## **Open Quantum Systems for Quarkonia in QGP**

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References: Akamatsu-Asakawa-Kajimoto-Rothkopf 1805.00167, Kajimoto-Akamatsu-Asakawa-Rothkopf (18), Akamatsu (15,13), Akamatsu-Rothkopf (12)

## **Motivation and Outline**



- 1. Basics of Open Quantum System
- 2. Application to Quarkonium in QGP
- ${\bf a}$ shapes. As in previous analyses  $[2,2,4]$ , possible  $[2,2,4]$  $\alpha$  dot-dashed blue lines) are also shown, as are the individual  $\alpha$ 3. Quantum State Diffusion Simulation for a Heavy Quark

 $\theta$  differences in Pb-Pb-Pb and pp acceptances due to physical pp acceptances due to physical **Can we understand the data in terms of in-medium QCD forces at high** *T***?** What do we learn from heavy-ion data?

# **Basics of Open Quantum System**

#### Open quantum systems



1. Total system consists of system (S) and environment (E)

$$
\mathcal{H}_{\mathsf{tot}}=\mathcal{H}_{S}\otimes\mathcal{H}_{E}
$$

2. Hamiltonian

$$
H_{\text{tot}} = H_S \otimes 1 + 1 \otimes H_E + H_I, \quad H_I = \sum H_I^{(S)} \otimes H_I^{(E)}
$$

3. Reduced density matrix & Master equation

$$
\rho_S(t) \equiv \text{Tr}_E \rho_{\text{tot}}(t), \quad i \frac{d}{dt} \rho_{\text{tot}} = [H_{\text{tot}}, \rho_{\text{tot}}] \quad \rightarrow \quad \underbrace{i \frac{d}{dt} \rho_S = ?}_{\text{Markovian limit}}
$$

#### 4. Theoretical methods

- $\blacktriangleright$  Influence functional path integral representation for the master equation
- $\triangleright$  Schwinger-Dyson equation time evolution equation for the density matrix

## Time scale hierarchies

### Three basic time scales

- ▶ Environment correlation time *τ<sup>E</sup>*
- ▶ System intrinsic time scale *τ<sup>S</sup>*
- ▶ System relaxation time *τ<sup>R</sup>*



#### Time scale hierarchies

▶ Quantum Brownian motion

 $\tau_E \ll \tau_R \; , \qquad \tau_E \ll \tau_S \quad \rightarrow$  good description in phase space Markov approx. derivative expansion

▶ Quantum optical system

*τE*  $\ll$  *τR τ* Markov approx. *, τ<sup>S</sup> ≪ τ<sup>R</sup> →* good description in eigenbasis rotating wave approx.

It is very important to estimate the relevant time scales We adopt QBM-type approximation scheme to study quarkonium

#### Time scales of a quarkonium quantum Brownian motion in QGP

#### ▶ Environment (QGP) correlation time *τ<sup>E</sup>*

1. Time scales of QGP



2. Heavy quarks mostly couple to electric field

$$
\tau_E \sim \frac{1}{gT}
$$

▶ System (Quarkonium) intrinsic time scale *τ<sup>S</sup>*

Orbital period  $=$  inverse energy gap  $=$  formation time



#### ▶ System relaxation time *τ<sup>R</sup>*

Kinetic equilibration / color relxation (for a single  $HQ$  / longer for a quarkonium)

$$
\tau_R^{\rm kin} \sim \frac{M}{T} \frac{1}{g^4 T \ln(1/g)}, \quad \tau_R^{\rm color} \sim \frac{1}{g^2 T}
$$

*⇒*Time scale hierarchy for quarkonium quantum Brownian motion

$$
\tau_E \ll \tau_R, \quad \tau_E \ll \tau_S \to g \underbrace{\ll}_{\text{color}} 1, \quad g^3 \ln(1/g) \underbrace{\ll}_{\text{kinetic}} \frac{M}{T} \underbrace{\ll}_{\text{potential}} \frac{g}{\alpha^2} \sim \frac{100}{g^3}
$$

