairo and P. Vander Griend  
*Transport coefficients from in medium quarkonium dynamics*  
arXiv:1903.08063
- (2) N. Brambilla, M.A. Escobedo, J. Soto and A. Vairo  
*Heavy quarkonium suppression in a fireball*  
Phys. Rev. **D97** (2018) 074009 arXiv:1711.04515
- (3) N. Brambilla, M.A. Escobedo, J. Soto and A. Vairo  
*Quarkonium suppression in heavy-ion collisions: an open quantum system approach*  
Phys. Rev. **D96** (2017) 034021 arXiv:1612.07248

# Quarkonium in a fireball

- After the heavy-ion collisions, heavy quark-antiquarks propagate freely up to 0.6 fm.
- From 0.6 fm to the freeze-out time  $t_F$  they propagate in the medium.
- We assume the medium infinite, homogeneous and isotropic.
- We assume the heavy quarks comoving with the medium.
- We assume the medium to be locally in thermal equilibrium,  
i.e., the temperature  $T$  of the medium changes (slowly) with time:

$$T = T_0 \left( \frac{t_0}{t} \right)^{v_s^2}, \quad t_0 = 0.6 \text{ fm}, \quad v_s^2 = \frac{1}{3} \text{ (sound velocity)}$$

## Quarkonium as a Coulombic bound state

The lowest quarkonium states (1S bottomonium and charmonium, 2S bottomonium) are the most tightly bound. For these we assume the hierarchy of energy scales

$$M \gg \frac{1}{r} \sim M\alpha_s \gg T \sim gT \gg \text{any other scale}, \quad v \sim \alpha_s$$

## Density matrices

- **Subsystem**: heavy quarks/quarkonium
- **Environment**: quark gluon plasma

We may define a **density matrix** in pNRQCD for the heavy quark-antiquark pair in a singlet and octet configuration:

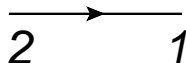
$$\begin{aligned} \langle \mathbf{r}', \mathbf{R}' | \rho_s(t'; t) | \mathbf{r}, \mathbf{R} \rangle &\equiv \text{Tr} \{ \rho_{\text{full}}(t_0) S^\dagger(t, \mathbf{r}, \mathbf{R}) S(t', \mathbf{r}', \mathbf{R}') \} \\ \langle \mathbf{r}', \mathbf{R}' | \rho_o(t'; t) | \mathbf{r}, \mathbf{R} \rangle \frac{\delta^{ab}}{8} &\equiv \text{Tr} \{ \rho_{\text{full}}(t_0) O^{a\dagger}(t, \mathbf{r}, \mathbf{R}) O^b(t', \mathbf{r}', \mathbf{R}') \} \end{aligned}$$

$t_0 \approx 0.6$  fm is the time formation of the plasma.

The system is in **non-equilibrium** because through interaction with the environment (quark gluon plasma) singlet and octet quark-antiquark states continuously transform in each other although **the number of heavy quarks is conserved**:  $\text{Tr}\{\rho_s\} + \text{Tr}\{\rho_o\} = 1$ .

# Evolution of the density matrix in pNRQCD. Singlet

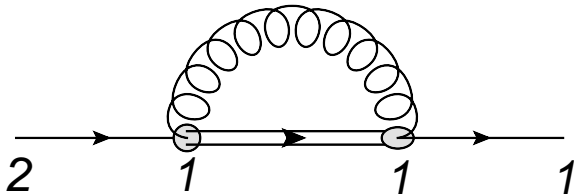
$$\partial_t \rho_s = -i[h_s, \rho_s] - i\Sigma \rho_s + i\rho_s \Sigma^\dagger + \mathcal{F}(\rho_o)$$



We use the non-equilibrium quantum field theory formalism to perform this computation. Tree level is equivalent to the  $T = 0$  LO evolution

# Evolution of the density matrix in pNRQCD. Singlet

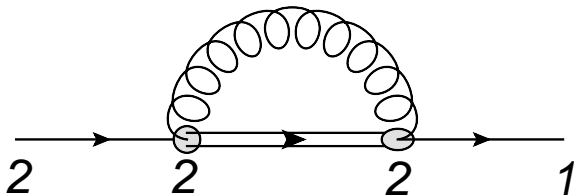
$$\partial_t \rho_S = -i[h_S, \rho_S] - i\Sigma \rho_S + i\rho_S \Sigma^\dagger + \mathcal{F}(\rho_S)$$



Self-energy diagram contributes to screening and to the decay width.

