The real-time finite-temperature static potential: a higher order calculation

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Motivation

- Check the robustness of our qualitative understanding of heavy quarkonium dissociation at finite temperature
 - ► J/ψ suppression (Matsui, Satz, 1986): heavy quarkonium states sequentially melt due to screening
 - ★ The Yukawa potential supports less and less bound states when the screening mass (~ T) increases
 - The potential develops an imaginary part, which is responsible for the melting rather than the screening (M. Laine, O. Philipsen, P. Romatschke, M. Tassler, JHEP 0703: 054, 2007)
 - * The melting temperature T_d can be parametrically estimated to be $T_d \sim -m_Q \alpha_s^{2/3} / \ln^{1/3} \alpha_s$ (M.A. Escobedo, JS, Phys. Rev. A 78, 032520 (2008))

• Provide additional physically motivated forms of the potential which may help extracting it from lattice data

The LO potential

 If p, m_D ≪ T the momentum space potential reads (G(p₀, p) is the Hard Thermal Loop (HTL) longitudinal gluon propagator, m_D ~ gT the screening mass, g the QCD coupling constant, C_F a color factor)

$$V_{\rm 1lo}(p)=g^2C_FG(0,p)$$

$$G(0,p) = -\frac{1}{m_D^2 + p^2} + \frac{i\pi T m_D^2}{p (m_D^2 + p^2)^2}$$

- ▶ When screening is important $p \leq m_D \Rightarrow \text{Im}(V_{1\text{lo}}) \gg \text{Re}(V_{1\text{lo}}) \Rightarrow$ No narrow resonance exist
- $\operatorname{Im}(V_{1\mathrm{lo}}) \sim \operatorname{Re}(V_{1\mathrm{lo}}) \Rightarrow p \equiv p_d \sim (m_D^2 T)^{\frac{1}{3}} \sim g^{\frac{2}{3}} T \Rightarrow m_D \ll p_d \ll T$
- The typical p must fulfill $p \gg p_d$ for a narrow resonance to exist

The LO calculation corresponds to the one-gluon exchange in HTL

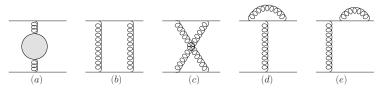
$$\delta \mathcal{L}_{HTL} = \frac{1}{2} m_D^2 F^{\mu\alpha} \int \frac{d\Omega}{4\pi} \frac{k_\alpha k_\beta}{-(k.\partial)^2} F^{\mu\beta} + m_f^2 \bar{\psi} \gamma^\mu \int \frac{d\Omega}{4\pi} \frac{k_\mu}{k.\partial} \psi$$

$$k = (1, \hat{\mathbf{k}}),$$
 $m_D^2 = g^2 T^2 (N_c + N_f/2)/3,$ $m_f^2 = g^2 T^2/16$
(Braaten, Pisarsky, 92)

- HTL ~ Integrating out the scale T ~ One-loop selfenergies at LO in the p/T expansion
- Power corrections to HTL ~ One-loop selfenergies at NLO in the p/T expansion (Manuel, JS, Stetina, 16; Carignano, Manuel, JS, 17; Carrington, Carignano, JS, 20; Ekstedt, 23; Gorda, Paatelainen, Säppi, Seppänen, 23)

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• We work in the Coulomb gauge. The relevant diagrams are



- Gluon lines in (b)-(e) = HTL longitudinal time-ordered gluon propagators
- Self-energy blob (a):
 - ★ Longitudinal gluon self-energy calculated in HTL
 - ★ Power correction calculated in QCD
- The hierarchy $m_D \ll p \ll T$ produces enormous simplifications $k \sim \text{loop momentum}$
 - (a), the longitudinal gluon self-energy calculated in HTL is dominated by k ~ p ⇒ HTL propagators and vertexes reduce to QCD ones up to m²_D/p² corrections
 - (b)-(e) are dominated by $k \sim m_D \Rightarrow$ can be expanded in $k/p, m_D/p$

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• The fully expanded result reads

$$\begin{split} V_{2,\exp}(p) &= -\frac{g^4 C_F N_c T}{16 \pi m_D p^2} \left(1 - \frac{3\pi^2}{16} + \frac{4\pi m_D}{p} + \frac{m_D^2}{p^2} \left(\frac{5\pi^2}{24} - \frac{4}{3} \right) \right) \\ &- i \frac{g^4 C_F T^2}{16 p^4} \left(N_c \left(\frac{56}{3\pi} - \left(1 - \frac{3\pi^2}{16} \right) \frac{m_D}{p} \right) - \frac{4}{\pi} \left(N_c - \frac{N_f}{2} \right) \frac{p}{T} \right) \end{split}$$

- Let $p \sim g^a T$, then the formula above holds up to corrections g^2 (g^{3a}, g^{2-a}) for the real (imaginary) part (1/3 < a < 2/3)
- We define the damped approximation by keeping factors $1/(p^2 + m_D^2)$ in the gluon propagators unexpanded
 - Expected to be more realistic for $m_D \lesssim p$
 - Reproduces the expanded formula above

