Quantum Hall Transition in Real Space: From Localized to Extended States

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Using scanning tunneling spectroscopy in an ultrahigh vacuum at low temperature (T = 0.3 K) and high magnetic fields ($B \le 12$ T), we directly probe electronic wave functions across an integer quantum Hall transition. In accordance with theoretical predictions, we observe the evolution from localized drift states in the insulating phases to branched extended drift states at the quantum critical point. The observed microscopic behavior close to the extended state indicates points of localized quantum tunneling, which are considered to be decisive for a quantitative description of the transition.

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Two-dimensional electron systems (2DES) are paradigms for quantum phase transitions [1]. For example, they exhibit metal-insulator, integer quantum Hall (IQH), and superconductor-insulator transitions [2], all described fairly successfully by a universal percolation-type model [3]. Some macroscopic aspects of this percolation have been verified, e.g., by transport experiments revealing critical exponents for the divergence of the localization length [4]. However, the most fundamental aspect, i.e., how wave functions change at the transition, has not been probed experimentally. Here, we directly observe the wave functions at the IQH transition.

The microscopic description of IQH transitions is theoretically well established [5-7]. The 2DES in magnetic B fields exhibits discrete kinetic energies which are called Landau levels (LL). The corresponding states are subject to random potential disorder, which has primarily a semiclassical effect [5]. Quasiclassically, the electrons perform a fast cyclotron rotation within the electrostatic disorder which leads to additional drift motion along the equipotential lines [8]. Quantum mechanically, so-called drift states meander along equipotential lines with a width of about the cyclotron radius r_c [6]. If the potential energy of the state is low (high), the drift states are closed trajectories around potential minima (maxima); i.e., they are localized and represent insulating electron phases. In the center of a LL, the adjacent trajectories merge at the saddle points of the potential leading to an extended state traversing the whole sample. It is known that this state is the quantum critical state of the IQH transition and responsible for the finite longitudinal resistance between quantized values of the Hall conductance [5-7,9].

Several experiments have addressed microscopic aspects of QH transitions [10,11]. However, none of these techniques revealed sufficient lateral resolution to image the drift states. Only scanning tunneling spectroscopy (STS) using the adsorbate-induced 2DES at *n*-InAs(110) surfaces found localized drift states within a LL tail [12]. This paves the path to observe the decisive transition from localized to extended states directly. Localized states in *B* field have also been observed on graphite, which, however, does not exhibit a pure 2DES [13].

Here, we probe the local density of states (LDOS) of the adsorbate-induced 2DES on *n*-InSb(110) by STS. The 2DES is prepared by depositing 0.01 monolayer of Cs atoms on cleaved *n*-InSb(110) at temperature T = 30 K and pressure $p \sim 10^{-10}$ mbar [14,15]. STS is performed *in situ* at T = 0.3 K and $B \le 12$ T [16] using a W tip carefully selected by trial and error to exhibit negligible tip-induced band bending [14,17]. The tip is stabilized at current I_{stab} and voltage V_{stab} , prior to measuring dI/dV as a function of sample voltage V_s directly by lock-in technique with modulation voltage V_{mod} .

First we sketch the general properties of the 2DES. Figure 1(a) shows a color scale plot of the *local* differential conductivity $dI/dV(V_s)$ of the 2DES measured along a straight line at B = 0 T and the corresponding spatially averaged dI/dV curve. They, respectively, represent the energy dependence of the LDOS (energy-position plot) and of the macroscopically averaged LDOS, i.e., the DOS [18]. The LDOS of the 2DES shows two apparent boundaries coinciding with two steplike features in the DOS at $V_s =$ -115 and -47 meV, which represent the first (E_1) and second (E_2) subband edges as indicated in Fig. 1(a). The



