Using Topological Insulator Proximity to Generate Perfectly Conducting Channels in Materials without Topological Protection

Sven Essert, Viktor Krueckl, and Klaus Richter

Institut für Theoretische Physik, Universität Regensburg, D-93040 Regensburg, Germany

(Dated: February 6, 2014)

We show that hybrid structures of topological insulators and materials without topological protection can be employed to create perfectly conducting channels hosted in the non-topological part. These states inherit the topological protection from the proximity of the topological insulator but are more fragile to time-reversal symmetry breaking because of their extended character. We explore their formation in the band structure of model hybrid systems as well as realistic heterostructures involving HgTe/CdTe-based two-dimensional topological insulators. Using numerical quantum transport calculations for the HgTe/CdTe material system we propose two experimental settings which allow for the detection of the induced perfectly conducting channels, both in the localized and diffusive regime, by means of magneto conductance and shot noise.

PACS numbers: 72.10.-d, 73.20.At, 73.21.Ac

One important feature of the topological classification of insulators [1] is the existence of gapless states at interfaces between materials which differ in their topological quantum number. In this way, one can understand the formation of quasi one-dimensional (1d) edge channels in quantum Hall systems and in two-dimensional (2d) topological insulators (TI) at the interface between the topologically non-trivial insulator and the trivial insulating vacuum. These edge channels carry the unique property of being perfectly conducting, even in presence of impurity scattering, corresponding to transmission eigenvalues of one. Such quantized, spin-polarized channels along the boundaries of a 2d TI crystal have already been detected experimentally [2] and give rise to peculiar conductance phenomena such as the quantum spin Hall effect [3, 4].

In this paper we show that perfectly conducting channels (PCCs) can be induced in materials without topological protection by the proximity of a 2d TI as sketched in Fig. 1. In such a hybrid structure the non-topological material could be a (disordered) metal, a gated semiconductor or even a conventional Anderson insulator. Such a setup exploits the hybridization of the upper TI edge state with the bulk states of the normal material, leading to an imbalance of left and right moving states in the normal material and thereby to a PCC therein [5, 6]. The emergence of PCCs in isolated non-topological systems has been earlier predicted for carbon nanotubes [6] and graphene nanoribbons [7, 8]. There the PCC was shown [7] to arise from an uneven separation of left and right moving states associated with specific (valley) symmetries [8]. However, since atomic defects or short-ranged impurity potentials break this symmetry, PCCs are rather fragile in these graphene-based systems and have not yet been observed therein. Contrarily, the proximity-induced PCCs reported here are robust against any kind of disorder as long as time-reversal symmetry is preserved, even in presence of spin-orbit coupling.

To our knowledge this phenomenon has not been discussed before in 2d heterostructures, only layers made of 2d TIs nested in 3d systems [9-11] and pure insulator heterostructures [12] have been considered very recently.

In the following, we first demonstrate the emergence of a proximity-induced PCC by means of a band structure analysis within a simplified model calculation. Subsequently, we will present numerical results that show the same features in HgTe/CdTe-based ribbons. Furthermore, we will study features of the PCC in quantum transport for experimentally realizable setups. In particular we reveal the creation of a PCC in an otherwise insulating disordered strip and how to disentangle the signature of a PCC from the diffusive modes in a disordered conductor.

Throughout this manuscript we model the electronic structure of a 2d HgTe/CdTe quantum-well by the four-band Bernevig-Hughes-Zhang (BHZ) Hamiltonian [3],

\[ H = \begin{pmatrix} h(k) & 0 & -\Delta \\ 0 & \Delta & 0 \\ -\Delta & 0 & h^*(-k) \end{pmatrix}, \] (1)

with spin-subblock Hamiltonians

\[ h(k) = \begin{pmatrix} M - (B + D)k^2 & -Ak_+ \\ Ak_- & -M + (B - D)k^2 \end{pmatrix}, \] (2)

where \( k_\pm = k_x \pm ik_y \) and \( k^2 = k_x^2 + k_y^2 \). In Eq. (1), the off-
diagonal coupling parameter $\Delta$ describes a Dresselhaus-type spin-orbit coupling of the two spin blocks [13].

