Strongly interacting Dirac liquid on the surface of a topological Kondo insulator

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A topological Kondo insulator (TKI) is a strongly correlated material, where hybridization between the conduction electrons and localized f electrons gives rise to a crossover from a metallic behavior at high temperatures to a topologically nontrivial insulating state at low temperatures. The existing description of the TKIs is based on a slave-boson mean-field theory, which neglects dynamic fluctuation phenomena. Here, we go beyond the mean-field theory and investigate the role of Kondo fluctuations on the topological surface states. We derive an effective theory of the Dirac surface states coupled to fluctuations and show that the latter mediate strong repulsive interactions between surface excitations. We show that these effects renormalize the plasmon spectrum on the surface. We also argue that Kondo-mediated interactions may drive a magnetic instability of the surface spectrum.

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Introduction. The last decade has witnessed a flurry of interest and tremendous advance in the studies of topological states of matter. By now, a new class of topological insulating state has been theoretically predicted [1–6] and experimentally discovered in both two- and three-dimensional material systems (see Refs. [7–9], and references therein). Moreover a complete topological classification of noninteracting band structures has been put together [10,11]. Despite this impressive progress, there still remain a number of open fundamental questions, and understanding the interplay between interactions and topological states is arguably the chief challenge among them on the theoretical side [12,13].

While some of these fundamental questions are academic at this stage, there is also an experimental motivation to look closer at the effect of interactions on the topological surface states, which comes from the recent discovery of a topological Kondo insulator in samarium hexaboride [14–25]. Indeed, these heavy-fermion topological insulators represent the first known class of material where the topological behavior arises not from the noninteracting band structure, but due to strong interactions and correlations [26,27]. Furthermore, a recent experiment by the Paglione group [28] on SmB₆ suggests that interactions in this topological Kondo insulator (TKI) are not a minor complication, but may be of the utmost importance, because they appear to drive the topological surface states into a ferromagnetic phase. The nature of the latter is not understood as of now and this calls for a critical revision of the underlying theoretical framework.

The theoretical description of a Kondo insulator is usually based on the Anderson model [see Eq. (1) below], which describes hybridization between the conduction d electrons and localized, strongly interacting f electrons. To solve the model, the slave-boson technique [29] is usually employed, which allows one to conveniently circumvent complications associated with strong (formally, infinite) repulsion between the f electrons by "splitting" the physical f electron into a product of a fermion and a slave boson, supplemented with a constraint to remove the double occupancy. A mean-field

analysis (reiterated in more detail below) involves condensing the boson field and neglecting all of its dynamics. This effectively leads to a noninteracting model of an insulator, where the gap is proportional to the condensate and hybridization parameter, and makes it amenable to a topological analysis. Since the electron states being hybridized have the opposite parities, the resulting Kondo insulator is topological [26,27] and contains its hallmark feature—the Dirac surface states

There is a good evidence that the mean-field theory is reasonable, and actually works at least at the qualitative level. In particular, the existing data on SmB_6 , including the observation of surface states there [15,16,19], appear to be in good agreement with it. However, there are also reasons, both experimental and theoretical, to study effects beyond mean field and this is the focus of our work.

There has been a large amount of prior work on the effect of Kondo fluctuations in heavy-fermion materials both metallic [30–37] and insulating [38]. The theory is well controlled in the large- $N_{\rm f}$ limit (where $N_{\rm f}$ is the degeneracy of the f level) and the mean-field approximation becomes exact for $N_{\rm f} \to \infty$. In realistic systems (with $N_{\rm f}=2$ or $N_{\rm f}=4$), the consensus appears to be that fluctuations play an important role in heavy-fermion metals, but are essentially irrelevant in (conventional) Kondo insulators. The case of a topological Kondo insulator considered here is of a mixed character, because the "matter fields," which the Kondo fluctuations are coupled to, are gapped in the bulk, but gapless on the surface.

To succinctly summarize the results of our theory detailed below, we find that Kondo fluctuations manifest themselves as strong short-range repulsive temperature-dependent interactions between surface Dirac states. Therefore, a proper description of the TKI's surface state is a *strongly interacting Dirac liquid*. The interaction can drive the magnetic instability on the surface, which is accompanied by opening of an insulating gap in the Dirac spectrum. In the absence of a phase transition, these interactions distort the spectrum of surface collective modes.

