Electron-electron interactions in antidot-based Aharonov-Bohm interferometers

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We present a microscopic picture of quantum transport in quantum antidots in the quantum Hall regime taking electron interactions into account. We discuss the edge state structure, energy-level evolution, charge quantization and linear-response conductance as the magnetic field or gate voltage is varied. Particular attention is given to the conductance oscillations due to Aharonov-Bohm interference and their unexpected periodicity. To explain the latter, we propose the mechanisms of scattering by point defects and Coulomb blockade tunneling. They are supported by self-consistent calculations in the Hartree approximation, which indicate pinning and correlation of the single-particle states at the Fermi energy as well as charge oscillation when antidot-bound states depopulate. We have also found interesting phenomena of antiresonance reflection of the Fano type.

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I. INTRODUCTION

An antidot is a potential energy hill in a two-dimensional electron gas (2DEG) formed in a GaAs/AlGaAs heterostructure by a negative voltage applied on a surface gate.¹ It is often regarded as an artificial repulsive impurity and thus considered to be the inverse of a quantum dot. By applying a magnetic field perpendicular to the 2DEG, antidots have been extensively and intensively studied to understand edge state transport in the quantum Hall regime, $2-8$ $2-8$ charging in open systems and its influence on Aharonov-Bohm (AB) interference, $4,8-11$ $4,8-11$ $4,8-11$ fractionally quantized charge of Laughlin quasiparticles, 12 non-Abelian statistics of the fractional quantum Hall $5/2$ state, 13 and others.¹ Though electron transport in antidots seemed to be well understood, recent experiments of Goldman *et al.* revealed new unexpected features such as multiple periodicity of the Aharonov-Bohm conductance oscillations[.14](#page-10-7)

In a magnetic field, quantum Hall edge channels form closed pathways encircling an antidot.^{1[,15](#page-10-8)} They are separated from extended edge channels propagating along device boundaries by quantum point-contact (QPC) constrictions (see the insets in Fig. 1). At a given magnetic field, there are *fleads* propagating edge states in the leads at the Fermi energy. The electron density in the constrictions is smaller than in the leads and, hence, only the lowest f_c states are fully transmitted, whereas the remaining highest $f_{leads}-f_{c}$ states are partially or fully reflected. A typical conductance of the AB interferometer as a function of magnetic field exhibits a steplike structure with plateaus separated by wide transitions regions[.3–](#page-10-9)[6](#page-10-10)[,14](#page-10-7) At very low magnetic fields, the conductances of the QPC constrictions are additive and behave like classical Ohmic resistors[.16](#page-10-11) When the magnetic field increases, the conductance evolves from classical to quantum behavior with the single-particle levels condensed into degenerated Landau levels (LLs). The steplike dependence of the conductance reflects successive depopulation of the LLs in the constrictions[.15](#page-10-8) The plateau regions correspond to the field range, where the QPC openings are fully transparent (the transmission coefficient through an individual QPC is integer

 $T \cong f_c$), and transition regions between these plateaus correspond to the partially transparent QPC openings (the transmission coefficient is noninteger $f_c < T < f_c + 1$). When a single-particle state of the antidot-bound edge channel coincides with the Fermi energy E_F , it provides a pathway for scattering from an edge channel on one side of the sample to an edge channel on the opposite side. Thus, it gives rise to pronounced AB conductance oscillations in the transition regions between the plateaus.

According to the conventional theory of the Aharonov-Bohm interferometer,¹⁵ its conductance shows a peak each time the enclosed flux $\phi = BS$ changes by the flux quantum $\phi_0 = h/e$. Thus, the conductance of the interferometer as a function of the magnetic field exhibits the periodicity

$$
\Delta B = \frac{\phi_0}{S}.\tag{1}
$$

Here $S = \pi r^2$ is the area enclosed by a circular antidot-bound state of radii *r*, which approximates the geometrical area the antidot gate (Fig. [1](#page-1-0)). The enclosed flux through the interferometer can also be varied at a fixed magnetic field by changing an antidot gate voltage *Vadot*. In the case when the area changes linearly with the change in the gate voltage ΔS $=\alpha \Delta V_{adot}$, the expected periodicity is

$$
\Delta V_{adot} = \frac{\phi_0}{\alpha B}.\tag{2}
$$

Since the experimental study of Hwang *et al.*, [3](#page-10-9) the inter-pretation of Aharonov-Bohm oscillations based on Eq. ([1](#page-0-0)) has been widely accepted.¹ Measuring the period ΔB , the radii of the edge states circulating around the antidot can be deduced. This gives valuable information about actual depletion region in 2DEG. However, in the recent experimental work of Goldman *et al.*,^{[14](#page-10-7)} it was reported that the periodicity of the AB oscillations as a function of the magnetic field depends on the number of fully transmitted states in the constriction f_c and is then well described by the dependence

FIG. 1. The AB conductance oscillations calculated in the Hartree (solid lines) and Thomas-Fermi (dashed lines) approximations for two antidot gate voltages *Vadot*=−0.25, −0.68 V, and for temperature *T*= 0.2 K. *B* marks the AB period, which is the same for any filling factor in the constriction f_c . Insets on the right show the wave-function modulus for different f_c . The arrows with stars mark positions where the former was calculated. Inset on the top shows schematic structure of the antidot-based AB interferometer. Top pattern denotes the metallic gates on the GaAs heterostructure. The radius of the antidot gate is *R*= 200 nm. 2DEG resides a distance *b* from the surface.

$$
\Delta B = \frac{1}{f_c} \frac{\phi_0}{S},\tag{3}
$$

which differs by a factor of $1/f_c$ from the conventional for-mula ([1](#page-0-0)). On the other hand, the back-gate charge period ΔV_{edot} was found to be the same for all f_c , independent of the magnetic field in stark contrast to Eq. (2) (2) (2) .^{[14](#page-10-7)} This departure from the conventional periodicity of the AB oscillations cannot be explained in a one-electron picture of noninteracting electrons. To account for it, it is necessary to consider electron interactions and/or Coulomb blockade (CB) charging effects. Coulomb interactions define a potentially important energy scale because even rough estimation gives Coulomb energy values that can exceed kinetic energy in magnetic field $\hbar \omega_c$; $\omega_c = eB/m^*$ and m^* is effective electron mass.¹⁴