We consider a 2d heterostructure of a usual conductor and a semi-infinite topological insulator plane. It is instructive to first look at a simplified model system which already captures the important physical features. The metal part is modeled by the Hamiltonian (1) with $A = D = M = \Delta = 0$, describing a free electron-hole gas. Confined to a finite-width strip by hard wall boundary conditions, this leads to an electron-hole symmetric parabolic band structure, as shown in Fig. 2(a). For the TI half plane, we use a minimal BHZ Hamiltonian $(D = \Delta = 0)$, which splits up into two electron-hole symmetric spin blocks. The band structure for one spin block shown in Fig. 2(a) exhibits the bulk band gap and a single edge state. A coupling between the two systems is implemented by replacing the central hard wall constraint by the continuity of the wave function $\Psi(y)$ and the flux $v_y \Psi(y)$ across the interface. The resulting band structure, shown in Fig. 2(b), resembles the one of the metallic strip, but contains an extra band (marked in blue) as a remainder of the topological edge state, which does not feature the linear momentum dependence because of band repulsion due to hybridization. The wave functions at various energies in Fig. 2(c) reveal that all states are predominantly localized in the metallic region and only small exponential tails remain in the TI part. Thus, we do not find evidence for an induced non-trivial topology as reported in Ref. 12, from which one would expect a new interface state forming on the outer hard wall boundary. Instead, the former edge state fully hybridizes with the metallic bands and covers the whole extended state region! This is in line with Ref. 10 where the authors also observe an extended metallic state in a 3d heterostructure, but differs from the findings of Ref. 11, where a protection from hybridization for topological edge states is reported for a 2d quantum Hall insulator on top of a 3d substrate.

The above result has important consequences. Counting the bands in Fig. 2(b) reveals an imbalance between left and right movers (up and down movers in Fig. 2), which leads to a PCC at any Fermi energy. Including the other spin block and spin-orbit interaction only slightly changes this picture. Since both spin blocks feature a PCC with opposite propagation direction in the decoupled system, the combined system has an equal but odd number of left and right moving channels. Due to the time-reversal symmetry of the Hamiltonian, the scattering matrix can be written in a basis where it is antisymmetric. Together with the odd number of channels, this implies a single PCC in each direction [15]. As the wave functions of all channels are predominantly localized in the metallic region, one can say that the proximity of the TI induces a single additional channel in each direction leading to one net PCC in the metallic region. In other words, the quasi 1d edge state of the TI migrates into the metallic region and forms an extended PCC. We believe that this is generally true for materials with extended states. It should provide a valuable tool to create and observe PCCs as they inherit the topological protection of the former edge channels and are expected to be stable for any disorder and interface type, as long as the latter allows for sufficient hybridization.

Given that the TI is a good insulator, i.e., that the decay length of the states inside the insulator is short compared to the width of the metallic ribbon, one expects that only a vanishing part of the probability density of a state is inside the TI. This allows for replacing the explicit consideration of the TI by an effective boundary condition for the wave function $\Psi(y)$ inside the metallic ribbon. For the right boundary of a heterostructure in $y$-direction it reads:

$$\Psi_2 = -\Psi_1,$$

$$i (v_y \Psi)_2 - i (v_y \Psi)_1 = 4Qk_x \Psi_1,$$  

where $(v_y \Psi)_i$ denotes the $i$-th component of the state obtained by applying the velocity operator normal to the boundary. Note that, Eq. (36) still contain a free parameter $Q$, which stems from the band curvature of the TI. The effect $Q$ is best seen by applying the boundary conditions to a topologically trivial insulator, like a gapped electron-hole gas: Then the group velocity of the emergent edge state band linearly depends on $Q$. Using the boundary conditions on the gapless electron-hole gas from the previous model heterostructure, Fig. 2(a), we obtain for $Q = B$ qualitatively the same band structure as in Fig. 2(b). While Eq. (3) is a simple Dirichlet boundary condition that fixes the phase of the two components, Eq. (36) is of the Robin type [16], i.e., it links

---

**FIG. 2.** (Color online) Band structure of (a) a non-topological strip (left) and a semi-infinite topological insulator plane (right), (b) the combined system. (c) Wave functions for parameters marked by dots in (b). Solid/Dashed lines show the first/second component of the wave functions.
FIG. 3. (Color online) (a) Scheme of the four-lead geometry used for PCC observation (b)–(d) Band structures of a nanoribbon ($W = 350$ nm, HgTe material parameters [17]) with different band topology [(b) $M = -10$ meV, (c) $M = 0$ meV, (d) $M = 10$ meV] in proximity to a TI ribbon (energy offset $E_0 = 30$ meV). Colors encode the position of the corresponding wave function (red: metallic part, blue: TI).

the wave function at the boundary to its derivative. This $k_x$-dependent mixing induces the PCC in the adjacent metal, as illustrated in the Supplemental Material [17].

