Fano resonances and electron spin transport through a two-dimensional spin-orbit-coupled quantum ring

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Abstract: Electron transport through a spin-orbit-coupled quantum ring is investigated within linear response theory. We show that the finite width of the ring results in the appearance of Fano resonances in the conductance. This turns out to be a consequence of the spin-orbit interaction that leads to a breaking of the parity of the states localized in the ring. The resonances appear when the system is close to maxima of Aharonov-Condorquantate conductance oscillations where spin transfer is heavily modified. When the spin-orbit coupling strength is detuned from the Aharonov-Condorequantate maxima the resonances are broadened resulting in a dependence of the spin transport on the electron Fermi energy in contrast to predictions from one-dimensional models.

DOI: 10.1103/PhysRevB.84.235319 PACS number(s): 73.23.Ad, 73.63.Nm

I. INTRODUCTION

Electrical manipulation of spin polarization of carriers is one of the key elements for semiconductor spintronics devices. Since the proposal of the spin-field-effect transistor by Das and Datta1 particular attention has been addressed to Rashba spin-orbit (SO) coupling.2 This interaction—a relativistic consequence of the presence of the electric field within the device—produces an effective magnetic field for the moving electrons which makes their spins precess. The Rashba coupling has been successfully implemented in quantum devices that operate on an electron spin through the control of voltages applied to the electrodes in the system, such as quantum gates3 and valves.5

Proposals of spin-operating devices concern also spin-orbit-coupled quantum rings as a realization of universal quantum gates,6 spin beam splitters,7 or spin filters.8 The electron transfer through quantum rings involves both the spin precession due to the SO interaction and the quantum interference effects related to Aharonov-Bohm9-10 and Aharonov-Condorequantate (AC) effects.11 The latter spin-interference effect12,13 results from the fact that the relative phase shifts for the wave function passing through both arms of the ring are spin dependent in the presence of SO interaction. The AC oscillations were probed experimentally in a HgTe single quantum ring,14 in a single15 and in an array of InGaAs quantum rings,16 or in an array of Bi2Se3 topological insulator quantum rings.17

Although the theoretical work on SO-coupled quantum rings is rich, it is based mostly on the idealized case of a ring with infinitesimal narrow channels, i.e., the one-dimensional approximation. This approach allowed one to obtain analytical description of charge18 and spin6 transport through the ring as function of the electron Fermi energy and the Rashba SO coupling strength. Theoretical studies concerning two-dimensional channels showed however that for an accurate description of transport through the spin-orbit-coupled ring, the finite width of the channels cannot be neglected.13,14,19 This is mainly due to the fact that for a finite-width ring the spin is no longer well defined. Nonetheless, full calculations are rather scarce. A ring with two-dimensional channels has been studied within a tight-binding formalism in Ref. 13 or in the framework of the scattering matrix method in Ref. 20. Reference 19 investigated the transport within the multiband Landauer-Büttiker formalism. This was later extended to describe the experimental data obtained in the presence of an external magnetic field.14

In this paper we develop a calculation scheme that allows us to study spin transport through a SO-coupled ring with two-dimensional channels. We show that the finite width of the channels along with SO interaction results in the appearance of Fano resonances in the conductance around the AC oscillation maxima. These sharp peak-dip structures have been previously studied in the presence of an external magnetic field, where they were the result of the broken symmetry of states localized in the ring,21 in systems of a quantum ring coupled to a quantum dot,22 or in one-dimensional quantum rings containing impurities23 and magnetic structures.24 Here we find that the Fano resonances originate from the coupling of the transmitted electron with the resonance states localized in the ring that have broken parity as a consequence of the SO interaction.

We find that in the resonance region the spin transport through the ring is strongly modified. We argue that the modification is caused by the coupling of the electron spin with the spin of the resonance states which is revealed by the application of an external magnetic field. When the SO coupling strength is detuned from the AC oscillation maxima the Fano resonances are broadened which results in a dependence of the spin transport on the electron Fermi energy. This particular result was not present in previous studies on spin transformations in one-dimensional rings6 and is of importance for spintronics devices based on SO-coupled rings.

