



## OPTICAL ALIGNMENT OF 2D-ELECTRON MOMENTA IN MULTIPLE QUANTUM WELL STRUCTURES

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A linear polarization of the recombination luminescence of photoexcited hot electrons in a multiple quantum well structure has been observed. The peculiarity of electron 2D-motion manifests itself in an observed increase of the polarization with 2D-motion energy. The time of optical-phonon emission by a 2D electron is determined by measuring a magnetic depolarization of the luminescence.

### 1. Experimental

Spectra and polarization of hot photoluminescence in a quantum-well structure have been first studied in this work. The results are presented for a p-GaAs/Al<sub>0.3</sub>Ga<sub>0.7</sub>As structure grown by molecular-beam epitaxy and composed of 100 periods of 70 Å GaAs quantum wells and 100 Å GaAlAs barriers. The acceptor (carbon) concentration in the region of quantum wells was of the order of 10<sup>16</sup> cm<sup>-3</sup>.

The photoluminescence spectrum under Kr-laser  $E_{ex} = 1.65$  eV excitation and directions of the exciting light and observation normal to the plane of the structure, is shown in Fig.1. For such a value of  $E_{ex}$ , the electrons are excited only into the first electron subband (see the scheme of the transitions in Fig.1). In the luminescence spectrum the 1e - 1hh excitonic transition as well as that from the first electron subband onto the acceptor level (1e - A) are revealed. The high-frequency tail of this latter component of the spectrum is just due to the recombination of hot electrons on the acceptor levels. The peak designated as "0" on the high-frequency edge of the spectrum corresponds to recombination of the electrons from the point of photocreation, i.e. prior to any energy relaxation. The kinetic energy of electrons in this point is  $E_0 = 60$  meV. The degree of linear polarization of the radiation from the point "0" measured under linearly polarized excitation has the value  $P_1 = 0.06$

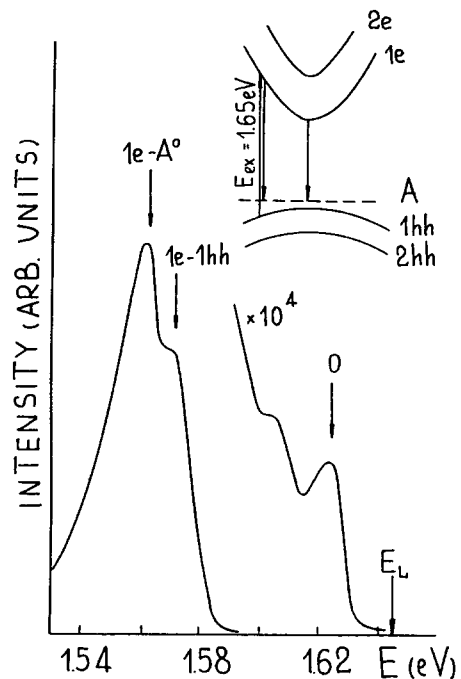


Fig.1. The photoluminescence spectrum of the quantum well structure (the parameters of the structure are given in the text,  $E_{ex} = 1.65$  eV). The right part of the curve is the spectrum of hot luminescence. "0" designates the zero-phonon peak recombination from the point of photocreation. The scheme of the transitions is in the right upper corner.

When  $E_{ex}$  is increased,  $P_1$  increases too amounting to 0.28 at  $E_{ex}=1.83$  eV (which corresponds to  $E_0 = 215$  meV). As seen in Fig.2, the degree of linear polarization decreases in the course of energy relaxation.

2. Discussion

The obtained results are explained by an optical alignment of electron momenta under linearly polarized excitation from the valence band of the GaAs-type semiconductors. The electrons created in the conduction band have anisotropic distribution of momenta. Let us first remind the result for the 3D-case [1,2]. For the sake of simplicity, we will use a spherical approximation for the valence band. We will also consider the acceptor wave function as formed mainly from the states of the heavy-hole subband [2]. The calculation of matrix elements of the dipole momentum shows that when excited from the heavy-hole subband, the electrons are created with the momenta  $\vec{p}$  directed preferentially perpendicular to the electric vector  $\vec{e}$  of the exciting light. The distribution function of these electrons is

$$F(p) = f(p)(p_y^2 + p_z^2), \quad (1)$$

where  $f(p)$  is a spherically symmetric function, the exciting light is directed along Z axis and  $\vec{e}$  along X axis.