For the practical observation of a PCC in such heterostructures the semi-infinite TI plane can be replaced by a TI-ribbon in a four-lead geometry, see Fig. 3(a). This geometry decouples the edge channel on the lower edge of the TI from the region with the PCC. This is necessary, as the lower edge state would add another channel to the scattering matrix, make it even dimensional and therefore lift the protection of the PCC. If the PCC in the upper metal is well coupled to the two metallic leads (e.g., by making the leads wide), the total transmission $T_M$ between the upper leads will be at least one. Even in presence of disorder there will be no complete localization independent of the system length.

It is not required to assemble the heterostructure presented in Fig. 3(a) from two different materials. For convenience the whole system can consist of a common HgTe/CdTe quantum well with inverted band order, where only the upper (metallic) part has a Fermi energy outside the topological band gap, such that propagating states exist in the bulk of the corresponding clean material. For strong disorder associated with a localization length shorter than the length $L$ of the ribbon connecting the left and right terminals, the total transmission $T_M$ between the upper leads and $T_{TI}$ between the lower leads is strongly suppressed. To achieve TI properties in the lower part of the strip, a gate is applied to locally shift the band structure of a region containing the edge and the entire connection between the strip and the lower leads [red region in the inset of Fig. 4(a)]. To illustrate the signatures of the PCC the transport properties of this setup are calculated for different gate potentials $V_g$ by means of a recursive Greens function algorithm [18]. If the gated part is tuned such
that the Fermi energy is in the TI bandgap, the suppressed transmission $T_{TI}$ of the localized regime takes the quantized value $T_{TI} \approx 1$ of the quantum spin Hall edge state. This behavior of $T_{TI}$ is shown in Fig. 4(a) by the solid (blue) line for gate voltages between 30 mV and 50 mV. Remarkably, the transmission $T_M$ between the two upper leads exhibits the same behavior and gets quantized, although there is no material with topological protection linking the two terminals. This quantized transmission in the otherwise localized regime is a clear-cut signature of and carried by the PCC.

The different nature of the PCC compared to the quantum spin Hall edge state can be probed by a perpendicular magnetic field $B$. Since the PCC state is extended over the entire spatial region $A$ of the upper strip, which can be seen in the local density of states [17], a weak field generating a flux $A \cdot B \approx \phi_0$, where $\phi_0$ is the flux quantum, is sufficient to effectively break the time-reversal symmetry. The numerical data presented in Fig. 4(b) shows that the conductance quantization of $T_M$ is fully destroyed for a magnetic field of $B = 2$ mT, corresponding to $\phi_0$ penetrating the device. This is not the case for the robust quantum spin Hall edge state, which is strongly localized at the lower boundary and accordingly quasi one-dimensional. As a result, the magneto transport is insensitive to magnetic fields in the mT regime, as shown for $T_{TI}$ in Fig. 4(c). As a side remark, the PCC turns out to be completely spin mixed already at small spin-orbit strengths Delta, in contrast to the quantum spin Hall edge channels, which remain spin-polarized up to moderate spin-orbit strengths [17].

If the electronic states of the disordered strip are not localized the total transmission $T_M = \sum_n t_n$ also contains contributions $0 < t_n < 1$ of non-perfectly conducting channels. For diffusive transport the distribution $P(t)$ of these transmission eigenvalues $t$ features a bimodal distribution [19, 20], see upper left panel of Fig. 5(a), whose intermediate transmission eigenvalues mask the PCC when considering $T_M$. In this case an additional tunnel barrier, depicted by the green bar in the sketch of Fig. 5, can be used to unravel the presence of the PCC. The probability for finding these intermediate transmission eigenvalues can be strongly reduced by the tunnel barrier (parametrized by $V_t$), as shown in Fig. 5(a) for $B = 0$ mT, $V_t = 40$ mV. Hence, for this configuration $T_M$ is mainly carried by the PCC. As in the localized case, protection of the PCC can be broken by a magnetic flux leading to a broader distribution for the high transmission eigenvalues [right panel in Fig. 5(a)].

