

Static hybrid potential in D dimensions at short distancesAntonio Pineda¹ and Maximilian Stahlhofen^{1,2}¹*Grup de Física Teòrica, Universitat Autònoma de Barcelona, E-08193 Bellaterra, Barcelona, Spain*²*University of Vienna, Faculty of Physics, Boltzmannngasse 5, A-1090 Wien, Austria*

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We compute the energy of a static hybrid, i.e., of a hybrid quarkonium with static quark and antiquark, at short distances in $D = 4, 3$, dimensions. The soft contribution to this energy is the static potential of a color-octet quark-antiquark pair at short distances, which is known at two loops for arbitrary D . We have checked this expression employing thermal field theory methods. Using the effective field theory potential nonrelativistic QCD we calculate the ultrasoft contributions to the hybrid (and singlet) static energy at the two-loop level. We then present new results for the static hybrid energy/potential and the hybrid decay width in three and four dimensions. Finally we comment on the meaning of the perturbative results in two space-time dimensions, where the hybrid does not exist.

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I. INTRODUCTION

In a recent paper [1] we have studied the potential and energy of a static color-singlet quark-antiquark state at short distances in D space-time dimensions (though with special emphasis on the three-dimensional case). It is the aim of this paper to perform a similar analysis for the static hybrid energy and the associated octet potential.

The energies of static hybrid systems are very interesting quantities. They can teach us a lot about the dynamics of QCD, and can be potentially relevant for the theoretical description of physical hybrids made of heavy quarks. Their behavior at long distances might contain information on the detailed dynamics responsible for confinement, whether it is due to strings and, if so, of which kind. Most relevant for us, however, is that the hybrid energy at short distances approximately equals the sum of the static octet potential energy and the respective gluelump mass. This relation was first quantified in Refs. [2,3] and recently exposed in a unified and model-independent framework [4,5] using the effective field theory “potential nonrelativistic QCD” (pNRQCD) [6] (for a review see [7]).

The physics of the static hybrid system in the short distance limit is governed by, at least, two physical scales. One is the soft scale $\sim 1/r$, the inverse distance between quark and antiquark, the other is the ultrasoft scale $\sim \Delta V \equiv V_o - V_s$, where V_o and V_s are the static octet and singlet potential, respectively. The effective theory pNRQCD is particularly suitable to study this limit, as it profits from the large-scale separation: $1/r \gg \Delta V$. For small distance r the static hybrid state can be understood to consist of a color-octet quark-antiquark state (acting as a static source in the octet representation) coupled to some ultrasoft and/or nonperturbative gluons to form an overall color-singlet state. A close relative of the static hybrid (as we will see below) is the gluelump, which consists of a static octet source (of whatever origin) attached to nonperturbative gluons forming a color-singlet state.

In order to gain a deeper understanding of the physics of the static hybrid system one can also consider how it is qualitatively affected by changing the number of dimensions from four (4D) to three (3D) or two (2D). The three-dimensional result is also important on its own. Four-dimensional thermal QCD effectively undergoes a dimensional reduction for large temperatures. Therefore, determining the renormalization group (RG) structure for the static potential in three dimensions might open the way to a resummation of logarithms at finite temperature. Three-dimensional space-time is moreover a good testing ground for renormalon issues, since the linear power divergences associated with renormalons in four dimensions become logarithmic divergences in three dimensions and can be traced back using dimensional regularization (see Ref. [1]). Last but not least the computations in less than four dimensions represent important consistency checks of the theoretical framework used to describe the 4D hybrid (and singlet) systems at short distances.

The color-octet static potential is obtained by integrating out the soft scale and has received quite some attention in the last decade. For $D = 4$, it was computed with two-loop precision in Ref. [8], the leading three-loop logarithms were obtained in Ref. [9], and the associated resummation of logarithms was carried out in Ref. [10]. The 4D two-loop result of Ref. [8] was confirmed in Ref. [11] by computing the correlator of two Polyakov loops at finite temperature, which is then related to the singlet and octet static potentials. From the terms given in Ref. [8] we have also been able to read off the general D -dimensional expression for the static octet potential, and verified it using the method developed in Ref. [11]. This expression is infrared-divergent for $D = 3$. The divergences must be canceled by the ultraviolet divergences of the ultrasoft contribution. We calculate the latter in pNRQCD for arbitrary D up to two loops and present the results in full detail on a diagram-by-diagram basis. We also compute the D -dimensional ultrasoft two-loop correction to the static

singlet energy providing an independent confirmation of the results obtained in Refs. [12,13]. The one-loop ultrasoft singlet and octet contributions are basically identical except for trivial modifications. The ultrasoft two-loop octet computation has previously been considered in Ref. [14], where a one-to-one correspondence to the singlet result in the temporal ($A_0 = 0$) gauge was claimed. This statement can be rather problematic due to the singular nature of the gluon propagator in this gauge. Actually, for $D = 4$, the anomalous dimension of the octet static potential we find differs from the one obtained using the relation between singlet and octet computation based on the $A_0 = 0$ gauge argument of Ref. [14]. In three dimensions the problem is even more acute, as the result obtained from the suggested singlet-octet correspondence does not have the correct divergence structure to make the computation renormalizable and to derive a meaningful anomalous dimension. Our explicit ultrasoft two-loop result in Feynman gauge, in contrast, does, and exactly cancels the infrared divergences of the static octet potential for $D = 3$. This represents a strong cross-check of our calculation.

The discussion in the present paper closely follows the lines of Ref. [1]. Therefore, we will skip some details by referring to that reference. The outline of the paper is as follows. In Sec. II we introduce the theoretical setup and present the bare two-loop soft and ultrasoft results for the color octet. The expressions for the relevant pNRQCD diagrams in Feynman gauge are shown in the Appendix, both for the color octet and singlet case. In Sec. III we obtain the static hybrid energy at next-to-next-to-next-to-leading logarithmic ($N^3\text{LL}$) order (apart from the soft three-loop matching constant) for $D = 4$. Sec. IV contains the complete RG-improved ultrasoft calculation up to NNLL level in three dimensions. In Sec. V we discuss the situation in two dimensions, and in Sec. VI we present our conclusions.

II. pNRQCD

Up to the next-to-leading order (NLO) in the multipole expansion (and irrespectively of the space-time dimension) the effective Lagrangian density of pNRQCD in the static limit takes the form¹

$$\begin{aligned} \mathcal{L} = & \text{Tr}\{S^\dagger(i\partial_0 - V_s(r))S + O^\dagger(iD_0 - V_o(r))O\} \\ & + gV_A(r)\text{Tr}\{O^\dagger \mathbf{r} \cdot \mathbf{E}S + S^\dagger \mathbf{r} \cdot \mathbf{E}O\} \\ & + g\frac{V_B(r)}{2}\text{Tr}\{O^\dagger\{\mathbf{r} \cdot \mathbf{E}, O\}\} + \mathcal{O}(r^2). \end{aligned} \quad (2)$$

¹In principle, to fully account for the nonperturbative contributions at $\mathcal{O}(r^2)$ we have to add the term

$$\delta\mathcal{L} = g\frac{V_C(r)}{8}r^i r^j \text{Tr}\{O^\dagger[\mathbf{D}^i \mathbf{E}^j, O]\} \quad (1)$$

to the Lagrangian in Eq. (2).

We define color-singlet and -octet fields for the quark-antiquark system by $S = S(\mathbf{r}, \mathbf{R}, t)$ and $O^a = O^a(\mathbf{r}, \mathbf{R}, t)$ respectively. $\mathbf{R} \equiv (\mathbf{x}_1 + \mathbf{x}_2)/2$ is the center position of the system. In order for S and O^a to have the proper free-field normalization in color space they are related to the fields in Eq. (2) as follows:

$$S \equiv \frac{11_c}{\sqrt{N_c}} S, \quad O \equiv \frac{T^a}{\sqrt{T_F}} O^a, \quad (3)$$

where $T_F = 1/2$ for the fundamental representation of $SU(N_c)$. All gluon fields in Eq. (2) are evaluated in \mathbf{R} and the time t , in particular, the chromoelectric field $\mathbf{E} \equiv \mathbf{E}(\mathbf{R}, t)$ and the ultrasoft covariant derivative $iD_0 O \equiv i\partial_0 O - g[A_0(\mathbf{R}, t), O]$.

In the following we will use the index “ B ” to explicitly denote bare quantities. Parameters without this index are understood to be renormalized. We will furthermore use the notation $D \equiv 4 + 2\epsilon \equiv n + 2\epsilon_n$, where $\epsilon_n = \frac{D-n}{2}$ parameterizes the (typically infinitesimal) difference to the closest integer dimension $n = 4, 3, 2$. The bare parameters of the theory are the coupling constant α_B (g_B) and the potentials $V_{\{s,o,A,B\},B}(r)$, generically denoted by V_B . The associated renormalized coefficients $\alpha(\nu)$ and $V_{\{s,o,A,B\}}(r; \nu)$ are the Wilson coefficients of the effective Lagrangian and depend on the renormalization scale (ν). They are typically fixed at a scale smaller than (or similar to) $1/r$ and larger than the ultrasoft and any other scale in the problem by matching the effective to the underlying theory, which in this case is QCD in the static limit.

