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OPEN Electric control of topological phase transitions in Dirac semimetal thin films

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Dirac semimetals host three-dimensional (3D) Dirac fermion states in the bulk of crystalline solids, which can be viewed as 3D analogs of graphene. Owing to their relativistic spectrum and unique topological character, these materials hold great promise for fundamental-physics exploration and practical applications. Particularly, they are expected to be ideal parent compounds for engineering various other topological states of matter. In this report, we investigate the possibility to induce and control the topological quantum spin Hall phase in a Dirac semimetal thin film by using a vertical electric field. We show that through the interplay between the quantum confinement effect and the field-induced coupling between sub-bands, the sub-band gap can be tuned and inverted. During this process, the system undergoes a topological phase transition between a trivial band insulator and a quantum spin Hall insulator. Consequently, one can switch the topological edge channels on and off by purely electrical means, making the system a promising platform for constructing topological field effect transistors.

The study of topological insulators (TIs) have been one of the most active research areas in the past ten years^{1,2}, which revolutionized our understanding of the electronic band structure. It is now understood that there could be nontrivial topologies encoded in the electronic wave-functions, characterized by various topological invariants according to the symmetry class of the system, and physically manifested by the appearance of topological boundary states. For example, two-dimensional (2D) TIs, also known as the quantum spin Hall (QSH) insulators, are characterized by a Z₂ topological invariant and have spin helical edge states on sample boundaries3, for which back-scattering is suppressed in the presence of time reversal symmetry^{1,2,4}. Hence they hold great promise for applications such as low-dissipation electronics, spintronics, and quantum computations. For these TIs, the nontrivial topology as well as the boundary states are protected by the finite energy gap, i.e., they are robust against perturbations as long as the insulating gap does not close.

It was later realized that the topological classification could be pushed beyond insulators to states without a gap⁵⁻⁷. In particular, a novel state called Dirac semimetal (DSM) has been proposed and successfully demonstrated in recent experiments for two crystalline materials Na₃Bi and $\hat{C}d_3\hat{A}s_2^{8-26}$. In these materials, the Fermi energy sits at two three-dimensional (3D) Dirac points—where the bands touch with a fourfold degeneracy—and the dispersion is linear along all three directions in reciprocal space. Each Dirac point can be viewed as consisting of two Weyl points of opposite chiralities and is protected by the crystalline symmetry^{8–10,27,28}. Such unusual electronic structure endows the system with many intriguing properties like the surface Fermi arcs and the quantum magnetoresistance 18,21,29,30. Perhaps more importantly, DSMs are expected to be an ideal parent compound for realizing other novel topological states such as Weyl semimetals, TIs, and topological superconductors¹¹. Particularly in this regard, DSMs offer a simple alternative to achieve the 2D TI phase through the quantum confinement effect 10,31.

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It has been shown that with increasing thickness, DSM thin films exhibit oscillations in the $2D \mathbb{Z}_2$ invariant whenever a quantum well state crosses the Dirac point 10,31 . Hence a QSH phase can be realized by a proper control of the film thickness. Since the QSH phase has so far been detected in only a few systems, given its fundamental and technological importance, the new approach to realize it using DSMs would be of great interest. Furthermore, the unique properties of DSMs may offer new methods to manipulate the properties of the QSH phase.

Motivated by these recent breakthroughs and by the great interest in utilizing DSMs for topological devices, in this work, we investigate the possibility of electric control of the topological phase transitions in a DSM thin film. We show that by using a vertical electric field, DSM thin films can be switched between a topological QSH phase and a trivial insulator phase. This topological phase transition is enabled by a combined effect of quantum confinement and field induced sub-band coupling. As a result, one can electrically manipulate the topological edge channels, and the charge and spin conduction through a finite sample can be readily switched on and off. This leads to a simple design of a DSM-based topological field effect transistor with advantages of fast-speed, low power consumption, and low dissipation, owing to the robust topological edge channels combined with full electric control.

