Dynamical charge susceptibility in nonequilibrium double quantum dots

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Double quantum dots are one of the promising two-state quantum systems for realizing qubits. In the quest of successfully manipulating and reading information in qubit systems, it is of prime interest to control the charge response of the system to a gate voltage, as filled in by the dynamical charge susceptibility. We theoretically study this quantity for a nonequilibrium double quantum dot by using the functional integral approach and derive its general analytical expression. One highlights the existence of two lines of maxima as a function of the dot level energies, each of them being split under the action of a bias voltage. In the low frequency limit, we derive the capacitance and the charge relaxation resistance of the equivalent quantum RC-circuit with a notable difference in the range of variation for $R$ depending on whether the system is connected in series or in parallel.

By incorporating an additional triplet state in order to describe the situation of a double quantum dot with spin, we obtain results for the resonator phase response which are in qualitative agreement with recent experimental observations in spin qubit systems. These results show the wealth of information brought by the knowledge of dynamical charge susceptibility in double quantum dots with potential applications for spin qubits.

I. INTRODUCTION

In double quantum dots (DQDs), the knowledge of the dynamical charge susceptibility (DCS), which measures the ability of a system to adapt its electronic charge to an ac gate-voltage, is of fundamental interest in the general context of circuit quantum electrodynamics with gated GaAs, silicon, and germanium semiconductor quantum dots. This field has become all the more important because of its expected implications in manipulation, control and readout of spin qubits\cite{1}. There certainly are some theoretical works on charge susceptibility in DQD but they are mainly restricted to the study at zero frequency\cite{2,3} and in the low frequency regime with the determination of mesoscopic admittance\cite{4} and quantum capacitance,\cite{5} or to calculations performed at the lowest order in dot-lead coupling amplitude.\cite{6} Electrical transport experiments in DQD systems are, however, restricted neither to the weak dot-lead coupling regime nor to the measurement at low frequencies. On the contrary, in the last ten years one has witnessed a considerable experimental effort\cite{7,8} with measurements performed by using either on-chip superconducting resonant detectors,\cite{9,10} or dispersive probed microwave spectroscopy via reflectometry techniques\cite{11,12}, all of them made in the high frequency regime. To accompany this growing experimental development, it becomes of primary importance to progress on the theoretical level in order to investigate circuit quantum electrodynamics at high frequency in nanoscale systems. Indeed, the interpretation of these experiments requires the precise knowledge of the DCS at any frequency and temperature range, in both equilibrium and out-of-equilibrium DQDs. This article is precisely devoted to this theoretical issue. It is organized as follows: in Sec. II we present the functional integral approach used to solve this problem and give the expression for the dynamical charge susceptibility, in Sec. III we give the results for both serial and parallel DQDs, in Sec. IV we focus on the characterization of the equivalent quantum RC-circuit to the DQD system. Finally, in Sec. V, we study the reflection phase of the system considered as a resonator embedded in an electromagnetic environment, shedding new light on recent measurements performed in microwave reflectometry experiments in spin qubit systems. We conclude in Sec. VI.

II. MODEL

A. Functional integral approach

Let us consider a DQD connected to two leads described by the Hamiltonian $\hat{H} = \hat{H}_{\text{leads}} + \hat{H}_{\text{dots}} + \hat{H}_{\text{hop}}$, with $\hat{H}_{\text{leads}} = \sum_{\alpha=L,R,k} \varepsilon_{\alpha k} \hat{c}_{\alpha k} \hat{c}_{\alpha k}^\dagger$, $\hat{H}_{\text{dots}} = \sum_{i=1,2} \varepsilon_{i} \hat{d}_{i} \hat{d}_{i}^\dagger + \sum_{i=1,2} \sum_{\alpha=L,R,k} \gamma_{\alpha k} \hat{c}_{\alpha k} \hat{d}_{i} \hat{d}_{i}^\dagger$, and $\hat{H}_{\text{hop}} = \sum_{n=L,R,k} \sum_{i=1,2} \sum_{\alpha=L,R,k} \gamma_{\alpha k} \hat{d}_{i} \hat{c}_{\alpha k} + H.c.$, where $\varepsilon_{i}$ is the energy level of the dot $i$, $\gamma_{12}$ is the interdot coupling, $\varepsilon_{k}$ is the energy of one electron with momentum $k$ in the lead $\alpha$, and $V_{i,\alpha k}$ is the hopping energy between the dot $i$ and the lead $\alpha$ for momentum $k$. The very general form considered in the expression for $\hat{H}$ allows one to describe the situations where the two dots are connected in series as well as in parallel (see Fig. 1). We use the functional integral approach to derive the expression for the DCS. The partition function of the system writes

$$Z = \int \prod_{i=1,2} \prod_{\alpha=L,R} \prod_{k=\alpha} d\hat{d}_{i} d\hat{d}_{i}^\dagger \prod_{\alpha=L,R} d\hat{c}_{\alpha k} d\hat{c}_{\alpha k}^\dagger \frac{Z_{0}}{Z} d\mathcal{L}(\tau),$$

where $\mathcal{L}(\tau) = \sum_{i=1,2} \partial_{\tau} \hat{d}_{i} \hat{d}_{i}^\dagger \partial_{\tau} \hat{c}_{\alpha k} \hat{c}_{\alpha k}^\dagger - \hat{H}$ is the Lagrangian, $d_{i}^{(1)}$ and $c_{\alpha k}^{(1)}$ are the Grassmann variables associated with the operators $\hat{d}_{i}^{(1)}$ and $\hat{c}_{\alpha k}^{(1)}$. 

\cite[university]{}
In the linear response theory it is related to a correlation function through a Kubo-type formula $\mathcal{X}_{ij}(t,t') = i\Theta(t-t')\langle [\Delta \hat{N}_i(t), \Delta \hat{N}_j(t')] \rangle$ with $\Delta \hat{N}_i = \hat{d}_i^{\dagger} \hat{d}_i$. Decomposing the correlation function in the eigenstate basis and taking the Fourier transform, for the DCS one gets: $\mathcal{X}_c(\omega) = \sum_{\alpha,\beta=\pm} C_{\alpha,\beta} \mathcal{X}_{\alpha,\beta}(\omega)$, where $C_{\alpha,\beta}$ are coherency factors defined in Ref [23]. $\mathcal{X}_{\alpha,\beta}(\omega)$ can be expressed as a function of the Green functions in the dots following

$$\mathcal{X}_{\alpha,\beta}(\omega) = i \int_{-\infty}^{\infty} \frac{d\epsilon}{2\pi} \left[ \mathcal{G}_{\alpha,\beta}(\epsilon) \mathcal{G}_{\alpha,\beta}(\epsilon - i\omega) + \mathcal{G}_{\alpha,\beta}^{\ast}(\epsilon + i\omega) \mathcal{G}_{\alpha,\beta}^{\ast}(\epsilon) \right], \quad (5)$$

where $\mathcal{G}_{\alpha,\beta}^{\ast}$ are retarded, advanced, and nonequilibrium Green functions in the eigenstate basis. $\mathcal{G}_{\alpha,\beta}^{\ast}$ are diagonal matrices of elements $\mathcal{G}_{\alpha,\beta}^{\ast}(\epsilon) = (\epsilon - \lambda_{\alpha})^{-1}$ and $\mathcal{G}_{\alpha,\beta}^{\ast}(\epsilon) = (\epsilon - \lambda_{\beta})^{-1}$. In the steady state, one has $\mathcal{G}_{\alpha,\beta}^{\ast}(\epsilon) = \mathcal{G}_{\alpha,\beta}^{\ast}(\epsilon)(1 - \lambda_{\alpha})^{-1}$, where $\lambda_{\alpha}$ is the transition matrix from the initial state basis to the eigenstate basis. $\mathcal{G}_{\alpha,\beta}^{\ast}(\epsilon) = i \sum_{\alpha=L,R} \lambda_{\alpha,\beta} F_{\alpha}(\epsilon,\mu_{\alpha})$, where $\lambda_{\alpha}$ is the nonequilibrium self-energy, and $F_{\alpha}(\epsilon,\mu_{\alpha}) = (1 + \exp((\epsilon - \mu_{\alpha})/k_B T))^{-1}$. The Fermi-Dirac distribution in the lead $\alpha$ of chemical potential $\mu_{\alpha}$ and temperature $T$. We have thus established a general expression for the DCS of a nonequilibrium DQD, valid whatever its geometry is.

### III. RESULTS AND DISCUSSION

#### A. DQD in series

The results obtained for a DQD symmetrically connected in series are shown in Fig. 2. The variation of the absolute value of the DCS, $|\mathcal{X}_{c}(\omega)|$, is plotted in the form of color-scale plots as a function of $\epsilon_1$ and $\epsilon_2$. Figures 2(a) and (b) show the presence of peaks in $|\mathcal{X}_{c}(\omega)|$ along four branches denoted as $B_L^c$, $B_R^c$, $B_L^R$ and $B_L^R$ corresponding to the alignment of the bonding and anti-bonding state energies with the chemical potential in the leads, occurring when $\text{Re}\{\lambda_{\pm}\} = \mu_{L,R}$.

