Imaginary Phases in Two-Level Model with Spontaneous Decay.

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Abstract

We study a two-level model coupled to the electromagnetic vacuum and to an external classic electric field with fixed frequency. The amplitude of the external electric field is supposed to vary very slow in time. Garrison and Wright [Phys. Lett. A128 (1988) 177] used the non-hermitian Hamiltonian approach to study the adiabatic limit of this model and obtained that the probability of this two-level system to be in its upper level has an imaginary geometric phase. Using the master equation for describing the time evolution of the two-level system we obtain that the imaginary phase due to dissipative effects is time dependent, in opposition to Garrison and Wright result. The present results show that the non-hermitian hamiltonian method should not be used to discuss the nature of the imaginary phases in open systems.

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1 Introduction

The natural approach to study quantum systems in contact with an environment is the density operator method. It allows one to study quantum states evolving from pure to mixed states. In general, exponential decay in the time evolution of the population at each excited state of the quantum system occurs due to the exchange between the system and its environment. An alternative approach to obtain those exponential decays in the probability is through complex energies obtained from a phenomenological non-hermitian Hamiltonian[1, 2, 3, 4]. The non-hermitian Hamiltonian method has been very fruitful in approaching various physical problems as, for example, the multiphoton ionization[2, 3, 4, 5] and free-electron laser theory[6, 7]. As mentioned by Baker in reference [3], it is expected that the non-hermitian Schrödinger equation be a bona fide description of the interaction among the parts of a system when the intervals of time are sufficiently short such that the coherence state of a subsystem is not destroyed by its interaction with the environment.

Since the disclosure of geometrical phases by M.V. Berry in 1984[8] in cyclic Hamiltonians evolving adiabatically, there has been significant search for geometrical phases in other physical contexts. For example, Joye et al.[9] and Berry[10] independently showed that the transition probability of instantaneous eigenstates of non-real Hamiltonians in the non-adiabatic regime gets an imaginary geometric phase. This imaginary geometric phase was measured by Zwanziger et al. in a two-level system[11]. A great part of the work done on geometric phases has been on pure states. Since the work by Uhlmann in 1986[12, 13], however, the study of holonomy has been extended to mixed states under unitary evolution[14, 15, 16]. More recently, Ericsson et al.[17] obtained the expression of the geometric phase of a quantum system interacting with its environment when the unitary evolution of the whole system (including the environment) is known. Certainly the discussion of holonomy in mixed states is a very interesting point, since the correct expression of the geometric phase for mixed states under unitary evolution is still under debate, as in a very recent work by Singh et al.[18]. However, that is not the issue of the present communication. Another equally interesting question about the imaginary phases in transition probabilities of open systems is the correctness of the application of the non-hermitian hamiltonian approach[19, 20, 21, 22] to discuss the nature of such phases.

In particular Garrison and Wright[19] used the non-hermitian Hamiltonian method to study a two-level model with linewidths in the presence of an external electric field, ob-
taining the adiabatic limit of the probability of this system being in its upper level, after
the external field has returned to its original configuration. They concluded that the de-
caying factor has an imaginary correction to Berry’s phases. Their result for the two-level
model coupled to an external classical electromagnetic field with spontaneous emissions
is already contained in the phases (75) and (76) of reference [3] for any open system de-
descibed by a non-hermitian Hamiltonian. It is the mathematical structure of these phases
in the non-hermitian Hamiltonian approach that makes the imaginary phase derived by
Garrison and Wright in reference [19] to have an imaginary geometric contribution.

Garrison and Wright mentioned in their conclusion of reference [19] that the nature
(time-dependent or path-dependent) of their imaginary phase due to dissipation effects
should be reexamined using the density matrix approach; this is the aim of the present
letter.