In our convention α_B has integer mass dimension, $[\alpha_B] = [\alpha] = M^{4-n}$ ($[g_B^2] = M^{4-D}$), and is related to g_B by

$$\alpha_B = \frac{g_B^2 \nu^{2\epsilon_n}}{4\pi}, \quad (4)$$

where ν is the renormalization scale. It has a special status since it does not receive corrections from other Wilson coefficients of the effective theory. Therefore, it can be renormalized multiplicatively

$$\alpha_B = Z_\alpha \alpha, \quad (5)$$

where

$$Z_\alpha = 1 + \sum_{s=1}^{\infty} Z_\alpha^{(s)} \frac{1}{\epsilon_n^s}. \quad (6)$$

The RG equation of α is

$$\nu \frac{d}{d\nu} \alpha \equiv \alpha\beta(\alpha; \epsilon_n) = 2\epsilon_n \alpha + \alpha\beta(\alpha; 0). \quad (7)$$

In the limit $\epsilon_n \rightarrow 0$

$$\nu \frac{d}{d\nu} \alpha \equiv \alpha\beta(\alpha; 0) \equiv \alpha\beta(\alpha) = -2\alpha \frac{d}{d\alpha} Z_\alpha^{(1)}. \quad (8)$$

For the octet potential we employ an additive renormalization convention

$$V_{o,B} = V_o + \delta V_o. \quad (9)$$

The counterterm δV_o generally depends on the Wilson coefficients of the effective theory, i.e., on α and V , and on the number of space-time dimensions. Using the minimal subtraction (MS) renormalization scheme in $D(n)$ dimensions we define

$$\delta V_o = \sum_{s=1}^{\infty} Z_{V_o}^{(s)} \frac{1}{\epsilon_n^s}. \quad (10)$$

From the scale independence of the octet bare potential

$$\nu \frac{d}{d\nu} V_{o,B} = 0, \quad (11)$$

one obtains the RG equation of V_o . It can schematically be written as

$$\nu \frac{d}{d\nu} V_o = B(V), \quad (12)$$

$$B(V) \equiv -\left(\nu \frac{d}{d\nu} \delta V_o\right). \quad (13)$$

Note that $B(V)$ is, in general, a function of all the potentials appearing in the Lagrangian. Note as well that Eq. (12) implies that all the $1/\epsilon_n$ poles disappear once the derivative with respect to the renormalization scale is performed. This imposes some constraints on δV_o :

$$\mathcal{O}(1/\epsilon_n): B(V) = -2\alpha \frac{\partial}{\partial \alpha} Z_{V_o}^{(1)}, \quad (14)$$

$$\mathcal{O}(1/\epsilon_n^2): B(V) \frac{\partial}{\partial V} Z_{V_o}^{(1)} + \alpha \beta(\alpha) \frac{\partial}{\partial \alpha} Z_{V_o}^{(1)} + 2\alpha \frac{\partial}{\partial \alpha} Z_{V_o}^{(2)} = 0, \quad (15)$$

and so on.

A. Bare results in D dimensions

This subsection summarizes the bare results of the calculations relevant in this work. In D -dimensional momentum space we can write the bare octet potential as

$$\tilde{V}_{o,B} = \frac{1}{2N_c} g_B^2 \frac{1}{\mathbf{k}^2} \sum_{n=0}^{\infty} g_B^{2n} \mathbf{k}^{2n\epsilon} \frac{\tilde{c}_n^{(o)}(D)}{(4\pi)^{nD/2}}, \quad (16)$$

where $\tilde{c}_0^{(o)}(D) = 1$ and $\tilde{c}_1^{(o)}(D) = \tilde{c}_1^{(s)}(D)$, $\tilde{c}_1^{(s)}(D)$ being the analogously defined one-loop coefficient for the static singlet potential, see Ref. [15]. The two-loop coefficient $\tilde{c}_2^{(o)}(D)$ differs from $\tilde{c}_2^{(s)}(D)$, which was computed in Ref. [16]. We denote this difference by

$$\delta \tilde{c}_2(D) = \tilde{c}_2^{(o)}(D) - \tilde{c}_2^{(s)}(D). \quad (17)$$

$\delta \tilde{c}_2(4)$ was first obtained in Ref. [8]. The result was confirmed in Ref. [11] using thermal QCD and effective theory methods to compute the correlator of two Polyakov loops and relate it to the singlet and octet potentials. Although no explicit expression is given one can also read the D -dimensional coefficient $\delta \tilde{c}_2(D)$ off the results in Ref. [8]. Using the approach developed in Ref. [11] we have been able to confirm that expression for arbitrary D . It reads

$$\delta \tilde{c}_2^{(o)}(D) = C_A^2 \frac{\pi^3 \Gamma(\epsilon + \frac{1}{2}) \sec^2(\pi\epsilon) (\epsilon(2\epsilon + 3) \Gamma(\epsilon - \frac{1}{2}) \Gamma(2\epsilon + 1) \csc(\pi\epsilon) + \pi \Gamma(3\epsilon + \frac{3}{2}) \sec(\pi\epsilon))}{\Gamma(\frac{3}{2} - \epsilon) \Gamma(2\epsilon + 1)^2 \Gamma(3\epsilon + \frac{1}{2})}. \quad (18)$$

After Fourier transformation to position space Eq. (16) becomes (see, e.g., Ref. [17])

$$\begin{aligned} V_{o,B} &= \frac{1}{2N_c} g_B^2 \sum_{n=0}^{\infty} \frac{g_B^{2n} r^{-2(n+1)\epsilon}}{r} \frac{\tilde{c}_n^{(o)}(D)}{(4\pi)^{nD/2}} \\ &\times \frac{\Gamma[1/2 + (n+1)\epsilon]}{2^{2-2n\epsilon} \pi^{3/2+\epsilon} \Gamma[1-n\epsilon]} \\ &\equiv \frac{1}{2N_c} g_B^2 \sum_{n=0}^{\infty} \frac{g_B^{2n} c_n^{(o)}(D) r^{-2(n+1)\epsilon}}{r}. \end{aligned} \quad (19)$$

We now consider the ultrasoft bare correction to the hybrid energy at $\mathcal{O}(r^2)$ in the multipole expansion. It can be determined from the calculation of the bare octet self-energy $\Sigma_B^{us}(E)$ in terms of (1PI) loop diagrams in pNRQCD. The octet propagator including the ultrasoft corrections (but neglecting nonperturbative ones) then takes the form

$$\int dt e^{iEt} \langle O^a(t) O^b(0) \rangle \sim \frac{i}{E - V_o^B - \Sigma_B^{us}(E)}. \quad (20)$$

To extract the ultrasoft correction to the static octet energy² at $\mathcal{O}(r^2)$ it is sufficient to set $E = V_o^B|_{\mathcal{O}(r^0)}$ in the bare expression for the self-energy,³ i.e., $\delta E_{o,B}^{us} = \Sigma_B^{us}(E = V_o^B) + \mathcal{O}(r^3)$.

The one-loop result equals the analogous singlet correction [9,18,19] except for a change of the color factor and the replacement $\Delta V \equiv V_o - V_s \rightarrow -\Delta V$

²This correction is understood to be complex as long as we do not explicitly disentangle the real part associated with the physical energy and the imaginary part associated with the decay width, see Sec. II B.

³In order to consistently treat the 3D infrared divergence at $\mathcal{O}(r^0)$ (which is canceled by an ultraviolet divergence of non-perturbative origin) in dimensional regularization it is important to set $E = V_o^B$ before performing the loop integrations of the self-energy diagrams in this ‘‘bare’’ approach. As a check we have computed the corresponding off shell diagrams and explicitly performed the renormalization of the potentials and the octet field. Finally taking the ‘‘renormalized’’ on shell limit $E \rightarrow V_o$ gives the same result for the ultrasoft correction to the static hybrid energy.

$$\begin{aligned} \delta E_{o,B}^{us}(1\text{-loop}) &= -g^2 \frac{1}{2N_c} V_A^2 (1 + \epsilon) \\ &\times \frac{\Gamma[2 + \epsilon] \Gamma[-3 - 2\epsilon]}{\pi^{2+\epsilon}} \mathbf{r}^2 (-\Delta V_B)^{3+2\epsilon}. \end{aligned} \quad (21)$$

In Feynman gauge it comes from the pNRQCD diagram shown in Fig. 1. The two-loop expression is new and reads

$$\begin{aligned} \delta E_{s,B}^{us}(2\text{-loop}) &= g^4 \frac{1}{2N_c} C_A V_A^2 \Gamma[-3 - 4\epsilon] \frac{1}{(2\pi)^2} \frac{1}{4\pi^{2+2\epsilon}} \\ &\times \Gamma^2[1 + \epsilon] [g(\epsilon) - (1 + 2\epsilon)g_1(\epsilon) \\ &+ g_o(\epsilon)] \mathbf{r}^2 (-\Delta V_B)^{3+4\epsilon}, \end{aligned} \quad (22)$$

$$g_o(\epsilon) = \frac{2^{-4\epsilon-5} (2\epsilon + 3)(2\epsilon^2 + 2\epsilon + 1) \Gamma(-2\epsilon - \frac{7}{2}) \Gamma^2(-\epsilon - \frac{3}{2}) \Gamma(2\epsilon + \frac{9}{2})}{\epsilon \Gamma(-4\epsilon - 3) \Gamma^2(\epsilon + 1)}. \quad (25)$$

It is the central result of this work. In the Appendix we separately present the analytic expressions for the two-loop self-energy diagrams that contribute to Eq. (22) in Feynman gauge. We have also computed the corresponding two-loop expressions for the color singlet, which provide an explicit confirmation of the result obtained in Refs. [12,13] and can also be found in the Appendix.

The computation of the ultrasoft octet self-energy has also been addressed in Ref. [14], where it was argued that the result could be related to the singlet one (which is gauge-invariant) in the temporal ($A_0 = 0$) gauge. In order to reach this conclusion, the effects involving the gluonic content of the hybrid in the soft and ultrasoft computation were neglected (see Eq. 17 in Ref. [14]). Then, after choosing the $A_0 = 0$ gauge, it is indeed easy to see that all nonvanishing self-energy diagrams at $\mathcal{O}(r^2)$ are, apart from the color factor and the sign of ΔV , the same for singlet and octet. Requiring gauge invariance of the final outcome one thus would obtain

$$\begin{aligned} V_A^2 \frac{T_F}{(D-1)(N_c^2-1)N_c} \mathbf{r}^2 \\ \times \int_0^\infty dt e^{t\Delta V_B} \langle vac | g \mathbf{E}_E^a(t) \phi(t, 0)_{ab}^{\text{adj}} g \mathbf{E}_E^b(0) | vac \rangle \end{aligned} \quad (26)$$

for the bare ultrasoft contribution to the static octet energy (in Euclidean space-time) at $\mathcal{O}(r^2)$ and to all orders in α .

