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# Intraband absorption of electromagnetic radiation by electrons with optical phonon participation in quantum dot superlattices

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**Abstract:** In this paper, we investigate the absorption of electromagnetic radiation by a free electron gas interacting with lattice vibrations in a quantum dot superlattice. It is assumed that the electron gas in the quantum dot superlattice is limited by the anisotropic parabolic potential. The absorption of light by free carriers with the participation of phonons is calculated in the second order of the perturbation theory. When calculating the absorption coefficient, the matrix elements of the electron-photon and electron-phonon interactions (with polar and nonpolar optical phonons) are used. We can expect three possible transitions in the absorption coefficient for electron- polar phonon scattering: (1) a transition due to the size subband levels for only the x direction, (2) a transition due to the size subband levels for only the y direction and (3) a transition due to the size subband levels for both the x direction and the y direction. For electron- nonpolar phonon scattering we can expect only one possible transitions in the, the absorption coefficient: a transition due to the size subband levels for both the x direction and the y direction.

**Keywords:** quantum dot, superlattice, miniband, intraband absorption

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## 1. INTRODUCTION

Advances in methods for growing crystals with sizes close to the interatomic distance, such as molecular beam epitaxy and metal-organic vapor deposition, have made it possible to obtain such low-dimensional systems as quantum wells, superlattices, quantum dots, etc. Since the beginning of the pioneering work of Esaki and Tsu [1], a large amount of research has been devoted to semiconductor superlattices due to their transport properties and their application in electronic devices such as generators, tunnel diodes, hot electron transistors, and optoelectronic devices [2,3]. Modern nanotechnology makes it possible to

manufacture quantum dots of various shapes, which are successfully used in various devices, quantum cascade lasers based on quantum dot superlattices [1–3]. Such systems have a significantly higher performance compared to existing quantum cascade lasers based on quantum well superlattices [4]. Superlattices of quantum dots attract the attention of researchers also due to their unusual optical and optoelectronic properties, which can be used to develop new optoelectronic devices [5-13].

It is well known that the electron-phonon interaction is an important factor affecting the physical properties of polar crystals, such as the binding energy of impurities, charge transfer, linear and nonlinear optical properties, especially in low-dimensional quantum systems [14,15].

In this paper, we develop a theory of light absorption by free carriers in a superlattice of quantum dots with an anisotropic parabolic potential upon scattering of carriers by polar and nonpolar optical phonons.

In bulk semiconductors, a quantum mechanical theory of light absorption by free carriers using the second-order perturbation theory was given by Dumke [16]. Later, it was studied in detail for acoustic phonons (through the deformation potential and piezoelectric coupling), non-polar and polar optical phonons, and impurities in bulk semiconductors [17]. These studies have shown a different power-law dependence of the absorption coefficient on wavelength to radiation for different scattering mechanisms. Moreover, these power laws differ from the classical dependence of the absorption coefficient. The absorption of light by free carriers in the scattering of electrons by phonons has been theoretically investigated in semiconductor quantum wells and wires [18-28] and in quantum dots [29,30].

## 2. THEORETICAL FRAMEWORK

It is known that the motion of an electron in a superlattice is limited and its energy spectrum is quantized. It is assumed that the quantization takes place in the  $z$ -direction. The electron-phonon interaction in a quantum dot superlattice with a periodic potential  $U(z)$  of period  $d$  along the  $z$ -direction of the form [31] is considered:

$$H = G_{\text{GAPL}}(v) = \text{Concat}_{j=1}^{\text{B}} \left( \frac{1}{\text{B}} \sum_{a=1}^{\text{B}} v_a \right) \in \mathbb{R}^{\text{B} \times \text{O}}. \quad (1)$$

where  $m^*$  is the effective mass and  $\omega_x$ ,  $\omega_y$  are the confinement frequencies in the  $x$ - and  $y$ -directions, respectively. In the strong-coupling approximation, the Hamiltonian for charge carriers in a quantum dot superlattice can be written as [8,9]:

$$H = \frac{(p_x^2 + p_y^2)}{2m^*} + \frac{m^*}{2} (\omega_x^2 x^2 + \omega_y^2 y^2) + \frac{\Delta}{2} (1 - \cos \frac{p_z d}{\hbar}), \quad (2)$$