Scale hierarchy satisfied/challenged at weak/strong coupling

Open quantum system by path integral

1. Path integral

$$
\rho_{\text{tot}}(t, \underbrace{x, y}_{\in S}, \underbrace{X, Y}_{\in E}) = \int dx_0 dy_0 dX_0 dY_0 \int_{x_0, y_0, X_0, Y_0}^{x, y, X, Y} \mathcal{D}[\bar{x}, \bar{y}, \bar{X}, \bar{Y}]
$$

$$
\times \underbrace{\rho_{\text{tot}}(0, x_0, y_0, X_0, Y_0)}_{\text{factorizable } \rho_S(0) \otimes \rho_E^{\text{eq}}} e^{iS_{\text{tot}}[\bar{x}, \bar{X}] - iS_{\text{tot}}[\bar{y}, \bar{Y}]}
$$

2. Influence functional  $S_{IF}$  [Feynman-Vernon (63)]

$$
\rho_S(t, x, y) = \underbrace{\int dX dY \delta(X - Y)}_{\text{trace out } E = \text{path closed at } t} \rho_{\text{tot}}(t, x, y, X, Y)
$$

$$
= \int dx_0 dy_0 \rho_S(0, x_0, y_0) \int_{x_0, y_0}^{x, y} \mathcal{D}[\bar{x}, \bar{y}] e^{iS_S[\bar{x}] - iS_S[\bar{y}] + iS_{\text{IF}}[\bar{x}, \bar{y}]}
$$

Influence functional contains all the information of the open system

#### Coarse graining for quantum Brownian motion

1. Influence functional up to quadratic order

$$
iS_{\text{IF}}[x,y]=-\,\frac{1}{2}\int_{0}^{t}dt_{1}dt_{2}(x,y)_{(t_{1})}\underbrace{\begin{pmatrix}G_{11}&-G_{12}\\-G_{21}&G_{22}\end{pmatrix}}_{\text{correlation function of }E} \begin{pmatrix}x\\y\end{pmatrix}_{(t_{2})}
$$

2. Choice of time after coarse graining

$$
t^> = \max(t_1, t_2), \quad s = |t_1 - t_2|
$$

3. Derivative expansion in *s* when  $\tau_S \gg \tau_E$ 

$$
iS_{\text{IF}}[x,y] = 2\gamma m T \int_0^t dt^>(x,y) \begin{pmatrix} -1 & 1\\ 1 & -1 \end{pmatrix} \begin{pmatrix} x\\ y \end{pmatrix}
$$
  
momentum diffusion (fluctuation)  

$$
+ i\pi m \int_0^t dt^>(x,y) \begin{pmatrix} -1 & -1\\ t & -1 \end{pmatrix} \begin{pmatrix} \dot{x} \\ \dot{x} \end{pmatrix}
$$

$$
+\underbrace{i\gamma m\int_0^t dt^>(x,y)\begin{pmatrix}-1 & -1\\ 1 & 1\end{pmatrix}\begin{pmatrix}\dot{x}\\ \dot{y}\end{pmatrix}}_{\text{drag force (dissipation)}}+\cdots
$$

Influence functional is single time integral after coarse graining

#### Caldeira-Leggett master equation

1. From path integral to differential equation

$$
\rho_S(t, x, y) = \int dx_0 dy_0 \rho_S(0, x_0, y_0) \int_{x_0, y_0}^{x, y} \mathcal{D}[\bar{x}, \bar{y}] e^{iS_S[\bar{x}]-iS_S[\bar{y}]+iS_{\text{IF}}[\bar{x}, \bar{y}]}
$$

$$
\rightarrow i \frac{\partial}{\partial t} \rho_S(t, x, y) = H(x) \rho_S(t, x, y) - H(y) \rho_S(t, x, y)
$$

$$
-i \gamma \left[ \underbrace{2mT(x-y)^2}_{\text{fluctuation}} + \underbrace{(x-y)(\partial_x - \partial_y)}_{\text{dissipation}} \right] \rho_S(t, x, y)
$$

▶ Equivalent to Fokker-Planck equation through Wigner transform

2. Ehrenfest equations

$$
\frac{d}{dt}\langle p\rangle=-2\gamma\langle p\rangle,\quad \frac{d}{dt}\langle H\rangle=-4\gamma\left(\langle H\rangle-\frac{T}{2}\right)
$$