At $\mathcal{O}(\alpha)$ (one-loop) Eq. (26) agrees with Eq. (21). However, at $\mathcal{O}(\alpha^2)$ (two-loop) Eqs. (22) and (26) are different, since $g_o(\epsilon) = 0$ in Eq. (26). This is in contradiction

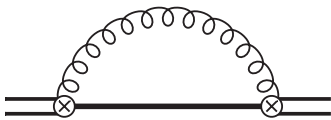


FIG. 1. One-loop octet self-energy diagram at $\mathcal{O}(r^2)$.

where

$$\begin{aligned} g(\epsilon) &= \frac{2\epsilon^3 + 6\epsilon^2 + 8\epsilon + 3}{\epsilon(2\epsilon^2 + 5\epsilon + 3)} \\ &- \frac{2\epsilon \Gamma(-2\epsilon - 2) \Gamma(-2\epsilon - 1)}{(2\epsilon + 3) \Gamma(-4\epsilon - 3)}, \end{aligned} \quad (23)$$

$$\begin{aligned} g_1(\epsilon) &= \frac{6\epsilon^3 + 17\epsilon^2 + 18\epsilon + 6}{\epsilon^2(2\epsilon^2 + 5\epsilon + 3)} + \frac{4(\epsilon + 1)n_f T_f}{\epsilon(2\epsilon + 3)N_c} \\ &+ \frac{2(\epsilon^2 + \epsilon + 1) \Gamma(-2\epsilon - 2) \Gamma(-2\epsilon - 1)}{\epsilon(2\epsilon + 3) \Gamma(-4\epsilon - 3)}, \end{aligned} \quad (24)$$

with our nonzero result for $g_o(\epsilon)$ in Eq. (25), which was obtained by an explicit two-loop calculation in Feynman gauge (see App. A 1). The existence of a nonzero value for g_o is required on a theoretical basis in Sec. IV. Otherwise $\delta E_{o,B}^{us}$ would not produce the right divergence structure in $D = 3$ dimensions to make the theory renormalizable nor lead to a consistent RG equation for the potential, as there would be some divergences proportional to logarithms of the ultrasoft scale left after the subtraction of the one-loop subdivergences. Our explicit computation, in contrast, is consistent with the renormalizability of the theory, and moreover perfectly cancels the infrared divergences of the soft computation. This represents a strong check of our computation. In Sec. III we will see, that with our result in Eq. (22) we also obtain a different NLL anomalous dimension of the octet potential V_o in four dimensions compared to Ref. [14]. The corresponding cancelation of soft infrared and ultrasoft ultraviolet divergences in four dimensions can however not be checked, since the required soft four-loop result is not known at present.

Actually, the fact that there are problems with the naive usage of $A_0 = 0$ gauge becomes apparent at several points of the calculation. In pNRQCD this already happens at $\mathcal{O}(r^0)$ in the multipole expansion for $D = 3$. At this order the bare soft result for V_o is infrared-divergent (cf. Equation (47)), but in the $A_0 = 0$ gauge there are no ultrasoft octet self-energy diagrams to cancel that divergence. It is also remarkable that in this gauge the spacelike strings of the rectangular Wilson loop contribute to the static potential, whereas in Feynman gauge they are expected to be negligible.⁴ This can be easily visualized for $D = 2$, since then time and space coordinates can be

⁴In Ref. [15] it was explicitly shown that the spacelike strings in Feynman gauge do not contribute to the 4D static singlet potential through two loops.

symmetrically interchanged (in the Euclidean), showing that in the $A_0 = 0$ gauge the contribution to the singlet potential comes exclusively from the spacelike strings. Actually that is true for arbitrary D at tree level. For a detailed analysis at one loop see, e.g., Ref. [20] and the references therein. Overall, we think it would be very interesting to quantitatively study the effect of the asymptotic gluonic degrees of freedom on the static singlet and hybrid energy in the $A_0 = 0$ gauge.

To the above discussion one should also add the $\mathcal{O}(r^2)$ ultrasoft contributions from octet self-energy diagrams without singlet-octet transitions. They are proportional to V_B^2 or V_C and may have logarithmic ultraviolet divergences, relevant for the determination of the counterterm (δV_o) and the anomalous dimension of V_o . These divergences can, however, not be proportional to positive powers of ΔV , because the ultrasoft scale ΔV does not appear in the diagrams with on shell external fields, and the corresponding contributions vanish in dimensional regularization. Therefore they cannot solve the $A_0 = 0$ gauge problem described above, since the first discrepancies between our two-loop result in Eq. (22) and the $\mathcal{O}(\alpha^2)$ term in Eq. (26) are linear in ΔV for $D = 3$ and cubic for $D = 4$.

B. Observables

At short distances we approximate the bare propagator of a generic hybrid system as follows:

$$\int dt e^{iEt} \langle H^a O^a(t) H^b O^b(0) \rangle \sim \frac{i}{E - V_o^B - \Sigma_B(E) + i\epsilon}. \quad (27)$$

H^a stands for the gluonic content of the hybrid ($H^a = B^a, E^a$, etc.) and $\Sigma_B(E)$ is the self-energy of the system expressed in terms of bare parameters (potentials, couplings). It accounts for effects at the ultrasoft and the nonperturbative scale.

Let us now construct some observables. In this paper we focus on the hybrid static energy and decay width. Their definition can become gauge-dependent for unstable particles. Fortunately this is not so in our case, since both the on shell and pole definition (see, e.g., Ref. [21] for the respective definitions) are identical at the precision of our computation. Close to the resonance we can expand $\Sigma_B(E)$ around the complex pole position $E_H + i\Gamma_H/2$ of the hybrid propagator in Eq. (27). In this paper we will consider ultrasoft self-energy contributions Σ_B^{us} up to two-loop level and NLO in the multipole expansion. It is therefore sufficient to approximate $\Sigma_B(E_H + i\Gamma_H/2)$ by $\Sigma_B(E_0)$ and set $E_0 = V_o^B|_{\mathcal{O}(\alpha^2)}$ for $D = 4$ and $E_0 = V_o^B|_{\mathcal{O}(\alpha)}$ for $D = 3$. Note that in the latter case $\mathcal{O}(\alpha^2)$ terms would produce subleading $\mathcal{O}(r^3)$ corrections to the energy and decay width.

We now choose to shift the energy origin by adding $-2m_B$, i.e., $E \rightarrow E - 2m_B$. The term $2m_B = 2m + 2\delta m_{(\text{soft})}$ will allow to absorb ultraviolet soft divergences

(into $\delta m_{(\text{soft})}$) that may appear in the computation. m has a clear physical meaning in the case of heavy quarkonium: it represents the heavy quark (pole) mass. At NLO in the multipole expansion we thus have

$$\frac{1}{E - 2m_B - V_o^B - \Sigma_B(E - 2m_B)} = \frac{Z_H}{E - E_H + i\Gamma_H/2} + \mathcal{O}((E - E_H + i\Gamma_H/2)^0), \quad (28)$$

where Z_H is the normalization of the hybrid propagator and

$$E_H - i\Gamma_H/2 = 2m_B + V_o^B + \Sigma_B(E_0) \quad (29)$$

is the complex position of the hybrid propagator pole. It can be understood to be composed of contributions from different energy regions. After factorization we can express it in the following way (either bare or renormalized):

$$\begin{aligned} E_H(r) - i\Gamma_H/2 &= 2m_B + V_{o,B} + \delta E_{o,B}^{us} + \delta E_{H,B}^{np} \\ &= 2m_{\overline{\text{MS}}}(\nu_s) + V_{o,\overline{\text{MS}}}(r; \nu_s, \nu) \\ &\quad + \delta E_{o,\overline{\text{MS}}}^{us}(\nu, \nu_{np}) + \delta E_{H,\overline{\text{MS}}}^{np}(\nu_{np}). \end{aligned} \quad (30)$$

The subscript $\overline{\text{MS}}$ of $m_{\overline{\text{MS}}}(\nu_s)$ denotes the scheme for possible hard infrared divergence subtractions (canceling the ultraviolet soft divergences), e.g., in three dimensions, and should not be confused with the ultraviolet (hard) renormalization scheme for the mass. In four dimensions we have $m_{\overline{\text{MS}}}(\nu_s) = m_{\text{OS}}$, where m_{OS} is the usual on shell mass of the heavy quarks. Similarly $V_{o,\overline{\text{MS}}}(r; \nu_s, \nu) \rightarrow V_{o,\overline{\text{MS}}}(r; \nu)$ in four dimensions, where V_o is the octet potential and encodes the effects due to soft degrees of freedom. δE_o^{us} and δE_H^{np} encode the effects due to the ultrasoft and nonperturbative degrees of freedom, respectively. Each term depends on the respective factorization/renormalization scales: ν_s separates the hard (m) from the soft scale ($1/r$), ν separates the soft from the ultrasoft scale (ΔV) and ν_{np} separates the ultrasoft from the nonperturbative scale (Λ_{QCD}). For real factorization scales ν_s, ν, ν_{np} , the decay width Γ_H is contained in the terms δE_o^{us} and δE_H^{np} . We give more details in the next sections.

III. RESULTS FOR $D = 4$

In four dimensions the renormalized coupling constant α satisfies a nontrivial RG equation. We will work here in the $\overline{\text{MS}}$ renormalization scheme, which is related to the MS scheme by a redefinition of the renormalization scale: $\nu_{(\text{MS})} \rightarrow \nu_{(\overline{\text{MS}})} c_{\overline{\text{MS}}}^{-1}$, where $c_{\overline{\text{MS}}} = e^{1/2(\ln(4\pi) - \gamma_E)}$. The counterterm coefficient

$$Z_\alpha^{(1)} = \frac{\alpha}{4\pi} \beta_0 + \frac{1}{2} \frac{\alpha^2}{(4\pi)^2} \beta_1 + \dots \quad (31)$$

is understood in the MS scheme, but the form of the anomalous dimension of α ,

$$\alpha\beta(\alpha) = -2\alpha\left(\beta_0\frac{\alpha}{4\pi} + \beta_1\frac{\alpha^2}{(4\pi)^2} + \dots\right) \quad (32)$$

is the same in both schemes, MS and $\overline{\text{MS}}$. The constants $\beta_0 = \frac{11}{3}C_A - \frac{4}{3}T_F n_f$ and $\beta_1 = \frac{34}{3}C_A^2 - 4C_F T_F n_f - \frac{20}{3}C_A T_F n_f$ are the standard $\overline{\text{MS}}$ (MS) one- and two-loop coefficients of the QCD beta function.

We also have to consider possible soft and ultrasoft corrections to V_A . They were obtained in Ref. [10] with leading logarithmic (LL) accuracy, in Ref. [13] with NLO accuracy and in Ref. [14] with NLL accuracy. The outcome is

$$\nu \frac{d}{d\nu} V_A = 0 + \mathcal{O}(\alpha^3) \quad (33)$$

for the anomalous dimension and $V_A = 1 + \mathcal{O}(\alpha^2)$ for the initial matching condition. We conclude that for the precision of our calculation we can use $V_A = 1$.

The counterterms of the octet potential, which subtract the ultraviolet divergences from the result of our ultrasoft two-loop computation in Eq. (22) in four dimensions, read

$$Z_{V_o}^{(1)} = -r^2 \Delta V^3 \frac{1}{2N_c} V_A^2 \left[\frac{\alpha}{3\pi} + \frac{\alpha^2 [C_A(47 - 12\pi^2) - 10T_F n_f]}{108\pi^2} \right], \quad (34)$$

$$Z_{V_o}^{(2)} = -r^2 \Delta V^3 \frac{1}{2N_c} V_A^2 \frac{2}{3} \beta_0 \frac{\alpha^2}{(4\pi)^2}. \quad (35)$$

The latter expression comes from the $1/\epsilon_4^2$ pole of the two-loop result and from the α^2/ϵ_4 divergence of α_B in the divergent term of the one-loop self-energy.

