

Lattice NRQCD Study of Thermal Sommerfeld Factor

Seyong Kim

Sejong University

based on

SK, M. Laine (ITP, U of Bern), arXiv:1602.08105

Outline

- 1 Introduction
- 2 Method
- 3 Result and Discussion

Thermal Sommerfeld Effect in QGP

- chemical equilibration rate of heavy quark in Quark-Gluon Plasma (QGP)

$$\Gamma_{\text{chem}} \simeq \frac{8\pi\alpha_s^2}{3M^2} \left(\frac{MT}{2\pi}\right)^{3/2} e^{-M/T} \left[\frac{\bar{S}_1}{3} + \left(\frac{5}{6} + Nf\right) \bar{S}_8 \right].$$

- Sommerfeld effect enhances the Born matrix elements

$$|\mathcal{M}_{\text{resummed}}|^2 = \mathcal{S} |\mathcal{M}_{\text{tree}}|^2$$

for the color singlet

$$\mathcal{S}_1 = \frac{X_1}{1 - e^{-X_1}}, X_1 = C_F \frac{g^2}{4V}$$

and for the color octet

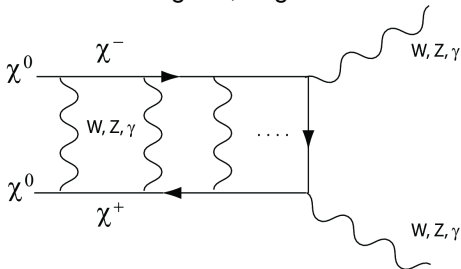
$$\mathcal{S}_8 = \frac{X_8}{e^{X_8} - 1}, X_8 = \left(\frac{N_c}{2} - C_F\right) \frac{g^2}{4V}$$

Thermal Sommerfeld Effect in QGP

- enhancement of co-annihilation for slowly moving particle :

$$\sim \frac{\alpha}{v}$$

per one rung in a ladder diagram, large for small v



How to calculate: Usual way

- consider Boltzmann equation*

$$\partial_t n \simeq -c(n^2 - n_{\text{eq}}^2) = \dot{n}_{\text{loss}} + \dot{n}_{\text{gain}}$$

with $\dot{n}_{\text{loss}} = -cn^2$

- in equilibrium, $n(t) = n_{\text{eq}}$ and $\delta\dot{n} \simeq -2cn\delta n$

$$\Gamma_{\text{chem}} = \left. \frac{\delta\dot{n}}{n} \right|_{\text{eq}} = -2 \frac{\dot{n}_{\text{loss}}}{n_{\text{eq}}}$$

- then

$$\Gamma_{\text{chem}} = \frac{2}{2N_c \int_{\mathbf{k}} f_F(E_k)} \int \int (2\pi)^4 \delta^4(P_1 + P_2 - K_1 - K_2) f_F(E_{k_1}) f_F(E_{k_2})$$

$$\left(\frac{1}{2} \sum |M_1|^2 [1 + f_B(\epsilon_{p_1})] [1 + f_B(\epsilon_{p_2})] + N_f \sum |M_2|^2 [1 - f_F(\epsilon_{p_1})] [1 - f_F(\epsilon_{p_2})] \right)$$

How to calculate: New way, Γ_{chem} as a transport coefficient

- chemical equilibration as a transport coefficient (D. Bödeker, M. Laine, JHEP07 (2012) 130, 01 (2013) 037)
- treat the approach to the equilibrium as a Langevin process

$$\delta\dot{n}(t) = -\Gamma_{\text{chem}}\delta n(t) + \xi(t)$$

$$\langle\langle \xi(t)\xi(t') \rangle\rangle = \Omega_{\text{chem}}\delta(t-t'), \quad \langle\langle \xi(t) \rangle\rangle = 0$$

where $\delta n(t)$ is the deviation from the equilibrium and $\xi(t)$ is a stochastic noise

$$\delta n(t) = \delta n(t_0)e^{-\Gamma_{\text{chem}}(t-t_0)} + \int_{t_0}^t dt' e^{\Gamma_{\text{chem}}(t'-t)}\xi(t')$$