Mean-field slave-boson theory of a TKI. Our starting point is the following Anderson lattice Hamiltonian:

$$\hat{H} = \sum_{ij\sigma} \epsilon_{ij}^{d} \hat{d}_{i\sigma}^{\dagger} \hat{d}_{j\sigma} + \sum_{ij\alpha} \epsilon_{ij}^{f0} \hat{f}_{i\alpha}^{\dagger} \hat{f}_{j\alpha}$$

$$+ \sum_{ij\sigma\alpha} \left(V_{ij}^{\sigma\alpha} \hat{d}_{i\sigma}^{\dagger} \hat{f}_{j\alpha} + \text{H.c.} \right) + \frac{U}{2} \sum_{i\alpha} \hat{f}_{i\alpha}^{\dagger} \hat{f}_{i\alpha} \hat{f}_{i\bar{\alpha}}^{\dagger} \hat{f}_{i\bar{\alpha}}, \quad (1)$$

where the first two terms describe hopping of d and felectrons, with $\epsilon_{\bf k}^{\rm d}$ and $\epsilon_{\bf k}^{\rm f0}$ being their bare dispersion laws as a function of the lattice momentum, k. As our goal is to extract new physics associated with the fluctuation effects in TKIs, we would like to separate them from nonuniversal phenomena associated with a particular compound. Hence, we will focus on a generic topological Kondo insulator with the simplest band structure and assume that the relevant f states form a Kramers doublet (labeled by the index $\alpha = \uparrow, \downarrow$ and $\bar{\alpha} = \downarrow, \uparrow$), $\epsilon_{\bf k}^{\rm d}=-E_{\rm d}+k^2/2m_{\rm d}$ and $\epsilon_{\bf k}^{\rm f0}=-E_{\rm f}^0-k^2/2m_{\rm f}$. The model requires an ultraviolet cutoff, $E_{\rm d}^{\rm m}$, which corresponds to the width of the conduction band. All energies are measured from the position of the unoccupied f levels. As shown below, the resulting theory is described by 4×4 matrices in the bulk and one Dirac cone on the surface (as opposed to an 8×8 bulk Hamiltonian and three surface Dirac cones for SmB₆ [39–41]). The third term in Eq. (1) describes hybridization between the fand d states. Since they have different parities, which dictates $\hat{V}_{-\mathbf{k}} = -\hat{V}_{\mathbf{k}}$ for their Fourier transforms, and the time-reversal symmetry is respected, the hybridization matrix element at small momenta can be approximated as $\check{V}_{\mathbf{k}} = V\mathbf{k} \cdot \check{\boldsymbol{\sigma}}$. This particular form of the hybridization is the source of nontrivial topology of the emergent bands. The last term in Eq. (1) is the Hubbard repulsion between two f electrons on a level. The Hubbard energy scale, U, much exceeds all other energy scales in the problem and hence the last term effectively enforces a no-double-occupancy constraint on each site. This constraint is ultimately a source of fluctuation effects we are studying.

To exclude the double occupation of the f states we use the slave-boson approach, with spinon \hat{f}_i and holon \hat{b}_i operators introduced in the following way:

$$\hat{f}_{i\alpha} \to \hat{f}_{i\alpha} \hat{b}_i^{\dagger}, \qquad \hat{f}_{i\alpha}^{\dagger} \to \hat{f}_{i\alpha}^{\dagger} \hat{b}_i.$$
 (2)

Singly occupied f states correspond to the occupation numbers $|1_{\uparrow s},0_{\downarrow s},0_{h}\rangle$ ($|0_{\uparrow s},1_{\downarrow s},0_{h}\rangle$) in terms of the spinons and holons, and an empty f-level corresponds to a state occupied by a holon $|0_{\uparrow s},0_{\downarrow s},1_{h}\rangle$. The doubly occupied f levels are excluded by the constraint on each site $\hat{Q}_{i}=0$ (operator identity), where \hat{Q}_{i} is given by

$$\hat{Q}_i = \sum_{\alpha} \hat{f}_{i\alpha}^{\dagger} \hat{f}_{i\alpha} + \hat{b}_i^{\dagger} \hat{b}_i - 1.$$
 (3)

