

String formation and chiral symmetry breaking in the heavy-light quark-antiquark system in QCD

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The effective quark Lagrangian is written for a light quark in the field of a static antiquark, explicitly containing field correlators as coefficient functions of products of quark operators. At large N_c the closed system of equations for the gauge-invariant quark Green's function in the field of static source is examined analytically. The formation of the string connecting the light quark to the static source is observed numerically. The scalar Lorentz nature of the resulting confinement is shown to hold for the considered case, implying chiral symmetry breaking. The resulting spectrum with and without perturbative gluon exchanges is obtained numerically and compared to the B and D meson masses and heavy quark effective theory.

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I. INTRODUCTION

The nature of QCD string formation between static sources was studied on the lattice [1,2] and analytically [2]. From these investigations it was shown that the string consists of a predominantly color-electric longitudinal field. At the critical temperature T_c this electric field disappears and at the same time the deconfined phase with color-magnetic condensate sets in. This effect was predicted theoretically in [3] and also seen in lattice measurements [4]. At the same temperature chiral symmetry breaking (CSB) for light quarks is found to disappear [5], which indicates that there is an intimate connection between the string formation and CSB. In the case of heavy quark systems CSB occurs due to the quark mass and confinement can be described as the area law of the Wilson loop. How CSB and confinement are explicitly realized for the light quark system and what equation describes its dynamics is an interesting and open problem.

It is the purpose of the present paper to study this issue in the simplest dynamical example—in the system of one light quark and a heavy antiquark. This allows us to describe the dynamics of light quark (its propagator) in a gauge-invariant manner, while physically the light quark is expected to be confined at the end of a string connected to the static source. Applying the formalism of field correlators (FC) [6,7], we derive the effective quark Lagrangian, containing any number of quark operators multiplied by field correlators. To proceed further one can use the limit of large N_c and write down the Dyson-Schwinger equations with the mass operator expressed through the Green's function. The resulting equations are nonlocal and nonlinear. It is not clear from the beginning how confinement and CSB would manifest themselves in the solution of these equations. Some hint was provided in Ref. [8] using a relativistic WKB analysis [9], where it was shown that at large distances from the heavy source the dynamics of the light quark is described by the Dirac equation with a scalar linear confining interaction, which leads to CSB.

In this paper we examine the properties of the Green's function of the color singlet $q\bar{Q}$ system, where the antiquark is treated in the static limit. In Sec. II we formulate the full form of nonlocal and nonlinear equations for the light quark propagator and its eigenfunctions and study the behavior at all distances. We also take into account both perturbative and nonperturbative contributions to the interaction kernel. As a result our equations contain both a confining interaction and color Coulomb part. Similar to the heavy quark situation we argue that the string formation for low angular momentum is of a color-electric nature. Moreover, the confinement of the light quark to the heavy one is shown to be of Lorentz scalar type. In Sec. III the resulting nonlinear equations are studied numerically. The energy spectrum and the structure of the low-lying eigenfunctions are presented. We in particular study the B and D meson spectra. We compare them with experimental data for B , D mesons and results of other calculations, exploiting for this purpose the expansion of heavy quark effective theory (HQET). In Sec. IV the chiral condensate of the light quarks $\langle\bar{q}q\rangle$ is determined by taking the limit of the Green's function $S(x,y)$, with both x and y tending to zero. In this limit the heavy quark is turned off and the condensate $\langle\bar{q}q\rangle\sim S(0,0)$ should have a value not depending on the presence of heavy quark.

II. DERIVATION OF EQUATIONS

A. Dyson-Schwinger equations

In this section we give an outline of the procedure to obtain the Dyson-Schwinger equations for the color white $q\bar{Q}$ system. Our starting point is the gauge-invariant light quark Green's function $S(x,y)$ in the presence of a static heavy antiquark placed in the origin. In the static limit, the heavy quark can be treated as an external source. Assuming the Euclidean metric and letting $T=(x-y)_4$, the heavy antiquark propagator in the modified Fock-Schwinger gauge [10] is proportional to the parallel transporter: namely,

$$S_{\bar{Q}}(A) = h(\mathbf{x}, \mathbf{y}) P \exp \left(ig \int_0^T A_4(\mathbf{r}=0, \tau) d\tau \right), \quad (1)$$

with

$$h(\mathbf{x}, \mathbf{y}) = \frac{i}{2} \delta^{(3)}(\mathbf{x} - \mathbf{y}) [(1 + \gamma_4) e^{-m_Q T} \Theta(T) + (T \rightarrow -T, \gamma_4 \rightarrow -\gamma_4)].$$

This acts as a static source situated at the origin. As a result the proper limit for $m_Q \rightarrow \infty$ of the Green's function of the $q\bar{Q}$ system can be defined as

$$S(x, y) = \langle \psi(x) \Pi(x, y) \psi^+(y) \rangle. \quad (2)$$

Here we have to average over the gluon fields A and light quark fields ψ , while Π contains the parallel transporters between the end points

$$\Pi(x, y) \equiv \phi(\mathbf{x}, x_4; 0, x_4) \phi(0, x_4; 0, y_4) \phi(0, y_4; \mathbf{y}, y_4),$$

with

$$\phi(\mathbf{x}, x_4; \mathbf{y}, y_4) = P \exp \left(ig \int_x^y A_\mu dz_\mu \right).$$

