

Confinement effects on decay rate of surface electron states over liquid helium

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The decay rate of excited states of surface electrons in liquid helium, trapped in a quantum dot system, is evaluated, taking into account the process of spontaneous radiation of two-rippbons with short wavelength. We find that the values of the decay rate in later process are rather higher than those for the one-rippon process previously calculated. The upper-bound limit lifetime of excited states of surface electrons in a quantum dot is found to be $\tau < 10^{-6}-10^{-7}$ s.

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Surface electrons (SEs) localized over the liquid helium surface have been an remarkable archetype for studying confined charged systems of low dimensionality [1,2]. In particular and more important, SE systems have been intensively investigated in the last years as potential candidate of qubit generators in quantum computation [3,4]. The main reasons for this are twofold. Firstly, the energy spectrum of SEs is quite simple and well-known. In the limit of zero holding electric field E_{\perp} oriented in the normal direction to the electron layer, the spectrum is given by hydrogenic-like subbands

$$\Delta_l = -\frac{m\Lambda_0^2}{2\hbar^2 l^2}; \quad l=1, 2, 3, \dots, \quad (1)$$

where $\Lambda_0 = (e^2/4)(\epsilon_{\text{He}} - 1)/(\epsilon_{\text{He}} + 1)$, $\epsilon_{\text{He}} \simeq 1.0572$ is liquid helium dielectric constant, e and m are electron charge and mass, respectively. The total electron energy is given by $E_{\mathbf{k},l} = E_{\mathbf{k}} + \Delta_l$; $E_{\mathbf{k}} = \hbar^2 k^2/(2m)$ where \mathbf{k} is 2D electron wave vector in the helium liquid–vapor interface. For nonzero holding field, the dependence $\Delta_l(E_{\perp})$ can be calculated within the variational approach

[5,6] (see Fig. 1 for $l=1$ and 2). The minimum energy gap between ground $l=1$ and first excited ($l=2$) subbands is $\Delta_{21} = \Delta_2 - \Delta_1 \simeq 5.7$ K for $E_{\perp} = 0$ and it is increasing

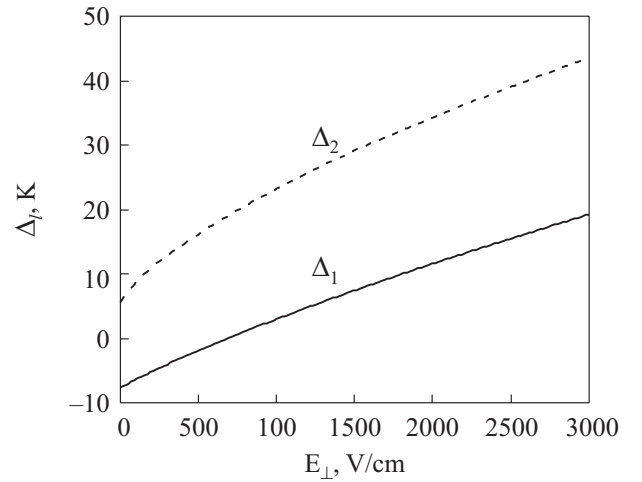


Fig. 1. The energies Δ_1 of the ground subband $l=1$ and Δ_2 of the first excited subband $l=2$.

function of E_{\perp} , attaining, for example, 8.7 (15.5) K for $E_{\perp} = 100$ (450) V/cm.