The reason behind this construction is that an equality is straightforward to enforce in the Lagrangian formalism by introducing a functional delta function (while the original inequality in terms of the physical f electrons is not as easy to implement). The Lagrangian for our system takes the

form

$$L = \sum_{i} \bar{b}_{i}(\tau) \partial_{\tau} b_{i}(\tau) + \sum_{i} i \lambda_{i}(\tau) Q_{i}(\tau)$$

$$+ \sum_{ij} \bar{\Psi}_{i}(\tau) \begin{pmatrix} \partial_{\tau} + \epsilon^{d}_{ij} & \check{V}_{ij} \bar{b}_{j}(\tau) \\ b_{i}(\tau) \check{V}^{\dagger}_{ij} & \partial_{\tau} + b_{i}(\tau) \epsilon^{f0}_{ij} \bar{b}_{j}(\tau) \end{pmatrix} \Psi_{j}(\tau),$$

$$(4)$$

where $\Psi_i = (d_{i\uparrow}, d_{i\downarrow}, f_{i\uparrow}, f_{i\downarrow})^{\rm T}$ is a column of Grassmann fields combining the d and f electrons, and $\lambda_i(\tau)$ is a time-dependent auxiliary real field, which enforces the no-double-occupancy constraint. \check{V} is a 2×2 hybridization matrix in the spin/pseudospin space and the diagonal components of the matrix in (4) are proportional to a 2×2 identity matrix (not shown).

The phase degree of freedom of slave particles is a redundant variable and can be integrated out after a gauge transformation [35–37], which separates the phase θ_i and amplitude of holon field b_i^a ,

$$b_i = b_i^{a} e^{i\theta_i}, \quad f_i \to f_i e^{i\theta_i}, \quad \lambda_i \to \lambda_i - \frac{d\theta_i}{d\tau}.$$
 (5)

In mean-field theory, the holon field condenses and it, together with the auxiliary field, acquires site- and time-independent values $b_i^a \to b_0$ and $i\lambda_i \to \delta E_{\rm f}$. The resulting mean-field Hamiltonian, $\hat{H}_{\rm MF}$, reads

$$\hat{H}_{\rm MF} = \sum_{\mathbf{k}} \hat{\Psi}_{\mathbf{k}}^{\dagger} \begin{pmatrix} \epsilon_{\mathbf{k}}^{\rm d} & b_0 \check{V}_{\mathbf{k}} \\ b_0 \check{V}_{\mathbf{k}}^{\dagger} & \epsilon_{\mathbf{k}}^{\dagger} \end{pmatrix} \hat{\Psi}_{\mathbf{k}}, \tag{6}$$

where $\epsilon_{\bf k}^{\rm f}=-E_{\rm f}^0+\delta E_{\rm f}-b_0^2k^2/2m_{\rm F}$ is a renormalized dispersion law of the f states. The spectrum consists of two bands $\epsilon_{\bf k}^\pm$, separated by direct gap $\Delta_{\bf k}=b_0({\rm Tr}[\check{V}_{\bf k}^\dagger\check{V}_{\bf k}]/2)^{1/2}$ and given by $\epsilon_{\bf k}^\pm=(\epsilon_{\bf k}^{\rm d}+\epsilon_{\bf k}^{\rm f})/2\pm\sqrt{(\epsilon_{\bf k}^{\rm d}-\epsilon_{\bf k}^{\rm f})^2/4+\Delta_{\bf k}^2}$. There is an indirect gap $\Delta_{\rm b}\sim\Delta_{\bf k}^2/E_{\rm d}^{\rm m}$ and it is assumed that the chemical potential of the TKI bulk lies within it. The set of self-consistent equations for b_0 and $\delta E_{\rm f}$ can be derived from the mean-field action and are given by

$$b_0^2 + \sum_{\alpha} \langle \hat{f}_{i\alpha}^{\dagger} \hat{f}_{i\alpha} \rangle = 1, \quad b_0 \delta E_{\rm f} = -\sum_{i\sigma\alpha} V_{ij}^{\sigma\alpha} \langle \hat{f}_{j\alpha}^{\dagger} \hat{d}_{i\sigma} \rangle.$$
 (7)