# Evolution of the density matrix in pNRQCD. Singlet

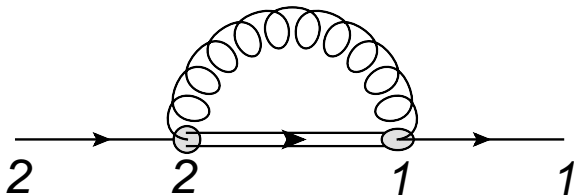
$$\partial_t \rho_S = -i[h_S, \rho_S] - i\Sigma \rho_S + i\rho_S \Sigma^\dagger + \mathcal{F}(\rho_o)$$



Hermitian conjugate of the previous diagram. We can reorganize this to diagrams in a redefinition of  $h_s$  and a term that represents the decay width.

# Evolution of the density matrix in pNRQCD. Singlet

$$\partial_t \rho_S = -i[h_S, \rho_S] - i\Sigma \rho_S + i\rho_S \Sigma^\dagger + \mathcal{F}(\rho_o)$$



The number of singlets is increased due to the octets in the medium that absorb a gluon.



# Evolution of the octet

Very similar reasoning.

$$\partial_t \rho_o = -i[H_o, \rho_o] - \frac{1}{2}\{\Gamma, \rho_o\} + \mathcal{F}_1(\rho_s) + \mathcal{F}_2(\rho_o)$$

- Computationally costly system of coupled equations. The density matrix contains many information, especially in the non-Abelian case (QCD).
- Total number of heavy particles is conserved.  $Tr(\rho_s) + Tr(\rho_o)$  is a constant of the evolution.
- Only if  $T \gg E$  the interaction with the medium can be considered local in time and the evolution is Markovian.

# Time scales

- Environment correlation time  $\tau_E \sim \frac{1}{T}$ .
- System intrinsic time scale  $\tau_S \sim \frac{1}{E}$ .
- System relaxation time  $\tau_R \sim \frac{1}{\Sigma} \sim \frac{1}{\alpha_s a_0^2 \Lambda^3}$  where  $\Lambda$  is either  $T$  or  $E$  depending on the temperature regime.

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- Because  $\frac{1}{r} \gg T, E$ , it follows that  $\tau_R$  is the longest time scale. Markovian approximation.
  - If  $T \gg E$  then  $\tau_S \gg \tau_E$ , the interaction with the medium is instantaneous from the point of view of the system. Lindblad equation limit.

# The $\frac{1}{r} \gg T, m_D \gg E$ regime

Brambilla, M.A.E., Soto and Vairo (2017-2018)

$$\partial_t \rho = -i[H(\gamma), \rho] + \sum_k (C_k(\kappa) \rho C_k^\dagger(\kappa) - \frac{1}{2} \{C_k^\dagger(\kappa) C_k(\kappa), \rho\})$$

$$\kappa = \frac{g^2}{6 N_c} \text{Re} \int_{-\infty}^{+\infty} ds \langle T E^{a,i}(s, \mathbf{0}) E^{a,i}(0, \mathbf{0}) \rangle$$

The heavy quark diffusion coefficient

$$\gamma = \frac{g^2}{6 N_c} \text{Im} \int_{-\infty}^{+\infty} ds \langle T E^{a,i}(s, \mathbf{0}) E^{a,i}(0, \mathbf{0}) \rangle$$

No lattice QCD information on  $\gamma$  but we observe that we reproduce data better if it is small. For comparison, in pQCD

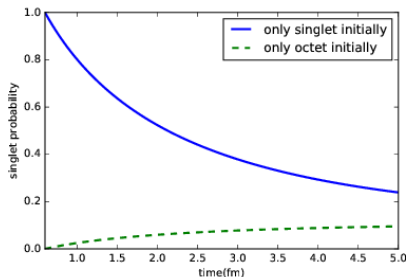
$$\gamma = -2\zeta(3) C_F \left( \frac{4}{3} N_c + n_f \right) \alpha_s^2(\mu_T) T^3 \approx -6.3 T^3$$

# Approach to equilibrium in the static limit

We assume that  $\frac{\kappa}{T^3}$  is a constant. Then,  $\Gamma(T) = \Gamma(T_0) \left(\frac{T}{T_0}\right)^3$ .