Quantum mechanical description for Brownian motion

Caldeira-Leggett master equation is NOT Lindblad

1. Positivity of the density matrix

$$
\forall |\alpha\rangle \rightarrow \langle \alpha|\rho_S|\alpha\rangle \ge 0
$$

2. Any Markovian positive map is written by the Lindblad equation  $[Lindblad (76)]$ 

$$
\frac{d}{dt}\rho_S(t) = -i[H, \rho_S] + \sum_{i=1}^N \gamma_i \left( L_i \rho_S L_i^{\dagger} - \frac{1}{2} L_i^{\dagger} L_i \rho_S - \frac{1}{2} \rho_S L_i^{\dagger} L_i \right)
$$

3. Lindblad form is obtained when higher order expansion is included  $[p_{\text{loss}}(93)]$ 

$$
S_{\text{IF}} = \underbrace{S_{\text{fluct}} + S_{\text{diss}}}_{\text{Calderia-Legget}} + \underbrace{S_{(2)}}_{\propto x \dot{x}} \underbrace{\times x \dot{x}}_{\text{diss}}
$$

 $If L ∼ x + \dot{x}, then L<sup>†</sup>L ⇒ \dot{x}\dot{x}$ 

Lindblad equation is not a must, but theoretically more complete

# **Application to Quarkonium in QGP**

#### Influence functional for heavy quarks

1. Heavy quarks in the non relativistic limit

$$
\mathcal{L}_I = -gA_0^a \left[ Q^{\dagger} t^a Q + Q_c t^a Q_c^{\dagger} \right] = -gA_0^a \rho^a
$$

2. Influence functional:  $-gA_0^a\rho^a$  is a source term for QGP

$$
e^{iS_{\text{IF}}[\rho]}\simeq \int \mathcal{D}[A,q]\rho_{\text{QGP}}^{\text{eq}}[A,q]\exp\Bigl[i\int_{x\in \text{CTP}}\left\{\mathcal{L}_{\text{QGP}}(A,q)-gA_0^a\rho^a\right\}\Bigr]
$$

- 3. Perturbative expansion in terms of gluon correlators in QGP
	- ▶ Choose  $t^> = \max(t_1, t_2)$  as a single time variable in  $S_{IF}$

$$
iS_{\text{IF}} = -g^2 \int_{t}^{t} \int_{x}^{t} (\rho_1^a, \ \rho_2^a)_{(t,x)} \int_{s>0} \left[ \begin{array}{cc} G^F & -G^< \\ -G^> & G^{\tilde{F}} \end{array} \right]_{(s,x-y)} \left( \begin{array}{c} \rho_1^a \\ \rho_2^a \end{array} \right)_{(t-s,y)}
$$

- 4. Derivative expansion based on hierarchy of time scales between *G* and *ρ*
	- ▶ Expand in *s*

$$
S_{\text{IF}} = \underbrace{S_{\text{pot}} + S_{\text{fluct}}}_{\propto \rho \rho} + \underbrace{S_{\text{diss}}}_{\propto \rho \dot{\rho}} + \underbrace{S_{(2)}}_{\propto \dot{\rho} \dot{\rho}} + \cdots
$$

#### More on influence functional for heavy quarks

1. Gluon correlators at low frequencies

$$
V(r) = g^2 G_R(\omega = 0, r), \quad D(r) = g^2 T \frac{\partial}{\partial \omega} \underbrace{\sigma(\omega = 0, r)}_{\text{spectral function}}
$$

2. Using the *ra*-basis:  $\rho_r = (\rho_1 + \rho_2)/2$ ,  $\rho_a = \rho_1 - \rho_2$ 

 $\blacktriangleright$  potenital

$$
S_{\text{pot}} = \int_{t} \int_{x y} V(x - y) \rho_{a}(x) \rho_{r}(y)
$$

▶ fluctuation

$$
S_{\text{fluct}} = \frac{i}{2} \int_t \int_{x y} D(\boldsymbol{x} - \boldsymbol{y}) \rho_a(x) \rho_a(y) \Leftrightarrow S_{\text{fluct}}^{CL} = 2i\gamma m T x_a^2
$$

