Phonon-induced relaxation and decoherence times of the hybrid qubit in silicon quantum dots

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We study theoretically the phonon-induced relaxation and decoherence processes in the hybrid qubit in silicon. A hybrid qubit behaves as a charge qubit when the detuning is close to zero and as a spin qubit for large detuning values. It is realized starting from an electrostatically defined double quantum dot where three electrons are confined and manipulated through only electrical tuning. By employing a three-level effective model for the qubit and describing the environment bath as a series of harmonic oscillators in the thermal equilibrium states, we extract the relaxation and decoherence times as a function of the bath spectral density and of the bath temperature using the Bloch-Redfield theory. For Si quantum dots the energy dispersion is strongly affected by the physics of the valley, i.e., the conduction band minima, so we also included the contribution of the valley excitations in our analysis. Our results offer fundamental information on the system decoherence properties when the unavoidable interaction with the environment is included and temperature effects are considered.

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I. INTRODUCTION

In solid-state physics, electrons or holes confined in 0dimensional nanostructures, i.e., quantum dots (QDs), represent promising platforms for the realization of qubits, for exploiting spin and/or charge, and for quantum computing applications [1–6]. However qubits are inevitably coupled to the degrees of freedom of the surrounding environment causing a loss of coherence that deeply affects the qubit operations [7,8]. Depending on the nature of the host materials, a source of noise could be predominant with respect to the others. For example, materials belonging to group IV, such as Si and Ge, possess isotopes with zero nuclear spin that allow them to reduce magnetic noise, while electrical noise remains an issue to be faced [9,10].

The dynamics of electron spin in a quantum dot is mainly affected by the interaction with two environments of different natures: the phonons in the lattice and the spins of atomic nuclei in the quantum dot. The spin-orbit interaction couples the spin and the orbital degrees of freedom that, being coupled to the phonons, provide an indirect coupling between the electron spin and the phonons. They constitute a large dissipative bosonic reservoir causing decoherence and relaxation. The short-time correlations in the phonon bath induce a Markovian dynamics of the qubit that is the subject of our study. Moreover the electron spin and the nuclear spins in the host material interact via the Fermi contact hyperfine interaction, which creates entanglement between them. It turns out that long-time correlations in the nuclear spin system induce a non-Markovian dynamics of the electron spin. However the presence of silicon isotopes with zero nuclear spin reduces

magnetic noise and non-Markovian dynamics becomes negligible.

Much progress has also been made in studying semiconducting QDs in III-V compounds, such as GaAs [11,12], that assures greater advantages in fabrication processes; however Si qubits have attracted recently a lot of attention also due to the immediate integrability with the existing Complementary Metal-Oxide Semiconductor (CMOS) technology of the microelectronic industry [13,14].

When Si QD-based qubits are considered, the sixfold degeneracy of the conduction band minima, which is due to the twofold degeneracy of the Δ valleys aligned along each one of the three main crystallographic directions, is an additional source of decoherence that may be overcome only if the typical qubit splitting energies are smaller with respect to the valley splittings [15,16]. Otherwise, in order to have a complete picture, it becomes indispensable to include valley effects in the Hamiltonian model.

The hybrid qubit (HQ) is realized by the electrostatic confinement of three electron spins in a double quantum dot [17,18]. We describe an HQ with an effective three-level model adopting a basis whose logical states are encoded in the S = 1/2 and $S_z = -1/2$ subspace, where S denotes the total angular momentum of the three electrons [19,20]. Then, we model the environment with which the HQ unavoidably interacts, by a bath consisting of a series of harmonic oscillators with frequencies ω_j . The effects of the bath temperature and of the bath spectral density on the qubit decoherence and relaxation times are studied.

In Ref. [21] the authors focus on ameliorating the dominant sources of decoherence in order to increase the coherence time in a Si/SiGe HQ. They measure $dE_Q/d\epsilon$, where E_Q is the qubit energy and ϵ is the detuning between the two QDs, and demonstrate that the HQ can be made resilient to charge noise by tuning appropriately the qubit parameters. More

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recently, in Ref. [22] atomic scale disorder at the quantum well interface has been put into direct connection with the dephasing of the HQ.

The study of the relaxation and decoherence processes is of fundamental as well as practical interest for quantum computation applications. For this reason, the aim of the present paper is to study theoretically the phonon-induced relaxation and decoherence times and how these times are affected by the HQ parameters as well as by the bath structure.

