Conductance, Valley and Spin polarization and Tunnelling magneto-resistance in ferromagnetic—normal—ferromagnetic junctions of silicene

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We investigate charge conductance and spin and valley polarization along with the tunnelling magneto-resistance (TMR) in silicene junctions composed of normal silicene and ferromagnetic silicene. We show distinct features of the conductances for parallel and anti-parallel spin configurations and the TMR, as the ferromagnetic—normal—ferromagnetic (FNF) junction is tuned by an external electric field. We analyse the behavior of the charge conductance and valley and spin polarizations in terms of the independent conductances of the different spins at the two valleys and the band structure of ferromagnetic silicene and show how the conductances are affected by the vanishing of the propagating states at one or the other valley. In particular, unlike in graphene, the band structure at the two valleys are independently affected by the spin in the ferromagnetic regions and lead to non-zero, and in certain parameter regimes, pure valley and spin polarizations, which can be tuned by the external electric field. We also investigate the oscillatory behavior of the TMR with respect to the strength of the barrier potential (both spin-independent and spin-dependent barriers) in the normal silicene region and note that in some parameter regimes, the TMR can even go from positive to negative values, as a function of the external electric field.

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I. INTRODUCTION

A close cousin of graphene, silicene has been attracting a lot of attention in recent times both theoretically and experimentally\textsuperscript{1–4}, due to the possibility of new applications, given its compatibility with silicon based electronics. Although earlier theoretical analyses\textsuperscript{5–7} had predicted the possibility of silicene and even germanene, stanene, etc, interest in this subject rose after experimentalists observed hexagonal structure in silicene sheets deposited on a silver substrate\textsuperscript{2–4}. Unlike graphene, silicene does not have a planar structure; instead it forms a buckled structure due to the large atomic radius of silicon\textsuperscript{8}. The possibility of using silicene as a spintronic device has pre-
dicted in a normal silicene junction\textsuperscript{2–4}. Moreover, interest in this area is essentially due to the realisation that devices based on the spin degree of freedom are almost dissipationless unlike those based on the charge degree of freedom. The possibility of using silicene as a spintronic device has been reported very recently in Ref.\textsuperscript{15–19} due to its strong spin-orbit coupling. Another important quantity in spintronics is the tunnelling magneto-resistance (TMR) which occurs at junctions between materials \textit{i.e.}, ferromagnet—normal metal—ferromagnet (FNF) junctions. The resistance of the junction is different for parallel and anti-parallel spin configurations, and this difference in resistance can be experimentally measured.

In recent years, spin based electronics or spintronics has become a prominent field of research both theoretically and experimentally\textsuperscript{14}. The upsurge of activity in this area is essentially due to the realisation that devices based on the spin degree of freedom are almost dissipationless unlike those based on the charge degree of freedom. The possibility of using silicene as a spintronic device has been reported very recently in Ref.\textsuperscript{15–19} due to its strong spin-orbit coupling. Another important quantity in spintronics is the tunnelling magneto-resistance (TMR) which occurs at junctions between materials \textit{i.e.}, ferromagnet—normal metal—ferromagnet (FNF) junctions. The resistance of the junction is different for parallel and anti-parallel spin configurations, and this difference in resistance can be experimentally measured.

In this paper, we study charge conductance and spin and valley polarizations along with the TMR in silicene junctions, in particular the FNF junctions, as the external electric field is tuned through the system. We model our FNF setup within the scattering matrix formalism\textsuperscript{20}. A similar setup has been utilised in graphene\textsuperscript{21} to study TMR, while valley polarization has been studied in a normal—ferromagnet—normal (NFM) junction in silicene\textsuperscript{22}. In Ref. \textsuperscript{22}, the author has investigated the con-
Ez Hamiltonian given by proximity effect when it is placed in proximity with a
in Fig. 1. Ferromagnetism is induced in silicene by the
conclusions.
the summary of our numerical results followed by the
various parameter regimes. Finally in Sec. V, we present
and valley polarizations and TMR in the FNF set up for
present our numerical results for the conductances, spin
polarizations (\(S = 1\)) and in right region for parallel
\(s = 1\) or \(P\) or \(AP\) spin configurations of magnetization re-