Results

Model and Analytic Analysis. Our analysis is based on a generic low-energy effective model describing the DSMs A_3 Bi (A = Na, K, Rb) and Cd_3As_2 as derived in previous works^{9,10}. In these materials, the states around Fermi energy can be expanded using a minimal four-orbital basis of $S_{1/2}$, $S_{1/2}$, $S_{1/2}$, and $S_{1/2}$, $S_{1/2}$, $S_{1/2}$, $S_{1/2}$, and $S_{1/2}$, $S_{1/2}$, $S_{1/2}$, $S_{1/2}$, and $S_{1/2}$, $S_{1/2}$, and $S_{1/2}$, and $S_{1/2}$, $S_{1/2}$, and $S_{1/2}$, and $S_{1/2}$, and $S_{1/2}$, $S_{1/2}$, and $S_{1/2}$, and S

$$\mathcal{H}(\mathbf{k}) = \varepsilon_0(\mathbf{k}) + \begin{bmatrix} M(\mathbf{k}) & Ak_+ & 0 & 0\\ Ak_- & -M(\mathbf{k}) & 0 & 0\\ 0 & 0 & M(\mathbf{k}) & -Ak_-\\ 0 & 0 & -Ak_+ & -M(\mathbf{k}) \end{bmatrix},$$
(1)

where $\varepsilon_0(\mathbf{k}) = C_0 + C_1 k_z^2 + C_2 \left(k_x^2 + k_y^2\right)$, $k_\pm = k_x \pm i k_y$, and $M(\mathbf{k}) = -M_0 + M_1 k_z^2 + M_2 \left(k_x^2 + k_y^2\right)$ with $M_0, M_1, M_2 > 0$ to reproduce the band inversion feature at Γ -point. The material-specific parameters A, C_b and M_i are determined by fitting the first-principles result or the experimental measurement. It has been shown that this model nicely captures the essential low-energy physics as compared with experiment N_1, N_2, N_3 .

For bulk DSMs, model (1) gives the energy dispersion $\mathcal{E}(\pmb{k}) = \varepsilon_0(\pmb{k}) \pm \sqrt{M(\pmb{k})^2 + Ak_\parallel^2}$, where $\pmb{k}_\parallel = \begin{pmatrix} k_x, k_y \end{pmatrix}$ is the 2D wave-vector in the k_x - k_y plane. The spectrum has two Dirac points located along the k_z -axis at $(0, 0, \pm k_{\rm D})$ with $k_{\rm D} = \sqrt{M_0/M_1}$. Each Dirac point is four-fold degenerate and can be viewed as consisting of two Weyl nodes with opposite chiralities (as represented by the two 2×2 diagonal sub-blocks in Hamiltonian (1)). The dispersion around each Dirac point is linear in all three directions, as can be seen by expanding $\mathcal{E}(\pmb{k})$ at $(0, 0, \tau k_{\rm D})$ ($\tau = \pm$ labels the two Dirac points): $\mathcal{E}(\pmb{k}) \simeq \sqrt{A^2k_x^2 + A^2k_y^2 + 4M_1^2k_{\rm D}^2(k_z - \tau k_{\rm D})^2}$. One notes that the low-energy spectrum is anisotropic as manifested in both the distribution of Dirac points as well as the different Fermi velocities along k_z versus that in the k_x - k_y plane (Fermi velocity along k_z is typically much slower), which leads to quite different behaviors when a DSM is confined along different directions³¹.

DSMs such as Na₃Bi and Cd₃As₂ have layered structures along crystal *c*-axis. Hence their thin film structures with confinement along *z*-direction can be more readily fabricated. Consider a DSM thin film with thickness *L* confined in the region $z \in [-L/2, L/2]$. For small *L*, the electron motion along *z* will be quantized into discrete quantum well levels due to quantum confinement effect. This generally turns the system from a semimetal to a semiconductor. Using quantum well approximation, each quantum well level has a quantized effective wave-vector k_z such that $\langle k_z \rangle_n = 0$ and $\langle k_z^2 \rangle_n \simeq (n\pi/L)^2$ with $n = 1, 2, \cdots$ counting the sub-bands and the angular bracket meaning the average over quantum well states.

One observes that for each sub-band n, Hamiltonian (1) has a similar form as the low-energy model describing the 2D QSH systems in HgTe/CdTe quantum wells³², with the term

$$M(\mathbf{k}) \to M(n, \mathbf{k}_{\parallel}) = \mathcal{M}_n + M_2(k_x^2 + k_y^2),$$
 (2)

where $\mathcal{M}_n \equiv -M_0 + M_1(n\pi/L)^2$ is a sub-band dependent mass which determines the gap of the sub-band at Γ -point of the 2D Brillouin zone. It's known that band inversion occurs (around $\mathbf{k}_{\parallel} = 0$) when $\mathrm{sgn}(\mathcal{M}_n/M_2) = -1^{32,33}$, i.e. when \mathcal{M}_n and M_2 have opposite signs, which signals a nontrivial $\mathbb{Z}_2 = 1$ character of the sub-band n. Given that M_0 , M_1 , $M_2 > 0$, this happens when $\tilde{k}_n \equiv \langle k_z^2 \rangle_n^{1/2} (=n\pi/L) < k_\mathrm{D}$ is satisfied. Therefore, for a thin-enough film such that $\tilde{k}_1 = \pi/L > k_\mathrm{D}$, all the sub-bands are topologically trivial with positive mass terms \mathcal{M}_n . With increasing film thickness, the