From Eqs. (34) and (35) we can derive the RG equation of the octet potential at two-loop order.⁵ We find

$$\nu \frac{d}{d\nu} V_{o,\overline{\text{MS}}} = r^2 \Delta V^3 \frac{1}{2N_c} V_A^2 \left[\frac{2\alpha}{3\pi} + \frac{\alpha^2 [C_A(47 - 12\pi^2) - 10T_F n_f]}{27\pi^2} + \mathcal{O}(\alpha^3) \right]. \quad (36)$$

This result holds in any momentum-independent renormalization scheme that is related to the MS scheme by a (D -independent) redefinition of the renormalization scale ν . Solving the RG equation we can write the RG-improved static octet potential as

$$V_{o,\overline{\text{MS}}}(r; \nu) = V_{o,\overline{\text{MS}}}^{i.c.}(r; \nu_i) + \delta V_{o,\overline{\text{MS}}}^{\text{RG}}(r; \nu_i, \nu), \quad (37)$$

where

⁵In our counting this translates to N³LL order, because ΔV comes at least with one power of α .

$$\delta V_{o,\overline{\text{MS}}}^{\text{RG}}(r; \nu_i, \nu) = -\frac{1}{2N_c} V_A^2 r^2 (\Delta V)^3 \frac{2\pi}{\beta_0} \left\{ \frac{2}{3\pi} \ln \frac{\alpha(\nu)}{\alpha(\nu_i)} - (\alpha(\nu) - \alpha(\nu_i)) \left(\frac{8}{3(4\pi)^2} \frac{\beta_1}{\beta_0} - \frac{C_A(47 - 12\pi^2) - 10T_F n_f}{27\pi^2} \right) \right\} \quad (38)$$

describes the ultrasoft RG evolution of V_o and the initial matching condition at the (soft) scale ν_i is given by

$$V_{o,\overline{\text{MS}}}^{i.c.}(r; \nu_i) = \frac{1}{2N_c} \frac{\alpha(\nu_i)}{r} \sum_{n=0}^3 \left(\frac{\alpha(\nu_i)}{4\pi} \right)^n a_n^{(o)}(r; \nu_i) \quad (39)$$

with coefficients ($a_0^{(o)}(r; \nu_i) = 1$)

$$\begin{aligned} a_1^{(o)}(r; \nu_i) &= a_1 + 2\beta_0 \ln(\nu_i e^{\gamma_E} r), \\ a_2^{(o)}(r; \nu_i) &= a_2^{(o)} + \frac{\pi^2}{3} \beta_0^2 + (4a_1 \beta_0 + 2\beta_1) \ln(\nu_i e^{\gamma_E} r) \\ &\quad + 4\beta_0^2 \ln^2(\nu_i e^{\gamma_E} r), \\ a_3^{(o)}(r; \nu_i) &= a_3^{(o)} + a_1 \beta_0^2 \pi^2 + \frac{5\pi^2}{6} \beta_0 \beta_1 + 16\zeta_3 \beta_0^3 \\ &\quad + \left(2\pi^2 \beta_0^3 + 6a_2^{(o)} \beta_0 + 4a_1 \beta_1 + 2\beta_2 \right. \\ &\quad \left. + \frac{16}{3} C_A^3 \pi^2 \right) \ln(\nu_i e^{\gamma_E} r) + (12a_1 \beta_0^2 \\ &\quad + 10\beta_0 \beta_1) \ln^2(\nu_i e^{\gamma_E} r) + 8\beta_0^3 \ln^3(\nu_i e^{\gamma_E} r). \end{aligned} \quad (40)$$

Explicit expressions for a_1 and $a_2^{(o)}$ can be found in the literature [16,22]. The constant $a_3^{(o)}$ is still unknown and represents the only missing piece in the N³LL result for the Wilson coefficient V_o , Eq. (37). Note that this expression is real for positive renormalization and matching scales ν and ν_i .

We can now determine the complex pole of the hybrid (octet) propagator in four dimensions. Up to nonperturbative effects, which we neglect here, it is given by

$$E_H(r) - i \frac{\Gamma_H}{2} \simeq E_o(r) - i \frac{\Gamma_o}{2} \simeq V_{o,\overline{\text{MS}}}^{i.c.}(r; 1/r) + \delta V_{o,\overline{\text{MS}}}^{\text{RG}}(r; 1/r, \Delta V) + \delta E_{o,\overline{\text{MS}}}^{\text{us}}(\Delta V). \quad (41)$$

Therefore, besides the octet potential, we also need the ultrasoft correction in the $\overline{\text{MS}}$ scheme, which we obtain from our bare two-loop result in Eq. (22) after subtraction of the divergences. It reads

$$\begin{aligned} \delta E_{o,\overline{\text{MS}}}^{us}(\nu) = & \frac{1}{2N_c} r^2 (-\Delta V)^3 V_A^2 \left[-\frac{\alpha}{9\pi} \left(6 \ln \left[\frac{-\Delta V}{\nu} \right] + 6 \ln 2 - 5 \right) + \frac{\alpha^2}{108\pi^2} \left(18\beta_0 \ln^2 \left[\frac{-\Delta V}{\nu} \right] - 6(N_c(13 - 8\pi^2) \right. \right. \\ & - 2\beta_0(-5 + 3 \ln 2)) \ln \left[\frac{-\Delta V}{\nu} \right] - 2N_c(-84 + 39 \ln 2 - 8\pi^2(-2 + 3 \ln 2) + 72\zeta(3)) \\ & \left. \left. + \beta_0(67 + 3\pi^2 - 60 \ln 2 + 18 \ln^2 2) \right) \right]. \end{aligned} \quad (42)$$

Note that this object is complex for positive ν . Since the octet potential is real, the imaginary part of $\delta E_{o,\overline{\text{MS}}}^{us}(\nu)$ directly gives the decay width. Choosing $\nu = \Delta V$ in Eq. (42) we find for the decay width of the hybrid (octet) system

$$\Gamma_H = \frac{4}{3} \alpha(\Delta V) \frac{1}{2N_c} r^2 \Delta V^3 \left[1 + \frac{\alpha(\Delta V)}{12\pi} (N_c(13 - 8\pi^2) - 2\beta_0(3 \ln 2 - 5)) + \mathcal{O}(\alpha^2) \right] + \mathcal{O}(r^4), \quad (43)$$

where we have to use the NLO expression for the ultrasoft scale ΔV :

$$\Delta V(r; \nu_i) = \frac{C_A}{2} \frac{\alpha(\nu_i)}{r} \sum_{n=0}^1 \left(\frac{\alpha(\nu_i)}{4\pi} \right)^n a_n^{(o)}(r; \nu_i), \quad (44)$$

which is scheme- and factorization-scale-independent at this order.

For E_H we need the real part of $\delta E_{o,\overline{\text{MS}}}^{us}(\nu)$:

$$\begin{aligned} E_H(r) = & V_{o,\overline{\text{MS}}}^{i.c.}(r; 1/r) + \delta V_{o,\overline{\text{MS}}}^{\text{RG}}(r; 1/r, \Delta V) \\ & + \text{Re} \left[\delta E_{o,\overline{\text{MS}}}^{us}(\Delta V) \right]. \end{aligned} \quad (45)$$

Replacing the first two terms in this equation by Eq. (37) and using the leading term of Eq. (42) in the last term, we reach N³LL accuracy. Equation (42) also provides the subleading ultrasoft correction to the hybrid energy relevant at N⁴LL order.

We will not attempt to compare our results with 4D lattice simulations. This would require the incorporation of nonperturbative effects and the proper treatment of renormalons, which goes beyond the scope of this paper.

IV. RESULTS FOR $D = 3$

The derivation of the static octet potential in three dimension is quite analogous to the color-singlet case, which has been discussed in detail in Ref. [1]. We therefore focus on the novel aspects of the analysis for the octet (hybrid) system in this section and refer to Ref. [1] for a more careful account on the universal issues related to the 3D static potential.

As argued in Ref. [1], the coefficients $V_{A/B}$ are not renormalized at $\mathcal{O}(r^0)$, i.e.

$$Z_{A/B} = 1 + \mathcal{O}(r). \quad (46)$$

The reason is that the potentials and α have to appear perturbatively (with positive powers) in the counterterms, otherwise the renormalizability of the theory at leading order of the multipole expansion would be spoiled. Moreover, we can set $V_A = V_B = 1$ (just like in 4D), as $\mathcal{O}(\alpha)$ soft corrections would be multiplied by factors of r and would move us away from the precision of V_o aimed for at this paper. Actually, from inspection of the possible diagrams that will contribute at the soft scale, we know that $V_{A/B} = 1 + \mathcal{O}(\alpha^2)$ [13].

From the bare soft computation we can completely fix δV_o through $\mathcal{O}(r^2)$. We obtain (for the ultrasoft counterterms)

$$\begin{aligned} Z_{V_o}^{(1)} = & \frac{C_A}{2} \alpha + r^2 \Delta V^2 \alpha \left(C_F - \frac{C_A}{2} \right) \frac{1}{4} + r^2 \Delta V C_A \alpha^2 \left(C_F - \frac{C_A}{2} \right) \frac{1}{2} - r^2 \alpha^3 \left(C_F - \frac{C_A}{2} \right) \\ & \times \frac{(13\pi^2 - 2304)C_A^2 + 8(19\pi^2 + 144)C_A T_F n_f - 48T_F n_f(4(\pi^2 - 10)C_F + \pi^2 T_F n_f)}{2304}, \end{aligned} \quad (47)$$

$$Z_{V_o}^{(2)} = r^2 \Delta V \alpha^2 \left(C_F - \frac{C_A}{2} \right) C_A \frac{1}{8} + r^2 \alpha^3 \left(C_F - \frac{C_A}{2} \right) C_A^2 \frac{1}{12}, \quad (48)$$

$$Z_{V_o}^{(3)} = r^2 \alpha^3 \left(C_F - \frac{C_A}{2} \right) C_A^2 \frac{1}{48}, \quad (49)$$

$$Z_{V_o}^{(n)} = 0 \quad \forall n > 3. \quad (50)$$

The soft calculation is organized in powers of αr . The tree-level computation gives the first term in Eq. (47) (once the ultraviolet divergences of the soft one-loop heavy quark self-energies have been subtracted). The soft one-loop contribution to the octet potential is infrared safe. The

two-loop computation leads to the remaining terms. These results are exact at $\mathcal{O}(r^2)$.

The fact that one can renormalize the potential with a finite number of terms at a given order in the multipole expansion, i.e., $Z_{V_o}^{(n)} = 0$ for $n > 3$ at $\mathcal{O}(r^2)$, reflects the super-renormalizability of the theory in three dimensions.