According to Eq. (5), one has

$$\lambda_{\pm} = \frac{1}{2} \left( \epsilon_1 + \epsilon_2 - i\Gamma \pm \sqrt{(\epsilon_1 - \epsilon_2)^2 + 4\epsilon V_{12}^2} \right),$$

making use of $\Gamma_{L,R,1} = \Gamma, \Gamma_{L,R,2} = \Gamma_{L,R,1} = 0$, and $\Gamma_{L,2} = \Gamma_{R,2} = 0$, and assuming $\alpha_1 = \alpha_2 = 1$ and $V_{12} = V_{21}$. At zero voltage, $B_L^c$ and $B_R^c$ coincide (the same for $B_L^R$ and $B_R^R$). $B_L^L$ and $B_L^R$ are two branches of a hyperbola of equations $\epsilon_1\epsilon_2 = |V_{12}|^2$, separated by a minimal distance along the diagonal $D$ of equation $\epsilon_1 = \epsilon_2$, taking the value of $2|V_{12}|$. From Eq. (6), the imaginary part of $\lambda_{\pm}$ are both equal to $-\Gamma/2$ and independent of $\epsilon_1$ and $\epsilon_2$. As a result, the width of the charge susceptibility peak arcs is uniform along the branches $B_L^c$ and $B_R^c$, as observed in Fig. 2(b). At finite voltage, Fig. 2(b) shows the splitting of the peak arcs into two branches $B_L^c$ and $B_R^c$, respectively $B_L^c$ and $B_R^c$, with the reduction of
According to Eq. (S37), one has

\[ B \] 

intensity along both halves of the branches 

\( B \) 

density at the intersections between the branches 

\( B \) 

with symmetrical couplings: \( \Gamma_{\alpha,ij} \).

\| \X_i(0) \|. At finite frequency, one observes a broadening of the peak arcs located around the \( B_{L,R}^{\perp} \) branches, together with the formation of an additional central peak at \( \epsilon_1 = \epsilon_2 = 0 \) (see Fig. 2(c) and (d)). An exact expression for \( \X_i(\omega) \) is derived from Eq. (S27) at \( T = 0 \) when \( \epsilon_1 = \epsilon_2 \) and \( \nu_{21} = \nu_{12} \) for a serial DQD. It writes

\[ \X_{\pm,\alpha}(\omega) = -\frac{\Gamma}{2h\omega(h\omega + i\Gamma)} \ln \left( 1 - \frac{h\omega(h\omega + i\Gamma)}{\Gamma} A_{\pm}(\mu_\alpha) \right), \]  

(7)

where \( A_{\pm}(\epsilon) = -\text{Im}\{\mathcal{G}_\epsilon^+(\epsilon)\}/\pi \) are the spectral function contributions from the anti-bonding and bonding states.

### B. DQD in parallel

Figure 3 shows the results obtained for a parallel DQD with symmetrical couplings: \( \Gamma_{\alpha,ij} \equiv \Gamma, \forall \alpha, i, j \), with \( \alpha_1 = \alpha_2 = 1 \) and \( \nu_{21} = \nu_{12} \). At zero frequency, one observes three differences towards the situation in series: (i) the intensity of \( |X_i(0)| \) is reduced on the branches \( B_{L,R}^{\perp} \); (ii) \( |X_i(0)| \) undergoes an extinction of its intensity at the intersections between the branches \( B_{L,R}^{\parallel} \) with the diagonal \( D \); (iii) at finite voltage, \( |X_i(0)| \) is of equal intensity along both halves of the branches \( B_{L,R}^{\perp} \) (respectively \( B_{L,R}^{\parallel} \)). These differences are understood as follows. According to Eq. (S37), one has

\[ \lambda_{\pm} = \frac{1}{2} \left( \epsilon_1 + \epsilon_2 - 2i\Gamma + \sqrt{(\epsilon_1 - \epsilon_2)^2 + 4(\nu_{12} - i\Gamma)^2} \right). \]  

(8)

The imaginary parts of \( \lambda_{\pm} \) depend on \( \epsilon_1 \) and \( \epsilon_2 \), contrary to what happens in the case in series where the imaginary parts of \( \lambda_{\pm} \) were equal to \(-\Gamma/2\) (see Eq. (6)). Typically the imaginary part of \( \lambda_+ \) for the parallel DQD is large along the branches \( B_{L,R}^{\perp} \) explaining the fact that the intensity of \( |X_i(0)| \) on the latter branches are reduced (property (i)). Along the diagonal \( D \) of equation \( \epsilon_1 = \epsilon_2 = \epsilon_0 \), one has \( \lambda_+ = \epsilon_0 - \nu_{12} \) whereas \( \lambda_- = \epsilon_0 + \nu_{12} - 2i\Gamma \). The imaginary part of \( \lambda_- \) is zero, meaning that the bonding state of the parallel DQD is disconnected from the leads, eliminating any dissipation effect through contacts to leads, and causing a significant reduction in the intensity of \( |X_i(0)| \) at the intersection of branches \( B_{L,R}^{\perp} \) and diagonal \( D \) (property (ii)). Finally, the property (iii) can be understood as follows: at finite voltage \( |X_i(0)| \) is maximal on the branches \( B_{L,R}^{\parallel} \) with equal intensity along both halves of the branches since each dot are equally connected to \( L \) and \( R \) leads for parallel DQD. At finite frequency, Figs. 3(c) and (d) show the broadening of the branches \( B_{L,R}^{\perp} \) and the widening of the gap in the branches \( B_{L,R}^{\parallel} \). At \( T = 0 \), \( \epsilon_1 = \epsilon_2 \) and \( \nu_{21} = \nu_{12} \), an exact expression for \( \X_{\pm,\alpha}(\omega) \) is derived from Eq. (S27) in the parallel geometry. It reads as

\[ \X_{\pm,\alpha}(\omega) = -\frac{2\Gamma}{h\omega(h\omega + 4i\Gamma)} \ln \left( 1 - \frac{h\omega(h\omega + 4i\Gamma)}{4\Gamma} A_{\pm}(\mu_\alpha) \right). \]  

(9)

### IV. EQUIVALENT QUANTUM RC-CIRCUIT

We now focus on the characterization of the equivalent quantum RC-circuit to the DQD whose capacitance and charge relaxation resistance are respectively given by

\[ C = e^2 |\X_i(0)| \]  

\[ R = e^2 \lim_{\omega \to 0} \text{Im} \{\X_i(\omega)\}/(\omega C^2) \]  

(22)

(23)
By employing similar arguments to those used for a serial DQD, and by noticing that \( n = 2 \) in the case in parallel since the bonding state is disconnected from the leads at \( \epsilon_1 = \epsilon_2 \), one shows that \( R \) varies from \( R_q/2 \) to \( R_q \), in agreement with what is observed in Fig. 4(d). It means that when \( \epsilon_1 = \epsilon_2 \), the number of quantum channels \( n \) allowing the charge to travel are equal to four in the case in series whereas it equals two in the case in parallel. The results given by Eqs. (10) and (11) can be viewed as the extension to a DQD of the results previously established in the cases of a single quantum dot or a quantum point contact. As in such systems, \( \chi_e(\omega) \) in a DQD obeys a generalized Korringa-Shiba relation according to which \( \lim_{\omega \to 0} \text{Im} \{ \chi_e(\omega) \}/\omega = \frac{1}{2} \sum x_{\alpha=L,R} \chi^2_{\alpha}(0) \), a characteristic of a Fermi liquid. It is worthwhile to explore the variation of \( \chi_e(\omega) \) at higher frequencies and to see how its frequency dependence deviates from this relation. Figure 5 reports the results obtained for \( \text{Im} \{ \chi_e(\omega) \}/\chi^2_{\omega}(0) \) as a function of \( \omega \). For a serial DQD, respectively parallel DQD, one observes a linear variation with frequency according to \( \omega/4 \), respectively \( \omega/2 \), at low frequencies in agreement with the generalized Korringa-Shiba relation, confirming the difference of a factor two found for the values of \( R \) between the cases in series and in parallel, whereas it shows strong deviations at higher frequencies.

V. REFLECTION PHASE

We end up by discussing the reflection phase of the system considered as a resonator embedded in an electromagnetic environment, which is defined as the phase shift between incoming and reflected microwaves, as measured in reflectometry experiment. A rapid calculation shows that this phase is related to the DCS via the relation \( \phi(\omega) = \arctan(\text{Re} \{ \chi_e(\omega) \}/\text{Im} \{ \chi_e(\omega) \}) \). To compare our predictions to the measurements performed in spin qubit systems, we generalize the description of the spinless DQD in series made above by taking spin into account.
account. This is simply done by considering triplet states in addition to the bonding and anti-bonding states which correspond to the singlet states in the case of a DQD with spin. The eigenenergy of the triplet state is given by 
\[ \lambda_T = (\varepsilon_d - i\Gamma)/2, \]
where \( \varepsilon_d = \varepsilon_2 - \varepsilon_1 \) is the detuning energy. Figure (6a) shows the \( \varepsilon_d \)-dispersion of the eigenenergies \( \lambda_\pm \) and \( \lambda_T \). At \( T = 0 \), we use a generalized expression for \( \chi_c(\omega) \) obtained from Eq. (7) by adding a term corresponding to the triplet state contribution according to \(-3\Gamma/(2\hbar\omega(\omega + i\Gamma)) \sum_{\alpha=L,R} \ln(1 - \hbar\omega(\omega + i\Gamma) A_T(\mu_\alpha)/\Gamma)\), where \( A_T(\varepsilon) = -\text{Im}(\mathcal{G}_T^\dagger(\varepsilon))/\pi \)
with \( \mathcal{G}_T(\varepsilon) = (\varepsilon - \lambda_T)^{-1} \). The results for \( \phi(\omega) \) as a function of detuning energy and frequency is shown in Fig. (6b). One observes a dip in phase inside two pockets spreading out symmetrically around the vertical axis \( \varepsilon_d = 0 \), the two pockets being separated by a gap area located around \( \hbar\omega = \nu_{12} \), spotted by the dashed line. The existence of this gap is a direct consequence of the presence of bonding and anti-bonding states whereas the formation of the low-frequency pocket below the gap results from the presence of the triplet state. Our prediction for \( \phi(\omega) \) is in good qualitative agreement with the experimental results obtained in spin qubit.

