

Transport across two interacting quantum dots: Bulk Kondo, Kondo box, and molecular regimesL. C. Ribeiro,^{1,2} I. J. Hamad,^{3,4} G. Chiappe,² and E. V. Anda³¹*Centro Federal de Educação Tecnológica Celso Suckow da Fonseca (CEFET-RJ/UnED-NI), RJ 26041-271, Brazil*²*Departamento de Física Aplicada, Unidad Asociada del Consejo Superior de Investigaciones Científicas and Instituto Universitario de Materiales, Universidad de Alicante, San Vicente del Raspeig, Alicante 03690, Spain*³*Departamento de Física, Pontificia Universidade Católica do Rio de Janeiro (PUC-Rio),**22452-970 Caixa Postal: 38071 Rio de Janeiro, Brazil*⁴*Instituto de Física Rosario, Universidad Nacional de Rosario, Bv. 27 de Febrero 210 bis, Rosario (2000), Argentina*

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We analyze the transport properties of a double quantum dot device with both dots coupled to perfect conducting leads and to a finite chain of N noninteracting sites connecting both of them. The interdot chain strongly influences the transport across the system and the local density of states of the dots. We study the case of a small number of sites, so that Kondo box effects are present, varying the coupling between the dots and the chain. For odd N and small coupling between the interdot chain and the dots, a state with two coexisting Kondo regimes develops: the bulk Kondo due to the quantum dots connected to leads and the one produced by the screening of the quantum dot spins by the spin in the finite chain at the Fermi level. As the coupling to the interdot chain increases, there is a crossover to a molecular Kondo effect, due to the screening of the molecule (formed by the finite chain and the quantum dots) spin by the leads. For even N the two Kondo temperatures regime does not develop and the physics is dominated by the usual competition between Kondo and antiferromagnetism between the quantum dots. We finally study how the transport properties are affected as N is increased. For the study we used exact multiconfigurational Lanczos calculations and finite- U slave-boson mean-field theory at $T = 0$. The results obtained with both methods describe qualitatively and also quantitatively the same physics.

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I. INTRODUCTION

Since the prediction of the occurrence of the Kondo effect [1] in a single quantum dot (QD) device [2,3] and its subsequent experimental observation [4], several single, double (DQD), and multiple quantum dots devices, or systems with atoms or molecules acting as magnetic impurities, have been studied both theoretically [5,6,8–11] and experimentally [12–16]. The interest in these systems stems from their potential applications to quantum and classical computing [17,18] and their usefulness as model systems to study the physics of strongly correlated electrons.

In a QD connected to leads the charge can be manipulated by means of an external gate potential. When the charge on the QD is close to an odd integer, the Kondo effect takes place and it results in transmission through the system at temperatures below the Kondo temperature T_K and for gate voltages that do not incorporate extra charge into the QD. The electrons in the QD and leads form a spin singlet, which is one of the most clear benchmarks of the Kondo effect [19]. The system constitutes an experimental realization of the single-impurity Anderson model [20].

Likewise, two QDs directly coupled between them amount to an experimental realization of the two-impurity Anderson model [21]. Here the physics is much richer, particularly in the regime where each dot contains an unpaired electron. In this case, the state of the system is characterized by the competition between the tendency of the conduction electrons on the leads to screen the spins localized in the QDs and the antiferromagnetic (AF) coupling between them [5,6] (also mediated by the reservoirs when the Coulomb repulsion U is very large [7]). The former favors the formation of two Kondo singlets, while the latter corresponds to a molecular state constituted by the two QDs. The resulting ground state

and transport properties of the system depend sensitively on the relative strength of the interactions and the topology of the system. In each scenario, phase transitions are predicted to occur between different quantum states. These states can be characterized by a single Kondo resonance, or by a Kondo peak with a very narrow dip at the Fermi level which is representative of a two-stage Kondo regime [22,23], or even by a situation in which there is a split resonance resultant of the dominance of the AF correlations between QDs. In this latter case the conductance changes from a large value to zero, in a wide region of the applied gate voltage. Moreover, it has been predicted that this system could suffer a quantum phase transition, which involves a non-Fermi-liquid fixed point [24], which in fact has proven extremely difficult to be observed [6,16,25]. These transitions in the two-impurity Anderson model have received wide attention in the theoretical and experimental literature in recent years [5,6,14]. An understanding of the physics of two interacting impurities is important and believed to play a role in the electronic properties of a wide range of strongly correlated materials, including spin glasses and heavy-fermion compounds [19].

It is known that when an impurity is coupled to a finite number of N noninteracting sites, the Kondo effect modifies its character in what is called the “Kondo box” effect, a phenomenon that has also been theoretically [26–30] and experimentally [31] studied. In this case, the finite system’s density of states (DOS) consists of a series of peaks separated by an energy Δ inversely proportional to v/N where v is the hopping matrix element among the sites and N is the number of them. The peaks could have a finite width δ if the quantum box is weakly connected to an infinite system. When N is very large these peaks overlap and the DOS resembles that of a continuum. As N is decreased or v augmented,

finite-size effects start taking place when $\Delta \approx T_K$ [26], if $\delta \ll \Delta$. Depending on whether one of these peaks coincides with the Fermi level or not (“at resonance” or “off resonance” case), there is a single or a split resonance in the impurity’s local DOS (LDOS) [26], and the Kondo temperature behaves in a different way [27].

The physics of a Kondo box described above can be realized in systems of two impurities coupled between them by a finite number of noninteracting sites. This problem is thus very interesting and has also been theoretically studied in various configurations [32–34]. Here, finite-size effects can take place together with a magnetic interaction between the impurities of the Ruderman-Kittel-Kasuya-Yosida (RKKY) type [35]. In this case, when the Kondo effect is present, the fourth-order RKKY interaction between the impurities is mediated by the electrons of the noninteracting sites, which are participating simultaneously in the Kondo screening of each impurity, as was studied in Ref. [32]. It has also been pointed out that other types of magnetic interactions arise such as the “Kondo correlated” or superexchange interactions studied in Ref. [36]. Also, using variational wave functions, it was predicted [37] that the interaction between the impurities is mainly due to an interference-enhanced hybridization that generates Kondo doublet states. This interaction can be more important than the RKKY type and of ferromagnetic character.