These equations should be supplemented by an additional one for the number of particles, $n_f + n_d = N_f$, which imposes full occupation of the band $\epsilon_{\mathbf{k}}^-.$ This equation fixes the position of the chemical potential for renormalized energy bands. The position of the chemical potential, and Fermi momentum are changing within gap opening, which can be expected from Luttinger's theorem arguments [42–46]. The phase diagram of the model, depending on the energy of localized states, degeneracy, and the lattice symmetry, has been extensively discussed recently [27,39,41,47–49]. Here we are interested in studying Kondo fluctuations relative to the mean-field saddle point. So, we treat the mean-field values b_0 and $\delta E_{\rm f}$ as parameters of the model and assume that the system is within TKI phase. Note the following hierarchy of energy scales: $\Delta_{\rm b} \ll \Delta_{\rm k} \ll E_{\rm d}^{\rm m}$. The shift of the f levels is of order $\delta E_{\rm f} \sim E_{\rm d}^{\rm m}$ and the Kondo temperature where the formation of an insulating band structure commences can be estimated as $T_{\rm K} \sim \Delta_{\rm b}$. The holon condensate b_0 appears at temperatures of order Δ_k , and smoothly increases with decreasing of temperature.

Kondo fluctuations in the TKI's bulk. Here we restore the dynamics to the fluctuations of holon condensate $\delta b_i^a(\tau) = b_i^a(\tau) - b_0$ and the constraint field $\delta \lambda_i(\tau) = \lambda_i(\tau) + \mathbf{i} \delta E_f$. A Lagrangian, describing Gaussian Kondo fluctuations and their interactions with electrons, is given by

$$L_{\text{fl}}^{\text{b}} = \sum_{ij} \bar{\Psi}_{i}(\tau) \begin{pmatrix} 0 & \check{V}_{ij} \delta b_{j}^{\text{a}}(\tau) \\ \delta b_{i}^{\text{a}}(\tau) \check{V}_{ij}^{\dagger} & \mathbf{i} \delta \lambda_{j}(\tau) \end{pmatrix} \Psi_{j}(\tau) + \sum_{i} \left\{ \delta E_{\text{f}} \left[\delta b_{i}^{\text{a}}(\tau) \right]^{2} + 2\mathbf{i} b_{0} \delta b_{i}^{\text{a}}(\tau) \delta \lambda_{i}(\tau) \right\}. \tag{8}$$

The "valence fluctuations" of the auxiliary field, λ_i (describing local fluctuations of f states' occupation) do not have a bare mass or bare field theory all together, and therefore its bare fluctuations are gapless. Note that an effective field theory for $\delta\lambda_i$ appears upon integrating out the matter fields—the f electrons. On the other hand, the fluctuations of the holon amplitude are massive and can be integrated out. Introducing Fourier components for fermionic and bosonic fields, we get a continuous low-energy part of the fluctuation Lagrangian

$$L_{\rm fl}^{\rm b} = \sum_{\mathbf{k}\mathbf{q}} \mathbf{i}\lambda_{\mathbf{q}}(\tau)\bar{f}_{\mathbf{k}+\mathbf{q}\alpha}(\tau)f_{\mathbf{k}\alpha}(\tau) + \sum_{\mathbf{q}} \frac{|\delta\lambda_{\mathbf{q}}(\tau)|^2}{2U_{\rm b}}.$$
 (9)

Here $U_{\rm b} = \delta E_{\rm f}/2b_0^2 n_{\rm b}$ corresponds to short-range repulsive interactions mediated by Kondo fluctuations and n_b is the three-dimensional concentration of f levels. It is very strong since the energy of interaction between two electrons on a cite is of order $\delta E_{\rm f}/b_0^2$ and considerably exceeds the gap, $\Delta_{\rm b}$. The value of the interaction decreases with the formation of the holon condensate. It can be shown that the interaction $U_{\rm h}$ is inverse proportional to the degeneracy N_f of the f states. So in the limit $N_{\rm f} \to \infty$ the slave-boson mean-field theory is exact. For a finite $N_{\rm f}$, the strong short-range interaction is important in both thermodynamics and transport if the bulk is metallic as was extensively discussed [30–34]. But if the bulk is insulating, electronic excitations at low temperatures $T \ll \Delta_b$ form a dilute gas, for which short-range interaction is unimportant. A TKI has an insulating bulk but topologically protected surface states. These Dirac surface states inherit interactions mediated by the Kondo fluctuations in the bulk.