The paper is organized as follows. In Sec. II we present the theoretical model describing the Si HQ, including valley degeneracy, the bath, and their interaction; moreover the relaxation and decoherence times are derived following the Bloch-Redfield theory. Section III is devoted to the analysis of the relaxation and decoherence processes when the effects of the bath are included and the space of the qubit parameters is explored. Finally, concluding remarks are reported in Sec. IV.

II. THEORY

This section is devoted to the description, through an effective Hamiltonian model, of the HQ interacting with a bath of harmonic oscillators when temperature effects are included. The analytical expressions for the evaluation of relaxation and decoherence times are presented.

A. Model of the silicon hybrid qubit in a thermal bath

We describe effectively the HQ adopting a threedimensional basis. The first state of the basis corresponds to a configuration with two electrons in the left dot and one in the other and consequently has a singlet charge form. The remaining basis states, on the contrary, correspond to the complementary configuration in which one electron is confined in the left dot and two electrons are confined in the right dot; they correspond to singlet and triplet charge configurations. The three-level matrix describing the qubit is

$$H_{S} = \begin{pmatrix} \frac{\epsilon}{2} & \Delta_{1} & \Delta_{2} \\ \Delta_{1} & -\frac{\epsilon}{2} & 0 \\ \Delta_{2} & 0 & -\frac{\epsilon}{2} + \Delta_{R} \end{pmatrix}, \tag{1}$$

where ϵ is the detuning between the two QDs; Δ_1 and Δ_2 refer to the tunnel couplings between different charge states from one dot to the other; and Δ_R corresponds to the low-energy splitting of the right dot, which reflects a valley excitation, an orbital excitation, or a combination. The detuning ϵ can be changed by varying the applied electrostatic potential in the left QD. The tunnel couplings Δ_1 and Δ_2 can be electrostatically modulated by changing the tunneling barrier between the two QDs. The Δ_R parameter can be effectively manipulated by varying the quantum confining energy profile of the right dot, exploiting the external electric field coming from a gate(s) close to the quantum dot. In particular, in the case of a Si/SiO₂ confining interface, the valley splitting depends on the electric field at that interface [23] so an additional gate(s) (for example, a back gate in a silicon on insulator structure) can be used to effectively modulate the electric field at the interface and the resulting Δ_R without



FIG. 1. A sketch of the HQ energy levels. The interdot tunnel couplings are Δ_1 and Δ_2 , ϵ is the detuning between the two QDs, and Δ_R corresponds to the low-energy splitting of the right dot.

affecting the confining energy potential. The estimation of such parameters is extractable from simulations adopting a tight-binding model as done in Ref. [22] for a strained Si quantum well sandwiched between strain-relaxed $Si_{0.7}Ge_{0.3}$. An illustrative sketch of the theoretical model describing HQ is reported in Fig. 1.

We model the surrounding environment by a series of N harmonic oscillators with the Hamiltonian

$$H_B = \sum_{j=1}^{N} \omega_j b_j^{\dagger} b_j, \qquad (2)$$

where $b_j(b_j^{\dagger})$ is the annihilation (creation) operator for the environment mode and ω_j is the frequency associated with each mode *j*.

The interaction Hamiltonian between HQ and the bath is written as [24]

$$H_I = \sum_{j=1}^N \lambda_j (b_j^{\dagger} + b_j) \otimes \frac{1}{2} \hat{O}_S, \qquad (3)$$

with λ_j representing the coupling qubit-bath, and the system operator \hat{O}_S is equal to

$$\hat{O}_S = \begin{pmatrix} 1 & 0 & 0 \\ 0 & -1 & 0 \\ 0 & 0 & -1 \end{pmatrix}.$$
 (4)