VI. CONCLUSION

We have developed a model to calculate the DCS in a nonequilibrium DQD. We have established a general expression for this quantity, which, at \( T = 0 \) is related to the spectral function contributions from the bonding and anti-bonding states, leading to the prediction of a splitting of the two branches of maxima of the DCS as a function of the energy levels of the dots resulting from the application of a finite bias voltage driving the DQD out-of-equilibrium. It would be interesting to check this prediction experimentally. In the low frequency regime, we have analyzed the results in terms of the capacitance and the resistance of the equivalent RC-circuit in both serial and parallel geometries. In the high frequency regime, by extending our results to take into account an additional triplet mode in order to describe spin qubits, we have deduced the evolution for \( \chi_c(\omega) \) as a function of \( \omega \) and \( \varepsilon_d \) and found a qualitative agreement with experimental results recently obtained. The approach presented here can be used in many other contexts: quantum dots with multiple energy levels, submitted to the application of a magnetic field, in the presence of exchange or Coulomb interactions, among others.

ACKNOWLEDGMENTS

We dedicate this article to Marc Sanquer †. We pay tribute to his memory for his role in the development of quantum nanoelectronics. Marc had played a major role in our interest in this topics and never ceased to provide us with encouragement, sharing with us his expertise and enthusiasm for this field of research. We would like to thank Romain Maurand for valuable discussions and help and Vyacheslavs Kashcheyevs for stimulating discussion.

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In this Supplemental Material, we give the details of the calculation concerning (i) the integration over Grassmann variables, (ii) the dynamical charge susceptibility $\chi_c(\omega)$ for both a DQD in series and in parallel, (iii) the analytical expression for $\chi_c(\omega)$ in the limit of zero temperature ($T = 0$), (iv) the capacitance $C$ and the charge relaxation resistance $R$ of the equivalent RC-circuit for aligned dot levels $\varepsilon_1 = \varepsilon_2$, (v) the determination of the phase response $\phi(\omega)$ of the corresponding resonator, and (vi) the expression for $\chi_c(\omega)$ in a DQD in series in the presence of additional triplet states.

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VII. FUNCTION INTEGRAL APPROACH

Within the functional integral approach, the partition function can be written as the integral over the Grassmann variables $d_i^\dagger$, $d_i$, $c_{ak}^\dagger$, $c_{ak}$, associated with the creation and annihilation operators $\hat{d}_{\alpha,i}$, $\hat{d}_i$ for the dots, and $\hat{c}_{\alpha k}^\dagger$, $\hat{c}_{\alpha k}$ for the leads. Indeed, one has

$$Z = \int \prod_{i=1,2} \mathcal{D} d_i \prod_{\alpha=L,R} \prod_{k<\alpha} \mathcal{D} c_{\alpha k} \exp \left( - \int_0^\beta d\tau \mathcal{L}(\tau) \right),$$

(S1)

where $\beta = 1/k_BT$, $\mathcal{D} d_i$ and $\mathcal{D} c_{\alpha k}$ are the differential elements defined as $\mathcal{D} d_i = dd_i^\dagger dd_i$, and $\mathcal{D} c_{\alpha k} = dc_{\alpha k}^\dagger dc_{\alpha k}$. $\mathcal{L}(\tau)$ is the Lagrangian of the system defined in three parts $\mathcal{L}(\tau) = \mathcal{L}_{\text{dots}}(\tau) + \mathcal{L}_{\text{leads}}(\tau) + \mathcal{L}_{\text{hop}}(\tau)$, where $\mathcal{L}_{\text{dots}}(\tau) = \sum_{i=1,2} d_i^\dagger \partial_\tau d_i - \hat{H}_{\text{dots}}$, $\mathcal{L}_{\text{leads}}(\tau) = \sum_{\alpha=L,R,k<\alpha} c_{\alpha k}^\dagger \partial_\tau c_{\alpha k} - \hat{H}_{\text{leads}}$ and $\mathcal{L}_{\text{hop}}(\tau) = -\hat{H}_{\text{hop}}$. One remarks that the sum of $\mathcal{L}_{\text{leads}}(\tau)$ and $\mathcal{L}_{\text{hop}}(\tau)$ can be written equivalently in the form of a perfect square from which a supplementary term is subtracted

$$\mathcal{L}_{\text{leads}}(\tau) + \mathcal{L}_{\text{hop}}(\tau) = \sum_{\alpha=L,R} \sum_{k<\alpha} \overline{c}_{\alpha k}^\dagger (\partial_\tau + \varepsilon_{\alpha k}) \overline{c}_{\alpha k} - \sum_{\alpha=L,R,k<\alpha} \sum_{i,j} V_{i,\alpha k}^\dagger (\partial_\tau + \varepsilon_{\alpha k})^{-1} V_{j,\alpha k} d_i^\dagger d_j,$$

(S2)

with $\overline{c}_{\alpha k} = c_{\alpha k} - \sum_{j=1,2} V_{j,\alpha k} (\partial_\tau + \varepsilon_{\alpha k})^{-1} d_j$. By integrating over the Grassmann variables $\overline{c}_{\alpha k}$ and $\overline{c}_{\alpha k}^\dagger$, one gets $Z = \int \mathcal{D} d_1 \mathcal{D} d_2 \exp(-\int_0^\beta d\tau \mathcal{L}_{\text{eff}}(\tau))$, within a constant multiplicative factor. $\mathcal{L}_{\text{eff}}(\tau)$ is an effective Lagrangian defined as

$$\mathcal{L}_{\text{eff}}(\tau) = \left( \begin{array}{cc} d_1^\dagger & d_2^\dagger \\ \partial_\tau + \varepsilon_1 + \Sigma_{11}(\tau) & \Sigma_{12}(\tau) \end{array} \right) \left( \begin{array}{c} d_1 \\ 0 \end{array} \right),$$

(S3)

where $\Sigma_{ij}(\tau) = \sum_{\alpha=L,R} \sum_{k<\alpha} V_{i,\alpha k}^\dagger g_{\alpha k}(\tau) V_{j,\alpha k}$, with $g_{\alpha k}(\tau) = - (\partial_\tau + \varepsilon_{\alpha k})^{-1}$ and $\Sigma_{ij}(\varepsilon) = \Sigma_{ij}(\varepsilon) + V_{i,\alpha k}^\dagger$. The retarded Green function in the dots can be obtained from $\mathcal{G}_{\alpha k}^R(\tau) = \int \mathcal{D} d_1 \mathcal{D} d_2 d_i^\dagger d_j \exp(-\int_0^\beta d\tau \mathcal{L}_{\text{eff}}(\tau))$ by performing the integration over the Grassmann variables $d_i$ and $d_j^\dagger$, and taking the Fourier transform, one gets

$$\mathcal{G}_{\alpha k}^R(\varepsilon) = \left( \begin{array}{cc} \varepsilon - \varepsilon_1 - \Sigma_{11}^R(\varepsilon) & -\Sigma_{12}^R(\varepsilon) \\ -\Sigma_{21}^R(\varepsilon) & \varepsilon - \varepsilon_2 - \Sigma_{22}^R(\varepsilon) \end{array} \right)^{-1}.$$

(S4)

From Eq. (S4) one extracts an effective Hamiltonian $\mathcal{H}_{\text{eff}}$ defined through the relation $\mathcal{G}_{\alpha k}^R(\varepsilon) = (\varepsilon - \mathcal{H}_{\text{eff}})^{-1}$. It is given by

$$\mathcal{H}_{\text{eff}} = \left( \begin{array}{cc} \varepsilon_1 - i \Gamma_{11}/2 & V_{12}^* - i \Gamma_{12}/2 \\ V_{21}^* - i \Gamma_{21}/2 & \varepsilon_2 - i \Gamma_{22}/2 \end{array} \right),$$

(S5)

since $\Sigma_{ij}(\varepsilon) = -i\Gamma_{ij}/2$ and $\Gamma_{ij} = \sum_{\alpha=L,R} \Gamma_{\alpha,ij}$, where $\Gamma_{\alpha,ij} = 2\pi V_{i,\alpha k}^* V_{j,\alpha k} \rho_{\alpha}$ does not depend on $k$ in the flat-wide-band approximation for electrons in the leads. The fact that $\mathcal{H}_{\text{eff}}$ is non-hermitian is the consequence of the integration over the lead Grassmann variables that has been performed. The coupling to the leads is thus explicitly taken into account in the Hamiltonian. It induces a dissipation since the eigenvalues of $\mathcal{H}_{\text{eff}}$ acquire an imaginary part

$$\lambda_{\pm} = \frac{1}{2} \left( \varepsilon_1 - i \Gamma_{11}/2 + \varepsilon_2 - i \Gamma_{22}/2 \pm \Delta \right),$$

(S6)

with $\Delta^2 = (\varepsilon_1 - i\Gamma_{11}/2 - \varepsilon_2 + i\Gamma_{22}/2)^2 + 4(V_{12}^* - i\Gamma_{12}/2)(V_{21}^* - i\Gamma_{21}/2)$. The eigenvalues $\lambda_-$ and $\lambda_+$ are respectively the energies of the bonding and anti-bonding states of the DQD system. The non-hermitian matrix $\Sigma$, of elements
$U_{ij}$, associated with this diagonalization is

$$U = \begin{pmatrix}
\frac{\mathcal{V}_{12}^* - i\Gamma_{12}/2}{D_1} & \frac{\delta}{2D_2} \\
-\frac{\delta}{2D_1} & \frac{\mathcal{V}_{21}^* - i\Gamma_{21}/2}{D_2}
\end{pmatrix},$$

(S7)

with

$$\delta = \varepsilon_1 - i \frac{\Gamma_{11}}{2} - \varepsilon_2 + i \frac{\Gamma_{22}}{2} - \Delta,$$

(S8)

$$D_1^2 = \left|\frac{\mathcal{V}_{12}^*}{2}\right|^2 + \left|\frac{\delta}{4}\right|^2.$$  

(S9)

