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Theory of the Interfacial Dzyaloshinskii-Moriya Interaction in Rashba Antiferromagnets

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In antiferromagnetic (AFM) thin films, broken inversion symmetry or coupling to adjacent heavy metals can induce Dzyaloshinskii-Moriya (DM) interactions. Knowledge of the DM parameters is essential for understanding and designing exotic spin structures, such as hedgehog Skyrmions and chiral Néel walls, which are attractive for use in novel information storage technologies. We introduce a framework for computing the DM interaction in two-dimensional Rashba antiferromagnets. Unlike in Rashba ferromagnets, the DM interaction is not suppressed even at low temperatures. The material parameters control both the strength and sign of the interfacial DM interaction. Our results suggest a route toward controlling the DM interaction in AFM materials by means of doping and electric fields.

Relativistic spin-orbit coupling (SOC) is the foundation of spin orbitronics, a rapidly developing branch of spintronics [1, 2]. Anisotropic magnetoresistance [3] and the anomalous Hall effect [4] are established SOC transport phenomena. More recent discoveries include the spin Hall and inverse spin Hall effects [5], topological surface states [2, 6, 7], spin-orbit torques [1, 8–11], and chiral domain walls and Skyrmions [12–18]. These phenomena are essential to enable novel ultrafast, nonvolatile, nanoscale spin-based storage and computation devices.

The Dzyaloshinskii-Moriya (DM) interaction between localized spins in a magnetic material is induced by SOC [19, 20]. The DM interaction is important for stabilizing noncollinear magnetic structures in ferromagnets. The original proposal by Dzyaloshinskii and Moriya, however, concerned antiferromagnetic (AFM) systems with weak magnetization, i.e., weak ferromagnets; this weak magnetism can be explained in terms of antisymmetric exchange, which is also referred to as the DM interaction.

The energy associated with the DM interaction between two neighboring spins in a lattice, $S_1$ and $S_2$, can be written in the form of the mixed product $\mathcal{H}_{\text{DM}} = -D_{12} \cdot S_1 \times S_2$, where $D_{12}$ is called the DM vector. The DM interaction, therefore, favors a perpendicular orientation of neighboring spins. By contrast, the Heisenberg exchange interaction $\mathcal{H}_{\text{ex}} = J_{\text{ex}} S_1 \cdot S_2$ favors a collinear magnetic order, which can be either ferromagnetic (FM) for $J_{\text{ex}} < 0$ or AFM for $J_{\text{ex}} > 0$. The competition between the Heisenberg exchange interaction and the DM interaction leads to the formation of exotic structures such as chiral domain walls, helices, and Skyrmions [2].

In the bulk of a noncentrosymmetric magnetic crystal, the DM vector points along one of the lattice vectors and gives rise to so-called Bloch-like structures [19, 20].

![FIG. 1: (a) The Rashba-AFM model is a model for systems with inversion asymmetry: an AFM/HM bilayer (left) and an AFM thin film asymmetrically embedded between two different insulator layers (right). (b), (c) Electronic dispersion relations in the limit of strong exchange coupling (b) and in the limit of high SOC (c). The superscript $\eta = \pm$ refers to the conduction (blue) and valence (red) bands, whereas the superscript $s = \pm$ specifies chiral bands. In these two extreme limits, the band structure is almost isotropic for an AFM system with in-plane anisotropy.](image-url)
candidates for use in the next generation of spin-based memory and processing devices [46, 47] with a nanoscale element base.

The coexistence of strong SOC and spin ordering at the interfaces of AFM/heavy-metal (HM) bilayers makes such heterostructures particularly promising for low-dimensional spin orbitronics applications [48]. An AFM thin film sandwiched between insulators is also a common functional geometry for spin-orbitronics due to interfacial SOC. The interfacial SOC in such a system is effectively described by Rashba SOC [39]. For magnetic films, the two-dimensional (2D) Rashba model captures the main physics and trends of SOC with a broken inversion symmetry [40], such as fieldlike and dampinglike spin-orbit torques [11, 9, 10], intrinsic spin Hall effects [5], intrinsic anomalous Hall effects [50], inverse Faraday effects [51], and magnetic anisotropy [52].