In reference [23] we studied the adiabatic limit of any periodic non-degenerate Hamil-
tonian using the density matrix approach (extending the discussion carried out by Born
and Fock in reference [24] to the density matrix in the basis of instantaneous eigenstates
of the Hamiltonian), and we concluded that for a quantum system to get an imaginary
correction to Berry’s phases in dissipative phenomena, the functions in the integrals of the
decay exponentials in the entries of the density matrix would have to satisfy special con-
ditions. Let \( e^{-\int_0^t dt' \varphi(t')} \) be a typical decreasing exponential originated by the presence of
the dissipative effects. For the integral in the exponential be written as a path-dependent
integral, the function \( \varphi(t) \) must have the form

\[
c(t) = \varphi_i(t) \frac{d}{dt}(\Psi_i(t)),
\]

and the functions \( \Psi_i(t) \) have to satisfy two conditions: \( i \) they must not be explicitly
time-dependent; \( ii \) the time-dependence of functions \( \Psi_i(t) \) must come only from their
dependence on the set of parameters \( \vec{k}(t) \equiv (k_1(t), k_2(t), \cdots, k_l(t)) \). We point out that
there is no restriction to the regime of the time variation of the set of parameters \( \vec{k}(t) \)
and that it has not to be a periodic function in time, i.e., the path in the \( \vec{k} \)-parameter
space has not to be closed (see details in reference [23]).

In reference [19], Garrison and Wright considered a two-level system interacting with
a classic external electromagnetic field

\[
\vec{E}_{\text{clas}}(t) = \Re \left[ \vec{e} \mathcal{E}(t)e^{i\mathcal{O}} \right],
\]

(2a)

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where the amplitude $\mathcal{E}(t)$ varies very slowly. The two energy states were supposed to have a linewidth. They used the time-dependent non-hermitian Schrödinger (Bethe-Lamb) equation in the rotating wave approximation (RWA)[19, 25]

$$i \frac{d}{dt} \begin{pmatrix} C_a(t) \\ C_b(t) \end{pmatrix} = \begin{pmatrix} -\frac{1}{2} i \gamma_a & V e^{i \Delta t} \\ V e^{-i \Delta t} & -\frac{1}{2} i \gamma_b \end{pmatrix} \begin{pmatrix} C_a(t) \\ C_b(t) \end{pmatrix},$$  \hspace{1cm} (2b)

where $\gamma_a$ and $\gamma_b$ are the decay rates for the upper and lower levels, respectively, and $C_a(t)$ and $C_b(t)$ are the corresponding probability amplitude of states $|a\rangle$ and $|b\rangle$. Moreover, $\Delta = \omega - \nu$, $\omega = (E_a - E_b)/\hbar$, $V = \vec{\mu} \cdot \vec{e} \mathcal{E}(t)/2\hbar$ and $\vec{\mu} = \vec{\mu}_{ab}$ is the electric dipole matrix element. Letting $\mathcal{E}(t) = \mathcal{E}_0 e^{i \phi(t)}$, where $\phi(0) = 0$ and $\phi(T) = 2\pi$, with $T \gg \frac{2\pi \hbar}{E_a - E_b}$ they obtained a complex Berry’s phase $\beta_-$ in the expression for the probability of the quantum system to be in the state $a$ at time $T$,

$$\beta_- = \frac{1}{2} [2\pi (1 - \cos(\theta_0))] = \frac{1}{2} [2\pi (1 - \frac{\Delta - i \delta}{\sqrt{|2V_0|^2 + (\Delta - i \delta)^2}})] \hspace{1cm} (2c)$$

where $\delta \equiv (\gamma_a - \gamma_b)/2$ and $V_0 = \vec{\mu} \cdot \vec{e} \mathcal{E}_0/2\hbar$.

In the present work we want to verify if the result about the path-dependence of the imaginary phases derived by non-hermitian method is faithful. To do so, we apply the density matrix approach presented in reference [23] to study the same physical problem considered in reference [19] with $\gamma_b = 0$. We wish to compare the nature (path or time-dependent) of the imaginary phases in the expression of the probability of the atomic electron, initially at state $|a\rangle$, being at the same state at time $t = T$, derived by the density matrix approach with the one obtained in the non-hermitian phenomenological hamiltonian framework.