II. THEORY

A. System

We consider a system described by the effective mass Hamiltonian

\[ H = \left( \frac{\hbar^2 k^2}{2m} + V_c(\mathbf{r}) \right) + \frac{1}{2} g \mu_B B \sigma_z + H_{SIA}, \]  

(1)
The dashed thick lines in the leads present the closed system of region where Rashba coupling is present is marked with blue color. The potential of the ring and the leads shown by the black curves. The \( V_c \)

M. P. NOWAK, B. SZAFRAN, AND F. M. PEETERS PHYSICAL REVIEW B 84, 235319 (2011)

in Fig. 1(a) by the black curve.

barrier). The contour of the confinement potential is depicted the potential that enters the Rashba Hamiltonian

include magnetic field \((0\ \hat{B})\) directed perpendicular to the plain of the device. We choose the Lorentz gauge \( A = (A_x, A_y, 0) = (0, Bx, 0) \).

We include Rashba SO interaction with the Hamiltonian \( H_{SIA} = \alpha \nabla \cdot (\sigma \times \mathbf{k}) \) resulting from the electric field felt by the propagating electron. In order to allow for a well-defined spin in the leads we assume that the Rashba coupling is present solely in the ring area by applying the electric field only therein. Experimentally this is realized by applying a voltage between the substrate and a gate that is restricted to the ring area.

In the considered two-dimensional system we describe the potential that enters the Rashba Hamiltonian \( H_{SIA} \) by \( V(x, y, z) = V_c(x, y) + |e| F_z f(y) \).

We include inhomogeneous electric field \( F_z f(y) \) that controls the coupling strength. In Fig. 1(a) in the blue region the electric field is approximately equal to \( F_z \).

The function that controls \( F_z f(y) \) has the form \( f(y) = |\text{arctan}(y - y_1) + \pi/2| - |\text{arctan}(y - y_2) + \pi/2|, \) where \( y_1 = 200 \) \( \text{nm} \) and \( y_2 = 500 \) \( \text{nm} \); see Fig. 1(a). The function is nearly steplike. We depict \( f(y) \) by the blue curve in Fig. 1(b).

Finally, we obtain the Rashba operator for the electric field in the growth direction that depends on the \( y \) position [we neglect the derivatives of \( V_c(r) \) as the wave function disappears in the proximity of the edges of the confinement potential]:

\[
H_{SIA} = \alpha |e| F_z \frac{\partial f(y)}{\partial y} \sigma_z k_z + a |e| F_z f(y) (\sigma_x k_y - \sigma_y k_x).
\]

Note that the first term in the latter equation guarantees that the two-dimensional Rashba Hamiltonian is Hermitian when the strength of the coupling is varied along the \( y \) direction. In this way we avoid artificial symmetrization needed in the previous work with nonhomogeneous Rashba coupling.

We employ material parameters for \( \text{In}_{0.5} \text{Ga}_{0.5} \text{As} \) alloy, i.e., \( m^* = 0.0465 m_0, g = -8.97 \). The Rashba coupling parameter \( \alpha = 0.572 \) \( \text{nm}^2 \) is adopted from Ref. 26.

### B. Lead eigenstates

We use a finite-difference approach with Hamiltonian Eq. (1) discretized on a grid with \( \Delta x = \Delta y = 4 \) \( \text{nm} \), employing gauge-invariant discretization of the kinetic energy operator for each of the spinor \( \Psi \) components:

\[
\frac{\hbar^2 k^2}{2m^*} \Psi_{i,j} = \frac{\hbar^2}{2m^* \Delta x^2} (4 \Psi_{i,j} - C_y \Psi_{i,j-1} - C_y^* \Psi_{i,j+1} - C_x \Psi_{i-1,j} - C_x^* \Psi_{i+1,j}),
\]

where \( \Psi_{i,j} = \Psi(x_i, y_j) \), \( C_y = \text{exp}[-i \frac{\hbar}{\Delta x} A_y] = \text{exp}[-i \frac{\hbar}{\Delta x} A_y] \) and \( C_x = \text{exp}[-i \frac{\hbar}{\Delta x} A_x] = 1 \). The derivatives in the SO Hamiltonian (3) are discretized straightforwardly.

