Coherence and Decoherence in Tunnelling between Quantum Dots

D. M. Cardamone, C. A. Stafford, and B. R. Barrett

Physics Department, University of Arizona, 1118 East 4th Street, Tucson, Arizona 85721

(Submitted October 12, 2001)

Subject classification: 73.20.Jc; 73.40.Gk; 73.63.Kv

Abstract

Coupled quantum dots are an example of the ubiquitous quantum double potential well. In a typical transport experiment, each quantum dot is also coupled to a continuum of states. Our approach takes this into account by using a Green’s function formalism to solve the full system. The time-dependant solution is then explored in different limiting cases. In general, a combination of coherent and incoherent behavior is observed. In the case that the coupling of each dot to the macroscopic world is equal, however, the time evolution is purely coherent.

The double-well potential is one of the simplest and best understood problems in modern quantum mechanics. Its utility is likewise unparalleled. Potential applications of double-well devices have been noted in Refs. [1-5]. For such devices to be useful, an understanding of the processes which couple the microscopic device to the macroscopic environment is paramount. That is to say, the decoherence processes of such systems must be well understood. To this end, we consider a simple exactly-solvable model of two coupled quantum dots. An excellent review of related systems and some approximate solutions is given in Ref. [6].

Each dot is coupled to an environment (a triple-barrier system) as in Fig. 1. The environment consists of a continuum of states, as would be appropriate for a macroscopic lead. The initial condition is a single electron localized in dot 1 at time \( t = 0 \). Only one state in each quantum dot is considered, which amounts to the assumption that the tunnelling parameters connecting our two states to any neglected state are much less than the energy differences with that state.

The Hamiltonian of the system can be written as the sum of two terms \( H = H_0 + H_c \). The first term, \( H_0 \), represents the dynamics of the simple two-state system

\[
H_0 = \varepsilon_1 d_1^\dagger d_1 + \varepsilon_2 d_2^\dagger d_2 + V (d_1^\dagger d_2 + d_2^\dagger d_1) \approx \begin{pmatrix} \varepsilon_1 & V \\ V & \varepsilon_2 \end{pmatrix},
\]

where \( d_i^\dagger \) creates an electron in dot \( i \), and the other symbols are defined in Fig. 1. \( V \) may be chosen real and positive without loss of generality. The second term, \( H_c \), reflects the coupling of each dot to its respective lead

\[
H_c = \sum_{i=1}^{2} \sum_k (V_{ki} c_k^\dagger d_i + V_{ki}^* d_i^\dagger c_k + \varepsilon_{ki} c_k^\dagger c_k),
\]

where the sum on \( k \) runs over the states in the continuum adjoining dot \( i \), \( \varepsilon_{ki} \) is the energy of the \( k \)th state of the appropriate continuum, and \( V_{ki} \) is the tunnelling parameter from the dot to the continuum. Throughout the paper, we use units with \( \hbar = 1 \).

1Corresponding author; e-mail: dmcard@physics.arizona.edu
In the following, we adopt a Green’s function approach which parallels that taken in Ref. [7]. The retarded Green’s function is defined by
\[ G_{ij}(t) = -i\Theta(t)\langle \{ d_i(t), d_j^\dagger(0) \} \rangle, \]
where \( \Theta(t) \) is the well-known step function and \( \{ \cdot, \cdot \} \) denotes the anticommutator. It is a simple matter to calculate \( G_0(t) \) from \( H_0 \) in the Heisenberg picture. From this, one can find the Fourier transform
\[ G_0^{-1}(E) = \frac{1}{2\pi} \int_{-\infty}^{\infty} dt \, G_0(t) e^{iEt}, \]
and finally the inverse
\[ G^{-1}_0(E) = G_0^{-1}(E) - \Sigma. \]
Here \( \Sigma \) is the self-energy due to interactions with the continua
\[ \Sigma = \begin{pmatrix} -i\Gamma_1/2 & 0 \\ 0 & -i\Gamma_2/2 \end{pmatrix}. \]
In the broad-band limit for the continua, the \( \Gamma_i \)'s are independent of energy, and are [8]
\[ \Gamma_i = 2\pi \sum_k |V_{ki}|^2 \delta(E - \varepsilon_{ki}). \]
Application of Dyson’s equation and inversion yields the total Green’s function of energy
\[ G(E) \approx \left( E - \varepsilon_1 + i\frac{\Gamma_1}{2} \right) \left( E - \varepsilon_2 + i\frac{\Gamma_2}{2} \right) - V^2 \]
\[ \approx \begin{pmatrix} E - \varepsilon_1 + i\frac{\Gamma_1}{2} & -V \\ -V & E - \varepsilon_2 + i\frac{\Gamma_2}{2} \end{pmatrix}. \]
The general solution is given by the inverse Fourier transform of \( G(E) \), but the initial conditions make this necessary for only the first row. From the first row of \( G(t) \), we can determine the probabilities to find an electron in each quantum dot as a function of time:
\[ P_1(t) = |G_{11}(t)|^2 = \frac{V^2}{\omega_1^2} e^{-\Gamma_1 t} \left( \frac{1}{\Gamma_1 - \omega_1} e^{\omega_1 t} \right. \]
\[ \left. + \frac{1}{\omega_1 + \Gamma_1} e^{-\omega_1 t} + \frac{i\omega_1 + \Gamma_1}{i\omega_1 - \Gamma_1} e^{i\omega_1 t} + \frac{i\omega_1 - \Gamma_1}{i\omega_1 + \Gamma_1} e^{-i\omega_1 t} \right), \]
\[ P_2(t) = |G_{12}(t)|^2 = \frac{2V^2}{\omega_1^2} e^{-\Gamma_1 t} (\cosh \omega_1 t - \cos \omega_1 t). \]
where
\[ \Gamma \equiv \frac{\Gamma_1 + \Gamma_2}{2}, \quad \Gamma' \equiv \frac{\Gamma_2 - \Gamma_1}{2}, \] (11)
and
\[ \omega \equiv \omega_r + i\omega_i \equiv \sqrt{4V^2 + (\varepsilon_2 - \varepsilon_1 - i\Gamma')^2}. \] (12)