The change of this probability distribution is also reflected in transport quantities, which can be experimentally probed. For example, upon raising the tunnel barrier $V_t$ the transmission at $B = 0$ exhibits a quantized minimum, which reflects the presence of a PCC, as shown by the red line in Fig. 5(b) for $V_t \in [30, 60]$ mV. As expected, this feature disappears at small magnetic fields.

A more pronounced PCC signature is found in the shot noise power [21], $P = \sum_n t_n(1 - t_n) \approx 11$, depending on the transmission eigenvalues $t_n$ between the upper leads. For conventional diffusive transport the bimodal distribution leads to a universal 1/3 suppression of the shot noise $P = T/3$, as marked by a dashed line in Fig. 5(c), independent of the universality class of the material [22]. In presence of a tunnel barrier all $t_n$ except the one of the PCC tend to zero, leading to a characteristic shot noise suppression, as shown by our numerics in Fig. 5(c) for $V_t \in [30, 60]$ mV. A $B$-field destroys the PCC, which in turn removes the shot noise suppression. At finite $B$, in absence of the PCC the shot noise signal even increases above $P = T/3$ due to the tunnel barrier. We suggest the tunnel conductance and shot noise as promising, experimentally accessible observables to verify the TI proximity induced PCC.

To summarize, we showed that the proximity of a 2d topological insulator (TI) creates a robust perfectly conducting channel (PCC) in an adjacent non-topological material. We believe this is a quite general effect that should work for almost all materials with extended states.
We expect that a similar phenomenon exists for a disordered conductor at the surface of a 3d TI. In addition, we showed that the proximity of a 2d TI to a metal can be understood in terms of a Robin type effective boundary condition, responsible for the creation of the PCC. This does not only allow simplifications in future theoretical studies, but might also pave the way to induce PCCs by artificially creating such boundary conditions, e.g., by using metamaterials for electromagnetic waves, without relying on TI heterostructures.

This work is supported by DFG (SPP 1666 and joined DFG-JST Research Unit FOR 1483). We thank I. Adagideli and M. Wimmer for useful conversations.

[5] Strictly speaking, the argument of channel imbalance only applies if the spin is conserved. However, the PCC also persists in presence of spin-orbit coupling as long as time-reversal symmetry holds, since such a setting leads to an odd number of channels in each direction.
[14] We use the term “metal” synonymously to “material with extended states” also referring to a gated semiconductor.
[17] See Appendix for the derivation of the effective boundary condition and details about the numerical model and spatial densities of the PCC.

**Band structure calculations**

This section provides some details on the band structure calculations shown in Fig. 2 of the main manuscript. We consider a metallic strip of finite width $W$, which borders vacuum on the one side and is in contact with a semi-infinite 2d-TI plane on the other. In the TI part $H_{\text{TI}}$ is given by a simplified 2-band BHZ-model

$$H_{\text{TI}} = \begin{pmatrix} M - Bk_x^2 & A(k_x + ik_y) \\ A(k_x - ik_y) & -M + Bk_x^2 \end{pmatrix},$$  

where in the metallic region $H_{\text{met}}$ is simply taken as a free electron-hole gas ($A = M = 0$)

$$H_{\text{met}} = \begin{pmatrix} -Bk_x^2 & 0 \\ 0 & Bk_x^2 \end{pmatrix}.$$  

For simplicity we assume the value of the parameter $B$ to be the same in the two regions. To solve for the transverse wave functions in this geometry, we require that the wave function $\psi_{k_y}(y)$ should vanish at the boundary to vacuum at $y = 0$ and that $\psi_{k_y}(y)$ should be continuous across the metal-TI boundary at $y = W$. We also require the probability current to be continuous across the boundary, i.e.,

$$\left( \begin{array}{cc} 2iB\partial_x & 0 \\ 0 & -2iB\partial_y \end{array} \right) \psi_{k_x}(W - \epsilon) = \left( \begin{array}{cc} 2iB\partial_x & iA \\ -iA & -2iB\partial_y \end{array} \right) \psi_{k_x}(W + \epsilon).$$