The infrared divergences of the bare potential have to cancel the ultraviolet ones of the ultrasoft computation. We can then use the explicit one- and two-loop ultrasoft contributions to partially check the structure of the octet potential counterterm in Eqs. (47)–(49).

At leading order in the multipole expansion the octet field has a residual interaction with the ultrasoft gluon field. The octet potential receives an ultraviolet-divergent correction from the one-loop self-energy diagram shown in Fig. 2, which yields the first term in Eq. (47). This is the only term at $\mathcal{O}(r^0)$, since higher loop diagrams cannot contribute because the coupling and the potentials have to appear perturbatively (with positive powers) in the Z 's. Since α has dimensions of mass, the potentials would appear with negative powers in multiloop diagrams at $\mathcal{O}(r^0)$. This is not allowed by renormalizability. By the very same reason the octet field does not require renormalization at $\mathcal{O}(r^0)$:

$$Z_o = 1 + \mathcal{O}(r^2). \quad (51)$$

We now consider higher orders in the multipole expansion. At two soft loops in dimensional regularization infrared poles up to $\mathcal{O}(1/\epsilon_3^3)$ appear. The ultrasoft computation in pNRQCD yields the following results for the counterterms:

- (1) The second term in Eq. (47) comes from the $1/\epsilon_3$ divergence of the ultrasoft one-loop correction in Eq. (21). It is scheme-independent and fixes, together with Eqs. (14) and (15), the first term of Eq. (48) and (49). It would also be possible to compute these $1/\epsilon_3^2$ and $1/\epsilon_3^3$ divergences directly. The $1/\epsilon_3^2$ term is confirmed by the ultrasoft two-loop result in Eq. (22). The $1/\epsilon_3^3$ term would however require an ultrasoft three-loop calculation, which has not been performed yet.
- (2) The third term in Eq. (47) follows from the remaining $1/\epsilon_3$ divergence in the ultrasoft two-loop correction, Eq. (22), once all subdivergences (associated with the $\mathcal{O}(r^0)$ octet potential) have been subtracted. This result combined with Eq. (15) then fixes the second term of Eq. (48).

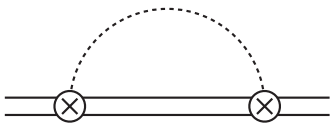


FIG. 2. One-loop contribution to the octet propagator at $\mathcal{O}(r^2)$. The dotted line represents the A^0 field.

In summary: the explicit ultrasoft computation allows us to check the three first terms of Eq. (47) and the first term of Eq. (48). We find perfect agreement. The use of Eqs. (14) and (15) allows us to completely check Eq. (48) and (49). We also find perfect agreement. Note that this can be understood as a nontrivial cross-check of two independent determinations of these terms.

The only remaining term is the last one in Eq. (47). This term would be canceled by the sum of the ultrasoft ultraviolet divergences from three-loop octet self-energy diagrams with and without singlet-octet transition. The latter diagrams are proportional to V_B^2 or V_C and scaleless (for on shell external fields), as they are insensitive to the ultrasoft scale ΔV . Therefore, the associated RG evolution will run down to the nonperturbative scale α . This is not so for the ultrasoft diagrams with singlet-octet transition, which get infrared regulated by the ultrasoft scale ΔV correspondingly producing logarithms of ΔV . Without an explicit ultrasoft three-loop calculation we are not able to distinguish among the divergences from the two types of diagrams nor check the last term of Eq. (47), which has been deduced from the soft result. Therefore, we turn the problem around and use the latter to fix the ultraviolet divergences of the ultrasoft three-loop computation for the next sections.

A. pNRQCD RG

We can now deduce the RG equations of V_o . Again the discussion is similar to the one of Ref. [1] to which we refer for extra details.

At leading order in the multipole expansion the RG equation of the octet potential reads

$$\nu \frac{d}{d\nu} V_o = -C_A \alpha. \quad (52)$$

The running of ΔV at this order reads (including the tree-level matching condition)

$$\Delta V_X(r; \nu) = -\alpha C_A \ln(r\nu d_X) + \mathcal{O}(\epsilon), \quad (53)$$

where X stands for the factorization scheme (e.g., MS or $\overline{\text{MS}}$)

$$\begin{aligned} d_{\text{MS}} &= e^{\gamma_E/2} \sqrt{\pi} \approx 2.36546, \\ d_{\overline{\text{MS}}} &= d_{\text{MS}} c_{\overline{\text{MS}}}^{-1} = e^{\gamma_E/2} \approx 0.890536, \end{aligned} \quad (54)$$

with $c_{\overline{\text{MS}}} = e^{1/2(\ln(4\pi) - \gamma_E)}$.

Now we consider the subleading contributions to V_o . The *complete* anomalous dimension of the static octet potential through $\mathcal{O}(r^2)$ takes the form

$$\begin{aligned} \nu \frac{d}{d\nu} V_o &= B(V_o) \\ &= -C_A \alpha + \sum_{n=0}^2 C_F \alpha V_A^2 r^2 \alpha B_n (C_A \alpha)^n (\Delta V)^{2-n}, \end{aligned} \quad (55)$$

and is explicitly given by

$$\begin{aligned} \nu \frac{d}{d\nu} V_o = & -C_A \alpha - r^2 \Delta V_X^2 \alpha \left(C_F - \frac{C_A}{2} \right) V_A^2 \frac{1}{2} - 2r^2 \Delta V_X C_A \alpha^2 \left(C_F - \frac{C_A}{2} \right) V_A^2 + r^2 \alpha^3 \left(C_F - \frac{C_A}{2} \right) \\ & \times V_A^2 \frac{(13\pi^2 - 2304)C_A^2 + 8(19\pi^2 + 144)C_A T_F n_f - 48T_F n_f (4(\pi^2 - 10)C_F + \pi^2 T_F n_f)}{384}. \end{aligned} \quad (56)$$

The form of the anomalous dimension in Eq. (56) is invariant under scheme transformations that amount to a redefinition of the renormalization scale: $\nu \rightarrow \nu c_X^{-1}$, where c_X is a ϵ_3 -independent constant ($c_{\text{MS}} = 1$). We indicate renormalized quantities in this class of momentum-independent renormalization schemes by an index “ X ” ($X = \text{MS}, \overline{\text{MS}}, \text{etc.}$) as, e.g., ΔV_X in the above equations.

By solving the RG equations in the MS scheme we obtain the static octet potential with $\mathcal{O}(\alpha^3 r^2)$ precision:

$$V_{o,\text{MS}}(r; \nu_s, \nu) = V_{o,\text{MS}}^{i.c.}(r; \nu_s, \nu_i) + \delta V_{o,\text{MS}}^{\text{RG}}(r; \nu_i, \nu), \quad (57)$$

where

$$\begin{aligned} \delta V_{o,\text{MS}}^{\text{RG}}\left(r; \frac{1}{r}, \nu\right) = & -C_A \alpha \ln(\nu r) + \left(C_F - \frac{C_A}{2} \right) r^2 \alpha^3 \left\{ -\frac{1}{6} C_A^2 \ln^3(r\nu) - \frac{1}{4} C_A^2 [-4 + \gamma_E + \ln\pi] \ln^2(r\nu) \right. \\ & + \left[C_A^2 \left(\frac{13\pi^2}{384} + \frac{1}{8} (-\gamma_E^2 + 8\gamma_E - 48 + (8 - 2\gamma_E) \ln\pi - \ln^2\pi) \right) + n_f T_F \left(C_A \left(3 + \frac{19\pi^2}{48} \right) \right. \right. \\ & \left. \left. + C_F \left(5 - \frac{\pi^2}{2} \right) \right) - (n_f T_F)^2 \frac{\pi^2}{8} \right] \ln(r\nu) \left. \right\} \end{aligned} \quad (58)$$

is the running and

$$\begin{aligned} V_{o,\text{MS}}^{i.c.}\left(r; \nu_s, \frac{1}{r}\right) = & C_F \alpha \ln(r^2 \nu_s^2 \pi e^{\gamma_E}) - \frac{C_A}{2} \alpha \ln(\pi e^{\gamma_E}) + \frac{\pi}{4} \left(C_F - \frac{C_A}{2} \right) (7C_A - 4n_f T_F) \alpha^2 r \\ & + \left(C_F - \frac{C_A}{2} \right) r^2 \alpha^3 \left\{ C_A^2 \left[\frac{\pi^2}{2304} (39\gamma_E - 715 + 564 \ln 2 + 39 \ln\pi) + \frac{25}{3} - \frac{31}{24} \zeta(3) \right. \right. \\ & - \frac{1}{48} (\gamma_E + \ln\pi) (144 - 12\gamma_E + \gamma_E^2 + \ln^2\pi - 12 \ln\pi + 2\gamma_E \ln\pi) \left. \right] + n_f T_F C_F \left[\frac{\pi^2}{24} (15 - 6\gamma_E) \right. \\ & - 8 \ln 2 - 6 \ln\pi + \frac{1}{2} (5\gamma_E - 8 + 5 \ln\pi) \left. \right] + n_f T_F C_A \left[\frac{\pi^2}{288} (57\gamma_E - 97 + 12 \ln 2 + 57 \ln\pi) \right. \\ & \left. \left. + \frac{1}{6} (9\gamma_E - 22 + 9 \ln\pi + 10 \zeta(3)) \right] + (n_f T_F)^2 \left[\frac{\pi^2}{48} (7 - 3\gamma_E - \ln(16\pi^3)) \right] \right\} \end{aligned} \quad (59)$$

is the initial matching condition. Note that the tree-level and one-loop matching conditions have been included. The tree-level result depends on the factorization scale $\nu_s \gg 1/r$, which separates the hard and the soft regime. The dependence on this factorization scale cancels the infrared scale dependence of the mass, cf. Equation (30).

We stress that we have obtained the *exact* contributions to the static potential through $\mathcal{O}(r^2)$. There is nothing left. Moreover, by setting $\nu \sim \Delta V$ large logarithms are resummed

up to the scale ΔV , i.e., we have determined all logarithms of $r\Delta V$ at $\mathcal{O}(\alpha^3 r^2)$.