 $\blacktriangleright$  dissipation

$$
S_{\text{diss}} = -\frac{1}{2T} \int_t \int_{x\mathbf{y}} D(\mathbf{x} - \mathbf{y}) \rho_a(x) \dot{\rho}_r(y) \Leftrightarrow S_{\text{diss}}^{CL} = -2\gamma m x_a \dot{x}_r
$$

▶ 2nd order

$$
S_{(2)} \simeq \frac{i}{4} \int_t \int_{xy} \frac{D(\boldsymbol{x} - \boldsymbol{y})}{8T^2} \dot{\rho}_a(x) \dot{\rho}_a(y)
$$

Fluctuation-dissipation theorem in QGP sector relates  $S<sub>fluct</sub>$  and  $S<sub>dis</sub>$ 

## Master equation from influence functional

## THIS IS THE MOST DIRTY PART

#### 1. From path integral to functional differential equation

▶ Analogous to "Schrödinger equation from path integral"

$$
\underbrace{\rho_S[t,Q_1^{\text{fin}},Q_2^{\text{fin}}]}_{\text{``wave function'' at }t} = \int dQ_{1,2}^{\text{ini}} \underbrace{\rho_S[0,Q_1^{\text{init}},Q_2^{\text{init}}]}_{\text{initial ``wave function''}} \int_{Q_{1,2}^{\text{init}}}^{Q_{1,2}^{\text{fin}}} \mathcal{D}[Q_{1,2}] e^{iS_S[Q_1] - iS_S[Q_2] + iS_{\text{IF}}[Q_1,Q_2]}\n \rightarrow \frac{\partial}{\partial t} \rho_S[t,Q_1,Q_2] = \mathcal{L}[Q_1,Q_2] \rho_S[t,Q_1,Q_2]
$$

#### 2. From functional density matrix to density matrix

(i) Recall that the basis of the functional space is the coherent state

$$
|Q\rangle\sim e^{-\int_{\mathbf{w}}Q(\mathbf{x})\hat{Q}^{\dagger}(\mathbf{x})}|\Omega\rangle
$$

(ii) Introduce a heavy quark by functional differentiation

$$
\rho_Q(t, \boldsymbol{x}, \boldsymbol{y}) \sim \frac{\delta}{\delta Q_1(\boldsymbol{x})} \frac{\delta}{\delta Q_2(\boldsymbol{y})} \rho_S[t, Q_1, Q_2]|_{Q=0}
$$

There must be several ways to derive the master equation from *S*IF

#### Lindblad equation for a quarkonium in QGP

$$
\frac{d}{dt}\rho_{Q\bar{Q}}(t) = -i[H, \rho_{Q\bar{Q}}] + \sum_{k} \left( L_k \rho_{Q\bar{Q}} L_k^{\dagger} - \frac{1}{2} L_k^{\dagger} L_k \rho_{Q\bar{Q}} - \frac{1}{2} \rho_{Q\bar{Q}} L_k^{\dagger} L_k \right)
$$

$$
L_k = \sqrt{D(k)} e^{ikx/2} \left[ 1 + \frac{ik \cdot \nabla_x}{\frac{4MT}{\Delta x_Q \sim k/MT}} \right] e^{ikx/2} + \text{heavy antiquark}
$$

- ▶ Scattering  $Qq \rightarrow Qq$  with momentum transfer *k* with rate  $D(k)$
- $\triangleright$  Momentum transfer without recoil = stochastic potential (no dissipation)

$$
L_k = \underbrace{\sqrt{D(k)}e^{ikx}}_{\Delta p_Q = k} + \text{heavy antiquark}
$$

- $\triangleright$  Quantum dissipation from heavy quark recoil during a collision
- $\triangleright$  Coefficient  $1/4MT$  fixed by fluctuation-dissipation theorem for QGP correlators

**Quantum State Diffusion Simulation for a Heavy Quark**

#### Quantum State Diffusion simulation for Lindblad equation

1. Lindblad equation

$$
\frac{d}{dt}\rho_S(t) = -i[H, \rho_S] + \sum_{i=1}^N \gamma_i \left( L_i \rho_S L_i^{\dagger} - \frac{1}{2} L_i^{\dagger} L_i \rho_S - \frac{1}{2} \rho_S L_i^{\dagger} L_i \right)
$$