How to calculate: New way, Γ_{chem} as a transport coefficient

- consider a correlator (classical)

$$\begin{aligned}
 \Delta_{cl}(t, t') &= \lim_{t_0 \rightarrow -\infty} \langle\langle \delta n(t) \delta n(t') \rangle\rangle \\
 &= \lim_{t_0 \rightarrow -\infty} \int_{t_0}^t dt_1 e^{\Gamma_{\text{chem}}(t_1 - t)} \int_{t_0}^{t'} dt_2 e^{\Gamma_{\text{chem}}(t_2 - t')} \langle\langle \xi(t_1) \xi(t_2) \rangle\rangle \\
 &= \Omega_{\text{chem}} \lim_{t_0 \rightarrow -\infty} \int_{t_0}^t dt_1 e^{\Gamma_{\text{chem}}(t_1 - t)} \int_{t_0}^{t'} dt_2 e^{\Gamma_{\text{chem}}(t_2 - t')} \delta(t_1 - t_2) \\
 &= \frac{\Omega_{\text{chem}}}{2\Gamma_{\text{chem}}} e^{-\Gamma_{\text{chem}}|t - t'|}
 \end{aligned}$$

and

$$\partial_t \partial_{t'} \Delta_{cl}(t, t') = -\frac{\Omega_{\text{chem}}}{2} \Gamma_{\text{chem}} e^{-\Gamma_{\text{chem}}|t - t'|} + \Omega_{\text{chem}} \delta(t - t')$$

How to calculate: New way, Γ_{chem} as a transport coefficient

- in frequency domain,

$$\tilde{\Delta}_{cl}(\omega) \equiv \int_{-\infty}^{\infty} dt e^{i\omega(t-t')} \Delta_{cl}(t, t') = \frac{\Omega_{\text{chem}}}{\omega^2 + \Gamma_{\text{chem}}^2}$$

and

$$\omega^2 \tilde{\tilde{\Delta}}_{cl}(\omega) = \int_{-\infty}^{\infty} dt e^{i\omega(t-t')} \partial_t \partial_{t'} \Delta_{cl}(t, t') = \frac{\omega^2 \Omega_{\text{chem}}}{\omega^2 + \Gamma_{\text{chem}}^2}$$

- then,

$$\langle (\delta n)^2 \rangle = \lim_{t_0 \rightarrow -\infty} \langle \langle \delta n(t) \delta n(t) \rangle \rangle = \frac{\Omega_{\text{chem}}}{2\Gamma_{\text{chem}}}$$

or

$$\Gamma_{\text{chem}} = \frac{\Omega_{\text{chem}}}{2\langle (\delta n)^2 \rangle}$$

How to calculate: New way, Γ_{chem} as a transport coefficient

- to access Ω_{chem} , consider correlators (for heavy quark, or non-relativistic QCD, the heavy quark number is replaced by hamiltonian) in imaginary time

$$\Omega(\tau) = \frac{1}{V} \langle \partial_t H(\tau) \partial_t H(0) \rangle_{\text{qm}}$$

or

$$\Delta(\tau) = \frac{1}{V} \langle H(\tau) H(0) \rangle_{\text{qm}}$$

then

$$\Omega_{\text{chem}} = \lim_{\Gamma_{\text{chem}} \ll \omega \ll \omega_{UV}} 2T \frac{\rho_{\Omega}(\omega)}{\omega} \quad \text{or} \quad = \lim_{\omega \ll T} 2T \omega \rho_{\Delta}(\omega)$$

and

$$\Gamma_{\text{chem}} = \frac{\lim_{\omega \rightarrow 0^+} 2T \frac{\rho_{\Omega}(\omega)}{\omega}}{2\chi_f M^2} \quad \text{or} \quad = \frac{\lim_{\omega \ll T} 2T \omega \rho_{\Delta}(\omega)}{2\chi_f M^2}$$