I. INTRODUCTION

Due to the buckled structure of silicene, the atoms in
the two sublattices respond differently to an externally
applied electric field \(E_z\). Thus \(iE_x\) is the potential
difference between the two sublattices A and B due to this
applied electric field where \(l\) is the separation between
the two sublattices. Hence, the potential difference is a tun-
able parameter and can be tuned by an external electric
field\(^1\). Also, when the Fermi energy is close to the Dirac
point, at the critical electric field \(E_z^c = \eta \sigma \lambda_{SO}\), one of
the valleys in silicene is up-spin polarized and the other
one is down-spin polarised\(^7\). Here \(\eta = \pm 1\) denotes the \(K\)
and \(K'\) valleys respectively and \(\sigma = \pm 1\) denotes the spin
indices. \(V(x)\) denotes the profile for the potential barrier
in the normal silicene region and \(h(x)\) corresponds to the
exchange splitting, or the energy difference between the
up and down spin electrons in the FS regions. Note how-
ever, that in real materials the proximity of a ferromagnet
to silicence can actually change the band structure of sil-
icene itself. In that case, the only way to proceed will be
to perform first principles calculations as have been done
in graphene\(^24\).

II. MODEL AND BAND STRUCTURE

We study the FNF junction set up in silicene as shown
in Fig. 1. Ferromagnetism is induced in silicene by the
proximity effect when it is placed in proximity with a
magnetic insulator, which we model by the following
Hamiltonian given by\(^22\)

\[
H = \hbar v_F (\eta k_x \hat{\tau}_z - k_y \hat{\tau}_y) + (\epsilon_l E_z - \eta \sigma \lambda_{SO}) \hat{\tau}_z + V(x) - h(x),
\]

(1)

where \(v_F\) is the Fermi velocity of the charge carriers in
silicene, \(e\) is the charge of the electron and \(\eta, \sigma\) correspond
to the valley and spin indices and \(\hat{\tau}\) corresponds to the
sublattice (pseudospin) Pauli matrices. \(\lambda_{SO}\) is the pa-
rameter that specifies the spin-orbit coupling in silicene.

We now consider the geometry shown in Fig. 1 and
assume that the system is translationally invariant along
the \(y\) direction. The interface between the normal and
the FS are located at \(x = 0\) and \(x = L\) where \(L\) the
length of the normal silicene region sandwiched between
the ferromagnetic patches. Here \(V(x) = U\Theta(x)\theta(L-x)\)
is the profile of the potential barrier modelled in the nor-
mal silicene region and \(h(x) = h[\sigma s\Theta(-x) + \sigma' s' \Theta(x-L)]\)
denotes the exchange field or Zeeman field in the two fer-
romagnetic regions with \(s = s'\) corresponding to the par-
allel (P) configuration and \(s = -s'\) corresponding to the
antiparallel (AP) spin configurations of magnetization re-
spectively. We show the schematic of up and down spin
in left region (for \(s = 1\)) and in right region for parallel
\((s' = 1\) or P or \(uw\)) and antiparallel \((s' = -1\) or AP or
\(ud\)) configurations in FNF junction for both \(E < h\) and
\(E > h\) regions in Fig. 2. The orientation of the exchange

FIG. 1. (Color online) Schematic of the FNF junction in
silicene in which two ferromagnetic patches (dark brown, dark
grey) have been deposited on two sides of a normal silicene
sheet (cyan, light grey) to induce ferromagnetism in it. Here
\(L\) is the length of the normal silicene region.

FIG. 2. (Color online) Schematic of the configurations with
P (\(uw\)) and AP (\(ud\)) spin polarizations for one of the valleys
of the FNF silicene junction.

We perform first principles calculations as have been done
in graphene\(^24\).
field in the left region is kept fixed by keeping \( s = 1 \) and then \( s' = 1 \) and \( s' = -1 \) in the right region corresponds to the parallel (P or \( uu \)) and antiparallel (AP or \( ud \)) configuration respectively. Note that the \( E < h \) line crosses the same band in the third region for the P configuration, but crosses the other band for the AP configuration. As reported earlier\(^ {21} \), this feature gives rise to negative TMR in graphene. In silicene, also, the same feature is responsible for negative TMR which we discuss later in Sec. IV.

