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Importance of non-local electron correlations in BaNiS₂ semimetal from quantum oscillations studies

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Abstract

By means of Shubnikov-de-Haas and de-Haas-van-Alphen oscillations, and *ab initio* calculations, we have studied the Fermi surface of high-quality BaNiS₂ single crystals, with mean free path $l \sim 400$ Å. The angle and temperature dependence of quantum oscillations indicates a quasi-two-dimensional Fermi surface, made of an electron-like tube centred at Γ , and of 4 hole-like cones, generated by Dirac bands, weakly dispersive in the out-of-plane direction. *Ab initio* electronic structure calculations, in the density functional theory framework, show that the inclusion of screened exchange is necessary to account for the experimental Fermi pockets. Therefore, the choice of the functional becomes crucial. A modified HSE hybrid functional with 7% of exact exchange outperforms both GGA and GGA+U density functionals, signalling the importance of non-local screened-exchange interactions in BaNiS₂, and, more generally, in 3d compensated semimetals.

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I. INTRODUCTION

After having been considered as a potential candidate for high-temperature superconductivity^{1,2}, which is however absent in the doping³ and pressure⁴ ranges explored, the BaNiS₂ compound has seen a renewed attention for its fascinating spin properties, discovered only very recently⁵. Indeed, angle resolved photoemission spectroscopy (ARPES) and *ab initio* electronic structure calculations based on the density functional theory (DFT) have shown that the spin-orbit (SO) interactions play an important role in this material^{5,6}. Despite the inversion symmetry of the crystal and the low Z number of the nickel atoms, an unexpected Rashba spin-orbit splitting was found in the bulk, with a Rashba spin-orbit coupling (SOC) as large as 26 eV Å. This value usually corresponds to much heavier elements, or to surface physics.

These surprising results can be related to the peculiar symmetry of crystal structure of BaNiS₂, reported in Fig. 1. The transition metal site is located at the center of edge-shared square pyramids, composed of sulfur atoms, whose orientation is staggered with respect to the basal plane. The BaNiS₂ structure is quasi-two-dimensional, with the barium atoms acting as a charge reservoir, intercalated between the electronically active layers, stacked along the c axis. The interlayer distance is approximately twice the metal-metal distance in the NiS plane. The presence of a large crystal field at the Ni sites in a locally asymmetric position enhances the SOC, and splits the electronic bands. Spin-orbit gaps and splittings of about 50 meV take place at the Fermi level, or in its proximity, paving the way to possible spintronic applications of this compound, which is a Pauli paramagnetic metal^{7,8}.

In view of exploiting the unique electronic properties of BaNiS₂ for future spintronic applications, it is important to have a detailed account of its band structure and Fermi surface topology. Following our previous ARPES experiment of Ref. 5, the Fermi surface can be further investigated by means of quantum oscillations (QOs). This complementary technique gives, by controlling the angle of the magnetic field, the topology of the Fermi sheets in the three-dimensional space, but without knowing their position in the first Brillouin Zone (FBZ). Except for the experimental investigations, the *ab initio* description of BaNiS₂ can also be improved. Previous DFT calculations of the band structure of BaNiS₂ did not consider SOC and the Hubbard U repulsion^{1,2}. Nevertheless, correlation effects in BaNiS₂ are non negligible, owing to the presence of a transition metal element at interme-

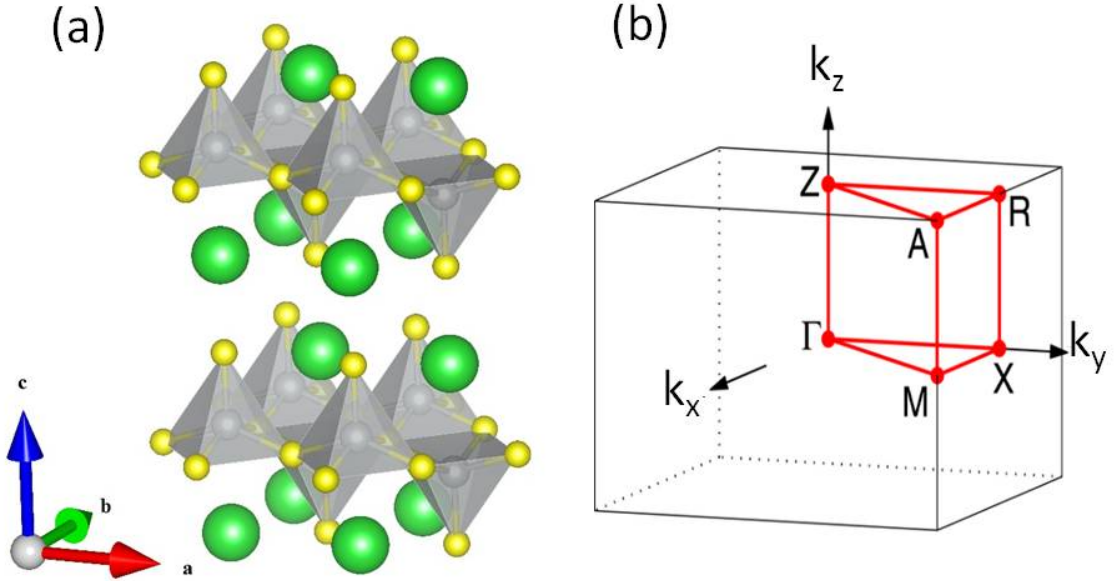


FIG. 1: (color online) (a) Crystallographic structure of BaNiS_2 . Large green and small yellow spheres stand for Ba and S, respectively. Ni is represented by medium size grey spheres at the center of the pyramids (b) First Brillouin zone of BaNiS_2 showing high-symmetry points.