The averaging over the gluon fields A has to be done over the perturbative and nonperturbative gluon fields a_μ and B_μ contributions, respectively, where the total gluonic field $A_\mu = B_\mu + a_\mu$. We now apply the method of field correlators, which was developed in a series of papers [6,7] to derive the effective quark Lagrangian from QCD. Let us first consider the effects of averaging over nonperturbative (NP) field B_μ . We may write for the partition function

$$\begin{aligned} \langle P e^{g \int \psi^+ \hat{B}(x) \psi dx} \rangle_B &\equiv e^{L_{\text{eff}}} \\ &= P \exp \left[\sum_n \frac{g^n}{n!} \int d^4 x_1 \dots d^4 x_n j(1) \dots j(n) \right. \\ &\quad \left. \times \langle \langle B(1) \dots B(n) \rangle \rangle \right], \quad (3) \end{aligned}$$

with $j(n) \equiv j_{\mu_n}(x_n) = \psi^+(x_n) \gamma_{\mu_n} \psi(x_n)$ and $B(n) = B_{\mu_n}(x_n)$. To write the correlator $\langle \langle B \dots B \rangle \rangle$ for the gauge-invariant situation corresponding to the color white $q\bar{Q}$ system one can use the modified Fock-Schwinger gauge [10] to express the correlator of $B(x)$ through FC:

$$\begin{aligned} N(1, \dots, n) &\equiv \langle \langle B(1) \dots B(n) \rangle \rangle \sim \int dx(1) \dots dx(n) \\ &\quad \times \langle \langle F(1) \dots F(n) \rangle \rangle. \quad (4) \end{aligned}$$

The effective interaction kernel in Eq. (3) can now be used to write a Dyson-Schwinger-type equation for the quark Green's function S . To simplify the matter, one can consider

the large N_c limit, in which case the connected self-energy kernel $M(x, y)$ is obtained from Eq. (3) by replacing any pair of adjacent ψ operators by

$$\psi_{a\alpha}(x) \psi_{a\beta}^+(y) \rightarrow N_c S_{\alpha\beta}(x, y), \quad (5)$$

where $a\alpha$ are color and Lorentz index, respectively (for details of derivation see [8]). As a result one obtains the equation for $S(x, y)$

$$(-i \hat{b}_x - im) S(x, y) - i \int M(x, z) S(z, y) d^4 z = \delta^{(4)}(x - y), \quad (6)$$

where the kernel M is expressed through N as

$$\begin{aligned} iM(x, y) &= N_{\mu\nu}^{(2)}(x, y) \gamma_\mu S(x, y) \gamma_\nu \\ &+ \sum_{n=3}^{\infty} \int d^4 x_2 \dots d^4 x_{n-1} \gamma_{\mu_1} S(x, x_2) \\ &\quad \times \gamma_{\mu_2} \dots \gamma_{\mu_{n-1}} S(x_{n-1}, y) \gamma_{\mu_n} N_{\mu_1 \dots \mu_n}^{(n)} \\ &\quad \times (x, x_2 \dots x_{n-1}, y). \quad (7) \end{aligned}$$

The system of equations (6), (7) is exact in the large N_c limit and is well defined provided all NP correlators $\langle \langle F(1) \dots F(n) \rangle \rangle$ are known.

Evidence has been found in recent accurate lattice calculations [11] of static potentials in different $SU(3)$ representations, that the contributions of the higher correlators $\langle \langle F(1) \dots F(n) \rangle \rangle$ for $n > 2$ to the planar Wilson loop are small. In particular, these terms are found to contribute only around a few percent of the dominant Gaussian correlator [12]. Hence, the Gaussian stochastic model, based on the lowest correlator $\langle \langle F(1) F(2) \rangle \rangle$, is expected to be a good approximation. In view of this we assume in this paper that the Gaussian approximation holds and we keep similar to Ref. [8] only the first term in Eq. (8). One can parametrize the Gaussian correlator according to [6] as

$$\begin{aligned} \langle \langle F(1) F(2) \rangle \rangle &= \frac{1}{N_c} \text{tr} \langle F_{\mu\lambda}(x) \phi(x, 0) F_{\nu\sigma}(0) \phi(0, x) \rangle \\ &= D(x) (\delta_{\mu\nu} \delta_{\lambda\sigma} - \delta_{\mu\sigma} \delta_{\nu\lambda}) + \Delta_{\mu\lambda\nu\sigma}^{(1)}, \quad (8) \end{aligned}$$

where only $D(x)$ is responsible for confinement and it contributes to string tension σ , while $\Delta^{(1)}$ is a full derivative. As a consequence, the latter contributes to the perimeter of Wilson loop and $\phi(x, 0) = P \exp ig \int_0^x B_\mu dz_\mu$. From this we find that the $N_{\mu\nu}^{(2)}$ can explicitly be written in the gauge [10] as

$$\begin{aligned} N_{\mu\nu}^{(2)}(x, y) &= (\delta_{\mu\nu} \delta_{ik} - \delta_{\mu k} \delta_{\nu i}) \int_0^x du_i \alpha_\mu(u) \\ &\quad \times \int_0^y dv_k \alpha_\nu(v) D(u - v), \quad (9) \end{aligned}$$

where $\alpha_4(u) = 1, \alpha_i(u) = u_i/x_i$.

In Eq. (9) only the nonperturbative confining piece of the Gaussian correlator (8) is retained, since the perturbative part and $\Delta^{(1)}$ do not produce either the string (confinement) or CSB [8]. In the case of a nonrotating string the terms in Eq. (7) with space components γ_k are suppressed by powers of velocity of the endpoints of the string. In what follows we shall keep for simplicity only the component $N_{44}^{(2)} \equiv J(x, y)$ of $N_{\mu\nu}^{(2)}$. Hence the kernel M is proportional to the FC of the color-electric field $E_i \equiv F_{i4}$. It is the dominant part of the string. Color-magnetic components are neglected in this first step and can be considered as a correction.