The Dirac surface states. Now we consider a TKI with open boundary conditions (requiring that the electron's wave function vanish on the surface), and find the topological surface states from the mean-field Hamiltonian (6). If the surface is perpendicular to the z axis and the TKI occupies the z < 0 halfspace, the contribution of surface states to the field operator $\hat{\Psi}_s(\mathbf{r},z)$ is given by

$$\hat{\Psi}_{s} = \sum_{\mathbf{p}} \psi(z) e^{i\mathbf{p}\mathbf{r}} \begin{bmatrix} \hat{\chi}_{\mathbf{p}}^{\uparrow} \begin{pmatrix} u \\ 0 \\ iw \\ 0 \end{pmatrix} + \hat{\chi}_{\mathbf{p}}^{\downarrow} \begin{pmatrix} 0 \\ u \\ 0 \\ -iw \end{pmatrix} \end{bmatrix}. \tag{10}$$

Here $\hat{\chi}_{\mathbf{p}}^{\uparrow,\downarrow}$ are a pair of annihilation operators for surface states with the in-plane momentum \mathbf{p} and up/down spin. Projection of the bulk mean-field Hamiltonian onto these states yields the Dirac Hamiltonian $\mathcal{H}_s = v_s[\mathbf{p} \times \boldsymbol{\check{\sigma}}]_z$ with the velocity $v_s = 2uwb_0V \approx 2b_0^2V\sqrt{m_d/m_f}$, which is temperature dependent. The coherence factors u^2 and w^2 define contributions to

the surface states from the conduction and f states and are given by $u^2 = b_0^2 m_{\rm d}/(m_{\rm f} + b_0^2 m_{\rm d})$ and $w^2 = m_{\rm f}/(m_{\rm f} + b_0^2 m_{\rm d})$. Since $m_{\rm d} \ll m_{\rm f}$, then $u^2 \ll w^2$ and Dirac surface states mostly consist of the f-electron states. We find the wave function of the surface states exponentially decaying in the TI's bulk and vanishing on its surface as follows $\psi(z) = N \left[e^{\eta - z} - e^{\eta + z} \right]$. Here N is a normalization factor, and η_{\pm} are given by

$$\eta_{\pm} = \sqrt{\kappa_1 - \kappa_2 \pm \sqrt{(\kappa_1 - \kappa_2)^2 - \kappa_2^2}},$$
(11)

with $\kappa_1 = 2m_d m_f V^2/\hbar^2$ and $\kappa_2 = 2m_d m_f (E_d - E_f)/\hbar (m_f + b_0^2 m_d)$.

The full field operator $\hat{\Psi} = \hat{\Psi}_s + \hat{\Psi}_b$ has contributions from both the surface, $\hat{\Psi}_s$, and the interior, $\hat{\Psi}_b$. To project the bulk model onto the surface we insert $\hat{\Psi}$ into the action (8) and neglect terms that contain the bulk states, $\hat{\Psi}_b$. This approximation is justified, if the energy of surface states is deep inside the gap of the bulk spectrum. Neglecting variation of the fluctuation field in the transverse direction (on the length scale of order penetration length of the surface states), integrating the action over z, and integrating out massive fluctuations of the holon field's amplitude, δb_i^a , we obtain the continuous low-energy model for surface Dirac states:

$$L_{s}(\tau) = \sum_{\mathbf{p}} \bar{\chi}_{\mathbf{p}}(\tau) (\partial_{\tau} + v_{s}[\mathbf{p} \times \check{\boldsymbol{\sigma}}]_{z} - \mu) \chi_{\mathbf{p}}(\tau)$$

$$+ \sum_{\mathbf{p},\mathbf{q}} i w^{2} \lambda_{\mathbf{q}}(\tau) \bar{\chi}_{\mathbf{p}+\mathbf{q}}(\tau) \chi_{\mathbf{p}}(\tau) + \frac{w^{4}}{2U_{s}} \sum_{\mathbf{q}} |\lambda_{\mathbf{q}}(\tau)|^{2}.$$
(12)