diate filling. For instance, the interplay of filling, correlation, and charge transfer leads to a Fermi liquid breakdown in BaCoS_2 , where Ni has been substituted by cobalt⁹. Although our simulations within the generalized gradient approximation (GGA) supplemented by the Hubbard repulsion (GGA+U)⁵ significantly improves the description of the band structure of BaNiS_2 as compared to previous studies, discrepancies in the k_z dispersion have been found with ARPES results.

In this paper, we bring about an improved and highly accurate description of the band structure and Fermi surface of BaNiS_2 . For the first time, we performed Shubnikov-de Haas (SdH) and de Haas-van Alphen (dHvA) oscillation measurements on BaNi_2 , which are direct probes for the bulk Fermi surface. The k -space resolution in the three crystallographic directions is high enough to determine Fermi surface cross-sections with a resolution of 10^{-3} of the FBZ. The resolution in energy is also very high ($\Delta E \approx \mu_B B = 0.6 \text{ meV}$ for $B = 10 \text{ T}$) so that one can access to the effective mass very close to the Fermi energy. To reach these goals, we worked at low temperatures and strong magnetic fields. Moreover, we have been able to grow BaNiS_2 samples with a large enough scattering time⁸, such that they can be investigated by means of quantum oscillations (QOs).

For the above reasons, QOs are a stringent probe to benchmark the quality of *ab initio* electronic structure methods. The measured Fermi pockets volumes can be directly compared with the outcome of band structure calculations. The corresponding Fermi surface is sensitive to bandwidth renormalization, band shifts, and chemical potential. Therefore, it is a delicate quantity to compute from first principles, as it results from the fragile balance of many contributions, whose magnitude is set by the electronic correlations. On the other hand, *ab initio* calculations add invaluable information to the QOs, because they allow to locate the measured Fermi pockets in the Brillouin zone, as long as a clear match can be made between measured and computed pockets.

We performed DFT band structure calculations within the GGA, GGA+U and hybrid functional approximations. We show that the local Hubbard repulsion included in a mean-field fashion in the GGA+U framework is not adequate to account for the data. Instead, we reach a qualitative agreement with the experiment only when a fraction of the exact non-local exchange is added in the functional. By appropriately tuning the Heyd-Scuseria-Ernzerhof (HSE) hybrid functional^{10,11}, we obtain a *quantitative* agreement with the measured QOs. This corresponds to a 7% of exact exchange, while the screening length l_0 is kept equal to the original HSE06 functional ($\omega = 1/l_0 = 0.200 \text{ \AA}^{-1}$)^{11,12}. The very good match between the measured and computed frequencies of the QOs allows us to provide a detailed picture of the Fermi surface topology and electronic properties of this compound. Moreover, the comparison between our three different DFT functionals against the QOs measurements reveals the importance of non-local correlation effects in this class of transition metal systems. The semimetallic character of BaNiS₂ is responsible for an enhanced screening of the exchange operator. However, its non-local nature cannot be neglected, because its impact on the low-energy properties and on the fermiology is strong. Indeed, the optimal value of exact exchange fraction results from electronic screening mechanisms acting on the non-local exchange operator. The stronger the metallic character, the smaller the exact exchange weight in the functional. Our 7% value, smaller than in the original HSE functional, but still non negligible, is thus a consequence of the semimetallic character of the material.

Similar effects have been observed in the iron pnictides and selenides superconductors^{13–16}, which share with BaNiS₂ the presence of *d* electrons at intermediate filling, a compensated (semi)metallic character, and the *P4/nmm* point group in the undistorted phases. Band shifts non explained by local correlations have been reported in both theory¹⁷ and

experiments¹⁸. Therefore, the interplay between local and non-local interaction effects is key to explain the rich physics of this class of transition metals. The present work demonstrates that QOs measurements and DFT calculations with advanced functionals can be fruitfully combined to assess and quantify in an unambiguous way the importance of these contributions.

The paper is organized as follows. In Sec. II, we detail the experimental methods (Subsec. II A) and report the results (Subsec. II B) of the QOs measurements. In Sec. III, we describe the first-principles methods (Subsec. III A) and present the outcome of the electronic structure calculations (Subsec. III B). We draw our conclusions in Sec. IV.