Let the two-level system that describes an atomic electron be represented by the Hamiltonian $\mathbf{H}_e$. The atomic electron interacts with a classic external electromagnetic field $\vec{E}_{\text{clas}}(t)$ (see eq.(2a)) and with the electromagnetic vacuum (the vacuum electromagnetic field operator being represented by $\vec{E}_0(\vec{x})$). The total Hamiltonian for this model in the Schrödinger picture is[26]

$$\mathbf{H}_T = \mathbf{H}_e + \mathbf{H}_f + \mathbf{H}_{\text{int}},$$  \hspace{1cm} (3)
\[ \mathbf{H}_e = \frac{\mathbf{P}^2}{2m} + V(r), \]  
\[ \mathbf{H}_f = \sum_{\vec{k}} \sum_{\lambda=1}^{2} \hbar \omega_{\vec{k}\lambda} \left( a_{\vec{k}\lambda}^\dagger a_{\vec{k}\lambda} + \frac{1}{2} \right), \]  
and
\[ \mathbf{H}_{int} = -e \left( \vec{E}_{\text{clas}}(t) + \vec{E}_0(\vec{x}) \right) \cdot \vec{r}, \]  
where \( \vec{P} \) is the momentum operator associated to the atomic electron, \( V(r) \) is the spherical interaction potential between the electron and the rest of atom and \( \mathbf{H}_f \) is the Hamiltonian of the electromagnetic energy operator of the electromagnetic vacuum. The operator \( a_{\vec{k}\lambda}^\dagger \) (\( a_{\vec{k}\lambda} \)) creates (destroys) a photon with momentum \( \vec{k} \) in the polarization state \( \lambda \). We also have: \( \omega_{\vec{k}} = c |\vec{k}| \).

Following reference [19], the classic external electromagnetic field \( \vec{E}_{\text{clas}}(t) \) is given by eq. (2a) assuming that \( \mathcal{E}(t) \) varies very slowly and \( \vec{e} \) is constant. The electromagnetic field operators in the vacuum \( \vec{E}_0(\vec{x}) \) in the Schrödinger picture have the expansion
\[ \vec{E}_0(\vec{x}) = \sum_{\vec{k}} \sum_{\lambda=1}^{2} \hat{e}_{\vec{k}\lambda} \sqrt{\frac{\hbar \omega_{\vec{k}\lambda}}{2e_0 V}} \left( -i a_{\vec{k}\lambda}^\dagger e^{-i\vec{k} \cdot \vec{x}} + i a_{\vec{k}\lambda} e^{i\vec{k} \cdot \vec{x}} \right). \]

The dynamics of the density operator of the complete system is driven by the Liouville von-Neumann equation. Taking the trace over the electromagnetic degrees of freedom we obtain the master equation for the reduced density matrix of the atomic electron \( \rho(t) \) written in the basis of the eigenstates of \( \mathbf{H}_e \) (\( \mathbf{H}_e |i\rangle = E_i |i\rangle \)). The dynamics of \( \rho(t) \) in the electric dipole approximation and in the RWA is [27, 28]
\[ \frac{d}{dt} \rho(t) = i \left[ (\omega_0 + \Omega_+ + \Omega_-) \sigma_z, \rho(t) \right] - \frac{1}{i\hbar} \overline{\mu}_{ab} \cdot \vec{E}_{\text{clas}}(t) [\sigma_x, \mathcal{A}_0(t)] + \gamma \left( 2\sigma_- \rho(t) \sigma_+ - \{ \sigma_+, \sigma_- \}, \rho(t) \right) + \frac{\overline{\mu}_{ab} \cdot \vec{E}_{\text{clas}}(t)}{\hbar^2} \left( -2A(t) \rho(t) \right) + 2A(t) \sigma_x \rho(t) \sigma_x + B(t) \left[ \{ \sigma_y, \rho(t) \}, \sigma_x \right], \]  
where \( \sigma_z \) and \( \sigma_\pm \) are the Pauli matrices with \( \sigma_\pm = \frac{1}{2} (\sigma_x \pm i\sigma_y) \), \( \overline{\mu}_{ab} \) is the electric dipole matrix (we are supposing \( \overline{\mu}_{ab} \) to be real) with \( \overline{\mu}_{ab} \equiv e \langle a | \vec{r} | b \rangle (e > 0) \).
\[ \gamma = \frac{\pi}{\hbar^2} \sum_{\lambda=1}^{2} \int d^3 \vec{k} \eta(\vec{k}) |g(\vec{k}, \lambda)|^2 \delta(\omega_0 - \omega_{\vec{k}}), \]  