VIII. GENERAL EXPRESSION FOR THE DYNAMICAL CHARGE SUSCEPTIBILITY

A. Calculation of $\chi_c(\omega)$

The charge susceptibility is given by $\chi_c(\omega) = \sum_{i,j=1,2} \alpha_i \alpha_j \chi_{ij}(\omega)$, where $\alpha_i$ are the level-arm coefficients measuring the asymmetry of the capacitive couplings, and $\chi_{ij}(\omega)$ is the Fourier transform of

$$\chi_{ij}(t, t') = i\Theta(t - t')\langle\{\Delta \hat{N}_j(t'), \Delta \hat{N}_i(t)\}\rangle,$$

(S10)

where $\Delta \hat{N}_i(t) = \hat{N}_i(t) - \langle\hat{N}_i\rangle$ with $\hat{N}_i(t) = \hat{d}_i^\dagger(t)\hat{d}_i(t)$, and $\langle, , \rangle$ denotes the anti-commutator. To calculate this quantity, one has first to calculate the time-ordered correlator defined as

$$\chi'_{ij}(\tau, \tau') = i \langle T_c[\Delta \hat{N}_j(\tau')\Delta \hat{N}_i(\tau)] \rangle,$$

(S11)

where $T_c$ is the time-ordered operator. By assuming that there is no Coulomb interaction, one can use the Wick theorem to calculate this correlator, thus

$$\chi'_{ij}(\tau, \tau') = i \langle T_c[\hat{d}_j^\dagger(\tau')\hat{d}_i(\tau)]\rangle \langle T_c[\hat{d}_j(\tau')\hat{d}_i^\dagger(\tau)]\rangle.$$

(S12)

By introducing the time-ordered Green functions defined as $G'_{ij}(\tau, \tau') = -i \langle T_c[\hat{d}_j^\dagger(\tau')\hat{d}_i(\tau)]\rangle$, one gets:

$$\chi_{ij}(\tau, \tau') = iG'_{ij}(\tau, \tau')G'_{ji}(\tau', \tau).$$

(S13)

In the basis of the eigenstates associated with the operators $\hat{d}_+$ and $\hat{d}_-$ in which $\mathcal{H}_{\text{eff}}$ is diagonal, one has

$$\hat{d}_1(t) = U_{11}\hat{d}_+(t) + U_{12}\hat{d}_-(t),$$  

(S14)

$$\hat{d}_2(t) = U_{21}\hat{d}_+(t) + U_{22}\hat{d}_-(t),$$  

(S15)

$$\hat{d}_1^\dagger(t) = U_{11}^*\hat{d}_+^\dagger(t) + U_{12}^*\hat{d}_-^\dagger(t),$$  

(S16)

$$\hat{d}_2^\dagger(t) = U_{21}^*\hat{d}_+^\dagger(t) + U_{22}^*\hat{d}_-^\dagger(t),$$  

(S17)

where $U_{ij}$ are the elements of the transition matrix $U$. Indeed, one has $\hat{d}_i = \sum_{s=\pm} \langle i | s \rangle \hat{d}_s$, with $\langle i | s \rangle = U_{is}$. It leads to

$$G'_{ij}(\tau, \tau') = U_{11}U_{j1}^*G_+^{\tau'}(\tau, \tau') + U_{12}U_{j2}^*G_-^{\tau'}(\tau, \tau'),$$  

(S18)

$$G'_{ji}(\tau', \tau) = U_{j1}U_{i1}^*G_+^{\tau}(\tau', \tau) + U_{j2}U_{i2}^*G_-^{\tau}(\tau', \tau),$$  

(S19)

where $G'_\pm(\tau, \tau') = -i\langle T_c[\hat{d}_\pm^\dagger(\tau')\hat{d}_\pm(\tau)]\rangle$. Finally, one has

$$\chi_{ij}(\tau, \tau') = iU_{11}^*U_{j1}^2G_+^{\tau'}(\tau, \tau')G_+^{\tau}(\tau', \tau) + iU_{12}^*U_{j2}^2U_{j2}G_-^{\tau'}(\tau, \tau')G_-^{\tau}(\tau', \tau),$$

(S20)

$$+iU_{11}^*U_{j1}U_{j2}G_+^{\tau'}(\tau, \tau')G_-^{\tau'}(\tau, \tau) + iU_{12}^*U_{j2}G_-^{\tau'}(\tau, \tau')G_-^{\tau'}(\tau, \tau),$$
which leads to
\[ \chi_\ell^\prime(\tau, \tau') = \sum_{s_1, s_2 = \pm} C_{s_1 s_2} \chi_{s_1 s_2}^\ell(\tau, \tau') , \] (S21)
where \( \chi_{s_1 s_2}^\ell(\tau, \tau') = i G_{s_1}^\ell(\tau, \tau') G_{s_2}^\ell(\tau', \tau) \) and \( C_{s_1 s_2} \) are coherence factors defined as
\[ C_{++} = \sum_{i,j=1,2} \alpha_i \alpha_j |U_{ij} U_{j' i'}|^2 , \] (S22)
\[ C_{+-} = \sum_{i,j=1,2} \alpha_i \alpha_j U_{i1} U_{j' 2} U_{1' j} U_{2' 2} , \] (S23)
\[ C_{-+} = \sum_{i,j=1,2} \alpha_i \alpha_j U_{i2} U_{j' 1} U_{1' j} U_{2' 2} , \] (S24)
\[ C_{--} = \sum_{i,j=1,2} \alpha_i \alpha_j |U_{i2} U_{j' 2}|^2 . \] (S25)

One underlines that \( C_{++} \) and \( C_{--} \) are both reals, and that \( C_{+-} \) is \( C_{-+} \). By performing an analytical continuation \( \Re \), one obtains the charge susceptibility \( \chi_c(\tau, \tau') = \sum_{s_1, s_2 = \pm} C_{s_1 s_2} \chi_{s_1 s_2}(\tau, \tau') \) with
\[ \chi_{s_1 s_2}(\tau, \tau') = i \left[ G_{s_1 s_2}^{<}(\tau, \tau') \xi_{s_2}^{<}(\tau) + \xi_{s_1}^{<}(\tau) \xi_{s_2}^{<}(\tau') + G_{s_1 s_2}^{<}(\tau, \tau') \right] \] (S26)

By taking the Fourier transform, one gets \( \chi_c(\omega) = \sum_{s_1, s_2 = \pm} C_{s_1 s_2} \chi_{s_1 s_2}(\omega) \) with
\[ \chi_{s_1 s_2}(\omega) = \int_{-\infty}^{\infty} \frac{d\epsilon}{2\pi} \left[ G_{s_1 s_2}^{<}(\epsilon) \xi_{s_2}^{<}(\epsilon - i\omega) + \xi_{s_1}^{<}(\epsilon + i\omega) \right] \] (S27)

The retarded and advanced Green functions \( G_{s_1}^{<}(\epsilon) = (\epsilon - \lambda_s)^{-1} \) and \( G_{s_2}^{<}(\epsilon) = (\epsilon - \lambda_s)^{-1} \) in the eigenstate basis associated with the operators \( \hat{d}_+ \) and \( \hat{d}_- \), where \( \lambda_s \) is given by Eq. (S6). In the steady state, the nonequilibrium Green function is given by
\[ \xi_{s_1}^{<}(\epsilon) = \xi_{s_2}^{<}(\epsilon) U^{-1}_1 e^{-\frac{\epsilon}{2T}} \xi_{s_2}^{<}(\epsilon) U_1 , \] (S28)
where \( \xi_{s_1}^{<}(\epsilon) = i \sum_{\alpha=L,R} f_\alpha(\epsilon) \Gamma_{s_2}^{-1} \) is the self-energy in the initial basis associated with the operators \( d_1 \) and \( d_2 \), \( f_\alpha(\epsilon) = [1 + \exp((\epsilon - \mu_\alpha)/k_B T)]^{-1} \) is the Fermi-Dirac distribution functions of the left (L) and right (R) leads, \( \mu_{L,R} \), the chemical potential, and \( T \), the temperature. \( G_{s_2}^{<}(\epsilon) \) is usually a non-diagonal matrix in the eigenstate basis.

### B. Coherence factors \( C_{s_1 s_2} \)

In case of symmetrical capacitive couplings, i.e. \( \alpha_1 = \alpha_2 = 1 \), one has
\[ C_{++} = (|U_{11}|^2 + |U_{12}|^2)^2 = 1 , \] (S29)
\[ C_{+-} = \sum_{i,j=1,2} U_{i1} U_{j' 2} U_{1' j} U_{2' 2} = C_{-+} , \] (S30)
\[ C_{--} = (|U_{21}|^2 + |U_{22}|^2)^2 = 1 . \] (S31)

In each of the possible geometries of the DQD, it leads to:

- In the case of a DQD in series with symmetrical real couplings, one has \( \Gamma_{L,11} = \Gamma_{R,22} \equiv \Gamma \), \( \Gamma_{L,22} = \Gamma_{R,11} = 0 \), \( \Gamma_{\alpha,12} = \Gamma_{\alpha,21} = 0 \), \( \forall \alpha \), and \( \mathcal{V}_{12} = \mathcal{V}_{21} \):
  \[ C_{+-} = C_{-+} = 0 , \] (S32)
so that only \( \chi_{+-}(\omega) \) and \( \chi_{-+}(\omega) \) contribute to the charge susceptibility according to \( \chi_c(\omega) = \chi_{+-}(\omega) + \chi_{-+}(\omega) \).