We then find the secular equation

$$4B^2 f_p \left( \lambda_2 u_1 v_2 - \lambda_1 u_2 v_1 \right) + (4B^2 - A^2) \lambda_1 \lambda_2 \left( u_2 v_1 - u_1 v_2 \right) - 2AB \left( \lambda_1 - \lambda_2 \right) \left( u_1 u_2 - v_1 v_2 \right) \frac{f_m\cosh(f_m W) \sinh(f_p W)}{f_m f_p \cosh(f_m W) \cosh(f_p W)}$$

$$= 4B^2 f_p f_m \left( u_1 v_2 - u_2 v_1 \right) + 4B^2 f_m \left( \lambda_2 u_2 v_1 - \lambda_1 u_1 v_2 \right) \frac{\sinh(f_m W) \sinh(f_p W)}{\sinh(f_m W) \cosh(f_p W)},$$

which may be solved numerically to yield the band structure. Here,

$$f_p = \sqrt{k_x^2 + \frac{E}{B}},$$  

$$f_m = \sqrt{k_x^2 - \frac{E}{B}},$$

and $u_{1/2}, v_{1/2}$ are the components of the decaying free solutions $\xi_{k_x}^{1/2}(y)$ in the TI-region,

$$\xi_{k_x}^{1/2}(y) = \begin{pmatrix} u_{1/2} \\ v_{1/2} \end{pmatrix} e^{\lambda_{1/2} y},$$

$$\lambda_{1/2} = \frac{\sqrt{B^2 + E^2}}{2B}.$$
where $H_{TI} \xi_{kx}^{1/2}(y) = E \xi_{kx}^{1/2}(y)$. The decay coefficients (which have a negative real part as long as $|E| < |M|$) are given by

$$\lambda_{1/2} = -\sqrt{\frac{A^2 + 2B^2k_x^2 - 2BM + \sqrt{A^2 + 4B^2E^2 - 4A^2BM}}{2B}}.$$

Plots of a typical band structure obtained from this model are shown in Fig. 2(b) of the main text together with some sample wave functions in Fig. 2(c). They were obtained using $A = 5, B = -100, M = -0.05, W = 1000$.

**Effective boundary condition**

**Derivation**

One can increase the level of abstraction and still observe the same physics, by not explicitly considering the semi-infinite 2d TI plane but instead replacing it by an effective boundary condition. This is especially justified if the extent of the evanescent wave component into the 2d-TI is small ($\frac{1}{|\text{Re}\lambda_{1/2}|} \ll 1$), such that the transverse wave function is to a good approximation only located inside the metallic region. As the considered model of the TI involves many parameters, there is no unique way of achieving this infinitely fast decay and hence there will be no unique set of boundary conditions. The limiting procedure presented in the following should be understood as one possible way, which will in the end provide a conceptually simple boundary condition, which will capture all important physical features.

For achieving the infinite decay constants we let $M \to -\infty$, i.e., we will consider a TI with an infinite band gap. In addition, we will have to rescale $A$ by choosing $A = X \cdot \sqrt{-M}$ with a parameter $X$ having the units of energy/length. If one now considers the limit $M \to -\infty$ keeping $X = \frac{A}{\sqrt{-M}}$ = const, one finds for the real parts of the decay constants $\lambda_{1/2}$:

$$\lim_{M \to -\infty} \text{Re} \lambda_1 = -\infty, \quad \lim_{M \to -\infty} \text{Re} \lambda_2 = -\infty,$$

meaning that we indeed expect the wave functions not to extend into the TI. For this limit, one can derive the appropriate boundary conditions that any attached two-band Hamiltonian should fulfill. We start from the continuity of the wave function of the metal region $\psi^M(y)$ and the probability current $v_y \psi^M(y)$ across the boundary at $y = W$,

$$\begin{align*}
(\psi_1^M(W)) &= c_1 (u_1) + c_2 (u_2) \\
(\psi_2^M(W)) &= c_1 (u_1) - c_2 (u_2) - i v_y (u_1)
\end{align*} \quad \ldots \quad (17)$$

$$\begin{align*}
(-i) v_y \psi_1^M(W) &= c_1 2B \lambda_1 (u_1) + c_2 2B \lambda_2 (u_2) \\
&+ A c_1 (v_1) + A c_2 (v_2)
\end{align*} \quad \ldots \quad (18)$$

for so far unknown constants $c_1$ and $c_2$. We are now interested in the behavior of the right hand sides for the limit discussed above. However, as