Secondly, the SE scattering along the liquid–vapor interface is limited for temperatures $T < 1$ K by ripples, the quantized capillary waves of the He surface, whose dispersion law is $\omega_{\mathbf{q}} \simeq (\alpha / \rho) q^{3/2}$ (α and ρ are the He surface tension coefficient and bulk density, respectively, q is wave number). The strength of the electron–ripple interaction decreases strongly with T which favours the experimental study of qubits produced by SE. One should emphasize that even though the electron–ripple interaction is essential to define both transport phenomena [2] and decay rate of excited electron states it is rather weak to influence the subband structure. The problem of the decay rate and the subband lifetime τ , which is inversely proportional to decay rate, was first discussed in Ref. 4. The authors concluded that the contribution of the one-ripple scattering processes predominates. In such a case the SE scattering is due to long wavelength ripples ($q \leq 10^6 - 10^7 \text{ cm}^{-1}$) and τ is inversely proportional to temperature attaining rather optimistic values $\tau \sim 10^{-5} - 10^{-4}$ s for T in the range of millikelvin. Furthermore a strong increase of the lifetime due to confinement effects in the plane of electron layer was predicted in presence of magnetic field normal to SE system.

However, the situation is quite different if one takes into account the contribution of SE scattering by short wavelength ripples [7]. This process occurs because the nonlinearity of the electron–ripple interaction. As it was shown in Ref. 7 even at very low temperature the SE can radiate a pair of short wavelength ripples in almost opposite directions. The total energy of the ripples $2\hbar\omega_{\mathbf{q}}$ can be of order of $E_{\mathbf{k},l}$ and the ripple-pair momentum satisfies the condition $s = |\mathbf{q} + \mathbf{q}'| \ll q$. The two-ripples process is responsible for the energy relaxation of SE system [7]. It was shown that its contribution to the decay rate of the excited level $l=2$ is two orders of magnitude higher than that of one-ripple process, i.e., $\tau \sim 10^{-6} - 10^{-7}$ s [8]. Furthermore confinement effects due to the presence of a perpendicular magnetic field, which now makes discrete the energy $E_{\mathbf{k}}$ (Landau spectrum), do not alter substantially the values of τ .

Recently, the trap of SE in a quantum dot with few electrons by specially designed electrostatic gates was realized [9]. Besides the fact that magnetic and electrostatic confinements have different physical and practical implications, the probability of SE transitions between in-plane energy levels in quantum dot do not depend on the SE momentum which leads to different matrix elements for the interaction potential. Accordingly the calculation of the SE lifetime and decay rate in quantum dot is required. The aim of present work is to perform the theoretical calculation of the decay rate in quantum dot due to the two-ripple creation process.

The discrete energy spectrum of the SE in a quantum dot, in the asymmetrical parabolic approximation with characteristic frequencies ω_x and ω_y , is written as

$$E_{n_x, n_y, l} = \hbar\omega_x \left(n_x + \frac{1}{2} \right) + \hbar\omega_y \left(n_y + \frac{1}{2} \right) + \Delta_l, \quad (2)$$

where $n_x, n_y = 0, 1, 2 \dots$. Here we assume the axis z to be normal to the electron layer located in the xy plane. In this approximation, the wave functions for the motion along x and y are given in terms of Hermite functions.

The decay rate and lifetime for both excited level $l=2$ and fundamental level $l=1$ is evaluated by considering the in-plane excited states with $n_x, n_y \geq 1$. Note that the decay rate defines the linewidth of the spectroscopic transitions between the SE energy levels [10].

We start from the basic relation for energy state decay rate defined by the interaction given by the Hamiltonian $\hat{H}(\mathbf{r}, z, t)$ [11]

$$\Gamma_{\{j\}} = \frac{1}{\hbar^2} \sum_{\{j'\}} \int_{-\infty}^{\infty} dt \exp \left[i \frac{(E_{\{j'\}} - E_j)t}{\hbar} \right] \langle U_{\{j'\}}(0) U_{\{j'\}}(t) \rangle, \quad (3)$$

where $U_{\{j'\}}(t) = \langle \{j'\} | \hat{H}(\mathbf{r}, z, t) | \{j\} \rangle$ is the ensemble-averaged matrix element for the states $\{j'\}$ and $\{j\}$, $\mathbf{r} = \{x, y\}$ is 2D position-vector in the plane of the electron layer. A straightforward calculation leads to the following expression for the decay rate in a quantum dot due to the two-ripple process:

$$\begin{aligned} \Gamma_{\{n_x, n_y, l\}}^{(2r)} &\simeq \frac{4\pi}{\hbar} \sum_{n'_x, n'_y, l'} \sum_{\mathbf{q}, \mathbf{s}} Q_{\mathbf{q}}^2 Q_{\mathbf{q}-\mathbf{s}}^2 |\langle l | V_{\mathbf{q}}^{(2r)} | l \rangle|^2 |J_{n_x, n'_x}(s_x)|^2 |J_{n_y, n'_y}(s_y)|^2 (N_{\mathbf{q}} + 1)(N_{\mathbf{q}-\mathbf{s}} + 1) \times \\ &\times \delta [\Delta_l - \Delta_{l'} + \hbar\omega_x(n'_x - n_x) + \hbar\omega_y(n'_y - n_y) + \hbar\omega_{\mathbf{q}} + \hbar\omega_{\mathbf{q}-\mathbf{s}}]. \end{aligned} \quad (4)$$

Here only radiation processes are included. In Eq. (4) and, $N_{\mathbf{q}}(T)$ is the ripple thermal distribution function $Q_{\mathbf{q}}^2 = \hbar q / (2\rho\omega_{\mathbf{q}})$. The matrix elements for in-plane electron motion are given by

$$|J_{n_x, n'_x}(s_j)|^2 = \frac{[\min(n_j, n'_j)]!}{[\max(n_j, n'_j)]!} \left(\frac{s_j^2 \lambda_j^2}{2} \right)^{|n_j - n'_j|} \exp \left(-\frac{s_j^2 \lambda_j^2}{2} \right) \left[L_{\min(n_j, n'_j)}^{|n_j - n'_j|} \left(\frac{s_j^2 \lambda_j^2}{2} \right) \right]^2, \quad (5)$$

where $L_V^\beta(x)$ are Laguerre polynomials and $\lambda_j^2 = \hbar / (m\omega_j)$ with $\omega_j = \omega_x$ if $j = x$ and $\omega_j = \omega_y$ if $j = y$. The structure of the matrix elements $\langle l | V_{\mathbf{q}}^{(2r)} | l \rangle$ for the two-ripplon process is rather interesting. As it was shown in Ref. 2 (see also Ref. 8) the matrix elements for the two-ripplon process in the electron-ripplon scattering should be calculated within the Bloch treatment for ripplon wave numbers exceeding $q \simeq 2 \cdot 10^7 \text{ cm}^{-1}$. For such q 's, typical for two-ripplon processes which are quadratic in the displacement $\xi(\mathbf{r})$ of the liquid-vapor interface from the equilibrium position $z = 0$, one has

$$\langle l | V_{\mathbf{q}}^{(2r)} | l \rangle = \kappa_0 \sqrt{\left(\frac{\partial v}{\partial z} \right)_{ll} \left(\frac{\partial v}{\partial z} \right)_{l'l'}}$$
 (6)

where $(\partial v / \partial z)_{ll} = \langle l | v(z) | l \rangle$, $v(z) = -\Lambda_0 / z + eE_{\perp} z$ and $\kappa_0 = \sqrt{2mV_0} / \hbar$. Indeed the matrix elements given by Eq. (6) depend explicitly on the height of the potential barrier $V_0 \simeq 1 \text{ eV}$ which prevents SE penetrating inside the liquid helium. The wave functions $f_l(z)$ for $l=1$ and $l=2$ are calculated within the variational approach [5,6].

Taking into account that $N_{\mathbf{q}} \ll 1$ and disregarding s in comparison with q in Eq. (4), we arrive to the following expressions for the decay rates for $l=2$ and $l=1$:

$$\Gamma_{(n_x, n_y, 2)}^{(2r)} = \frac{\kappa_0^2}{12\pi^2 \alpha^{3/2} \rho^{1/2} \lambda_x \lambda_y} \times \sum_{n'_x n'_y} I_{n_x n'_x} I_{n_y n'_y} \left[\frac{\left(\frac{\partial v}{\partial z} \right)_{22}^2}{q_0^{1/2}} + \frac{\left(\frac{\partial v}{\partial z} \right)_{11} \left(\frac{\partial v}{\partial z} \right)_{22}}{(q_0^*)^{1/2}} \right];$$
 (7)