In this way one arrives at the system of equations [8] where we keep the same notation $M(x, y)$ for the retained piece of the kernel,

$$iM(x, y) = J(x, y) \gamma_4 S(x, y) \gamma_4, \quad (10)$$

$$(-i\hat{b}_x - im)S(x, y) - i \int M(x, z) S(z, y) d^4z = \delta^{(4)}(x - y). \quad (11)$$

B. Partial wave reduced equations

The kernel $J(x, y)$ in Eq. (10) depends on the time as $(x_4 - y_4)/T_g$. Therefore in the limit of $T_g \rightarrow 0$ it becomes local in time. As a result, the Fourier transform of M in the fourth coordinate does not depend on the momentum p_4 for $T_g \rightarrow 0$. The corresponding eigenfunctions $\psi_n(\mathbf{x}, p_4) = \psi_n(\mathbf{x})$ and eigenvalues $\varepsilon_n = \varepsilon_n(p_4)$ of the light quark in the white light-heavy configuration can readily be obtained from Eqs. (10), (11) in the discussed instantaneous limit. We find

$$\left(\frac{\boldsymbol{\alpha}}{i} \frac{\partial}{\partial x} + \beta m \right) \psi_n(\vec{x}) + \beta \int M(\mathbf{x}, \mathbf{z}) \psi_n(\mathbf{z}) d^3z = \varepsilon_n \psi_n(\mathbf{x}), \quad (12)$$

$$M(\mathbf{x}, \mathbf{z}) = J(\mathbf{x}, \mathbf{z}) \beta \Lambda(\mathbf{x}, \mathbf{z}), \quad \Lambda(\mathbf{x}, \mathbf{z}) = \sum_k \psi_k(x) \text{sgn } \varepsilon_k \psi_k^+(z). \quad (13)$$

The spherical-spinor decomposition of ψ_n for the total and orbital angular momentum channel $j, l = j \pm 1/2$,

$$\psi_n(r) = \frac{1}{r} \begin{pmatrix} G_n(r) \Omega_{j l M} \\ i F_n(r) \Omega_{j l' M} \end{pmatrix}, \quad l' = 2j - 1 \quad (14)$$

yields equations for the partial waves

$$\frac{dF_n}{dr} - \frac{\kappa}{r} F_n + (\varepsilon_n - m) G_n - M_{11} G_n - i M_{12} F_n = 0, \quad (15)$$

$$\frac{dG_n}{dr} + \frac{\kappa}{r} G_n - (\varepsilon_n + m) F_n - M_{22} F_n + i M_{21} G_n = 0, \quad (16)$$

with $\kappa = \pm(j + 1/2)$. Clearly the kernels M_{ik} are nonlocal in space, i.e.,

$$M_{ik} G_n = \int_0^\infty M_{ik}^{jj}(r, r') G_n(r') r r' dr', \quad (17)$$

with $M_{ik}^{jj} = \langle \Omega_{j l M} | M_{ik} | \Omega_{j l' M} \rangle$ and M_{ik} is given by Eq. (13).

Equations (15), (16) are invariant under the transformation

$$\varepsilon_n \rightarrow -\varepsilon_n, \quad \kappa \leftrightarrow -\kappa, \quad G_n \leftrightarrow F_n \quad (18)$$

which also yields $M_{11} \leftrightarrow M_{22}$, $M_{12} \leftrightarrow -M_{21}$. The symmetry (18) implies that the spectrum is symmetric in $\varepsilon_n \leftrightarrow -\varepsilon_n$, which is a property of a scalar interaction. The Lorentz scalar nature of the confining interaction has the nice feature that it does not lead to instability problems in the Dirac equation [14]. Moreover, nontrivial solutions of Eqs. (15), (16), if they exist, signify spontaneous CSB.

We solve Eqs. (15), (16) using the relativistic WKB approximation for the kernel M_{ik} . To simplify the calculations the Gaussian form for $D(x)$ was used (since all observables are integrals of FC, its explicit form is not essential at large distances, provided the FC have a finite range T_g and it yields the same value of the string tension σ)

$$D(u) = D(0) \exp(-u^2/4T_g^2), \quad D(0) = \frac{\sigma}{2\pi T_g^2} \quad (19)$$

with $\sigma = 0.2 \text{ GeV}^2$, and $T_g = 0.25 \text{ fm}$, taken in accordance with the lattice measurements [4,13].

As the reference basis we take the WKB solutions of the Dirac equation for the local linear confining potential σr , and the WKB computed kernels \tilde{M} and $\tilde{\Lambda}(\mathbf{x}, \mathbf{y})$ in Eq. (13) are determined by explicit summation over eigenstates. It was checked by an independent calculation [15] that the relativistic WKB procedure yields eigenvalues of the linear potential with accuracy better than 1%.

The general structure of M and \tilde{M} can be derived from Eqs. (13), (14). One can write M as a 4×4 matrix as follows [16]

$$M = M^{(0)} I + M^{(i)} \hat{\sigma}_i + M^{(4)} \gamma_4 + M^{(i)} \gamma_i, \quad (20)$$

where γ_i, γ_4 are the usual Dirac matrices, $i = 1, 2, 3$ and $\hat{\sigma}_i = \begin{pmatrix} \sigma_i & 0 \\ 0 & \sigma_i \end{pmatrix}$.

The same representation holds for the WKB approximated \tilde{M} . From the WKB analysis [8,16] we find that $\tilde{M}^{(0)}$ is the only growing kernel. It behaves asymptotically as

$$M^{(0)}(x, y) = \sigma \frac{|\mathbf{x} + \mathbf{y}|}{2} \tilde{\delta}^{(3)}(\mathbf{x} - \mathbf{y}), \quad (21)$$

where $\tilde{\delta}^{(3)}(\mathbf{r})$ is a smeared δ function with the range of non-locality decreasing asymptotically with growing $|\mathbf{x}|, |\mathbf{y}|$. The term $M^{(i)}$ is proportional to the angular momentum L and asymptotically it behaves as $O(1/x)$. One can also prove that $M^{(4)}, M^{(i)}$ do not grow at large x, y [16]. Hence in all problems where large distances are dominant one can consider only the first term in Eq. (20). Using this approximation we get, for the kernel (13),

$$\tilde{M} = \tilde{M}^{(0)}(\mathbf{x}, \mathbf{y}) I = J(\mathbf{x}, \mathbf{y}) \tilde{\Lambda}(\mathbf{x}, \mathbf{y}) I, \quad (22)$$

with

$$J(\mathbf{x}, \mathbf{y}) = \sigma \frac{\mathbf{xy}}{\sqrt{\pi}} f(\mathbf{x}, \mathbf{y}), \quad f(\mathbf{x}, \mathbf{y}) = \int_0^1 ds \int_0^1 dt e^{-(xs-yt)^2/4T_g^2} \quad (23)$$

and where $\tilde{\Lambda}$ is now a scalar quantity.