system becomes nontrivial once the first (n=1) sub-band has its mass \mathcal{M}_1 inverted when $\tilde{k}_1 < k_D$. The inverted sub-band contributes a $\mathbb{Z}_2 = 1$ and in the inverted band gap, there appears a pair of spin-helical edge states protected by time reversal symmetry on each edge of the quasi-2D system. Further increasing L would invert \mathcal{M}_2 of the second sub-band, leading to two pairs of edge states. However, for \mathbb{Z}_2 group: 1+1=0, hence this state is topologically trivial. Physically, it is because backscattering can occur between edge states from different time reversal pairs. Following this logic, the topological properties as well as the bulk band gap show oscillatory behavior as a function of the film thickness 10,31 .

Since the sub-band dependent mass \mathcal{M}_n plays the key role in determining the topological properties of the system, we shall focus on the change of \mathcal{M}_n by a vertical electric field, aiming to achieve an electric control of the topological phase of DSM thin films. To proceed, one notes that the lower diagonal block of Hamiltonian (1) is formally the time reversal counterpart of the upper block, which share the same energy spectrum and the E field does not mix the two. (This can also be argued by observing that the low-energy Hamiltonian possesses the unitary symmetry $\mathcal{U} = I \otimes \sigma_z$ and the anti-unitary symmetry $\mathcal{U}' = -i\sigma_y \otimes IK$, which may be regarded as emergent symmetries for the low-energy physics. Here σ 's are the Pauli matrices, I is the 2×2 identity matrix, and K is the complex conjugation operator.) Hence to study the change of \mathcal{M}_n , it is enough to consider only the upper block denoted by $h(\mathbf{k})$. Modeling with the hard-wall boundary condition for the confinement potential, we have for sub-band n,

$$h_n(\mathbf{k}_{\parallel}) = \varepsilon_0(n, \mathbf{k}_{\parallel})I + Ak_x\sigma_x - Ak_y\sigma_y + M(n, \mathbf{k}_{\parallel})\sigma_z, \tag{3}$$

where σ 's are the Pauli matrices, I is the 2 × 2 identity matrix, and $\varepsilon_0(n, \mathbf{k}_{\parallel}) = C_0 + C_1(n\pi/L)^2 + C_2k_{\parallel}^2$. The energy eigenstates are given by

$$\left\langle \mathbf{r} \middle| \Psi_{n\alpha} \Big(\mathbf{k}_{\parallel} \Big) \right\rangle = \frac{1}{\sqrt{S}} e^{ik_{x}x + ik_{y}y} \psi_{n}(z) \chi_{n\alpha}, \tag{4}$$

with eigen-energies

$$\mathcal{E}_{n\alpha}(\mathbf{k}_{\parallel}) = \varepsilon_0(n, \mathbf{k}_{\parallel}) + \alpha \sqrt{A^2 k_{\parallel}^2 + M(n, \mathbf{k}_{\parallel})^2}, \tag{5}$$

where S is the area of the thin film, $\alpha = \pm$, $\chi_{n\alpha}$ are the two eigen-spinors along the quantization direction $\left(Ak_x, -Ak_y, M\left(n, \mathbf{k}_{\parallel}\right)\right)$, and

$$\psi_n(z) = \sqrt{\frac{2}{L}} \sin\left[\frac{n\pi}{L}\left(z + \frac{L}{2}\right)\right] \tag{6}$$

is the quantum well state for the *n*-th sub-band. The vertical electric field is modeled by adding a diagonal potential energy term V(z) = eEz where (-e) is the electron charge and E is the effective field strength which may be considered as including the static screening effects.

For a qualitative analysis, we assume small field and treat V perturbatively. Because V(z) is odd in z, it is easy to see that the first order perturbation in energy vanishes. The leading order perturbation comes at the second order, with

$$\delta \mathcal{E}_{n\alpha}(\mathbf{k}_{\parallel}) \approx \sum_{m\beta \neq n\alpha} \frac{\left| \left\langle \Psi_{m\beta}(\mathbf{k}_{\parallel}) \middle| V \middle| \Psi_{n\alpha}(\mathbf{k}_{\parallel}) \right\rangle \right|^{2}}{\mathcal{E}_{n\alpha}(\mathbf{k}_{\parallel}) - \mathcal{E}_{m\beta}(\mathbf{k}_{\parallel})}, \tag{7}$$

where the summation is over all other states different from $\Psi_{n\alpha}$, and in reality, it has a physical cutoff for which the low-energy description is no longer valid. One notes that in order to analyze the renormalized \mathcal{M}_n , it is sufficient to focus on the change at Γ -point of the 2D Brillouin zone by setting $\mathbf{k}_{\parallel} = 0$.