The above results have been presented in the MS scheme. We can easily transform them to another momentum-independent scheme by redefining the renormalization scale. In particular, if we write the expression in terms of ΔV , most of the scheme dependence gets encapsulated in ΔV . For instance, Eq. (58) can be reexpressed as

$$\begin{aligned} \delta V_{o,\text{MS}}^{\text{RG}}\left(r; \frac{1}{r}, \nu\right) = & \Delta V_X(r; \nu) - \Delta V_X(r; 1/r) + \frac{1}{6} \left(\frac{C_F}{C_A} - \frac{1}{2} \right) r^2 \alpha V_A^2 (\Delta V_X^3(r; \nu) - \Delta V_X^3(r; 1/r)) + r^2 C_A \alpha^2 \left(\frac{C_F}{C_A} - \frac{1}{2} \right) V_A^2 (\Delta V_X^2(r; \nu) \\ & - \Delta V_X^2(r; 1/r)) - r^2 \alpha^2 \left(\frac{C_F}{C_A} - \frac{1}{2} \right) V_A^2 \left(\frac{(13\pi^2 - 2304)C_A^2 + 8(19\pi^2 + 144)C_A T_F n_f - 48T_F n_f (4(\pi^2 - 10)C_F + \pi^2 T_F n_f)}{384} \right) \\ & \times (\Delta V_X(r; \nu) - \Delta V_X(r; 1/r)). \end{aligned} \quad (60)$$

B. Ultrasoft contributions (δE_o^{us})

The general structure of the ultrasoft contribution δE_o^{us} at $\mathcal{O}(r^2)$ is

$$\delta E_o^{us}(\nu, \nu) = -\frac{1}{2N_c} \alpha V_A^2 r^2 \Delta V^2 \sum_{n=0}^{\infty} \left(\frac{C_A \alpha}{\Delta V} \right)^n \times \sum_{s=0}^{n+1} c_{n,s} \ln^s \left[\frac{\Delta V}{\nu} \right], \quad (61)$$

where we have set $\nu_{np} = \nu$. The dependence on ν_{np} first appears at three loops in this $\mathcal{O}(r^2)$ contribution to δE_o^{us} . At present, concrete results for the ultrasoft corrections are available at one, $\mathcal{O}(g^2)$, and two-loop level, $\mathcal{O}(g^4)$, given in Eqs. (21) and (22). After MS subtraction they read

$$\delta E_{o,MS}^{us}(1\text{-loop}) = \frac{1}{8N_c} \alpha V_A^2 r^2 \Delta V_{MS}^2 \left(1 + \gamma_E - \ln(4\pi) + 2 \ln \left[\frac{-\Delta V_{MS}}{\nu} \right] \right), \quad (62)$$

$$\delta E_{o,MS}^{us}(2\text{-loop}) = \frac{1}{4N_c} \alpha^2 V_A^2 r^2 \Delta V_{MS} \left(-C_A \ln^2 \left[\frac{-\Delta V_{MS}}{\nu} \right] + C_A (2 - \gamma_E + \ln(4\pi)) \ln \left[\frac{-\Delta V_{MS}}{\nu} \right] - \frac{1}{2} C_A \left(\pi^2 - 7 + \frac{1}{2} (2 - \gamma_E + \ln(4\pi))^2 \right) - 2n_f T_F \right). \quad (63)$$

Note that these expressions have imaginary parts. Setting $\nu \sim \Delta V_X(\nu)$ resums large logarithms into the potential V_o and minimizes them in the ultrasoft contribution δE_o^{us} . It also simplifies the determination of the decay width of the octet (hybrid) system

$$\Gamma_H = \frac{1}{2N_c} \alpha V_A^2 r^2 \Delta V_{MS}^2 (\bar{\nu}_{us}) \pi + \frac{1}{2N_c} \alpha^2 V_A^2 r^2 \Delta V_{MS} (\bar{\nu}_{us}) \times C_A \pi (2 - \gamma_E + \ln(4\pi)) + \mathcal{O}(\alpha^3 r^2 \Delta V^0), \quad (64)$$

where

$$\bar{\nu}_{us} \equiv \Delta V_X(\bar{\nu}_{us}) = C_A \alpha W(1/(C_A \alpha d_X r)) \quad (65)$$

and $W(z)$ is the Lambert function (here: $X = MS$).

The precision of Eq. (64) is limited by the existence of some unknown single logarithms that appear in the three-loop ultrasoft self-energy (see next subsection). The imaginary parts of those logarithms produce terms like $\delta\Gamma \sim \alpha^3 r^2$.

C. Subleading ultrasoft and nonperturbative effects

Even without an explicit computation we can obtain some information on the ultrasoft three-loop terms. At three loops we start to have contributions from diagrams

with no singlet-octet vertices. They are proportional to V_B^2 or V_C and vanish in dimensional regularization (for on shell external fields), as they are insensitive to the scale ΔV . Therefore, any ultraviolet divergence proportional to V_B^2 or V_C should be canceled by an infrared one, which signals a sensitivity to the nonperturbative scale α . We quantify this statement with the following equation

$$\nu \frac{d}{d\nu} \delta E_o^{us}(\nu, \nu) = -B(V) - \nu \frac{d}{d\nu} \delta E_H^{np}, \quad (66)$$

where the RG structure of the nonperturbative term is the following

$$\nu \frac{d}{d\nu} \delta E_H^{np} = C_A \alpha + B r^2 \alpha^3 (1 + \mathcal{O}(\alpha/\Delta V)). \quad (67)$$

Solving this equation we obtain

$$\delta E_H^{np}(\nu_{np}) = C_A \alpha \left(\ln \left(\frac{\nu_{np}}{\alpha} \right) + c_H \right) + B \alpha^3 r^2 \ln \left(\frac{\nu_{np}}{\alpha} \right) + \mathcal{O}(\alpha^3 r^2), \quad (68)$$

where c_H is a nonperturbative constant that depends on the specific hybrid (gluelump) we consider. Actually, the $\mathcal{O}(r^0)$ term is nothing but the gluelump mass

$$\Lambda_H(\nu_{np}) = C_A \alpha \left(\ln \left(\frac{\nu_{np}}{\alpha} \right) + c_H \right). \quad (69)$$

It can be related to a gauge-invariant correlator as

$$\Lambda_H = \lim_{T \rightarrow \infty} \frac{i}{T} \ln \langle H^a(T/2) \phi_{ab}^{\text{adj}}(T/2, -T/2) H^b(-T/2) \rangle, \quad (70)$$

where H^a is the field operator associated with the gluonic component of the hybrid and $\phi_{ab}^{\text{adj}}(T/2, -T/2)$ is a Wilson line in the adjoint representation, see Ref. [4].

Note on the other hand that B is independent of the hybrid type and can be obtained from perturbation theory. It is however unknown at present and constrains, besides the nonperturbative gluelump mass Λ_H , the precision of our result.

At this point we could also analyze subleading ultrasoft effects in the $\alpha/\Delta V$ expansion along the lines of Ref. [1], as some of them, namely, the logarithmic terms proportional to V_A^2 , are fixed by the RG structure. In view of the dominant uncertainties discussed above, and because there are more unknowns than in the singlet case [1], we refrain from performing that analysis here.

Combining Eqs. (57), (62), (63), and (68) we can write down the static hybrid energy $E_H = 2m_{MS} + V_{o,MS} + \delta E_{o,MS}^{us} + \delta E_{H,MS}^{np}$. Expressed as a double expansion in αr and $1/\ln(r\Delta V)$ it reads with $\mathcal{O}(\alpha^2 r^2)$ and NNLL accuracy

$$\begin{aligned}
 E_H &= 2m_{\text{MS}}(\nu_s) + C_F \alpha \ln(r^2 \nu_s^2 \pi e^{\gamma_E}) + \frac{\pi}{4} \left(C_F - \frac{C_A}{2} \right) (7C_A - 4n_f T_F) \alpha^2 r + \delta V_{o,\text{MS}}^{\text{RG}} \left(r; \frac{1}{r}, \bar{\nu}_{us} \right) \\
 &\quad + \delta E_{o,\text{MS}}^{us}(\bar{\nu}_{us}, \nu_{np}) + \delta E_{H,\text{MS}}^{np}(\nu_{np}) + \mathcal{O}(\alpha^3 r^2 \ln^0) \\
 &= 2m_{\text{MS}}(\nu_s) + C_F \alpha \ln(r^2 \nu_s^2 \pi e^{\gamma_E}) - \frac{C_A}{2} \alpha \ln(r^2 \alpha^2 \pi e^{\gamma_E}) + C_A \alpha c_{H,\text{MS}} \\
 &\quad + \frac{\pi}{4} \left(C_F - \frac{C_A}{2} \right) (7C_A - 4n_f T_F) \alpha^2 r - \left(C_F - \frac{C_A}{2} \right) \alpha^3 r^2 \left\{ \frac{1}{6} C_A^2 \ln^3(r \Delta V_{\text{MS}}) + \frac{1}{4} C_A^2 (2\gamma_E - 3 - 2 \ln 2) \ln^2(r \Delta V_{\text{MS}}) \right. \\
 &\quad \left. + \left[C_A^2 \left(\frac{83\pi^2}{384} + \frac{1}{8} (4\gamma_E^2 + 4 \ln^2 2 - \gamma_E (10 + \ln 256) + 38 + 8 \ln 2 - 2 \ln \pi) \right) \right. \right. \\
 &\quad \left. \left. - n_f T_F \left(C_A \left(2 + \frac{19\pi^2}{48} \right) + C_F \left(5 - \frac{\pi^2}{2} \right) \right) + (n_f T_F)^2 \frac{\pi^2}{8} \right] \ln(r \Delta V_{\text{MS}}) \right\} + \mathcal{O}(\alpha^3 r^2 \ln(\Delta V_{\text{MS}}/\alpha)). \tag{71}
 \end{aligned}$$

We have checked that the explicit scheme dependence of $\delta V_{o,\text{MS}}^{\text{RG}}$ and $\delta E_{o,\text{MS}}^{us}$ and the implicit scheme dependence of Eq. (71) through the logarithms of ΔV cancel up to $\mathcal{O}(\alpha^3 r^2 \ln^0)$. Note also that $c_{H,\text{MS}}$ is scheme-dependent.

The leading uncertainty of Eq. (71) comes from the nonperturbative gluon mass Λ_H . The constant Λ_H is independent on the distance r and therefore drops out in the force, i.e., the derivative of the potential. In that case the leading uncertainty is due to the coefficient B , which is unknown but, unlike Λ_H , of perturbative origin and could be determined by an ultrasoft three-loop calculation.

Equation (71) represents one of the main results of this paper. One could try to perform some comparison with the (quenched) lattice data existing in the literature [23–27]. It is not clear though that these simulations reach short enough distances that our results can be tested quantitatively (on the other hand it should not be too costly to perform a dedicated short-distance simulation to test our expression). Moreover, it would be as well very interesting to study the accidental symmetries and degeneracies among different hybrid energies that should appear at short distances along the lines of the studies in four dimensions performed in Refs. [4,5]. To this end one could consider energy differences where the perturbative terms (the non-analytic, i.e., logarithmic dependence in r) cancels. The noncanceled terms should be produced by the nonperturbative terms (note that the scheme dependence of $c_{H/H'}$ cancels in the difference)

$$E_H - E_{H'} = C_A \alpha (c_H - c_{H'}) + C_{HH'} r^2 + \mathcal{O}(r^3). \tag{72}$$

All this would require a dedicated analysis, which will be carried out elsewhere.