Experimentally, these kind of configurations have also been studied, as was done recently [15] with a system of two Co impurities interacting through N Cu atoms placed between them, constituting a linear CoCu_NCo cluster deposited on a surface of $\text{Cu}(1,1,1)$. It was observed that the number N of Cu in the linear chain strongly influences the Kondo temperature T_K of each impurity, in general lowering it and having an oscillatory behavior for $N > 3$ compatible with RKKY interactions [15].

In this paper, we analyze a DQD (or double impurity) system simultaneously connected to metallic leads and between themselves through a finite number of noninteracting sites (i.e., with no Coulomb repulsion). The study emphasizes the fact that, as mentioned above, the impurities in such a system are coupled to, and interact through, a noninteracting linear chain (NILC) that constitutes a quantum box, whose electrons can participate in the Kondo screening. This implies that the physics corresponding to this problem is one in which there is an interplay between a bulk continuous Kondo regime and a two-impurity Kondo box. We analyze the transport properties for different values of the number N of sites of the NILC and for different couplings of the QDs with it. We study the system at $T = 0$, but the results are valid also for temperatures well below the characteristic single-impurity Kondo temperature. In particular, we present a detailed analysis of NILC with $N = 1$ and $N = 2$, which are limiting cases of a more general situation in which discrete systems present resonances at the Fermi level or not, respectively. We ask ourselves what is the effect of varying the connection of the impurities to the finite system. In doing so, we are able to characterize a crossover, which occurs for $N = 1$ (odd N in general), from a two Kondo temperature regime to a molecular Kondo regime, in which three resonances rise up as the coupling between the NILC and QDs is increased. This is reflected in the LDOS and in the transmission of the system. For even N , a crossover

from a single central peak to a split peak is observed in the LDOS, in the transmission, and in the conductance. In this case, the system behaves in a similar way as one with a direct connection between the impurities [9–11]. Finally, we also present results for larger N , but small enough so as to guarantee that finite-size effects are still predominant. We conclude that the characteristics of the LDOS and the conductance (or transmission) are largely influenced by the spectrum of the quantum box formed by the two dots plus the NILC.

The paper is organized as follows. In Sec. II we present the model and the Hamiltonian we used to study it. In Sec. III we briefly describe the methods used to study the system, while in Sec. IV we present the results. Finally, we make the concluding remarks in Sec. V.

II. MODEL AND HAMILTONIAN

The two-QD system is sketched in Fig. 1. It is described by an Anderson Hamiltonian composed of three terms, $H = H_0 + H_t + H_{lcc}$. The first contribution carries the local physical information of the QDs and is given by

$$H_0 = \sum_{\substack{i=\alpha;\beta \\ \sigma}} \epsilon_i f_{i\sigma}^\dagger f_{i\sigma} + \sum_{i=\alpha;\beta} U f_{i\downarrow}^\dagger f_{i\downarrow} f_{i\uparrow}^\dagger f_{i\uparrow}, \quad (1)$$

where ϵ_i , U , and $f_{i\sigma}^\dagger$ ($f_{i\sigma}$) represent, respectively, the local energy state, the electron-electron Coulomb interaction, and the operator that creates (annihilates) an electron with spin σ in the i th QD. We assume $\epsilon_i = 0$. The local energy of the dots is tuned by a gate potential V_g that, for simplicity, is considered to be the same for both QDs. The second term can be written as

$$H_t = t' \sum_{\substack{i=L;R \\ j=\alpha;\beta \\ \sigma}} (c_{i\sigma}^\dagger f_{j\sigma} + \text{H.c.}) + t'' \sum_{\substack{i=1;N \\ j=\alpha;\beta \\ \sigma}} (f_{j\sigma}^\dagger c_{i\sigma} + \text{H.c.}) \quad (2)$$

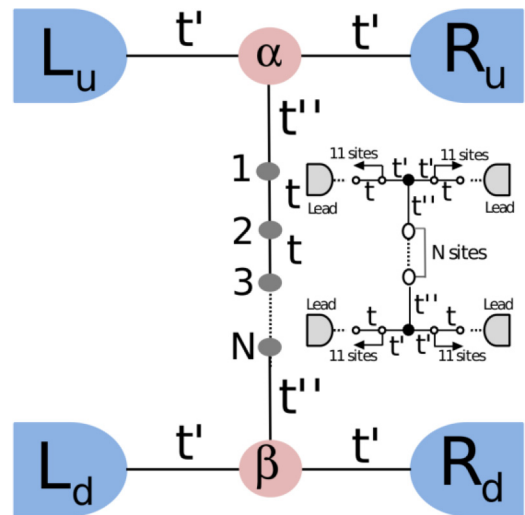


FIG. 1. (Color online) The structure studied in this work. The sites named α and β represent the two dots, with finite intrasite Coulomb repulsion U . The sites in between the dots have no Coulomb repulsion. Inset: The scheme used for the multiconfigurational Lanczos calculations (see Sec. III B).

and describes the connections of the QDs to both the metallic leads and the NILC through the hopping terms t' and t'' , respectively. As the physics depends on the relative values between these two parameters plus the Coulomb repulsion U , we fix t' and U and concentrate on the effects of varying t'' . The third term,

$$H_{lcc} = t \sum_{\substack{i \in L_u, R_u, L_d, R_d \\ \sigma}}^{\infty} (c_{i\sigma}^\dagger c_{i+1,\sigma} + \text{H.c.}) + t \sum_{\substack{i=1 \\ \sigma}}^N (c_{i\sigma}^\dagger c_{i+1,\sigma} + \text{H.c.}), \quad (3)$$

describes the leads and the NILC. The Fermi level in the reservoirs is adjusted to zero, $E_f = 0$, while the hopping t is adopted as the energy unit.

III. METHODS

A. Finite- U slave-boson mean-field approximation

The main idea concerning the slave-boson approach [38,39] consists of enlarging the Hilbert space by introducing in the Hamiltonian a set of bosonic operators which incorporate into the system the physics underlying the Kondo regime. These operators, e_i , $p_{i\sigma}$, and d_i , with i corresponding to the i th impurity, are responsible for projecting the system on a state of zero, single, and double occupation on the impurity and are introduced by the hybridization of the fermion operator which creates (annihilates) an electron with spin σ in the i th impurity,

$$f_{i\sigma}^\dagger \rightarrow Z_{i\sigma}^\dagger f_{i\sigma}^\dagger, \quad (4)$$

where

$$Z_{i\sigma}^\dagger = [1 - d_i^\dagger d_i - p_{i\sigma}^\dagger p_{i\sigma}]^{-\frac{1}{2}} (e_i^\dagger p_{i\sigma} + p_{i\bar{\sigma}}^\dagger d_i) \times [1 - e_i^\dagger e_i - p_{i\bar{\sigma}}^\dagger p_{i\bar{\sigma}}]^{-\frac{1}{2}} \quad (5)$$

is an operator that, in the mean-field approximation, becomes a parameter \bar{Z} that reproduces correctly the results expected in the noninteracting limit $U = 0$ [39] and is responsible for the renormalization of the connections of the impurities to the NILC and to the metallic leads. It is observed that this parameter becomes less than 1, $\bar{Z} < 1$, as the system enters into the Kondo regime.