We are most interested in the case in which the mass (gap) of the first sub-band can be inverted by the electric field, because then the two sides of the topological phase transition have the most salient contrast: absence or presence of the topological edge channels, hence leading to the best on-off ratio when considering a topological transistor based on it. For such case, we consider a thickness L such that $\mathcal{M}_1 = -M_0 + M_1(\pi/L)^2 > 0$, i.e. an initially trivial system with $\mathbb{Z}_2 = 0$ in the absence of E field. At $k_\parallel = 0$, we have $\chi_{n+} = |\uparrow\rangle$ and $\chi_{n-} = |\downarrow\rangle$ for all n, where $|\uparrow\rangle$ and $|\downarrow\rangle$ are the two eigenstates of σ_z . For the n=1 sub-band, we have at $k_\parallel = 0$,

$$\delta \mathcal{E}_{1+} \approx -e^2 E^2 \sum_{m>1} \frac{\left| \left\langle \psi_m(z) | z | \psi_1(z) \right\rangle \right|^2}{\mathcal{E}_{m+} - \mathcal{E}_{1+}} = -e^2 E^2 \frac{64L^4}{\pi^6} \frac{1}{(M_1 + C_1)} \eta. \tag{8}$$

where

$$\eta = \sum_{m \in \text{even}} \frac{m^2}{(m^2 - 1)^5}.$$
(9)

In the first equality of (8) we used the fact that the state $|\Psi_{1+}(0)\rangle$ does not mix with the $|\Psi_{m-}(0)\rangle$ states from the valence bands by the E field because their pesudo-spin parts χ are orthogonal. Also note that one has $M_1>|C_1|$ in order for the model (1) to describe a semimetal phase, hence $\mathcal{E}_{m+}>\mathcal{E}_{1+}$ for m>1, hence the perturbation due to the coupling between $|\Psi_{1+}(0)\rangle$ and $|\Psi_{m+}(0)\rangle$ (with m>1) generally pushes down the energy level of $|\Psi_{1+}(0)\rangle$, making $\delta\mathcal{E}_{1+}<0$. In the expression (9) of the constant factor η (with a rapidly converging value \simeq 0.0165), the summation only includes the even integer numbers, because V(z) only couples states with opposite parities in z.

Similarly, the energy shift for state $|\Psi_{1}(0)\rangle$ can be calculated,

$$\delta \mathcal{E}_{1-} \approx e^2 E^2 \frac{64L^4}{\pi^6} \frac{1}{(M_1 - C_1)} \eta,$$
 (10)

which is positive, showing that the coupling induced by the E field pushes up the energy of $|\Psi_1(0)\rangle$. Therefore, in the Hilbert sub-space of the first sub-band, the E field renormalizes the value of the mass:

$$\mathcal{M}_1 \to \mathcal{M}_1 + \delta \mathcal{M}_1,$$
 (11)

with the correction

$$\delta \mathcal{M}_1(E) = \frac{\delta \mathcal{E}_{1+} - \delta \mathcal{E}_{1-}}{2} \approx -e^2 E^2 \frac{64L^4}{\pi^6} \frac{2M_1}{M_1^2 - C_1^2} \eta. \tag{12}$$

Using this estimation, one observes that the gap of the first sub-band would decrease with increasing E field, and closes at

$$E_c \approx \frac{\pi^3}{8e^2L^2} \sqrt{\frac{M_1^2 - C_1^2}{2\eta M_1}} \mathcal{M}_1,$$
 (13)

which signals a topological phase transition point and beyond which the gap reopens with the system turned into a QSH insulator phase characterized by $\mathbb{Z}_2 = 1$.

For large E field with eEL being comparable or even larger than \mathcal{M}_1 , the perturbative calculation is no longer expected to be accurate. Nevertheless, the general physical pictures from the above discussion still applies: the level repulsion due to higher sub-bands would generally decrease and invert the gap of the first sub-band, generating a topological phase transition. We shall explicitly demonstrate this in the next section through numerical calculations.