where  $\Delta$  is the miniband width. The normalized eigenwave functions of the electron  $\Psi_{n,e,k_z}(r)$  and the eigenvalues of the energy  $E_{n,l}(k_z)$  in the conduction band are set, respectively, in the form [10]:

$$\Psi_{n,l,k_z}(r) = \frac{1}{\sqrt{L_z}} \Psi_n(x) \Psi_l(y) \xi_{k_z}(z), \quad (3)$$

$$E_{n,l}(k_z) = (n + \frac{1}{2}) \hbar \omega_x + (l + \frac{1}{2}) \hbar \omega_y + \frac{\Delta}{2} (1 - \cos k_z d) = \varepsilon_{n,l} + \varepsilon(k_z), \quad (4)$$

where  $n = 0, 1, 2, \dots$  and  $l = 0, 1, 2, \dots$  are the level indices of the electronic subbands,  $k_z$  is the component of the wave vector in the  $z$ -direction,  $\Psi_n(x) = \sqrt{1/\sqrt{\pi} l_x} 2^n n! \exp[-x^2/2l_x^2] H_n(x/l_x)$ ,  $\Psi_l(y) = \sqrt{1/\sqrt{\pi} l_y} 2^l l! \exp[-y^2/(2l_y^2)] H_l(y/l_y)$  are the eigenwave functions of the harmonic oscillator,  $H_n(x)$  is the Hermite polynomial,  $l_x = \sqrt{\hbar/(m^* \omega_x)}$ ,  $l_y = \sqrt{\hbar/(m^* \omega_y)}$ ,  $\xi_{k_z}(z)$  is the Bloch function in the approximation of strong coupling in the  $z$ -direction, and  $L_z$  is the normalized length in the direction.

The absorption of light by free carriers with the participation of phonons is calculated in the second order of the perturbation theory. The absorption coefficient is determined by the well-known formula [20]:

$$\alpha = \frac{\varepsilon^{1/2}}{n_0 c} \sum_i W_i f_i, \quad (5)$$

where  $\varepsilon$  is the dielectric constant,  $n_0$  is the number of photons in the radiation field,  $f_i$  is the distribution function of free current carriers,  $W_i$  is the transition probability, determined by the following expression:

$$W_i = \frac{2\pi}{\hbar} \sum_{f,i} \left[ |\langle f | M_+ | i \rangle|^2 \delta(E_f - E_i - \hbar\Omega - \hbar\omega_q) + |\langle f | M_- | i \rangle|^2 \delta(E_f - E_i - \hbar\Omega + \hbar\omega_q) \right], \quad (6)$$

where  $E_i$  and  $E_f$  are the energies of the initial and final states of electrons, respectively,  $\hbar\Omega$  is the photon energy,  $\hbar\omega_q$  is the phonon energy, and  $\langle f | M_{\pm} | i \rangle$  are the elements of the matrix of transition from the initial state to the final state for the interaction between electrons, phonons, and photons. Elements of the transition matrix can be represented as:

$$\langle f | M_{\pm} | i \rangle = \sum_{\alpha} \left( \frac{\langle f | H_R | \alpha \rangle \langle \alpha | V_s | i \rangle}{E_i - E_{\alpha} \mp \hbar\omega_q} + \frac{\langle f | V_s | \alpha \rangle \langle \alpha | H_R | i \rangle}{E_i - E_{\alpha} - \hbar\Omega} \right), \quad (7)$$

where  $i, \alpha, f$  indices denote the initial, intermediate and final states of an electron with quantum numbers  $k, n, l$ ,  $H_R$  is the operator of electron-photon interaction,  $V_s$  is the operator of electron-phonon interaction.

Substituting expressions for wave functions (3) into expressions (7) for matrix elements of electron-photon interaction, we obtain:

$$\langle n'l'k'_z | H_R | nlk_z \rangle = \frac{e\Delta d \sin(k_z d)}{2\hbar} \left( \frac{2\pi\hbar n_0}{V\Omega \epsilon} \right)^{1/2} \delta_{nm} \delta_{l'l'} \delta_{kk_z}, \quad (8)$$

where  $V$  is the volume of the crystal, the radiation field is polarized along the  $\hat{z}$ -direction,  $\epsilon$  – is the vector of polarization.