![Image](https://via.placeholder.com/150)

**FIG. 3.** (Color online) (a) to (d) gives the schematic of the band structure for \( s = +1 \), at \( K \) (for both \( \uparrow \) and \( \downarrow \) spin) and \( K' \) (for both \( \uparrow \) and \( \downarrow \) spin) valleys for ferromagnetic silicene for four different values of the dimensionless parameter \( E_z/h \) (red \((0.1)\), green \((0.5)\), blue \((1.0)\) and magnetica \((2.0)\)). On the other hand \((e)\) to \((h)\) gives the same for \( s = -1 \). We use these diagrams to qualitatively explain the dependence of conductances on the electric field as mentioned in the text.

In Fig. 3 we demonstrate the spin polarization of both the \( K \) and \( K' \) valleys for FS with the magnetisation directions up \((u\) defined by \( s = +1\)) and down \((d\) defined by \( s = -1\)) via an energy band diagram. The diagram clearly shows that different spin orientations behave differently in the \( K \) and \( K' \) valleys at different values of the exchange field \( h \). To visualize this picture, we fix \( \lambda_{SO}/h = 0.5 \) and show the dispersion for \( K \uparrow, K \downarrow, K' \uparrow \) and \( K' \downarrow \) for four different values of the tunable parameter \( E_z/h \), for the cases \( u \) and \( d \). Note that we use \( \uparrow \), \( \downarrow \) to denote the spins of the incoming (also reflected and transmitted) charge carriers and we use \( u, d \) to denote the orientation of magnetic exchange. For the \( u \) or \( s = +1 \) case, at \( E_z/h = \lambda_{SO}/h = 0.5 \) both \( K \uparrow \) and \( K' \downarrow \) have a vanishing band gap. On the other hand, for \( E_z/h = 0.1 \), \( E_z/h = 1 \), and \( E_z/h = 2 \), the valleys at \( K \uparrow, K \downarrow, K' \uparrow, K' \downarrow \) are all gapped. Also note that due to the exchange splitting \( h \), the \( K' \) valley is shifted upwards for the \( \uparrow \) spin while the \( K \) valley is shifted downwards for the \( \uparrow \) spin. The case for \( d \) or \( s = -1 \) is the other way around. Hence, it is clear that unlike in graphene, the contributions to the conductances from various spin configurations will not be identical for the \( K \) and \( K' \) valleys.

### III. THE SCATTERING MATRIX

We model our FNF setup within the scattering matrix formalism\(^ {20} \) where we match the wave functions at each ferromagnet−normal interface to obtain the scattering matrix and find the conductances and the TMR. The wave functions for the valley \( \eta \) in each of the three regions, \( x < 0, 0 < x < L \) and \( x > L \) can be written as

\[
\psi_i = a_i e^{ik_i x} \begin{pmatrix} \eta k_i S_i e^{i\eta \theta_i} \\ \tau_i \end{pmatrix} + b_i e^{-ik_i x} \begin{pmatrix} -\eta k_i S_i e^{-i\eta \theta_i} \\ \tau_i \end{pmatrix},
\]

(2)

where \( a_1 = 1, b_1 = r \) for \( x < 0, a_2 = a, b_2 = b \) for \( 0 < x < L \) and \( a_3 = t, b_3 = 0 \) for \( x > L \). Note that we keep track of the sign of the charge carriers by including the index \( S_i \) in all the regions (the sign of the charge carriers changes from electron-type to hole-type when \( S_i \) is negative in any region). This actually happens for the anti-parallel configuration when energy of the incident charge carrier is below the induced magnetic field energy \( \sigma \) \( i.e. \ (E < h) \). This charge reversal actually qualitatively changes the conductances, as was shown in graphene\(^ {21} \).

We obtain the scattering matrix both for \( E > h \) and \( E < h \) by matching the wave functions (see Eq.(2)) at \( x = 0 \) and \( x = L \) and solving Eq.(3) numerically.