II. QUANTUM OSCILLATIONS MEASUREMENTS

A. Methods

We have grown samples by a self-flux technique similar to that reported in Ref. 19, yielding high-quality platelet-like single crystals with dimensions of $\sim 1 \times 1 \times 0.1$ mm³. We selected those crystals with the best residual resistivity ratios, $\rho_{ab}(300K)/\rho_{ab}(10K) \approx 17$ for experiments in magnetic fields. We detected SdH oscillations by a four probe resistivity measurement using a lock-in detection in an Oxford Kelvinox 400 dilution refrigerator equipped with a 16-T superconducting magnet. We applied a current along the ab -plane and the magnetic field perpendicular to it. We swept the magnetic field at a rate of 10 G/s. dHvA oscillations were deduced from magnetic torque measurement in pulsed magnetic fields of up to 58 T, using a commercial piezoresistive microcantilever²⁰ at the LNCMI in Toulouse. We measured variations of the piezoresistance of the cantilever with a Wheatstone bridge with an ac excitation at a frequency of 63 kHz. We varied the angle θ between the normal to the ab -plane and the magnetic field by a one-axis rotating sample holder with a precision of $\pm 1^\circ$.

B. Results

Figures 2 (a) and (b) show the oscillating parts of the longitudinal magnetoresistance and of the torque for various temperatures. Fourier analysis of the oscillatory magnetoresistance (Fig. 2(c)) displays peaks at frequencies $F_\alpha = 37$ T, $F_\beta = 60.5$ T, $F_{\beta'} = 122$ T, $F_\gamma = 184$ T

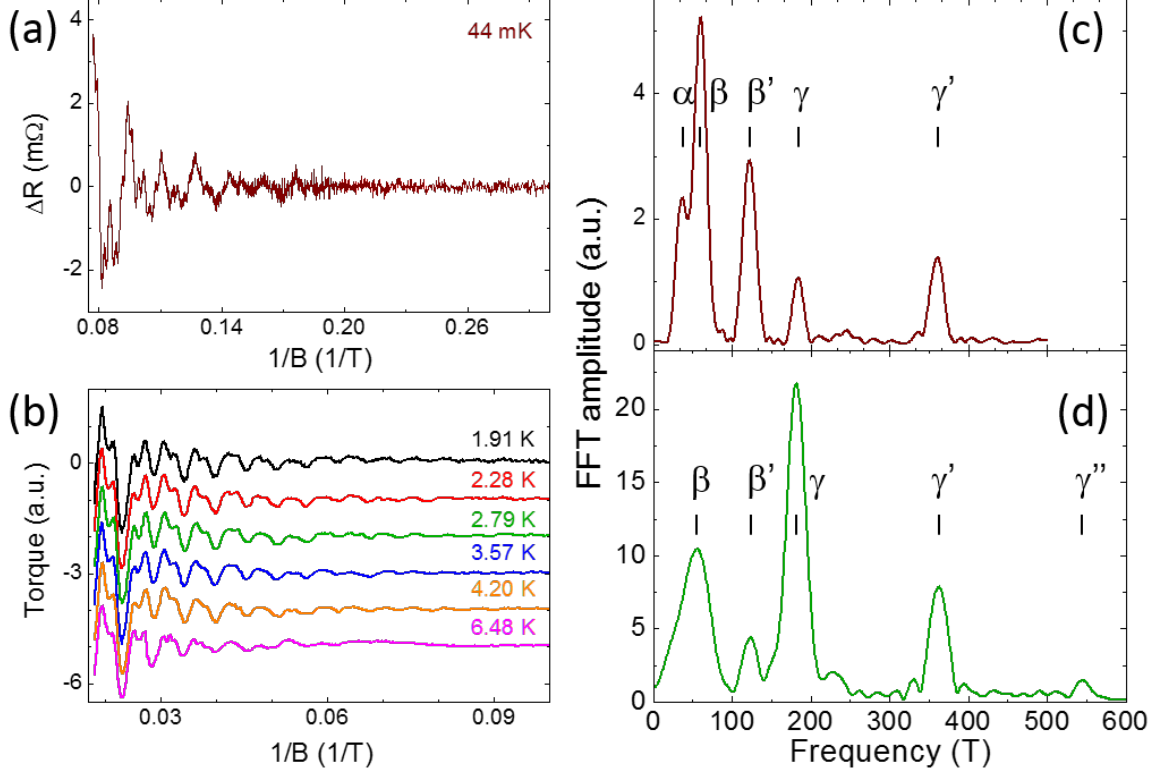


FIG. 2: (color online). (a) Oscillating part of the longitudinal magnetoresistance, ΔR measured at $T = 44$ mK. (b) Oscillating part of the torque, $\Delta\tau$ at various temperatures. θ was set to $17(1)^\circ$. (c) and (d) display the Fourier transforms of ΔR and $\Delta\tau$ ($T = 2.28$ K) taken in the range [5 - 13 T] and [10 - 55.2 T], respectively.

and $F_{\gamma'} = 361$ T. These data are in agreement with Fourier analysis of the oscillatory torque (Fig.2(d)) which displays peaks at frequencies $F_\beta = 60(6)$ T, $F_{\beta'} = 120(3)$ T, $F_\gamma = 182(1)$ T, $F_{\gamma'} = 363(3)$ T and $F_{\gamma''} = 549(8)$ T.