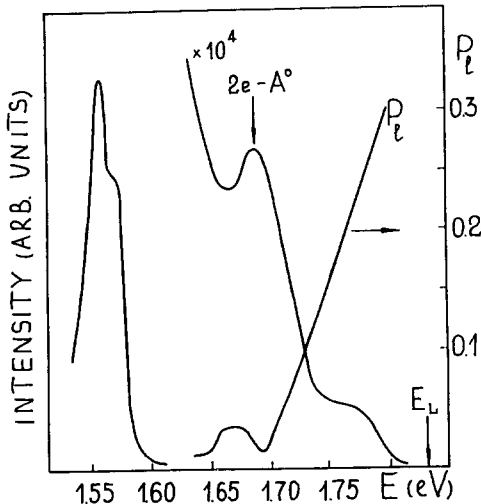


Fig.2. The spectra of photoluminescence intensity and linear polarization for the same structure as in Fig.1.  $E_{ex} = 1.83$  eV. The minimum on  $P_1$  curve is due to contribution of weakly polarized transitions.

A transition probability for recombination has the same matrix elements as for excitation. Consequently the intensity of the radiation polarized along the X and Y axes is respectively

$$I_x \sim \int F(p)(p_y^2 + p_z^2) d\vec{p}, \quad (2)$$

$$I_y \sim \int F(p)(p_x^2 + p_z^2) d\vec{p}.$$

The integration in the 3D-case must be performed in p-space over the angles  $\psi$  and  $\varphi$  ( $p_x = p \sin \psi \cos \varphi$ ,  $p_y = p \sin \psi \sin \varphi$ ,  $p_z = p \cos \psi$ ), because to the fixed energy of the recombination quantum corresponds a fixed value of  $p^2$ .

Thus, for the recombination from the conduction band into the band of heavy holes or on the acceptor levels it is easy to get the degree of polarization  $P_1 = (I_x - I_y) / (I_x + I_y) = 1/7$ .

Note that  $P_1$  does not depend on the energy of the recombining electrons.

In the 2D-case one must take into account a dimensional quantization of electron and hole energies. In a very simplified model, in which the  $p_z$  component of the momentum is considered as fixed and is determined by energy  $E_1$  of the bottom of the 1e subband

$$(p_z^2 = 2m_c E_1),$$

the integration in Eq. (2) must be performed only over  $\psi$ . Then we have for the degree of polarization under the excitation from the subband of heavy holes

$$P_1 = \frac{(E_0/E_1)^2}{8(1+E_0/E_1) + 2(E_0/E_1)^2}, \quad (3)$$

where the energy of the in-plane motion of the electrons in the point of photocreation is  $E_0 = (p_x^2 + p_y^2) / 2m_c$ .

Equation (3) shows that in the 2D-case, contrary to the 3D-one, there exists an essential dependence of  $P_1$  on  $E_0$ .  $P_1$  increases from 0 to 0.5 with an increase of  $E_0$ . (For electrons excited from the subband of light holes and recombining on the acceptor levels,  $P_1$  changes in the range from 0 to -0.3). Such a dependence of  $P_1$  on  $E_0$  results from the fact that at small  $E_0$  the electrons are moving nearly normal to the plane of the structure and consequently for the direction of observation along Z their radiation is almost unpolarized. For  $E_0 \gg E_1$ , the

movement of the electrons proceeds mainly in-plane and recombination radiation is strongly polarized.

The increase of  $P_1$  with the increasing electron energy  $E_0$ , observed in the experiment, is consistent with the presented ideas. For a studied structure,

the value  $E_0 = 60$  meV corresponds to  $E_0/E_1 = 1.2$  and in accordance with Eq. (3) to  $P_1 = 0.075$  (the measured value was  $P_1 = 0.06$ ). The value  $E_0 = 215$  meV corresponds to  $E_0/E_1 = 4.3$  and the calculated and measured values of  $P_1$  were 0.23 and 0.28 respectively. The minimum on the spectral dependence of  $P_1$  in Fig. 2 is due to the contribution of a weakly polarized transitions from the bottom of the second electron subband ( $2e - \bar{A}$ ).