The term ‘‘slave boson’’ comes from the constraints

$$e_i^\dagger e_i + \sum_{\sigma} p_{i\sigma}^\dagger p_{i\sigma} + d_i^\dagger d_i - 1 = 0 \quad (6)$$

and

$$f_{i\sigma}^\dagger f_{i\sigma} - p_{i\sigma}^\dagger p_{i\sigma} - d_i^\dagger d_i = 0 \quad (7)$$

that are imposed on the boson operators in order to eliminate the nonphysical states, assuring that the impurity is occupied with zero, one, or two electrons and establishing a correspondence between bosons and fermions. They are incorporated into the Hamiltonian through the Lagrange multipliers λ_1^i and $\lambda_{2\sigma}^i$.

Within the slave-boson mean-field approximation at finite U (finite- U SBMFA) we write the effective

Hamiltonian,

$$H_{\text{eff}} = \sum_{\substack{i=\alpha,\beta \\ \sigma}} \epsilon_i f_{i\sigma}^\dagger f_{i\sigma} + t'' \sum_{\substack{j=\alpha,\beta \\ i=1;N \\ \sigma}} \bar{Z}_j (f_{j\sigma}^\dagger c_{i\sigma} + \text{H.c.}) + \sum_{i=\alpha,\beta} U_i \langle d_i \rangle^2 + t' \sum_{\substack{i=L;R \\ j=\alpha,\beta \\ \sigma}} \bar{Z}_j (c_{i\sigma}^\dagger f_{j\sigma} + \text{H.c.}) + \sum_{i=\alpha,\beta} \lambda_1^i (\langle e_i \rangle^2 + \langle p_{i\sigma} \rangle^2 + \langle d_i \rangle^2 - 1) + \sum_{\substack{i=\alpha,\beta \\ \sigma}} \lambda_{2\sigma}^i (f_{i\sigma}^\dagger f_{i\sigma} - \langle p_{i\sigma} \rangle^2 - \langle d_i \rangle^2) + H_{lcc}, \quad (8)$$

where we observe that the local energy levels $\epsilon_i (i = \alpha; \beta)$ are renormalized by the Lagrange multiplier $\lambda_{2\sigma}^i (i = \alpha; \beta)$, $\epsilon_i + \lambda_{2\sigma}^i$, and the connections t' and t'' of the QDs by the multiplicative slave-boson parameter \bar{Z} . These are the two renormalizations that, in the context of the finite- U SBMFA, carry the system into the Kondo regime. At $T = 0$ the effective Hamiltonian H_{eff} is minimized with respect to the Lagrange multipliers and to the mean values of the bosons operators, resulting in a nonlinear set of ten equations that have to be solved self-consistently in order to obtain the numerical values of these quantities.

B. Multiconfigurational Lanczos

In the multiconfigurational Lanczos (MCL) calculation [40,41] the system is described by the same tight-binding-based Hamiltonian discussed before. The process requires the definition of a subsystem, a cluster of an arbitrary number of M sites that includes the DQD, the NILC, and a portion of the connecting leads. Due to numerical reasons, each part of the four leads is taken to be constituted by 11 noninteracting sites, as is shown in the inset of Fig. 1. As N is the number of sites of the NILC, the size of the cluster taken is given by $M = N + 2 + 44$. In each lead we do a transformation from left/right sites to symmetric/antisymmetric channels. Only symmetric channels are coupled with the central structure (DQD + NILC).

In order to numerically obtain the ground state of the cluster we use a mean-field solution as the starting point. The Hilbert space is increased by successive applications of H . In each step we obtain the ground state within the Hilbert subspace generated by the application of H . In each step, the ground state energy is lowered, and this process continues until the desired convergence is reached in the ground state energy. With the ground state so obtained, we proceed to calculate the Green's functions of the finite cluster, using a Lanczos procedure. Finally, the cluster is embedded into the rest of the system, within the LDECA and ECA formalism [23,42,43], using a Dyson equation to obtain the Green's function of the entire system. LDECA allows to obtain the value of the conductance exactly, but the behavior of transmission or density of states (DOS) as function of the frequency is poorly described. To get it we use ECA, but previously we verify the results at the Fermi level with LDECA.

IV. RESULTS

In this section we study the transport properties of the system. To this purpose we calculate, using both methods described above, the local density of states (LDOS) at the QDs, the transmission $T(\omega)$ as a function of ω , and the conductance $G = T(0)$ as a function of the gate potential V_g , or, equivalently, the energy levels of the impurities. Conductance or transmission, magnitudes that can be experimentally measured, are calculated from the upper left to the upper right lead. The transmission, using the Landauer-Buttiker formalism [44], is calculated as

$$T(\omega) = 4\pi^2 t^4 \text{Im}G_l(\omega) \text{Im}G_r(\omega) [|G_{\alpha\beta}(\omega)|^2 + |G_{\alpha\alpha}(\omega)|^2], \quad (9)$$

where $G_{l(r)}$ is the Green's function of the upper left (right) lead, and $G_{\alpha\beta}$ ($G_{\alpha\alpha}$) is the propagator from dot α to dot β (α). We also present finite- U SBMFA results for the parameters \bar{Z} and $\bar{\epsilon}$, that help to understand the physics involved. The results were obtained for the system in different regions of the parameter space and are divided into $N = 1$, $N = 2$, and large- N limit. The Coulomb interaction in the QDs and the connection with the leads are always $U = 0.5$ and $t' = 0.2$, respectively, in units of t .