Within the reported uncertainties, it can be considered that the frequencies $F_{\beta'}$, F_γ , $F_{\gamma'}$ and $F_{\gamma''}$ are multiple of F_β . Hence, it cannot be excluded that these frequencies are harmonics of either F_β or F_γ . As discussed in the following this point can be checked through the analysis of the temperature dependence of the relevant Fourier amplitudes. Indeed, according to the Lifshitz-Kosevich (LK) formula, the amplitude of the first harmonic of a dHvA Fourier component with frequency F can be written ²¹:

$$\Delta M \propto R_T R_D R_S \sin\left(\frac{2\pi F}{B} + \phi\right) \quad (1)$$

where $R_T = \kappa T m^*/B \sinh[\kappa T m^*/B]$, $R_D = \exp(-\kappa T_D m^*/B)$ and R_S are the thermal,

Dingle and spin damping factors, respectively, with $\kappa = 2\pi^2 k_B m_0 / e\hbar \simeq 14.69$ T/K and $T_D = \hbar / 2\pi k_B \tau$ the Dingle temperature. m^* is the cyclotron effective mass in unit of the bare electron mass, m_0 . τ is the scattering time and ϕ is the phase. For the p^{th} harmonic, with frequency pF , the effective mass is pm^* . Temperature dependence of the Fourier amplitudes, an example of which is given in figure 3, are studied for three different directions of the magnetic field ($\theta = 17^\circ$, 34.5° , and 68°). The deduced effective masses are given in table I.

θ (°)	m_α^*	m_β^*	m_γ^*	$m_{\gamma'}^*$	$m_{\gamma''}^*$
17			0.39(12)	0.77(8)	
34.5	~ 0.4	~ 0.6	0.41(12)	0.87(11)	1.7(3)
68		0.77(13)	0.86(4)		

TABLE I: Effective masses $m^*(\theta)$ in bare electron mass units for few of the Fourier components observed in figure 2.

Within the error bars, these data are compatible with the relation $m_\gamma^* = m_{\gamma'}^* / 2 = m_{\gamma''}^* / 3$ whereas $m_\gamma^* \neq 3m_\beta^*$. It can therefore be supposed that three cyclotron orbits with frequencies F_α , F_β and F_γ enter the Fermi surface while $F_{\beta'}$, $F_{\gamma'}$, and $F_{\gamma''}$ are harmonics. Given the uncertainty on the amplitude of $F_{\beta'}$, it was not possible to determine its effective mass.

Dingle plots for the γ orbit at $\theta = 34.5^\circ$ are reported in figure 4 in which solid lines are best fits of Eq. 1 to the data with $T_D = 7$ K. Hence, the deduced scattering time is $\tau = 8.4 \times 10^{-14}$ s from which, by assuming a circular orbit with $v_F = \hbar k_F / m^*$, a mean free path, $l \sim 400$ Å can be deduced. The latter value is close to the longest mean free paths measured in high- T_C cuprates²² and iron-based pnictides²³.

The dimensionality of the Fermi surface can be studied by analyzing the angle dependence of the frequencies. A strictly 2D Fermi surface would result in a θ -independent value of $F \cos(\theta)$. As shown by figure 5, this is not the case for β and γ , since $F \cos(\theta)$ decreases as θ increases. Figure 5 also indicates that β and γ orbits belong to two different pockets because otherwise the two curves would cross at a given Yamaji angle²⁴. One may also wonder if the α orbit belongs to the same Fermi surface as the β or γ orbits. In fact, the data in Table I show similar effective masses for the three orbits. As a consequence, it is