Before proceeding, we mention that our above analysis based on the low-energy effective model around Γ -point is valid, because for these materials: the band inversion only occurs around Γ -point; while the bands near the Brillouin zone boundary (away from Γ -point) have normal band ordering and are higher in energy, to invert them would require very large applied field if not impossible. Hence for the consideration of topological properties, we can focus on the band evoluation around Γ -point. In addition, in our treatment of confinement we are focusing on the thickness range where the relevant sub-band gap is small and close to band inversion, which means the corresponding effective wave-vector \tilde{k}_n is close to k_D , hence its low-energy behavior and especially the topological property are well-captured by this effective model in Eq.(1). Similar treatment based on the low-energy effective models has been successfully applied in the study of quantum confined structures, e.g. in semiconductor quantum wells³⁴, 3D topological insulator thin films³⁵, and also in previous studies of Dirac semimetal thin films^{10,31}. The above analysis can also be applied to higher sub-bands when a thicker film with the n-th (n > 1) sub-band most close to transition is considered. We will discuss this later in the discussion section.

Numerical Results. For numerical investigation, we discretize the model (1) on a 3D lattice with lattice constants $a_x = a_y = 0.5448 \,\mathrm{nm}$ (1.264 nm), and with $a_z = 0.4828 \,\mathrm{nm}$ (2.543 nm) being set to the interlayer separation pertinent to Na₃Bi (Cd₃As₂). The standard substitutions

$$k_i \to \frac{1}{a_i} \sin(k_i a_i), \quad k_i^2 \to \frac{2}{a_i^2} [1 - \cos(k_i a_i)]$$
 (14)

are adopted (i=x,y,z) for lattice discretization around Γ -point. Since we require the initial state at E=0 is of a trivial insulator phase, we need the number of layers $\ell < \lfloor \pi M_1/(M_0a_z) \rfloor + 1$, where $\lfloor \cdots \rfloor$ is the floor function. And in order for the band gap to be inverted by a relatively small E field, one may wish to have the initial gap size $2M_1$ not too large.

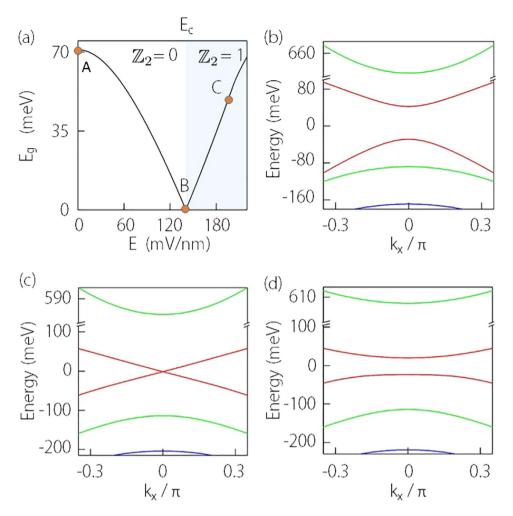


Figure 1. Field induced topological phase transition in Na₃Bi thin film. (a) Variation of energy gap as a function of the vertical field E showing the gap closing and reopening process, marking a topological phase transition between $\mathbb{Z}_2=0$ and $\mathbb{Z}_2=1$ phases. (b–d) Energy spectra corresponding to A, B, and C as marked in (a) plotted versus k_x (with $k_y=0$). The first, second, and third sub-bands are marked using red, green, and blue colors respectively. The parameters for used in the calculation are $\ell=5$, $C_0=-63.82\,\mathrm{meV}$, $C_1=87.538\,\mathrm{meV}$ nm², $C_2=-84.008\,\mathrm{meV}$ nm², $C_2=-84.008\,\mathrm{meV}$ nm², $C_2=-84.008\,\mathrm{meV}$ nm², $C_2=-84.008\,\mathrm{meV}$ nm², and $C_2=-84.008\,\mathrm{meV}$ nm², arigid energy shift is applied to make the gap center at zero energy.

Let's consider Na₃Bi first. The model parameters we use are listed in the caption of Fig. 1, which have been extracted from the first-principles calculations and compared well with experiment. For Na₃Bi thin films, the critical thickness for which the first sub-band undergoes band inversion is around $\ell_c \simeq 6^{31}$. Hence we take a film with $\ell=5$ layers ($L\simeq 2.414\,\mathrm{nm}$) for demonstration. Figure 1(a) shows the variation of the band gap $E_{\rm g}$ as a function of the E field. The result is symmetric between positive and negative values of E, so only the positive part is shown here. Initially, at E=0, the system has a confinement gap about 71 meV (marked by point A). With increasing E field, the gap decreases and closes at a critical value $E_c\simeq 140\,\mathrm{mV/nm}$ (marked by point B), and then reopens and increases with E. This is consistent with our previous analytic analysis. The value of E_c is also not far from our estimation in Eq.(13) which is about 208 mV/nm. We also plot in Fig. 1(b–d) the energy spectrum of the system corresponding to the three representative states marked by A, B, and C in Fig. 1(a). It shows that on both sides of the gap closing point, the system is insulating with the gap belong to the first sub-band. At the critical value E_c , the band gap closes at Γ -point, marking the topological phase boundary which separates the topologically trivial and nontrivial phases.