A. $N = 1$

We take the $N = 1$ case as representative of a quantum box with states at the Fermi level. For this case, it is instructive to think of the two-dot system plus the central noninteracting site as a three-atom molecule, with three molecular energy levels. One level has zero energy and weight only in the QDs. The other two levels have energy $\pm\sqrt{2}t''\bar{Z}$ (where \bar{Z} is the finite- U SBMFA renormalization parameter that, when squared, represents the weight of the Kondo state), whose larger weight falls on the central site. These levels exist since each dot is in a Kondo state due to its coupling to the respective leads, and hence has a resonance at zero energy. Therefore, the charge of the molecule is three electrons. In Fig. 2 we present the LDOS calculated (a) in the QDs and (b) in the noninteracting central site, obtained with the gate potential V_g adjusted in the particle-hole symmetric position $V_g = -U/2$. We can clearly identify two quantum regimes in the system. The first, for $t'' \lesssim 0.04$, in which we observe a peak with a narrow dip just at the Fermi level, is characterized by the existence of two energy scales that correspond to two Kondo temperatures, T_{K1} and T_{K2} . These energy scales naturally emerge from the figure and are associated, respectively, with the widths of the dip and of the peak in the LDOS of the QDs. In this regime the particles, and hence the spins, are equally distributed in the three quasidegenerated molecular energy states that exist next to $E_f = 0$. Then, the first Kondo temperature can be thought of as coming from the partial screening of the QDs spin by the free spin allocated at the central noninteracting site. The remanent spin is then completely screened by the leads, characterizing the second Kondo temperature.

One of the key points to obtain this double Kondo temperature behavior is that the NILC (in this case composed of a single atom) shows, for low values of t'' , a resonance in

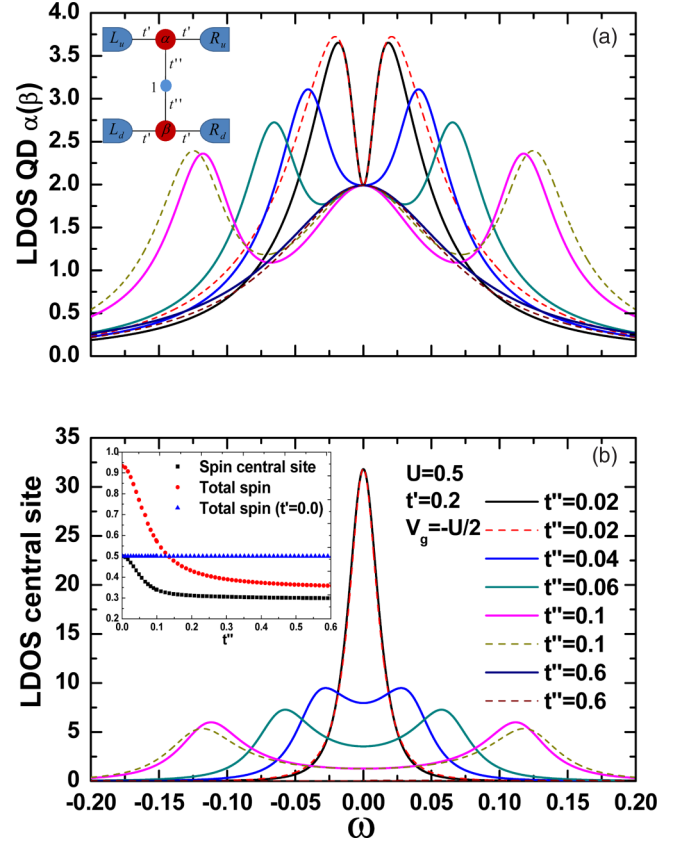


FIG. 2. (Color online) LDOS calculated (a) in the QDs and (b) in the noninteracting central site 1 as a function of the energy ω for the $N = 1$ case, and for different magnitudes of the connection t'' between the dots and the noninteracting central site at $V_g = -U/2$. The continuous lines corresponds to the finite- U SBMFA results, while the dashed lines correspond to MCL results. In the inset of panel (b) are shown the spin in the noninteracting central site (black squares), the total spin (red circles), and the total spin of the isolated molecule (blue triangles). See discussion in the text.

the LDOS just at the Fermi level. Then, there is a free spin allocated in this level that screens the spins at the QDs. The width of this resonance is similar to the width of the dip in the LDOS at the QDs. As already explained, when the molecule is disconnected from leads it has three levels: one at zero energy and weights only on the QDs and another two at energies $\pm\sqrt{(2)}t''\bar{Z}$ with most of its weight on the central site. When connected, the peaks are broadened by their hybridization with the leads. While t'' is small enough as compared with the connection between the dots and the leads (t'), these two levels superpose creating a single resonance at the central site. The analysis of the central site's Green's function shows that this occurs while $t'' \lesssim T_K / (2\sqrt{(2)}\bar{Z})$, where we have assumed that the Kondo peaks have a width of T_K . For the parameters used in Fig. 2, this occurs when $t'' \lesssim 0.02$, above which a crossover begins to a molecular Kondo regime, discussed below, and a structure with two peaks appears at the LDOS of the central site. It is interesting to observe that the resonance at the central site opens a tunneling channel, at the Fermi level, between the upper left and the lower leads. As a consequence, the transmission at the Fermi level between the upper left and

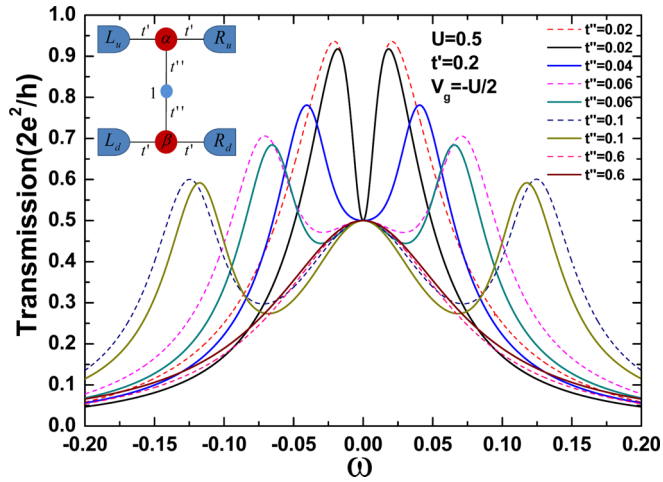


FIG. 3. (Color online) Transmission $T(\omega)$ as a function of ω for the system with $N = 1$, $V_g = -U/2$, and different magnitudes of the connection t'' with the central noninteracting site. The continuous lines correspond to the finite- U SBMFA results, while the dashed lines correspond to MCL results.

upper right leads, presented in Fig. 3, reaches the value e^2/h which corresponds to half of its maximum $2e^2/h$, obtained when the transmission is analyzed displaced by a small amount outside the Fermi level [45]. A similar result can be obtained by studying the behavior of the conductance shown in Fig. 4: conductance through the upper wire is exactly half of the value of that obtained when the dots are disconnected ($t'' = 0$). When the connection between the dots is established the electronic flow can be equally divided between the interdot channel and the upper left-right one, independently of the value of t'' [46].