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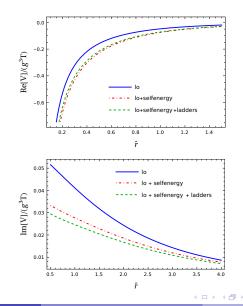
• Upon Fourier transforming, the coordinate space potential in the damped approximation reads $(\hat{r} = rm_D)$

$$\operatorname{Re}[V_2] = \frac{g^4 N_c C_F T}{64\pi^2 \hat{r}} \left\{ 8 \left(I_2(\hat{r}) - I_1(\hat{r}) \right) + \frac{e^{-\hat{r}}}{16} \left(3\pi^2 - 16 + \frac{\hat{r}}{6} \left(16 - \pi^2 \right) \right) \right\}$$

$$i \text{Im}[V_2] = -i \frac{g^3 C_F T}{16\pi^2 \hat{m}_D} \left\{ \frac{3\pi^2 - 16}{32 \,\hat{r}} l_2(\hat{r}) + \frac{7}{3} \, N_c e^{-\hat{r}} - \frac{2g \,\hat{m}_D}{\pi \hat{r}} \left(N_c - \frac{N_f}{2} \right) \left(l_1(\hat{r}) - l_2(\hat{r}) \right) \right\}$$

$$I_j(\hat{r}) = \int_0^\infty d\hat{p} \, rac{\sin\left(\hat{p}\hat{r}
ight)}{(\hat{p}^2+1)^j} \quad , \quad m_D = gT\,\hat{m}_D$$

The potential beyond LO (g = 1.8)



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- The coordinate space potential for $r \ll 1/m_D$ gets contributions from $p \sim 1/r$ and $p \lesssim m_D$
 - ► The contribution from p ~ 1/r can be obtained by Fourier transforming V_{2,exp}
 - The contribution from $p \leq m_D$ is a polynomial in r^2 , since the $\exp(i\vec{p}\vec{r})$ in the Fourier transform can be expanded
 - ▶ The damped approximation reshuffles part of the $p \lesssim m_D$ contribution out of the polynomial

Comparison with lattice data

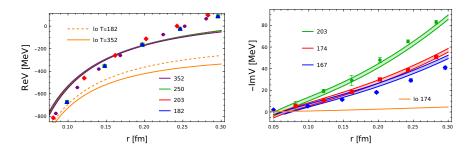
- $g \text{ small} \Rightarrow \text{short distance observables}$
 - ► Data for the potential up to 0.3 fm (Bazavov, Hoying, Kaczmarek, Larsen, Mukherjee, Petreczky, Rothkopf, Weber, 23) = [1]
 - Mass shift and decay width of Ŷ(1s) (Larsen, Meinel, Mukherjee, Petreczky, 20)≡ [2]
- We fix g from the best fit to T = 0 data of the first reference (g = 1.8)
- The soft contribution to the coordinate space potential is accounted for by including

 $\delta \text{Re}(V) = q_0 g^3 T$, $\delta \text{Im}(V) = i_0 g^3 T + i_2 g^5 r^2 T^3$

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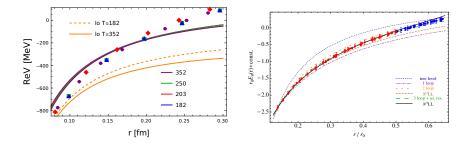
and fitting q_0 , i_0 and i_2 to the lattice data

Comparison with [1]



- Re(V) depends very little on T like the lattice data
 - The slight difference in the shape may be accounted for by higher order T-independent corrections
- The dramatic improvement in Im(V) is due to the inclusion of the soft contribution $(q_0, i_0, i_2) = (0.049, -0.021 \pm 0.002, 0.205 \pm 0.001)$

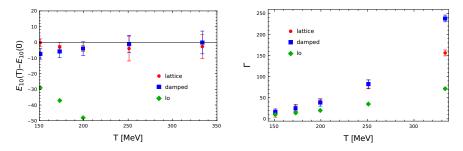
Comparison with [1]



- The rhs plot shows the T = 0 potential at different orders of perturbation theory, $r_0 = 0.468$ fm (Bazavov, Brambilla, Garcia i Tormo, Petreczky, Soto, Vairo, 12)
- The slightly different shape between our results and data in the lhs plot is similar to the one between tree level and data in the rhs plot
 It may be fixed by a higher order calculation

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Comparison with [2]



- Dramatic (considerable) improvement in the binding energy (decay width)
- The inclusion of soft contributions is crucial $(q_0, i_0, i_2) = (0.078 \pm 0.004, -0.026 \pm 0.009, 0.053 \pm 0.002)$
- The value of *i*₂ is not consistent with the one obtained from [1]
- The same size for (q_0, i_0, i_2) is expected. This is fulfilled except for i_2 from [1]

Dissociation temperature

- We define the dissociation temperature (T_d) as the temperature for which the binding energy equals the decay width
 - At LO we obtain: $T_d = 193.2 \text{ MeV}$
 - Damped approximation, fit to [1]: $T_d = 151.8 \pm 1.2$ MeV
 - Damped approximation, fit to [2]: $T_d = 225 \pm 10$ MeV
- Fit to [1] produces a very low T_d and an unnatural value of $i_2 \Rightarrow$ there might be a problem with the data

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Conclusions

- We have calculated higher order corrections to the momentum space potential, both to the real and to the imaginary parts, when $m_D \ll p \ll T$
- We have proposed an approximation (damped) that partially includes soft contributions
- We have pointed out that the soft contributions $p \lesssim m_D$ to the coordinate space potential are universal for $r \ll 1/m_D$, and can be described by a polynomial in r^2 (up to logs)
- We get a reasonable description of lattice data at short distances
 - We have been able to identify an inconsistency between two sets of lattice data
- Our results provide additional inputs for the Bayesian methods to obtain the potential from Euclidean lattice data.