The possible importance of CB charging was suggested by earlier experiments of Ford *et al.*[4](#page-10-3) and Kataoka *et al.*, [8](#page-10-2) where doubled frequency of the AB oscillations was observed. For the filling factor $f_c = 2$, one may anticipate that resonances from one spin species should occur halfway between the neighboring resonances of the second spin species. This is, however, not the case. In a later experiment, using selective injection and detection of spin-resolved edge channels, it was shown that the antidot states with up spin do not provide resonant paths in the *h*/2*e* AB oscillations[.11](#page-10-4) No model of noninteracting electrons can explain this because in such models the $h/2e$ oscillations should be a simple composition of the two *h*/*e* oscillations coming from the two spin species, and the phase shift between the two *h*/*e* oscillations is determined by the ratio between the Zeeman energy and the single-particle level spacing. Since the ratio depends on the antidot potential at the Fermi level and the magnetic field, the noninteracting model cannot provide an explanation of the sample-independent π phase shift. Moreover, in the absence of interactions, both spins should participate in the resonant scattering, contradicting the experimental observation that only the spin species with the larger Zeeman energy contributes to the resonances. $4,8,11$ $4,8,11$ $4,8,11$ Thus, experimental findings gave strong motivation for a model that takes electron interactions into account.

To explain double-frequency Aharonov-Bohm oscillations, models accounting for the formation of compressible rings around the antidot⁸ and capacitive interaction between excess charges were introduced.⁹ The first model is based on the assumption that there are two compressible regions encircling an antidot separated by an insulating incompressible ring. Screening in compressible regions, and Coulomb blockade, then force the resonances through the outer compressible region to occur twice per h/e cycle.⁸ In the capacitiveinteraction model, $\frac{9}{5}$ two antidot-bound edge states are assumed to localize excess charges that are spatially separated from each other and from extended edge channels by incompressible regions. This allows one to include in the antidot Hamiltonian the capacitive coupling of excess charges. In a regime of weak coupling, Coulomb blockade prohibits relaxation of the excess charges unless one of the antidot states accumulates exactly one electron or spin-flip cotunneling between them is allowed. Analyzing the evolution of the excess charges as a function of magnetic field, it was proposed that the process responsible for doubling of the AB oscillations comprises two consecutive tunneling events of spin-down electrons and one intermediate Kondo resonance[.9](#page-10-12) This type of process agrees with the experiment, $\frac{11}{11}$ where only the electrons with spin down contribute to the *h*/2*e* AB oscillations.

A related topic of $1/f_c$ periodic AB oscillations has been recently investigated both experimentally $17,18$ $17,18$ and theoretically^{19,[20](#page-10-16)} for the case of quantum dot-based interferometers. The experiments of Camino *et al.*[17](#page-10-13) clearly demonstrated that the magnetic-flux AB period is described by Eq. ([3](#page-1-1)) and the gate voltage period stays constant for any filling factor f_c . Moreover, the authors reanalyzed the existing experimental data and showed that all of it can be also well described by Eq. ([3](#page-1-1)). To explain $1/f_c$ scaled period of AB oscillations in quantum dots, Coulomb blockade theory was introduced in Ref. [19.](#page-10-15) Assuming that a compressible island exists inside the quantum dot, the AB period is caused by charging of f_c fully occupied LLs in the dot. The validity of this assumption for the compressible island as well as the electrostatics of the AB interferometer has been discussed in Ref. [20](#page-10-16) by two of the present authors. It was shown and explained why the scattering theory based on Landauer formula^{[1](#page-0-0)5} predicts the conventional AB periodicity [Eqs. (1) and ([2](#page-0-1))]. A very recent experiment of Zhang *et al.*^{[18](#page-10-14)} pointed out that $1/f_c$ period oscillations that are caused by the CB effect hold in an AB interferometer with a small quantum dot. However, as the dot size increases, the charging energy becomes an unimportant energy scale and the conventional AB oscillations are restored.

In the present paper, motivated by the experiment of Goldman *et al.*^{[14](#page-10-7)} we address electron transport through the antidot AB interferometer from different standpoints. Starting from a geometrical layout of the device, we calculate self-consistently the edge state structure, energy-level evolution, charge quantization, and linear-response conductance in the Thomas-Fermi (TF) and Hartree approximations. We find that the AB periodicity is well described by the conventional formulas (1) (1) (1) and (2) (2) (2) in the case of the ideal structure without impurities. The conductance is dictated by the highestoccupied (f_c+1) th state in the constrictions, but the f_c antidot-bound states are well localized and do not participate in transport. Electron interactions in the Hartree approximation pin the f_c single-particle states to the Fermi level and force their mutual positions to be correlated. The Hartree approach also predicts that the number of electrons around the antidot oscillates in a saw-tooth manner reflecting sequential escape of electrons from the f_c edge states. For low temperatures, we have found an interesting phenomenon that we call antiresonance reflection of the Fano type. While the Hartree and Thomas-Fermi approximations do not reproduce experimental $1/f_c$ AB periodicity,¹⁴ we explore two mechanisms that might be relevant, namely, scattering by impurities and Coulomb blockade tunneling.

II. MODEL

We consider an antidot AB interferometer defined by split gates in the GaAs heterostructure similar to those studied experimentally.^{1[,3](#page-10-9)[–5,](#page-10-17)[10](#page-10-18)[,14](#page-10-7)} A schematic layout of the device is illustrated in Fig. [1.](#page-1-0) Charge carriers originating from a fully ionized donor layer form the 2DEG, which is buried inside a substrate at the GaAs/Al_xGa_{1-*x*}As heterointerface situated at a distance *b* from the surface. Metallic gates placed on the top define the antidot and the leads at the depth of the 2DEG.