The electron distribution function for a nondegenerate electron gas has the form:

$$f_0(E_{nk_z}) = \frac{2n_{1D} d e^{\frac{\Delta}{k_B T}} \sinh\left(\frac{\hbar\omega_x}{2k_B T}\right) \sinh\left(\frac{\hbar\omega_y}{2k_B T}\right)}{M\left(\frac{1}{2}, 1, \frac{\Delta}{k_B T}\right)} \exp\left(-\frac{E_{nk_z}}{k_B T}\right), \quad (9)$$

where  $n_{1D}$  is the number of electrons per unit length and  $M(a, c, x)$  is the confluent hypergeometric function [31]

$$M(a, c; x) = \sum_{n=0}^{\infty} \frac{(a)_n x^n}{(c)_n n!}. \quad (10)$$

The matrix element of the electron-phonon interaction has the following form:

$$\langle k'_z n' l' | V_s | k_z n l \rangle^2 = C_j'^2 J_{m'}(x) J_{l'}(y) I(q_z), \quad (11)$$

where  $V_s$  is the operator of the interaction energy of an electron with a phonon,  $C_j$  is a function that characterizes the interaction between electrons and phonons,

$$\begin{aligned} J_{m'}(q_x) &= \int_{-\infty}^{\infty} e^{iq_x x} dx \Psi_m(x) \Psi_{m'}(x), \\ J_{l'}(q_y) &= \int_{-\infty}^{\infty} e^{iq_y y} dy \Psi_l(y) \Psi_{l'}(y), \\ I(q_z) &= \int_0^d \xi_{k_z}(z) \xi_{k'_k}(z) e^{iq_z z} dz, \end{aligned} \quad (12)$$

For the interaction of an electron with polar optical phonons, we have:

$$\begin{aligned} \tilde{N}_{POL}^2 &= 2\pi e^2 \hbar \omega_0 \left\{ \frac{1}{\epsilon_{\infty}} - \frac{1}{\epsilon_0} \right\}, \\ F_{POL} &= \frac{N_0^{\pm}}{q^2}, \quad N_0 = \left[ \exp\left(\frac{\hbar\omega_0}{K_B T}\right) - 1 \right]^{-1}, \end{aligned} \quad (13)$$

where  $N_0^- = N_0$ ,  $N_0^+ = N_0 + 1$ .

Here  $\epsilon_{\infty}$  and  $\epsilon_0$  are the high-frequency and static dielectric constants of the semiconductor, respectively. The phonon energy is taken  $\hbar\omega_q = \hbar\omega_0 = \text{const}$

$$N_0 = \left[ \exp\left(\frac{\hbar\omega_0}{K_B T}\right) - 1 \right]^{-1}, \quad N_0^- = N_0, \quad N_0^+ = N_0 + 1, \quad (14)$$

where  $N_0^-$  ( $N_0^+$ ) describes the annihilation and creation of a phonon.

For the interaction of an electron with non-polar optical phonons

$$C_{np}^2 = \frac{\hbar D}{2\rho\omega_0 V}, \quad F_{np}(q) = N_0^{\pm}, \quad (15)$$

where  $D$  is the constant of the non-polar optical deformation potential.

Taking into account formulas (6) and (11) in (5) and summing over  $k'_z$ , for the absorption coefficient we obtain:

$$\begin{aligned} \alpha &= \left( \frac{\pi e \Delta d}{\hbar^2} \right)^2 \frac{1}{cV\Omega^3 \epsilon^{3/2}} \times \\ &\times \sum_{n'e' n k_z} \sum_{m' C_j} \sum_q |I_{m'}(q_x)|^2 |I_{e'}(q_y)|^2 (\sin(k_z + q_z) d - \sin k_z d)^2 \times \\ &\times \left\{ N_0 \delta \left( (n'-n)\hbar\omega_x + (e'-e)\hbar\omega_y + \frac{\Delta}{2} (\cos k_z d - \cos(k_z + q_z) d - \hbar\Omega + \hbar\omega_q) \right) + \right. \\ &\left. + (N_0 + 1) \delta \left( (n'-n)\hbar\omega_x + (l'-l)\hbar\omega_y + \frac{\Delta}{2} (\cos k_z d - \cos(k_z + q_z) d) - \hbar\Omega - \hbar\omega_q \right) \right\}. \end{aligned} \quad (16)$$

