Statistical properties of the charmonium spectrum and a new mechanism of J/ψ suppression

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The statistical properties of the charmonium energy spectrum determined by the Bethe-Salpeter equation are investigated. It is found that the regular motion of the $c\bar{c}$ system can be expected at a small value of color screening mass but the chaotic motion is expected at a large one. It is shown that the level mixing due to color screening serves as a new mechanism resulting in the J/ψ suppression. Moreover, this kind of suppression can occur before the color screening mass reaches its critical value for J/ψ dissociation. This implies that a strong J/ψ suppression is possible in the absence of dissociation of J/ψ . [S0556-2813(99)03509-8]

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Strongly interacting matter of sufficiently high density is predicted to undergo a transition to a state of deconfined quarks and gluons. Deconfinement occurs when color screening shields a given quark from the binding potential of any other quarks or antiquarks. Bound states of heavy quarks, such as the J/ψ or the Y, whose radii are much smaller than those of the usual mesons and nucleons, can survive in a deconfined medium until the temperature or density becomes so high that color screening prevents their tighter binding. It was theoretically predicted that a suppression of J/ψ or ψ' production due to the dissociation can be found in relativistic heavy ion collision, which can serve as a clear signature for the formation of quark-gluon-plasma [1]. Subsequently this suppression was observed by the NA38 Collaboration [2]. However, successive research pointed out that such suppression could also exist in hadronic matter, even though caused by a completely different mechanism [3]. Recently, an anomalously strong J/ψ suppression has been observed by the NA50 Collaboration [4] and there have been a number of attempts (see, for example, Refs. [5-7]) to explain it. However, the mechanism of J/ψ suppression is still a question of debate.

It is known that, for a complex system, the statistical properties of the quantum spectrum and the complexity of the eigenfunctions reveal frequently new features of the system. Because of the complexity of the interaction and the temperature and density effect, the heavy $q\bar{q}$ systems in hot matter can be regarded as complex systems. Nevertheless, the spectral statistical properties and the complexity of eigenfunctions of heavy $q\bar{q}$ systems have not yet been discussed.

In this paper we study the spectral statistical properties and the complexity of eigenfunctions of the $c\bar{c}$ system. The influence of color screening on the statistical properties and the mixing of eigenfunctions which characterizes the complexity of eigenfunctions will be examined. And an alternative interpretation of the strong J/ψ suppression is proposed.

To investigate the statistical properties of the $c\bar{c}$ system a reasonable energy spectrum is crucial. Usually the bound

states of the $c\bar{c}$ system were studied in the nonrelativistic formulation [8,9]. However, the motion of a quark and an antiquark in a meson should be relativistic. As pointed out in Ref. [10], the kinetic energy of the $c\bar{c}$ system is about 13% of the total energy and the ratio of relativistic corrections to the quark mass will not decrease with the increase of quark mass. As a result the bound state equation for J/ψ should be relativistic. Since the Bethe-Salpeter (BS) equation [11] is the unique effective relativistic equation of two-body bound states and is consistent with quantum field theory, we take the BS equation to determine the energy spectrum of charmonium in this paper.

It is known that the full bound state BS equation, written in the two-sided notation, reads [12]

$$(\eta_1 \mathbf{P} + \mathbf{p} - m_1) \chi_P(p) (\eta_2 \mathbf{P} - \mathbf{p} + m_2) = \frac{i}{(2\pi)^4} \int d^4 p' V(p, p'; P) \chi_P(p'), \qquad (1)$$

where $\eta_i = m_i / (m_1 + m_2)$ (*i*=1,2), χ_P is the wave function for the quark-antiquark system with total four momentum *P*, *p* is the relative four momentum. *V* is the interaction kernel that acts on χ_P and formal products $V\chi_P(p')$ take the form $V\chi_P(p') = V_s\chi_P(p') + \gamma_\mu \otimes \gamma^\mu V_v\chi_P(p')$, in which V_s and V_v are the scalar, vector potential, respectively.