The Hamiltonian of the whole system, including the semiinfinite leads, can be written in the form $H = H_0 + V(\mathbf{r})$, where

$$
H_0 = -\frac{\hbar^2}{2m^*} \left\{ \left(\frac{\partial}{\partial x} - \frac{eiBy}{\hbar} \right)^2 + \frac{\partial^2}{\partial y^2} \right\}
$$
(4)

is the kinetic energy in the Landau gauge, and the total confining potential

$$
V(\mathbf{r}) = V_{conf}(\mathbf{r}) + V_H(\mathbf{r}),\tag{5}
$$

where $V_{conf}(\mathbf{r})$ is the electrostatic confinement (including contributions from the top gates, the donor layer, and the Schottky barrier), $V_H(\mathbf{r})$ is the Hartree potential,

$$
V_H(\mathbf{r}) = \frac{e^2}{4\pi\epsilon_0 \epsilon_r} \int d\mathbf{r}' n(\mathbf{r}') \left(\frac{1}{|\mathbf{r} - \mathbf{r}'|} - \frac{1}{\sqrt{|\mathbf{r} - \mathbf{r}'|^2 + 4b^2}} \right),\tag{6}
$$

where $n(r)$ is the electron density, the second term corresponds to the mirror charges situated at the distance *b* from the surface, and $\varepsilon_r = 12.9$ is the dielectric constant of GaAs. Equation ([6](#page-2-0)) accounts for the average electrostatic potential generated by the total charge density. Though it is in general three dimensional, the thickness of the 2DEG is assumed to be small such that electrons occupy only the lowest-energy level in the *z* direction. Typically, the electrons reside within a well 10 nm thick, 15 which is about the lattice constant of our discretization mesh. Thus, the three-dimensional nature of the 2DEG is not expected to affect the computational results significantly. Another assumption underlying Eq. (6) (6) (6) is the low-frequency approximation for image charges. The antidot and the leads are treated on the same footing, i.e. the electron interaction and the magnetic field are included both in the lead and in the antidot regions. 21

We calculate the self-consistent electron densities, potentials, and the conductance on the basis of the Green's function technique. The description of the method can be found in Refs. [21–](#page-10-19)[23](#page-10-20) and thus the main steps in the calculations are only briefly sketched here. First, we compute the selfconsistent solution for the electron density, effective potential, and the Bloch states in the semi-infinite leads by the technique described in Ref. [24.](#page-10-21) Knowledge of the Bloch states allows us to find the surface Green's function of the semi-infinite leads. We then calculate the Green's function of the central section of the structure by adding slice by slice and making use of the Dyson equation on each iteration step. Finally, we apply the Dyson equation in order to couple the left and right leads with the central section and thus compute the full Green's function $G(E)$ of the whole system. The electron density is integrated from the Green's function (in the real space),

$$
n(\mathbf{r}) = -\frac{1}{\pi} \int_{-\infty}^{\infty} \Im[\mathcal{G}(\mathbf{r}, \mathbf{r}, E)] f_{FD}(E - E_F) dE, \tag{7}
$$

where f_{FD} is the Fermi-Dirac distribution. This procedure is repeated many times until the self-consistent solution is reached; we use a convergence criterion $|n_i^{out} - n_i^{in}|/(n_i^{out})$ $+n_i^{in}$ < 10⁻⁵, where n_i^{in} and n_i^{out} are input and output densities on each iteration step *i*.

Finally, the conductance is computed from the Landauer formula, which in the linear-response regime is 15

$$
G = -\frac{2e^2}{h} \int_{-\infty}^{\infty} dET(E) \frac{\partial f_{FD}(E - E_F)}{\partial E},
$$
 (8)

where the transmission coefficient $T(E)$ is calculated from the Green's function between the leads. $21-23$

To clarify the role of the electron interaction we also calculate the conductance in the TF approximation, where the self-consistent electron density and potential are given by the semiclassical TF equation at zero field

$$
\frac{\pi \hbar^2}{m^*} n(\mathbf{r}) + V(\mathbf{r}) = E_F.
$$
\n(9)

This approximation does not capture effects related to electron-electron interaction in quantizing magnetic field such as the formation of compressible and incompressible strips and, hence, it corresponds to the noninteracting oneelectron approach where, however, the total confinement is given by a smooth fixed realistic potential. $21,25$ $21,25$

While the present approach is not expected to account for single-electron tunneling in the conductance (leading to the Coulomb blockade peaks), 26 26 26 one can expect that it correctly reproduces a global electrostatics of the interferometer and microscopic structure of the quantum-mechanical edge states regardless whether the conductance is dominated by a singleelectron charging or not. This is because the interferometer is an open system with a large number of electrons surrounding the antidot and, thus, the electrostatic charging caused by a single electron hardly affects the total confining potential of the interferometer. Thus, the results of the self-consistent Hartree approach provide accurate information concerning the locations of the propagating states and the structure of compressible/incompressible strips in the interferometer. Our calculations are also expected to provide detailed information concerning the coupling strengths between the states in the leads and around the antidot.

III. RESULTS AND DISCUSSION

We calculate the magnetotransport of a quantum antidot AB interferometer with the following parameters representa-tive of a typical experimental structure.^{1,[3](#page-10-9)-5[,10,](#page-10-18)[14](#page-10-7)} The 2DEG is buried at $b = 50$ nm below the surface (the widths of the cap, donor, and spacer layers are 14 nm, 26, nm and 10 nm, respectively), the donor concentration is 1.02×10^{24} m⁻³. The width of the quantum wire and semi-infinite leads is 700 nm. The radius of the antidot gate is $R = 200$ nm (see Fig. [1](#page-1-0)). The gate voltage applied to the lead gates is $V_{lead} = -0.4$ V. With these parameters of the device, there are 25 channels available for propagation in the leads and the electron density in the center of the leads is $n_{lead} = 2.5 \times 10^{15} \text{ m}^{-2}$.