V. RESULTS FOR $D = 2$

It is also interesting to investigate static hybrid systems in two space-time dimensions. In the following we set $n_f = 0$ for simplicity. In exactly two dimensions physical hybrids do not exist, as there are no propagating (physical) gluons. It is instructive to see how this finding arises in an

explicit calculation. For a typical hybrid, the object to be computed is, e.g.,

$$\begin{aligned}
 \langle W_H \rangle &= \left\langle \frac{1}{N_c} \text{Tr} \mathcal{P} \mathbf{E}(T/2, \mathbf{R}) \cdot \mathbf{E}(-T/2, \mathbf{R}') \right. \\
 &\quad \left. \times \exp \left(-ig \oint_{\Gamma} A_{\mu} dx^{\mu} \right) \right\rangle, \tag{73}
 \end{aligned}$$

where the contour Γ of the integral in Eq. (73) is a rectangle with spatial extension r and temporal extension T in Minkowski space and we have inserted two chromoelectric fields on each spacelike end-string of the rectangular Wilson loop. We could have also chosen other gluonic configurations instead, but this would only make the discussion more complicated without changing the physical outcome.

Equation (73) is gauge-invariant. Choosing axial ($A_1 \equiv 0$) gauge in exactly two dimensions, the only non-vanishing component of the gluon field-strength tensor is $F_{01} = -F_{10} = -\partial_1 A_0$. Hence, (A_0) gluons neither interact among themselves, nor propagate in time, since no time derivative acting on the gluon field is left in the Lagrangian. Therefore only⁶ planar “ladder” diagrams (with “potential” gluons) contribute to $\langle W_H \rangle$. In the large T limit the final result takes the form

$$\langle W_H \rangle \sim \langle E^a(T/2) E^a(-T/2) \rangle e^{-iV_o T} \tag{74}$$

where $V_o(r) = \frac{1}{2N_c} \alpha r$. Note that this octet potential is the same as obtained from a leading order, tree-level, computation in 2D static QCD. Therefore the nonexistence of the hybrid does not arise from the fact that the octet potential diverges in two dimensions (generating an infinite hybrid mass) but rather from the fact that

$$\langle E^a(T/2) E^a(-T/2) \rangle \sim \delta(T) \tag{75}$$

is local in time. That is because the gluon does not propagate in time.

⁶We do not include in this discussion possible interactions of \mathbf{E} with the gluons in the timelike strings. They would produce terms proportional to the singlet: $\sim e^{-iV_s T}$, but not contribute to the hybrid energy.

For $D > 2$ one expects $\langle W_H \rangle$ to converge towards the previous result in Eqs. (74) and (75) as $D \rightarrow 2$, provided the limit is smooth. The discussion parallels to some extent the one for the singlet potential in Ref. [1]. Like in that case we cannot make definite statements, since the calculation of the Wilson loop becomes intrinsically nonperturbative. The reason is that the proper ultrasoft expansion parameter is now $\Delta V/g \sim gr$, see Ref. [1]. Thus the nonperturbative ultrasoft ($\Delta V/g$) expansion and the perturbative multipole (gr) expansion mix.

All we can say is that the cancelation of the soft and ultrasoft one-loop contributions we observed for the singlet potential [1] also occurs for the hybrid potential, because the computation differs only in the prefactor. For the chromoelectric correlator we find at leading order in α

$$\langle E^a(T/2)E^a(-T/2) \rangle \sim \mathcal{O}(\epsilon). \quad (76)$$

This is what we would expect from the decoupling of gluons in two dimensions. Nevertheless at NLO in α one obtains a finite contribution. Again we can not draw any definite conclusion, as this result is obtained in standard perturbation theory, whereas the correct expansion entails inverse powers of the coupling constant g .

VI. CONCLUSIONS

We have computed the energy of a static hybrid in the weak coupling limit in four and three space-time dimensions. Employing finite temperature theory methods we have checked the static potential of a color-octet quark-antiquark pair with distance r at two loops for an arbitrary number of dimensions D . The result represents the soft contribution to the static hybrid energy. Using the effective theory pNRQCD we have explicitly calculated the missing ultrasoft two-loop contributions. We have also confirmed the respective two-loop result for the color singlet. The cancelation of soft infrared and ultrasoft ultraviolet divergences is a strong cross-check of our results.

For $D = 4$ we have determined the static octet potential and the static hybrid energy through N^3 LL order, up to the unknown three-loop soft matching condition. The result

disagrees with an earlier result obtained in Ref. [14]. We have also given the ultrasoft contribution to the static hybrid energy at N^4 LL and computed the decay width of the hybrid/octet system at NLL order.

In $D = 3$ dimensions our result for the static hybrid energy reaches $\mathcal{O}(\alpha^3 r^2)$ accuracy in the (soft) multipole expansion. At this order we have determined the complete expression at NNLL order in the ultrasoft $\alpha/\Delta V$ expansion. The precision of the result is only limited by unknown terms of order $\mathcal{O}(\alpha^3 r^2 \ln(\Delta V/\alpha))$ and the nonperturbative gluelump mass. The former can in principle be obtained by a perturbative three-loop computation, whereas the latter requires lattice simulations (note, however, that the gluelump mass vanishes in derivatives of the potential, like the force). Besides the energy we have determined the 3D hybrid/octet decay width through $\mathcal{O}(\alpha^3 r^2)$ in the multipole expansion and NLO in the $\alpha/\Delta V$ expansion.

We have also studied the two-dimensional case, where the exact result is known: hybrids do not exist. This is due to the temporal locality of the gluonic correlators as there are no propagating (physical) gluons in *exactly* two space-time dimensions, see Eqs. (74) and (75), rather than due to an infinite octet potential or hybrid mass. In the $D \rightarrow 2$ limit this result is more complicated to obtain, because already the ultrasoft contribution is intrinsically nonperturbative. Strong cancellations have to occur among the different contributions from the soft, the ultrasoft and the nonperturbative (g) scale. We can only report partial and inconclusive results on this limit, particularly because they are based on perturbation theory, which is not reliable for $D \rightarrow 2$. Nevertheless, it might be worth exploring the $D \rightarrow 2$ limit in more detail, as it could provide nontrivial information about the dependence of the (D -dimensional) perturbative results on $(D - 2)$.

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APPENDIX A: TWO-LOOP PNRQCD SELF-ENERGY DIAGRAMS IN FEYNMAN GAUGE

1. Color octet

$$\begin{aligned} \text{Diagram} &= C_o^{(d)} \frac{\Gamma(5-2d)\Gamma(\frac{d}{2})^2}{4(d-4)(d-3)} \\ &= \begin{cases} C_o^{(3)} \left[\frac{1}{8\epsilon_3^2} + \frac{2L+1}{4\epsilon_3} + L^2 + L + \frac{11\pi^2}{48} + 1 + \mathcal{O}(\epsilon_3) \right] \\ C_o^{(4)} \left[-\frac{1}{12\epsilon_4^2} + \frac{-6L+11-6\ln 2}{18\epsilon_4} - \frac{2L^2}{3} - \frac{2L(6\ln 2-11)}{9} - \frac{\pi^2}{8} - \frac{349}{108} - \frac{2\ln^2 2}{3} + \frac{22\ln 2}{9} + \mathcal{O}(\epsilon_4) \right] \end{cases} \end{aligned} \quad (A1)$$

$$\begin{aligned} \text{Diagram 1} &= C_o^{(d)} \frac{\Gamma(5-2d)\Gamma\left(\frac{d}{2}-1\right)^2}{16} \end{aligned} \quad (\text{A2})$$

$$= \begin{cases} C_o^{(3)} \left[-\frac{1}{4\epsilon_3} - L + 1 + \mathcal{O}(\epsilon_3) \right] \\ C_o^{(4)} \left[-\frac{1}{24\epsilon_4} - \frac{L}{6} + \frac{11-6\ln 2}{36} + \mathcal{O}(\epsilon_4) \right] \end{cases}$$

$$\begin{aligned} \text{Diagram 2} &= C_o^{(d)} \frac{(d-2)\Gamma(5-2d)\Gamma\left(\frac{d}{2}-1\right)^2}{8(d-3)} \end{aligned} \quad (\text{A3})$$

$$= \begin{cases} C_o^{(3)} \left[-\frac{1}{4\epsilon_3^2} + \frac{1-2L}{2\epsilon_3} - 2L^2 + 2L - \frac{11\pi^2}{24} - 2 + \mathcal{O}(\epsilon_3) \right] \\ C_o^{(4)} \left[-\frac{1}{6\epsilon_4} - \frac{2L}{3} + \frac{25}{18} - \frac{2\ln 2}{3} + \mathcal{O}(\epsilon_4) \right] \end{cases}$$

$$\begin{aligned} \text{Diagram 3} &= C_o^{(d)} \left[\frac{2^{1-2d} (d^2 - 6d + 10) \Gamma\left(\frac{1}{2}-d\right) \Gamma(1-d)^2 \Gamma\left(d+\frac{1}{2}\right) \Gamma(d)^2}{(d-4)\Gamma\left(\frac{d-1}{2}\right) \Gamma\left(\frac{d+1}{2}\right)} \right. \\ &+ \left. \frac{\sqrt{\pi} 2^{-d-2} (-3d^4 + 30d^3 - 111d^2 + 181d - 112) \Gamma(5-2d)\Gamma\left(\frac{d}{2}-2\right) \Gamma(d-4)}{\Gamma\left(\frac{d+1}{2}\right)} \right] \end{aligned} \quad (\text{A4})$$

$$= \begin{cases} C_o^{(3)} \left[-\frac{1}{2\epsilon_3} - 2L + \frac{1}{6} (15 - \pi^2) + \mathcal{O}(\epsilon_3) \right] \\ C_o^{(4)} \left[-\frac{1}{8\epsilon_4^2} + \frac{-9L-2\pi^2+15-9\ln 2}{18\epsilon_4} - L^2 - \frac{1}{9}L(-30 + 4\pi^2 + 18\ln 2) \right. \\ \left. + \frac{1}{6}(8\zeta(3) - 27 + 20\ln 2 - 6\ln^2 2) - \frac{\pi^2}{432}(192\ln 2 - 47) + \mathcal{O}(\epsilon_4) \right] \end{cases}$$