The total Hamiltonian given by the sum of the three contributions, i.e., $H = H_S + H_B + H_I$, is transformed by adopting the unitary transformation $U = [\mathbf{m}_0, \mathbf{m}_1, \mathbf{m}_2]$. Each column of U contains the eigenvectors \mathbf{m}_k of H_S , in such a way that transforming H_S through U results in a diagonal form. After making explicit calculations for $\tilde{H} = U^{\dagger}HU$, we finally obtain

$$\tilde{H} = \begin{pmatrix} E_0 & 0 & 0\\ 0 & E_1 & 0\\ 0 & 0 & E_2 \end{pmatrix} + \frac{\zeta}{2} \otimes \begin{pmatrix} \chi_{00} & \chi_{01} & \chi_{02}\\ \chi_{10} & \chi_{11} & \chi_{12}\\ \chi_{20} & \chi_{21} & \chi_{22} \end{pmatrix} + \tilde{H}_B,$$
(5)

where E_i (i = 0, 1, and 2) are the eigenvalues of H_S , $\zeta = \sum_{j=1}^N \lambda_j (b_j^{\dagger} + b_j)$, and $\chi_{ij} = \chi_{ji}$ are the transformed matrix elements of \hat{O}_S through U and $\tilde{H}_B = H_B$.

B. Relaxation and decoherence times

We determine the explicit expressions for the relaxation and the decoherence times. First, we calculating the power spectrum $S_{\zeta}(\omega)$ for a bath in thermal equilibrium at temperature *T*. In the framework of the Linblad master equation describing the HQ, we trace over the environmental bath degrees of freedom, obtaining

$$S_{\zeta}(\omega) = \frac{1}{2\pi} \int_{-\infty}^{+\infty} \langle f(t)f(0) \rangle_{\beta} e^{i\omega t} dt.$$
 (6)

The correlator $\langle f(t)f(0)\rangle_{\beta}$, where $\beta \equiv (k_BT)^{-1}$, is evaluated analytically giving as a result

$$\begin{aligned} \langle f(t)f(0) \rangle_{\beta} &\equiv \operatorname{Tr}_{B}[e^{-\beta H_{B}}f(t)f(0)] \\ &= \operatorname{Tr}_{B}(e^{-\beta H_{B}}e^{iH_{B}t}\zeta e^{-iH_{B}t}\zeta) \\ &= \sum_{j=1}^{N} \lambda_{j}^{2} \bigg[\cos(\omega_{j}t) \coth\left(\frac{\beta\hbar\omega_{j}}{2}\right) - i\sin(\omega_{j}t) \bigg], \end{aligned}$$

$$\tag{7}$$

and the following relations have been exploited:

$$\langle b_{j}^{\dagger}b_{j}\rangle_{\beta} = \frac{1}{e^{\beta\omega} - 1} = \frac{1}{2}\coth\left(\frac{\beta\hbar\omega_{j}}{2}\right) - \frac{1}{2},$$
$$\langle b_{j}b_{j}^{\dagger}\rangle_{\beta} = \langle b_{j}^{\dagger}b_{j} + 1\rangle_{\beta} = \frac{1}{2}\coth\left(\frac{\beta\hbar\omega_{j}}{2}\right) + \frac{1}{2}.$$
 (8)

Inserting Eq. (7) into Eq. (6) we obtain

$$S_{\zeta}(\omega) = \frac{\sum_{j=1}^{N} \lambda_j^2}{2} \left\{ \delta(\omega + \omega_j) \left[\coth\left(\frac{\beta \hbar \omega_j}{2}\right) - 1 \right] + \delta(\omega - \omega_j) \left[\coth\left(\frac{\beta \hbar \omega_j}{2}\right) + 1 \right] \right\}.$$
(9)

In the hypothesis that *N* is large, the sum over *j* can be approximated by the frequency integral $\sum_{j=1}^{N} \lambda_j^2 \approx \int_0^{+\infty} J(\omega) d\omega$, where $J(\omega)$ is the spectral function of the oscillator bath, and Eq. (9) is rewritten in the following way:

$$S_{\zeta}(\omega) = \frac{1}{2}J(\omega) \left[\coth\left(\frac{\beta\hbar\omega}{2}\right) + 1 \right], \ \omega > 0, \tag{10}$$

$$S_{\zeta}(\omega) = \frac{1}{2}J(\omega) \left[\coth\left(\frac{\beta\hbar\omega}{2}\right) - 1 \right], \ \omega < 0.$$
(11)

The spectral density is supposed to be of the following general form: $J(\omega) = (\eta \omega^s / \omega_c^{s-1}) e^{-\omega/\omega_c}$, with a high-energy cutoff ω_c and where η is an effective dimensionless coupling that determines the overall strength of the electron-phonon coupling. The parameter *s* distinguishes among s = 1 ohmic, s > 1 superohmic, and s < 1 subohmic baths.