- In the case of a DQD in parallel with symmetrical real couplings, one has \( \Gamma_{\alpha,i j} \equiv \Gamma \), \( \forall \alpha, i, j \) and \( \mathcal{V}_{12} = \mathcal{V}_{21} \):
  \[ C_{+-} = C_{-+} = \frac{|(\mathcal{V}_{12} - i\Gamma)\delta|^2 - \text{Re}((\mathcal{V}_{12} + i\Gamma)^2\delta^2)}{2D^4} , \] (S33)
where \( D^2 = |\mathcal{V}_{12} - i\Gamma|^2 + |\delta|^2/4 \), \( \delta = \varepsilon_1 - \varepsilon_2 - \Delta \), and \( \Delta^2 = (\varepsilon_1 - \varepsilon_2)^2 + 4(\mathcal{V}_{12} - i\Gamma)^2 \). One takes note that at either \( \Gamma = 0 \) or \( \varepsilon_1 = \varepsilon_2 \), Eq. (S33) leads to \( C_{+-} = C_{-+} = 0 \).
FIG. S1. Coherence factor $C_{++} = C_{--}$ for a DQD coupled in parallel at $V_{12} = \Gamma = 0.1$.

**IX. EXPRESSION FOR $\chi_c(\omega)$ AT $T = 0$**

For a DQD system, one has $X_c(\omega) = \sum_{s_1, s_2 = \pm} C_{s_1 s_2} X_{s_1 s_2}(\omega)$ with

$$X_{s_1 s_2}(\omega) = i \int_{-\infty}^{\infty} \frac{d\varepsilon}{2\pi} \left[ G^<_{s_1 s_2}(\varepsilon) G^a_{s_2}(\varepsilon - \hbar \omega) + G^r_{s_1}(\varepsilon + \hbar \omega) G^<_{s_1 s_2}(\varepsilon) \right],$$  

(S34)

where $G^r_{\pm}(\varepsilon)$ and $G^a_{\pm}(\varepsilon)$ are the retarded and advanced Green functions in the dots given by $G^r_{\pm}(\varepsilon) = (\varepsilon - \lambda_{\pm})^{-1}$ and $G^a_{\pm}(\varepsilon) = (\varepsilon - \lambda_{\pm}^*)^{-1}$, and $G^<_{\pm}(\varepsilon)$ are the nonequilibrium Green functions in the dots which in the steady state are given by

$$G^<_{\pm}(\varepsilon) = G^r_{\pm}(\varepsilon) U^{-1} \Sigma^<_{\pm}(\varepsilon) U, \quad (S35)$$

with $\Sigma^<_{\pm}(\varepsilon) = i \sum_{\alpha = L, R} f_{\alpha}(\varepsilon) \Gamma_{\alpha}$. The matrices $G^r_{\pm}(\varepsilon)$ and $G^a_{\pm}(\varepsilon)$ are diagonal while $G^<_{\pm}(\varepsilon)$ may be non-diagonal.

We set the constant $\hbar$ to 1 in the details of the calculation but we will restore this omitted factor $\hbar$ in the final expressions of the results.

**A. DQD in series**

For a DQD in series with symmetrically capacitive couplings $\alpha_1 = \alpha_2 = 1$, one has $\Gamma_{L, 11} = \Gamma_{R, 22} = \Gamma$ and $\Gamma_{L, 22} = \Gamma_{R, 11} = \Gamma_{\alpha, 12} = 0$, thus:

- Coherence factors: $C_{++} = C_{--} = 1$ and $C_{+-} = C_{-+} = 0$.
- $\Gamma_L$ and $\Gamma_R$ are diagonal matrices:

$$\Gamma_L = \begin{pmatrix} \Gamma & 0 \\ 0 & 0 \end{pmatrix}, \quad \Gamma_R = \begin{pmatrix} 0 & 0 \\ 0 & \Gamma \end{pmatrix} \Rightarrow \Sigma^<_{\pm}(\varepsilon) = \begin{pmatrix} i \Gamma f_L(\varepsilon) & 0 \\ 0 & i \Gamma f_R(\varepsilon) \end{pmatrix},$$  

(S36)

- Eigenvalues:

$$\lambda_{\pm} = \frac{1}{2} \left( \varepsilon_1 + \varepsilon_2 - i \Gamma \pm \sqrt{(\varepsilon_1 - \varepsilon_2)^2 + 4|V_{12}|^2} \right).$$  

(S37)

- Diagonalization matrix:

$$U = \frac{1}{\sqrt{|V_{12}|^2 + \delta^2/4}} \begin{pmatrix} V_{12}^* & \delta \\ \frac{\delta}{2} & V_{12} \end{pmatrix} \Rightarrow U^{-1} = \frac{1}{\sqrt{|V_{12}|^2 + \delta^2/4}} \begin{pmatrix} V_{12}^* & \frac{\delta}{2} \\ \frac{\delta}{2} & V_{12} \end{pmatrix},$$  

(S38)

with $\delta = \varepsilon_1 - \varepsilon_2 - \Delta$, $\Delta^2 = (\varepsilon_1 - \varepsilon_2)^2 + 4|V_{12}|^2$, and $V_{21}^* = V_{12}$. 


It leads to

\[
\mathcal{G}_G^\alpha(\varepsilon) = \frac{\Gamma}{|V_{12}|^2 + \delta^2/4} \left( \begin{array}{ccc} G_\alpha^G(\varepsilon) & 0 & f_L(\varepsilon) \\ 0 & G_\alpha^G(\varepsilon) & 0 \\ \frac{\delta}{2} & \frac{\delta}{2} & V_{12} \end{array} \right) \left( \begin{array}{ccc} \frac{\delta}{2} & \frac{\delta}{2} & V_{12} \\ 0 & f_R(\varepsilon) & 0 \\ V_{12} & 0 & 0 \\ \frac{\delta}{2} & \frac{\delta}{2} & V_{12} \end{array} \right) \left( \begin{array}{ccc} G_\alpha^G(\varepsilon) & 0 & 0 \\ 0 & G_\alpha^G(\varepsilon) & 0 \\ \frac{\delta}{2} & \frac{\delta}{2} & V_{12} \end{array} \right) \left( \begin{array}{ccc} \frac{\delta}{2} & \frac{\delta}{2} & V_{12} \\ 0 & f_R(\varepsilon) & 0 \\ V_{12} & 0 & 0 \\ \frac{\delta}{2} & \frac{\delta}{2} & V_{12} \end{array} \right) \right).
\]

By using this result, one gets at \( T = 0 \)

\[
\mathcal{X}_{++}(\omega) = -\frac{\Gamma}{|V_{12}|^2 + \delta^2/4} \int_{-\infty}^{\infty} \frac{d\varepsilon}{2\pi} |V_{12}|^2 f_L(\varepsilon) + \delta^2 f_R(\varepsilon)/4) |G_\alpha^G(\varepsilon)|^2 \left[ G_\alpha^G(\varepsilon - \omega) + G_\alpha^G(\varepsilon + \omega) \right]
\]

and

\[
\mathcal{X}_{--}(\omega) = -\frac{\Gamma}{|V_{12}|^2 + \delta^2/4} \int_{-\infty}^{\infty} \frac{d\varepsilon}{2\pi} |V_{12}|^2 f_L(\varepsilon) + \delta^2 f_R(\varepsilon)/4) |G_\alpha^G(\varepsilon)|^2 \left[ G_\alpha^G(\varepsilon - \omega) + G_\alpha^G(\varepsilon + \omega) \right]
\]

One has to calculate the following integral

\[
I_{\alpha,\pm}(\omega) = \int_{-\infty}^{\mu_\alpha} \frac{d\varepsilon}{2\pi} |G_\alpha^G(\varepsilon)|^2 \left[ G_\alpha^G(\varepsilon - \omega) + G_\alpha^G(\varepsilon + \omega) \right].
\]

From Eq. [S37], one can write

\[
|G_\alpha^G(\varepsilon)|^2 = |(\varepsilon - \lambda_{\pm})^{-1}|^2 = \frac{1}{(\varepsilon - \text{Re}\{\lambda_{\pm}\})^2 + \Gamma^2/4},
\]

and

\[
G_\alpha^G(\varepsilon - \omega) + G_\alpha^G(\varepsilon + \omega) = \frac{1}{\varepsilon - \omega - \text{Re}\{\lambda_{\pm}\} - i\Gamma/2} + \frac{1}{\varepsilon + \omega - \text{Re}\{\lambda_{\pm}\} + i\Gamma/2}.
\]

Thus

\[
I_{\alpha,\pm}(\omega) = \int_{-\infty}^{\mu_\alpha} \frac{d\varepsilon}{2\pi} (\varepsilon - \text{Re}\{\lambda_{\pm}\})^2 + \Gamma^2/4 \left( \frac{1}{\varepsilon - \omega - \text{Re}\{\lambda_{\pm}\} - i\Gamma/2} + \frac{1}{\varepsilon + \omega - \text{Re}\{\lambda_{\pm}\} + i\Gamma/2} \right).
\]

By performing this integral explicitly, one gets

\[
I_{\alpha,\pm}(\omega) = \frac{1}{4\pi\omega(T - i\omega)} \left[ 2 \arctan \left( \frac{\Gamma/2}{\text{Re}\{\lambda_{\pm}\} - \mu_\alpha + \omega} \right) - 2 \arctan \left( \frac{\Gamma/2}{\text{Re}\{\lambda_{\pm}\} - \mu_\alpha - \omega} \right) + i \ln \left( \frac{(\mu_\alpha - \text{Re}\{\lambda_{\pm}\})^2 + \Gamma^2/4}{(\mu_\alpha + \omega - \text{Re}\{\lambda_{\pm}\})^2 + \Gamma^2/4} \right) \right].
\]
which can be written as
\[ I_{\alpha, \pm}(\omega) = -\frac{1}{2\pi\omega(\omega + i\Gamma)} \ln \left( \frac{(\mu_\alpha - \text{Re}(\lambda_\pm))^2 + \Gamma^2/4}{(\mu_\alpha - \omega - \text{Re}(\lambda_\pm) - i\Gamma/2) (\mu_\alpha + \omega - \text{Re}(\lambda_\pm) + i\Gamma/2)} \right), \]  
(S47)

or as well
\[ I_{\alpha, +}(\omega) = -\frac{1}{2\pi\omega(\omega + i\Gamma)} \ln \left( \frac{|\mu_\alpha - \lambda_\pm|^2}{(\mu_\alpha - \omega - \lambda_\pm) (\mu_\alpha + \omega - \lambda_\pm)} \right) = \frac{1}{2\pi\omega(\omega + i\Gamma)} \ln \left( \frac{(1 - \omega G^+_\pm(\mu_\alpha))(1 + \omega G^-_\pm(\mu_\alpha)))}{(1 - \omega^2|G^+_\pm(\mu_\alpha)|^2)} \right). \]  
(S48)