$$\Gamma_{(n_x, n_y, 1)}^{(2r)} = \frac{\kappa_0^2}{12\pi^2 \alpha^{3/2} \rho^{1/2} \lambda_x \lambda_y} \times \sum_{n'_x n'_y} I_{n_x n'_x} I_{n_y n'_y} \left[\frac{\left(\frac{\partial v}{\partial z} \right)_{11}^2}{q_0^{1/2}} + \frac{\left(\frac{\partial v}{\partial z} \right)_{11} \left(\frac{\partial v}{\partial z} \right)_{22}}{(q_0^{**})^{1/2}} \right],$$
 (8)

where the characteristic ripplon wave numbers are

$$q_0 = (\rho / 4\alpha)^{1/3} [\omega_x (n_x - n'_x) + \omega_y (n_y - n'_y)]^{2/3};$$

$$q_0^* = (\rho / 4\alpha)^{1/3} [\Delta_{21} / \hbar + \omega_x (n_x - n'_x) + \omega_y (n_y - n'_y)]^{2/3};$$

$$q_0^{**} = (\rho / 4\alpha)^{1/3} [-\Delta_{21} / \hbar + \omega_x (n_x - n'_x) + \omega_y (n_y - n'_y)]^{2/3}$$

and

$$I_{n_j, n'_j} = \frac{[\min(n_j, n'_j)]!}{[\max(n_j, n'_j)]!} \times \int_0^\infty x^{|n_j - n'_j| - 1/2} \exp(-x) \left[L_{\min(n_j, n'_j)}^{|n_j - n'_j|}(x) \right]^2 dx.$$

The sums which appear in Eqs. (7) and (8) are carried out over the values of n'_x and n'_y satisfying the condition of positiveness of the expressions in square brackets of q_0 , q_0^* , and q_0^{**} , respectively. One emphasizes that the decay rate does not depend on temperature which is a direct consequence of neglecting $N_{\mathbf{q}}(T)$ in Eq. (4) because it is very small for short wavelength ripples. On the contrary, for long wavelength ripples $N_{\mathbf{q}} \simeq T / \hbar\omega_{\mathbf{q}} \gg 1$. Therefore a linear T -dependence of the decay rate is obtained when only the one-ripplon process is considered [4].

The first term in the sums of Eqs. (7) and (8) correspond to intrasubband scattering contributions whereas the second term is related to intersubband scattering.

Looking over the expressions for the decay rates one must emphasize that the matrix elements given by Eq. (5) depend on $s_j^2 \lambda_j^2 / 2 \sim 1$ for $s \ll q$. Likewise the calculation for the one-ripplon process leads to the matrix elements which depend on $q_j \lambda_j \gg 1$. Consequently the matrix elements in Eq. (5) being proportional to $\exp(-q_j^2 \lambda_j^2 / 2) \ll 1$ are quite small. The conclusion is that the contribution from one-ripplon process for the decay rate of SE in a quantum dot is absolutely negligible in comparison with that from two-ripplon process.

In Fig. 2 we depicted the dependence of the lifetimes $\tau_2^{(2r)} = [\Gamma_{\{n_x, n_y, 2\}}^{(2r)}]^{-1}$ and $\tau_1^{(2r)} = [\Gamma_{\{n_x, n_y, 1\}}^{(2r)}]^{-1}$ on the holding electric field for $n_x = 1$ and $n_y = 0$. The energies of the parabolic confinement are $\hbar\omega_x = 0.5 \text{ K}$ and $\hbar\omega_y = 0.6 \text{ K}$, respectively. We see that the order of magnitude of the