Some understanding of the physical meaning of the solution may be arrived at by examining the limiting case of identical quantum dots: i.e., \( \varepsilon_1 = \varepsilon_2 \). In that case, we have
\[ \omega = \sqrt{4V^2 - \Gamma'^2}, \] (13)
where \( \omega \) must be either purely real or purely imaginary. Working with \( P_2(t) \) for algebraic simplicity, we find that if \( 2V > |\Gamma'| \), \( \omega \) is purely real, and
\[ P_2(t) = \frac{4V^2}{\omega^2} e^{-\Gamma t} \sin^2 \frac{\omega t}{2}. \] (14)

In this limit, the solution exhibits the characteristic coherent behavior of Rabi oscillations, as one would expect from a simple two-well problem. The presence of the environment is visible in the exponentially decaying envelope function: electrons leave/enter the two-state system at an overall rate \( \bar{\Gamma} \).

In the case that \( \varepsilon_1 = \varepsilon_2 \) and \( 2V < |\Gamma'| \), on the other hand, we find a purely imaginary \( \omega \), and
\[ P_2(t) = \frac{4V^2}{|\omega|^2} e^{-\Gamma t} \sinh^2 \frac{|\omega| t}{2}. \] (15)

The hyperbolic function and the resulting lack of Rabi oscillations demonstrate that interdot tunneling is completely incoherent in this regime.

In the more general case that \( \varepsilon_1 \neq \varepsilon_2 \), neither the circular nor the hyperbolic functions vanish, giving a combination of both coherent and incoherent behavior. For \( \Gamma' = 0 \), however, a purely real \( \omega \) is recovered, and the coherent evolution of Eq. (14) is obeyed, with Rabi frequency
\[ \omega = \sqrt{4V^2 + (\varepsilon_1 - \varepsilon_2)^2}. \] (16)

This somewhat startling result is especially gratifying when viewed in the context of measurement theory: \( \Gamma' = 0 \) indicates that the coupling of the two-well system to the environment does not distinguish between the dots, and thus the coherence between the dots is in no way destroyed by the macroscopic world.

Additional insight may be gained by examining the poles of the Green’s function, which can be interpreted as complex energy levels of the system, or resonances [9]. Defining complex energy levels
\[ \tilde{\varepsilon}_i = \varepsilon_i - i\Gamma_i/2 \] (17)
for each dot by including the local self-energy, the poles of \( G(E) \) may be written as
\[ \varepsilon_{\pm} = \frac{\tilde{\varepsilon}_1 + \tilde{\varepsilon}_2}{2} \pm \sqrt{\left( \frac{\tilde{\varepsilon}_2 - \tilde{\varepsilon}_1}{2} \right)^2 + V^2}. \] (18)

This is the standard formula for the hybridization of two energy levels coupled by a matrix element \( V \), except that the energy levels here are complex (see Figure 2). That the poles of \( G(E) \) should have this form can be understood from Dyson’s equation, if the interdot tunneling is taken as the perturbation, rather than the coupling to the environment. Neglecting interdot tunneling, the Green’s function is diagonal, and consists of two independent Breit-Wigner resonances at \( \tilde{\varepsilon}_1 \) and \( \tilde{\varepsilon}_2 \). The interdot tunneling operator then hybridizes these
Figure 2: Hybridization of the complex energy levels $\tilde{\varepsilon}_i$ yields resonances $\varepsilon \pm$ with complex level-splitting $\omega$, the nature of which determines the behavior of the system.

resonances via Dyson’s equation, leading to the poles displayed in Eq. (18). From Eqs. (12) and (18), it is easy to see that the energy difference

$$\varepsilon_+ - \varepsilon_- = \omega.$$  \hspace{1cm} (19)

Thus $\omega$ can be interpreted as the complex level splitting between the hybridized states, corresponding to a complex Rabi frequency: $\omega_r$ is the frequency of real Rabi oscillations, while $\omega_i$ is the rate of incoherent tunneling.

We have presented an exactly-solvable model for the tunneling dynamics of a two-level quantum system coupled to a macroscopic environment, appropriate to describe a double quantum dot connected to conducting leads. Despite its simplicity, this model displays many of the essential features of more complicated models [6]. The tunneling dynamics is determined by the complex Rabi frequency given in Eq. (12). Our results can be interpreted in terms of measurement theory: interdot coherence is suppressed to the degree that the environment can distinguish between the two orbitals.

Acknowledgements The authors thank Ned S. Wingreen for useful conversations. This work was supported in part by NSF grants PHY0070858 and DMR0072703.

References