- 2. Stochastic unravelling
	- ▶ Equivalent to a nonlinear stochastic Schrödinger equation [Gisin-Percival (92)]

$$
\rho_S(t) = \lim_{N \to \infty} \frac{1}{N} \sum_{i=1}^N \frac{|\phi_i(t)\rangle \langle \phi_i(t)|}{||\phi_i(t)||^2} = \mathsf{M} \left[ \frac{|\phi(t)\rangle \langle \phi(t)|}{||\phi(t)||^2} \right],
$$
\n
$$
|d\phi\rangle = -iH|\phi(t)\rangle dt + \sum_n \Big( \underbrace{2\langle L_n^{\dagger} \rangle_{\phi} L_n}_{\text{nonlinear in } \phi} - L_n^{\dagger} L_n \Big) |\phi(t)\rangle dt + \sum_n L_n |\phi(t)\rangle d\xi_n,
$$
\n
$$
\underbrace{\langle d\xi_n d\xi_m^* \rangle}_{\text{complex noise}} = 2\delta_{nm} dt
$$

#### Apply this technique to heavy quark Lindblad equation

Nonlinear stochastic Schrödinger equation for a heavy quark

▶ Nonlinear stochastic Schrödnger equation

$$
d\phi(x,t) = \phi(x, t + dt) - \phi(x,t)
$$
  
\n
$$
\simeq \left(i\frac{\nabla^2}{2M} - \frac{1}{2}D(0)\right)\phi(x)dt + d\xi(x)\phi(x)
$$
  
\n
$$
+ \frac{dt}{||\phi(t)||^2}\int d^3y D(x - y)\phi^*(y)\phi(y)\phi(x) + \mathcal{O}(T/M)
$$

▶ Correlation of complex noise field

$$
\langle d\xi(x)d\xi^*(y)\rangle = D(x-y)dt, \quad \langle d\xi(x)d\xi(y)\rangle = \langle d\xi^*(x)d\xi^*(y)\rangle = 0
$$

 $\blacktriangleright$  Density matrix for a heavy quark

$$
\rho_Q(x,y,t)=\mathsf{M}\left[\frac{\phi(x,t)\phi^*(y,t)}{||\phi(t)||^2}\right]
$$

What is the equilibrium solution of the Lindblad equation? How does a heavy quark approach equilibrium?

QSD simulation for a single heavy quark in an external potential

Numerical setups

$$
V_{\text{ext}}(x) = 0, \quad \frac{1}{2}M\omega^2 x^2, \quad -\frac{\alpha}{\sqrt{x^2 + r_c^2}}
$$

$$
D(x) = \gamma \exp\left[-x^2/l_{\text{corr}}^2\right]
$$



$$
\Delta x = \frac{1}{M} \ll l_{\text{corr}} = \frac{10}{M} \ll N_x \Delta x = \frac{128}{M}
$$

Do the density matrix approach *∝* exp(*−H/T*)?

## Solitonic wave function in one sampling



Wave function is localized because of the nonlinear evolution equation

Equilibration of a heavy quark:  $V_{\text{ext}} = 0$ 

Time evolution of momentum distribution

▶ Relaxation time of corresponding classical system *Mτ*relax *∼* 300



Equilibrium momentum distribution is the Boltzmann distribution!

## Equilibration of a heavy quark:  $V_{\text{ext}} = V_{\text{HO/Coulomb}}$

Time evolution of eigenstate occupation (lowest 3 levels)

 $\blacktriangleright$  Harmonic potential (left), regularized Coulomb potential (right)



Eigenstate occupation relaxes to a static state Relaxation time depends on the initial state and rate equation is inapplicable  $\frac{22}{22}$ 

Equilibrium distribution of a heavy quark:  $V_{\text{ext}} = V_{\text{HO/Coulomb}}$ 

Equilibrium distribution of eigenstates (lowest 10 levels)

▶ Harmonic potential (top), regularized Coulomb potential (bottom)



We also checked that off-diagonal part is 0 within statistical fluctuation

Eigenstate distribution in the external potential is also the Boltzmann distribution

#### $QSD$  simulation without quantum dissipation (= stochastic potential)

Heavy quark is overheated because energy increases without dissipation

 $\blacktriangleright$  Neglect  $\mathcal{O}(T/M)$  terms in the nonlinear stochastic Schrödinger equation



#### Dissipation is more important for smaller bound state because decoherence is ineffective

## Summary and outlook

### Quantum State Diffusion simulation for Lindblad equation

- ▶ Equivalent to nonlinear stochastic Schrödinger equation (integro-differential equation)
- ▶ Numerically confirm the equilibration of a heavy quark *→* Can be shown analytically?