$$\lim_{M \to -\infty} |u_{1/2}| = \lim_{M \to -\infty} |v_{1/2}| = 0,$$

one needs to rescale the constants to obtain a finite value of the wave function at the border to the TI:

$$c_{1/2}' = \frac{c_{1/2}}{\sqrt{-M}}.$$

This way, we find

$$\begin{align*}
\psi_1^M(W) &= c_1' \sqrt{-M} u_1 + c_2' \sqrt{-M} u_2 \\
\psi_2^M(W) &= c_1' \sqrt{-M} u_1 + c_2' \sqrt{-M} v_2,
\end{align*} \quad \ldots \quad (21)$$

$$\ldots \quad (22)$$

from Eq. (17). We can now do the limiting process on the right hand side always keeping $X = \frac{A}{\sqrt{-M}}$ fixed, using

$$\begin{align*}
\lim_{M \to -\infty} u_1 \sqrt{-M} &= \frac{1}{2} (X + \sqrt{4B + X^2}) \\
&\quad \equiv q_1(B, X) \\
\lim_{M \to -\infty} v_1 \sqrt{-M} &= -\frac{1}{2} (X + \sqrt{4B + X^2}) \\
&\quad \equiv q_1(B, X) \\
\lim_{M \to -\infty} u_2 \sqrt{-M} &= \frac{1}{2} (X - \sqrt{4B + X^2}) \\
&\quad \equiv q_2(B, X) \\
\lim_{M \to -\infty} v_2 \sqrt{-M} &= -\frac{1}{2} (X - \sqrt{4B + X^2}) \\
&\quad \equiv q_2(B, X)
\end{align*} \quad \ldots \quad (23)$$

$$\ldots \quad (26)$$

where, on the way, we made use of the fact that $B < 0$ and $X > 0$. To simplify the writing we introduced the functions $q_1(B, X)$ and $q_2(B, X)$. From this, we can already find the first boundary condition, as now

$$\begin{align*}
\psi_1^M(W) &= c_1' q_1(B, X) + c_2' q_2(B, X) \\
\psi_2^M(W) &= -c_1' q_1(B, X) - c_2' q_2(B, X), \quad \ldots \quad (27)$$

$$\psi_1^M(W) = -\psi_2^M(W). \quad \ldots \quad (28)$$

Note that this is independent of the choice of the parameters $B$ and $X$. The second boundary condition which links the currents is, however, more complicated and does depend on the choice of $B$ and $X$. It can be simplified by additionally choosing $X = 2\sqrt{-B}$, which will then imply

$$q_1 \left(B, X = 2\sqrt{-B}\right) = q_2 \left(B, X = 2\sqrt{-B}\right) = \sqrt{-B}. \quad \ldots \quad (29)$$

$$\ldots \quad (30)$$
If we now subtract the two components of Eq. (18),
\[
(i (v_y \psi^M)_2 (W) - i (v_y \psi^M)_1 (W)) = c'_1 (2B \lambda_1 + A) (u_1 + v_1) \sqrt{-M} + c'_2 (2B \lambda_2 + A) (u_2 + v_2) \sqrt{-M}.
\]
(31)
and again perform the limits,
\[
\lim_{M \to -\infty} (2B \lambda_1 + A) (u_1 + v_1) \sqrt{-M} = 4B \sqrt{-B} k_x (32)
\]
\[
\lim_{M \to -\infty} (2B \lambda_2 + A) (u_2 + v_2) \sqrt{-M} = 4B \sqrt{-B} k_x (33)
\]
we obtain
\[
i (v_y \psi^M)_2 (W) - i (v_y \psi^M)_1 (W) = 4B k_x \left(c'_1 \sqrt{-B} + c'_2 \sqrt{-B}\right).
\]
\[
= 4B k_x \psi^M_1 (W)
\]
\[
= 2B k_x (\psi^M_1 (W) - \psi^M_2 (W))
\]
(3)
Here, we already made use of the first boundary condition. In total we find, that the boundary to a TI in the limit discussed above can be approximately described by the following boundary conditions, which still contain one free parameter,
\[
\psi_1 (W) = -\psi_2 (W)
\]
\[
i (v_y \psi)_1 (W) - i (v_y \psi)_2 (W) = 2Q k_x (\psi_1 (W) - \psi_2 (W)) (36)
\]
which we renamed from \(B\) to \(Q\), to distinguish it from the band curvature of the material on which we enforce this boundary condition. Like this we can generalize the case discussed in Sec. and need not restrict ourselves to heterostructures of materials with the same band curvature. Fig. (6) shows band structure calculations using the boundary conditions with different choices for the parameter \(Q\). One notes that for \(Q = B\) one almost recovers the band structure from Fig. 2(b) of the main text which was obtained by a full calculation which explicitly includes the TI half-plane.