It is also important to notice that in this regime, the magnetic moment of the noninteracting site assumes its maximum value and as a consequence is capable of screening the dot's spins [see inset of Fig. 2(b)]. To reinforce this image, note also that while the total spin for the free molecule is $S = 0.5$, it changes to $S = 1$ as it is connected to the leads, reducing its value as t'' increases. This $S = 1$ value for $t'' \approx 0$ is in fact a mean value between $S = 0.5$ and $S = 1.5$, the two possible values of the total spin of three independent spins. It shows that, as is natural for $t'' \approx 0$, the molecule is not formed as a whole, something that happens when t'' increases, but all the same connecting the molecule to the leads changes the internal spin correlations. This is a quite general fact also observed in other systems, as was shown for example in Ref. [47], where the spin of MnPc changes from 1.5 to 1 by depositing it on a Bi(110) surface.

To emphasize the points mentioned above we have verified with MCL that, in the two Kondo temperatures regime, when there is a peak with a narrow dip in the LDOS of the dots, a magnetic field applied at the central noninteracting site produces the disappearing of the dip and the single Kondo peak is recovered. It is clear that for the central site to be able to participate in the screening of the QDs spins, its electron has to be at the Fermi level and with a fluctuating spin. The external magnetic field freezes the spin by opening a spin-dependent Zeeman splitting at the central site, eliminating its screening capabilities. The recovering of the central Kondo peak as the

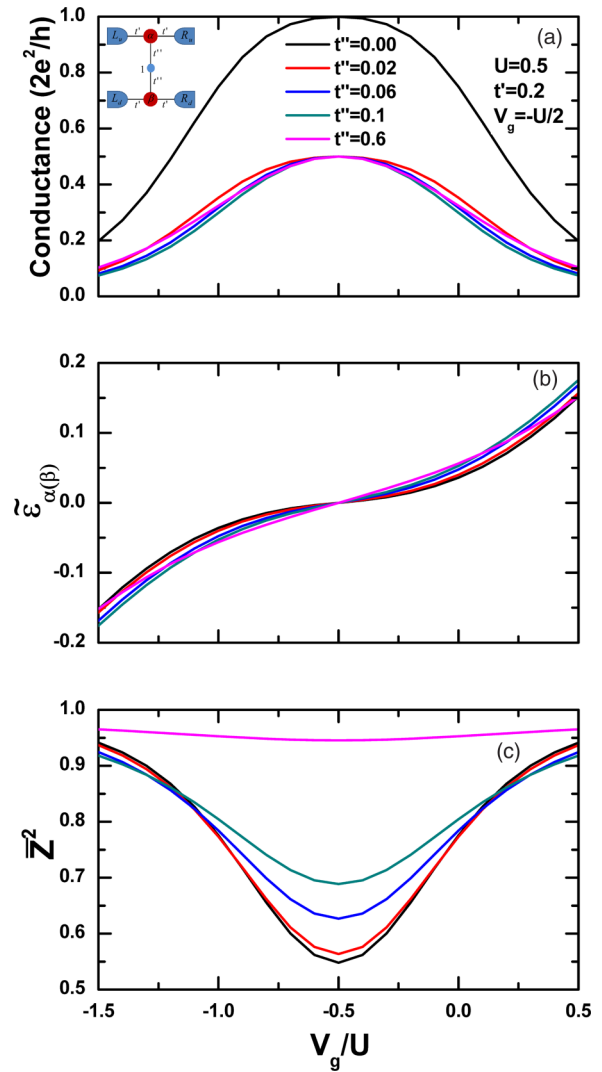


FIG. 4. (Color online) Conductance through the upper leads (a), renormalized energy level $\tilde{\epsilon}_{\alpha(\beta)}$ (b), and the parameter \tilde{Z}^2 (c), as a function of the gate potential V_g applied in the basis of the QDs for the system with different magnitudes of the connection t'' with the central noninteracting site.

central dip disappears is a confirmation of this process and of the role played by the interdot site.

In our effective one-body finite- U SBMFA Hamiltonian, after combining the upper and lower channels with a symmetric-antisymmetric transformation, a further symmetric-antisymmetric transformation with respect to the central site results in a symmetric pseudodot that is coupled to a reservoir *and* to the single noninteracting central site, plus another independent antisymmetric pseudodot that is only coupled to a reservoir. Let us concentrate on the symmetric pseudodot. Following Ref. [48], we can calculate the ground state of a simple system consisting of a site with energy ϵ_f and Coulomb repulsion U , which we take as infinite for simplicity, weakly coupled ($V \ll -\epsilon_f$) to a single site with energy $\epsilon_l = 0$, without U . From these we can calculate the Green's function of this simple system; i.e., $\langle GS | f \frac{1}{\omega + E_{GS} - H} f^\dagger | GS \rangle + \langle GS | f^\dagger \frac{1}{\omega - E_{GS} + H} f | GS \rangle$. The first term gives

$\langle GS|f \frac{1}{\omega+E_{GS}-H} f^\dagger|GS\rangle = \frac{2V^2}{\epsilon_f^2(\omega+2V^2/\epsilon_f)}$. This term has a pole at $\omega = -\frac{2V^2}{\epsilon_f}$, that since $V/\epsilon_f \ll 1$ is located very near zero energy, and has a weight of $2V^2/\epsilon_f^2 \ll 1$. This is basically the Kondo resonance. Due to the hybridization, the ground state energy is a singlet, which energy is lowered in comparison to the $V = 0$ case by an amount $2V^2/\epsilon_f$, which can be taken, divided by k_B , as the Kondo temperature of the system [48]. The other term gives poles at higher energies and is not important for this analysis. It can be easily checked, using equations of motion, that the LDOS of the symmetric pseudodot, which is coupled to both an infinite reservoir and a single separated site, consists of a peak with a dip at the Fermi level, due to interference effects between the Kondo peak due to the coupling to the reservoir and the one corresponding to the single isolated site. On the other hand, the LDOS of the antisymmetric one is just a Lorentzian peak. The LDOS at the QDs can be written as the sum of the LDOS corresponding to each pseudodot, which gives the results of Fig. 2, which, we emphasize, have a very good agreement with the MCL results. Note that our results differ from a study of a similar system in the so called “at resonance” situation [33]. The two Kondo temperature regime is manifested as a peak with a dip in the LDOS, rather than only one Kondo resonance that results from the sum of two peaks with different widths, each one reflecting its corresponding Kondo temperatures, as proposed in the mentioned study.