It can be written equivalently in the form
\[ I_{\alpha, \pm}(\omega) = \frac{1}{2\pi\omega(\omega + i\Gamma)} \ln \left( 1 - \frac{2\pi\omega(\omega + i\Gamma)}{\Gamma} A_{\pm}(\mu_\alpha) \right). \]  
(S49)

since one has \( A_{\pm}(\varepsilon) = -\text{Im}\{G^\pm_\pm(\varepsilon)\}/\pi \) and for a DQD in series
\[ G^+_\pm(\mu_\alpha) = \frac{1}{\mu_\alpha - \lambda_\pm} = \frac{1}{\mu_\alpha - \text{Re}(\lambda_\pm) + i\Gamma/2} = (\mu_\alpha - \text{Re}(\lambda_\pm) - i\Gamma/2)|G^+_\pm(\mu_\alpha)|^2 \]
\[ \Rightarrow \text{Im}\{G^\pm_\pm(\mu_\alpha)\} = -\frac{\Gamma}{2}|G^+_\pm(\mu_\alpha)|^2 \Rightarrow 2\text{Im}\{G^+_\pm(\mu_\alpha)\} + \Gamma|G^+_\pm(\mu_\alpha)|^2 = 0. \]  
(S50)

Finally, one gets \( \chi_c(\omega) = \chi_{++}(\omega) + \chi_{--}(\omega) \) with
\[ \chi_{++}(\omega) = -\frac{\Gamma}{2\pi\omega(\omega + i\Gamma)(|\mathcal{V}_{12}|^2 + \delta^2/4)} \left( |\mathcal{V}_{12}|^2 \ln \left( 1 - \frac{2\pi\omega(\omega + i\Gamma)}{\Gamma} A_+(\mu_L) \right) + \frac{\delta^2}{4} \ln \left( 1 - \frac{2\pi\omega(\omega + i\Gamma)}{\Gamma} A_+(\mu_R) \right) \right), \]  
(S51)
\[ \chi_{--}(\omega) = -\frac{\Gamma}{2\pi\omega(\omega + i\Gamma)(|\mathcal{V}_{12}|^2 + \delta^2/4)} \left( |\mathcal{V}_{12}|^2 \ln \left( 1 - \frac{2\pi\omega(\omega + i\Gamma)}{\Gamma} A_-(\mu_R) \right) + \frac{\delta^2}{4} \ln \left( 1 - \frac{2\pi\omega(\omega + i\Gamma)}{\Gamma} A_-(\mu_L) \right) \right). \]  
(S52)

The expression for \( \chi_c(\omega) \) can be simplified in the following two special cases (we restore the omitted \( \hbar \) factor):
- For \( \varepsilon_1 = \varepsilon_2 = \varepsilon_0 \), one has \( \delta = -2|\mathcal{V}_{12} | \), which leads to
\[ \chi_c(\omega) = -\frac{\Gamma}{2\hbar\omega(\hbar\omega + i\Gamma)} \sum_{\pm} \sum_{\alpha = L, R} \ln \left( 1 - \frac{\hbar\omega(\hbar\omega + i\Gamma)}{\Gamma} A_{\pm}(\mu_\alpha) \right). \]  
(S53)

- For \( \mu_L = \mu_R = 0 \) (zero-voltage), one gets whatever \( \varepsilon_1 \) and \( \varepsilon_2 \) are:
\[ \chi_c(\omega) = -\frac{\Gamma}{\hbar\omega(\hbar\omega + i\Gamma)} \sum_{\pm} \ln \left( 1 - \frac{\hbar\omega(\hbar\omega + i\Gamma)}{\Gamma} A_{\pm}(0) \right). \]  
(S54)

### B. DQD in parallel

For a DQD in parallel with symmetrically capacitive couplings: \( \alpha_1 = \alpha_2 = 1 \), identical couplings between each dot and each lead: \( \Gamma_{\alpha, ij} \equiv \Gamma \) for \( \forall \alpha, i, j \), and assuming coupling \( \mathcal{V}_{12} \) to be real which implies \( \mathcal{V}_{12} = \mathcal{V}_{21} \), one has:
- Coherence factors: \( C_{++} = C_{--} = 1 \) and \( C_{+-} = C_{-+} = 0 \) since one has from Eq. (S33)
\[ C_{+-} = C_{-+} = \frac{|(\mathcal{V}_{12} - i\Gamma)\delta|^2 - \text{Re}(\mathcal{V}_{12} + i\Gamma)^2\delta^2}{2D^4}, \]  
(S55)

where \( D^2 = |\mathcal{V}_{12} - i\Gamma|^2 + |\delta|^2/4 \), \( \delta = \varepsilon_1 - \varepsilon_2 - \Delta \), and \( \Delta^2 = (\varepsilon_1 - \varepsilon_2)^2 + 4|\mathcal{V}_{12} - i\Gamma|^2 \).
For \( \varepsilon_1 = \varepsilon_2 = \varepsilon_0 \), it leads to \( \Delta = 2|\mathcal{V}_{12} - i\Gamma| \), \( \delta = -\Delta \) and \( D^2 = 2|\mathcal{V}_{12} - i\Gamma|^2 \), thus \( C_{+-} = C_{-+} = 0 \).
\(\Gamma_L\) and \(\Gamma_R\) are non-diagonal matrices equal to:

\[
\Gamma_L = \begin{pmatrix} \Gamma & \Gamma \\ \Gamma & \Gamma \end{pmatrix}, \quad \Gamma_R = \begin{pmatrix} \Gamma & \Gamma \\ \Gamma & \Gamma \end{pmatrix} \Rightarrow \sum^< (\varepsilon) = i \sum_{\alpha=L,R} f_{\alpha}(\varepsilon) \begin{pmatrix} \Gamma & \Gamma \\ \Gamma & \Gamma \end{pmatrix}.
\] (S56)

- Eigenvalues:

\[
\lambda_\pm = \frac{1}{2} \left( \varepsilon_1 + \varepsilon_2 - 2i\Gamma \pm \sqrt{(\varepsilon_1 - \varepsilon_2)^2 + 4(\nu_{12} - i\Gamma)^2} \right).
\] (S57)

- Diagonalization matrix:

\[
U = \frac{1}{D} \begin{pmatrix} \nu_{12} - i\Gamma & \frac{\delta}{2} \\ -\frac{\delta}{2} & \nu_{12} - i\Gamma \end{pmatrix}, \quad U^{-1} = \frac{D}{(\nu_{12} - i\Gamma)^2 + \delta^2/4} \begin{pmatrix} \nu_{12} - i\Gamma & -\frac{\delta}{2} \\ \frac{\delta}{2} & \nu_{12} - i\Gamma \end{pmatrix}.
\] (S58)

It leads to

\[
\sum^< (\varepsilon) = \frac{i \sum_{\alpha=L,R} f_{\alpha}(\varepsilon)}{(\nu_{12} - i\Gamma)^2 + \delta^2/4} \begin{pmatrix} G^p_\alpha(\varepsilon) & 0 \\ 0 & G^\sigma_\alpha(\varepsilon) \end{pmatrix} \begin{pmatrix} \nu_{12} - i\Gamma & -\frac{\delta}{2} \\ \frac{\delta}{2} & \nu_{12} - i\Gamma \end{pmatrix} \begin{pmatrix} 1 & 1 \\ 1 & 1 \end{pmatrix} \begin{pmatrix} \nu_{12} - i\Gamma & \frac{\delta}{2} \\ -\frac{\delta}{2} & \nu_{12} - i\Gamma \end{pmatrix} \begin{pmatrix} G^p_\alpha(\varepsilon) & 0 \\ 0 & G^\sigma_\alpha(\varepsilon) \end{pmatrix}
\]

\[
= \frac{i \sum_{\alpha=L,R} f_{\alpha}(\varepsilon)}{(\nu_{12} - i\Gamma)^2 + \delta^2/4} \begin{pmatrix} G^p_\alpha(\varepsilon) & 0 \\ 0 & G^\sigma_\alpha(\varepsilon) \end{pmatrix} \begin{pmatrix} \nu_{12} - i\Gamma & -\frac{\delta}{2} \\ \frac{\delta}{2} & \nu_{12} - i\Gamma \end{pmatrix} \begin{pmatrix} 1 & 1 \\ 1 & 1 \end{pmatrix} \begin{pmatrix} \nu_{12} - i\Gamma & \nu_{12} - \frac{\delta}{2} \\ -\nu_{12} - \frac{\delta}{2} & \nu_{12} - i\Gamma \end{pmatrix} \begin{pmatrix} G^p_\alpha(\varepsilon) & \frac{\delta}{2}G^\sigma_\alpha(\varepsilon) \\ \frac{\delta}{2}G^\sigma_\alpha(\varepsilon) & (\nu_{12} - i\Gamma)G^p_\alpha(\varepsilon) \end{pmatrix}
\]