FIG. 1 (color). (a) Color scale plot of $dI/dV(V_s)$ measured along a straight line at B = 0 T (left); corresponding spatially averaged $dI/dV(V_s)$ (right); dashed lines mark the Fermi level E_F and the subband energies E_i (i = 1, 2); $V_{stab} = 150$ mV, $I_{stab} = 0.13$ nA, $V_{mod} = 2$ mV. (b) Self-consistently calculated band bending at the InSb surface (black line) and electrondensity distribution of *i*th subband (red areas); E_i and electron areal densities n_i are marked. (c) $dI/dV(V_s)$ at different *B* recorded at the same lateral position; E_i, E_F , (dI/dV = 0) lines on the right, LLs and spin directions are marked; $V_{stab} =$ 150 mV, $I_{stab} = 0.13$ nA, $V_{mod} = 2.0$ mV (0 T), 1.3 mV (6 T), 1.0 mV (9 T), 0.9 mV (12 T). (d) Experimentally determined Landau fan diagram; green (red) dashed lines: spin-down LLs of the 1st (2nd) subband; arrows: spin levels of the lowest LL; $V_{stab} = 150$ mV, $I_{stab} = 0.10$ nA, $V_{mod} = 1.5$ mV.

subband energies are excellently reproduced by a selfconsistent calculation [Fig. 1(b)] [14,19]. The spatial irregularity of the onset at E_1 in Fig. 1(a) (left) is a signature of the potential disorder [18]. Figure 1(c) shows a set of dI/dV curves measured at the same position at different B. Starting at B = 6 T, the dI/dV curves exhibit distinct LLs with a pronounced twofold spin splitting. The LLs are separated by regions of $dI/dV \approx 0$ evidencing complete quantization. Indeed, spin-resolved IQH plateaus up to filling factor 6 were recently observed by magnetotransport on an adsorbate-induced 2DES on InSb(110) [20]. Repeating the measurement of Fig. 1(c) using smaller B-field steps highlights the continuous evolution of the spin-split LLs, i.e., the LL fan diagram in Fig. 1(d). The green (red) dashed lines mark the four (two) spin-down LLs of the 1st (2nd) subband. The accompanying spin-up LLs are visible at higher energies as marked by blue arrows for the lowest LL. Partly, LLs of different subbands cross without anticrossing indicating orthogonality and, thus, negligible interaction between the subbands. From the peak distances, we deduce the effective mass m^* and the absolute value of the *g* factor |g| via the separation of LLs $\hbar |e|B/m^*$ (\hbar , Dirac's constant; *e*, electron charge) and of spin levels $|g|\mu_B B$ (μ_B , Bohr magneton). For the lowest peaks at B = 6 T, we find $m^*/m_e = 0.019 \pm 0.001$ (m_e , free-electron mass) and $|g| = 39 \pm 2$. This is close to the known values at the band edge $m^*/m_e = 0.014$ and |g| = 51 with slight deviations due to nonparabolicity and energy dependent spin-orbit coupling as known for InSb [21,22].

Figures 2(a)-2(g) show our central result, the LDOS across an IQH transition, i.e., dI/dV images recorded in the lowest spin-down LL of the 2DES (LL0 \downarrow) at B =12 T. The corresponding, spatially averaged dI/dV curve is shown in Fig. 2(h). The continuous change of the LDOS with energy is available as a movie in [14]. The LDOS in the low-energy tail of $LL0 \downarrow$ is shown in Fig. 2(a). It exhibits spatially isolated closed-loop patterns with averaged FWHM $\simeq 6.9$ nm close to $r_c = 7.4$ nm. Thus, we attribute the closed patterns to localized drift states of LL0 | aligning along equipotential lines around a potential minimum. Accordingly, at slightly higher energy [Fig. 2(b)], the area encircled by the drift states increases indicating that the drift states probe a longer equipotential line at higher energy within the same valley. In contrast, the ring patterns at the high-energy tail, marked by green arrows in Figs. 2(f) and 2(g), encircle an area decreasing in size with increasing voltage. These states are attributed to localized drift states around potential maxima. Notice that the structures in Figs. 2(a) and 2(b) appear nearly identically in Figs. 2(f) and 2(g) as marked by white arrows. The latter structures are the LL0 [↑] states localized around potential minima, which energetically overlap with the high-energy LL0 | states localized around potential maxima. When the voltage is close to the LL center [Figs. 2(c) and 2(e)], adjacent drift states coalesce and a dense network is observed directly at the LL center [Fig. 2(d)]. This is exactly the expected behavior of an extended drift state at the IOH transition [6,7]. Figure 2(i) shows a calculated extended state at B = 12 T using the parameters of a 2DES of InSb [23,24] and the potential disorder provided by the known dopant density of the sample [14]. Good qualitative agreement with the measurement is achieved supporting the interpretation of the coalesced LDOS patterns as extended states. Figure 2(j) shows another extended state recorded on a larger area at different B demonstrating that the coalesced pattern is not restricted to small length scales.