Here $\chi_{\mathbf{p}} = (\chi_{\mathbf{p}}^{\uparrow}, \chi_{\mathbf{p}}^{\downarrow})^{\mathrm{T}}$ and μ is the chemical potential of surface states, $U_{\mathrm{s}} = \delta E_{\mathrm{f}} w^4 / 2 b_0^2 n_{\mathrm{s}}$ is the Fourier transform of an effective short-range interaction mediated by the Kondo fluctuations, and n_{s} is the surface concentration of f sites. As seen from Eq. (12), the valence fluctuations mediate repulsive interactions between the emergent Dirac excitations in the density-density channel, whose strength depends on the properties of the underlying mean-field state. With the formation of the holon condensate, Kondo fluctuations and interactions mediated by them are decreasing.

Now we discuss possible manifestations of these fluctuations/interactions in the observables. Since our theory is generic we cannot provide universal predictions, but we can discuss two distinct scenarios: either interactions lead to a phase transition or not.

Modification of the plasmon spectrum. If the effective repulsion is sufficiently weak and no phase transition takes place, Kondo fluctuations do not induce a reconstruction of the surface spectrum, but have only qualitative consequences (much like in the theory of Landau Fermi liquid). One such effect is a correction to the spectrum of collective modes on the surface. Since the two-dimensional Dirac electrons are charged, there is a long-range Coulomb interaction, which gives rise to collective charge oscillations—plasmons. This Coulomb interaction between Dirac surface states can be introduced by inserting an additional term in the action (12):

$$L_{\phi} = \sum_{\mathbf{p},\mathbf{q}} i\phi_{\mathbf{q}}(\tau)\bar{\chi}_{\mathbf{p}+\mathbf{q}}(\tau)\chi_{\mathbf{p}}(\tau) + \sum_{\mathbf{q}} \frac{|\phi_{\mathbf{q}}(\tau)|^2}{2U_{\mathbf{q}}}.$$
 (13)

Here $\phi_{\bf q}(\tau)$ is a bosonic field, which originates from the Hubbard-Stratonovich decoupling of the Coulomb interaction and plays the role of a scalar electrical potential; $U_{\bf q}=2\pi e^2/\epsilon q$ is the Fourier transform of the Coulomb potential in two dimensions and ϵ is an effective dielectric permittivity of the TKI surface. Fields $\phi_{\bf q}$ and $\lambda_{\bf q}$ can be incorporated in a single field $\Phi_{\bf q}=\phi_{\bf q}+w^2\lambda_{\bf q}$ and one of them is redundant and can be integrated out. After the integration the action for $\Phi_{\bf q}$ is given by

$$L_{\Phi} = \sum_{\mathbf{p}, \mathbf{q}} i \Phi_{\mathbf{q}}(\tau) \bar{\chi}_{\mathbf{p}+\mathbf{q}}(\tau) \chi_{\mathbf{p}}(\tau) + \sum_{\mathbf{q}} \frac{|\Phi_{\mathbf{q}}(\tau)|^2}{2(U_{\mathbf{q}} + U_{\mathbf{s}})}.$$
 (14)

In the random-phase approximation, the dispersion law of collective excitations satisfies the equation