$$\begin{aligned} \text{Diagram 4} &= C_o^{(d)} \frac{3(3-d)\Gamma(5-2d)\Gamma\left(1-\frac{d}{2}\right) \Gamma\left(\frac{d}{2}-1\right) \Gamma\left(\frac{d}{2}\right)}{16\Gamma\left(3-\frac{d}{2}\right)} \end{aligned} \quad (\text{A5})$$

$$= \begin{cases} C_o^{(3)} \left[-3 + \mathcal{O}(\epsilon_3) \right] \\ C_o^{(4)} \left[\frac{1}{8\epsilon_4^2} + \frac{3L-4+3\ln 2}{6\epsilon_4} + L^2 + \frac{2L(3\ln 2-4)}{3} + \frac{3\pi^2}{16} + \frac{26}{9} + \ln^2 2 - \frac{8\ln 2}{3} + \mathcal{O}(\epsilon_4) \right] \end{cases}$$

$$\begin{aligned} \text{Diagram 5} &= C_o^{(d)} \frac{C_A(3d-2) - 4(d-2)n_f T_F}{C_A} \times \\ &\times \frac{\sqrt{\pi}(d-3)\Gamma(5-2d)\Gamma\left(1-\frac{d}{2}\right) \Gamma(d-2)\Gamma\left(\frac{d}{2}\right)}{2^{d+3}\Gamma\left(3-\frac{d}{2}\right) \Gamma\left(\frac{d+1}{2}\right)} \end{aligned} \quad (\text{A6})$$

$$= \begin{cases} C_o^{(3)} \left[\frac{7}{4} - \frac{n_f T_F}{C_A} + \mathcal{O}(\epsilon_3) \right] \\ C_o^{(4)} \left[\left[\frac{4n_f T_F}{9C_A} - \frac{5}{9} \right] \frac{1}{8\epsilon_4^2} + \frac{1}{2\epsilon_4} \left(\left[\frac{4n_f T_F}{9C_A} - \frac{5}{9} \right] (L - \frac{5}{4} + \ln 2) + \frac{1}{18} \right) + \frac{L}{9} - \frac{1}{6} + \frac{\ln 2}{9} \right. \\ \left. + \left[\frac{4n_f T_F}{9C_A} - \frac{5}{9} \right] \left(L^2 + L(2\ln 2 - \frac{5}{2}) + \frac{3\pi^2}{16} + \frac{95}{36} + \ln^2 2 - \frac{5\ln 2}{2} \right) + \mathcal{O}(\epsilon_4) \right] \end{cases}$$

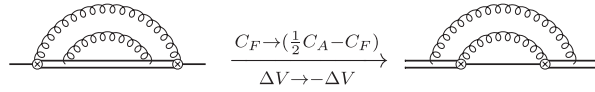
$$L := \ln \left[-\frac{\Delta V}{\nu_{\overline{\text{MS}}}} \right]; \quad \nu_{\overline{\text{MS}}} := \nu e^{-(1/2)(\gamma_E - \ln(4\pi))} \quad (\text{A7})$$

$$C_o^{(d)} := -16iC_A\left(\frac{C_A}{2} - C_F\right)\pi^{2-d}r^2V_A^2\alpha^2(-\Delta V)^{2d-5}\nu_{\text{MS}}^{-4\epsilon_n}e^{-2\epsilon_n(\gamma_E - \ln(4\pi))} \quad (\text{A8})$$

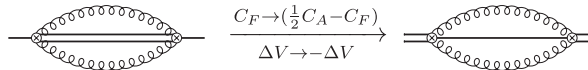
$$C_o^{(3)} := -iC_A\left(\frac{C_A}{2} - C_F\right)\Delta Vr^2V_A^2\alpha^2 \quad (\text{A9})$$

$$C_o^{(4)} := -\frac{iC_A\left(\frac{C_A}{2} - C_F\right)\Delta V^3r^2V_A^2\alpha^2}{\pi^2} \quad (\text{A10})$$

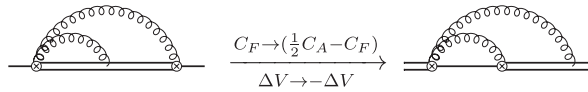
2. Relation between singlet and octet diagrams



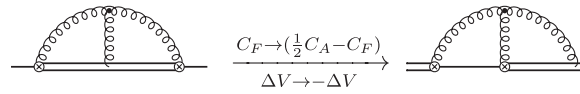
$$\quad \xrightarrow[\Delta V \rightarrow -\Delta V]{C_F \rightarrow (\frac{1}{2}C_A - C_F)} \quad (\text{A11})$$



$$\quad \xrightarrow[\Delta V \rightarrow -\Delta V]{C_F \rightarrow (\frac{1}{2}C_A - C_F)} \quad (\text{A12})$$



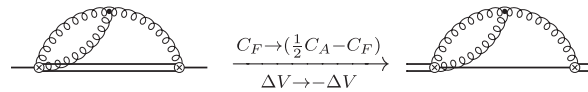
$$\quad \xrightarrow[\Delta V \rightarrow -\Delta V]{C_F \rightarrow (\frac{1}{2}C_A - C_F)} \quad (\text{A13})$$



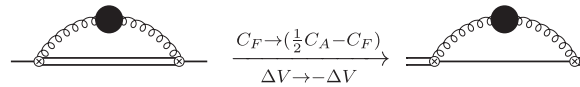
$$\quad \xrightarrow[\Delta V \rightarrow -\Delta V]{C_F \rightarrow (\frac{1}{2}C_A - C_F)} \quad + \Delta_{so} \quad (\text{A14})$$

$$\Delta_{so} = -C_o^{(d)} \frac{4^{-d}(d^2 - 6d + 10)\Gamma(\frac{1}{2} - d)\Gamma(\frac{1}{2} - \frac{d}{2})^2\Gamma(\frac{d}{2} + \frac{1}{2})^2\Gamma(d + \frac{1}{2})}{(d - 4)\Gamma(\frac{d-1}{2})\Gamma(\frac{d+1}{2})} = \quad (\text{A15})$$

$$= \begin{cases} C_o^{(3)} \left[\frac{1}{4\epsilon_3^2} + \frac{4L+1}{4\epsilon_3} + 2L^2 + L + \frac{13\pi^2}{24} + \frac{7}{4} + \mathcal{O}(\epsilon_3) \right] \\ C_o^{(4)} \left[\frac{\pi^2}{6\epsilon_4} + \frac{2\pi^2 L}{3} + \frac{2}{9}\pi^2(3\ln 2 - 2) + \mathcal{O}(\epsilon_4) \right] \end{cases} \quad (\text{A16})$$



$$\quad \xrightarrow[\Delta V \rightarrow -\Delta V]{C_F \rightarrow (\frac{1}{2}C_A - C_F)} \quad (\text{A17})$$



$$\quad \xrightarrow[\Delta V \rightarrow -\Delta V]{C_F \rightarrow (\frac{1}{2}C_A - C_F)} \quad (\text{A18})$$

[1] A. Pineda and M. Stahlhofen, *Phys. Rev. D* **81**, 074026 (2010).
 [2] I. H. Jorjusz and C. Michael, *Nucl. Phys.* **B302**, 448 (1988).
 [3] M. Foster and C. Michael (UKQCD Collaboration), *Phys. Rev. D* **59**, 094509 (1999).
 [4] N. Brambilla, A. Pineda, J. Soto, and A. Vairo, *Nucl. Phys.* **B566**, 275 (2000).
 [5] G. S. Bali and A. Pineda, *Phys. Rev. D* **69**, 094001 (2004).
 [6] A. Pineda and J. Soto, *Nucl. Phys. B, Proc. Suppl.* **64**, 428 (1998).
 [7] N. Brambilla, A. Pineda, J. Soto, and A. Vairo, *Rev. Mod. Phys.* **77**, 1423 (2005).
 [8] B. A. Kniehl, A. A. Penin, Y. Schroder, V. A. Smirnov, and M. Steinhauser, *Phys. Lett. B* **607**, 96 (2005).
 [9] N. Brambilla, A. Pineda, J. Soto, and A. Vairo, *Phys. Rev. D* **60**, 091502 (1999).

- [10] A. Pineda and J. Soto, *Phys. Lett. B* **495**, 323 (2000).
- [11] N. Brambilla, J. Ghiglieri, P. Petreczky, and A. Vairo, *Phys. Rev. D* **82**, 074019 (2010).
- [12] M. Eidemuller and M. Jamin, *Phys. Lett. B* **416**, 415 (1998).
- [13] N. Brambilla, X. Garcia i Tormo, J. Soto, and A. Vairo, *Phys. Lett. B* **647**, 185 (2007).
- [14] N. Brambilla, A. Vairo, X. Garcia i Tormo, and J. Soto, *Phys. Rev. D* **80**, 034016 (2009).
- [15] Y. Schroder, Ph.D. thesis, University of Hamburg, 1999, <http://www.physik.uni-bielefeld.de/~yorks/www/papers/t2.pdf>.
- [16] Y. Schroder, *Phys. Lett. B* **447**, 321 (1999); M. Peter, *Phys. Rev. Lett.* **78**, 602 (1997).
- [17] P. Pascual and R. Tarrach, *Lect. Notes Phys.* **194**, 1 (1984).
- [18] A. Pineda and J. Soto, *Phys. Lett. B* **420**, 391 (1998).
- [19] B. A. Kniehl and A. A. Penin, *Nucl. Phys.* **B563**, 200 (1999).
- [20] B.J. Hand and G. Leibbrandt, [arXiv:hep-th/9511140](https://arxiv.org/abs/hep-th/9511140).
- [21] B.A. Kniehl and A. Sirlin, *Phys. Rev. Lett.* **81**, 1373 (1998).
- [22] W. Fischler, *Nucl. Phys.* **B129**, 157 (1977).
- [23] O. Philipsen, *Phys. Lett. B* **535**, 138 (2002).
- [24] K.J. Juge, J. Kuti, and C. Morningstar, [arXiv:hep-lat/0312019](https://arxiv.org/abs/hep-lat/0312019).
- [25] M. Caselle, M. Pepe, and A. Rago, *J. High Energy Phys.* **10** (2004) 005.
- [26] J. Kuti, *Proc. Sci.*, LAT2005, 001 (2006); HW2005 (2006) 009 [[arXiv:hep-lat/0511023](https://arxiv.org/abs/hep-lat/0511023)].
- [27] B.B. Brandt and P. Majumdar, *Phys. Lett. B* **682**, 253 (2009).
- [28] D. Binosi and L. Theussl, *Comput. Phys. Commun.* **161**, 76 (2004).