Following the theory [25,26], in which the Bloch-Redfield master equation has been used to describe the dynamics of the qubit interacting with the phonon bath, the relaxation time T_1 , the pure dephasing time T_{ϕ} , and the decoherence time T_2 are



FIG. 2. T_1 (red lines) and T_2 (blue lines) as a function of the bath temperature *T* for three different regimes: s = 1 (solid lines), s = 2 (dot-dashed lines), and s = 1/2 (dashed lines). The Si/SiGe HQ parameters are set to $\epsilon = 225 \ \mu eV$, $\Delta_1 = 19.27 \ \mu eV$, $\Delta_2 = 12.20 \ \mu eV$, and $\Delta_R = 54.18 \ \mu eV$ [22]. The bath parameters are set to $\eta = 0.5$, $\omega_c = 10\Delta_1$, and $\omega_{cutoff}/2\pi = 1$ Hz [27].

directly linked to the power spectrum $S_{\zeta}(\omega)$ by the following relations:

$$\frac{1}{T_1} = \frac{\pi}{2} \chi_{10}^2 S_{\zeta}(E_Q), \tag{12}$$

$$\frac{1}{T_{\phi}} = \frac{\pi}{4} (\chi_{11} - \chi_{00})^2 S_{\zeta}(0), \tag{13}$$

$$\frac{1}{T_2} = \frac{1}{2T_1} + \frac{1}{T_{\phi}},\tag{14}$$

where $E_Q \equiv E_1 - E_0$ is the qubit energy. The power spectrum $S_{\zeta}(0)$ is calculated at the first order of the Taylor expansion giving as a result $S_{\zeta}(0) \approx \eta/(\hbar\beta)$ for the ohmic regime and $S_{\zeta}(0) \approx \eta \omega/(\hbar\omega_c\beta)$ for the superohmic regime. Then in the extreme subohmic case, $J(\omega \approx \eta \omega_{\text{cutoff}}$ (i.e., a nonzero constant at low frequency), which corresponds to $S_{\zeta}(0) \approx \eta \omega_{\text{cutoff}}/(\hbar\omega\beta)$.

III. RESULTS

In this Section we report a detailed analysis on the relaxation and decoherence times when different experimental parameters related to the bath as well as to the HQ are varied.

In Fig. 2 the behavior of the relaxation T_1 (red lines) and the decoherence T_2 (blue lines) times calculated through Eqs. (12)–(14) is reported as a function of the bath temperature *T* for the three different regimes: s = 1 ohmic bath (solid lines), s = 2 superohmic bath (dot-dashed line), and s = 1/2subohmic bath (dashed line). The parameter η is chosen in such a way to assure the Hamiltonian-dominated regime, that is, $E_Q \gg \eta k_B T$, and at the same time to obtain relaxation and dephasing times compatible with experimental results recently obtained [21] for the parameter values analyzed. The parameters of the HQ defined in Si/SiGe QDs as well as the bath parameters are taken from the literature [22,27].

The relaxation time, as well as the decoherence, increases when the bath passes from a subohmic to a superohmic regime and decreases when the bath temperature grows.



FIG. 3. T_1 (top panels), T_{ϕ} (middle panels), and T_2 (bottom panels) as a function of ϵ and Δ_R in correspondence to an ohmic bath (s = 1) at three different temperatures: T = 0.1, 0.3, and 1.6 K. The other qubit and bath parameters are the same as those in Fig. 2.

Guided by experimental results in which the coherence times are estimated in the range of hundreds of nanoseconds [21,22], we choose to focus on the ohmic regime. We analyze in Fig. 3 the relaxation T_1 , the pure dephasing T_{ϕ} , and the decoherence T_2 times as a function of two significant qubit parameters that are the detuning ϵ that is tunable from external control voltages and the low-energy splitting of the right dot Δ_R that is linked to the qubit fabrication. We explore larger values of Δ_R with respect to the Si/SiGe case in order to include the valley splitting achievable in Si-MOS HQ [3]. For the two-dimensional (2D) plots we select three significant temperatures for experimentalists: T = 0.1, 0.3, and 1.6 K.