We obtain the asymptotic states in the leads far away from the ring (i.e., for \( y = 0 \)), where \( -\hbar \frac{\partial}{\partial y}, H \) = 0 and the spinor of propagating wave can be written as

\[
\Psi(x, y) = e^{i k_y y} \begin{pmatrix} \Psi^+_1(x) \\ \Psi^-_1(x) \end{pmatrix},
\]

We insert this form of the spinor in the discretized Hamiltonian (1) and obtain a one-dimensional eigenproblem for transverse quantization in the lead. We plot the dispersion relation in Fig. 1(c). The energies of the (split by the Zeeman energy \( 2E_z \)) spin-up and spin-down states are plotted with the red and blue curves, respectively.
In the present work we consider the range of Fermi energies such that $E_f$ lies below the energy of the third subband [the dashed red curve in Fig. 1(c)]. In this case there are four possible values of the electron wave vector for a given electron Fermi energy. It can either belong to the lowest subband having $k_1$ or $-k_1$ and spin oriented along the $z$ direction or belong to the second subband having $k_1$ or $-k_1$ and spin oriented antiparallel to the $z$ direction.

We solve the transport problem through the solution of the stationary Schrödinger equation $H \Psi = E \Psi$ assuming boundary conditions such that the electron enters the system from the bottom electrode and can either be backscattered or be transmitted through the system. Details on the boundary conditions and on the approach applied to solve the transport problem can be found in the Appendix.

Unless stated otherwise in the calculations we lift the spin degeneracy of the states in the leads by applying a residual magnetic field with $B = 0.1$ mT, which does not induce any observable orbital effects.

III. RESULTS

A. Fano resonances

Let us start with the case of no SO coupling. In the ring there are two types of localized states: states with positive and negative parity with respect to the $y = 0$ axis. We inspect these states by diagonalization of Hamiltonian (1) for a closed system with varied length of the leads given by $L$ [see Fig. 1(c)] and plot the energy spectrum with black dotted curves in Fig. 2(b). Note that in fact each curve corresponds to the energies of two spin-opposite states. In Fig. 2(c) we plot the real part of the wave function of the spin-up states whose energies we mark by A, B, and C. The A and C states are the states of positive parity and due to the fact that their wave function is nonzero in the leads their energy changes with $L$. On the other hand the wave function of negative-parity state $B$ is zero in the leads and its energy is independent of $L$.

When the electron is transmitted through the ring the current-carrying state from the lead (which is a state of positive parity, i.e., the ground state of transverse excitation) couples to the localized states with positive parity. The conductance of the ring as function of electron Fermi energy [see black dotted curve in Fig. 2(a)] exhibits wide resonances due to this coupling. In the absence of SO interaction the localized states of the negative parity are bound; i.e., their lifetime is infinite in spite of the fact that their energy lies in the energy continuum—above the lowest subband transport threshold. On the other hand the energy of those states is still below the transport threshold for the next subband with wave functions of negative parity with respect to the axis of the channel.

When the SO coupling is introduced (we discuss first the case of weak Rashba coupling with $F_z = 50 \text{ kV/cm}$) the parity of the localized states is no longer well defined. For instance mean values of parity operator for states A, B, C are $0.939, -0.962, 0.957$, respectively. Due to the broken symmetry, the current-carrying state from the lead couples now to all the localized states. This results in the appearance of sharp peaks in the conductance, plotted by the red curve in Fig. 2(a), in addition to the wide resonances. These sharp peaks are Fano resonances with characteristic asymmetric dip-peak structures. Their energy corresponds to the energy of states localized purely in the ring. The small width of the resonances is reflected in a finite but long lifetime of the resonance state. In Fig. 2(b) the red curves present the energy spectrum of a closed system as function of $L$ in the presence of SO coupling. Note that now due to the fact that the states lack a well-defined parity anticrossings appear in the spectrum.

B. Spin transport

Let us now inspect spin transport through the ring. We consider the transmission of the electron with spin initially polarized along the $z$ direction (from the $k_1$ subband) and study the spin state at the output of the system.\(^{27}\) However, for the considered residual magnetic field $B = 0.1$ mT the transfer probabilities from both subbands are exactly the same and the spin at the output of the ring is exactly opposite.

In Fig. 3 we plot the conductance (black curve) and mean spin components at the output lead by solid colored curves. Notice that outside the resonance regions the output spin orientation remains unchanged when $E_f$ is varied—see the red and green curves—which is in agreement with the results of Ref. 6. However, when the electron Fermi energy is tuned to a resonance value the spin at the output is modified, as seen clearly in Figs. 3(b) and 3(c).