A curious feature of the considered equations is the way the string connecting the light quark to the source is being created. Actually, if one takes only lowest partial waves inside the kernel M [i.e., in $\Lambda(\mathbf{x}, \mathbf{y})$, Eq. (13)], then the effective potential in Eqs. (15), (16) is not confining. If one, however, sums up over all *angular states and radial excitations* in Λ , then the resulting $\Lambda(x, y)$ is a smeared δ function leading to the quasilocal confining kernel \tilde{M} . For example, using eigenfunctions for the local case, $\tilde{\Lambda}$ can be computed quasi-classically to be

$$\tilde{\Lambda}(\mathbf{x}, \mathbf{y}) = \frac{\sigma^2 xy}{2\pi^2} \frac{K_1[\sigma\sqrt{xy}\sqrt{(x-y)^2 + \theta^2 xy}]}{\sqrt{(x-y)^2 + \theta^2 xy}}, \quad \mathbf{xy} = xy \cos \theta. \quad (24)$$

In Eq. (24) one can clearly see that $\tilde{\Lambda}(x, y)$ is a normalized smeared δ function, with smearing radius in $|x-y|$ being $1/\sigma\sqrt{xy}$. For large distances it is nonvanishing only in the forward direction.

Insertion of this $\tilde{\Lambda}$ into \tilde{M} , Eq. (22) produces linear confinement due to the kernel $\tilde{\Lambda}(\mathbf{x}, \mathbf{y})$, as is given by Eq. (24) [while simply averaging the contribution from each individual orbital in Eq. (13) over the angle between \mathbf{x} and \mathbf{y} would produce no confinement at all]. The computed kernel $\tilde{M}(x, y)$ turns out to be nonlocal, but very close to the linear potential at large distances. Indeed, the effective localized potential defined as

$$V_{\text{eff}}(r) = \int M^{(0)}(r, x) r x dx \quad (25)$$

approaches at very large distances, i.e., for $\sigma^{1/2}r > 200$, a linear dependence with a slope given by σ . However at shorter distances V_{eff} also looks linear over a relatively large

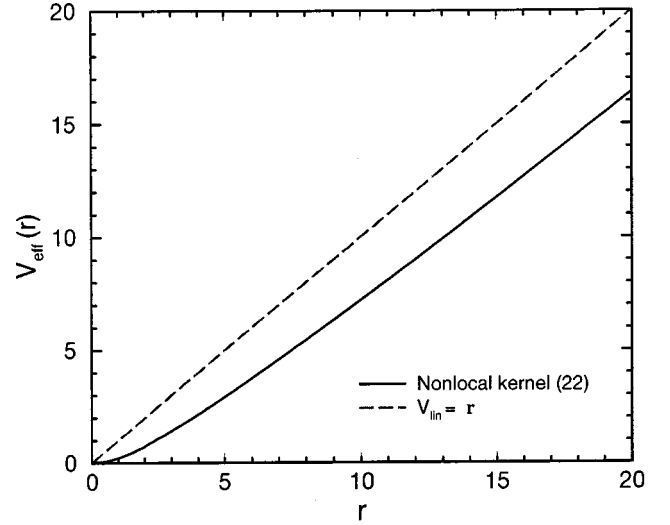


FIG. 1. Behavior of the effective potential, $V_{\text{eff}}(r) = \int \tilde{M}_{11}(r, x) r x dx$, as a function of r (solid line). The kernel (24) and units of $\sqrt{\sigma}$ have been used. For comparison, the linear local potential $V_{\text{lin}}(r) = r$ (dashed line) is also plotted.

region with a slope almost the same as σ , reflecting the presence of a small local curvature. In particular, we find that in the region $5 < \sigma^{1/2}r < 20$ the effective potential can reasonably well be described by $V_0(r) = 0.9\sigma r - 1.8\sigma^{1/2}$. In Fig. 1 are shown the results up to $r = 20$ in units of $\sigma^{1/2}$.

Since it is only the higher states and large distances which are important in the creation of this δ -function-type behavior of $\Lambda(x, y)$, and since the WKB method does well for high states and at large distances, one can clearly conclude that linear confinement should be obtained if one sums over all exact solutions of Eqs. (15), (16). Hence this should be a property of the exact solution.

The property that the kernel has a focusing effect in the forward direction can be used to get a somewhat simpler form. For this purpose we may also use [8]

$$\tilde{\Lambda}(\mathbf{x}, \mathbf{y}) = \frac{\sigma}{\pi^2 \sqrt{xy}} K_0(a) \delta(\cos \theta_x - \cos \theta_y), \quad a = \sigma \sqrt{xy} |x - y|. \quad (26)$$

TABLE I. Energy eigenvalues for Eqs. (13), (14) with $J = \frac{1}{2}$ and the kernels, given by Eq. (24) and Eq. (26), as compared with those of the Dirac equation for the linear local potential $V_{\text{lin}}(r) = r$.

n	$V = r$		kernel (24)		kernel (26)	
	$\kappa = -1$	$\kappa = 1$	$\kappa = -1$	$\kappa = 1$	$\kappa = -1$	$\kappa = 1$
0	1.619	2.294	0.925	1.472	0.969	1.516
1	2.603	3.031	1.719	2.056	1.765	2.113
2	3.291	3.626	2.277	2.541	2.334	2.608
3	3.855	4.138	2.740	2.964	2.809	3.042
4	4.345	4.594	3.144	3.342	3.226	3.432
5	4.784	5.008	3.507	3.685	3.603	3.790
6	5.186	5.334	3.838	4.002	3.950	4.123