In the derivation, we assumed the TI half plane to extend in the positive \(y\)-direction, which is why we matched the exponentially decaying wave functions on the TI side. Of course the same calculation can in principle be redone for an arbitrary interface orientation. E.g., choosing the TI plane to extend in the negative \(y\)-direction with a boundary at \(y = 0\), leads to the following set of boundary conditions:
\[
\psi_1 (0) = \psi_2 (0) \quad (37)
\]
\[
i (v_y \psi)_1 (0) + i (v_y \psi)_2 (0) = -2B k_x (\psi_1 (0) + \psi_2 (0)) (38)
\]
i.e., we find an additional phase factor in front of one wave function component.

Effect of Robin boundary condition

To show that it is indeed the Robin boundary condition, Eq. (36) or (38), which is responsible for the appearance of a perfectly conducting channel, we consider a one-component free electron gas Hamiltonian
\[
H = B (k_x^2 + k_y^2)
\]
subject to hard wall boundary conditions at \(y = 0\) \((\psi(0) = 0)\) and a one-component version of the boundary
FIG. 7. Band structure of a wave guide with a free electron gas Hamiltonian, Eq. (39), subject to hard wall boundary conditions at \( y = 0 \), and the Robin boundary condition from Eq. (41) at \( y = W \) using \( \frac{1}{2}Q = B = 1 \).

condition from Eq. (36) at \( y = W \):

\[
\begin{align*}
  i (v_y \psi) (W) &= 2B \partial_y \psi (W) = -2Qk_x \psi (W) \\
  \rightarrow \partial_y \psi (W) &= -\frac{Q}{B} k_x \psi (W).
\end{align*}
\]

For \( |Q| > |B| \) one obtains a band structures similar to the one shown in Fig. 7, which has been calculated by setting \( \frac{1}{2}Q = B = 1 \). It resembles the simple parabolic band structure that one obtains by putting hard wall (or Neumann) boundary conditions on both sides, but it contains an extra band which tends to negative energies with increasing \( k_x \) and thereby creates a PCC. While the emergence of large negative eigenvalues is known for Robin problems \( \partial_y \psi = -\alpha \psi \) with large negative parameters \( \alpha \) [23], as far as we know there has not been a discussion on the emergence of perfectly conducting channels in one-sided Robin wave guides so far.

By implementing this boundary condition, or a similar one (e.g. by replacing \( k_x \) in Eq. (41) by a different but still odd function of \( k_x \)), on one side of a wave guide, one should be able to create PCCs for many scenarios, i.e., for quantum as well as electromagnetic wave guides.

Transport Calculations

In the manuscript we use numerical quantum transport calculations to investigate the transport signatures of a PCC, which is induced in a conventional metal by the proximity of a TI. To this end, we propose two different setups, which will be introduced in more detail in the following. Both of them are based on a common HgTe/CdTe heterostructure, described by the BHZ Hamiltonian (1) using the material parameters summarized in Table I.

As first transport example, we simulate a H-shaped structure consisting of a long strip, which is connected to four leads, as sketched in Fig. 8(a). The whole system consists of a single spatially constant material, which is discretized on a square grid with a lattice spacing of 4 nm. We assume the material hole doped and set the Fermi energy to \( E_F = -40 \text{ meV} \), such that no quantum Hall edge states are involved in transport and the sample behaves like a conventional metal. Furthermore, Anderson disorder is added on each lattice site. The strength of this disorder is set to \( U_0 = 30 \text{ meV} \) and the length of the strip to \( L = 8 \mu \text{m} \), which leads to a strong localization and a very low transmission between the right and the left side. The signatures of the strong localization are visible in the non-equilibrium local density of states (LDOS) in Fig. 8(b), which is calculated for a potential gradient using higher chemical potentials for the two leads at the left side. The states entering the system at the left are not able to propagate through the whole sample, which leads to a gradient in the LDOS from high densities (dark colors) at the left side to low densities (light colors) at the right side.