(6a)

being \( g(\vec{k}, \lambda) = -i \vec{e}(\vec{k}, \lambda) \cdot \vec{\mu}_{ab} \sqrt{\hbar \omega_{\vec{k}}/2a_0} \) and \( \eta(\vec{k}) \) is the density of states introduced in the integration,

\[ \omega_0 \equiv \frac{E_a - E_b}{2\hbar} = \frac{\omega}{2}, \]  

(6b)

\[ \Lambda_{ij}^0(t) \equiv e^{-\frac{t}{2}(E_i-E_j)} \rho_{ij}(0), \quad i, j = a, b, \]  

(6c)

and

\[ \Omega_+ = -\frac{\gamma}{\pi} \ln \left[ \left| \frac{\omega_c}{\omega_0} - 1 \right| \left( \frac{\omega_c}{\omega_0} + 1 \right) \right], \]  

(6d)

where \( \omega_c \) is the cutoff frequency that preserves the dipole approximation (\( \omega_c < c/a_0 \), with \( a_0 \) being the atomic Bohr radius). The cutoff \( \omega_c \) becomes a parameter of the effective model to be determined by fitting to experimental data. The term \( \Omega_+ \) corresponds to the frequency shift[27]. The elements \( \rho_{ij}(0) \) are the initial values of the entries of matrix \( \tilde{\rho}(t) \). The functions \( A(t) \) and \( B(t) \) appearing in eq.(5) are defined as:

\[ A(t) \equiv \int_0^t dt' \vec{\mu}_{ab} \cdot \vec{E}_{\text{clas}}(t') \cos \left( \frac{E_a - E_b}{\hbar}(t - t') \right) \right) and \( B(t) \equiv \int_0^t dt' \vec{\mu}_{ab} \cdot \vec{E}_{\text{clas}}(t') \sin \left( \frac{E_a - E_b}{\hbar}(t - t') \right). \]

We remind that the density matrix of a two-level model must satisfy two conditions:

(i) \( Tr(\tilde{\rho}(t)) = 1\) and \( ii) \rho_{ab}(t) = (\rho_{ab}(t))^*, \) where \( \rho_{ab}(t) \equiv \langle a | \tilde{\rho}(t) | b \rangle \). As a consequence of those conditions, the density matrix has only two independent elements; we choose \( \rho_{aa}(t) \) and \( \rho_{ab}(t) \) to be such elements. ¿From eq.(5) the time equations for those two elements are

\[ \frac{d}{dt} \rho_{aa}(t) = \left[ -2\gamma - 4\vec{\mu}_{ab} \cdot \vec{E}_{\text{clas}}(t) \Re \left( e^{i(E_a-E_b)t} G(t) \right) \right] \rho_{aa}(t) \]

\[ + 2\vec{\mu}_{ab} \cdot \vec{E}_{\text{clas}}(t) \Im \left[ e^{-i(E_a-E_b)t} \rho_{ab}(0) \right] + 2\vec{\mu}_{ab} \cdot \vec{E}_{\text{clas}}(t) \Re \left[ e^{i(E_a-E_b)t} G(t) \right] \]