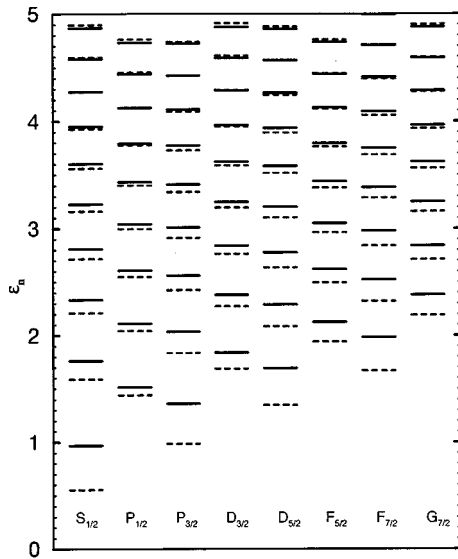


FIG. 2. Level structure calculated with the Dirac equation using the kernel (26) (solid line) in comparison with the predictions of the linear potential $V_0(r) = 0.9r - 1.8$ (dashed line). States carry the quantum numbers L_J , with L, J being the orbital and total angular momentum.

The eigenvalues and eigenfunctions in Eqs. (15), (16) have been determined using the kernel $\tilde{M}(x, y)$, given by Eqs. (24) and (26). Some results are shown in Table I and Fig. 2. In our present study we have taken $m = 0$. Note that $\tilde{M}(x, y)$ is approaching a local linear potential at large distances, $x, y \gtrsim \sigma^{-1/2}$, which justifies *a posteriori* our choice of the reference basis. Due to the nonlocality of the interaction the predicted spectrum is found to be different from that of the linear potential, valid at large distances. Moreover, comparing the level structures of the $J = 1/2$ channel as obtained using the kernels (24) and (26) we see from Table I that they are qualitatively very similar, corroborating that there is indeed a strong forward focusing effect in the quark propagator. From Fig. 2 we see that for all L values the higher radially excited levels are close to the predictions of linear potential $V_0(r) = 0.9\sigma r - 1.8\sigma^{1/2}$, in agreement with the fact that the interaction at large distance can indeed be described by a local linear potential. On the other hand, the nonlocal kernel predictions for the low-lying states clearly deviates strongly from those of the (shifted) linear potential. Hence the nonlocal nature of the force does affect the spectrum in an essential way.

The eigenfunctions for the nonlocal kernels look qualitatively similar to the corresponding ones of the shifted linear potential. In Fig. 3 are shown the ground state and first excited state for the $J = 1/2$ channels. Although the differences

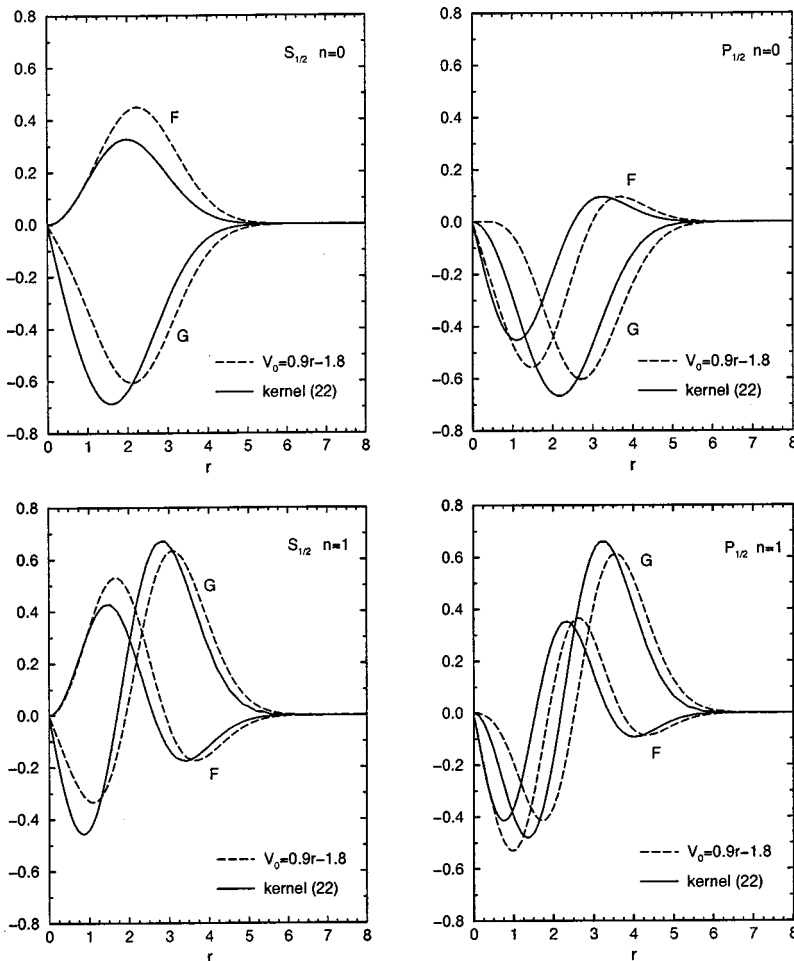


FIG. 3. Eigenfunctions of the ground state and the first excited state for the $S_{1/2}$ and $P_{1/2}$ channels. The solutions correspond to using the kernel (26) (solid line) and the shifted linear potential $V_0(r) = 0.9r - 1.8$ (dashed line).

are substantial for these low-lying states, the agreement for higher excited states is considerably better. Moreover, we find that the large distances and high states of the WKB states agree well with the corresponding eigenfunctions.

C. Inclusion of perturbative exchanges

Till now we have considered only the NP part of the gluonic field, B_μ . In this section we include the perturbative part, a_μ , and neglect for simplicity the interference terms. Therefore the effect of a_μ is accounted for in the appearance of an additional factor in the partition function (3), namely,

$$Z = Z_{NP} Z_{\text{pert}}, \quad Z_{\text{pert}} = \langle e^{g \int dx \psi^\dagger \hat{a}(x) \psi(x) + i g \int dz_4 a_4(z_4)} \rangle_a, \quad (27)$$

where the second term in the exponent of Eq. (27) corresponds to the interaction of the perturbative part of the gluon field with the static antiquark. We have used in Eq. (27) that due to the 't Hooft identity [17] one can average independently over B_μ and a_μ .