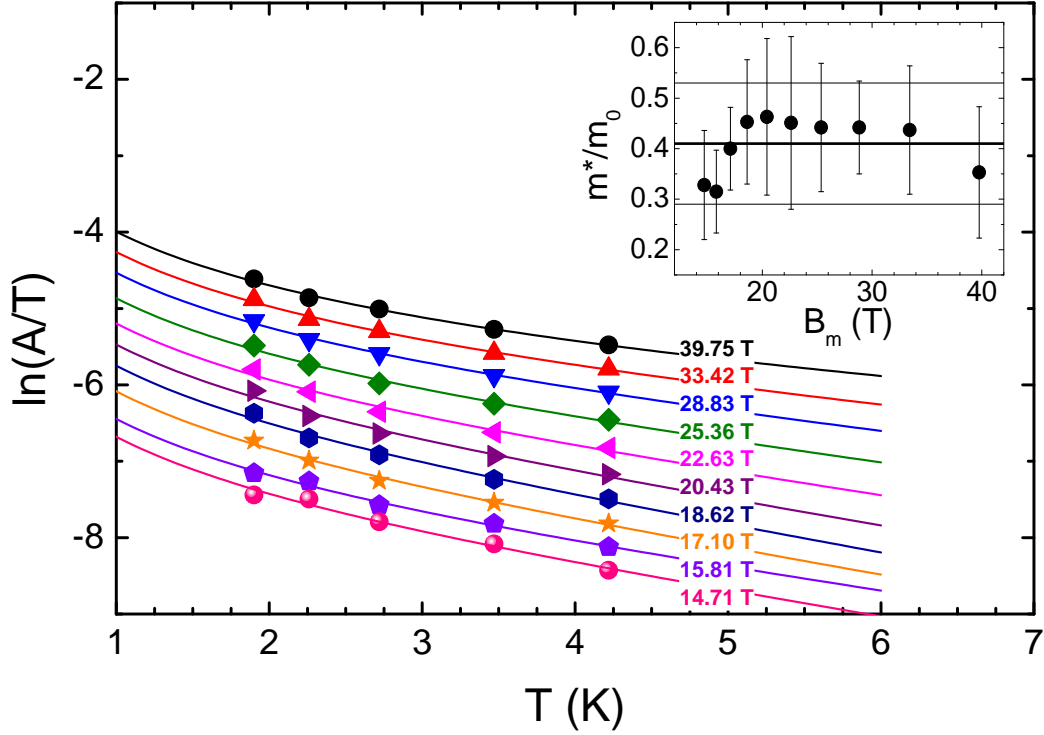


FIG. 3: (color online). Temperature dependence of $\ln(A/T)$ where A is the Fourier amplitude of the component with the frequency F_γ determined at different mean magnetic fields, B_m for $\theta = 34.5^\circ$. Solid lines are the best fits of Eq. 1 to the data. Inset shows the normalized effective mass, m^*/m_0 issued from these fits. The mean value of m^*/m_0 is shown by the thick straight line.

possible that two of these orbits are extreme orbits of the same sheet. This point will be discussed later based on the angle dependence of the frequencies and on band calculations. It can be noticed that the angle dependence of F_β and F_γ is nicely accounted for by the expression :

$$F(\theta) \cos(\theta) = \frac{F(0)}{\sqrt{1 + \xi^2 \tan^2(\theta)}} \quad (2)$$

obtained for Fermi surface pockets with a bi-axial ellipsoidal shape derived from standard parabolic bands, where $\xi = k_F^{xy}/k_F^z$ is the ratio of the major to minor semi-axes. Here $k_F^z > k_F^{xy}$ because $F \cos(\theta)$ decreases with increasing θ . Assuming a circular cross section in the $k_x k_y$ - plane, k_F^{xy} is related to $F(0)$ through the Onsager relation : $k_F^{xy} = \sqrt{\frac{2e}{\hbar} F(0)}$.