In order to further explore the spin changes in the narrow resonance regions let us increase the magnetic field, such the spin splitting of the localized resonance states is pronounced; namely, we apply $B = 0.02$ T. Now the transfer probabilities for an electron incoming from the subband $k_1$ and $k_2$ are no longer the same. In Fig. 4(a) with the purple curve we plot the transfer probability for an electron incoming from the lowest subband in the lead, i.e., $k_1$ with spin polarized parallel to the $z$ direction. With the green curve we show the transfer

![Figure 2](https://example.com/figure2.png)

FIG. 2. (Color online) (a) Conductance of the ring as function of the Fermi energy. (b) Energy spectrum for a closed system of a ring with leads of length $L$. In (a) and (b) black dotted curves represent results obtained for $F_z = 0$ and red solid curves for $F_z = 50 \text{ kV/cm}$. (c) Real part of the spin-up wave function corresponding to the states A, B, and C marked with blue arrows on (b) obtained without SO coupling for $L = 80 \text{ nm}$. Green contours depict the profile of the confinement potential.

235319-3
probability for an electron incoming from the second subband, i.e., \( k_y \) with spin polarized antiparallel to the \( z \) direction. We find that the Fano resonances from Figs. 2(a) and 3(a) are now converted to sharp peaks in the transfer probabilities occurring separately in both \( T_{k_x} \) and \( T_{k_y} \). In Fig. 4(c) we plot the energy spectrum of the closed system with varied length of the leads. Notice that the magnetic field splits the spin doublets of the localized states. The states have nonzero average spin components in the \( z \) direction (with \( |\langle s_z \rangle| \leq 0.9 \)). The states lying lower in energy have \( \langle s_z \rangle < 0 \); the states with higher energies have \( \langle s_z \rangle > 0 \). The splitting energy of the doublet is not equal to the Zeeman splitting as the spin and orbital parts of the wave function of the states are mixed by the Rashba coupling present in the ring. Namely, the energy of twice the Zeeman splitting is 

\[ 2E_z = |g\mu_B B| = 10 \mu eV \]

and the energy difference between the states whose energies we mark with orange and light blue curves is 12 \( \mu eV \) and for the pair plotted with blue and red curves the energy difference is 17 \( \mu eV \). The mean values of the spin operators in the \( x \) and \( y \) direction are zero.

The peaks in the transport probabilities for an electron with initially spin polarized parallel to the \( z \) direction [see purple curve in Fig. 4(a)] appear for energies equal to those of the resonance states marked with orange, blue, and green colors in Fig. 4(c), with positive \( \langle s_z \rangle \). For opposite spin orientation peaks are present for energies corresponding to the second state from the spin doublet (with energies marked with light blue, red, and purple curves), with negative \( \langle s_z \rangle \). This indicates that resonances in the transfer probability appear when the spin of the localized state matches the orientation of the spin of the incoming electron.

Let us now inspect the average spin components at the output of the system. In Fig. 4(b) we present the mean \( x \), \( y \), and \( z \) spin components multiplied by the conductance with blue, green, and red curves, respectively. Solid curves correspond to initial spin-up polarization and the dashed one to initial spin-down polarization. We observe that when the transported electron couples to the resonance states localized in the ring the spin at the output is close to the average spin of the resonance state (see the peaks marked with vertical dashed lines). Outside of the resonances we observe that the spin of the transferred electron deviates from the \( z \) or \(-z \) direction.

C. Dependence of the spin orientation on the Fermi energy for increased SO strength

Let us go back to the case of residual magnetic field \( B = 0.1 \) mT and inspect the dependence of the spin transport on the electron Fermi energy for increased strength of the SO coupling. In Fig. 5 we plot the mean spin components for four \( F_z \) values. We observe distinct broadening of the spin changes at the Fano resonances. The broadening of the resonances corresponds to a reduction of the lifetime of the resonance states in the ring.

Although the Fermi energy affects the spin evolution in the ring, the spin measured at the output (i.e., multiplied by the \( G \)) remains unchanged as function of \( E_f \) for \( F_z < 200 \) kV/cm. This is made clear by the contour maps of the mean spin components multiplied by the conductance presented in Fig. 6(a) for the \( y \) component and in Fig. 6(d) for the \( z \) component. Outside the narrow Fano region spin changes are marked by the blocked transport through the ring.