How to calculate: New way, Γ_{chem} as a transport coefficient

- for non-relativistic QCD), the change in the number of heavy quark/anti-quark can be treated as a “perturbation” since the pair annihilation/creation of heavy quark–anti-heavy quark has “high energy” ($\sim 2M$) and in NRQCD, $p \sim M$ is integrated out
- again, consider an imaginary time correlator

$$\Delta(\tau) = \frac{1}{V} \langle H(\tau)H(0) \rangle_{\text{qm}}$$

- and its leading order perturbation by four-fermion vertices

$$O_1(^1S_0) = \psi^\dagger \chi \chi^\dagger \psi, \quad S_M = \frac{f_1(^1S_0) O_1(^1S_0)}{M^2}$$

and

$$O_8(^1S_0) = \psi^\dagger T^a \chi \chi^\dagger T^a \psi, \quad S_M = \frac{f_8(^1S_0) O_8(^1S_0)}{M^2}$$

How to calculate: New way, Γ_{chem} as a transport coefficient

- We need to be careful about the analytic structure of these four-fermion vertices since we are interested in imaginary time formulation for finite temperature physics.
- we need to consider, a correlator

$$\Delta(\tau) = \text{Im} \frac{f_1(^1S_0)}{\pi M^2} \varepsilon(\tau)$$

with

$$\varepsilon(\tau) = \int_{\mathbf{x}, \mathbf{y}} \int_0^\beta d\tau_1 \int_0^\beta d\tau_2 \frac{\langle H(\tau, \mathbf{x}) H(0, \mathbf{y}) (\theta^\dagger \chi)(\tau_1, \mathbf{0}) (\chi^\dagger \theta)(\tau_2, \mathbf{0}) \rangle}{|\tau_1 - \tau_2|}$$

How to calculate: New way, Γ_{chem} as a transport coefficient

- numerator becomes

$$\begin{aligned}
 & \int_{\vec{x}, \vec{y}} \langle H(\tau, \vec{x}) H(0, \vec{y}) (\psi^\dagger \chi)(\tau_1, \vec{0}) (\chi^\dagger \psi)(\tau_2, \vec{0}) \rangle |_{\tau_1 > \tau_2} [\theta(\tau - \tau_1) + \theta(\tau_2 - \tau)] \\
 &= \frac{1}{Z} \text{Tr} \left[e^{-\beta \hat{\mathcal{H}}} (\hat{\psi}^\dagger \hat{\chi})(\tau_1, \vec{0}) (\hat{\chi}^\dagger \hat{\psi})(\tau_2, \vec{0}) \int_{\vec{x}, \vec{y}} \hat{H}(0, \vec{x}) \hat{H}(0, \vec{y}) \right] \\
 &= \frac{1}{Z} \sum_{m, n} \langle q\bar{q}, m | e^{-\beta \hat{\mathcal{H}}} (\hat{\psi}^\dagger \hat{\chi})(\tau_1, \vec{0}) | n \rangle \langle n | (\hat{\chi}^\dagger \hat{\psi})(\tau_2, \vec{0}) \int_{\vec{x}, \vec{y}} \hat{H}(0, \vec{x}) \hat{H}(0, \vec{y}) | \\
 &= \frac{4M^2}{Z} \sum_{m, n} e^{-\beta E_m} e^{(\tau_1 - \tau_2)(E_m - \varepsilon_n)} \langle q\bar{q}, m | \hat{\psi}^\dagger \hat{\chi} | n \rangle \langle n | \hat{\chi}^\dagger \hat{\psi} | q\bar{q}, m \rangle
 \end{aligned}$$