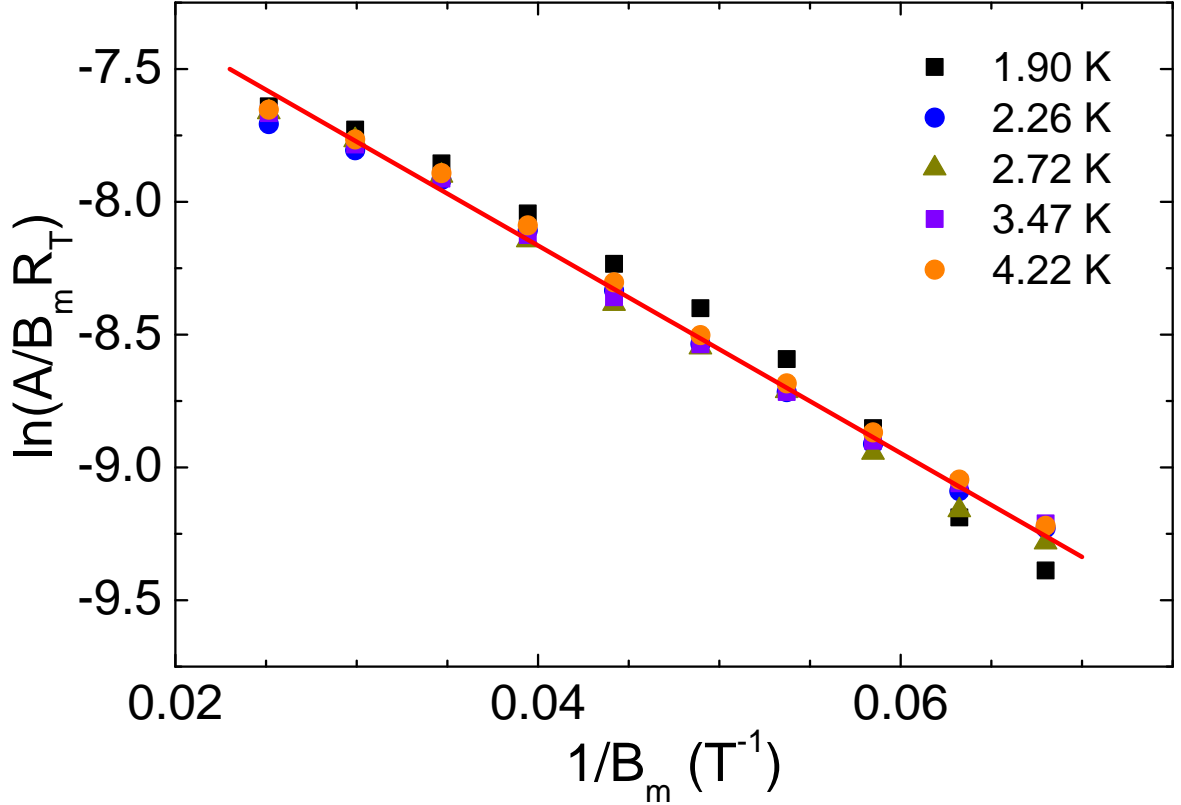


FIG. 4: (color online). Dingle plot for the frequency F_γ , at $\theta = 34.5^\circ$. The solid line is the best fit of Eq. 1 with $m^*(34.5^\circ) = 0.41m_0$ and $T_D = 6.3$ K.

At small enough θ values, in particular for $\theta = 17^\circ$ and 34° , both frequencies and effective masses are reasonably well approximated by a 2D Fermi surface, i.e. $F(0) = F(\theta) \cos(\theta)$ and $m^*(0) = m^*(\theta) \cos(\theta)$. While for $\theta = 68^\circ$ it is necessary to consider the ellipsoidal shape:

$$m^*(\theta) \cos(\theta) = \frac{m^*(0)}{\sqrt{1 + \xi^2 \tan^2(\theta)}} \quad (3)$$

introducing ξ values found in Eq. 2. $F(0)$, deduced cross section areas, $m^*(0)$ and ξ values are given in table II. All the orbits have very small area which is in agreement with the low carrier densities of BaNiS₂ given by band structure calculations¹ and magnetotransport measurements⁸. ξ values of 0.24(2) and 0.179(4) correspond to k_F^z/c^* ratios of 0.51(6) and 1.15(4) for the β and γ pockets, respectively. Therefore the β pocket can be modeled by an elongated ellipsoid. On the other hand, $k_F^z/c^* > 1$ obtained for γ is not consistent with an ellipsoid. It probably means that the ellipsoid model describes the dispersion only up to $\theta \approx 80^\circ$ but not more. Instead, it can be assumed that γ is the belly orbit of a two-

dimensional warped tube running along c^* . Within this picture, α could be the neck of this warped tube.

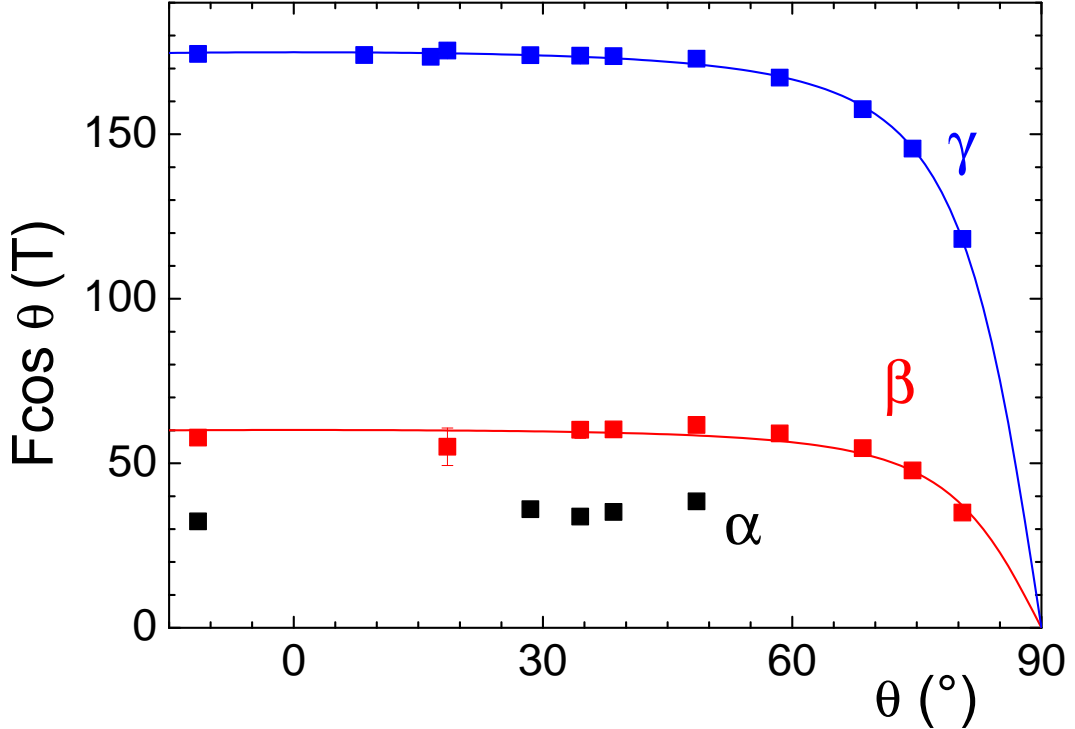


FIG. 5: (color online). Observed dHvA frequencies plotted as $F \cos(\theta)$ vs θ . Solid lines are best fit of Eq. 2 to the data relevant to β and γ orbits.