There is a second Kondo regime that appears when t'' is increased. In this new regime two lateral peaks appear in the LDOS of the noninteracting central site, together with an important reduction of the LDOS at E_f [see Fig. 2(b)]. As explained above, this is because the energy broadening of the molecular states with weight in it is now lower than their energy separation $\Delta E \sim t''$. These molecular orbitals of higher and lower energy are empty and double occupied, respectively. Remembering that the central site has weight mostly on these levels, we can conclude that with the noninteracting central site occupied with (zero) or two electrons no internal screening is now possible. The molecule has a net spin $S \approx 1/2$, as shown in Fig. 2(b), inset, equally distributed between the two dots, which become Kondo correlated with the spins in the leads. The net spin is not exactly $1/2$ since the molecule is connected to the leads through t' . It can be checked that, decreasing t' , the value of the total spin tends to $1/2$ as t'' increases. Although the molecule is connected to four leads, it represents all the same a traditional bulk one-channel SU(2) Kondo effect, as the electrons can flow freely from one channel to the other [49]. In Fig. 3 we observe a three-peak structure in the transmission for $t'' \gtrsim 0.04$, the central one being associated with a molecular Kondo effect. The lateral peaks correspond to tunneling resonances in which the molecule has a charge that fluctuates around 4 (left peak) or 2 (right peak) and the total spin is close to zero. This molecular Kondo regime represents a system that has a state at the Fermi level located mainly at the QDs and providing a channel for the electrons to go from one dot to the other. So, similarly to the previously analyzed two Kondo temperature regime, the conductance through the upper left and upper right leads is e^2/h , half of its maximum value $2e^2/h$ reached when $t'' = 0$, as shown in Fig. 4(a).

In Figs. 4(b) and 4(c) we present, respectively, finite- U SBMFA results showing the renormalized energy level $\tilde{\epsilon}_i$ of

the local state in the QDs and the renormalization parameter \tilde{Z}^2 as a function of the gate potential V_g applied on the QDs and for some of the values of the connections t'' used in Figs. 2 and 3. For all values of t'' we observe a plateau structure in $\tilde{\epsilon}_i$ which, in the context of the finite- U SBMFA, is the fingerprint of the Kondo effect. This plateau is an indication of the existence of a resonance, the Kondo peak, in the LDOS of the QDs, which remains fixed near the Fermi level as the potential V_g varies. This result, together with the renormalization observed in \tilde{Z}^2 , corroborates the Kondo nature of the two states described in this section. It is important to observe that the two Kondo temperature regime, obtained for $t'' \lesssim 0.04$, is associated with a larger plateau structure in $\tilde{\epsilon}_{\alpha(\beta)}$ and with a stronger renormalization (lower \tilde{Z}^2) of the connections if compared to the Kondo molecular regime.

The magnetic RKKY interaction between impurities was carefully studied in a system similar to ours in Ref. [32], taking into account the discrete character of the spectrum, with the conclusion that the equivalent to our $N = 1$ (odd N in general) case (“at resonance”) implies a ferromagnetic correlation. Keeping in mind what was pointed out in the Introduction, that besides the RKKY interaction there can be other types of magnetic interactions in systems like these, we have evaluated the spin-spin correlations through a MCL calculation. The results show that the correlation between the dots and the central site is AF, while between them a ferromagnetic correlation is established. Taking into account the results presented in the inset of Fig. 2(b), we can conclude that what determines the physics of our system is the fact that, as t'' increases, the two QDs plus the central site behave as a single entity, a molecule with $S \approx 1/2$ as we explained above, and this results in the molecular Kondo regime, which is reflected in the LDOS, transmission, and conductance of the system.

B. $N = 2$

The two-dot system defined with $N = 2$ corresponds to a four-atom molecule with two noninteracting intermediate sites. In this case the molecule has four levels, two well above and below the Fermi level, bonding and antibonding states created by the strong connection between the two sites forming a singlet. The other two levels are separated from the Fermi level by an amount proportional to t'' . In Fig. 5 we present the transmission $T(\omega)$ between the leads left (up) and right (up) for the system with $V_g = -U/2$ and, therefore, for a charge occupation of four electrons in the molecule. Results for the LDOS, as for the $N = 1$ case, are very similar to the transmission, and are not presented to avoid repetition. Again, the agreement between finite- U SBMFA and MCL is very good. The results do not present a dip in the LDOS of the QDs for small values of t'' and therefore it is not observed any signal of two coexisting Kondo temperatures, like in the $N = 1$ case. Instead, a single Kondo peak is observed. This is so because no internal screening is possible as the two free electrons in the center occupy the antibonding state described before. Each QD has a spin which is only screened by the leads, giving rise to a Kondo effect at each QD characterized by the resonance observed in the transmission $T(\omega)$ for $t'' \leq 0.3$. Consequently,

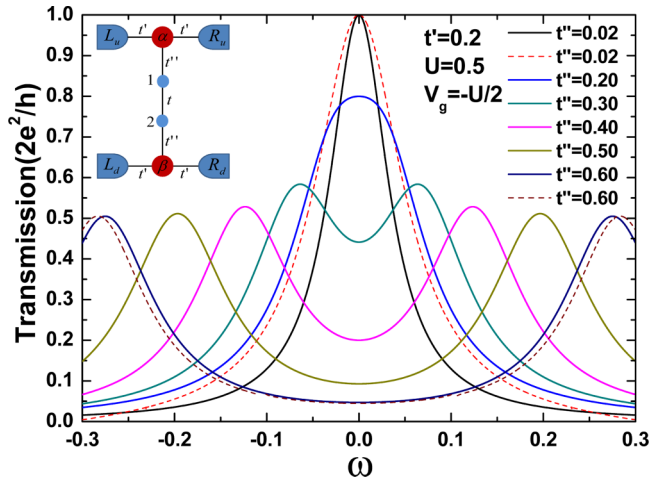


FIG. 5. (Color online) Transmission $T(\omega)$ as a function of ω for different magnitudes of the connection t'' for the $N = 2$ case with $V_g = -U/2$. The continuous lines correspond to the finite- U SBMFA results, while the dashed lines correspond to MCL results. As for the $N = 1$ case, the agreement between both methods is very good.

the conductance, presented in Fig. 6(a), reaches its maximum value $2e^2/h$.