With the standard reduction and spin-independent treatment being implemented, the three-dimensional spinindependent reduced Salpeter equation can be written as [13],

$$(M - E_1 - E_2)\phi(\vec{p}) = \int \frac{d^3p'}{(2\pi)^3} \sum_{i=s,v} F_i^{\text{SI}}(\vec{p}, \vec{p}') V_i(|\vec{p} - \vec{p}'|)\phi(\vec{p}'), \quad (2)$$

where $\phi(\vec{p}) = \int dp^0 \chi_P(p^0, \vec{p})$ is the three-dimensional equaltime BS wave function. *M* is the mass of the $q\bar{q}$ bound state. $E_i = (\vec{p}^2 + m_i^2)^{1/2}$, i=1,2 represent a quark and an antiquark,

TABLE I. The calculated spin-averaged mass spectra M_{nl} of the $c\bar{c}$ bound states and the comparison with experimental data and previous calculations.

			Cal. (GeV)		
	nl	Exp. (GeV)	Ref. [8]	Ref. [9]	Ours
$c\overline{c}$	1S	$J/\psi(3.068)$	3.070	3.070	3.067
	2S	$\psi'(3.663)$	3.698	3.686	3.663
	35	$\psi''(4.025)$	4.170	4.081	4.019
	1 <i>P</i>	$\chi_c(3.525)$	3.500	3.505	3.526

respectively. The functions F_v^{SI} and F_s^{SI} are the transformation coefficients of the spin-independent vector, scalar potential, respectively, when the four-dimensional BS equation is reduced to the three-dimensional spin-independent BS equation. Their expressions have been given explicitly in Ref. [13].

It is definite that the momentum dependence of the interaction is treated exactly in Eq. (2). Equation (2) gives a welldefined eigenvalue problem for the masses M of the $q\bar{q}$ bound states in the momentum space. To solve Eq. (2), we take the color screening potential between the quark and antiquark [8]. In momentum space the color screening potential is written as

In our calculation, the reduced BS equation which has been described well in Refs. [12,13] and a color screened quark-antiquark potential [13] are implemented. In a momentum space, the potential with the color screening can be written as

$$V_{s}(|\vec{p}-\vec{p'}|) = \frac{\sigma}{\mu} \delta^{3}(\vec{p}-\vec{p'}) - \frac{\sigma}{\pi^{2}} \frac{1}{[(\vec{p}-\vec{p'})^{2} + \mu^{2}]^{2}}$$
(3)

and

$$V_{v}(|\vec{p}-\vec{p}'|) = -\frac{2}{3\pi^{2}}\frac{\alpha_{s}}{[(\vec{p}-\vec{p}')^{2}+\mu^{2}]},$$
 (4)

where V_s , V_v refer to the scalar, vector potential respectively, and σ is the string tension, α_s the effective coupling constant, and μ the Debye screening mass. The parameters including the value of the μ at zero temperature (denoted as μ_0 hereafter) are determined by a least square fitting to the experimental data of 1*S*, 2*S*, and 3*S* states of the $c\bar{c}$ system. The fixed parameters are $\sigma = 0.22 \text{ GeV}^2$, $\mu_0 = 0.06 \text{ GeV}$, $m_c = 1.474 \text{ GeV}, \alpha_s(c\bar{c}) = 0.47$, which are within the scope of customary usage. Some of the obtained spin-averaged energy (mass) [13] spectrum of the $c\bar{c}$ bound states at zero temperature (T=0) are listed in Table I. From Table I, one can observe that the mass spectrum obtained at present is more consistent with the experiment than the previous ones [8,9]. With the parameters σ , m_c , and $\alpha_s(c\bar{c})$ fixed above, we evaluate the mass spectrum of the $c\bar{c}$ bound states for different μ , so that the temperature dependence of the spectrum can be investigated. The calculation indicates that the critical value of the screening mass for J/ψ to be dissociated is about $\mu_c = 0.90$ GeV.

As the spectrum has been determined, we study the spectral statistics. To this end we exploit the well-established random matrix theory (RMT) [14,15]. The RMT was initially used to describe the statistical properties of complex nuclear spectra in the early of 1950's [14]. In the RMT, the matrix of the Hamiltonian of the system, which is sufficiently complex and (or) governed by a very complicated or even unknown dynamics, is replaced by a random matrix with well-defined symmetry properties. Any desired quantity of the system is determined by performing an average over an ensemble of the random matrices of which the elements are distributed according to a certain probability distribution. The distribution is usually taken to be a Gaussian with a parameter d,