E_m are the states with heavy quarks, ε_n are states without heavy quarks

How to calculate: New way, Γ_{chem} as a transport coefficient

- with $C_{mn} = \frac{4M^2}{Z} e^{-\beta E_m} \langle q\bar{q}, m | \hat{\psi}^\dagger \hat{\chi} | n \rangle \langle n | \hat{\chi}^\dagger \hat{\psi} | q\bar{q}, m \rangle$

$$\varepsilon(\tau) = \sum_{m,n} C_{mn} \int_0^\beta d\tau_1 \int_0^\beta d\tau_2 \frac{e^{(\tau_1 - \tau_2)(E_m - \varepsilon_m)}}{\tau_1 - \tau_2} [\theta(\tau - \tau_1) + \theta(\tau_2 - \tau)]$$

- then with $\omega, \varepsilon_m \ll E_m$

$$\rho_\Delta(\omega) = \frac{\text{Im}f_1(1S_0)}{M^2} \sum_{m,n} C_{mn} \frac{e^{\beta\omega} - e^{-\beta\omega}}{\omega^2}$$

How to calculate: New way, Γ_{chem} as a transport coefficient

- finally,

$$\begin{aligned}
 \Omega_{\text{chem}} &= \frac{4\text{Im}f_1(1S_0)}{M^2} \sum_{m,n} C_{mn} \\
 &= 16\text{Im}f_1(1S_0) \frac{1}{Z} \sum_{m,n} e^{-\beta E_m} \langle q\bar{q}, m | \hat{\psi}^\dagger \hat{\chi} | n \rangle \langle n | \hat{\chi}^\dagger \hat{\psi} | q\bar{q}, m \rangle \\
 &= 16\text{Im}f_1(1S_0) \frac{1}{Z} \text{Tr} \left[e^{-\beta \hat{\mathcal{H}}} (\hat{\psi}^\dagger \hat{\chi})(0^+, \vec{0}) (\hat{\chi}^\dagger \hat{\psi})(0, \vec{0}) \right] \\
 &= 16\text{Im}f_1(1S_0) \langle (\psi^\dagger \chi)(0^+, \vec{0}) (\chi^\dagger \psi)(0, \vec{0}) \rangle
 \end{aligned}$$

How to calculate: New way, Γ_{chem} as a transport coefficient

$$P_1 \equiv \frac{1}{2N_c} \text{Re} \langle G_{\alpha\alpha;ij}^\theta(\beta, \vec{0}; 0, \vec{0}) \rangle ,$$

$$P_2 \equiv \frac{1}{2N_c} \langle G_{\alpha\gamma;ij}^\theta(\beta, \vec{0}; 0, \vec{0}) G_{\gamma\alpha;ji}^{\theta\dagger}(\beta, \vec{0}; 0, \vec{0}) \rangle ,$$

$$P_3 \equiv \frac{1}{2N_c^2} \langle G_{\alpha\alpha;ij}^\theta(\beta, \vec{0}; 0, \vec{0}) G_{\gamma\gamma;ji}^{\theta\dagger}(\beta, \vec{0}; 0, \vec{0}) \rangle .$$

- singlet Sommerfeld factor

$$\bar{S}_1 = \frac{P_2}{P_1^2} .$$

- octet Sommerfeld factor

$$\bar{S}_8 = \frac{N_c^2 P_3 - P_2}{(N_c^2 - 1) P_1^2} .$$

Lattice setup

- anisotropic Euclidean lattices (i.e., the time direction lattice spacing is different from the space direction lattice spacing, $a_s/a_t = 3.5$),
 $a_s = 0.1227(8)$ fm
- $N_f = 2 + 1$ light quark flavors ($M_\pi \simeq 400$ MeV, $M_K \simeq 500$ MeV)
- $24^3 \times N_t$ lattices
- $T_c = 185$ MeV, $a_s M = 2.92$
- lattices used for bottomonium at $T \neq 0$ study (G. Aarts et al, JHEP07 (2014) 097) and electric conductivity of QGP (G. Aarts et al, JHEP-2 (2015) 186)