(7a)
and
\[
\frac{d}{dt} \rho_{ab}(t) = \left( -2i(\omega_0 + \Omega_+) - \gamma - 2\mu_{ab} \cdot \vec{E}_{\text{clas}}(t) G(t) e^{i(E_a - E_b)t} \right) \rho_{ab}(t) \\
+ 2\mu_{ab} \cdot \vec{E}_{\text{clas}}(t) G^*(t) e^{-i(E_a - E_b)t} \rho_{ba}(t) + i\mu_{ab} \cdot \vec{E}_{\text{clas}}(t)(1 - 2\rho_{aa}(0)).
\]

(7b)

The expression of the function \( G(t) \) is: \( G(t) \equiv \int_0^t dt' \mu_{ab} \cdot \vec{E}_{\text{clas}}(t') e^{-i(E_a - E_b)t'} \). We see from eq.(7b) that the real and imaginary parts of \( \rho_{ab}(t) \) are coupled due to non-linear effects in the classic external electromagnetic field. From eqs.(7) on, we will be using natural units \( (\hbar = c = 1) \).

In the same manner as in references [23, 29], in order to study the \( T \to \infty \) limit of eqs.(7) we apply the transformation
\[
\tilde{\rho}_{ij}(t) \equiv e^{i(E_i - E_j)t} \rho_{ij}(t), \quad i, j = a, b
\]

(8)

and change the time scale to \( s = t/T; \) in the limit \( T \to \infty, \) one obtains
\[
\frac{d}{ds} \tilde{\rho}_{aa}(s) \approx T \left[ -2\gamma - 2\mu_{ab} \cdot \vec{e} \mathcal{E}_0 \text{Re}(\tilde{G}(s)e^{-i(\phi(s) - \Delta Ts)}) \right] \tilde{\rho}_{aa}(s) \\
+ T\tilde{\rho}_{ab} \cdot \vec{e} \mathcal{E}_0 \left[ \text{Im}(e^{i(\phi(s) - \Delta Ts)} \tilde{\rho}_{ab}(0)) + \text{Re}(\tilde{G}(s)e^{-i(\phi(s) - \Delta Ts)}) \right]
\]

(9a)

and
\[
\frac{d}{ds} \tilde{\rho}_{ab}(s) \approx - \left( 2i\Omega_+ + \gamma + \mu_{ab} \cdot \vec{e} \mathcal{E}_0 \tilde{G}(s)e^{-i(\phi(s) - \Delta Ts)} \right) T \tilde{\rho}_{ab}(s) \\
+ \mu_{ab} \cdot \vec{e} \mathcal{E}_0 \tilde{G}^*(s)e^{-i(\phi(s) - \Delta Ts)} T \tilde{\rho}_{ba}(s) + iT\mu_{ab} \cdot \vec{e} \frac{\mathcal{E}_0}{2}(1 - 2\tilde{\rho}_{aa}(0)) e^{-i(\phi(s) - \Delta Ts)} .
\]

(9b)

As we write eqs.(9), we are assuming that \( \Delta \equiv 2\omega_0 - \nu \sim \frac{2\pi}{T} \). The terms proportional to \( e^{\pm i((2\omega_0 + \nu)T s + \phi(s))} \) do not contribute in the limit \( T \to \infty \) (as shown in references [23, 29]) and the terms proportional to \( e^{\pm i(\Delta Ts - \phi(s))} \) contribute to the dynamics of the density matrix in the resonance region when \( \Delta \sim \frac{2\pi}{T} \). The function \( \tilde{G}(s) \) appearing in eqs.(9) is given by \( \tilde{G}(s) = \mu_{ab} \cdot \vec{e} \frac{\mathcal{E}_0 T}{2} \int_0^T ds e^{i(\phi(s') - \Delta Ts')} \).

