

## Electron delocalization in the 2D electron gas at a silicon (100) surface

V. T. Dolgoplov, A. A. Shashkin, N. B. Zhitenev, and S. I. Dorozhkin  
*Institute of Solid State Physics, Academy of Sciences of the USSR*

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The number of localized electrons,  $N_l$ , has been studied experimentally as a function of the total concentration  $N_s$  in metal-insulator-semiconductor structures on the (100) face of silicon. The number  $N_l$  is found to decrease with increasing  $N_s$ , because of an improved screening of the potential relief by free electrons.

In an ideal 2D electron system the state density  $D(E)$  would not depend on the energy. In a real system, in contrast, there are always fluctuations of the potential which result in a localization of electrons with energies below some value  $E_c$  (Ref. 1). In this case the state density has a tail which falls off exponentially at low energies

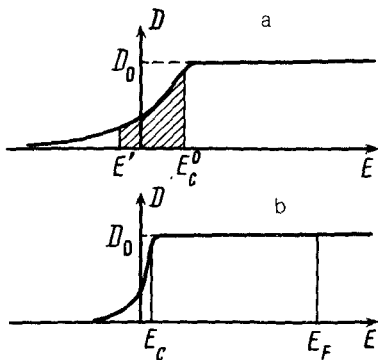


FIG. 1. Schematic diagram of the energy dependence of the state density. a—The Fermi level is the region of localized states; b—it is in the region of delocalized states. The hatching shows the region of delocalized electrons.

( $E < E_c$ ; Fig. 1). The existence of a tail in the state density has been verified experimentally in measurements of the activation energy ( $E_c - E_F$ ) as a function of the concentration ( $N_s$ ) of 2D electrons on low densities<sup>2,3</sup> ( $E_F < E_c$ ) and in optical experiments.<sup>4</sup> At comparatively high electron concentrations ( $E_F > E_c$ ), no evidence of any sort for a tail on the state density has so far been found.<sup>5</sup>

In the present letter we report an experimental study of the number of localized electrons,  $N_l = \int_{-\infty}^{E_c} D(E)dE$ , as a function of the concentration  $N_s$  in the 2D electron gas near the (100) surface of silicon at  $E_F > E_c$ . The results show that with increasing concentration of the 2D electrons, the number of localized electrons falls off, and the  $D(E)$  curve approaches the shape we would expect for an ideal system.

We measured the magnetotransport characteristics of samples with a high mobility  $\mu$  at magnetic fields  $\mu H \lesssim 1$ . We studied the dependence  $\sigma_{xx}(H)$  in a sample in the Corbino geometry. Using the Drude formula

$$\sigma_{xx} = \frac{\sigma_0}{1 + (\mu H)^2}, \quad (1)$$

whose applicability to our sample was verified over a broad range of magnetic fields and concentrations, we found  $\mu(N_s)$ . Assuming  $\sigma_0 = (N_s - N_l)e\mu$ , we then determined  $N_l(N_s)$ . The total electron concentration is found independently, from measurements of the capacitance of the structure and the threshold voltage  $V_T$ , determined from the positions of the minima of Shubnikov-de Haas oscillations in magnetic fields 5–10 T. The dependence  $N_l(N_s)$  found in this manner is shown in Fig. 2 for concentrations corresponding to  $\sigma > 10\sigma_{\min}$ .

Measurements in the Corbino geometry are carried out by a two-contact method, so that there is always the danger of finding an effect similar to that described above but resulting from a resistance of the contacts. In our case, the resistance of the contacts was apparently negligible, since the experimental curve of  $\sigma_{xx}(H)$  is described well by expression (1). We also made an attempt to carry out similar measurements in a sample in the Corbino geometry, whose conductivity was measured by a capacitance-coupling technique.<sup>6</sup> We were not able to measure an  $N_l(N_s)$  curve for this sample, since the accuracy of the method for measuring  $\sigma_{xx}$  falls off with increas-

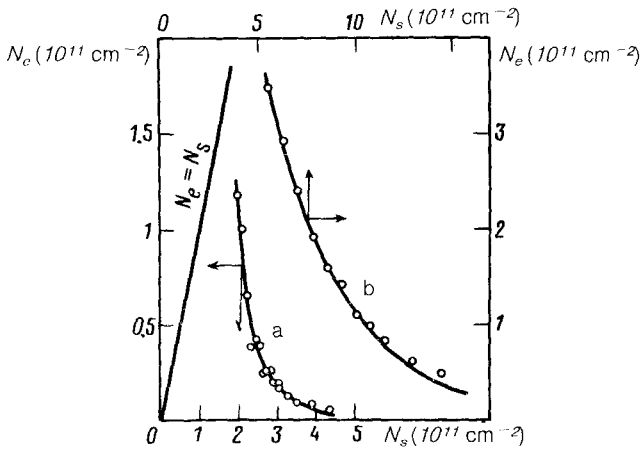


FIG. 2. Number of localized electrons,  $N_l$ , versus the total number of electrons in the inversion layer,  $N_s$ . *a*—Sample in the Corbino geometry [the maximum mobility  $\mu_{\max} = 20\,400\text{ cm}^2/(\text{V} \cdot \text{s})$  is reached at  $N_s = 8.6 \times 10^{11}\text{ cm}^{-2}$ ,  $T = 1.46\text{ K}$ , and  $V_{SB} = 0$ ]; *b*—Hall transistor [ $\mu_{\max} = 21\,000\text{ cm}^2/(\text{V} \cdot \text{s})$  at  $N_s = 4.5 \times 10^{11}\text{ cm}^{-2}$ ,  $T = 1.42\text{ K}$ , and  $V_{SB} = 0$ ].

ing conductivity. However, again in the case of this sample, the electron concentration calculated from the conductivity  $\sigma_0$  turned out to be lower than  $N_s$  at  $\sigma \gtrsim 10\sigma_{\min}$ , and the difference in concentrations was clearly greater than the experimental error.