In this Letter, we develop a framework for computing the interfacial DM interaction in AFM layers with inversion asymmetry. The typical system illustrated in Fig. 1 is described on the basis of an effective 2D AFM-Rashba Hamiltonian. We find that both the sign and the magnitude of the DM vector depend on the ratio relating three energy scales: the chemical potential, the $s$-$d$ exchange interaction, and the Rashba SOC strength. In particular, the strong dependence on the chemical potential suggests that the DM interaction can be tuned by modifying the electron density by means of doping or voltage gating [54].

A generic effective 2D Hamiltonian describing itinerant electrons in the AFM layer [see Fig. 1(a)] can be written as

$$
\mathcal{H} = \mathcal{H}_{\text{kin}} + \mathcal{H}_{\text{sd}} + \mathcal{H}_{\text{so}},
$$

where $\mathcal{H}_{\text{kin}}$ is the kinetic energy of the electrons, $\mathcal{H}_{\text{sd}} = \int_0^{\beta} \mathcal{J}_{\text{sd}} \Gamma d\tau \mathbf{n}$, $\Gamma$ describes an effective interaction with a strength $J_{\text{sd}}$ between the spins of the itinerant $s$ electrons and the localized $d$ electrons [50], and $\mathcal{H}_{\text{so}}$ describes the SOC. The operator $\Gamma$ is the direct product of the electron spin operator and the sublattice position operator, which, in the case of an AFM system, accounts for the effects of sublattice staggering. The unit vector $\mathbf{n}$ is the order parameter, which can represent either the total magnetization in a ferromagnet or the staggered magnetization in an antiferromagnet.

We compute the electronic contributions to the DM interaction parameter $D$ and to the exchange stiffness $A$ in the following way. First, we evaluate how the itinerant electrons influence the magnetic subsystem by finding an effective action. We expand the effective action up to linear order with respect to the deviation of the spins from their equilibrium direction. The corresponding susceptibility tensor describes the influence of the electronic degrees of freedom on the localized magnetic moments. A linear expansion of the susceptibility tensor in spatial gradients of $\mathbf{n}$ defines the DM interaction strength $D$. The contribution of the itinerant electrons to the exchange stiffness $A$ is extracted from the second-order expansion in spatial gradients.

The action $S$ defines the system partition function $Z = \int d[\Phi^*] d[\Phi] d[n] e^{-S[\Phi^*,\Phi,n]/\hbar}$, where $\Phi$ is the Grassmannian coherent-state spinor and $\hbar$ is Planck’s constant. In the $s$-$d$ approach to magnetic systems, the action is decomposed into the sum $S = S_F + S_B$, where $S_F[\Phi^*,\Phi,n]$ is the fermionic action corresponding to the Hamiltonian of Eq. (1), which also includes the $s$-$d$ coupling, and $S_B[n]$ is the bosonic action describing the dynamics of the localized spins (magnons) in the absence of itinerant electrons. In our model, it is the coupling between the itinerant electrons and the local moments that determines the DM interaction, which is also directly linked to the SOC of the itinerant electrons. We will not specify the bosonic part of the action $S_B[n]$ since it is irrelevant for the subsequent discussion.

The fermionic action reads

$$
S_F = \int_0^{\beta} d\tau d\tau' \int d\mathbf{r} d\mathbf{r}' \Phi^*_{\tau',\tau} [\mathbf{G}^{-1}_{\mathbf{r},\tau';\mathbf{r}',\tau} \Phi_{\mathbf{r}',\tau'}],
$$