Equation (9a) gives us the probability that the atomic electron be at state \( |a\rangle \). Its solution after one period \( T \) is
\[ \rho_{aa}(T) = \left\{ \rho_{aa}(0) + i \int_0^T dt' e^{i\left(\frac{dG(t')}{dt'} - \frac{dG(t')}{dt} \rho_{ba}(0) - \frac{dG(t')}{dt} \rho_{ab}(0)\right)} e^{2(\gamma T + |\bar{G}(t')|^2)} \right\}^I_J + \frac{1}{2} \int_0^T dt' e^{2\gamma t'} \left\{ e^{2|\bar{G}(t')|^2} \right\}, \]

Since the function \( \bar{G}(t) \) is an explicit time-dependent function, the condition \( i \) in eq.(1) is not satisfied and none of the integrals in eq.(10) is a time-independent integral. We have an overall exponential decay (imaginary factor), but from the definition of the constant \( \gamma \) (see eq.(6a)), we obtain that the exponential \( e^{-2\gamma t} \) is not a geometric (path-dependent) imaginary phase, but a time-dependent one as well as the contribution to the decreasing exponential coming from \( |\bar{G}(T)|^2 \).

For the sake of completeness, we should also examine the solution of eq.(9b) after a period \( T \). Differently from reference [29], here the dynamical equation of \( \rho_{ab}(t) \) couples its real and imaginary parts under the regime of strong classic external electric field. We get a \( SU(2) \) structure for the solution of \( \rho_{ab}(t) \). Calling \( \rho_1(t) = \rho_{ab}(t) \) and \( \rho_2(t) = \rho_{ba}(t) \), the solution of eq.(9b) is

\[ \rho_I(T) = \left[ T(e^{\int_0^T dt' \vec{B}(t') \cdot \vec{\sigma}}) \right]_{IJ} \left\{ \int_0^T dt' \left[ T(e^{\int_0^{t'} dt'' \vec{B}(t'') \cdot \vec{\sigma}}) \right]_{JK} \times \right. \]

\[ \left. e^{(\gamma t + \text{Re}^2(\bar{G}(t')) - \text{Im}^2(\bar{G}(t')))} d_K(t') + \rho_J(0) \right\} e^{-i^{(2\omega_0T)} e^{-(\gamma T + \text{Re}^2(\bar{G}(T)) - \text{Im}^2(\bar{G}(T))} \right), (11) \]

with \( I = 1, 2 \) and \( J, K = 1, 2 \). In these two last indices we are using the implicit sum notation on the r.h.s. of eq.(11). The symbol \( T \) means the time-ordering integrals[30], and \( \vec{\sigma} \) are the Pauli matrices. The elements of the column \( d_J(t) \) are: \( d_1(t) = d_2(t) = i \vec{\mu}_{ab} \cdot \vec{e}^{\frac{\Delta t}{2}[1 - 2 \rho_{aa}(0)]} e^{-i(\phi(t') - \Delta t)} \). The components of the vector \( \vec{B} \) are: \( B_x(t) = \frac{d|\bar{G}(t)|^2}{dt} \), \( B_y(t) = 2 \left[ \text{Im}(\bar{G}(t)) \frac{d(\text{Re}(\bar{G}(t)))}{dt} - \text{Re}(\bar{G}(t)) \frac{d(\text{Im}(\bar{G}(t)))}{dt} \right] \) and \( B_z(t) = -2i \left[ \Omega_+ + \frac{d}{dt} \left[ \text{Re}(\bar{G}(t)) \text{Im}(\bar{G}(t)) \right] \right] \).

We have again the function \( \bar{G}(t) \) that is explicitly time-dependent and then the time-ordering integrals cannot be converted to path-ordering integrals. Therefore the integrals and the phases on the r.h.s. of eq.(11) are all time-dependent, as well as the functions with the overall decreasing exponential.