A. Magnetic-flux periodicity of Aharonov-Bohm oscillations

Figure [1](#page-1-0) shows the conductance of the AB interferometer as a function of magnetic field calculated for the quantummechanical Hartree and semiclassical Thomas-Fermi approximations. At very low magnetic field, the conductances of two parallel QPC constrictions are additive and behave like Ohmic resistors.¹⁶ At high magnetic field, in the quantum Hall regime, the total conductance exhibits steplike dependence caused by gradual depopulation of the LLs.¹⁵ All channels for electron propagation are fully open or fully blocked in the plateau regions, while a partly transmitted channel is present in the transition regions. The conductance oscillations due to Aharonov-Bohm interference are clearly seen in the transition regions, which indicates that they are caused by the (f_c+1) th partly transmitted channel.

The Aharonov-Bohm oscillations reveal the same period $\Delta B = 15$ mT independent of f_c for both the Hartree and Thomas-Fermi approximations as is seen in Fig. [1.](#page-1-0) This periodicity is in excellent agreement with the conventional AB formula ([1](#page-0-0)). It corresponds to an area enclosed by skipping orbits of $0.28 \mu m$, which gives a circle radius of r = 300 nm. The latter is slightly larger that the geometrical radius of the antidot gate $R = 200$ nm and is in agreement with the wave-function maps in the insets in Fig. [1.](#page-1-0) The plotted wave functions show how the antidot-bound states depend on the number of Landau levels f_c (and therefore on the value of magnetic field) and also on the antidot gate voltage.

The AB oscillations can be related to the evolution of the corresponding energy spectrum when single-electron states cross the Fermi level each time the flux, through the antidot, increases by the flux quantum. This is illustrated in Fig. $2(b)$ $2(b)$, which shows an evolution of the resonant levels as a function of magnetic field in the vicinity of the Fermi energy. To obtain the evolution of the resonant levels, we draw an imaginary ring and analyze the density of states (DOS) there at each given *B*. The inner radius of the ring is chosen to be the antidot radius and the outer radius is 150 nm larger, i.e. we account for all states in the range 200–350 nm from the center. When DOS has been calculated, its peaks are searched for and their positions are plotted as illustrated in Fig. $2(b)$ $2(b)$]. Each conductance minimum of the AB interferometer seen in Fig. $2(a)$ $2(a)$ corresponds to a resonant-level aligns with the Fermi energy E_F in Fig. [2](#page-4-0)(b), implying a condition of resonants reflection of the extended edge state in the lead by the antidot.² However, two out of every three states present in Fig. $2(b)$ $2(b)$ are associated with neither minimum nor maximum of the AB conductance oscillations at 0.[2](#page-4-0) K [the dashed curve in Fig. $2(a)$]. Only if the temperature is lowered from 0.2 to 0.02 K does the conductance start to reveal a feature caused by their presence at E_F . Inspection of the DOS at $B = 0.814$ T [see the inset in Fig. [2](#page-4-0)(b)] shows that these extra states produce very narrow and sharp resonances (with broadening $\Gamma \ll k_B T$ for $T=0.2$ K), while the resonance due to the conductance-mediating state is broad and low (Γ) $\sim k_B T$). Let us call two former states "antidot-bound states" and refer to the latter one as a "transport state." The singleparticle transport state originates from the partly transmitted (f_c+1) th LL in the constriction. It is strongly coupled to the extended edge states in the leads. By contrast, all f_c antidotbound states are weakly coupled to the extended edge states because they are situated closer to the antidot hill and surrounded by an incompressible strip. Figure $2(c)$ $2(c)$ displays local density of states (LDOS) maps clearly showing the presence of $f_c = 1$ and $f_c = 2$ antidot-bound states at corresponding resonance peaks in DOS. Note that the evolution of both the

FIG. 2. (a) The conductance calculated in the Thomas-Fermi approximation for different temperatures and *Vadot*=−0.4 V. Solid curve: 0.02 K; dashed curve: 0.2 K. (b) Evolution of the resonant energy levels near E_F . The inset shows DOS in a ring around the antidot for the specified value of magnetic field $B = 0.814$ T. The ring is chosen so that contains the antidot-bound states and has outer and inner radii equal to 350 nm and 200 nm, respectively (note that the antidot gate radius is $R = 200$ nm). The evolution of the energy levels was obtained from the peak positions of the DOS at each given value of B . (c) LDOS at three peaks corresponding to three different energies at $B = 0.814$ T.

antidot-bound and transport states is not correlated between each other. All of them cross the Fermi energy at different magnetic fields although the ΔB interval for single-particle states belonging the same LL stays constant. This is a characteristic feature of noninteracting electrons in the Thomas-Fermi approximation.

FIG. 3. The AB resonant reflection and adiabatic transmission at f_c =2 as calculated in the Thomas-Fermi approximation and marked by arrows in Fig. $2(a)$ $2(a)$. The left column corresponds to minimum and right column to maximum of the conductance $B=0.81$ T and 0.802 T, respectively. Plots (a) - (h) show the wave-function modulus $|\Psi_i|^2$ of *i*th edge state and (i)–(l) are for LDOS integrated over transverse y direction. Panels (i) and (j) illustrate resonant levels in the vicinity of E_F in an enlarged scale of (k) and (l). Fat solid lines in (k) and (l) are the total confinement potential along $y=0$.

B. Antiresonance reflection

To understand features that are seen to be superimposed on the AB conductance oscillations in Fig. $2(a)$ $2(a)$, let us first look at the wave functions and LDOS at minima and maxima of the conventional AB oscillations due to the (f_c+1) th transport state. These are shown in Fig. [3:](#page-4-1) the wave functions for f_c lead edge states show perfect transmission through the device, while the state $f_c + 2$ as well as all higher states exhibit perfect reflection of the incoming state from the left lead back into the left lead. The only state modulating the transport through the antidot is the transport state $f_c + 1$. It

FIG. 4. (a) and (b) Fragment of energy structure and DOS at $B = 0.814$ T from Fig. [2](#page-4-0)(b). (c) Transmission coefficient $T_3 = \sum_i T_{i3}$ from the third mode in the left lead to all available *i*th modes in the right lead. There is a reflection antiresonance at *E*= 0.014 meV that manifests itself as a zigzag jump in T_3 . (d) Derivative of the Fermi-Dirac function for $T=0.02$ and 0.2 K. (e) and (f) Third wavefunction squared modulus $|\Psi_3|^2$ at two sides of the antiresonance.

effectively provides a resonant tunneling pathway between incoming and outgoing edge channels via the antidot-bound state. Inspection of LDOS in Figs. $3(i)-3(1)$ $3(i)-3(1)$ shows the presence of the single-particle state at the Fermi energy during resonant reflection via (f_c+1) th transport state.