where $\beta = 1/k_B T$ is the inverse temperature and $\tau$ is the imaginary time. The inverse Green’s function operator is $\mathbf{G}^{-1}_{\mathbf{r},\tau';\mathbf{r}',\tau'} = - (\hbar \partial_{\tau'} + \mathcal{H}) \delta(\mathbf{r} - \mathbf{r}') \delta(\tau - \tau')$ in terms of the Hamiltonian of Eq. (1). We compute the effective theory for the vector field $\mathbf{n}(\mathbf{r})$ by integrating out the fermionic degrees of freedom. This standard procedure results in an additional, effective contribution to the bosonic action of the form $\Delta S_F^{\text{eff}}[\mathbf{n}] = \int_0^{\beta} d\tau \int d\mathbf{r} \left(-\hbar \text{Tr}[\ln(\mathbf{G}^{-1})]\right)$. Below, we analyze $\Delta S_F^{\text{eff}}[\mathbf{n}]$ and its influence on the magnet in the AFM-Rashba model.

In our system, the symmetry-breaking direction (z direction) is perpendicular to the plane [81]. Without the loss of generality, we choose the $x$ axis to be in the direction of the in-plane vector field $\mathbf{n}$. A small deviation of the unit vector from its equilibrium direction is, then, parameterized by $\mathbf{n} = \hat{x} + \delta\mathbf{n}$, where $\delta\mathbf{n} = (-\delta n_y^2 + \delta n_z^2)/2, \delta n_y, \delta n_z$. The effective action is obtained from a perturbation with respect to $\delta\mathbf{n}$, [53] that holds irrespective of the value of the exchange

\[\text{FIG. 2: (a) The Feynman diagram corresponding to Eq. (1) for a single-loop polarizability tensor. (b) The Feynman diagram corresponding to Eq. (5), describing the interfacial DM interaction for a system with inversion symmetry breaking in the z direction and an order parameter vector n in the x direction.}\]
strength $J_{sd}$. The effective action is conveniently established from the Dyson equation, $G^{-1} = G^0^{-1} - \Sigma$, where the unperturbed Green’s function refers to $\delta n = 0$ and the self-energy $\hbar \Sigma = J_{sd} \delta n \cdot \mathbf{q} \delta(k - k') \delta(\tau - \tau')$ is introduced.

Computing the self-energy up to the second order in $\delta n$ yields [57]

$$\Delta \Sigma^{\text{eff}}[n] = \hbar \beta \sum_{k \neq 0, l} \delta n_{k, \nu}^a \Pi_{k, \nu}^{a,b} \delta n_{-k, -\nu}, \quad (3)$$

where the indices $a, b = (y, z)$ denote the transverse vector components with respect to the equilibrium $\hat{x}$ direction, $\nu = 2\pi n / \beta$ denotes the bosonic Matsubara frequencies, and

$$\Pi_{k, \nu}^{b} = J_{sd}^{2} \frac{2\hbar \beta}{\gamma} \sum_{q, n} \frac{1}{\xi_{k}^{2}} \left[ \Gamma_{x} G_{q, i\omega_{n}}^{0} \Gamma_{y} G_{q-k, i\omega_{n}+i\nu}^{0} \right] \quad (4)$$

is the dynamical susceptibility tensor, pictured schematically in Fig. 2a. Here, $G_{q, i\omega_{n}}^{0} = (i\hbar \omega_{n} - \mathcal{H})^{-1}$ is the equilibrium Green’s function, and the $\omega_{n} = (2n + 1)\pi / \beta$ are the fermionic Matsubara frequencies. We compute the sum over the fermionic Matsubara frequencies by using the identity $\sum_{\nu} (i\hbar \omega_{\nu} - E)^{-1} / \beta = f(E)$, where $f(E)$ is the Fermi distribution.

By expanding the static limit, $\nu = 0$, on the spin susceptibility to the second order in the wave vector $k$, we find both the electronic contribution to the symmetric Heisenberg exchange stiffness, which is determined by the symmetric terms in the diagonal elements of the susceptibility tensor, and the antisymmetric exchange interaction (DM interaction), which is determined by the antisymmetric terms in the off-diagonal elements. From the partition function of the canonical ensemble, $Z = \int d[n] e^{-\delta S^{\text{eff}}[n]} / \hbar = \int d[n] e^{-\beta F[n]}$, we obtain

$$\delta n \delta S^{\text{eff}} / \hbar = \beta \delta n \delta F[n],$$

where the micromagnetic free energy, including the stiffness and the DM interaction, is $F[n] = \int d^2r (A(\nabla \mathbf{n})^2 - D \mathbf{n} \cdot (\hat{x} \times \nabla) \times \mathbf{n})$. By comparing the microscopic free energy with the expression for the effective action, Eq. (3), we define the micromagnetic parameters $A$ and $D$, which characterize the free carrier contributions to the exchange stiffness and the DM interaction, respectively.