\[
\begin{bmatrix}
\sqrt{2E \tau_1} e^{i\eta \theta_1} & -\eta k_2 S_2 e^{-i\eta \theta_2} & \eta k_2 S_2 e^{-i\eta \theta_2} & 0 \\
2\eta k_2 S_2 e^{i\eta \theta_2} e^{ik_2 x} & -\tau_2 & -\tau_2 & 0 \\
0 & 2\eta k_2 S_2 e^{-i\eta \theta_2} e^{-ik_2 x} & -\tau_2 & 0 \\
\end{bmatrix}
\begin{bmatrix}
\begin{bmatrix} r \\ a \\ b \\ t \end{bmatrix}
\end{bmatrix} =
\frac{1}{\sqrt{2E \tau_1}}
\begin{bmatrix}
\begin{bmatrix} -\eta k_1 S_1 e^{i\eta \theta_1} \\ 0 \\ 0 \\ 0 \end{bmatrix}
\end{bmatrix}
\]

(3)

Further,

\[
k_i = \sqrt{E_i^2 - (eE_z - \eta \sigma_i \lambda_{SO})^2}
\]

and

\[
\tau_i = E_i - (eE_z - \eta \sigma_i \lambda_{SO})
\]

(4)

with \( \sigma_1 = \sigma, \sigma_3 = \sigma', E_1 = E + sh, E_3 = E + s'h \) and \( S_i = \text{sgn}[E_i - (eE_z - \eta \sigma_i \lambda_{SO})] \). Since momentum is conserved in the \( y \) direction and does not change, it is convenient to write the \( x \)-component of the wave-vectors
$$k_{x_1} = \sqrt{k_{x_1}^2 - k_y^2}. \quad (5)$$

where $k_y$ is the conserved momentum in the $y$ direction.

In Eqs. (4 and 5) we consider $i = 1, 3$ only. For simplicity we assume that $E_2 = 0$ in the middle region which makes the momentum in the middle region independent of valley and spin ($\sigma_2$). Hence for the central region,

$$k_2 = \sqrt{E_2^2 - \lambda_{SO}^2},$$

$$k_{x_2} = \sqrt{k_{x_2}^2 - k_y^2},$$

and $\tau_2 = E_2$. \quad (6)

where $E_2 = E - U$ and $U$ is the height of the potential barrier in the normal silicene region.

IV. NUMERICAL RESULTS

In this section we present our numerical results for the FNF junction for different parameter regimes. We first study the model described in Eq.(1) which has a spin-independent barrier in the normal silicene region. We compute the conductance using the transmission coefficients obtained in Eq.(3) for both the parallel ($s = s'$) and the anti-parallel ($s = -s'$) configurations of spins, using the Landauer–Buttiker formalism. We use the scattering matrix to compute the total transmission probability $T^{ss'}(\theta_1) = |t|^2k_{x_3}/k_{x_1}$ for parallel and anti-parallel configurations by choosing the spins appropriately and for a particular incident angle $\theta_1$ (which fixes the angles in the other regions as well). The factor of $k_{x_3}/k_{x_1}$ in the transmission function is needed because the probability flux density includes a factor of the velocity which is essentially $\hbar k/m$. Since experimentally, it is not easy to control the angle of incidence of the impinging electron, we then compute the conductance by integrating over the possible angles of incidence and multiplying by the number of modes within the width $W$ of the silicene sample. At zero temperature, this leads us to a conductance given by

$$G^{ss'} = \frac{e^2}{h} W k_1 \left[ \frac{\theta_C}{\pi} \right] \int_0^{\theta_C} T^{ss'}(\theta_1) \cos \theta_1 d\theta_1. \quad (7)$$

Here, $\theta_C$ is the critical angle of the incident particles which is needed to ensure propagating particles in the first and third regions and is given by $\theta_C = \pi/2$ for $k_1 \leq k_3$ and $\theta_C = \arcsin(k_1/k_3)$ for $k_1 > k_3$. Note that unlike the case for graphene, in silicene, the contributions at the two valleys are not identical and hence, we do not get the degeneracy factor of two. Instead, the contributions at both the valleys have to be computed independently and added to obtain the total conductance through the junction. Thus we define the total charge conductances $G_c$, valley and the spin polarizations ($P_v, P_S$) and TMR through the FNF junction in terms of the following constituent conductances— $G^{ss'}_{K'K}$, $G^{ss'}_{K'K'}$. Here, $ss'$ denotes $uu$ (P) or $ud$ (AP) spin configurations, $V$ denotes the valley (K or K') and $\sigma$ denotes the spin of the incoming charge carrier in region 1 and $\sigma'$ denotes the spin of the outgoing charge carrier in region 3, which can be different, because we have spin-orbit coupling in the system. These conductances have been shown in Fig. 4 for both $E > h$ and $E < h$. It is now easy to realize, in reference to the band diagrams given in Fig. 3 that the conductances go to zero when there is a gap in the density of states either in region 1 or 3. The maxima can also be understood by noting that the density of states at those values of $E_z/h$ are maximum and reduce both when $E_z/h$ is reduced or increased. This can be checked for each of the various conductances on a case by case
The total charge conductance \( G_c^{s's'} \) and the valley \( P_v^{s's'} \) and spin \( P_s^{s's'} \) polarizations for both the P (\( s = s' \)) and AP (\( s = -s' \)) configurations are now defined as