Lattice setup

$$G^\theta(0, \mathbf{x}; \cdot) = \frac{\delta_{\mathbf{x}, \vec{0}}}{a_s^3},$$

$$G^\theta(a_t, \mathbf{x}; \cdot) = \left(1 - \frac{a_t \mathcal{H}_0}{2n}\right)^n U_0^\dagger(0, \mathbf{x}) \left(1 - \frac{a_t \mathcal{H}_0}{2n}\right)^n G^\theta(0, \mathbf{x}; \cdot),$$

$$G^\theta(\tau + a_t, \mathbf{x}; \cdot) = \left(1 - \frac{a_t \mathcal{H}_0}{2n}\right)^n U_0^\dagger(\tau, \mathbf{x}) \left(1 - \frac{a_t \mathcal{H}_0}{2n}\right)^n (1 - a_t \delta \mathcal{H}) G^\theta(\tau, \mathbf{x}; \cdot),$$

where U_0 is a time-direction gauge link. The lowest-order Hamiltonian reads

$$\mathcal{H}_0 = -\frac{\Delta^{(2)}}{2M},$$

where $\Delta^{(2)}$ is a discretized gauge Laplacian.

Lattice setup

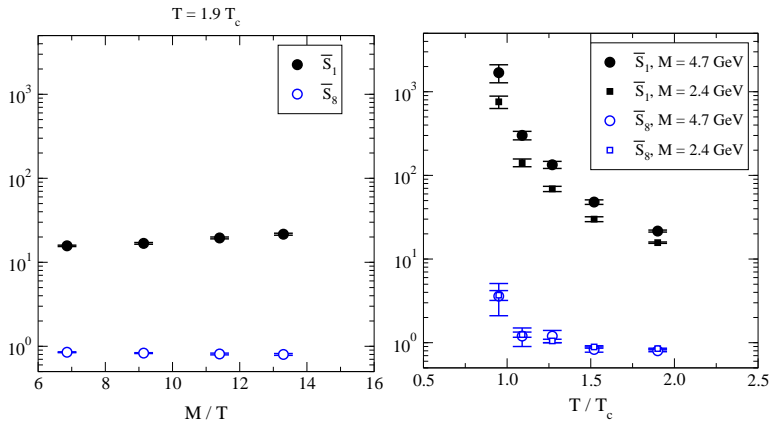
The higher order correction is

$$\delta\mathcal{H} = -\frac{(\Delta^{(2)})^2}{8M^3} + \frac{ig_0(\nabla \cdot \mathbf{E} - \mathbf{E} \cdot \nabla)}{8M^2} - \frac{g_0 \boldsymbol{\sigma} \cdot (\nabla \times \mathbf{E} - \mathbf{E} \times \nabla)}{8M^2} \\ - \frac{g_0 \boldsymbol{\sigma} \cdot \mathbf{B}}{2M} + \frac{a_s^2 \Delta^{(4)}}{24M} - \frac{a_t (\Delta^{(2)})^2}{16nM^2},$$

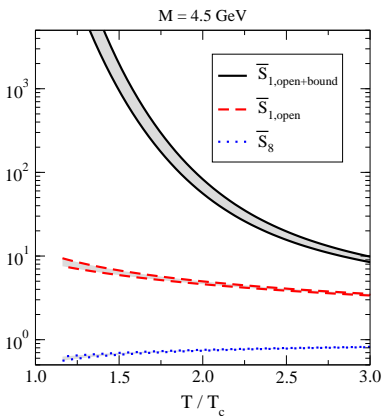
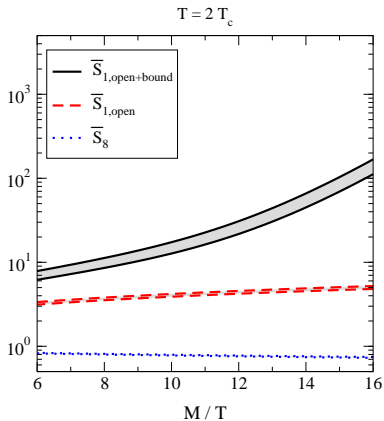
Lattice setup