$$P(H) \propto \exp(-\operatorname{Tr}\{HH^+\}/2d^2), \tag{5}$$

and the corresponding random matrix ensemble is called the Gaussian ensemble. The underlying space-time symmetries obeyed by the system impose some important restrictions on the admissible ensemble. If the Hamiltonian is time-reversal and rotational invariant, the Hamiltonian matrices can be chosen to be real symmetric. The corresponding ensemble is called the Gaussian orthogonal ensemble (GOE). If the Hamiltonian is not time-reversal invariant, irrespective of its behavior under rotations, the Hamiltonian matrices are complex Hermitian. The corresponding ensemble is called the Gaussian unitary ensemble (GUE). If the system is time-reversal invariant but not invariant under rotations, and it has half-odd-integer total angular momentum, the matrices are quaternion real. The corresponding ensemble is called the Gaussian symplectic ensemble (GSE).

If the Hamiltonian possesses other invariances, the corresponding ensemble can be the Gaussian unitary ensemble, or the Gaussian symplectic ensemble. The RMT has been applied successfully in many areas of physics, such as nuclear physics [16], quantum field theory [17], and the physics of disordered systems [18]. The success of the RMT is rooted in the fact that there exist so-called universal quantities which do not depend on the details of the dynamics but only on the underlying symmetries. This enables one to separate generic features from properties that do depend on the details of dynamics. The universal quantities usually mean the nearest neighbor level spacing distribution P(s) and the spectral rigidity $\Delta_3(L)$ which are powerful to analyze the spectral statistical properties of complex systems.

In present case the $c\bar{c}$ system is time-reversal and rotational invariant. It is then a member of the GOE and its spectrum should be analyzed by comparing with the predictions of the GOE. To make comparison with the universal and dimensionless results of the RMT, it is necessary to make a transformation or normalization on the spectrum. This operation is called the unfolding. The procedure is as follows. For a spectrum $\{E_i\}$, one separates its smoothed average part from its fluctuating part at first, then counts the number of the levels below E so that one can define a staircase function N(E) of the spectrum

$$N(E) = N_{\rm av}(E) + N_{\rm fluct}(E).$$

The unfolded spectrum can be finally obtained with the mapping

$$\widetilde{E}_i = N_{\mathrm{av}}(E_i)$$
.

This unfolded levels \tilde{E}_i are obviously dimensionless and have a constant average spacing of 1, but the actual spacings exhibit frequently strong fluctuations. The nearest neighbor level spacings are defined as $s_i = \tilde{E}_{i+1} - \tilde{E}_i$. The distribution P(s) is defined as that P(s) ds is the probability that s_i lies within the infinitesimal interval [s, s+ds].

It has been shown that the nearest neighbor level spacing distribution P(s) measures the level repulsion (the tendency of levels to avoid clustering) and short-range correlations between levels, which mean that, for a given level, its repulsion to the neighbor levels is limited within several level spacings [18,19]. For the GOE, the distribution can be expressed explicitly as $P(s) = (\pi/2)\exp(-\pi s^{2}/4)$, which can be deduced from Eq. (5) [15]. With the Brody parameter ω in the Brody distribution $P_{\omega}(s) = (1+\omega)\alpha s^{\omega}\exp(-\alpha s^{1+\omega})(\alpha = \{\Gamma[(2+\omega)/(1+\omega)]\}^{1+\omega}, \Gamma(x)$ is the Γ function), the transition from regularity to chaos can be measured quantitatively. It is evident that $\omega = 1$ corresponds to the GOE distribution, while $\omega = 0$ to the Poisson-type distribution. A value $0 < \omega < 1$ means an interplay between the regular and the chaotic.

As to the spectral rigidity $\Delta_3(L)$, it is defined as

$$\Delta_3(L) = \left(\min_{A,B} \frac{1}{L} \int_{-L/2}^{L/2} [N(x) - Ax - B]^2 dx \right), \quad (6)$$

where N(x) is the staircase function of a unfolded spectrum in the interval [-L/2,x]. The minimum is taken with respect to the parameters A and B. The average denoted by $\langle \cdots \rangle$ is taken over a suitable energy interval over x. Thus from this definition $\Delta_3(L)$ is the local average least square deviation of the staircase function N(x) from the best fitting straight line. For the GOE the expected value of $\Delta_3(L)$ can only be evaluated numerically, but it approaches the value $\Delta_3(L)$ $\approx 1/\pi^2(\ln L - 0.0687)$ for large L. It has also been shown that the spectral rigidity $\Delta_3(L)$ signifies the long-range correlations of quantum spectra [18,19] which make it possible that for a chaotic spectrum very small fluctuation of the staircase function around its average can be found in an interval of given length (the interval may cover dozens of level spacings). It has been generally accepted that if the spectral statistical properties of a quantum system approach to the GOE, the motion of the system is chaotic, and if the statistical properties appear as Poisson-types $\left[P(s) = \exp(-s)\right]$ and $\Delta_3(L) = L/15$, the motion is regular [18].