\[
G_c^{s's'} = \sum_{V_{\sigma\sigma'}} G_{c,V_{\sigma\sigma'}}^{s's'},
\]
\[
P_v^{s's'} = \sum_{\sigma\sigma'} (G_{c,V_{\sigma\sigma'}}^{s's'} - G_{c,V_{\sigma\sigma'}}^{s's'}),
\]
and
\[
P_s^{s's'} = \sum_{\sigma\sigma'} (G_{c,V_{\sigma\sigma'}}^{s's'} - G_{c,V_{\sigma\sigma'}}^{s's'}). \tag{8}
\]

Note that the indices \( s's' \) in Eq. (8) gives rise to four possible spin configurations \( uu, ud, du, dd \) for a FNF junction, of which \( uu \) and \( dd \) imply P configurations with \( s = s' \) and \( ud \) and \( du \) denote AP configurations with \( s = -s' \).

Now from the knowledge of all the possible conductances, one can define the tunneling magnetoresistance (TMR) through the FNF geometry as

\[
\text{TMR} = \frac{G_{c,uu} - G_{c,ud}}{G_{c,uu}}. \tag{9}
\]

Note that the standard definition of TMR has \( G_{c,ud} \) in the denominator. However it is also sometimes defined with \( G_{c,uu} \) and we choose this definition because in our case, \( G_{c,ud} \) vanishes at \( E_z = h \). This implies a singularity in the TMR which is avoided in our definition. Note that for \( E_z > h \), the difference in TMR between the two definitions is negligible. For \( E_z < h \), there are numerical differences, but no qualitative difference in the behavior of TMR with the two definitions.

The results are different for the energy regimes \( E > h \) and \( E < h \), because of the difference in band structure, which has band gaps and hence no propagating states available for transport (see Fig. 3), for certain ranges of \( E_z/h \) for \( E < h \). Since the main difference of silicene from graphene is the fact that the gap in silicene is tunable by the external electric field \( E_z \), we choose to focus on the dependence of conductances on \( E_z \). In cases, where we study the conductances as functions of other parameters such as the barrier strength \( U \) or the exchange splitting \( \Delta \), we present our results for three different values of \( E_z/h \).

Note that when \( E_z = 3\Delta \), silicene is actually coplanar, i.e. the two sublattices are in same plane like in graphene. But the spin-orbit coupling in silicene is much stronger than in graphene\(^3\). This increases or decreases the momentum of the incident charge carrier (see Eq. (4)) depending on the spin-polarization of the ferromagnetic graphene. Hence, we do not expect to reproduce the results of FNF junctions in graphene in the gapless regime.

### A. Spin-independent Barrier

Here we discuss the case where we have a finite spin-independent (scalar) barrier in the normal silicene region. The energy of the incident electron can be in the regime \( E > h \) or \( E < h \) and we present the behavior of the conductances and the TMR below. To carry out our numerical analyses, we have chosen to normalise all our energy scales by the Zeeman energy \( h \), so that all our results are in terms of dimensionless quantities. We also choose to measure conductances in units of \( e^2 W/\pi h \).
We first present the results for the various constituent conductances for the P and AP configuration in Fig. 4 in order to understand the behaviour of the conductances, the valley and spin polarizations and the TMR in various parameter regimes. The conductances are shown independently at the \( K \) and \( K' \) valleys as well as independently for the incoming (\( \sigma \)) and the outgoing (\( \sigma' \)) spins of the charge carriers. In Figs. 4(a–h), we show the behaviour of the conductances at the \( K \) and \( K' \) valleys, with respect to the dimensionless parameter \( E_z/h \) for the four possibilities (\( \uparrow \uparrow, \uparrow \downarrow, \downarrow \uparrow, \downarrow \downarrow \)) in the \( E > h \) regime for both the \( uu \) and \( ud \) configurations. Here the \( uu \) configuration corresponds to the majority spin density of states in the left and right FS regions being up spin (parallel to each other) and the \( ud \) configuration corresponds to the majority spin being opposite (anti-parallel) in the two regions. When \( \uparrow \) charge carrier comes in from the left then it can either go to \( \uparrow \) state or \( \downarrow \) state in right region. So \( \uparrow \uparrow \) etc, denote spins of the incoming and scattered charge carriers. The behaviour of the various conductances in Figs. 4(a–h) can now be understood easily when analysed in terms of the band structure presented in Fig. 3.