At low temperatures, the f_c antidot-bound states produce a sharp zigzaglike feature in the conductance as shown in Fig. $2(a)$ $2(a)$ at $T=0.02$ K. This effect is most pronounced for *B* $= 0.79$ T and might be attributed to a Fano-type resonance.²⁷ It differs qualitatively from the common resonance dip due to (f_c+1) th transport states. In order to get insight into its origin, let us look at the transmission coefficient T_{f+1} vs energy plot [Fig. $4(c)$ $4(c)$]. Note that the conductance is an integral over the total transmission weighted by the Fermi-Dirac derivative $[Eq. (8)]$ $[Eq. (8)]$ $[Eq. (8)]$ and, thus, it can not resolve all of the details unless the temperature is extremely low. Overall, the dependence of T_{f_c+1} is quite smooth showing a deep minimum associated with resonant reflection of (f_c+1) th state. This state can also be monitored in DOS as the wide and low peak [Fig. $4(b)$ $4(b)$]. However, there is a superimposed zigzaglike jump in T_{f_c+1} , which is totally different and caused by scattering between the (f_c+1) th state and the f_c one. It happens at $E=0.014$ $E=0.014$ meV in Fig. 4(c) and is caused by $3 \rightarrow 2$ scattering. At an energy slightly less than *E*= 0.014 meV, there is a sharp dip, which is associated with the third extended state being well coupled to the second antidot-bound state [Fig. $4(e)$ $4(e)$]. However, as energy increases and slightly exceeds $E = 0.014$ meV, it turns into a sharp peak with the antidot-bounded state becoming perfectly isolated Figs. $4(f)$ $4(f)$]. The fact that it is isolated can be seen as an absence of "bridges" to the extended lead state and almost perfect circled shape. Because the shake reveals in the transmission, this feature is not indent in the conductance at temperatures exceeding $k_B T > \Gamma$. We refer to this type of resonance as an antiresonance of the Fano type $2⁷$ since it results from quantum interference between two processes: one involving

FIG. 5. (a) Number of electrons around quantum antidot *N*, (b) the AB conductance oscillations, and evolution of the resonant energy levels near E_F [(c) and (d)] calculated in the Hartree approximation for different temperatures and *V_{adot}*=−0.4 V. *N* and energy levels are calculated in an annulus around the antidot, whose outer and inner radii are equal to 350 nm and 200 nm, respectively (note that the antidot gate radius is $R = 200$ nm). Evolution of the energy levels in (c) is calculated at $T=0.2$ K, while (d) is for $T=0.1$ K. The multiple levels marked by smaller circles are due to bulk states in the leads, which inevitably captured in the region of interest, i.e., the annulus around the antidot.

strongly localized state, the f_c antidot orbital, and the other being the less strongly localized $f_c + 1$ orbital. When the incident energy exactly coincides with the antiresonance energy, the AB phase $2\pi\phi/\phi_0$ flips by $\phi = \pi$. It is worth mentioning that the antiresonance here differs from the phase change in the AB conductance oscillations observed in Refs. [4–](#page-10-3)[6,](#page-10-10) where a phase flip between consecutive oscillations accompanied by the change in oscillation period. 5.6 5.6

C. Effect of electron interaction in Hartree approximation

Accounting for electron interactions within the quantummechanical Hartree approximation brings qualitatively new features to the AB oscillations [Fig. 5]. First, the conductance does not show perfect smooth regular oscillations any longer: their shapes become very distorted, which is especially pronounced as temperature lowers. Secondly, the energy levels from both $f_c + 1$ and f_c edge states become correlated: their positions at E_F become more equally spaced. This is attributed to the effect of Coulomb interaction that favors only one single-particle state being depopulated at a given magnetic field. In other words, electrons escape from localized f_c states one by one. Third, as the temperature decreases the energy levels due to f_c edge states become pinned to E_F [Figs. $5(c)$ $5(c)$ and $5(d)$]. We define the pinning as a lower slope at E_F (see Ref. [21](#page-10-19) for a detailed discussion of the pinning effect). When a state is pinned to E_F , it easily adjusts its position or occupation in response to any external perturbation. The role of the perturbation might be played by either magnetic field or applied gate voltage. Therefore, the screening of the antidot, and related metalliclike behavior of the system, is provided solely by the f_c states not by the (f_c) $+1$)th transport state. The latter crosses E_F steeply and mediates the conductance oscillations in a similar fashion to that in the semiclassical Thomas-Fermi approximation. Fourth, the number of electrons *N* in an annulus around the antidot shows saw-tooth oscillations that reflects pinning and depopulation of f_c edge states. Intervals of magnetic field with linear negative slopes of *N* and pinned states are marked by the shaded regions in Fig. [5.](#page-5-1) The negative slopes of *N* in Fig. $5(a)$ $5(a)$ are caused by gradual depopulation when the corresponding single-particle state is pushed up in energy and its occupation decreases. Note that the change in electron number is less than unity, which we explain by the finite temperature and bulk states captured in the region of interest, i.e. in the annulus of 200–350 nm size from the center. It also worth noting that the saw-tooth dependence of *N* is in excellent agreement with the experimental findings in Ref. [10.](#page-10-18)

