Fast and robust spin manipulation in a quantum dot by electric fields

Yue Ban,¹ Xi Chen,¹, ² E. Ya Sherman,¹, ³ and J. G. Muga¹, ²

¹Departamento de Química-Física, UPV/EHU, Apdo 644, 48080 Bilbao, Spain
²Department of Physics, Shanghai University, 200444 Shanghai, People’s Republic of China
³IKERBASQUE, Basque Foundation for Science, 48011 Bilbao, Spain

We apply an invariant-based inverse engineering method to control by time-dependent electric fields electron spin dynamics in a quantum dot with spin-orbit coupling in a weak magnetic field. The designed electric fields provide a shortcut to adiabatic processes that flips the spin rapidly, thus avoiding decoherence effects. This approach, being robust with respect to the device-dependent noise, can open new possibilities for the spin-based quantum information processing.

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Introduction.- Coherent spin manipulation in quantum dots (QDs) [1, 2] is the key element in the state-of-the-art spintronics and solid-state quantum information [10, 11]. Accurate spin manipulation can be achieved by several techniques. One of them is the conventional electron spin resonance (ESR) induced by magnetic field oscillating at the Zeeman transition frequency [1]. A more robust technique is the spin manipulation with geometric Berry phases during the adiabatic motion [2, 3]. Nowadays, there is also a growing interest in the electric control of spin using spin-orbit (SO) coupling [12]. It has been applied to high-fidelity spin manipulation on the 100 ns time scale [7, 8]. This highly efficient all-electrical method has other advantages. For example, it is easy to generate time-dependent electric fields on the nanoscale by adding local electrodes and produce spin manipulation by making them Zeeman-resonant [7]. As a result, Rabi spin oscillations appear at a frequency much smaller than the Zeeman frequency making the flip relatively slow and prone to decoherence. We shall propose here another all-electrical technique to flip spin with high fidelity via “shortcuts to adiabaticity”, in a time that can be much shorter than any decoherence time.

Recently, several shortcuts to adiabaticity have been put forward to speed up the adiabatic passage of quantum systems, and achieve a robust and fast adiabatic-like control [12–20]. The transitionless or counter-diabatic control algorithms proposed by Demirplak, Rice [13], and Berry [14], are designed to add supplementary time-dependent interactions that cancel the diabatic couplings of a reference process. The system then follows exactly the adiabatic trajectory of the original unperturbed process, in principle in an arbitrarily short time. Transitionless quantum drivings have been implemented in two-level systems: spins in a magnetic field [14], atoms [15], and Bose-Einstein condensates in optical lattices [16]. A different shortcut is provided by inverse engineering the transient Hamiltonian [17–18] using Lewis-Riesenfeld invariants [21]. This method has been used for time-dependent traps [17, 21], atomic transport [22], and other applications [23, 24]. Although these two methods are potentially equivalent [25], their implementations and results can be quite different. Here we choose the invariant-based inverse engineering approach, since it is better suited than the transitionless driving to be produced by the desired all-electrical means.

Model.- We consider the electric control of electron spin in a QD formed in the $x$-$y$ plane of a two-dimensional electron gas confined in the $z$-direction by the coordinate-dependent material composition, under a weak magnetic field $B_0 \parallel z$, as shown in Fig. 1. Here the total Hamiltonian $H$ of the electron interacting with the external electric field $\mathcal{E}(t) = -\partial A/\partial t$ is $H = H_0 + H_{\text{so}} + H_{\text{int}}$, with [12]

$$H_0 = \frac{p_x^2 + p_y^2}{2m} + U(x, y) + \frac{\Delta_z}{2} \sigma_z, \quad (1)$$

$$H_{\text{so}} = (-\alpha \sigma_y + \beta \sigma_z)p_x + \alpha \sigma_x p_y, \quad (2)$$