Upon expanding the off-diagonal elements of the tensor $\Pi$ to the first order in the wave vector $k$ [see Fig. 2b], we obtain the relation

$$D = \frac{i}{\beta} \left. \frac{\partial \Pi_{k, \nu}^{y}}{\partial k_x} \right|_{k=0} = - \left. \frac{i}{\beta} \frac{\partial \Pi_{k, \nu}^{y}}{\partial k_y} \right|_{k=0} \quad (5a)$$

$$= \frac{J_{sd}^{2}}{2\beta} \sum_{q, n} \frac{1}{\xi_{k}^{2}} \left[ \Gamma_{y} G_{q, i\omega_{n}}^{0} \Gamma_{x} G_{q-k, i\omega_{n}+i\nu}^{0} \right], \quad (5b)$$

where $v_y = \hbar^{-1} \partial \mathcal{H}_{q} / \partial q_y$ is the y component of the velocity operator. From the second-order terms, we obtain the electron contribution to the exchange stiffness:

$$A = - \left. \frac{\partial^{2} \Pi_{k, \nu}^{y}}{\partial k_{x}^{2}} \right|_{k=0} = - \left. \frac{\partial^{2} \Pi_{k, \nu}^{y}}{\partial k_{y}^{2}} \right|_{k=0}. \quad (6)$$

In our model, $A$ describes the contribution to the AFM exchange interaction from a superexchange-type interaction between the localized spins in the AFM layer via the itinerant spins.

We should emphasize here that in this approach, we have ignored the spin fluctuations of the localized AFM spins, which is a valid omission as long as the system temperature is much less than the critical Néel temperature.

To model an AFM system with interfacial SOC, we use the 2D AFM-Rashba Hamiltonian [10 11] on a square lattice:

$$\mathcal{H} = \gamma_{k} \tau_{z} \sigma_{0} + J_{sd} \tau_{x} \mathbf{n} - \alpha_{R} \tau_{z} (\sigma \times k) \cdot \hat{z}, \quad (7)$$

where $\sigma$ and $\tau$ are the vectors of Pauli matrices representing the spin and AFM sublattice degree of freedom, respectively; $\sigma_{0}$ is the identity matrix; $\mathbf{n}$ is the staggered order parameter (the normalized Néel vector); and $\alpha_{R}$ is the strength of the Rashba SOC. The kinetic energy of the itinerant electrons is $\gamma_{k} = a^{2}(t(k^{2} - k_{z}^{2}))$, where $t$ is the nearest-neighbor hopping energy and $k_{0} = 2/a$, with $a$ being the lattice constant.

The band structure of the AFM-Rashba Hamiltonian of Eq. (7) is, in general, anisotropic. It is convenient to parameterize the four spectral branches as follows:

$$E_{k}^{\pm} = \pm \sqrt{s^{2} + J_{sd}^{2} + \alpha_{R}^{2}k^{2} + 2s \alpha_{R}k_{x}}, \quad (8)$$

where $s, \eta = \pm 1$ are the spin chirality and electron/hole band indices, respectively. We also introduce $\xi_{k} = (\gamma_{k}^{2} + J_{sd}^{2} \cos^{2} \phi)^{1/2}$ with the in-plane wave vector $k$ parameterized by the angle $\phi$, such that $k = k(\cos \phi, \sin \phi, 0)$.