In formula (16), passing from summation over  $k_z$  and  $q$  to integration over  $k_z$  and  $q$  using the expression:

$$\sum_q \frac{V}{(2\pi)^3} \int_{-\infty}^{\infty} \int_{-\infty}^{\infty} \int_{-\pi/d}^{\pi/d} dq_x dq_y dq_z, \quad (17)$$

$$\sum_{k_z} (\dots) = \left( \frac{L_z}{2\pi} \right) \int_{-\pi/d}^{\pi/d} (\dots), \quad (18)$$

and using the properties of the Dirac  $\delta$ -function:

$$\delta[f(x)] = \sum_i \frac{\delta(x - x_i)}{|f'(x_i)|}, \quad (19)$$

where  $x_i$  are simple zeros of the function  $f(x)$ ,

$$\int_{-\pi/d}^{\pi/d} dq_z f(q_z) \delta(q_z - a) = f(a). \quad (20)$$

for the absorption coefficient for scattering by polar and non-polar optical phonons, we obtain:

$$\begin{aligned} \alpha_{pol} &= \frac{4\pi e^4 \Delta d \omega_0 L_z}{c\Omega^3 \epsilon^{3/2} \hbar^3} \left( \frac{1}{\epsilon_{\infty}} - \frac{1}{\epsilon_0} \right) \times \\ &\times \sum_{n'e' n} \sum_{\pm} \int_{-\pi/d}^{\pi/d} dk_z f_{nek_z} \left( N_0 + \frac{1}{2} \pm \frac{1}{2} \right) \frac{\sqrt{1 - \frac{4}{\Delta^2} \theta_{\pm}^2 - \sin k_z d}}{\sqrt{1 - \frac{4}{\Delta^2} \theta_{\pm}^2}} \times \end{aligned} \quad (21)$$

$$\times \int_0^{\infty} \int_0^{\infty} \frac{|I_{m'}(q_x)|^2 |I_{e'}(q_y)|^2}{(q_x^2 + q_y^2) + a_{\pm}^2} dq_x dq_y,$$

$$\alpha_{n,pol} = \frac{D^2 e^2 \Delta d}{\pi c \rho \omega_0 \Omega^3 L_x L_y \epsilon^{3/2}} \times \sum_{n'e'} \sum_{ne} \sum_{\pm} \int_{-\pi/d}^{\pi/d} dk_z f_{nek_z} \left( N_0 + \frac{1}{2} \pm \frac{1}{2} \right) \frac{\sqrt{1 - \frac{4}{\Delta^2} \Theta_{\pm}^2 - \sin^2 k_z d}}{\sqrt{1 - \frac{4}{\Delta^2} \Theta_{\pm}^2}} \times \frac{1}{n' - n} \cdot \frac{1}{l' - l}, \quad (22)$$

where

$$\Theta_{\pm}(k_z d) = (n' - n) \hbar \omega_x + (l' - l) \hbar \omega_y \pm \hbar \omega_0 + \hbar \Omega + \frac{\Delta}{2} \cos k_z d, \quad (23)$$

$$a_{\pm}^2 = \frac{1}{d^2} \left( k_z d - \cos^{-1} \left( \frac{2}{\Delta} \Theta_{\pm}(k_z d) \right) \right)^2,$$

When deriving formula (22), it was taken into account that  $\int_0^{\infty} \frac{|I_{nn'}(u)|^2 du}{u} = \frac{1}{n' - n}$ .

### 3. RESULTS

From expression (22) it is seen that the absorption coefficient diverges whenever the condition  $1 - 4\Theta_i^2/\Delta^2 = 0$  is satisfied. In addition, from the fact that  $1 - 4\Theta_i^2/\Delta^2 = 0$  are real and positive, we can obtain an energy range in which the absorption coefficient are allowed.

As follows from equations (21-22), peaks are observed at certain frequencies of the incident photon. The resonance behavior of the absorption coefficient for electron-phonon scattering appears for frequencies satisfying the relation