As is seen in Fig. 3, the relaxation time increases when the detuning is large and, in the region where ϵ is smaller, the valley splitting Δ_R has to be kept small in order to assure larger times. Looking at the pure dephasing time, it increases in the large bias region and presents also high values in a narrow section for high Δ_R where $(\chi_{11} - \chi_{00}) \simeq 0$ [see Eq. (13)]. When the relaxation time is combined with the pure dephasing time, it then gives a contribution to the total decoherence time in the large-bias region that is smaller than the contribution in the narrow section. The overall result is that the decoherence time in the large-bias region rises above its values at the narrow section, albeit the latter remains a local section of maximum for the coherence time. All the characteristic times generally reduce as the temperature is increased.

A. Relaxation time

We now focus our attention on T_1 . In Fig. 4 we report on how the two ingredients composing the T_1^{-1} behave against the detuning ϵ and Δ_R : χ_{10}^2 is plotted in Fig. 4(a), whereas the power spectrum of the bath $S_{\zeta}(E_Q)$ is presented in Fig. 4(b). Both the functions are calculated in the same range as used in Fig. 3 and $S_{\zeta}(E_Q)$ is presented for the three different



FIG. 4. (a) χ_{10}^2 as a function of ϵ and Δ_R . (b) The power spectrum $S_{\zeta}(E_Q)$ in the same (ϵ, Δ_R) range of (a) for T = 0.1, 0.3, and 1.6 K.

temperatures considered, while χ^2_{10} depends uniquely on the qubit parameters.

Note how χ_{10}^2 can be heavily modulated in the range studied, suggesting that driving an HQ in the region of small ϵ and featuring high Δ_R has to be avoided to obtain high T_1 . On the contrary, $S_{\zeta}(E_Q)$ does not present the same tunability, even if it depends on the E_Q through $J(\omega)$ and $\coth(\hbar\omega/2kT)$ with $\omega = E_Q/\hbar$ [see Eq. (10)]. Obviously $S_{\zeta}(E_Q)$ augments as *T* is increased [see again Eq. (10)], contributing to magnifying the detrimental effects of χ_{10}^2 on the relaxation times as highlighted in the corresponding plots of the first row in Fig. 3.

B. Pure dephasing and decoherence times

When the working point is set by choosing a value for ϵ , it is interesting to analyze how T_1 , T_{ϕ} , and T_2 are affected by the tunnel couplings, partially defined by the geometry of the HQ.

Our study is focused on the effect of (Δ_2, Δ_R) variations on HQ characteristic times. The Δ_1 values, which can be less effectively modulated with respect to Δ_2 values, produce small variations on T_1 , T_{ϕ} , and T_2 when compared to those caused by Δ_2 and Δ_R .

To show how this analysis is strictly connected to the HQ eigenvalues trend, in Fig. 5(a) we plot the eigenvalues of H_S and in Fig. 5(b) we plot the qubit energy E_Q (solid black line) and its derivative with respect to ϵ , that is, $dE_Q/d\epsilon$ (dashed red lines), both as a function of the detuning. All these quantities are calculated in correspondence to four different sets of Δ_2 and Δ_R at the range boundaries explored in Figs. 5(c) and 5(d) and are marked in the plots with different symbols (circle, triangle, square, and star). We also add two vertical lines highlighting the values of the detunings set to $\epsilon = 50 \ \mu eV$ (cyan) and 225 μeV (green), which are the values



FIG. 5. (a) Eigenvalues of H_S as a function of the detuning ϵ when Δ_2 and Δ_R are set. The symbols at the corners of the subplots (circle, triangle, square, and star) denote different qubit parameters (i.e., set of (Δ_2, Δ_R) values) in correspondence to the range boundaries explored in panels (c) and (d). The colored vertical lines highlight the values of the detuning $\epsilon = 50 \ \mu eV$ (cyan) and $\epsilon = 225 \ \mu eV$ (green) chosen for the plots in panels (c) and (d), respectively. (b) Energy qubit E_Q (solid black lines) and $dE_Q/d\epsilon$ (dashed red lines) as a function of ϵ for the same qubit parameter sets. (c) T_1 (top panel), T_{ϕ} (middle panel), and T_2 (bottom panel) as a function of Δ_2 and Δ_R in correspondence to an ohmic bath (s = 1) at T = 0.1 K and $\epsilon = 50 \ \mu eV$. The other parameters are the same as those in Fig. 2. (d) The same as panel (c) at $\epsilon = 225 \ \mu eV$.

chosen in Figs. 5(c) and 5(d), respectively. Figures 5(c) and 5(d) show 2D plots in which T_1 , T_{ϕ} , and T_2 are reported as a function of Δ_2 and of the low-energy splitting of the right dot Δ_R at a fixed temperature T = 0.1 K, which is of interest in experiments, for the two detunings chosen. We conclude that, independently of the value of the detuning, in correspondence to the small value of the qubit energy we have smaller relaxation and coherence times. This corresponds also to the condition in which Δ_2 is small and quite closer to Δ_1 (circle and square) for T_1 , while for T_{ϕ} and T_2 this corresponds to the condition in which Δ_2 is larger than Δ_1 (triangle and star).