$$H_{\text{int}} = -\frac{e}{c} A(t) \cdot \mathbf{v}, \quad (3)$$

where $m$ is the electron effective mass and $\sigma_i \ (i = x, y, z)$ are the Pauli matrices. $H_0$ represents the kinetic energy,
the potential \( U(x, y) \), and the Zeeman \( \Delta_z = g \mu_B B_0 \), where \( \mu_B \) is the Bohr magneton, and \( g \) is the Landé factor. The eigenfunctions of \( H_0 \) are \( \psi_j(x, y)|\sigma\rangle \), where \( |\sigma\rangle = |\pm 1\rangle \) is the eigenstate of \( \sigma_z \), and the spectrum is given by \( E_j \pm \Delta_z / 2 \), where \( E_j \) are the orbital eigenenergies in the confinement potential.

The SO coupling is the sum of structure-related Rashba \((\alpha)\) and bulk-originated Dresselhaus \((\beta)\) terms for \([110]\) growth axis. The vector potential \( \mathbf{A}(t) \) is in the \((x, y)\)-plane and corresponding spin-dependent velocity operators are

\[
v_x = \frac{i}{\hbar} [H_0 + H_{so}, x] = p_x + m/\hbar \sigma_x - \alpha \sigma_y, \quad (4)
v_y = \frac{i}{\hbar} [H_0 + H_{so}, y] = p_y + m/\hbar \sigma_x. \quad (5)
\]

We focus on the doublet \( \Psi_1 = \psi_1|1\rangle, \Psi_2 = \psi_1|-1\rangle \), include the higher orbitals by Löwdin partition \([28, 29]\), and reduce the full Hamiltonian into an effective \( 2 \times 2 \) matrix,

\[
H_{\text{eff}} = \frac{g \mu_B}{2} \left( \begin{array}{cc}
Z & X + iY \\
-X - iY & -Z
\end{array} \right), \quad (6)
\]

where \( X = B_2(1 + \xi_y), Y = (\alpha/\beta)(1 + \xi_x)B_1, Z = B_0 + (1 + \xi_z)B_1 \), with the effects of higher states characterized by \( \xi_x \) and \( \xi_y \), and the components of the designed electric field are renormalized by the factors of \( 1/(1 + \xi_z) \). The effective magnetic fields are expressed with the SO coupling parameters as \( B_1 = -2e\alpha A_y / cg \mu_B \), and \( B_2 = -2e\alpha A_y / cg \mu_B \). The resulting electric fields are:

\[
E_x(t) = \frac{g \mu_B}{2c \beta} \frac{\partial B_1}{\partial t}, \quad E_y(t) = \frac{g \mu_B}{2c \alpha} \frac{\partial B_2}{\partial t}. \quad (7)
\]

In practice, some slowly varying electric fields can be applied to drive the state from \( \Psi_1 \) to \( \Psi_2 \) adiabatically along an instantaneous eigenstate of the Hamiltonian in Eqs. \([11]\). To accelerate the driving using the transitionless algorithm, counter-diabatic fields should be provided \([22]\).

However, the common dependence of \( Y \) and \( Z \) on \( B_1 \) precludes the implementation of the fast driving terms only by electric fields. In contrast, invariant-based inverse engineering naturally leads to an all-electrical driving.

**Dynamical invariant and spin-flip example.** We shall design the time dependence of the external electric fields to guarantee the state transfer in some fixed time \( t_f \) by using the dynamical \( 2 \times 2 \) invariant \( I(t) \) satisfying the condition \( dI(t)/dt = \iota \hbar \partial I(t)/\partial t - [H_{\text{eff}}(t), I(t)] = 0 \). Parametrizing the Bloch sphere (Fig. \(1\), by the angles \( \theta \) and \( \varphi \), we construct yet unknown orthogonal eigenstates \( |\chi_{\pm}(t)\rangle \) of \( I(t) \) as

\[
|\chi_{+}(t)\rangle = \left( \begin{array}{c}
\cos \theta e^{i \varphi} \\
\sin \theta/2
\end{array} \right), |\chi_{-}(t)\rangle = \left( \begin{array}{c}
\sin \theta/2 \\
-\cos \theta e^{-i \varphi}
\end{array} \right). \quad (8)
\]