Increasing t'' the system goes into a crossover regime characterized by the interplay between the Kondo ground state and the AF correlation between the spins of the QDs. These states are associated, respectively, with two energy scales: the Kondo temperature T_K and the value of the effective AF interaction I . In Fig. 5 we can identify the crossover region for $0.02 < t'' < 0.3$ and observe, for $t'' > 0.3$, the splitting in the transmission $T(\omega)$ (or in the LDOS, which has the same form as has been mentioned) due to the increasing of I . This splitting characterizes the establishment of the AF regime in the system [9,10]. We have calculated with MCL the spin-spin correlations (not shown) between each dot and the rest of the sites of the system, verifying that, as t'' increases, the (Kondo) AF correlation of each dot with the reservoirs to which it is directly connected decreases, while the also AF correlation between the two dots increases, giving support to the well established image of the Kondo-AF crossover mentioned. Hence, in this case, we can conclude that as t'' increases, the two QDs plus the two noninteracting central sites also form a molecule, but with net spin $S = 0$.

In Fig. 6 the conductance (a) and the slave-boson parameters $\tilde{\epsilon}_i$ (b) and \bar{Z}^2 (c) are presented as a function of V_g . We observe that with the transition from the Kondo to the AF regime the conductance goes to zero in $V_g = -U/2$. We also conclude from an inspection of Figs. 6(b) and 6(c) that $\bar{Z} \mapsto 1$ together with the destruction of the plateau in $\tilde{\epsilon}$, indicating, in the context of the finite- U SBMFA, that the system is driven out of the Kondo regime with the increasing of t'' . Considering the dependence of the conductance on V_g , we observe in (a) the formation of two lateral peaks. These peaks correspond to the molecular Kondo resonances [10,23] associated with a charge occupation of approximately five (left peak) or three (right peak) electrons and a total spin close to 0.5. In this regime, transmission across the upper channel has a maximum value of e^2/h , half of the quantum of conductance, because

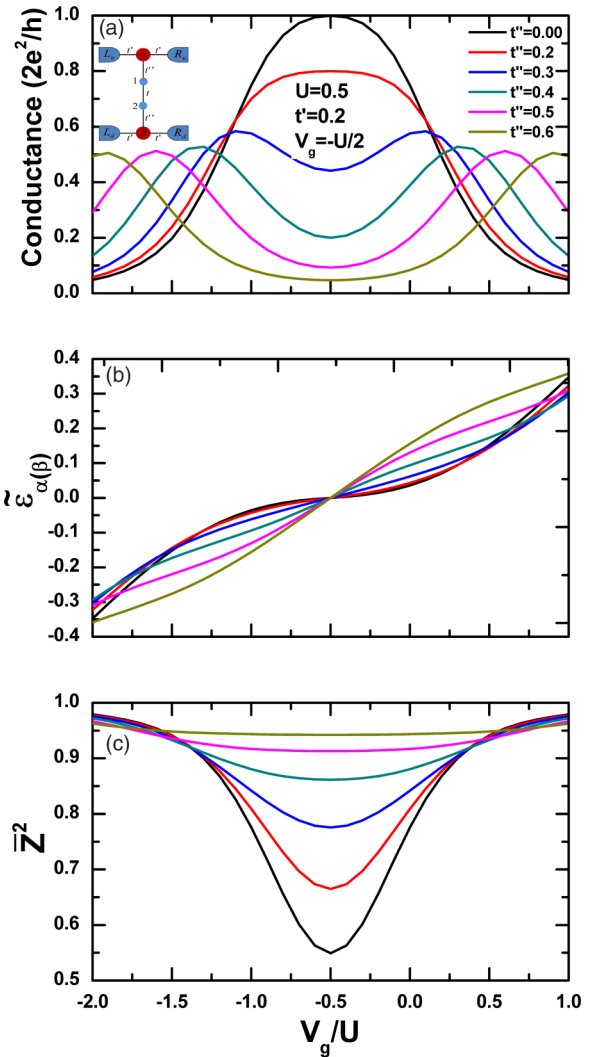


FIG. 6. (Color online) Conductance through the upper leads (a), renormalized energy level $\tilde{\epsilon}_{\alpha(\beta)}$ (b), and the parameter \bar{Z}^2 (c), as a function of the gate potential V_g applied in the basis of the QDs for the system with different magnitudes of the connection t'' with the central NILC.

there is an interdot flow of electrons through the molecular state that is half populated. The coherence existing between both dots in this Kondo molecular regime provides a channel for the conduction. Through the analysis of Fig. 6(b), within the context of the finite- U SBMFA, we observe that, as t'' increases, $\tilde{\epsilon}_i$ tends to form a double-plateau structure [50] in the regions of V_g corresponding to a charge occupation of approximately three or five electrons, corroborating the molecular Kondo regime, while the suppression of the plateau for four electrons certifies that the system does not have a Kondo ground state. We can conclude that this system behaves in a similar way to a two-dot structure with a direct connection between them [10,11].

C. Large N

In Fig. 7 we present the transmission $T(\omega)$ between the leads L_u and R_u for the system with the gate potential adjusted