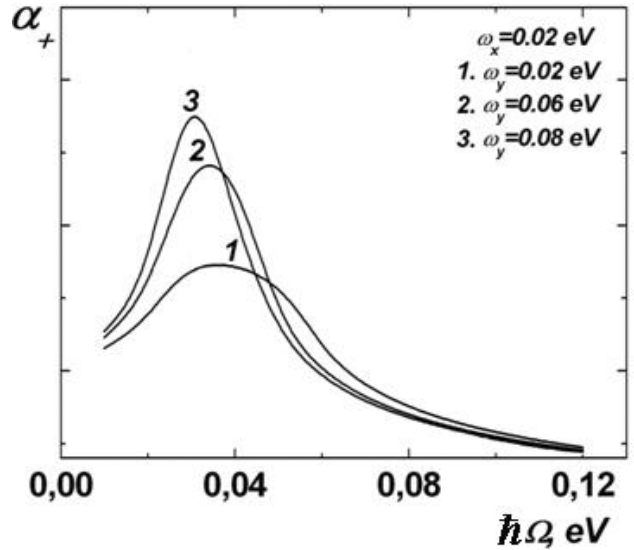
$$N \hbar \omega_x + P \hbar \omega_y \pm \hbar \omega_0 = \hbar \Omega, \quad (24)$$

here  $N = n' - n = 1, 2, 3, \dots$ , and  $P = l' - l = 1, 2, 3, \dots$

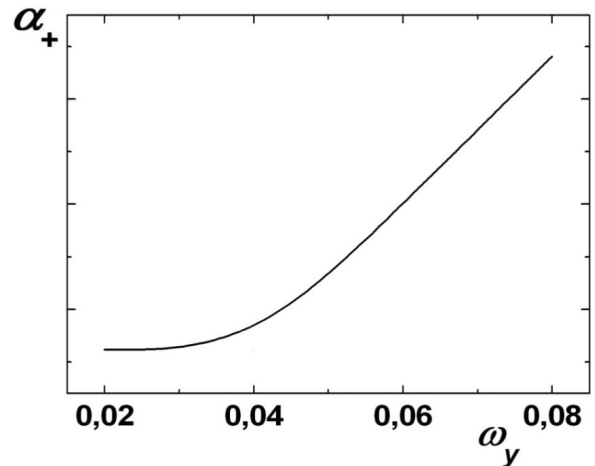
From Eq. (24), in the course of scattering events, the electrons in the subband levels specified by the level index  $n(l)$  can make transitions to one of the subband levels  $n'(l')$  by absorbing and/or emitting a photon of energy  $\hbar \Omega$  during the absorption of a LO phonon of energy  $\hbar \omega_0$ .

Eq. (24) is the basic equation for the absorption coefficient spectral lineshape, which enable us to analyze resonant effects in semiconductors.

In **Fig. 1**  $\alpha_+$  dependence on  $\hbar \Omega$  for  $\hbar \omega_x = 0.03 \text{ eV}$ , and  $\hbar \omega_y = 0.02 \text{ eV}, 0.06 \text{ eV}, 0.08 \text{ eV}$  is given. **Fig. 2** is presented  $\alpha_{+max}$  dependence on  $\omega_y$  for  $\Delta = 0.0024 \text{ eV}$ .



**Fig. 1.**  $\alpha_+$  dependence on  $\hbar \Omega$  at  $\hbar \omega_x = 0.03 \text{ eV}$ , and  $\hbar \omega_y = 0.02 \text{ eV}, 0.06 \text{ eV}, 0.08 \text{ eV}$ .



**Fig. 2.**  $\alpha_{+max}$  dependence on  $\omega_y$  at  $\Delta = 0.0024 \text{ eV}$ .

As seen, with increasing  $\omega_y$  the  $\alpha_{+max}$  grows. The  $\omega_y$  values, affecting the change of  $\alpha_{+max}$ , as well as the frequency at which it is achieved ( $\hbar \Omega$ ), shift towards small frequencies.

### 4. CONCLUSION

In conclusion, we calculated free-carrier absorption of electromagnetic radiation with polar and non-polar optical phonons. It should be noted that the summations of Eq. (21) over the size subband levels contain three types of contributions: (i)  $n' \neq n, l' = l$ , (ii)  $n' = n, l' \neq l$ , and (iii)  $n' \neq n, l' \neq l$ . From the selection rules, we can expect three possible transitions in the absorption coefficient for electron-polar phonon scattering: (1) a transition due to the size subband levels for only the  $x$ -direction, (2) a transition due to the size subband levels for only the  $y$ -direction

and (3) a transition due to the size subband levels for both the  $x$ -direction and the  $y$ -direction.

From the Eq. (22) we can expect only one possible transitions in the, the absorption coefficient for electron-nonpolar phonon scattering: a transition due to the size subband levels for both the  $x$ -direction and the  $y$ -direction.

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