Orbit	$F(0)$ (T)	% of FBZ	ξ	$m^*(0)$
α	36(7)	0.17(4)		~ 0.3
β	63(3)	0.30(2)	0.24(2)	0.34(6)
γ	176(1)	0.839(5)	0.179(4)	0.35(8)

TABLE II: Parameters relevant to the three detected orbits projected in the $k_x k_y$ -plane. The cross section areas are given in percentage of the FBZ area. The ratio of the major to minor semiaxes (ξ) are deduced from the data in figure 5 assuming ellipsoidal Fermi surfaces (see Eq. 2). Effective masses are in bare electron mass units.

In summary, the quantum oscillations data of BaNiS₂ are consistent with a Fermi surface composed of two pockets. The larger one, named S_1 is a constricted tube with neck and

belly orbits α and γ , respectively. The angle dependence of the corresponding frequencies is in agreement with a sinusoidal dispersion along k_z , which at small enough angle between the field direction and the normal to the conducting plane can be approximated by an ellipsoid. A better resolution of F_α and a more detailed exploration of quantum oscillations as a function of θ would be necessary to investigate the existence of a Yamaji angle. The total volume of S_1 assuming such an energy dispersion is $V_1 = 6.4(6) \times 10^{-3} \text{ \AA}^{-3}$. The smaller pocket, denoted as S_2 is an elongated ellipsoid which is responsible of a single frequency F_β in quantum oscillations and which covers half of the FBZ in the k_z direction for a total volume $V_2 = 1.5(4) \times 10^{-3} \text{ \AA}^{-3}$.

In order to establish a model of Fermi surface, we cross-check this analysis with previously published results of photoemission spectroscopy and magnetotransport measurements^{5,8}. According to the Hall effect data⁸, BaNiS₂ is a semimetal very close to compensation, with a charge imbalance of only 1% in favor of electrons. For the compensation to be fulfilled, S_1 and S_2 must have different electronic characters and, since $V_1 \approx 4V_2$ and owing to the tetragonal symmetry of BaNiS₂, the former must be located at the center of the FBZ while the latter has to be repeated four times. Within this picture, the total number of holes and electrons is of the order of $\sim 5 \times 10^{19} \text{ cm}^{-3}$, in agreement with magnetotransport data. These statements are also in agreement with ARPES measurements⁵ which, for a given k_z value, reveal two pockets: an electron-like one at the center of the FBZ and a hole-like one along Γ -M at mid-distance between the two high-symmetry end points.

III. *AB INITIO* ELECTRONIC STRUCTURE CALCULATIONS

A. Methods

In order to identify the Fermi pockets detected by QOs, we are going to use three different density functionals, namely the GGA with the Perdew-Burke-Ernzerhof (PBE) functional²⁵, the PBE supplemented by local Hubbard interactions (GGA+U)^{26,27}, and a modified hybrid Heyd-Scuseria-Ernzerhof (HSE) functional^{10,11} with an optimized 7% of exact exchange. Nowadays, the hybrid functionals, introduced for molecular systems¹², are getting more and more popular in solid state physics as well²⁸⁻³⁰, because they correct for some deficiencies of the regular density functionals, thanks to a reduced self-interaction error, and the restoration

of the derivative discontinuity. The hybrid functionals are non-local and orbital dependent. In all three cases, we included nickel spin-orbit (SO) interactions, whose importance has been recently demonstrated in this system^{5,6}. Indeed, despite the quite low Z number of the nickel atoms, the peculiar crystal structure of BaNiS₂ enhances the SOC, with sizable splittings showing up in certain regions of the Brillouin zone, and for some bands crossing the Fermi level⁵. Therefore, we expect that the shape of some Fermi pockets will be affected by SOC.

All *ab initio* DFT calculations have been carried out by using the Quantum ESPRESSO package^{31,32}. The geometry of the cell and the internal coordinates are taken from the experimental values reported previously³³. Ni, Ba and S atoms are described by norm-conserving pseudopotentials. The Ni pseudopotential is fully relativistic, with 10 valence electrons ($4s^2 3d^8$) and non-linear core corrections. For Ba, the semi-core states have been explicitly included in the calculations. The S pseudopotential is constructed with the $3s^2 3p^4$ in-valence configuration. The k -point sampling convergence is reached for a $8 \times 8 \times 8$ electron-momentum grid and a Methfessel-Paxton smearing of 0.01 Ry. The plane-waves cutoff has been set to 120 Ry for the wave function in the GGA and GGA+U calculations. In the HSE calculations, we lowered this cutoff to 60 Ry in order to reduce the computational burden of the non-local exchange potential evaluation. We checked that this lower cutoff does not affect the results within a target accuracy of 0.01 eV in the band structure.

For the GGA+U calculations performed in the fully rotational invariant framework²⁷, we took the same Hubbard parameters as in Ref. 5, namely $U=3$ eV, $J=0.95$ eV, and with the F4/F2 Slater integral ratio equal to its atomic value. This is in agreement with experimental values³⁴ and recent cRPA estimates⁹.