They satisfy \( I(t)|\chi_{\pm}(t)\rangle = \lambda_{\pm} |\chi_{\pm}(t)\rangle \). Introducing \( \lambda_{\pm} = \pm g \mu_B B_c / 2 \), we construct the invariant as \([22]\)

\[
I(t) = \frac{g \mu_B}{2} B_c \left( \begin{array}{cc}
\cos \theta & \sin \theta e^{i \varphi} \\
-\sin \theta e^{-i \varphi} & -\cos \theta
\end{array} \right), \quad (9)
\]

where \( B_c \) is an arbitrary constant magnetic field to keep \( I(t) \) with units of energy. According to the Lewis-Riesenfeld theory, the solution of the Schrödinger equation, \( i \hbar \partial \Psi = H_{\text{eff}}(t)|\Psi\rangle \), is a superposition of orthonormal “dynamical modes”, \( \Psi(t) = \sum_n C_n e^{i \omega_n(t)} |\chi_n(t)\rangle \quad ([27]) \), where \( C_n \) are time-independent amplitudes and \( \alpha_n(t) \) is Lewis-Riesenfeld phase,

\[
\alpha_n(t) = \frac{1}{\hbar} \int_0^t \langle \chi_n(t')| \hbar \partial/\partial t' - H_{\text{eff}}(t') |\chi_n(t')\rangle dt'. \quad (10)
\]

From the invariant condition, \( dI(t)/dt = 0 \), the angles \( \theta \) and \( \varphi \) are related to \( X, Y \) and \( Z \) by auxiliary equations

\[
\dot{\theta} = \eta (X \sin \varphi - Y \cos \varphi), \quad (11)
\]

\[
\dot{\varphi} = \eta (X \cos \varphi \cot \theta + Y \sin \varphi \cot \theta - Z), \quad (12)
\]

where \( \eta = g \mu_B / \hbar \). Since \( X \) is a function of \( B_2 \), while \( Y \) and \( Z \) are functions of \( B_1 \), once \( \theta \) and \( \varphi \) are fixed, Eqs. \([11]\) and \([12]\) give the effective magnetic fields

\[
B_1 = \frac{-\beta \dot{\theta} \cot \theta \cos \varphi + \beta (\dot{\varphi} + \eta B_0) \sin \varphi}{\eta(1 + \xi_\beta)(\alpha \cot \theta - \beta \sin \varphi)}, \quad (13)
\]

\[
B_2 = \frac{\alpha \dot{\varphi} \cos \varphi + \alpha (\dot{\varphi} + \eta B_0) \cos \varphi - \beta \dot{\theta}}{\eta(1 + \xi_\beta)(\alpha \cot \theta - \beta \sin \varphi)}, \quad (14)
\]

from which the electric fields are calculated using Eq. \([7]\). During the spin-flip process, there exist some time instants \( t = t_s \) which satisfy

\[
\alpha \cot \theta(t_s) = \beta \sin \varphi(t_s), \quad (15)
\]

and make the denominators of \( B_1 \) and \( B_2 \) zero. To get rid of such singularities we impose the conditions

\[
\beta \sin \varphi(t_s) \left[ \dot{\varphi}(t_s) + \eta B_0 - (\beta/\alpha) \dot{\theta}(t_s) \cos \varphi(t_s) \right] = 0, \quad (16)
\]

\[
\alpha \cos \varphi(t_s) \left[ \dot{\varphi}(t_s) + \eta B_0 - (\beta/\alpha) \dot{\theta}(t_s) \cos \varphi(t_s) \right] = 0, \quad (17)
\]

which make the numerators of \( B_1 \) and \( B_2 \) zero simultaneously. In the following example, we will show how this works.