While the shape of the AB conductance oscillations is strongly nonsinusoidal in the Hartree approximation, their periodicity still described by the conventional Eq. (1) (1) (1) that, in turn, disagrees with the experiment of Goldman *et al.*[14](#page-10-7) The Hartree approach is known to describe well the electrostatics of the system at hand. This is confirmed by the good agreement with numerous experiments including, for example, formation of compressible/incompressible strips in quantum wires²⁸ and the statistics of conductance oscillations in open quantum dots[.25](#page-10-22) Thus, the validity of the energy-level evolution as well as the electron number oscillation presented in Fig. [5](#page-5-1) is indeed qualitatively correct. The conductance, however, may be incorrect. It is calculated using the Landauer-Buttiker formalism that has been shown to fail in the regime of weak coupling, when the conductance is less than the conductance quantum $G_0 = 2e^2/h$.^{[26](#page-10-23)} Though the total conductance is larger than G_0 , the antidot is in the weak-coupling regime (a related discussion of quantum dot-based AB inter-ferometers is given in Ref. [20](#page-10-16)). This is because of the adiabatic character of the transport when the lowest f_c states pass through the interferometer without any reflection (see Fig. [3](#page-4-1)). The highest (f_c+1) th transport edge state, giving rise to the AB oscillations in the transition regions between the plateaus, becomes thus effectively decoupled from the remaining f_c states. Therefore, because of well-localized f_c states and partly localized (f_c+1) th state, the electron charge in the antidot may become quantized and transport through the interferometer strongly affected by the Coulomb blockade effect.

FIG. 6. The AB conductance oscillations calculated in the Hartree (solid lines) and Thomas-Fermi (dashed lines) approximations for fixed magnetic field $B=0.8$ T. ΔV_{adot} is the AB period, which scales linearly with the filling factor in the constriction f_c . Temperature $T = 0.2$ K.

D. Gate voltage periodicity of Aharonov-Bohm oscillations

The Aharonov-Bohm oscillations can be also observed when the gate voltage varies for a fixed magnetic field. Figure [6](#page-6-0) shows the conductance as a function of the antidot gate voltage calculated in the Thomas-Fermi and Hartree approximations for $B=0.8$ T. As in the case of magnetic field de-pendence [Fig. [1](#page-1-0)], the AB conductance oscillations are pronounced in the transition regions between plateaus. However, the period ΔV_{adot} scales linearly with the filling factor f_c , in agreement with the conventional formula ([2](#page-0-1)).

Both magnetic field and gate voltage periodicity of the AB oscillations are well described by conventional formulas ([1](#page-0-0)) and ([2](#page-0-1)). However, it disagrees with the experiment (Ref. [14](#page-10-7)). To overcome this discrepancy, we consider two effects in the following, namely, the scattering by random impurity potentials and the Coulomb blockade theory.

E. Scattering by random impurity potentials

It is known that impurity scattering might substantially modify electron transport in AlGaAs heterostructures.¹⁵ Depending on the nature of scatters, the scattering potential varies widely. Charged impurities, such as ionized donors, have a long-range potential, whereas neutral impurities have short-range potentials. These two cases have different effects on electron transport. The first leads to the localization of edge states in the quantum Hall regime and strong modification of the transition regions between quantum Hall plateaus[.29](#page-10-26) The edge states circulating around the antidot might change their locations and, as a result, the AB oscillations might experience sudden period changes[.5,](#page-10-17)[6](#page-10-10) Because $1/f_c$ periodicity in the experiment of Goldman *et al.*^{[14](#page-10-7)} is robust and measured for different cool down cycles, we are skeptical that the long-range scattering is responsible for $1/f_c$ periodicity and concentrate in the following on the short-range scattering.

A physical realization of the short-range disorder potential occurs if some neutral impurity, such as Al atom, has diffused out of a AlGaAs barrier into a GaAs well where the

FIG. 7. (a) The conductance and (b) resonant energy structure calculated in the Thomas-Fermi approximation for various disorder potentials V_{imp} . (c) The conductance calculated in the Hartree approximation. $V_{adot} = -0.4$ V.

2DEG resides. The Al atoms are the scattering centers with a potential *Vimp*. Because we are interested in scattering between the extended and bound states and between different bound and partly bound states, we generated random point scatters in the vicinity of the antidot. Figures $7(a)$ $7(a)$ and $7(b)$ show the conductance and energy-level structure calculated in the Thomas-Fermi approximation for the case of 500 random scatters, which corresponds to concentration *nimp* $= 10^{15}$ m⁻². As the magnitude of *V_{imp}* increases, the scattering between different edge states becomes stronger. The AB oscillations due to resonant reflection of the (f_c+1) th state are gradually suppressed with a new oscillation pattern emerging from resonant transmission via the f_c antidotbound states. The positions of the AB conductance peaks are clearly correlated with the antidot-bound states crossing E_F [see Figs [7](#page-7-0)(a) and 7(b)]. The transport (f_c+1) th state, which governs the conductance in the ideal case without impurities, becomes easily destroyed because it is half-filled in the QPC constrictions and, thus, slight potential fluctuations effectively block its propagation. However, in contrast to the ideal case with no defect scattering, for strong defect scattering a resonant transmission peak is visible in Fig. $7(a)$ $7(a)$ whenever any antidot-bound state crosses E_F in Fig. [7](#page-7-0)(b) at the higher values of the magnetic field that are shown.

The Hartree approximation [Fig. $7(c)$ $7(c)$] does not clearly recover resonant transmission triggered by the disorder potential, as it occurs in the Thomas-Fermi approach. In this case, there are only faint remnants of resonant transmission

via the f_c antidot-bound states at V_{imp} = 20 meV. We attribute this to inadequate modeling of point defects due to Al atoms. The minimal area that can be occupied by a defect in our simulations is 5×5 nm². This is two orders of magnitude larger than the realistic cross section of the Al atom. On the other hand, the height of the realistic defect potential is also much larger and on the order of $V_{imp}^{Al} \approx 1$ eV. Note that the screened potential after self-consistent calculation is about ten times smaller than the input V_{imp} in Fig. [7](#page-7-0)(c). If we perform simulation for higher potentials, the self-consistent density becomes quickly washed out preventing electron transport through the region occupied by the defects. Based on our calculations in the TF approximation that are described above, we conclude that the short-range disorder is a plausible source of $1/f$ AB periodicity, but a quantitative comparison based on the more sophisticated Hartree approximation remains to be done and is beyond the scope of the present work.