To perform a meaningful RMT analysis one has to sort the spectrum in symmetry sectors corresponding to the symmetry of the Hamiltonian. Since the symmetry invariant subspaces are orthogonal to one another, each such symmetry



FIG. 1. The spacing distribution P(s) and the spectral rigidity Δ_3 of the *s*-wave charmonium energy spectrum at different values of color screening masses μ measured in GeV. In (a) the dashed-line for Poisson distribution, the solid one for the GOE distribution, and the histograms for our numerical results.

invariant subspace is characterized by a specific set of quantum numbers. The RMT analysis is performed on sets of the eigenlevels having the same quantum numbers. In view of the above consideration we take the first 200 levels out of 300 s-wave (spin-weighted) eigenvalues which have the same quantum numbers (e.g., l=0). We have noticed that *p*-wave eigenvalues (spin-weighted) have the same statistical properties as those of the *s*-wave ones. The spacing distribution P(s) and spectral rigidity Δ_3 for several color screening masses μ are shown in Fig. 1. The figure shows obviously that, when μ is small, for instance 0.06 GeV, both the P(s)and Δ_3 display the Poisson-type distribution. This indicates that the spectrum is regular and the levels are uncorrelated. The motion of the $c\bar{c}$ system is then regular. With the increase of μ (for example, $\mu = 0.20$ GeV), the P(s) and Δ_3 depart from the Poisson types. Then the spectrum becomes less regular and the level correlation and repulsion increase gradually. As μ is large enough, say μ =0.60 GeV, both the P(s) and the Δ_3 appear the GOE behavior. This means that the spectrum becomes extreme irregular (chaotic) and the levels are strongly correlated. Strong level repulsion can thus be expected and the motion of the $c\bar{c}$ system is chaotic.

Fitting to the Brody distribution, we get the Brody parameter $\omega = 0.22, 0.51, 0.89$ for $\mu = 0.06, 0.20, 0.60$ (GeV), respectively. This also indicates that a transition from regular to chaotic motion can be induced in the $c\bar{c}$ system by the increasing of the color screening and the chaotic motion takes place when the hot matter is of high temperature. Recalling the idea that the P(s) and Δ_3 measures the correlation among the levels [18,19], one can recognize that the correlation of the $c\bar{c}$ resonances can be manifested by the P(s)and Δ_3 . The calculated result shows thus that the $c\bar{c}$ resonances are strongly correlated at a high temperature. This



FIG. 2. The probability density $P(c^2)$ of J/ψ versus c^2 for various values of μ . The solid histogram for μ =0.90 GeV, the dotted one for μ =0.60 GeV, the dash-dotted one for μ =0.20 GeV, and the dashed one for μ =0.10 GeV.

information may be useful for us to analyze the relevant experiments.

As a matter of fact, both the level correlation and level repulsion stem from level mixing [20] which characterizes the complexity of eigenfunctions. We analyze then the influence of the color screening on the level mixing. For this purpose, we take the *s*-wave eigenfunctions of the $c\bar{c}$ system at temperature zero $|k\rangle$ (J/ ψ , ψ' , ψ'' , ..., μ =0.06 GeV, equivalent to T=0) as a basis and expand the eigenfunctions at a finite temperature $|i\rangle$ in this basis as

$$|i\rangle = \sum_{k} c_{k}^{i} |k\rangle , \qquad (7)$$

with $\Sigma_k (c_k^i)^2 = 1$.