In general, the eigenstates of the invariant are not the same as the instantaneous eigenstates of the Hamiltonian, since \( I(t) \) and \( H_{\text{eff}}(t) \) do not commute. If we impose for \( \theta \) at \( t = 0 \) and \( t_f \) that

\[
\theta(0) = 0, \quad \theta(t_f) = \pi, \quad \dot{\theta}(0) = 0, \quad \dot{\theta}(t_f) = 0, \quad (18)
\]

then \( [H_{\text{eff}}(0), I(0)] = 0 \) and \( [H_{\text{eff}}(t_f), I(t_f)] = 0 \), which guarantees common eigenstates at initial and final times. Moreover the state obeying Eq. \([13]\) will flip from
\[ \psi(t_f) = \pi/2, \quad \varphi(t_f) = \pi/2. \tag{19} \]

These conditions are not sufficient though, since we still have to deal with the singularities and their cancellation. As \( \cot \theta(t_f/2) = 0 \), we may satisfy Eq. \( \text{(15)} \) and impose zeros of the denominators for \( B_{1,2} \) at \( t_s = t_f/2 \), if \( \sin \varphi(t_f/2) = 0 \). Imposing the two conditions

\[ \varphi(t_f/2) = 0, \tag{20} \]

\[ \dot{\varphi}(t_f/2) = (\beta/\alpha)\dot{\theta}(t_f/2) - \eta B_0, \tag{21} \]

to satisfy Eqs. \( \text{(10)} \) and \( \text{(17)} \) at \( t_s = t_f/2 \), we cancel the singularity there. With the conditions in Eqs. \( \text{(19)}-\text{(21)} \), we solve the third order polynomial ansatz \( \varphi = \sum_{j=0}^{3} b_j t^j \) to determine \( \varphi(t) \).

Explicit calculations demonstrate that for the third-order polynomial ansatz and boundary conditions imposed here, there is only one (removable) singularity at \( t_s = t_f/2 \) when the field \( B_0 \) is smaller than certain upper limit \( B_0^{\text{max}} \), shown in Fig. \( 2 \) as a function of \( t_f \). For \( B_0 > B_0^{\text{max}} \), more solutions of \( \text{(15)} \) appear [\( B_0 \) and \( \sin \varphi(t) \) are coupled by Eq. \( \text{(21)} \), which cannot be canceled with the third order polynomial. To satisfy Eqs. \( \text{(10)} \) and \( \text{(17)} \) at more than one zero of the denominators of \( B_{1,2} \), one may set higher order polynomials for \( \varphi \) and further conditions. This increases the bound \( B_0^{\text{max}} \), but also complicates the driving fields. In the present examples, we just apply the third order polynomial ansatz with the boundary conditions in Eqs. \( \text{(18)}-\text{(21)} \), so that the applied magnetic field \( B_0 \) should not go beyond the upper limit in Fig. \( 2 \). As the upper limit field grows for smaller times \( t_f \), this is not a problem in practice. Figure \( 3 \) shows examples of spin flip for different values of \( B_0 \).

With the functions \( \theta \) and \( \varphi \) fixed [see Fig. \( 3 \) (a)], the designed electric fields, \( \mathcal{E}_x(t) \) and \( \mathcal{E}_y(t) \), corresponding to \( B_0 = 0.01 \text{ T} \) and \( B_0 = 0.07 \text{ T} \) (close to the upper limit), are depicted in Fig. \( 3 \) (b) and (c). The populations (not shown) of the two spin states, given by \( P_1 = \cos^2(\theta/2) \) and \( P_{-1} = \sin^2(\theta/2) \), cross each other smoothly as \( \theta \) goes from \( 0 \) to \( \pi \). The choice of \( B_0 \) determines the trajectory on the Bloch sphere for a given \( t_f \). When \( B_0 \) approaches
the upper limit, the electric fields exhibit sharp peaks, see Fig. 3 (c). The smooth time-dependence in Fig. 3 (b) is well suited for the applications, while the complicated dependence in Fig. 3 (c) should be avoided. Undesirable excitation of the orbital modes does not occur here since the spin flip $t_f \sim 1$ ns, while the energy split of the orbital states in typical QDs exceeds 0.1 meV. Therefore, regarding the orbital motion, our perturbation is strongly adiabatic, and no orbital excitation occurs.