3. Depolarization in a magnetic field and determination of  $\tau_{p0}$

The study of the depolarization of hot luminescence in a magnetic field makes it possible to determine characteristic relaxation times of hot electrons [2]. The polarization of recombination luminescence in a magnetic field  $\vec{B}$  directed along the light beam (Faraday configuration) is

$$P_1(B)/P_1(0) = (1 + 4 \omega_c^2 \tau_0^2)^{-1/2}, \quad (4)$$

where  $\omega_c$  is the cyclotron frequency, and  $\tau_0$  is the "life time" of electrons in the point of photocreation  $E_0$ . At moderate doping level and pumping density  $\tau_0$  has the meaning of the LO-phonon emission time  $\tau_{p0}$ . Indeed, the distinct peak "0" observed in the spectrum (Fig.1) confirms that energy relaxation occurs by LO-phonon emission. It follows from  $P_1(B)$  measurements at  $E_0 = 60$  meV (Fig.3) that  $\tau_{p0} = 150 \pm 10$  fs. This value agrees with the probability of intraband transitions which was calculated in this work. The applied technique can be used in a wide range of electron energies to study kinetics of intra- and interband transitions in quantum wells. This technique is not influenced by the essentially complicating factors (such as, for instance, phonon heating), which were a reason of large spread in  $\tau_{p0}$  values determined in earlier works on 2D-structures (see, for instance, reviews [3,4]).

4. Luminescence under circularly polarized excitation

We would like to draw attention to some peculiarities of the 2D-case under excitation with circularly polarized light. Let us remind that in the 3D-case in GaAs-type semiconductors the momenta and spins of the electrons created via interband transitions are correlated [1]. The component of the density matrix responsible for the correlation is proportional to  $(\hat{\sigma}_p)(\hat{n}_p)$ , where  $\hat{\sigma}$  are Pauli matrices and  $\hat{n}$  is the

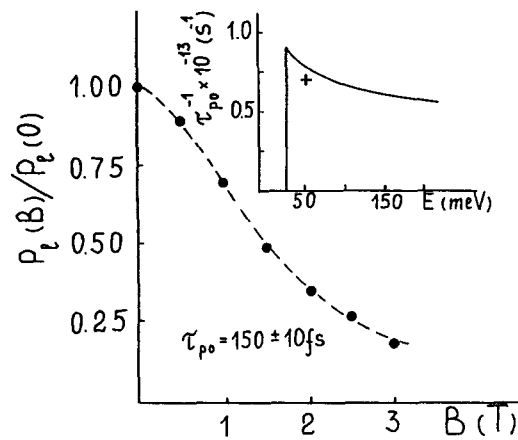


Fig.3. The depolarization of hot photoluminescence in a magnetic field is described by the Lorentz contour (4) with  $\tau_{p0} = 150$  fs. The calculated energy dependence  $\tau_{p0}^{-1}(E_0)$  for the studied structure is shown in the right upper corner, the experimental value corresponding to  $E_0 = 60$  meV is shown by a cross.

unit vector in the direction of the light beam. Such a correlation does not exist in the 2D-case. This is due to the fact that after averaging over the motion along Z the nondiagonal elements of the matrix  $(\hat{\sigma}_p)(\hat{n}_p)$  become zero. A very convincing magneto-optical manifestation of the correlation in the 3d-case is the decrease of circular polarization of the luminescence in a longitudinal magnetic field due to the destroying of the aforementioned correlation [5]. In the 2d-case we did not observe any decrease of circular polarization for magnetic fields as high as 7T, which is a direct evidence of the absence of the correlation.

Let us also attract attention to the dependence of the degree of circular polarization,  $P_c$ , on the energy of recombining electrons. Within the framework of approximation used in p.2 the mean spin of the electrons photo-excited from the subband of heavy holes is

$$S = -\hat{n}(2 + E_0/E_1)^{-1}.$$

And the degree of polarization for recombination on the acceptor levels is

$$P_c = -\hat{S}\hat{n} = (2 + E_0/E_1)^{-1}.$$

Thus, under excitation close to the bottom of the first electron band the photocreated electrons have com-

pletely oriented spins and  $P_0=0.5$ . As  $E_0$  increases, the degree of polarization decreases from 0.5 to 0. We have observed the decrease of  $P_1$  from 0.3 at  $E_0 = 60$  meV to 0.22 at  $E_0 = 200$  meV.

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