For \( E > h \), the results have been presented for \( E/h = 4 \). Using the band diagram, it is easy to check that both at \( K \) and \( K' \), there are always electron states available for conductance for both the P and AP configurations and for all possible incoming and outgoing spins. The differences in the magnitude both at \( K \) and \( K' \) valleys stems from the decrease in the momentum of the electrons at the Fermi energy, as can be seen from the band diagram.

The results for \( E < h \) is shown in Figs. 4(i–p) for \( E/h = 0.5 \). Consider the \( uu \) case for the valleys \( K \) and \( K' \). For \( G_{K,K' \uparrow \uparrow} \), (red lines shown in Fig. 4(i) and Fig. 4(m)), the band diagram (see Fig. 3) shows that if we start with a spin up electron at the \( K \) valley, (Fig. 4(ii)) then there is a non-zero density of states for \( \uparrow \) electrons in the third region for all values of \( E_z/h \) until it reaches the value of \( 2 \) (the magenta line goes above the value of \( E_z/h = 0.5 \)). On the other hand, for the \( K' \) valley, (shown in Fig. 4(m)), there is no density of states for the \( \uparrow \) electrons beyond \( E_z/h = 1 \) (the blue line goes above the value \( E_z/h = 0.5 \)) in the third region. This explains why beyond \( E_z/h = 2 \) for the \( K \) valley and beyond \( E_z/h = 1 \) for the \( K' \) valley, the conductances \( G_{K,K' \uparrow \uparrow}^{uu} \) and \( G_{K,K' \uparrow \uparrow}^{uu} \) are zero. It is also clear from the band diagram that for \( G_{K,K' \uparrow \uparrow}^{uu} \), its value increases from the value at \( E_z/h = 0 \), because the momenta of the electrons at \( E/h = 0.5 \) grows (comparing the red and green lines) and beyond that it decreases (comparing the green, blue and magenta lines). On the other hand, for \( G_{K,K' \uparrow \uparrow}^{uu} \), it is clear that the momentum of the electrons at \( E_z/h = 0.5 \) decreases as a function of \( E_z/h \) (comparing the red, green and blue lines). This explains why the conductance rises initially and then falls beyond \( E_z/h = 0.5 \) for \( G_{K,K' \uparrow \uparrow}^{uu} \) and why it falls monotonically for \( G_{K,K' \uparrow \uparrow}^{uu} \).

A similar detailed analysis can also be made for the \( ud \) case as well as each of the other graphs in Figs. 4(j,k,l) and Figs. 4(n,o,p), which explains each feature of the graph. However, since the method is similar to what has been described above, we will not go through each one of the graphs in detail. The behaviour of the charge conductance, the valley and spin polarizations and the TMR are also now understandable, since we can explain how each of the constituents \( G_{V,\sigma \sigma'}^{ss} \) behave as a function of \( E_z/h \) from the band diagram.

FIG. 7. (Color online) In the left panels, we show the total charge conductance \( G_c \) in units of \( e^2W/\pi h \), valley and spin polarizations \( (P_{\sigma}^{\sigma'}, P_{\sigma}^{ss}) \) for \( E > h \) and on the right panels for \( E < h \) respectively. The dashed and solid lines correspond to two different exchange splittings (\( \Delta/h \)) due to the spin-dependent barrier in the normal silicene region. The value of the other parameters are chosen to be \( \lambda_{SO}/h = 0.5 \), \( U/h = 30 \). Energy of the incident electron, for \( E > h \), is \( E/h = 4.0 \) and for \( E < h \), \( E/h = 0.5 \). The small difference in magnitude of \( P_{\sigma}^{\sigma'} \) and \( P_{\sigma}^{ss} \) is highlighted in the insets for \( E > h \) regime.