To further demonstrate the topological nature of the transition and to visualize the edge states, we compute the surface local density of states (LDOS) for the side surface. Due to the isotropy in the k_x - k_y plane of the low-energy model (1), without loss of generality, we choose the surface perpendicular to y-direction of the quasi-2D system. The surface LDOS $\rho(k_x)$ can be calculated for each k_x from the

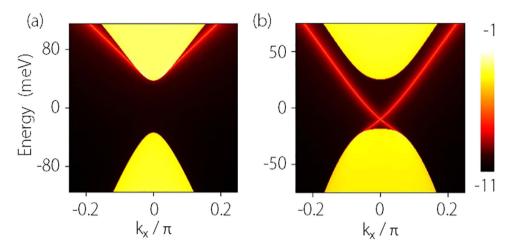


Figure 2. Calculated LDOS of a side surface for Na_3Bi thin film. (a) is for point A (trivial insulator) and (b) is for point C (QSH insulator) as marked in Fig. 1(a). The calculation is for a slab which is semi-infinite along y-direction and the parameters are the same as for Fig. 1.

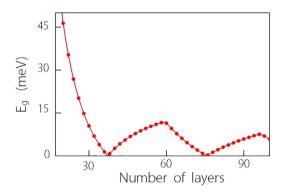


Figure 3. Confinement induced gap size versus thickness ℓ for Cd₃As₂ thin films. The parameters used in the calculation are $C_0 = -219 \,\text{meV}$, $C_1 = -300 \,\text{meV}$ nm², $C_2 = -160 \,\text{meV}$ nm², $M_0 = 10 \,\text{meV}$, $M_1 = 9600 \,\text{meV}$ nm², $M_2 = 180 \,\text{meV}$ nm², and $A = 275 \,\text{meV}$ nm.

surface Green's function $\rho\left(k_x\right)=-\mathrm{Tr}\left[\mathrm{Im}\;G_{00}(k_x)\;\right]/\pi$, where G_{00} is the retarded Green's function for the surface layer (labled by index 0) of the lattice³⁶. G_{00} can be evaluated by the transfer matrix through a standard numerical iterative method³⁷. The obtained surface LDOS for states before and after the phase transition (for state A and C) are plotted in Fig. 2. One observes that for both cases, the confinement-induced bulk gap can be clearly identified. Before the topological phase transition ($E < E_c$), there is no states inside the gap. In contrast, after transition ($E > E_c$), there appear two bright lines crossing the gap, corresponding to the spin helical edge states for the \mathbb{Z}_2 nontrivial QSH phase. As long as time reversal symmetry is preserved, these gapless modes are protected and carriers in these channels cannot be backscattered^{1,2}. Therefore transport through these channels is in principle dissipationless. In a two-terminal measurement, this would lead to a quantized conductance, which has been confirmed experimentally³.

Similar analysis applies to Cd_3As_2 as well. In Fig. 3, we plot the variation of its confinement-induced gap versus the film thickness, which clearly shows the oscillation behavior of the gap³¹. One observes that the critical thickness ℓ_c is at about 37 layers. Here we choose a film thickness of $\ell=20$ layers ($L=50.86\,\mathrm{nm}$) for demonstration. The variations of the gap with respect to the E field as well as representative energy spectra are shown in Fig. 4. Again the gap closing and reopening process similar to Fig. 1(a) is observed. The critical value of $E_c \simeq 5.26\,\mathrm{mV/nm}$ also agrees well with the estimation $\approx 4.17\,\mathrm{mV/nm}$ from Eq.(13). The energy spectra also coincide with our previous analysis. Figure 5 shows the side surface LDOS plots for states A and C (marked in Fig. 4(a), clearly showing the appearance of topological edge states across the transition. These results show qualitatively the same features as those for Na₃Bi.

Our numerical results discussed above thus confirm our analytical analysis. A vertical electric field can be used to control the topological phase transitions and the topological edge channels in a DSM thin film.