Figures 6(a) and 6(b) show that as the strength of the Rashba coupling is varied the conductance of the ring is changed due to the phase shift of the wave functions traveling in the left and right arm of the ring—the AC effect that modifies the conductance by a factor of

\[ G_{AC} = \frac{e^2}{h} \left[ 1 - \cos \left( \pi \frac{2Rm^2a|e|F_z}{\hbar^2} \right) \right]. \] (6)

For appropriately chosen strength of the SO coupling transport through the ring is quenched. For parameters taken in the present calculation the first AC oscillation minimum \( (G_{AC} = 0) \) is present around \( F_z = 200 \) kV/cm which can be observed.

FIG. 3. (Color online) (a) Conductance through the ring (black curve) and mean spin components (blue, green, and red solid curves) at the output lead for the electron with initial spin polarization along the \( z \) direction. (b), (c) Close-ups of the resonances in plot (a).

FIG. 4. (Color online) (a) Transfer probability for electron incoming with spin polarized parallel (purple curve) and antiparallel (green curve) to the \( z \) direction. (b) Mean values of the spin components in the \( x \), \( y \), and \( z \) directions at the end of output lead multiplied by the transfer probabilities. Solid curves are obtained for transport from \( k_y \), dashed curves for \( k_x \). (c) Eigenergies of closed system with leads of length \( L \). The results are obtained for \( B = 0.02 \) T.
in Figs. 6(a) and 6(d). On the other hand the first maxima of the AC oscillations appear for \( F_z = 0 \) and around \( F_z = 340 \text{ kV/cm} \) (note that the quenching of the conductance is found at slightly lower values of \( F_z \) for higher Fermi energy).

Similarly to the dependence of the spin on the electron Fermi energy observed before the first AC oscillation minimum \( (F_z < 200 \text{ kV/cm}) \) we find that the spin changes for \( F_z > 200 \text{ kV/cm} \)—after the first AC oscillation minimum [see Figs. 6(b) and 6(e)]. Similarly to the spin dependence in Fig. 5, they originate in the Fano resonances. Only now the resonances are associated with the second AC oscillation maximum around \( F_z = 350 \text{ kV/cm} \). Moreover now the dependence of the spin on the Fermi energy is no longer masked by the quenched transfer probabilities and it is visible in the contour maps of Figs. 6(a) and 6(d).

We conclude that the changes in the spin orientation originate at the Fano resonances appearing around and in the AC oscillation maxima and are broadened for SO coupling strength detuned from the AC oscillation maxima.

D. Comparison with one-dimensional model

In Ref. 6 it was found that when the electron is transferred through a one-dimensional spin-orbit-coupled ring

\[
E_f = \frac{2}{h} \left( \frac{\alpha R^2}{m^*} \right) \theta \]

the spin precesses around the \( y \) direction by an angle \( 2\theta = 2 \arctan(-\frac{F_z e}{\alpha 2m^* R^2/h^2}) \) which is independent of the electron Fermi energy. We calculated the spin after the rotation (taking \( R \) as the external radius of the ring, the area where the SO is present) about the angle \( 2\theta \) and plot the spin components in Figs. 6(c) and 6(f). We observe that before the first AC oscillation minimum both of the present results—of the two-dimensional model and the results for the one-dimensional ring—are similar with the exception of the spin resonances that are present in the first case. However, as the SO coupling strength is increased—the entanglement of the orbital and spin part of the wave function present for the system with finite-width channel increases—the discrepancy between our calculation and the results of the one-dimensional model increases. Namely, we find positive values of the \( y \) spin component, which is not seen in the one-dimensional approximation [note that the predictions of Eq. (6) for AC oscillation maxima and minima still hold for this strength of SO coupling].

E. Impact of the channel width and the ring radius

Let us now inspect the influence of the channel width on the spin behavior. We perform calculations in which we keep the mean radius of the ring constant and modify the channel width \( W \). We observe that the dependence of the spin orientation on the Fermi energy is changed. In Fig. 7 we plot the spin components at the output of the ring for different channel width. The wide resonances of the spin at the output as function of Fermi energy found for the wide-channel ring [see Fig. 7(c)] are transformed to sharp peaks for a ring with narrow channels [see Fig. 7(a)]. As the strength of the Rashba coupling and the mean radius of the ring are kept constant, we conclude that the dependence of the spin on the electron Fermi energy is an effect of the two-dimensional character of the channels which gets weaker for decreased \( W \); i.e., the resonances become narrower. This is in agreement with the intuition that for infinitesimal narrow channels the dependence of \( E_f \) should vanish (the peaks should be infinitesimally narrow) and the spin changes...
should not depend on the electron Fermi energy as found in the one-dimensional model. Note that in Fig. 7(c) we plot only the Fermi energy range below the energy of the third subband in the lead.