F. Coulomb blockade model for Aharonov-Bohm oscillations

The quantum antidot does not confine electrons electrostatically, but sufficiently large magnetic field causes the formation of localized bound states where charge might be quantized. Direct evidence of the charging effect in the antidot was given in the experiment of Kataoka *et al.*[10](#page-10-18) Placing a noninvasive voltage probe in close proximity to the antidot, they detected steady accumulation followed by sudden relaxation of a localized excess charge nearby. The saw-tooth resistance oscillations measured by the detector coincide with the resonances monitored in the antidot conductance. Therefore, it was concluded that a source of the excess charge is the antidot, and its conductance is mediated by Coulomb charging. Additional evidence for the CB effect follows from the measurement of the conductance as functions of both magnetic field and source-drain bias, where clear and regular Coulomb diamonds were observed[.10](#page-10-18) The doubled frequency conductance oscillations measured in Refs. [4](#page-10-3) and [8](#page-10-2) are also a strong indication of the Coulomb charging in the antidot. Thus, we conclude that $1/f_c$ periodicity observed by Goldman *et al.*^{[14](#page-10-7)} might be a result of the CB effect. As an indirect support, it is worth mentioning that a rough estimation of the electron Coulomb interaction energy gives values exceeding kinetic energy in magnetic field.¹⁴

Motivated by these experimental arguments as well as our calculation results presented above, we develop a simple phenomenological model based on the CB orthodox theory.^{30[,31](#page-10-28)} Let us consider a case of f_c fully occupied LLs in the QPC constrictions, when the conductance is near a plateau region. Notice that this case is represented in the experiment of Goldman *et al.*^{[14](#page-10-7)} Figure [8](#page-8-0)(a) illustrates schematically the edge states existing around the antidot. For a given magnetic field and antidot gate voltage, we may draw a closed curve of area *S* that encompasses all f_c bound states. There is some background number of electrons *N* inside the area *S*. If we fix the position of the curve and increase the field by amount $\Delta B = \phi_0 / S$, the total flux through the reference area ϕ will increase by one flux quantum ϕ_0 . One single-particle state in each f_c LL gets pushed up in energy,

FIG. 8. (a) Schematic illustration of the edge states in a quantum antidot when the filling factor in constrictions is $f_c = 3$. An electron from extended edge states may tunnel into the outermost antidotbound state, but not into the bound states due to the low-lying LLs. (b) Equivalent single-electron scheme for the antidot edge states. Dotted lines mark paths for sequential tunneling. (c) Excess charge on the LLs Δq and (d) Coulomb blockade oscillations calculated within the orthodox theory.

crosses the Fermi energy, and becomes depopulated. The number of electrons in the reference area drops by f_c . However, it is known that the magnetic field does not change the number of electrons¹⁵ and, thus, electrons cannot just disappear. There must be a balancing influx of electrons into LLs in a way that the degeneracy of each LL increases exactly by one. For the case of a voltage applied to the antidot gate, a background charge in the reference area can be increased and *Ngate* additional electrons can be attracted to the area *S*. Hence, we obtain a total charge imbalance inside the area *S* given by $eN + \text{frac}(ef_c\phi) - eN_{gate}$, which leads to a charging energy

$$
E = \frac{e^2}{2C} [N + \text{frac}(ef_c \phi) - N_{gate}]^2, \tag{10}
$$

where *C* is the capacitance of the island and frac (x) means the fractional part of x . Note that the charging energy (10) (10) (10) takes the same value when ϕ changes by $1/f_c$. Therefore, magnetic field periodicity satisfies Eq. (3) (3) (3) as measured in the experiment of Goldman *et al.*^{[14](#page-10-7)} The antidot gate period, however, is one-electron charge in the reference area independent of the filling factor. It is also worth noting a similar approach to CB charging for the quantum dot-based interferometer (Ref. [19](#page-10-15)).

We have several comments about the CB theory proposed above. First, the change in magnetic field or gate voltage is supposed not to be large so that the reference area *S* always encloses a fixed number f_c of edge states. Secondly, all edge states encircle the antidot at about the same radii from the center and the extent of their wave functions is an unimportant length scale. This is evidently correct for large antidot radii or large magnetic fields. Third, the charging energy ([10](#page-8-1)) does not depend on which particular edge state builds the charge imbalance at a given field and gate voltage. Equation (10) (10) (10) rather treats all f_c states as one single-electron island. However, our calculations within both the Thomas-Fermi and Hartree approaches identify the highest LL as the most important for electron transport. On the other hand, experimental data clearly indicate that transport occurs via the highest LL (the outermost edge state) when the antidot in the CB regime.¹¹ Thus, we assume that a single-electron island implied in Eq. (10) (10) (10) has an internal structure and functions as a single-electron trap³⁰ [see Fig. $8(b)$ $8(b)$]. Electrons sequentially hop from the low-lying LLs into the extended edge states and vice versa via highest (f_c) th LL. This is supported by electrostatics shown in Fig. $5(a)$ $5(a)$, where states depopulate one by one.

To solve the electron-transport problem in the Coulomb blockade regime as governed by Eq. (10) (10) (10) , we employ the standard orthodox theory.^{30,[31](#page-10-28)} It describes an evolution of system via a "master" equation for $p(N)$, the probability that there are *N* electrons in the island

$$
\Gamma_{N-1 \to N} p(N-1) + \Gamma_{N+1 \to N} p(N+1) = [\Gamma_{N \to N-1} + \Gamma_{N \to N+1}] p(N).
$$
 (11)

Here $\Gamma_{N' \to N}$ is the sum of the transition rates through tunnels barriers, which change the electron number N' to N . Each transition rate treats tunneling of a single electron through a tunnel barrier as a random event and depends on the reduction in the electrostatic energy of the system $[Eq. (10)]$ $[Eq. (10)]$ $[Eq. (10)]$, resulting from such a tunneling event. For example, the transition rate from the source electrode to island reads as

$$
\Gamma^{s \to i}(N) = \frac{\Delta E^{s \to i}(N)}{e^2 R_s} (1 - e^{\Delta E^{s \to i}(N)/k_B T})^{-1},
$$
(12)

where $\Delta E^{s\rightarrow i}(N)$ is the energy change after tunneling from the source to the island and R_s is the resistance of the tunneling barrier between the source and island. We solve Eq. (11) (11) (11) and then calculate the average current as

$$
I = \sum_{N} \left[\Gamma^{s \to i}(N) - \Gamma^{i \to s}(N) \right] p(N). \tag{13}
$$