Interestingly, the drift states around the LL center indicate quantum tunneling at the saddle points. Within the classical percolation model, the adjacent drift states are connected at a singular energy at each saddle eventually leading to a localization exponent $\nu = 4/3$. However, experimental [4] and numerical [7] results for IQH tran-



FIG. 2 (color). LDOS of lowest LL. (a)–(g) Measured dI/dV (x, y); B = 12 T, $V_s = -116.3$ mV (a), -111.2 mV (b), -104.4 mV (c), -100.9 mV (d), -99.2 mV (e), -92.4 mV (f), -89.0 mV (g), $I_{stab} = 0.1$ nA, $V_{stab} = 150$ mV, $V_{mod} = 1.0$ mV; same dI/dV color scale in each image; white (green) arrows in (a), (b), (f), and (g) mark drift states encircling potential minima (maxima); red, yellow arrows in (c) and (e) mark tunneling connections existing at identical positions; crosses in (d) mark extended LDOS areas at saddle points. (h) Spatially averaged dI/dV curve with circles at the V_s used in (a)–(g). (i) Calculated LDOS at the center of LL0 \downarrow at B = 12 T; red arrows mark tunneling connections at the saddle points; white crosses mark extended areas. (j) dI/dV image close to the center of LL0 \uparrow at B = 6 T, $V_s = -99$ mV; image includes the area of (a)–(g) within the marked rectangle.

sitions revealed $\nu \approx 2.3$. The discrepancy is attributed to quantum tunneling between classically localized drift states [6,7]. Tunneling connections are indeed visible in our data spreading over an energy range of 5 meV. As an example, the red (yellow) arrows in Figs. 2(c) and 2(e) mark the same connection point at $V_s = -104.4$ mV and -99.2 mV. LDOS is faintly visible at both positions in both images and sharply rotates by about 90° between the images. The reason is simply that the tunneling interconnection mediates between valley states at low energies and between hill states at high energies, which are connected via two nearly orthogonal lines. Such weak links are reproduced by the calculation as marked by red arrows in Fig. 2(i) (see also Fig. S3 of [14]). Note that the intrinsic energy resolution of the experiment is 0.1 meV [16], while peaks in the LL fan diagram exhibit a FWHM of 2.5 meV probably due to lifetime effects. Thus, broadening due to the energy resolution can hardly account for intensity at the saddles within an energy range of 5 meV. Another intriguing observation is the LDOS areas larger than r_c around the saddles. They are again visible in experiment [crosses in Fig. 2(d)] and calculation [crosses in Fig. 2(i)] and are probably due to the flat potential at the saddles leading to slow drift speed and, thus, extended LDOS intensity. Notice that the spreading of LDOS intensity at the saddles in energy and position is consistent with previous quantum mechanical calculations [6,7].