As expected from Eq. (6) the position of the AC oscillation minima and maxima with respect to the Rashba coupling strength is inversely proportional to the ring radius. In Fig. 8 we present the results obtained for a ring with mean radius \( R = 220 \) nm. First, we observe that spin components oscillate more frequently as function of the Fermi energy which can be attributed to the lower spacings between the energies of the localized states of a larger ring. Nevertheless, we observe a similar qualitative spin behavior with respect to the Fermi energy as in the case of a smaller ring [compare Figs. 8(b) and 8(e) with Figs. 6(b) and 6(e)]. Only now we are able to observe more periods of AC oscillations. We again find that the spin changes originate from the Fano resonances, which appear when the SO strength is tuned to the AC conductance oscillation maxima, i.e., for \( F_z \simeq 185 \) kV/cm and \( F_z \simeq 320 \) kV/cm. However, in terms of the mean spin values multiplied by the conductance, we observe a dependence on the Fermi energy only for strong SO coupling (namely \( F_z > 200 \) kV/cm) similarly to the case of a smaller ring. Also the correspondence with the one-dimensional results [see Figs. 8(c) and 8(f)] remains the same as for smaller ring radius.

**IV. SUMMARY AND CONCLUSIONS**

In summary, we have studied the spin and charge transport through a spin-orbit-coupled quantum ring with two-dimensional channels. We found Fano resonances of the conductance which are present for nonzero SO coupling strength tuned to the maxima of AC oscillations. These narrow resonances are an effect of coupling of the current-carrying states from the leads with the localized states in the ring that have a long lifetime. The coupling is possible due to the breaking of the parity of these states by the SO interaction.

In the Fano resonances spin transport through the ring is modified. We argue that the modification is due to the coupling of the transferred electron spin with the spin of the resonance states which we investigated in the presence of finite external magnetic field.

When the SO coupling strength is such that system is outside the AC oscillation maxima the Fano resonances are broadened. In that case the spin modification is translated into a wide dependence of the spin transport on the Fermi energy. The latter result is in contrast to the findings of the one-dimensional model\(^6\) which employed spin transformations (independent of \( E_f \)) performed in a quantum ring to realize a universal set of quantum gates. However, when the width of the channels is decreased the resonances that result in the dependence of spin transport on the Fermi energy become narrower; the results tend toward the prediction of the one-dimensional model.

Moreover, by the direct comparison of the results of the one-dimensional model and the two-dimensional calculation we found that for strong SO coupling the spin evolution proves to behave in a way exceeding the predictions for a one-dimensional ring even outside the resonance region.

**ACKNOWLEDGMENTS**

This work was supported by the “Krakow Interdisciplinary PhD Project in Nanoscience and Advanced Nanostructures” operated within the Foundation for Polish Science MPD Programme cofinanced by the EU European Regional Development Fund, Project No. N N201103938 supported by the Ministry of Science and Higher Education (MNiSW) for 2010–2013, the Belgian Science Policy (IAP), and the Flemish
From the discretized form of the derivative vector of incoming electron $i k$ the electron enters the system with a given wave vector and the method which are used for a solution of the electron transport through system described by Hamiltonian (1).

**APPENDIX**

This Appendix describes the applied boundary conditions and the method which are used for a solution of the electron transport in the output lead. Let us subtract from the derivative positive wave vectors (as we assume no backscattered waves in such channel is a combination of channel eigenstates with energy). (corresponding to a given energy) and can exit the system in part by PL-Grid Infrastructure.