#### Possible application

- $\triangleright$  Quarkonium evolution in heavy-ion collisions [Akamatsu et al, in progress]
- ▶ Dark matter bound state in early universe? [Kim-Laine (17)]
- ▶ Cold atomic gases? [Braaten-Hammer-Lepage (16)]

# Back Up

## Explicit form of gluon correlators in HTL approximation

$$
G_R(\omega = 0, r) = -\frac{e^{-m_D r}}{4\pi r},
$$

$$
\frac{\partial}{\partial \omega} \sigma_{ab,00}(0, \vec{r}) = \int \frac{d^3k}{(2\pi)^3} \frac{\pi m_D^2 e^{i\vec{k} \cdot \vec{r}}}{k(k^2 + m_D^2)^2},
$$

$$
m_D^2 = \frac{g^2 T^2}{3} \left( N_c + \frac{N_f}{2} \right)
$$

#### Example 1 – Quantum optical master equation



 $\triangleright$  A two-level atom in a photon gas

$$
i\frac{d}{dt}\rho_A = \gamma \underbrace{(N(\omega_0) + 1)}_{\text{emission}} \left[ \sigma_{-}\rho_A \sigma_{+} - \frac{1}{2}\sigma_{+}\sigma_{-}\rho_A - \frac{1}{2}\rho_A \sigma_{+}\sigma_{-} \right]
$$

$$
+ \gamma \underbrace{N(\omega_0)}_{\text{absorption}} \left[ \sigma_{+}\rho_A \sigma_{-} - \frac{1}{2}\sigma_{-}\sigma_{+}\rho_A - \frac{1}{2}\rho_A \sigma_{-}\sigma_{+} \right]
$$

▶ Approximations

$$
\underbrace{\rho_{\text{tot}}(t) \simeq \rho_A(t) \otimes \rho_B^{\text{eq}}}_{\text{Born approx. (weak coupling) }}, \quad \underbrace{\tau_B \ll \tau_R \equiv 1/\gamma}_{\text{Markov approx.}}, \quad \underbrace{\tau_A \equiv 1/\omega_0 \ll \tau_R}_{\text{rotating wave approx.}}
$$

- ▶ Environment correlation time *τ<sup>B</sup>*
- ▶ System intrinsic time scale *τA*, system relaxation time *τ<sup>R</sup>*

Master equation is an effective description at  $\tau_R \gg \tau_B$  for  $\tau_A \ll \tau_R$ 

## Example 2 – Quantum Brownian motion



Caldeira-Leggett model [Caldeira-Leggett (83)]

 $\triangleright$  Brownian particle linearly coupled to harmonic oscillators

$$
i\frac{d}{dt}\rho_A=[H_A,\rho_A]+\underbrace{\gamma[x,\{p,\rho_A\}]}_{\text{drag force}}-\underbrace{2i\gamma m T[x,[x,\rho_A]]}_{\text{momentum diffusion}}
$$

▶ Approximations

$$
\underbrace{\rho_{\text{tot}}(t) \simeq \rho_A(t) \otimes \rho_B^{\text{eq}}}_{\text{Born approx. (weak coupling) }}, \quad \underbrace{\tau_B \ll \tau_R \equiv 1/\gamma}_{\text{Markov approx.}}, \quad \underbrace{\tau_B \ll \tau_A}_{\text{derivative expansion}}
$$

- ▶ Environment correlation time *τ<sup>B</sup>*
- ▶ System intrinsic time scale *τA*, system relaxation time *τ<sup>R</sup>*

Master equation is an effective description at  $\tau_R \gg \tau_B$  for  $\tau_A \gg \tau_B$