Focusing on T_2 , we observe that it is marginally affected by variations of the qubit parameters in the low-bias regime [Fig. 5(c)], whereas the coherence time can be much more improved in the high-bias regime [Fig. 5(d)], especially for low values of the Δ_2 parameter. In fact, in this region (including green circles and squares), T_2 is enhanced thanks to the rise of T_{ϕ} for ϵ points with low $|dE_Q/d\epsilon|$ values, as highlighted in Fig. 5(b). To complete our analysis, we report in Fig. 6 the pure dephasing rate T_{ϕ}^{-1} as a function of $(dE_Q/d\epsilon)^2$ for the different temperatures studied.

As it can be seen, the pure dephasing rate of the qubit shows a linear dependence on $(dE_Q/d\epsilon)^2$, with higher temperatures leading to higher slopes. Note that the configurations where ϵ assumes high values (green symbols), assuring low $dE_Q/d\epsilon$, produce dephasing rates lower than those of the cases with high $dE_Q/d\epsilon$, when ϵ is low (cyan symbols). This is due to the relation $(\chi_{11} - \chi_{00}) \propto dE_Q/d\epsilon$ [24] that when inserted in Eq. (13) gives $1/T_{\phi} \propto (dE_Q/d\epsilon)^2$.

IV. CONCLUSIONS

The phonon-induced relaxation and decoherence processes are studied in the hybrid qubit in silicon quantum dots. We extract the relaxation, pure dephasing, and decoherence times as a function of the bath spectral density and of the bath temperature using the Bloch-Redfield theory. For Si quantum dots the energy dispersion is strongly affected by the physics of the valleys, so the contribution of the valley excitations



FIG. 6. Pure dephasing rate as a function of $(dE_Q/d\epsilon)^2$. The symbols correspond to different qubit parameters highlighted in Fig. 5 (green symbols correspond to $\epsilon = 50 \ \mu eV$ and cyan symbols to $\epsilon = 225 \ \mu eV$). The bath temperatures are $T = 1.6 \ K$ (solid line), $T = 0.3 \ K$ (dashed line), and $T = 0.1 \ K$ (dotted line).

- [1] D. Loss and D. P. DiVincenzo, Phys. Rev. A 57, 120 (1998).
- [2] D. P. DiVincenzo, D. Bacon, J. Kempe, G. Burkard, and K. B. Whaley, Nature (London) 408, 339 (2000).
- [3] M. Veldhorst, J. C. C. Hwang, C. H. Yang, A. W. Leenstra, B. de Ronde, J. P. Dehollain, J. T. Muhonen, F. E. Hudson, K. M. Itoh, A. Morello, and A. S. Dzurak, Nat. Nanotechnol. 9, 981 (2014).
- [4] J. J. L. Morton, D. R. McCamey, M. A. Eriksson, and S. A. Lyon, Nature (London) 479, 345 (2011).
- [5] E. Kawakami, P. Scarlino, D. R. Ward, F. R. Braakman, D. E. Savage, M. G. Lagally, M. Friesen, S. N. Coppersmith, M. A. Eriksson, and L. M. K. Vandersypen, Nat. Nanotechnol. 9, 666 (2014).
- [6] E. Ferraro, M. Fanciulli, and M. De Michielis, J. Phys. Commun. 2, 115022 (2018).
- [7] E. Ferraro, M. Fanciulli, and M. De Michielis, Quantum Inf. Process. 17, 130 (2018).
- [8] E. Ferraro, M. Fanciulli, and M. De Michielis, Adv. Quantum Technol. 1, 1800040 (2018).
- [9] C. B. Simmons, J. R. Prance, B. J. VanBael, T. S. Koh, Z. Shi, D. E. Savage, M. G. Lagally, R. Joynt, M. Friesen, S. N. Coppersmith, and M. A. Eriksson, Phys. Rev. Lett. **106**, 156804 (2011).
- [10] M. Xiao, M. G. House, and H. W. Jiang, Phys. Rev. Lett. 104, 096801 (2010).
- [11] J. M. Kikkawa and D. D. Awschalom, Phys. Rev. Lett. 80, 4313 (1998).
- [12] C. Barthel, J. Medford, C. M. Marcus, M. P. Hanson, and A. C. Gossard, Phys. Rev. Lett. **105**, 266808 (2010).
- [13] D. Rotta, M. De Michielis, E. Ferraro, M. Fanciulli, and E. Prati, Quantum Inf. Process. 15, 2253 (2016).
- [14] D. Rotta, F. Sebastiano, E. Charbon, and E. Prati, npj Quantum Inf. 3, 26 (2017).