$$1 - (U_{\mathbf{q}} + U_{\mathbf{s}})\Pi(\omega, \mathbf{q}) = 0, \tag{15}$$

where $\Pi(\omega, \mathbf{q})$ is the polarization operator of a Dirac electron gas, which has been calculated in Refs. [50] and [51] (see also Ref. [52], and references therein). The dispersion law of collective excitations depends on two dimensionless coupling constants $\alpha_s = e^2/\hbar v_s \epsilon$ and $\lambda_s = v_F U_s$, which describe the strength of the interactions. Here $v_F = \mu/2\pi(\hbar v_s)^2$ is the density of Dirac states at the Fermi level. For $v_s q \ll \omega \ll \mu$ the polarization operator can be approximated as $\Pi(\omega, \mathbf{q}) =$ $v_{\rm F}(v_{\rm s}q)^2/2\omega^2$. In the long-wavelength limit, the long-range Coulomb interaction dominates, but at $q \gtrsim q_0 = p_F \alpha / \lambda$, where $p_{\rm F}$ is the Fermi momentum, the short-range interaction takes over. For $q \lesssim q_0$, the plasmon dispersion is the standard in two dimensions square-root law, $\omega = v_s \sqrt{\alpha_s p_F q/2}$, while for $q_0 \lesssim q \ll p_{\rm F}$ it is a linear zero-sound-like mode, $\omega = v_{\rm s} \sqrt{\lambda_{\rm s}/2} q$. For estimates, we have used the following set of parameters, relevant to SmB₆ [16,39], $v_s \approx 0.5 \times 10^5$ m/s, $\delta E_{\rm f} \approx 0.15 \text{ eV}, \ b_0 \approx 0.05, \ \epsilon \approx 50, \ \Delta_{\rm b} \approx 1 \text{ meV}, \ n_{\rm s} \approx 4 \times 1 \text{ meV}$ 10^{14} cm⁻², which corresponds to the coupling constants $\alpha_{\rm s} \approx$ 0.7 and $\lambda_s \approx$ 1.3. The dispersion law of surface plasmons for $\alpha_s = 0.7$ and different values of parameter $\lambda_s = 0.1, 4, 12$ is presented in Fig. 1. If $\omega < v_s q$ or $\omega > 2\mu - v_s q$ the plasmons are strongly damped due to intraband (Landau damping) or interband transitions (the corresponding borders are shown by dotted lines in the figure). Since plasmons are undamped only for $q < p_{\rm F}$, the fluctuation-mediated interaction is important in that case only if $q_0 \ll p_F$, which is satisfied if $\alpha_{\rm s} \ll \lambda_{\rm s}$.

Magnetic phase transition in the Dirac liquid. While the gapless Dirac spectrum on a TKI's surface is protected by the time-reversal symmetry, this symmetry can be spontaneously broken due to sufficiently strong interactions (with the appearance of an out-of-plane spin polarization [53] or spin-density wave [54] order parameter). An out-of-plane spin polarization opens a gap in the Dirac spectrum, which can be energetically favorable since it reduces both the kinetic

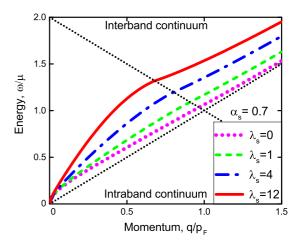


FIG. 1. (Color online) Dispersion law of plasmons on the surface of a TKI, calculated from (15), for dimensionless Coulomb coupling constant $\alpha_s = e^2/\hbar v_s \epsilon = 0.7$ and for different values of dimensionless short-range coupling constant $\lambda_s = v_F U_s = 0, 1, 4, 12$. Black dotted lines are borders of continuums $\omega < v_s q$ and $\omega > 2\mu - v_s q$, corresponding to intraband and interband single-particle excitations.

energy and exchange interaction energy of Dirac electrons. This state can also be interpreted as an excitonic insulator of electrons and holes from the conduction and valence bands of the surface Dirac spectrum, the possibility of which has been widely discussed in the context of graphene [55–57]. The critical value of short-range repulsion is [53] $U_s^* = 4\pi v_F \hbar/k_A$, where $k_{\Lambda} \approx \Delta_{\rm b}/\hbar v_{\rm F}$ is an ultraviolet cutoff of the linear spectrum and Δ_b is the indirect gap in the TKI's bulk. Using the same set of parameters as above, we get $U_{\rm s}/U_{\rm s}^* \approx 0.6$ and conclude that the short-range repulsion mediated by Kondo fluctuations is in the neighborhood of the critical value in realistic materials and may potentially drive the system into a magnetic phase. Furthermore, since the interaction strength is decreasing with lowering temperature, a highly unusual scenario of a double magnetic transition (where a ferromagnetic state would appear as an intermediate phase in a finite temperature window) is conceivable. Apart from this speculative scenario, another possibility is that the Dirac liquid on the TKI's surface remains magnetically ordered in the ground state, which may be consistent with the experiment of the Paglione group [28] in samarium hexaboride (where, however, the situation is complicated by the presence of three Dirac cones, whose Dirac points are offset relative to one another [39,40,58]).

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