Let us now analyze the expressions of Eqs. (6) and (7) for the AFM-Rashba Hamiltonian of Eq. (7) in the limit of weak spin-orbit interaction, $\left\{ k_{0}\alpha_{R}, J_{sd} \right\} \ll \min\{t, \epsilon_{F}\}$, and at zero temperature. Using the relation $\Gamma_{x} = \tau_{z} \sigma_{0}$, we obtain the relation

$$D = - \frac{A k_{0}^{2}}{4t} \alpha_{R}, \quad (9)$$

where the stiffness parameter is

$$A = \frac{t_{sd}^{2}}{2\pi \epsilon_{F}^{2}} \begin{cases} 2 - \epsilon_{F}^{2} / 16t^{2}, & \epsilon_{F} < 4t, \\ 1, & \epsilon_{F} > 4t, \end{cases} \quad (10)$$

which is manifestly independent of the SOC strength in the limit of weak spin-orbit interaction. Here, the Fermi energy $\epsilon_{F}$ is measured with respect to the center of the energy gap [see Figs. 10(b) and (c)]. In this regime, the
FM chiral bands are occupied (11) with those for a Rashba-FM system [58]. If both
\( \epsilon \)s-d for different values of the
energy gaps at \( k = 0 \) and \( k = \pm k_0 \) are equal to \( 8t \) and
\( 2J_{sd} \), respectively.
For the case in which the Fermi energy lies within the
band gap, i.e., for \( |\epsilon_F| < J_{sd} \), and in the limit of \( k_0 \alpha_R \ll J_{sd} \ll t \), we obtain
\[ D = \left( \frac{k_0^2}{2\pi} \right) \alpha_R, \] (11)
which is of the opposite sign compared with the metallic
regime of Eqs. (9) and (10).
It is instructive to compare the results of Eqs. (9)-
(11) with those for a Rashba-FM system [58]. If both
FM chiral bands are occupied (\( \epsilon_F > J_{sd} \)), the interfacial
DM interaction in the Rashba-FM model vanishes at
zero temperature due to the exact cancellation between
the Fermi surface and Fermi sea contributions [58]. How-
ever, such a cancellation is absent in the AFM-Rashba
model, giving rise to a finite result for Eq. (9), by virtue
of an additional contribution from the valence bands. For
\( \epsilon_F < J_{sd} \), i.e., when only the lowest FM chiral band is
occupied, the DM interaction in the weak SOC limit and
at zero temperature is finite and linearly proportional to
\( A \) [20] [55] [56] [58].
On the other hand, in the Dirac model of an
FM/topological-insulator bilayer characterized by the
Hamiltonian of the Rashba SOC symmetry, the DM in-
teraction vanishes inside the gap but remains finite out-
side the gap, even at zero temperature [59] [60]. Thus,
we conclude that the DM interaction exhibits a quali-
tatively different behavior in antiferromagnets compared
with that in ferromagnets.
In Fig. 3 we illustrate the behavior of the DM interaction
on the basis of a numerical analysis of Eq. (5) beyond
the weak SOC regime of Eqs. (9) [11], where typical
material parameters are assumed: \( t = 3 \) eV, \( \epsilon_F = 2 \) eV,
\( J_{sd} = 1 \) eV, \( a = 0.4 \) nm and \( k_0 \alpha_R = 0.1 \) eV [10]. The
results indicate a rather large interfacial DM interaction,
with \( D \approx -1.6 \) pJ/m. Moreover, when the Fermi energy
lies within the gap, i.e., \( |\epsilon_F| < J_{sd} \), we find an order-
of-magnitude enhancement of the DM interaction with
the opposite sign, \( D \approx 12.7 \) pJ/m. Importantly, the esti-
mated strength of the DM interaction is at least an order
of magnitude larger than that for a Rashba-FM layer
with the same parameters [55].
For small Fermi energies, the DM interaction is pos-
tive and almost independent of the Fermi energy (see
the top panel in Fig. 5), as might be expected from
Eq. (11). In the metallic regime (i.e., for Fermi ener-
gies well above the gap), the DM interaction is nega-
tive, and its strength decreases in proportion to \( \epsilon_F^{-2} \),
in good agreement with Eqs. (9) and (10). The sign inver-
sion of the DM interaction is rather sharp and occurs
at \( \epsilon_F = \sqrt{J_{sd} \alpha_R^2 + J_{sd}^2|\alpha_R|} \). The bottom panel of Fig. 3 also
confirms that the strength of the DM interaction is lin-
early proportional to \( \alpha_R \) in the weak SOC regime, in
agreement with Eq. (9). Thus, we conclude that the DM
interaction in an AFM material may vary by orders of
magnitude depending on the material parameters, as il-
ustrated in Fig. 3.
Our calculations also show that, unlike in FM-Rashba
systems [58], the temperature dependence of the DM in-
teraction in the AFM-Rashba model of Eq. (7) is weak
due to a large contribution to the DM interaction from
the valence bands (not shown). As we have already dis-
cussed, this is correct if the system temperature is much
less than the Néel temperature, i.e., in a regime in which
the spin fluctuations of the AFM layer are suppressed.