Finally, we discuss the possible influence of the tip. It is known that a mismatch of tip and surface potential leads to band bending within the sample [17,18]. To avoid this, we used only W tips exhibiting a minimum of tip-induced band bending. By analyzing the dI/dV data at B = 0 T with and without adsorbates, we can safely rule out a work function mismatch between tip and sample larger than 15 meV [14,17,18]. The applied V_s leads to an additional tip-induced band bending with a lever arm of 10 as determined from experiments described in [14]. Thus, we get an additional band bending of less than 12 meV. The influence of such small band bendings is tested by the theoretical calculations of a disordered 2DES in *B* field [14,23,24]. The tip-induced potential is added to the disorder potential of the 2DES as a Gaussian with 50 nm FWHM and amplitude $|V_{tip}| < 20$ meV [17,18]. The LDOS is calcu-



FIG. 3 (color online). (a) Calculated LDOS (*E*, *x*) at B = 6 T; $V_{\text{tip}} = 0$ and ± 20 meV as marked; $m^*/m_e := 0.02$, |g| := 28[23] (b), (c) Measured $dI/dV(V_s, x)$ at B = 6 and 12 T; *x*: lateral position along straight line; $I_{\text{stab}} = 0.13$ nA, $V_{\text{stab}} = 150$ mV, $V_{\text{mod}} = 1.3$ mV (b), 0.9 mV (c). (d),(e) Calculated LDOS (*E*, *x*) at B = 6 and 12 T; $V_{\text{tip}} = 0$ meV.

lated at each lateral tip position and its intensity below the tip is displayed as a function of tip position. Figure 3(a) shows the resulting energy-position plots for three different V_{tip} at B = 6 T. Obviously, V_{tip} leads only to a rigid energetic shift of the patterns. This result is robust to changes in *B* and FWHM of the tip-induced potential. Thus, the spatial dependence of the LDOS patterns is not influenced by a small potential mismatch $|V_{\text{tip}}| < 20$ meV.

To substantiate this result, Figs. 3(b)-3(e) show experimentally determined voltage-position plots in comparison with calculated energy-position plots [14,23] at B = 6 and 12 T. Spin-split LLs fluctuating in energy are visible as meandering pairs of lines. The theoretical calculations do not require free parameters by neglecting the tip-induced potential. They reproduce the experimental trends of spatially fluctuating spin-resolved LLs. The fluctuation length is 80% larger and the fluctuation amplitude 40% lower than in the experiment. This is most likely due to the neglected inhomogeneous charge distribution of the adsorbates, which would increase the potential disorder given by the bulk dopants. Indeed, as shown in Fig. S1 of [14], the spatial distribution of adsorbates is related to the resulting drift states, and the correspondence is improved by adding surface charges to the calculation (not shown). However, since many details of the adsorbate potential are not known, we omit its inclusion, thereby avoiding the use of further parameters. Note that the parameter-free calculation and the experiment quantitatively exhibit the same trends with B field and energy, i.e., a 30% increase of fluctuation amplitude between 6 and 12 T and a 60% decrease of fluctuation amplitude between LL0 and LL3 at B = 6 T. Both effects are explained straightforwardly by the fact that drift states can only probe potential fluctuations down to length scales of $r_c = \sqrt{(2n+1)\hbar/(|e|B)} (n,$ LL index).

In summary, our STS experiments provide the first direct observation of wave functions across an IOH transition. They reveal the development from localized to extended states including indications of the theoretically predicted quantum tunneling at saddle points of the disorder potential. We probe states well away from the Fermi level, and thus artificially exclude the influence of electron-electron interaction, thereby considering a theoretically pure IQH transition [6-9]. Indeed, single-particle calculations can largely reproduce our experimental results. In principle, the experiments can be extended to 2DES states at E_F by using *p*-type samples. Thus, we could add the electronelectron interaction, which reveals a wealth of further quantum phase transitions, in future experiments [9,11]. However, already the present results go far beyond previous results [12,13] by providing the states at the quantum phase transition, a LL fan diagram, and a detailed theoretical reproduction of the observed experimental LDOS features.

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