N_t	T/T_c	$a_s M$	M/T	$10^4 P_1$	$10^5 P_2$	$10^6 P_3$	\bar{S}_1	\bar{S}_8
32	0.95	2.92	26.7	0.75(7)	0.938(5)	1.06(1)	1690(410)	3.6(15)
28	1.09	2.92	23.4	3.0(1)	2.78(1)	3.19(3)	301(35)	1.2(3)
24	1.27	2.92	20.0	8.0(3)	8.64(3)	10.3(1)	134(13)	1.2(2)
20	1.52	2.92	16.7	24.4(5)	28.5(1)	36.1(4)	48(3)	0.83(6)
16	1.90	2.92	13.3	68.9(9)	102.2(3)	147(1)	21.6(6)	0.80(2)
32	0.95	1.50	13.7	1.31(7)	1.308(8)	1.51(1)	758(128)	3.7(5)
28	1.09	1.50	12.0	4.6(1)	3.04(2)	3.62(3)	142(15)	1.25(9)
24	1.27	1.50	10.3	10.4(2)	7.38(4)	9.21(8)	69(5)	1.05(5)
20	1.52	1.50	8.57	25.7(3)	19.7(1)	27.2(2)	30(2)	0.89(3)
16	1.90	1.50	6.86	61.9(6)	60.1(3)	95.8(8)	15.7(3)	0.85(1)
16	1.90	2.00	9.14	57.6(6)	55.7(2)	86.4(8)	16.8(4)	0.83(1)
16	1.90	2.50	11.4	62.5(7)	76.0(2)	113(1)	19.5(5)	0.81(2)

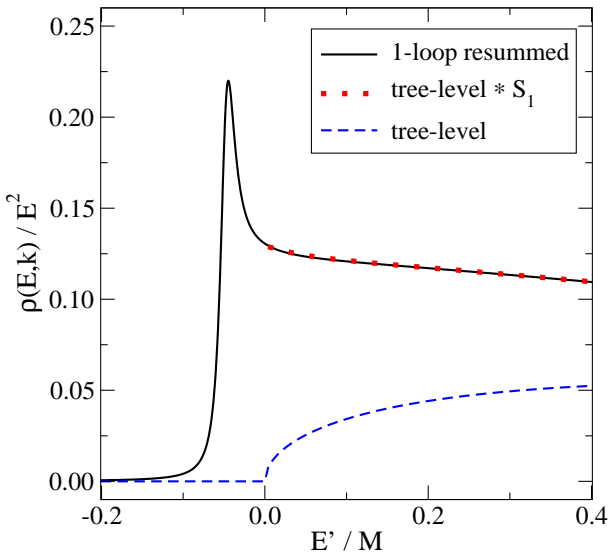
Lattice result of thermal Sommerfeld factor



Analytic estimate of thermal Sommerfeld factor



Analytic estimate of thermal Sommerfeld factor

 $M = 4.5 \text{ GeV}, T = 2 T_c$ 

Conclusion

- a real time quantity, **chemical equilibration rate**, is calculated non-perturbatively using Euclidean lattice **without analytic continuation**
- in the current relativistic heavy ion collision experiments, charm quark and bottom quarks do not chemically equilibrate
- thermal Sommerfeld effect for bottomonium co-annihilation is calculated using lattice NRQCD and is found to be **two orders of magnitude larger than perturbative estimate**
- **bound state effect** is the cause of large enhancement
- similarly, in “strongly interacting” dark matter scenario, the bound state effect is expected to enhance co-annihilation of dark matter far beyond a naive perturbative estimate.