In order to measure the level mixing and delocalization of eigenfunctions at a finite temperature, we define the following probability:

$$P(c^2) = \frac{\Delta N(c^2)}{N_t \delta},\tag{8}$$

where c^2 is the probability for the eigenfunctions at zero temperature to appear. $\Delta N(c^2)$ is the number having the probability between c^2 and $c^2 + \delta$ with δ being a small interval of c^2 . N_t is the total number of the eigenstates involved (at present $N_t = 200$). For the J/ψ state (the ground state) at a finite temperature, the $P(c^2)$ versus c^2 for a few values of μ are shown in Fig. 2. It is evident that, when μ is small, for example 0.10 GeV, the components with nonzero probabilities are limited within the lowest states. This means that the components with small probabilities prevail and the distribution width of $P(c^2)$ is rather small. In other words, the level mixing is localized. With the increase of μ , the number of the components with nonzero probabilities goes up. This leads to the increase of $P(c^2)$ in the region of large probabilities (see the case $\mu = 0.20$ GeV). As μ approaches to μ_c , the components with nonzero probability emerge on a large scale and become rather random (as shown by the cases μ =0.60, 0.90 GeV). Then, a wide distribution of $P(c^2)$ appears and the level mixing is strong and delocalized. Referring to Fig. 1, one can find that the quantum chaotic motion occurs when a strong level mixing emerges. It can also be found that, for higher eigenstates at a finite temperature, the level mixing changes with μ in a similar way as J/ψ state. Furthermore the strong level mixing appears for these eigenstates even if μ takes a smaller value.

The feature of the effect of the level mixing induced by the variance of μ is that the probability for the J/ψ state at zero temperature to appear in the J/ψ state at a finite temperature descends with the increasing of μ . Present result shows that at $\mu = 0.10, 0.20$, the probability is about 90, 70%, respectively. As μ increases, the probability decreases distinctly. For instance, when $\mu = 0.60$ GeV, the probability goes down to 10%. Referring to Eq. (7) one can know that the neighboring levels (e.g., ψ', ψ'', \dots) at a finite temperature will also contribute to the J/ψ at zero temperature. However, their production probability is much smaller than that of J/ψ at a finite temperature (e.g., in p-p collisions, the ratio of ψ' to J/ψ is about 1/7 [21], and in Pb-Pb collision it is predicted to be 4% [22]). Then their contribution to the J/ψ at zero temperature should be very small. One can hence easily recognize that the production of J/ψ at zero temperature, which means in fact the probability for the charmonium states at a finite temperature to include the J/ψ state at zero temperature, will be suppressed with the growth of μ and a strong suppression can be expected when μ is large enough. On the other hand, It has been well known that, in the deconfinement scheme, the suppression due to the dissociation of J/ψ occurs only if μ reaches or exceeds the critical value μ_c [1]. Present analysis indicates that, the suppression due to the level mixing can take place before μ reaches the critical value for the dissociation to happen. In addition, we find that for *p*-wave energy spectrum of the $c\bar{c}$ system, the transition to the chaos and the suppression due to the level mixing occur too. Moreover, the similar behavior can be found in the $b\overline{b}$ system.

To apply the above discussion to an actual physical situation, one needs to know the specific dependence of the Deby screening mass μ on the temperature T. Even though several studies on the properties of QCD at finite temperatures and densities have been accomplished (see for example Refs. [23-27]) and the general feature of the temperature dependence $\mu(T)$ has been well established, the explicit expression $\mu(T)$ has not yet been determined uniquely because a temperature-dependent running coupling constant is involved. Then the density dependence of μ is not clear now. Any way, it is certain that the $\mu(T)$ increases with T increasing. The above calculated result manifests that, as the statistical properties of the mass spectrum of the $c\bar{c}$ system is taken into account, the critical value μ_c for the strong suppression to take place gets lowered obviously from that in the deconfinement scheme (from 0.9 to 0.6 GeV). Then the corresponding critical temperature T_c decreases too. Consequently, much stronger suppression than that predicted in the deconfinement scheme can emerge in the same experimental status. The level mixing, or the chaotic motion, due to the

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color screening is then an alternative mechanism to induce the strong J/ψ suppression.

In conclusion, in this paper we have investigated the level statistics of the charmonium energy spectrum in the framework of the BS equation and found that, with the increasing of the color screening mass, the transition from regular to chaotic motion can be induced for the $c\bar{c}$ system. The level mixing due to the color screening is then an alternative mechanism resulting in the strong J/ψ suppression, and furthermore this kind of suppression can occur before μ reaches its critical value for the dissociation of $c\bar{c}$. This implies that a strong J/ψ suppression is possible in the absence of dissociation of the J/ψ .

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