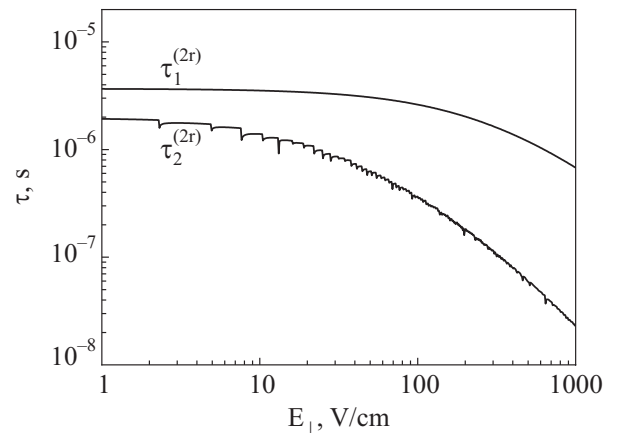


Fig. 2. The lifetime $\tau_2^{(2r)}$ for the SE excited states $\{n_x = 1; n_y = 0, l = 2\}$ and $\tau_1^{(2r)}$ for $\{n_x = 1; n_y = 0; l = 1\}$ limited by the two-ripplon scattering process.

lifetimes is 10^{-6} s for $E_{\perp} = 0$ and both $\tau_2^{(2r)}$ and $\tau_1^{(2r)}$ decrease strongly as the holding electric field increases. Note that the values of $\tau_2^{(2r)}$ are close to those obtained in Ref. 8 both for free electron motion in xy plane and in presence of magnetic field along z . We also observe the values of $\tau_1^{(2r)}$ are higher than those of $\tau_2^{(2r)}$. We should remark that only the first term of the intrasubband scattering in the sum of Eq. (8) contributes to $\Gamma_{\{n_x, n_y, l\}}^{(2r)}$

and consequently to $\tau_1^{(2r)}$ for the chosen $\hbar\omega_x = 0.5$ K and $\hbar\omega_y = 0.6$ K. The sharp difference between the values of $\tau_2^{(2r)}$ and $\tau_1^{(2r)}$ may be favourable in measurements of the spectroscopic linewidth in a quantum dot. Indeed the signals of the transitions from the excited SE subband $l=2$ and from excited states of in-plane motion with $n_x, n_y \geq 1$ for $l=1$ should be clearly distinguished. Interestingly enough are the oscillations appearing in $\tau_2^{(2r)}(E_{\perp})$ whose origin comes from the denominator of the second term of the sum in Eq. (7). One can easily check that there is a square root singularity for Δ_{21} / \hbar close to $\omega_x(n'_x - n_x) + \omega_y(n'_y - n_y)$. For example for $E_{\perp} = 2.3$ V/cm ($\Delta_{21} \simeq \simeq 5.8$ K) the q_0^* trends to zero for $\{n'_x = 3; n'_y = 8\}$ and $\{n'_x = 9; n'_y = 3\}$. Similar oscillations were observed for SE decay rate under magnetic field [8]. One should stress that the Eq. (4), obtained for the process of creation of two short wavelength ripples with $q > 10^7$ cm $^{-1}$, becomes invalid for $q \rightarrow 0$. For this reason the neighborhood of this singularity should be treated appropriately by taking into account on equal footing both two-ripples and one-ripple processes. On the other hand, note that the first term in the sum of Eq. (8) has no singularities. In this case, only the state $n'_x = n'_y = 0$ does contribute to $\Gamma_{(n_x, n_y, l)}^{(2r)}$ for all holding fields. As a result, a smooth dependence $\tau_1^{(2r)}(E_{\perp})$ is obtained.

In conclusion we determined the decay rate of excited states of surface electrons on liquid helium trapped in a

quantum dot. We showed that both decay rate and lifetime of the excited states do not depend on temperature and are limited by the process of spontaneous creation of a pair of short wavelength ripples (with wave numbers exceeding 10^7 cm $^{-1}$). The values of the decay rate are essentially higher and therefore the lifetimes are much smaller than those limited by one-ripple process [4]. The results obtained here should be essential for the experimental study of qubits based on surface electrons on helium.

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