In reference [19] Garrison and Wright present in eq.(4.9) the value of the geometric imaginary phase in the limit of weak electric field and \( \gamma_b \gg \gamma_a \). In the weak electric field
limit, the \( SU(2) \) structure in eq.(11) disappears and the solutions of eqs.(10) and (11) become simpler,

\[
\rho_{aa}(t) = \left\{ \rho_{aa}(0) + i \int_0^t dt' \frac{\mathcal{E}}{2} \mu_{ab} \cdot e \left[ e^{i(\Delta t' - \phi(t'))} \rho_{ba}(0) - e^{-i(\Delta t' - \phi(t'))} \rho_{ab}(0) \right] e^{2\gamma t'} \right\} e^{-2\gamma t} 
\]

(12a) and

\[
\rho_{ab}(t) = \left\{ \rho_{ab}(0) + i \int_0^t dt' \frac{\mathcal{E}}{2} \mu_{ab} \cdot e \left( 1 - 2\rho_{aa}(0) \right) e^{i(\Delta t' - \phi(t'))} e^{(2i\Omega + \gamma)t'} \right\} e^{-(2i(\Omega + \omega_0) + \gamma)t} .
\]

(12b)

Equations (12) have terms with exponential decay (imaginary factors) but they are time-dependent, as was obtained in references [23, 29], in opposition to the results of reference [19] (see eq.(2c)).

Strictly speaking, geometric phases appear in adiabatic processes that happen only in the limit \( T \to \infty \). However, this limit is experimentally implemented by taking \( T \gg 2\pi/\omega \). The experimentally measured path-dependent phases are pretty much independent of the particular value of \( T \), once the previous inequality is true. In eq.(4.3) of reference [19] we have two imaginary phases: one of them \( e^{-2Im(\beta)} \) does not depend on the value of \( T \) and consequently is path-dependent, whereas the other one depends on the particular value of \( T \). In our results (eqs.(10) and (11)), valid for arbitrary intensities of the classical external electric field, all the imaginary phases depend on the chosen value of \( T \). It means that the greater the period \( T \) of \( \phi(t) \), the smaller the probability that the electron be in the upper level; also, the smaller the correlation \( \rho_{ab}(t) \).

In summary, we study the nature of the imaginary phase acquired by the probability of a two-level model coupled to an external electric field with fixed frequency due to its interaction with the electromagnetic vacuum. The interaction among the quantum system (two-level system), the environment (vacuum) and the external classical electric field is taken in the electric dipole aproximation and in the RWA. Garrison and Wright in reference [19] discussed this same model using the non-hermitian Schrödinger equation approach, concluding that the probability has an imaginary phase that is path-dependent. In this brief report we apply the matrix density formalism to study its limit of \( T \to \infty \) and from eqs.(10) and (11) we conclude that all imaginary phases are time-dependent and consequently depend on the chosen value of \( T \). We are considering the situation when
the effects due to dissipation and time variation of Hamiltonian are of the same order and consequently the coherence of initial states are not preserved along the whole period $T$. It is not surprising that those distinct approaches (master equation and non-hermitian Hamiltonian) give different results.

Finally we showed in the present work that even thought the non-hermitian Hamiltonian method has been a very important and useful tool in describing open systems, it should not be applied to discuss the nature (time-dependent or path-dependent) of imaginary phases.

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References


[30] We define the time-ordering operator as

\[ T \left( e^{\int_t^{\tau} A(t') dt'} \right) = 1 + \int_t^{\tau} A(t_1) dt_1 + \int_t^{\tau} A(t_1) \int_t^{t_1} A(t_2) dt_2 + \]
\[ + \int_t^{\tau} A(t_1) \int_t^{t_1} A(t_2) \int_t^{t_2} A(t_3) dt_3 + \cdots \]

where \( \tau \) can be smaller or bigger than \( t \).