To check the stability with respect to initialization errors, we assume now the initial state as $(\sqrt{T - e^{i\phi_0}}, \sqrt{\epsilon})^T$ with an arbitrary phase $\phi_0$, and find $\theta$ and $\varphi$ from Eqs. (11) and (12) for the same designed electric fields (Fig. 4). The final value $\theta(t_f)$ depends on the error $\epsilon$, but insensitive to the initial phase $\phi_0$, while $\varphi(t_f)$ is sensitive to both initial conditions. Since our goal is to realize the spin flip, the final $\varphi$ is irrelevant, and the experimental effort should focus on achieving a small error $\epsilon$.

*Decoherence and noise effects.* To show feasibility of our approach, we study the effects of noise and decoherence on the spin-flip fidelity. We begin with a generic approach for coupling to the incoherent environment, based on the conventional Lindblad formalism as can arise, e.g., from interaction with the conduction electron bath. The master equation reads

$$\dot{\rho} = -\frac{i}{\hbar}[H_{\text{eff}}, \rho] - \frac{\gamma}{2} [\{\sigma_x, [\sigma_x, \rho]\} + [\sigma_y, [\sigma_y, \rho]] + [\sigma_z, [\sigma_z, \rho]]].$$

(22)

FIG. 4: (Color online) Time evolution of $\cos \theta$ (a) and $\sin \varphi$ (b) for the same Hamiltonian with the designed electric fields and $B_0 = 0.01$ T, where $\epsilon = 0$, $\varphi_0 = \pi/2$ (solid blue); $\epsilon = 0.01$, $\varphi_0 = \pi/4$ (dashed black). Other parameters are the same as in Fig. 3.

FIG. 5: (Color online) Fidelity as a function of $\gamma$ for different times $t_f = 0.1$ ns (solid red) and $t_f = 1$ ns (dotted black). The fidelity estimated from perturbation theory is also compared, for $t_f = 0.1$ ns (dashed blue, undistinguished) and $t_f = 1$ ns (dot-dashed orange). $B_0 = 0.01$ T and other parameters are the same as in Fig. 2.

where $\gamma$ is the dephasing rate. We introduce the Bloch vector with components $u = \rho_{11} - \rho_{10}$, $v = -i(\rho_{10} - \rho_{01})$, and $w = \rho_{00} - \rho_{11}$, and obtain from Eq. (23)

$$\begin{pmatrix} \dot{u} \\ \dot{v} \\ \dot{w} \end{pmatrix} = \begin{pmatrix} -4\gamma & \eta Z & -\eta Y \\ -\eta Z & -4\gamma & \eta X \\ -\eta Y & -\eta X & -4\gamma \end{pmatrix} \begin{pmatrix} u \\ v \\ w \end{pmatrix}.$$ 

(23)

We solve Eq. (23) numerically and calculate fidelity $F = |\langle -1|\Psi(t_f)\rangle|$, see Fig. 5. For $\gamma t_f \ll 1$ the time-dependent perturbation theory [21] yields the bound $F \gtrsim 1 - 2\gamma t_f$. Since the induced flip occurs very fast, it can overcome the main danger for the low-temperature spin manipulation in QDs coming from the hyperfine coupling to the nuclear spins, where the decoherence times exceed 100 ns [7].

Another source of decoherence is the device-dependent noise in the electric field acting on the spin. This can be important when the relatively weak electric fields are applied. We analyze in detail the effect of this noise and find that our method is robust to this randomness in [32].

*Conclusion and outlook.* We have proposed a fast and robust method to flip electron spin in a QD with SO coupling and weak perpendicular magnetic field. The spin-flip process, designed by Lewis-Riesenfeld invariants, is faster than the decoherence for any known low-temperature dephasing mechanism. This method can be further complemented by optimal control theory for time-and energy-minimization subjected to different physical constraints [33]. Implementation of this technique will allow for high-fidelity spin-manipulation for quantum information processing.