has been effectively included in our analysis. It is found that the characteristics of both the spectral density of the bath and the energy spectrum of the qubit play an essential role. The contribution of phonons to relaxation and pure dephasing effects is bias dependent, leading to the conclusion that the coherence time can be higher in the large-bias region than in the small-bias region, due to stronger relaxation at small bias. We also observed that the relaxation time is much more affected by the qubit energy spectrum than by the bath power spectrum. The pure dephasing rate exhibits a linear dependence on the square of the derivative of the qubit energy with respect to the detuning. This demonstrates a strong inverse proportionality, small values of the derivative of the qubit energy correspond to larger dephasing times, as confirmed by the experiments. Moreover, the higher the temperature is, the higher the slope is, meaning that at higher temperature there is a large variability of the pure dephasing time with respect to the energy qubit derivative.

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- [15] M. De Michielis, E. Prati, M. Fanciulli, G. Fiori, and G. Iannaccone, Appl. Phys. Exp. 5, 124001 (2012).
- [16] M. Friesen and S. N. Coppersmith, Phys. Rev. B 81, 115324 (2010).
- [17] Z. Shi, C. B. Simmons, J. R. Prance, J. K. Gamble, T. S. Koh, Y.-P. Shim, X. Hu, D. E. Savage, M. G. Lagally, M. A. Eriksson, M. Friesen, and S. N. Coppersmith, Phys. Rev. Lett. 108, 140503 (2012).
- [18] D. Kim, D. R. Ward, C. B. Simmons, D. E. Savage, M. G. Lagally, M. Friesen, S. N. Coppersmith, and M. A. Eriksson, npj Quantum Inf. 1, 15004 (2015).
- [19] E. Ferraro, M. De Michielis, G. Mazzeo, M. Fanciulli, and E. Prati, Quantum Inf. Process. 13, 1155 (2014).
- [20] M. De Michielis, E. Ferraro, M. Fanciulli, and E. Prati, J. Phys. A 48, 065304 (2015).
- [21] B. Thorgrimsson, D. Kim, Y.-C. Yang, L. W. Smith, C. B. Simmons, D. R. Ward, R. H. Foote, J. Corrigan, D. E. Savage, M. G. Lagally, M. Friesen, S. N. Coppersmith, and M. A. Eriksson, npj Quantum Inf. 3, 32 (2017).
- [22] J. C. Abadillo-Uriel, B. Thorgrimsson, D. Kim, L. W. Smith, C. B. Simmons, D. R. Ward, R. H. Foote, J. Corrigan, D. E. Savage, M. G. Lagally, M. J. Calderón, S. N. Coppersmith, M. A. Eriksson, and M. Friesen, Phys. Rev. B 98, 165438 (2018).
- [23] C. H. Yang, A. Rossi, R. Ruskov, N. S. Lai, F. A. Mohiyaddin, S. Lee, C. Tahan, G. Klimeck, A. Morello, and A. S. Dzurak, Nat. Commun. 2, 2069 (2013).
- [24] X.-K. Qin, Europhys. Lett. 114, 37006 (2016).
- [25] M. Borhani, V. N. Golovach, and D. Loss, Phys. Rev. B 73, 155311 (2006).
- [26] V. Kornich, C. Kloeffel, and D. Loss, Quantum 2, 70 (2018).
- [27] L. M. Cangemi, G. Passarelli, V. Cataudella, P. Lucignano, and G. De Filippis, Phys. Rev. B 98, 184306 (2018).