\[
= \frac{i \sum_{\alpha=L,R} f_{\alpha}(\varepsilon)}{(\nu_{12} - i\Gamma)^2 + \delta^2/4} \begin{pmatrix} G^p_\alpha(\varepsilon) & 0 \\ 0 & G^\sigma_\alpha(\varepsilon) \end{pmatrix} \begin{pmatrix} \nu_{12} - i\Gamma & -\frac{\delta}{2} \\ \frac{\delta}{2} & \nu_{12} - i\Gamma \end{pmatrix} \begin{pmatrix} 1 & 1 \\ 1 & 1 \end{pmatrix} \begin{pmatrix} \nu_{12} - i\Gamma & (\nu_{12} - \frac{\delta}{2})G^p_\alpha(\varepsilon) \\ -(\nu_{12} - \frac{\delta}{2})G^p_\alpha(\varepsilon) & (\nu_{12} - \frac{\delta}{2})G^\sigma_\alpha(\varepsilon) \end{pmatrix}
\]

\[
= \frac{i \sum_{\alpha=L,R} f_{\alpha}(\varepsilon)}{(\nu_{12} - i\Gamma)^2 + \delta^2/4} \begin{pmatrix} G^p_\alpha(\varepsilon) & 0 \\ 0 & G^\sigma_\alpha(\varepsilon) \end{pmatrix} \begin{pmatrix} \nu_{12} - i\Gamma & -\frac{\delta}{2} \right)^2 G^p_\alpha(\varepsilon) \\ (\nu_{12} - i\Gamma)^2 - \frac{\delta^2}{4} \right)G^\sigma_\alpha(\varepsilon) \end{pmatrix}
\]

which gives when \(\varepsilon_1 = \varepsilon_2 = \varepsilon_0\)

\[
\sum^< (\varepsilon) = 2i\Gamma \sum_{\alpha=L,R} f_{\alpha}(\varepsilon) \begin{pmatrix} |G^p_\alpha(\varepsilon)|^2 & 0 \\ 0 & 0 \end{pmatrix},
\] (S60)

since one has \(\delta = -\Delta = -2(\nu_{12} - i\Gamma)\) in that case.

Using this result, one gets for \(\mathcal{X}_c(\omega)\) at \(T = 0\)

\[
\mathcal{X}_c(\omega) = -2\Gamma \sum_{\alpha=L,R} \frac{d^\alpha f_{\alpha}(\varepsilon) |G^p_\alpha(\varepsilon)|^2 [G^p_\alpha(\varepsilon - \omega) + G^p_\alpha(\varepsilon + \omega)]}{d\varepsilon}.
\]

\[
-2\Gamma \sum_{\alpha=L,R} \left[ (\nu_{12} - i\Gamma)^2 - \frac{\delta^2}{4} \right] |G^\sigma_\alpha(\varepsilon)|^2 + (\nu_{12} - i\Gamma)^2 - \frac{\delta^2}{4} \right] G^\sigma_\alpha(\varepsilon) G^p_\alpha(\varepsilon)
\]

By performing this integral explicitly, one gets

\[
\mathcal{X}_c(\omega) = -\frac{2\Gamma}{2\pi\omega(\omega + 4i\Gamma)} \sum_{\alpha=L,R} \ln \left( (1 - \omega G^\sigma_\alpha(\mu_\alpha)) (1 + \omega G^\sigma_\alpha(\mu_\alpha)) \right)
\]

\[
= -\frac{2\Gamma}{2\pi\omega(\omega + 4i\Gamma)} \sum_{\alpha=L,R} \ln \left( 1 - \omega \left( G^\sigma_\alpha(\mu_\alpha) - G^\sigma_\alpha(\mu_\alpha) \right) - \omega^2 |G^\sigma_\alpha(\mu_\alpha)|^2 \right).
\] (S62)
which can be written equivalently in the form

\[
\mathcal{X}_c(\omega) = -\frac{2\Gamma}{\hbar(\hbar\omega + 4i\Gamma)} \sum_{\alpha=L,R} \ln \left( 1 - \frac{\hbar\omega(h\omega + 4i\Gamma)}{4\Gamma} A_\pm(\mu_\alpha) \right) \tag{S63}
\]

since one has for a DQD in parallel

\[
G_+^\prime(\mu_\alpha) = \frac{1}{\mu_\alpha - \lambda_+} = \frac{1}{\mu_\alpha - \text{Re}\{\lambda_+\} + 2i\Gamma} = (\mu_\alpha - \text{Re}\{\lambda_+\} - 2i\Gamma)|G_+^\prime(\mu_\alpha)|^2

\Rightarrow \text{Im}\{G_+^\prime(\mu_\alpha)\} = -2\Gamma|G_+^\prime(\mu_\alpha)|^2 \Rightarrow \text{Im}\{G_+^\prime(\mu_\alpha)\} + 2\Gamma|G_+^\prime(\mu_\alpha)|^2 = 0 . \tag{S64}
\]

X. EQUIVALENT RC-CIRCUIT AT \( T = 0 \) AND \( \varepsilon_1 = \varepsilon_2 = \varepsilon_0 \)

In this section, we calculate the capacitance and the resistance of the equivalent quantum RC-circuit which can be used to account for the behavior of the DQD system in the low-frequency regime.\(^3\)

A. DQD in series

1. \( C \) for a DQD in series

One has \( C = e^2\mathcal{X}_c(\omega = 0) \). One gets from Eq. (S53)

\[
C = -\lim_{\omega \to 0} \frac{e^2\Gamma}{2\hbar(\hbar\omega + i\Gamma)} \sum_{\pm} \sum_{\alpha=L,R} \ln \left( 1 - \frac{\hbar\omega(\hbar\omega + i\Gamma)}{\Gamma} A_\pm(\mu_\alpha) \right) \Rightarrow C = \frac{e^2}{2} \sum_{\pm} \sum_{\alpha=L,R} A_\pm(\mu_\alpha) \tag{S65}
\]

using \( \ln(1 + x) \approx x \).

2. \( R \) for a DQD in series

One has

\[
R = \frac{e^2}{C^2} \lim_{\omega \to 0} \frac{\text{Im}\{\mathcal{X}_c(\omega)\}}{\omega} . \tag{S66}
\]

By using \( \ln(1 + x) \approx x - x^2/2 \), one gets from Eq. (S53)

\[
\mathcal{X}_c(\omega) = -\frac{\Gamma}{2\hbar(\hbar\omega + i\Gamma)} \sum_{\pm} \sum_{\alpha=L,R} \left( \frac{\hbar\omega(\hbar\omega + i\Gamma)}{\Gamma} A_\pm(\mu_\alpha) - \frac{\hbar^2\omega^2(i\Gamma)^2}{2\Gamma^2} A_\pm^2(\mu_\alpha) \right) . \tag{S67}
\]

It leads to

\[
\mathcal{X}_c(\omega) = -\frac{\Gamma(\hbar\omega - i\Gamma)}{2\hbar(\hbar^2\omega^2 + \Gamma^2)} \sum_{\pm} \sum_{\alpha=L,R} \left( \frac{\hbar\omega(\hbar\omega + i\Gamma)}{\Gamma} A_\pm(\mu_\alpha) + \frac{\hbar^2\omega^2(i\Gamma)^2}{2\Gamma} A_\pm^2(\mu_\alpha) \right) . \tag{S68}
\]

The imaginary part of \( \mathcal{X}_c(\omega) \) is then given by

\[
\text{Im}\{\mathcal{X}_c(\omega)\} = \frac{\Gamma^2}{4\hbar(\hbar^2\omega^2 + \Gamma^2)} \sum_{\pm} \sum_{\alpha=L,R} (\hbar^2\omega^2 A_\pm^2(\mu_\alpha)) = \frac{\hbar\omega\Gamma^2}{4(\hbar^2\omega^2 + \Gamma^2)} \sum_{\pm} \sum_{\alpha=L,R} A_\pm^2(\mu_\alpha) . \tag{S69}
\]

Thus

\[
\lim_{\omega \to 0} \frac{\text{Im}\{\mathcal{X}_c(\omega)\}}{\omega} = \frac{\hbar}{4} \sum_{\pm} \sum_{\alpha=L,R} A_\pm^2(\mu_\alpha) . \tag{S70}
\]
By using $C = (e^2/2) \sum_{\pm} \sum_{\alpha=L,R} A_{\pm}(\mu_\alpha)$, one finally gets

$$R = \frac{\hbar}{e^2} \frac{\sum_{\pm} \sum_{\alpha=L,R} A_{\pm}^2(\mu_\alpha)}{(\sum_{\pm} \sum_{\alpha=L,R} A_{\pm}(\mu_\alpha))^2}$$

(S71)

The Cauchy-Schwarz inequality \((\sum_{i=1}^n x_i^2)(\sum_{i=1}^n y_i^2) \geq (\sum_{i=1}^n x_i y_i)^2\) for \(x_i = A_{\pm}(\mu_\alpha)\) and \(y_i = 1\) leads to

$$\left( \sum_{\pm} \sum_{\alpha=L,R} A_{\pm}^2(\mu_\alpha) \right) \left( \sum_{\pm} \sum_{\alpha=L,R} 1^2 \right) \geq \left( \sum_{\pm} \sum_{\alpha=L,R} A_{\pm}(\mu_\alpha) \times 1 \right)^2$$

$$\Rightarrow \frac{\sum_{\pm} \sum_{\alpha=L,R} A_{\pm}^2(\mu_\alpha)}{(\sum_{\pm} \sum_{\alpha=L,R} A_{\pm}(\mu_\alpha))^2} \geq \frac{1}{4} \Rightarrow R \geq \frac{\hbar}{4e^2},$$

(S72)

Moreover, in all generality one has \(\sum_{i=1}^n x_i^4 \leq (\sum_{i=1}^n x_i^2)^2\), thus

$$\sum_{\pm} \sum_{\alpha=L,R} |G_{\pm}(\mu_\alpha)|^4 \leq \sum_{\pm} \sum_{\alpha=L,R} |G_{\pm}^r(\mu_\alpha)|^2$$

$$\Rightarrow \frac{\sum_{\pm} \sum_{\alpha=L,R} A_{\pm}^2(\mu_\alpha)}{(\sum_{\pm} \sum_{\alpha=L,R} A_{\pm}(\mu_\alpha))^2} = \frac{\sum_{\pm} \sum_{\alpha=L,R} |G_{\pm}^r(\mu_\alpha)|^4}{(\sum_{\pm} \sum_{\alpha=L,R} |G_{\pm}^r(\mu_\alpha)|^2)^2} \leq 1 \Rightarrow R \leq \frac{\hbar}{e^2},$$

(S73)

since one has for a DQD in series: \(A_{\pm}(\varepsilon) = \left(-\text{Im}\{G_{\pm}^r(\varepsilon)\}/\pi = (\Gamma/2\pi)|G_{\pm}^r(\varepsilon)|^2\right)\).