In the static limit we can solve the equation analytically using Bjorken evolution, the approach to equilibrium is power like.

$$\rho_s(t) = \frac{1}{9} - \frac{8}{9} \left( \frac{\rho_o(t_0)}{8} - \rho_s(t_0) \right) \left( \frac{t_0}{t} \right)^{\frac{9}{8}\Gamma(T_0)t_0}$$



# Initial conditions

At any point in time, our density matrix is diagonal by boxes in colour space

$$\rho(t) = \begin{pmatrix} \rho_s(t) & 0 \\ 0 & \rho_o(t) \end{pmatrix} .$$

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## Physical insight

- The production of a heavy quark pair is a high energy process.
- The production of a pair in a singlet state is suppressed by a factor  $\alpha_s(m)$  with respect to that of the octet.

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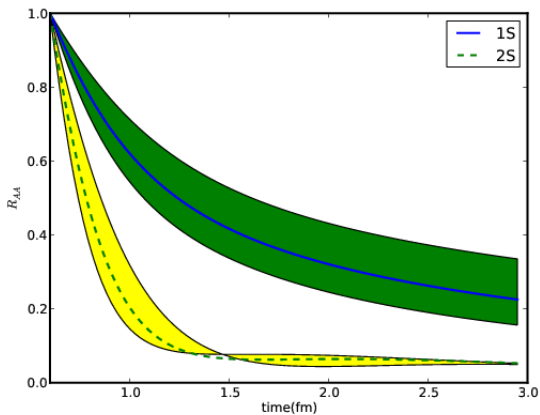
## Initial conditions

- We take as initial condition  $\rho_s(0) = A|\mathbf{r} = \mathbf{0}\rangle\langle\mathbf{r} = \mathbf{0}|$  and  $\rho_o(0) = \frac{\delta}{\alpha_s(m)}\rho_s(0)$ . Then, we evolve the density matrix in the vacuum from  $t = 0$  to  $t = t_0 = 0.6 \text{ fm}$ .
- $\delta$  is a free parameter. However, we observe that the results depend very little on its value.



We compute  $R_{AA}$  (before feed-down) using the formula

$$R_{AA}(n) = \frac{\langle n | \rho(t_f) | n \rangle}{\langle n | \rho(t_0) | n \rangle}$$



- The evolution equations that we are able to deduce by systematically exploiting the separation between energy scales are quantum.
- Approximations that can lead to simplifications (like semi-classical equations), as for example the rotating wave approximation or multiple-scale analysis, are not fully compatible with a systematic expansion of ratios of energy scales.
- At the moment, it is difficult to determine quantitatively the importance of quantum coherence.
- Quantum coherence might force us to change how we think about initial conditions. Not in terms of probabilities, but in terms of density matrix.

## Beyond Bjorken evolution

- Due to the high numerical cost of our computations, we have only implemented Bjorken evolution. This overestimates suppression.
- We are studying how to reduce this cost. When this is done we will implement more realistic hydro.
- We have calculated the binding energy and decay width in an anisotropic plasma. (Biondini, Brambilla, Escobedo and Vairo (2017)).

# Future improvements

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## Quarkonium in a hot wind

We have also studied how the potential, the binding energy and the decay width are modified when quarkonium is moving with a finite velocity with respect to the medium. This might be important for the  $R_{AA}$  dependence with transverse momentum. (Escobedo, Giannuzzi, Mannarelli and Soto (2013)).

# Future improvements II

- Use more realistic initial conditions.
- Use more recent determinations of  $\kappa$  and  $\gamma$  and include a more complex temperature dependence.
- Relax the assumption  $T \gg E$ .