(a) \( E > h \)

In Figs. 5(a), 5(c) and 5(e), we show the behaviour of the charge conductance \( (G_c^{ss}) \) and the valley and spin polarizations \( (P_{\sigma}^{ss} \) and \( P_{\sigma}^{ss}) \) with respect to the dimensionless parameter \( E_z/h \) for the P and AP configurations \((uu \) and \( ud \)) in the \( E > h \) regime. Note that in Fig. 5(a), \( G_c^{uu} \) and \( G_c^{ud} \) are both finite at \( E_z/h = 0 \) and start decreasing as we increase the value of \( E_z/h \). \( G_c \) is obtained by summing \( G_c^{ss} \) and \( G_c^{ss} \), which in turn are obtained as

\[
G_c^{\sigma \sigma'} = \sum_{\sigma''} G_{K+\sigma''}^{\sigma \sigma'} + G_{K-\sigma''}^{\sigma \sigma'}
\]

and similarly for \( G_c^{ss} \). In other words, the total charge conductance \( G_c^{ss} = G_c^{\uparrow \uparrow} + G_c^{\downarrow \downarrow} \) is obtained by summing over all the conductances in the panels (a) to (h) in Fig. 4.

In Fig. 5(c), \( P_{\sigma}^{uu} \) and \( P_{\sigma}^{ud} \) are plotted which are close to zero on the scale of the charge conductance. How-
However, they are not identical, as shown in the inset. But it appears that in this regime, silicene has negligible valley polarization, similar to graphene, which in fact has no valley polarization at all, since the two valleys are identical. This can be understood because the valley polarization $\mathcal{P}_v$ is simply proportional to $G_{uu}^s - G_{ud}^s$, and as can be seen from Fig. 4 that the magnitudes of the conductances at $K$ and $K'$ are almost the same for $E > h$. In Fig. 5(e), the behaviour of the spin polarization has been shown for both P ($uu$) and AP ($ud$) configurations, which is also very small in this regime. As shown in the inset, the spin polarization is positive for the P and negative for the AP configurations and increases as a function of $E_z/h$. This difference is due to the spin-orbit coupling in silicene, whereas in graphene, they are much smaller, since the spin-orbit coupling is vanishingly small. Finally, in Fig. 6(a), the behavior of the TMR is shown as a function of $E_z/h$ by the green solid line in the $E > h$ regime. In this regime, the TMR is close to zero.

(b) $E < h$

The right panels in Fig. 5 shows the behaviour of the charge conductance $G_{uu}^s$ and the valley and spin polarizations $\mathcal{P}_v^s$, $\mathcal{P}_v^s$ with respect to the dimensionless parameter $E_z/h$, in the $E < h$ regime, for the same spin and polarization configurations mentioned earlier. These are just the appropriate sums and differences of the constituent conductances in Figs. 4(i-p). Here also, their behaviour is easy to understand by comparing each of the graphs in Figs. 4(i-p) with the band diagrams in Fig. 3 and noting when there is no density of states for the configuration in either the incoming or the outgoing spin configuration of the charge carriers. For instance, for the $uu$ case, there is only one contribution for $E_z/h > 1$ in Fig. 4(i) and for the $ud$ case, there is no contribution for $E_z/h > 1$. This is because we have chosen $E/h = 0.5$ and the blue line ($E_z/h = 1.0$) in the band diagram goes above that line either for the incoming or scattered region for all cases in the $ud$ configuration and all but one case in the $uu$ configuration. In other words, their behaviour follows what is expected from the availability or non-availability of propagating states at the $K$ and $K'$ valleys as explained above in the discussion of Fig. 4.

The most interesting point to note is that the valley polarization and the spin polarization for the parallel or $uu$ configuration is unity for $E_z/h > 1$ in $E < h$ regime. This is simply because of the entire contribution to the conductance in this regime originates from $G_{uu}^s$. So the conductance is both fully valley and spin polarised and would be an important regime to achieve by tuning the incident electron energy $E/h < 1$ and the electric field $E_z/h > 1$. In the anti-parallel or $ud$ regime, the spin polarization can be tuned to negative values when $E_z/h < 1$, but without any valley polarization.