Figures $8(c)$ $8(c)$ and $8(d)$ show a representative calculation for the case of $f_c = 3$. We use capacitances and resistances of the tunnel barriers estimated from our self-consistent calculations given above. For a realistic quantum antidot shown in Fig. [1,](#page-1-0) approximate parameters are $C_1 = C_2 \approx 10^{-18}$ F and $R_1 = R_2 \approx 125 \text{ k}\Omega$. R_1 and R_2 are chosen to yield peak widths in Fig. $8(d)$ $8(d)$ similar to those in the experimental measurements of Goldman *et al.*^{[14](#page-10-7)} The excess charge Δq as a func-

tion of magnetic flux is shown in Fig. $8(c)$ $8(c)$. It oscillates in a saw-tooth manner within a window $-0.5 < Δ_q < 0.5$. Each time the value +0.5 is approached, an electron tunnel through the potential barriers with no energy cost and substantial current starts flowing through the antidot. When the magnetic flux increases further, one of the LLs gets recharged by one electron and the process repeats for another LL. For $f_c = 3$, the increase in magnetic field by the flux quantum ϕ_0 generates three successive rises of Δq and three related peaks of the current. Note that the current is proportional to the resistance measured in a two-terminal setup.

IV. CONCLUSIONS

In the present paper, we provide a microscopic physical description of the edge states existing in the quantum antidot focusing on the related conductance oscillations due to the Aharonov-Bohm interference. Motivated by recent experiment of Goldman *et al.*, [14](#page-10-7) we discuss different mechanisms that might be a source of measured $1/f_c$ periodicity of the AB conductance oscillations. Our findings are summarized as follows.

(1) Approaches based of the Hartree and Thomas-Fermi models for an ideal antidot structure predict the conventional AB conductance oscillations, as described by formulas (1) (1) (1) and (2) (2) (2) , i.e., magnetic field period does not depend on f_c and gate voltage period scales linearly with f_c . This is caused by transport isolation of the f_c edge states circulating around the antidot and the conductance being modulated solely by the highest-occupied (f_c+1) th state in the constrictions.

(2) Electron interactions in the quantum-mechanical Hartree approximation bring qualitatively new features to the electrostatics of the antidot AB interferometer. The singleparticle states originating from the f_c edge states become pinned to the Fermi energy. Their mutual positions at the Fermi energy are correlated, such that there is only one single-particle state depopulating at a given magnetic field and gate voltage. The number of electrons around the antidot shows a related saw-tooth dependence. It reflects the fact that particles escape sequentially, i.e., one by one.

(3) As the temperature decreases, both the Hartree and Thomas-Fermi approximations start to reveal a reflection antiresonance of the Fano type. This manifests itself as zigzag jumps in the conductance, with the (f_c+1) th transport state scattered into f_c antidot-bound state and vice versa. On the lower-energy side of the antiresonance, these states are strongly coupled with each other, but they are perfectly isolated on the high-energy side.

(4) The experimentally measured $1/f_c$ periodicity might be recovered if some disorder is introduced around the antidot. This can be naturally realized due to Al atoms diffused into the well where the 2DEG is. A short-ranged potential forces different edge states to mix and all f_c states might start to participate in transport. This is accompanied by changing of the AB conductance from resonant reflection to resonant transmission.

(5) A simple Coulomb blockade theory might also explain the $1/f_c$ periodicity. Using information about edge states and their occupancy from the self-consistent calculations, we

FIG. 9. Transmission coefficients $T_i = \sum_i T_{ij}$ from *i*th mode in the left lead to all available *j*th modes in the right leads calculated in the Hartree approximation for $f_c = 2$. Dashed lines present ideal structure, while solid lines are given for impurities with *Vimp* $= 20$ meV.

write down the charging energy of the system and calculate the transport in the sequential tunneling regime. When the magnetic field changes by a flux quantum, the system experiences recharging by f_c electrons.

While the experiment¹⁴ as well as our present study show presumably that $1/f_c$ AB periodicity is caused by CB tunneling, we suggest an experiment to verity that. If the size of the antidot were increased many times, the capacitance of the system would increase proportionally. The charging energy, therefore, would decrease and CB tunneling becomes unimportant.¹⁸ In this situation, the conventional AB oscillations, independent of f_c if magnetic field changes, should be restored.

From the theoretical point of view, it would be interesting to calculate the conductance by the Hartree-Fock approach, where the self-interaction problem is eliminated. One might also try density-functional theory with a proper exchangecorrelation functional that avoids self-interaction errors as well.

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APPENDIX: EXPLANATION OF INCREASED CONDUCTANCE DUE TO PRESENCE OF IMPURITIES IN HARTREE APPROXIMATION

Figure [9](#page-9-0) shows the partial transmission coefficients for the antidot without and with impurities as calculated in the Hartree approximation. Each curve gives the transmission probability for an electron incident from the left lead and passing into the right lead. For the ideal antidot structure, transmission is near unity or zero except for the (f_c+1) th

state, i.e., the third state in Fig. [9.](#page-9-0) When impurities are present, they effectively mix all states and prevent perfect transmission or reflection with unity or zero transmission. As a result, electrons from LLs in the leads that were perfectly reflected become partly transmitted through the antidot and the conductance, which is the sum of all transmission coefficients, rises [see Fig. $7(c)$ $7(c)$]. This feature of the selfconsistent calculation in the Hartree approach is in contrast to the simpler Thomas-Fermi approach. The latter, as expected, predicts the conductance being decreased when impurities are introduced because in the Thomas-Fermi model the increase in the transmission of modes 4 and 5 when impurities are introduced is insufficient to compensate for the decrease in the transmission of modes 1 and 2.

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