The resistance of the contacts plays no role in measurements by a four-contact method. We used a four-contact bridge arrangement to carry out measurements with Hall transistors.<sup>7</sup> The results on  $N_s - N_l = H(ec\rho_{xy})^{-1}$  as a function of the transistor gate voltage from these measurements are shown in Fig. 3; the corresponding curve for  $N_l(N_s)$  is shown in Fig. 2.

Figure 3 also shows the positions of the minima of the quantum oscillations of  $\rho_{xx}$

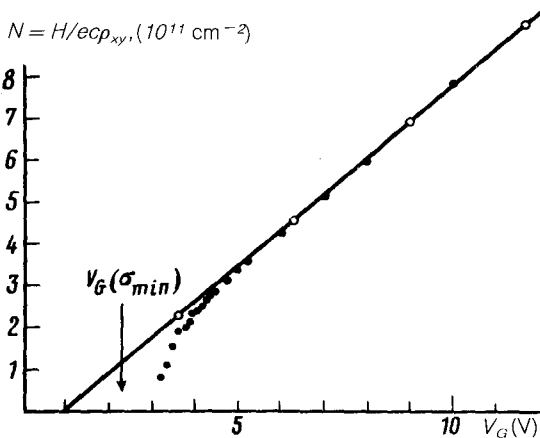


FIG. 3. Filled circles—measurements of  $N = H/ec\rho_{xy}$  as a function of the gate voltage in magnetic fields  $\mu H \lesssim 1$ ; open circles—concentrations determined from the positions of the minima of the quantum oscillations in a magnetic field  $H = 4.77$  at  $T = 1.42\text{ K}$ .

in a magnetic field  $H = 4.77$  T. We see from this figure that the positions of the minima of the Shubnikov-de Haas oscillations are determined by the total electron concentration  $N_s$ . We obtained a similar result for a sample in the Corbino geometry.

The decrease in the number of localized electrons with increasing  $N_s$  results from an improvement of the screening. The effective screening parameter  $q_s$  ( $q$ ) is constant at low temperatures, with  $q \ll 2k_F$ , and it falls off sharply at  $q > 2k_F$ , so that wells with different scale dimensions  $a$  are screened at different concentrations of free electrons. Let us assume that the width distribution of the wells is described by<sup>5</sup>

$$a = a_0 \left( \frac{E_c^0}{E_c^0 - E} \right)^s \quad (1)$$

[the origin on the energy scale is chosen in such a way that we have  $N_l = \int_{-\infty}^{E_c} D(E) dE = D_0 E_c D_0 = g_v m / \pi \hbar^2$ ] and that the state density is  $D = D_0 \exp((E - E_c)/E_c)$  if  $E < E_c$ . Upon the appearance of free electrons with a wave vector  $k_F$ , all the wells with  $a > k_F^{-1}$  are screened. Electrons with energies  $E' < E < E_c$  ( $k_F = 0$ ) =  $E_c^0$  become delocalized, and the energy  $E_c$  becomes dependent on the electron concentration (Fig. 1). We then have

$$E_F - E_c(N_s) = \frac{\hbar^2}{2ma_0^2} \left( \frac{E_c^0 - E'}{E_c^0} \right)^{2s}, \quad (2)$$

$$D_0(E_c^0 - E_c) = D_0 \int_{E'}^{E_c^0} dE \exp\left(\frac{E - E_c^0}{E_c^0}\right). \quad (3)$$

The first of these relations determines the Fermi energy  $E_F$ , while the second corresponds to the case in which the number of delocalized electrons is equal to the change in the number of free electrons. Eliminating  $E'$  from (2) and (3), we find

$$N_l(N_s) = D_0 E_c(N_s) = D_0 E_c^0 \exp[-\{\pi a_0^2 (N_s - N_l)\}^{1/2s}]. \quad (4)$$

For our samples, we find the best agreement with the experimental data with the values  $s = 0.35 \pm 0.05$  and  $a_0 = 66$  Å for a sample in the Corbino geometry or with  $\alpha_0 = 130$  Å for the Hall transistor (solid lines in Fig. 2). The values found for  $s$  by Pollitt<sup>3</sup> range from 1 to 0.75, depending on the difference  $E - E_c$ . The improvement in the screening by free electrons thus makes it possible to explain the observed dependence  $N_l(N_s)$  with completely reasonable values of the parameters  $s$  and  $a_0$ .

The difference between the mobilities  $\mu$  and  $\mu_H$ , which can be seen in Fig. 1 of Ref. 9, and which disappears with increasing  $N_s$ , can apparently be explained in terms of the presence of localized electrons. The temperature dependence of  $\rho_{xy}$  found in the 2D electron gas on <sup>10</sup>Si(111) is probably also due to an effect of this sort, since the corrections to  $\rho_{xy}$  for the electron-electron interaction in the dirty limit<sup>11</sup> vanish in the limit  $k_F q_S(0) \rightarrow 0$ . [The parameter  $k_F q_S(0)$  in Ref. 10 was  $\sim 0.1$ , and even smaller in our own case.]

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