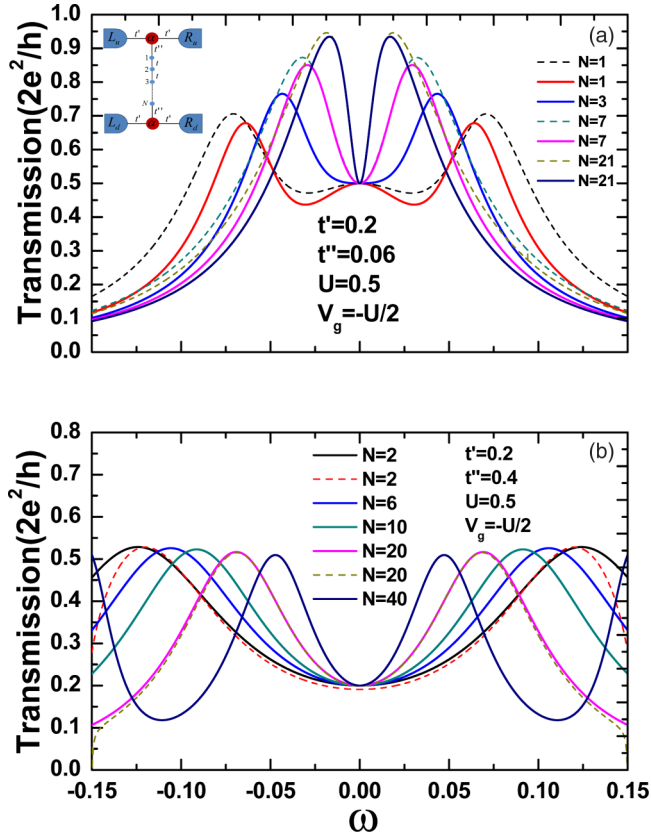


FIG. 7. (Color online) Transmission $T(\omega)$ as a function of ω for different lengths of the NILC for the system with $V_g = -U/2$. The upper panel (a) corresponds to odd N with $t'' = 0.06$ while the lower panel (b) to the even- N case with $t'' = 0.4$. The continuous lines correspond the finite- U SBMFA results while the dashed lines correspond to MCL results.

to $V_g = -U/2$, and different lengths of the NILC. We present results for odd N with $t'' = 0.06$ and for even N with $t'' = 0.4$, since with these values of the connections the effects of varying N are most clearly seen, as the energies involved are different for each case. The LDOS in the QDs (not shown) behaves in the same way as the transmission. The description for large odd or even N is qualitatively similar to the cases $N = 1$ and $N = 2$, respectively. This occurs since N is low enough so as to preserve finite-size effects, which happens if $\Delta \ll T_{K_0}$, T_{K_0} being the isolated-dot Kondo temperature. The behavior of the Kondo resonance alternates between even and odd N , as expected.

In Fig. 7(a), we clearly identify the existence of a regime of two coexisting Kondo temperatures for odd N with the lowering of the width of the dip (related to the first Kondo temperature T_{K_1}) as N is increased. Since the central noninteracting sites are linked by a hopping $t \gg t'' = 0.06$, the level structure consists of one level at E_f , two Kondo peaks at E_f , which splitting is controlled by t'' and N , and $(N-1)/2$ levels above and below E_f . As in the case of $N = 1$, T_{K_1} is due to the existence of a NILC state at the Fermi level, which allows us to localize two spins in the DQD and one free spin in the NILC. Hence, the two Kondo temperature regime develops in the same way as was described in the $N = 1$ case. As N is

increased, the weight of the state at the Fermi energy is spread along the odd sites of the NILC, reducing the weight of the local wave function, which is directly connected to the dots through t'' . This effect reduces the splitting of the two Kondo peaks and hence the width of the dip decreases. Increasing N produces, in some extent, a similar result to reducing the effective value of t'' .

For even N a similar phenomenon to the $N = 2$ case takes place as N is increased since there is no energy level at E_f . The two lateral peaks that exist for a chosen value of t'' are pushed to energies nearer to the Fermi level, as can be seen in Fig. 7(b). Eventually, as N increases, additional peaks appear in the transmission close to the Fermi level, as can be observed in the $N = 40$ case, where two additional lateral peaks are observed. A simple Green's function calculation allows us to verify the fact that the LDOS or the transmission $T(\omega)$, at $\omega = 0$, are independent of N , as long as N is low enough so as to keep finite-size effects predominant. This is also observed in the mentioned figure. Naturally, for a large number of intermediate sites N , the level separation in the NILC is comparable to the isolated-dot Kondo temperature and the system begins to behave as in the continuum limit [26], with a typical Kondo peak at LDOS of the dots.

V. CONCLUDING REMARKS

We have studied a system of two QDs connected with infinite leads and between them through a NILC. We have emphasized the quantum box character of this channel connecting the two dots. These kinds of systems are very interesting since they can be used as quantum gates and they are experimentally feasible. Nowadays it is possible to manipulate QDs and even single atoms or molecules that act as magnetic impurities, and to control precisely the intermediate number of noninteracting sites through which the interaction between them is mediated [15]. For the particular geometry assumed, we analyzed the dependence of the transport properties and the LDOS of each QD as its connection to the NILC t'' is varied. We studied in detail the $N = 1$ and $N = 2$ cases, representative of the odd and even N , respectively, and also presented results for larger N . Our calculations were done using both the finite- U SBMFA and MCL, which showed a remarkable qualitative and quantitative agreement, giving support to our results.

For the $N = 1$ case, we could identify two regimes. For small t'' , there is a double Kondo temperature regime, one being representative of the bulk SU(2) Kondo regime of each dot spin screened by the spins of the leads to which it is connected, and the other being a Kondo box regime in which the screening is done by the spin of the electron occupying the level at the Fermi energy of the interdot chain. This is reflected in the LDOS and transmission as a peak with a dip at the Fermi level. As t'' increases, there is a crossover to a molecular Kondo regime where the two QDs plus the central site act as a whole entity with spin $S = 1/2$, Kondo correlated with the conduction electron spins of the leads. The conductance across the upper leads reflects this behavior.

For the $N = 2$ case, increasing t'' there is a crossover from a single-impurity Kondo resonance to a split one that eventually disappears when the antiferromagnetic correlations between

the dots, mediated by the intermediate sites, is dominant. This crossover is reflected in the transmission through the upper leads.

For a NILC with larger N , the behavior obtained is similar to the $N = 1$ (for odd N) and $N = 2$ (for even N) cases. For odd number of sites in the NILC and low values of the connection t'' with the dots, a two Kondo temperature regime is developed. The width of the dip (related to one of the Kondo temperatures) is reduced as N increases. In the even- N case, and for values of t'' big enough so as to have a split peak in the LDOS or transmission, the splitting is reduced as N increases. In both cases the behavior is a consequence of the renormalization of the energy spectrum as N is increased.

In general it can be concluded that, for the type of system we have studied in this work, it is the structure of energy levels in the molecule composed by the two dots plus the

central site(s), together with the fact that the QDs are initially Kondo correlated with the leads to which they are connected, that determines the structure of the LDOS and the transport properties of the system. Varying the connection t'' of the dots with the NILC changes the energy levels and produces crossovers that are reflected in the LDOS and in the transport properties.

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