Controlling the DM interaction is essential for engi-
neering chiral magnetic structures. If the DM interaction
parameter exceeds a certain critical value, which is de-
termined by the Heisenberg exchange interaction and the
uniaxial anisotropy, then the ground state changes from
a collinear configuration to either a helimagnetic state or
a Skyrmion lattice. A weaker DM interaction enables
the stabilization of isolated Skyrmions in a metastable
state [15] [18]. In chiral magnets, the sign of the DM inter-
action determines the direction, or handedness, of spin
rotation.

The asymmetry of the spin-wave dispersion in spin-
polarized electron energy-loss spectroscopy and Brillouin
light scattering is a measure of the DM interaction in FM
systems [27,30,42]. Although the interfacial DM interactions in a few FM/HM bilayers have been experimentally studied in recent years [29,42], we are not aware of similar measurements in AFM heterostructures. However, the first observation of a bulk DM interaction in an AFM system, namely, the noncentrosymmetric $\alpha - Cu_2V_2O_7$, has recently been reported in Ref. [43] based on inelastic neutron scattering. Very recently, a large DM interaction in Fe/Ir bilayers on Rh(001) has been predicted on the basis of ab initio calculations [61]. We hope that our work will stimulate new experiments and ab initio works on such AFM heterostructures.

In summary, we have computed the DM interaction, both analytically and numerically, using an effective model for AFM layers with interfacial Rashba SOC. In the AFM-Rashba model, the induced interfacial DM interaction appears as a Lifshitz-type invariant term. Our results show that both the sign and the strength of the DM interaction may be tuned by modifying the electron density, e.g., by applying a gate voltage or through doping. This tunability implies that a rich variety of chiral magnetic structures can emerge in layered AFM/HM systems with different interfacial charge densities, $s$-$d$ exchange interactions, and SOC interactions.

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In general, the effective action has two extra terms related to the spin susceptibilities with a long wave-length nature, i.e., \( k = 0 \), which have no contributions in the exchange stiffness and DM interaction: 
\[
\Delta S_{\text{eff}}[\mathbf{n}] = \hbar \delta n \left( \delta n \chi^{s} + \sum_i \delta n_i \Pi_{\alpha\beta}^{sd} \delta n^a_{\nu_i} \right).
\]
The static spin susceptibility related to the first-order term in \( \delta n^a \) is 
\[
\chi^a = \left( J_{sd}/\hbar \beta \right) \sum_{q,n} \frac{1}{\nu} \text{Tr}[\Gamma_a G^0_{q,i\omega_n} ],
\]
and the dynamic spin susceptibility related to the second-order term in \( \delta n^a \) is 
\[
\Pi_{\alpha\beta}^{sd} = \sum_{q,n} \left\{ - \left( J_{sd}/2\hbar \beta \right) \text{Tr}[\Gamma_a G^0_{q,i\omega_n} \delta_{ab} + \left( J_{sd}/2\hbar \beta \right) \text{Tr}[\Gamma_a G^0_{q,i\omega_n} \Gamma_b G^0_{q,i\omega_n+i\omega} ]] \right\}, \text{ where } \delta_{ab} \text{ is the Kronecker delta function.}
\]