In conclusion, one has \(R \in [\hbar/4e^2, \hbar/e^2]\) for a DQD in series at \(T = 0\) and \(\varepsilon_1 = \varepsilon_2\).

### B. DQD in parallel

#### 1. \(C\) for a DQD in parallel

One has \(C = e^2 \mathcal{X}_c(\omega = 0)\). One gets from Eq. (S63)

$$C = \frac{2e^2 \Gamma}{\hbar \omega (\hbar \omega + 4i \Gamma)} \sum_{\alpha=L,R} \ln \left( 1 - \frac{\hbar \omega (\hbar \omega + 4i \Gamma)}{4 \Gamma} A_{\pm}(\mu_\alpha) \right) \Rightarrow C = \frac{e^2}{2} \sum_{\alpha=L,R} A_{\pm}(\mu_\alpha)$$

(S74)

using \(\ln(1 + x) \approx x\).

#### 2. \(R\) for a DQD in parallel

One has

$$R = \frac{e^2}{C} \lim_{\omega \to 0} \frac{\text{Im}\{\mathcal{X}_c(\omega)\}}{\omega}.$$  

(S75)

By using \(\ln(1 + x) \approx x - x^2/2\), one gets from Eq. (S63)

$$\mathcal{X}_c(\omega) = -\frac{2 \Gamma}{\hbar \omega (\hbar \omega + 4i \Gamma)} \sum_{\alpha=L,R} \left( \frac{\hbar \omega (\hbar \omega + 4i \Gamma)}{\Gamma} A_{\pm}(\mu_\alpha) - \frac{\hbar^2 \omega^2 (4i \Gamma)^2}{32 \Gamma^2} A_{\pm}^2(\mu_\alpha) \right).$$  

(S76)

It leads to

$$\mathcal{X}_c(\omega) = -\frac{2 \Gamma (\hbar \omega - 4i \Gamma)}{\hbar \omega (\hbar \omega^2 + 16 \Gamma^2)} \sum_{\alpha=L,R} \left( -\frac{\hbar \omega (\hbar \omega + 4i \Gamma)}{\Gamma} A_{\pm}(\mu_\alpha) + \frac{\hbar^2 \omega^2}{2} A_{\pm}^2(\mu_\alpha) \right).$$  

(S77)

The imaginary part of \(\mathcal{X}_c(\omega)\) is then given by

$$\text{Im}\{\mathcal{X}_c(\omega)\} = \frac{4 \Gamma^2}{\hbar \omega (\hbar^2 \omega^2 + 16 \Gamma^2)} \sum_{\alpha=L,R} (\hbar^2 \omega^2 A_{\pm}^2(\mu_\alpha)) = \frac{4 \hbar \omega \Gamma^2}{(\hbar^2 \omega^2 + 16 \Gamma^2)} \sum_{\alpha=L,R} A_{\pm}^2(\mu_\alpha).$$  

(S78)
Thus
\[ \lim_{\omega \to 0} \frac{\text{Im}\{\mathcal{X}_c(\omega)\}}{\omega} = \frac{\hbar}{4} \sum_{\alpha=L,R} A^2_\alpha(\mu_\alpha) . \] (S79)

By using \( C = (e^2/2) \sum_{\alpha=L,R} A(\mu_\alpha) \), one finally gets
\[ R = \frac{\hbar}{e^2} \frac{\sum_{\alpha=L,R} A^2_\alpha(\mu_\alpha)}{\left( \sum_{\alpha=L,R} A_\alpha(\mu_\alpha) \right)^2} \] (S80)

The Cauchy-Schwarz inequality \((\sum_{i=1}^n x_i^2)(\sum_{i=1}^n y_i^2) \geq (\sum_{i=1}^n x_i y_i)^2\) for \( x_i = A_\alpha(\mu_\alpha) \) and \( y_i = 1 \) leads to
\[ \left( \sum_{\alpha=L,R} A^2_\alpha(\mu_\alpha) \right) \left( \sum_{\alpha=L,R} (1)^2 \right) \geq \left( \sum_{\alpha=L,R} A_\alpha(\mu_\alpha) \times 1 \right)^2 \]
\[ \Rightarrow \frac{\sum_{\alpha=L,R} A^2_\alpha(\mu_\alpha)}{(\sum_{\alpha=L,R} A_\alpha(\mu_\alpha))^2} \geq \frac{1}{2} \Rightarrow R \geq \frac{\hbar}{2e^2} , \] (S81)

Moreover, one has in all generality \( \sum_{i=1}^n x_i^4 \leq (\sum_{i=1}^n x_i^2)^2 \), thus
\[ \sum_{\alpha=L,R} |G_\alpha'(\mu_\alpha)|^4 \leq \left( \sum_{\alpha=L,R} |G_\alpha'(\mu_\alpha)|^2 \right)^2 \]
\[ \Rightarrow \frac{\sum_{\alpha=L,R} A^2_\alpha(\mu_\alpha)}{(\sum_{\alpha=L,R} A_\alpha(\mu_\alpha))^2} = \frac{\sum_{\alpha=L,R} |G_\alpha'(\mu_\alpha)|^4}{(\sum_{\alpha=L,R} |G_\alpha'(\mu_\alpha)|^2)^2} \leq 1 \Rightarrow R \leq \frac{\hbar}{e^2} , \] (S82)

since one has for a DQD in parallel: \( A_\alpha(\varepsilon) = -\text{Im}\{G_\alpha'(\varepsilon)\}/\pi = (2\Gamma/\pi)|G_\alpha'(\varepsilon)|^2 \).

In conclusion, one has \( R \in [h/2e^2, h/e^2] \) for a DQD in parallel at \( T = 0 \) and \( \varepsilon_1 = \varepsilon_2 \).

**XI. PHASE RESPONSE \( \phi(\omega) \) OF THE RESONATOR**

By definition, the phase response of the resonator is equal to
\[ \phi(\omega) = \arctan \left( \frac{\text{Im}\{Z(\omega)\}}{\text{Re}\{Z(\omega)\}} \right) , \] (S83)

where \( Z(\omega) \) is the impedance defined as \( Z(\omega) = dV_{\text{gate}}(\omega)/dI(\omega) \). Given that the dynamical charge susceptibility is \( \mathcal{X}_c(\omega) = dN(\omega)/d(eV_{\text{gate}}(\omega)) \)\(^\text{[3]}\) and that the electrical current is defined as \( I(t) = -dQ(t)/dt = -edN(t)/dt \Rightarrow I(\omega) = -ie\omega N(\omega) \), one gets
\[ \mathcal{X}_c(\omega) = \frac{dN(\omega)}{d(eV_{\text{gate}}(\omega))} = \frac{1}{e} \frac{dN(\omega)}{dI(\omega)} \frac{dI(\omega)}{dV_{\text{gate}}(\omega)} = -\frac{1}{ie^2\omega Z(\omega)} = \frac{iZ^*(\omega)}{e^2|\omega|Z(\omega)|^2} . \] (S84)

Thus
\[ \text{Re}\{Z(\omega)\} = e^2\omega|Z(\omega)|^2\text{Im}\{\mathcal{X}_c(\omega)\} \quad \text{and} \quad \text{Im}\{Z(\omega)\} = e^2\omega|Z(\omega)|^2\text{Re}\{\mathcal{X}_c(\omega)\} . \] (S85)

By incorporating these latter expressions into Eq. (S83), one gets
\[ \phi(\omega) = \arctan \left( \frac{\text{Re}\{\mathcal{X}_c(\omega)\}}{\text{Im}\{\mathcal{X}_c(\omega)\}} \right) . \] (S86)

**XII. EXPRESSION FOR \( \mathcal{X}_c(\omega) \) IN A DQD IN SERIES IN THE PRESENCE OF ADDITIONAL TRIPLET STATES**

In order to compare with experiments performed in spin qubits, one needs to take triplet states into account in addition to the singlet states. To do this, one includes an additional term to the expression for the dynamical charge
susceptibility given in (S54) for a DQD in series at $T = 0$ and when $\varepsilon_1 = \varepsilon_2$, corresponding to the triplet state contribution, as follows

$$\chi_c(\omega) = -\frac{\Gamma}{2\hbar\omega(\hbar\omega + i\Gamma)} \sum_{\alpha=L,R} \left( \sum_{\pm} \ln \left( 1 - \frac{\hbar\omega(\hbar\omega + i\Gamma)}{\Gamma} A_{\pm}(\mu_{\alpha}) \right) + 3 \ln \left( 1 - \frac{\hbar\omega(\hbar\omega + i\Gamma)}{\Gamma} A_T(\mu_{\alpha}) \right) \right), \quad (S87)$$

where $A_T(\varepsilon) = -\text{Im}\{G_T^R(\varepsilon)\}/\pi$ is the spectral function associated with the Green function $G_T^R(\varepsilon) = (\varepsilon - \lambda_T)^{-1}$ where $\lambda_T = (\varepsilon_2 - \varepsilon_1 - i\Gamma)/2$ is the energy of the triplet states. The factor 3 in front of the second term in the r.h.s. of Eq. (S87) results from the fact that the number of states in the triplet states is equal to 3.

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