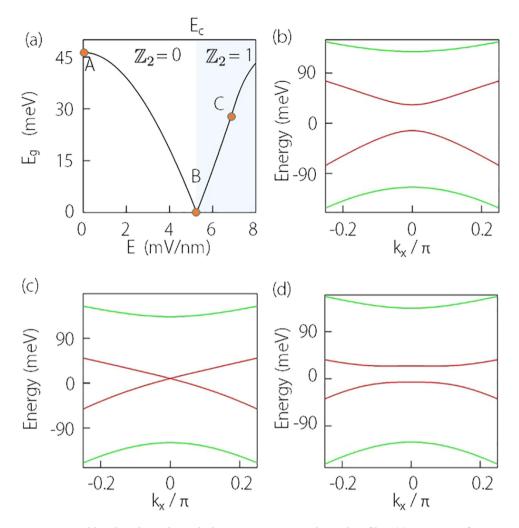


Figure 4. Field induced topological phase transition in Cd₃**As**₂ **thin film.** (a) Variation of energy gap as a function of the vertical field E. (**b-d**) Energy spectra corresponding to A, B, and C as marked in (a) plotted versus k_x (with $k_y = 0$). The first and the second sub-bands are marked using red and green colors respectively. The parameters for used in the calculation are the same as for Fig. 3 and $\ell = 20$ is taken.

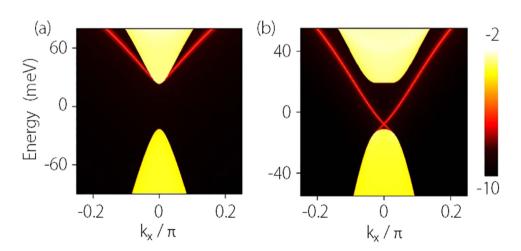


Figure 5. Calculated LDOS of a side surface for Cd_3As_2 thin film. (a) is for point A (trivial insulator) and (b) is for point C (QSH insulator) as marked in Fig. 4(a). The calculation is for a slab which is semi-infinite along *y*-direction and the parameters are the same as for Fig. 4.

Discussion

This work theoretically demonstrates the possibility to electrically control the $\mathbb{Z}_2=0/1$ topological phase transitions in a DSM thin film. Since the bulk topology is tied to the existence of topological edge channels and hence to the charge/spin conductance, this indicates that one can achieve a full electric control of the on/off charge/spin conductance of such a system, making it a suitable candidate for a topological field effect transistor. The electric field can be generated by the standard top and bottom gates setup. Compared with the traditional MOSFET which works by injection and depletion of charge carriers in the channel region and has a response timescale depending on factors such as the charge concentration and the carrier mobility, the operating mechanism for a topological transistor is expected to have a high on/off speed with electronic response timescale and a better power efficiency³⁸. In addition, multiple conducting channels in a transistor can be achieved by designing a multilayer structure with alternating DSM layers and insulating layers, similar to the structure as in Ref. 38.

For device design, we have seen that a proper film thickness can be chosen such that the starting confinement gap is small hence can be easily inverted by a small applied field. However, there is a tradeoff because if the gap is too small, then the thermally populated carriers in the bulk could strongly contribute to the transport. Therefore, a balance needs to be achieved for the device to have an optimal performance with relatively low power consumption.

In our analysis, we have focused on the phase transition in the first sub-band. Similar analysis can also be extended to higher sub-bands if a thicker film with its n-th (n>1) sub-band close to phase transition is considered. For example, consider a film thickness such that its second sub-band is just before the gap-closing. In this case, we would have $\mathcal{M}_2 > 0$ and $\mathcal{M}_1 < 0$, and the system is in a $\mathbb{Z}_2 = 1$ phase. Perturbation to second order in the field strength gives the energy correction of

$$\delta \mathcal{E}_{2+} \approx -e^2 E^2 \frac{256L^4}{\pi^6} \left[\frac{1}{M_1 + C_1} \eta' - \frac{1}{81[3C_1 + 5M_1 - 2M_0(L/\pi)^2]} \right], \tag{15}$$

for the $|\Psi_{2+}(0)\rangle$ state of the second sub-band, with $\eta' = \sum_{m \in \text{odd}, m \geq 3} m^2/(m^2-4)^5$. A similar expression can be obtained for $\delta \mathcal{E}_{2-}$ as well. The first term in the parenthesis of (15) is from the coupling with the m>2 sub-bands, while the second term is from the coupling with the first sub-band. The sign of this energy shift (and hence the correction of \mathcal{M}_2) would depend on the competition between the two terms and is not necessarily negative. Nevertheless, for such higher sub-band case, even if a field-induced top-ological phase transition can be realized, the topologically trivial phase would in fact still possess edge states. Although these (even number of pairs of) channels are not topologically robust, their presence would make the trivial state not completely 'off' hence is detrimental to the performance of a transistor.