![FIG. 8. (Color online) a) Sketch of the setup showing the perfectly conducting channel in transport through a strongly disordered HgTe “H-bar” used in Fig. 4 of the main manuscript (\( L = 8 \mu \text{m}, W_M = 350 \text{ nm}, W_T = 150 \text{ nm}, W_L = 800 \text{ nm} \)). b) Typical carrier distribution (local density of states) arising for a small bias with additional gating (\( E_F = -40 \text{ meV} \), dark colors represent higher densities). c) Typical carrier distribution for a gate potential of \( V_g = 40 \text{ mV} \). d) Corresponding spin polarization.]

![FIG. 8. (Color online) a) Sketch of the setup showing the perfectly conducting channel in transport through a strongly disordered HgTe “H-bar” used in Fig. 4 of the main manuscript (\( L = 8 \mu \text{m}, W_M = 350 \text{ nm}, W_T = 150 \text{ nm}, W_L = 800 \text{ nm} \)). b) Typical carrier distribution (local density of states) arising for a small bias with additional gating (\( E_F = -40 \text{ meV} \), dark colors represent higher densities). c) Typical carrier distribution for a gate potential of \( V_g = 40 \text{ mV} \). d) Corresponding spin polarization.]

TABLE I. Material parameters for HgTe/CdTe quantum wells (units in meV and nm), taken from Ref. [24].

<table>
<thead>
<tr>
<th>A</th>
<th>B</th>
<th>D</th>
<th>M</th>
<th>Δ</th>
</tr>
</thead>
<tbody>
<tr>
<td>354.5</td>
<td>−686</td>
<td>−512</td>
<td>−10</td>
<td>1.6</td>
</tr>
</tbody>
</table>
FIG. 9. (Color online) Tunnel setup for the detection of the perfectly conducting channel in the diffusive regime using an “H-bar” ($L = 1.2 \mu m$, $W_M = 350 \, nm$, $W_{TI} = 250 \, nm$, $W_L = 400 \, nm$, $W_T = 120 \, nm$) made of a HgTe/CdTe heterostructure. The same setup is used for the results presented in Fig. 5 of the main manuscript. Typical carrier distribution arising for a small bias with higher chemical potentials applied to the left leads ($E_F = -40 \, meV$, dark colors represent higher densities). Lower part is tuned into the topological bandgap by $V_g = 40 \, mV$ and the upper strip is divided into two parts by an additional tunnel barrier with $V_t = 40 \, mV$.

In the following, a quantum spin Hall effect is induced by gating a part of the material into the TI bandgap at the lower edge of the structure, as depicted by the blue region in Fig. 8(a). This results in a quantized conductance between the lower contacts, carried by a quantum spin Hall edge state, which shows up as the thin dark connection between the lower leads in the LDOS presented in Fig. 8(c). At the same time a PCC arises between the two upper leads and gives rise to quantized transport, as shown by the data in Fig. 4 of the manuscript. In comparison to the quantum spin Hall state this state is completely spread out over the whole upper part. Furthermore, the spin polarization of the PCC differs strongly from the spin polarization of the edge state. In the quantum spin Hall regime the edge state shows a very strong spin polarization, visible by the constant blue color in the spin density at the lower edge in Fig. 8(d). On the contrary the PCC has no polarization, which can be seen by the patchy pattern in the upper part of the spin density.

In the previous setup a quantized conductance was used to demonstrate the emergence of a PCC in an otherwise localized region. Therefore a very long system was needed, such that all channels, except the PCC, are sufficiently localized by the impurity potential. Although the PCC is still present if $L$ is smaller, its transport signature is masked by the additional diffusive non-localized modes. In this case an additional tunnel barrier can be used to suppress the influence of the non-perfectly conducting channels as shown in the manuscript. Therefore, we investigate the transport properties in presence of such an additional tunnel barrier, depicted by the green bar in the sketch of Fig. 9. Similar to the previous setup, a gate $V_g$ is used to induce a quantum spin Hall state between the lower leads and a PCC in the upper part, which can be seen in the LDOS in Fig. 9. The tunnel barrier $V_t$ disconnects the upper two leads and reduces the contributions of the non-perfectly conducting channels. As shown in Fig. 5 of the main manuscript, this leads to a vanishing shot noise and a quantized conductance.