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I. Hamiltonian reduction

Here we give the details on the derivation of the effective $2 \times 2$ Hamiltonian in the paper. For simplicity, we consider a 4-level system. The total Hamiltonian is $H = H_0 + H_{so} + H_{int}$. In the basis of two lowest spin-split orbital states,

$$H_0 = \begin{pmatrix} E_1 + \Delta_z/2 & 0 & 0 & 0 \\ 0 & E_1 - \Delta_z/2 & 0 & 0 \\ 0 & 0 & E_2 + \Delta_z/2 & 0 \\ 0 & 0 & 0 & E_2 - \Delta_z/2 \end{pmatrix},$$

where $\Delta_z$ is the Zeeman splitting, and

$$H_{so} + H_{int} = \begin{pmatrix} -\frac{e}{c} \beta A_x & -\frac{e}{c} \alpha (iA_x + A_y) & (\beta - \tilde{A}_x)\bar{p}_x - \tilde{A}_y\bar{p}_y & \alpha (\bar{p}_x + \bar{p}_y) \\ \frac{e}{c} \alpha (iA_x - A_y) & -\frac{e}{c} \beta A_x & \alpha (\bar{p}_x - \bar{p}_y) & -\alpha (\bar{p}_x + \bar{p}_y) \\ -(\beta + \tilde{A}_x)\bar{p}_x + \tilde{A}_y\bar{p}_y & -\alpha (\bar{p}_x - \bar{p}_y) & \frac{e}{c} \beta A_x & -\frac{e}{c} \alpha (iA_x + A_y) \\ \frac{e}{c} (iA_x - A_y) & (\beta + \tilde{A}_x)\bar{p}_x + \tilde{A}_y\bar{p}_y & -\frac{e}{c} \beta A_x & \frac{e}{c} \alpha (iA_x - A_y) \end{pmatrix},$$

with $\bar{p}_1 = \langle \psi_1 | \bar{p}_1 | \psi_2 \rangle$ and $\tilde{A}_i \equiv eA_i/mc$.

Our aim is to flip the spin between two levels taking into account the effects from orbital motion and the influence of the other levels. As the energy gap of the orbital states is much larger than the Zeeman splitting $\Delta_z$, the Löwdin partition \[28, 29\] will enable us to obtain an effective $2 \times 2$ matrix Hamiltonian. We split the total Hamiltonian $H$ into four blocks $\hat{Q}, \hat{B}, \hat{C}, \hat{C}^\dagger$, each of which is a $2 \times 2$ matrix,

$$H = \begin{pmatrix} \hat{Q} & \hat{C} \\ \hat{C}^\dagger & \hat{B} \end{pmatrix}. \tag{26}$$

The time-independent Schrödinger equation can be formally written as

$$\hat{Q}\bar{F} + \hat{C}\bar{G} = E\bar{F},$$
$$\hat{C}^\dagger\bar{F} + \hat{B}\bar{G} = E\bar{G}. \tag{27}$$

Substituting the formal solution of Eq. \[28\], $G = (E - B)^{-1}C^\dagger F$, into Eq. \[27\], we get a closed equation for $F$,

$$\hat{Q}\bar{F} + \hat{C}(E - \tilde{B})^{-1}C^\dagger \bar{F} = E\bar{F}. \tag{29}$$

Therefore, the effective Hamiltonian for this subspace is given by

$$H_{\text{eff}} \rightarrow \hat{Q} + \hat{C}(E - \tilde{B})^{-1}C^\dagger$$

including the effects from $\hat{B}, \hat{C}$ and $\hat{C}^\dagger$.