The behaviour of TMR is demonstrated in Figs. 6(a) and 6(b) for the $E < h$ regime. For $E_z/h = 0$, the TMR is negative and reaches its maximum negative value. Then the TMR increases as we increase the value of $E_z/h$ and changes sign and reaches saturation at $E_z/h \simeq 1$ since the conductance becomes fully spin polarised at that point. Note that we need to restrict the value of $E_z/h$ below two as the TMR takes an indeterminate form at $E_z/h = 2$ due to the vanishing of both $G_{uu}^s$ and $G_{ud}^s$ (see Fig. 5(c) and Eq. (9)) for all spin configurations. This is a consequence of our choice of the incident energy at $E/h = 0.5$.

The striking feature of positive to negative transition in the TMR also arises as we vary the strength of the potential barrier $U$ in the middle normal silicene region for different values of $E_z/h$. This feature is shown in Fig. 6(b) where the TMR oscillates between positive and negative values with respect to $U$ for $E_z/h = 0.1$. Such oscillations of the TMR from positive to negative values have been reported earlier in Ref. 21 in graphene, due to the change in the type of the charge carrier in the third region. Note also there is no significant qualitative change in the behavior of TMR even if we choose $U \sim h$. This extra tunability of the TMR with respect to an external electric field is a unique feature of silicene that we wish to emphasize here in this manuscript.

B. Spin dependent Barrier

Here we discuss the effect of a spin-dependent barrier on the total charge conductance, valley and spin polarizations and the TMR. The barrier is modelled in the normal silicene region as $U_c = U - \pi \Delta$, which is shown in Fig. 2. Here positive (negative) $\Delta$ represents the anti-parallel spin orientation of the FS in the first region.

In Figs. 7(a-f) we show the behaviour of the total charge conductances $G_{uu}^s$, spin and valley polarizations $\mathcal{P}_v^s$ and $\mathcal{P}_v^s$ in the $E > h$ and $E < h$ regimes, for $\pm \Delta/h$. Since the qualitative behavior of all the con-

FIG. 8. (Color online) The variation of the TMR with respect to $E_z/h$ and $U/h$ is shown for two values of $\Delta/h$. In the panels, the dashed and the solid lines correspond to $\Delta/h = 0.5$ and $\Delta/h = -0.5$ respectively. We choose the same value of the other parameters as mentioned in Fig. 7.
ductances remain similar to the spin independent barrier case, we do not show the behaviour of the conductances at the $K$ and $K'$ valleys independently, or analyse the graphs in detail via the band structure. Similarly, in Figs 8(a-c), we show the behaviour of the TMR as a function of $E_z/h$ and as a function of $U/h$ as well (for $E_z < h$), for both $\pm \Delta/h$. We find that the results are fairly similar to the spin-independent barrier case.

V. SUMMARY AND CONCLUSIONS

To summarize, in this paper, we have investigated the transport properties (charge conductance as well as spin and valley polarizations) and the TMR through a FNF junction in silicene. Here we have adopted the Landauer-Buttiker formalism to carry out our analysis. We show that the conductances and the TMR in this geometry can be tuned by an external electric field that the conductances and the TMR in this geometry can be tuned by an external electric field. We also find that the TMR can be tuned to $100\%$ in this geometry via the electric field. This is another conclusion of our analysis. Hence, from the application point of view, our FNF geometry may be a possible candidate for making future generation spintronic devices out of silicene.

As far as the practical realization of such a FNF structure in silicene is concerned, it should be possible to fabricate such a geometry with the currently available experimental techniques. Ferromagnetic exchange in silicene may be achieved via proximity effect using a magnetic insulator, for instance, EuO\(^{24,26}\). The typical spin orbit energy in silicene is $\lambda_{SO} \sim 4$ meV\(^3\). For an incident electron with energy $E \sim 4$ meV and exchange energy $h \sim 8$ meV, the maximum value of the spin and valley polarization as well as sign change of TMR from positive to negative value occur at an electric field $E_z \sim 0.03$ VÅ\(^{-1}\), potential barrier of height $U \sim 160$ meV, exchange splitting in the normal silicene region $\Delta \sim 4$ meV and width of the barrier $L \sim 100$ nm.

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