**1. Boundary conditions**

The boundary conditions applied in the method assume that the electron enters the system with a given wave vector $k_{inc}$ and the method which are used for a solution of the electron in the output lead). Let us subtract from the derivative waves with negative wave vector. For instance, for transport condition in the bottom of the computational box. Only now we have $i k_{inc} \Psi(x, y)$ and the sum includes a positive wave vector of incoming electron $k_{inc}$ and the two backscattered waves with negative wave vector. For instance, for transport of an electron with wave vector $k_{inc}$,

$$\frac{\partial \Psi(x, y)}{\partial y} = \sum_{k > 0} i k c_{k}^{\text{out}} \exp[i k y](\psi_{1}^{\dagger}(x)) \psi_{1}(x)$$  \hspace{1cm} (A1)

$i k_{inc} \Psi(x, y)$, obtaining

$$\frac{\partial \Psi(x, y)}{\partial y} = \sum_{k > 0} i (k - k_{inc}) c_{k}^{\text{out}} \exp[i k y](\psi_{1}^{\dagger}(x)) \psi_{1}(x) + i k_{inc} \Psi(x, y).$$ \hspace{1cm} (A2)

From the discretized form of the derivative

$$\frac{\partial \Psi(x, y)}{\partial y} = \frac{\psi(x, y + \Delta y) - \psi(x, y - \Delta y)}{2\Delta y},$$ \hspace{1cm} (A3)

we obtain

$$\psi(x, y + \Delta y) = 2\Delta y \sum_{k > 0} i (k - k_{inc}) c_{k}^{\text{out}} \exp[i k y](\psi_{1}^{\dagger}(x)) \psi_{1}(x) + \psi(x, y - \Delta y) + 2\Delta y i k_{inc} \Psi(x, y).$$ \hspace{1cm} (A4)

The same procedure leads to the form of the boundary condition in the bottom of the computational box. Only now we add $i k_{inc} \Psi(x, y)$ and the sum includes a positive wave vector of incoming electron $k_{inc}$ and the two backscattered waves with negative wave vector. For instance, for transport of an electron with wave vector $k_{inc}$,

$$\psi(x, y - \Delta y) = -2\Delta y \sum_{k} i (k + k_{inc}) c_{k}^{\text{in}} \exp[i k y](\psi_{1}^{\dagger}(x)) \psi_{1}(x) + \psi(x, y + \Delta y) + 2\Delta y i k_{inc} \Psi(x, y).$$ \hspace{1cm} (A5)

We use the above forms of the wave functions to obtain the boundary conditions, i.e., $\psi(x, y + \Delta y)$ for the top edge of the computational box and $\psi(x, y - \Delta y)$ at the bottom of the computational box (at the left and right edge of the mesh we assume $\psi = 0$). The used boundary conditions are transparent; i.e., the transport results do not depend on the length of the leads.

**2. Solving the transport problem**

We solve the system of equations produced by the discretization of the Schrödinger equation with the boundary conditions described above. In the present method on the one hand the amplitudes $c_{k}^{\text{in}}$ and $c_{k}^{\text{out}}$ are required for the boundary condition and on the other hand they can be obtained from the solution of the Schrödinger equation. Thus we assume initial values of the amplitudes (namely, $c_{k}^{\text{in}} = c_{k}^{\text{out}} = 1/\sqrt{2}$ for $k_{inc} = k_{1}$; however, we checked that the particular choice of the initial values does not change the final result) and put them into the boundary conditions. Then we solve the Schrödinger equation. From the solution we extract new values of the amplitudes by projection (in the input and output leads) of the function

$$\psi(x, y) = \sum_{k} c_{k} \exp[i k y](\psi_{1}^{\dagger}(x)) \psi_{1}(x)$$ \hspace{1cm} (A6)

(which accounts all possible wave vectors for a given energy) onto the solution and solve again the Schrödinger equation. Such procedure is repeated until convergence is reached—the extracted amplitudes do not change in the subsequent iterations, and the amplitudes $c_{k}$ are such that in the input channel there is only one incoming wave and in the output lead there are no backscattered waves.

We calculate transport probability from the ratio of the probability currents $j_{k}$ in the leads for respective wave vectors:

$$T_{k \rightarrow k'} = \frac{|c_{k}^{\text{out}}|^{2} j_{k'}}{j_{k}}.$$ \hspace{1cm} (A7)

The conductance $G$ is calculated as a sum of the transmission probabilities over available subbands; i.e., $G = \sum_{k} \sum_{k'} j_{k} T_{k \rightarrow k'}$. The described approach allows one to study the electron transport for a given Fermi energy in contrast to the methods involving transmission of a wave packet that consists of the superposition of $H$ eigenstates. Also, as the approach is based on an exact solution of the Schrödinger equation it naturally includes evanescent modes that can appear in the ring.

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27Mean values of the x, y, and z spin components are calculated on a normalized wave function on a stripe at the end of the output lead.