Assuming that the electric field is not strong enough to excite other orbital states, namely, $e \max(|A_x\bar{p}_x|, |A_y\bar{p}_y|)/mc \ll E_2 - E_1$, and keeping the first order term of the spin-orbit coupling constant, the effective Hamiltonian becomes

$$H_{\text{eff}} = \begin{pmatrix} E_1 + h_{11}^0 + h_{12}^0 & h_{12}^0 + h_{22}^0 & h_{11}^0 + h_{12}^0 & h_{22}^0 + h_{22}^0 \\ h_{21}^0 + h_{21}^0 & E_2 - E_1 & h_{21}^0 + h_{21}^0 & h_{22}^0 + h_{22}^0 \end{pmatrix},$$

where

$$h_{11}^0 = -h_{22}^0 = -\frac{e}{c} \beta A_x,$$
$$h_{12}^0 = h_{21}^0 = -\frac{e}{c} \alpha (iA_x + A_y),$$
$$h_{11} = -h_{22} = 2\frac{e\beta \bar{p}_x (A_x\bar{p}_x + A_y\bar{p}_y)}{mc (E_2 - E_1)},$$
$$h_{12} = h_{21} = 2\frac{e\alpha (A_x\bar{p}_x + A_y\bar{p}_y)}{mc (E_2 - E_1)}. \tag{30}$$

By a simple shift we may ignore the common term $E_1$ on the diagonal, and finally express the Hamiltonian as

$$H_{\text{eff}} = \frac{g\mu_B}{2} \begin{pmatrix} Z & X + iY \\ X - iY & -Z \end{pmatrix},$$

where $X = B_2(1 + \xi_y), Y = (\alpha/\beta)(1 + \xi_x)B_1, Z = B_0 + (1 + \xi_x)B_1$ with $B_1 = -2e\beta A_x/(eg \mu_B), B_2 = -2e\beta A_y/(eg \mu_B)$. According to perturbation theory, the effects from other states can be characterized by $\xi_x$ and $\xi_y, \xi_i = 2\sum_{n>1} |\langle \psi_1 | p_i | \psi_n \rangle|^2 / |m(E_n - E_1)|$, summing up all contributions.

II. Apparatus Noise Effect

Besides the decoherence resulting from the interaction between the system and the environment introduced in the main text, we consider apparatus noise, i.e. the
FIG. 6: (Color online) Dependence of fidelity on $\lambda_0^2$, given by $t_f = 1$ ns (solid black) and $t_f = 0.1$ ns (dashed red). Other parameters are the same as in Fig. 2 (b) in the main text: $\hbar\alpha = 2 \times 10^{-6}$ meV·cm, $\hbar\beta = 10^{-6}$ meV·cm, $g = -0.44$ for GaAs, $B_0 = 0.01$ T, $\xi_x = \xi_y = 0$.

Hamiltonian $H_{\text{eff}}$ [Eq. (35)] perturbed by a stochastic part $H^{\text{noise}}$ describing noise from the electric field source. Therefore, the stochastic Schrödinger equation is

$$i\hbar\frac{d}{dt}\Psi(t) = (H_{\text{eff}} + H^{\text{noise}})\Psi(t),$$

(36)

where $H^{\text{noise}} = \lambda H'\zeta(t)$, $\langle \zeta(t) \rangle = 0$, $\langle \zeta(t)\zeta(t') \rangle = \delta(t-t')$ and $\lambda = \lambda_0\sqrt{t_f}$, where $\lambda_0$ is the strength of noise. $H'$ is the part of $H_{\text{eff}}$ contributing from the time-dependent electric field. Then the density matrix [34] is

$$\dot{\rho} = -\frac{i}{\hbar}[H_{\text{eff}}, \rho] - \frac{\lambda^2}{2\hbar^2}[H', [H', \rho]]$$

(37)

Following the same procedure to derive the Bloch equation in the main text after Eq. (22), we obtain the corresponding equation with source noise. In order to analyze the effects from randomness induced by the electric field, we calculate numerically the fidelity $F = |\langle -1 | \Psi(t_f) \rangle|$, see Fig. 0. The high fidelity shows that our method is robust to the source noise.