The result of averaging yields a new additive term in L_{eff} , Eq. (3),

$$L_c = g \int dx \psi^\dagger(x) \hat{A}^{(c)}(x) \psi(x), \quad (28)$$

where we have defined

$$A_\mu^{(c)}(x) = -i g \int dz_4 \langle a_4(x) a_4(z_4) \rangle = \delta_{\mu 4} \frac{(-i) g C_2}{4 \pi |\mathbf{x}|}. \quad (29)$$

The presence of L_c in Eq. (3) does not influence the derivation of basic Eqs. (10), (11). The only difference is that Eq. (11) assumes the form

$$[-i \hat{\partial} - g \hat{A}^c(x) - im] S(x, y) - i \int M(x, z) S(z, y) d^4 z = \delta^{(4)}(x - y). \quad (30)$$

Equation (10) does not change and the kernel $J(x, y)$ contains as before only nonperturbative contributions. Note, however, that $S(x, y)$ in $M(x, y)$ in Eq. (10) now contains also perturbative gluon exchanges. This is a new type of interference of perturbative and NP terms, which appears irrespective of our neglect of this interference within the av-

eraging procedure over B_μ and a_μ . In other words, another class of diagrams is responsible for this interference.

Correspondingly in the static equations (12) one should replace

$$\beta m \rightarrow \beta m - \frac{C_2 \alpha_s}{|\mathbf{z}|}. \quad (31)$$

The equations for the partial waves (15), (16) are modified due to the presence of the color Coulomb potential $V(r)$ in a simple way. Since

$$V(r) = - \frac{C_2 \alpha_s}{r} \quad (32)$$

is local and a Lorentz vector, it always appears in the combination $\varepsilon_n - V(r)$. Hence one has instead of Eqs. (15), (16)

$$\frac{dF_\nu}{dr} - \frac{\kappa}{r} F_\nu + [\varepsilon_\nu - V(r) - m] G_\nu - M_{11} G - i M_{12} F = 0, \quad (33)$$

$$\frac{dG_\nu}{dr} + \frac{\kappa}{r} G_\nu - [\varepsilon_\nu - V(r) + m] F_\nu - M_{22} F + i M_{21} G = 0, \quad (34)$$

where we have denoted

$$M_{ik} \left(\frac{G}{F} \right) \equiv \int \langle \nu | M_{ik} | \nu' \rangle \left(\frac{G_{\nu'}(w)}{F_{\nu'}(w)} \right) r w dw. \quad (35)$$

Here M_{ik} is defined as in Eq. (13) and the matrix Λ_{ik} in Eq. (13) involves the sum over all states, including positive and negative ε_n . There in Sec. II B we have exploited the symmetry (18). However, Eqs. (33), (34) are invariant under another transformation, namely,

$$\varepsilon_n \leftrightarrow -\varepsilon_n, V(r) \leftrightarrow -V(r), \kappa \leftrightarrow \kappa, G_n \leftrightarrow F_n. \quad (36)$$

Now the sum over negative ε_n can be expressed through the corresponding sum over positive ε_n with exchange $G_n \leftrightarrow F_n$ as before, but also with the inversion of sign of Coulomb interaction, i.e., Coulomb attraction for positive ε_n is replaced by Coulomb repulsion for negative ε_n .

In what follows we shall denote wave functions of the positive energy states with repulsive Coulomb with the sign of tilde: $\tilde{G}_\nu, \tilde{F}_\nu$. Then using Eq. (13) the matrix $\beta \Lambda_{ik}$ can be written as a sum over only positive ε_n as follows:

$$\beta \Lambda_{ik}^{\mu\mu'} = \frac{1}{xy} \sum_{jIM, n>0} \begin{pmatrix} G_\mu G_{\mu'}^* - \tilde{F}_\mu \tilde{F}_{\mu'}^*, & -i(G_\mu F_{\mu'}^* - \tilde{F}_\mu \tilde{G}_{\mu'}^*) \\ -i(G_\mu G_{\mu'}^* - \tilde{G}_\mu \tilde{F}_{\mu'}^*), & \tilde{G}_\mu \tilde{G}_{\mu'}^* - F_\mu F_{\mu'}^* \end{pmatrix}. \quad (37)$$

Since $\beta \Lambda$ is exactly the combination which enters the mass matrix (13), one can list in Eq. (37) scalar and vector (proportional to β) parts:

$$M = M_s I + M_v \beta + \Delta M, \quad (38)$$

TABLE II. Ground state energy eigenvalue (in units of $\sqrt{\sigma}$) for Eqs. (33), (34) with $\sigma_s=0, 0.3$, and 0.39 and quark masses $m = 5$ MeV, 0.15 GeV, and 0.2 GeV (upper, middle, and lower entry) for different values of T_g , $T_g=0, 0.25, 0.5$, and 1 (in units of $1/\sqrt{\sigma}$).

T_g α_s	0	0.25	0.5	1
	1.628	0.985	0.979	0.907
0	1.886	1.225	1.217	1.145
	1.978	1.314	1.305	1.233
	1.163	0.684	0.679	0.628
0.3	1.378	0.884	0.877	0.826
	1.456	0.959	0.951	0.900
	1.004	0.585	0.580	0.536
0.39	1.201	0.768	0.761	0.717
	1.272	0.837	0.830	0.786

where ΔM contains spin-dependent terms, which can be considered as in Sec. II B, while M_s, M_v are

$$M_{s,v} = C \sum_{j|lM_{\mu,n}>0} [G_{\mu}G_{\mu}^* - \tilde{F}_{\mu}\tilde{F}_{\mu}^* \pm (\tilde{G}_{\mu}\tilde{G}_{\mu} - F_{\mu}F_{\mu}^*)], \quad (39)$$

where

$$C = \frac{1}{4} \sqrt{\pi} T_g D(0) \frac{\mathbf{x} \cdot \mathbf{y}}{xy} f(\mathbf{x}, \mathbf{y}),$$

$$f(\mathbf{x}, \mathbf{y}) = \int_0^1 ds \int_0^1 dt \exp\left(-\frac{(\mathbf{x}s - \mathbf{y}t)^2}{4T_g^2}\right). \quad (40)$$

From Eq. (39) it is clear that the vector part M_v is only due to the presence of Coulomb interaction. Corrections at large distances due to the vector part can be treated again in the relativistic WKB. A rough estimate of M_v at large r yields

$$\frac{M_v}{M_s} \sim \frac{\alpha_s}{\sigma r^2} \quad (41)$$

and hence can be neglected at large enough r .