Finally, in real systems, there could be other perturbation terms, e.g. Rashba spin-orbit couplings from possible structrual inversion asymmetry due to field or substrate effects. Such terms could generate a trivial gap competing with the QSH gap. Since a topological phase is protected by the bulk gap, the QSH phase would survive and is robust as long as the corresponding QSH gap dominates over the trivial gap due to other mechanisms³⁹.

In summary, we have demonstrated that full electric control of topological phase transitions in a DSM thin films can be achieved through the interplay between the quantum confinement effect and the coupling between sub-bands induced by a vertical electric field. As a result, a topological field effect transistor can be constructed in which carriers are conducted through the topological edge channels. Given that several DSM materials have been experimentally demonstrated and that the progress in material fabrication technology such as molecular beam epitaxy has allowed film growth with atomic precision, it is quite promising for the physical effect and the DSM-based topological transistor proposed here to be realized in the near future.

Methods

Lattice model. To investigate the effect of a vertical electric field, we discretize a generic low-energy effective model in Eq.(1) on a 3D lattice with lattice constants a_x , a_y , and a_z along the three orthogonal directions. The low-energy effective Hamiltonian for a 3D DSM around the Γ -point in the Brillouin zone is given by

$$H = H_0 + H_x + H_y + H_z. (16)$$

Here,

$$H_{0} = \begin{pmatrix} M_{+} & 0 & 0 & 0 \\ 0 & M_{-} & 0 & 0 \\ 0 & 0 & M_{+} & 0 \\ 0 & 0 & 0 & M_{-} \end{pmatrix} \sum_{i} c_{i}^{\dagger} c_{i}, \tag{17}$$

$$H_{x} = \begin{pmatrix} -C_{2}^{a} + M_{2}^{a} & -\frac{A^{a}}{2}i & 0 & 0\\ -\frac{A^{a}}{2}i & -C_{2}^{a} - M_{2}^{a} & 0 & 0\\ 0 & 0 & -C_{2}^{a} + M_{2}^{a} & \frac{A^{a}}{2}i\\ 0 & 0 & \frac{A^{a}}{2}i & -C_{2}^{a} - M_{2}^{a} \end{pmatrix} \sum_{i} (c_{i}^{\dagger}c_{i+x}) + H. c.,$$
(18)

$$H_{y} = \begin{pmatrix} -C_{2}^{a} + M_{2}^{a} & \frac{A^{a}}{2} & 0 & 0 \\ -\frac{A^{a}}{2} & -C_{2}^{a} - M_{2}^{a} & 0 & 0 \\ 0 & 0 & -C_{2}^{a} + M_{2}^{a} & \frac{A^{a}}{2} \\ 0 & 0 & -\frac{A^{a}}{2} & -C_{2}^{a} - M_{2}^{a} \end{pmatrix} \sum_{i} (c_{i}^{\dagger} c_{i+y}) + H. c.,$$
(19)

and

$$H_{z} = \begin{pmatrix} -C_{1}^{a} + M_{1}^{a} & 0 & 0 & 0 \\ 0 & -C_{1}^{a} - M_{1}^{a} & 0 & 0 \\ 0 & 0 & -C_{1}^{a} + M_{1}^{a} & 0 \\ 0 & 0 & 0 & -C_{1}^{a} - M_{1}^{a} \end{pmatrix} \sum_{i} (c_{i}^{\dagger} c_{i+z}) + H. c.,$$

$$(20)$$

where $M_{\pm}=C_0\pm M_0+(2C_1^a+4C_2^a)\mp(2M_1^a+4M_2^a)$. Using this lattice Hamiltonian, the low energy spectrum and the wavefunctions for the DSM slab under a perpendicular electric field can be constructed.

Surface LDOS. The surface LDOS can be derived from $\rho(k) = -\frac{1}{\pi} \text{TrIm} G_{00}(k)$, where G_{00} is the retarded Green's function for the surface layer of a semi-infinite 3D lattice model. The surface Green's function can be obtained through the transfer matrix method, with

$$G_{00} = (E - H_{00} - H_{01}T)^{-1}$$
(21)

$$T = (E - H_{00} - H_{01}T)^{-1}H_{01}^{\dagger}$$
(22)

where H_{00} and H_{01} are Hamiltonian matrix elements for a single layer and for the interlayer coupling, and T is the transfer matrix. Generally, Eq.(22) can be solved by iterative calculations until T converges with the help of a fast iteration algorithm³⁷.

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Author Contributions

S.A.Y. conceived the idea. M.W. and H.P. performed the numerical calculation. Y.L. and S.A.Y. did the analytic calculation and analysis. All authors contribute to the data analysis and interpretation. S.A.Y. and H.P. supervised the project and wrote the manuscript. All authors reviewed the manuscript.

Additional Information

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