III. NUMERICAL SOLUTIONS OF EQUATIONS AND COMPARISON TO B, D MESONS

We have performed numerical studies of Eqs. (33), (34) with the kernel (24) for different values of the quark mass m and different values of T_g . To simplify calculations only the dominant part of the mass operator M_{ik} was retained, i.e., $\hat{M}_{11} = M_{22} = M^{(0)}$, while M_{12}, M_{21} have been neglected. For $M^{(0)}$ the representation (22) was used

$$M^{(0)}(\mathbf{x}, \mathbf{y}) = J(\mathbf{x}, \mathbf{y}) \tilde{\Lambda}(\mathbf{x}, \mathbf{y}) I,$$

where $\tilde{\Lambda}$ is taken to be the kernel (24). Results of our calculations for the ground state energy are listed in Table II. One can see a rather sharp change of energy when α_s changes

from 0 to 0.3 and when T_g is changing from 0 to 0.25, while further increase of α_s or T_g does not produce such a strong dependence.

Solutions of our equations (33), (34) can be compared with physical states of B, D and B_s, D_s mesons. To this end one should have in mind that in Eqs. (33), (34) the static approximation for the heavy quark b, c was used, and hence all corrections $O(1/m_Q^n)$ with $n \geq 1$ are neglected.

One can exploit at this point the HQET expansion for the mass m_H of heavy-light boson [18,19]

$$m_H = m_Q \left[1 + \frac{\bar{\Lambda}}{m_Q} + \frac{1}{2m_Q^2} (\lambda_1 + d_H \lambda_2) + O(1/m_Q^3) \right], \quad (42)$$

where λ_n are free parameters, depending on dynamics, and d_H is the hyperfine splitting parameter. It is clear from the preceding that eigenvalues of Eqs. (33), (34) yield the value $\bar{\Lambda}$, which depends on the quantum numbers of the state,

$$\bar{\Lambda}(j, l, n_r) = \varepsilon_n(j, l). \quad (43)$$

Consider now the results of the present approach, i.e., solutions of Dirac-type equations (33), (34). In the local case ($T_g \rightarrow 0$) when the kernel M reduces to the linear potential σr , we have

$$\bar{\Lambda}_D^{(loc)} = 0.690 \text{ GeV} \quad (\alpha_s = 0, \sigma = 0.18 \text{ GeV}^2) \quad (44)$$

and

$$\bar{\Lambda}_D^{(loc)} = 0.493 \text{ GeV} \quad (\alpha_s = 0.3, \sigma = 0.18 \text{ GeV}^2). \quad (45)$$

This should be compared to the nonlocal case

$$\bar{\Lambda}_D^{(nonloc)} = 0.415 \text{ GeV} \quad (\alpha_s = 0, \sigma = 0.18 \text{ GeV}^2) \quad (46)$$

and

$$\bar{\Lambda}_D^{(nonloc)} = 0.288 \text{ GeV} \quad (\alpha_s = 0.3, \sigma = 0.18 \text{ GeV}^2). \quad (47)$$

These latter values are in general agreement with the results of the QCD heavy-flavor sum rules [20,21]

$$\bar{\Lambda} = 0.57 \pm 0.07 \text{ GeV} \quad (48)$$

and more recent analysis from semileptonic B decays [22]

$$\bar{\Lambda} = 0.39 \pm 0.11 \text{ GeV}. \quad (49)$$

Another interesting comparison is with the experimental values of the B -meson mass [the term λ_2 in Eq. (42) can be determined from the $B^* - B$ mass difference]. Using Eq. (47) and $\bar{M}_B = (3M_{B^*} + M_B)/4 = 5.312 \text{ GeV}$ one can estimate (neglecting λ_1) the pole mass of the b quark to be $m_b(\text{pole}) \cong 5.0 \text{ GeV}$, which is in reasonable agreement with the analysis of the quarkonium spectra in [23].

TABLE III. Energy eigenvalues $\bar{\Lambda}$ of the heavy-light system in the static heavy quark approximation obtained in different approaches.

Refs.	Method	$\bar{\Lambda}$ (GeV)
[20]	QCD sum rules	>0.5
[21]	QCD sum rules	$0.4-0.5$
[24]	Lattice	0.18 ± 0.03
[22]	experim.	0.33 ± 0.11
[25]	QCM	0.35
[26]	QCM	$0.5-0.6$
[27]	Rel. QCM	0.386
This work	Nonlin. Dirac eq.	0.287

A similar analysis can be done for the B_s meson; the corresponding values for $\bar{\Lambda}_s$ with $m_s=0.15$ and 0.20 GeV, are $\bar{\Lambda}_s - \bar{\Lambda} = 0.084$ and $\bar{\Lambda}_s - \bar{\Lambda} = 0.115$ for $\alpha_s = 0.3$. One can compare these values with the mass difference $B_s, B^0, \Delta M_s(B) = (0.090 \pm 0.0038)$ GeV. These numbers for $\bar{\Lambda}$ can be compared with those in Table III, where also results of lattice calculations [24] and of the constituent quark model (CQM) [25–27] are given.

IV. CHIRAL CONDENSATE

As a check of CSB in our Eqs. (15), (16) we have computed the chiral condensate, which can be expressed through the eigenfunctions as in [8] (to simplify the matter we disregard in this section perturbative contributions).

$$\langle \bar{q}q \rangle = -\frac{N_c}{2\pi} \sum_{n=0}^{\infty} [(A_n^-)^2 - (B_n^+)^2], \quad (50)$$

where $A_n^- = (G_n(r)/r)_{r=0}$, $B_n^+ = (F_n(r)/r)_{r=0}$, and G_n, F_n refer to solutions with $\kappa = -1, l=0$, and $\kappa = +1, l=1$, respectively. In the local linear potential case the values of A_n^-, B_n^+ have been computed in the WKB method [8] and shown to yield a monotonically divergent series $\langle \bar{q}q \rangle = -N_c/2\pi \sum_n \text{const}/\sqrt{n}$.

TABLE IV. The difference $s_n = |A_n|^2 - |B_n|^2$ for $n=0, 1, \dots, 6$, in the case of the nonlocal kernels (24) and (26) and corrected for a normalized Gaussian nonlocality (51) with a range of $a=0.3$ and $a=0.5$. For comparison the results are shown for a local linear potential $V_{lin}=r$.

n	$V=r$	kernel (24)	$A_n^2 - B_n^2$		
			kernel (26)	$a=0.3$	$a=0.5$
0	0.79	0.42	0.50	0.15	0.23
1	0.51	0.21	0.34	0.04	0.12
2	0.41	0.12	0.19	0.02	0.10
3	0.35	0.10	0.16	0.02	0.09
4	0.31	0.08	0.11	0.01	0.09
5	0.27	0.07	0.11	0.01	0.08
6	0.26	0.06	0.09	0	0.07

It can be argued [using Eqs. (8), (9) from [8]], that the nonlocality of the kernel M in spacetime, present by definition in Eq. (8), improves the convergence of the series and yields a finite result for $\langle \bar{q}q \rangle$. We have found A_n^-, B_n^+ from the solutions of the nonlocal equations (15), (16) with the kernel \tilde{M} and compared them with the local case, when \tilde{M} reduces to the local linear potential. Results are shown in Table IV.

One can see from the results that in the nonlocal case the magnitude of $s_n \equiv (A_n^-)^2 - (B_n^+)^2$ is clearly diminished as compared to the reference local case, and is of reasonable order of magnitude. From the obtained sequence of s_n we get that $\langle \bar{q}q \rangle = -0.5\sigma^{3/2}$ and $-0.7\sigma^{3/2}$ in the nonlocal cases of the kernels (21) and (22), respectively. Adopting a value of $\sigma = 0.2 \text{ GeV}^2$ we find $\langle \bar{q}q \rangle = -(350 \text{ MeV})^3$ and $-(400 \text{ MeV})^3$, respectively, to be compared with the usually acceptable value of $-(250 \text{ MeV})^3$. However convergence is still slow as seen from Table IV and the converged values are somewhat higher. We have checked that convergence is somewhat improved when one takes into account the intrinsic nonlocality of the kernel M in \mathbf{x}, \mathbf{y} . To this end we have modified the kernel \tilde{M} obtained from WKB analysis, replacing $\tilde{\delta}$ in Eq. (21) by a Gaussian factor

$$N \exp\left(-\frac{(x-y)^2}{a^2}\right) \delta(\cos \theta_x - \cos \theta_y) \quad (51)$$

and studied the sequences of s_n as functions of the nonlocality range a . Results are shown in Table IV for two values of $a = 0.3\sigma^{1/2}$ and $0.5\sigma^{1/2}$. The strength N is chosen such that numerically the slope of $\sigma = 0.2 \text{ GeV}^2$ is reproduced for large distances. The condensate values vary in the considered region substantially, showing that effects of nonlocality are important.

The slope of the effective potential $V_{\text{eff}}(r)$, determined by N , strongly depends on a for $a \cong T_g$. We believe that the reason for this lies in the fact that the chiral condensate $\langle \bar{q}q \rangle$ depends crucially on the nonlocality both in time components of $M(\mathbf{x}, \mathbf{y}; x_4, y_4)$ and in spatial components. The first nonlocality was, however, disregarded in Eqs. (12), (13), when the p_4 dependence was omitted in M and ψ_n (the static limit). It was indeed shown in Ref. [8] that taking this de-

pendence into account significantly improves convergence of the sum in Eq. (50). The full account of this effect requires solution of time-dependent Eqs. (10), (11), which is numerically a much more difficult problem.

V. CONCLUSION

We have studied the confining and CSB properties in the system of one light quark and one static antiquark. The effective mass operator is written explicitly for large N_c , as a sum over vacuum field correlators. Keeping only the Gaussian field correlator, we have obtained a closed system of equations in the limit of large N_c . Our results support the presence of a Lorentz scalar linear confinement for the light quark, which signifies CSB for this system, and yield eigenfunctions and eigenvalues for the heavy-light system containing both confinement and CSB.

As a direct evidence of CSB we have computed the chiral condensate, which appears to be of the correct sign and having the proper large N_c dependence. Our result yields a reasonable order of magnitude of $\langle \bar{q}q \rangle$, provided convergence of the sum is achieved. At this point it is useful to compare the CSB picture of the Nambu–Jona–Lasinio model and our

approach. In the NJL model confinement and string are absent and CSB may occur due to the condensation of $q\bar{q}$ pairs in the scalar channel. In our case, being the large N_c approximation of the real QCD, a string is built up between light and heavy quark, which depends not only on light quark coordinates \mathbf{x} , \mathbf{y} , but also on the distance from them to the heavy antiquark. In the presence of confinement, the phenomenon of CSB is due to the spontaneous creation of the scalar string, which is forbidden by chiral symmetry.

Eigenvalues ε_n and eigenfunctions obtained numerically for lowest states represent the leading contributions of the HQET expansion in powers of $1/m_Q$. Results for the energies ε_n in our method are compared of the lattice and QCD sum rule calculations, and also with experimental extraction of $\varepsilon_n = \bar{\Lambda}(n)$, showing an overall agreement with the B and D meson masses.

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