For the HSE functional, we used the latest fast implementation of the exact Fock exchange energy³², based on the adaptively compressed exchange (ACE) scheme³⁵. The significant CPU time reduction, allowed by the ACE implementation, makes hybrid functional calculations of BaNiS₂ possible in a plane-waves framework, which guarantees a systematic and controlled basis-set convergence. The q -integration of the non-local Fock operator has been performed in a downsampled grid made of $8 \times 8 \times 2$ q -points. To minimize the number of non-equivalent momenta in the $k + q$ grid, we found that shifting the k -grid by half-a-grid step in the z direction further speeds up our calculation. The loss of precision caused by the q -downsampling is critical for metals. However, given the quasi-two-dimensional nature of

BaNiS₂, reducing the q grid by a factor of 4 in the out-of-plane direction does not deteriorate the results, which still fulfill our target accuracy, while leading to a significant gain in CPU time.

As a final step, we obtained an accurate determination of the band structure, chemical potential, Fermi surface, and Fermi pocket cross-sections by the Wannier interpolation of the *ab initio* bands, performed with the Wannier90 code³⁶ for all functionals taken into account. The tight-binding model comprises 22 bands, originating from all Ni d and S p states in the BaNiS₂ unit cell.

B. Results

The band structures are reported in Fig. 6, while the corresponding Fermi surfaces are plotted in Fig. 7. Four cones are present in the BZ, whose axes are located at the proximity of the $(-1/4, -1/4, z)$, $(-1/4, 1/4, z)$, $(1/4, -1/4, z)$, and $(1/4, 1/4, z)$ directions. At $k_z = 0$, the Fermi level crosses the cones at their apex. The conical energy dispersion drifts as a function of k_z , being lifted up when we move away from $k_z = 0$. The dispersion along k_z is caused by the necessity of charge-carrier compensation. This leads to 4 Fermi surfaces that are also conical, instead of being Dirac lines, with vertices lying on the $k_z = 0$ plane. These four conical surfaces are a very robust feature of the system, present in *all* DFT calculations performed here. Therefore, they are insensitive to the correlation level, and to the quality of the density functional used. Nevertheless, their related QOs signal, evaluated from the area of the extremal cross-section at $k_z = \pi/c$, varies from functional to functional, as reported in Fig. 8. However, their F fluctuation is much milder than for any other Fermi pocket, and it is in a quite good agreement with the experiment for all functionals. Thus, from the *ab initio* values, we can safely attribute the β orbit to these 4 cones. The linear dispersion close to the cone apex is not captured by our experimental results, because it would require the angle dependence of both frequencies and effective mass, as stated in Ref. 37. However, it is reassuring that our conclusions drawn in Sec. IIB regarding the carrier concentration remain valid within this picture since, according to the DFT calculations, the volume of the conic part is small. The highly mobile holes observed in magnetotransport measurements are assigned to this conic portion.

In contrast to the robustness of the conical Fermi surfaces, the most fragile Fermi structure

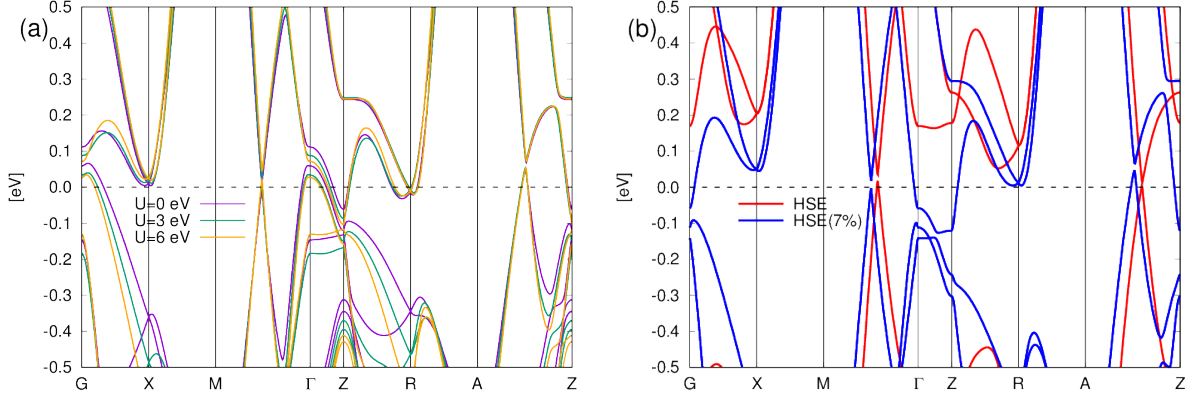


FIG. 6: (color online). Panel (a): Band structure obtained with the GGA and GGA+U functionals. The U values are reported in the key. The main difference between the functionals is in the band dispersion along the Γ -Z path. The larger the U , the lower the position of the bands that cross the Fermi level. Panel (b): Band structure obtained with the HSE functional with two fractions of exact exchange. The regular 25% of exchange is used in “HSE”, while only the 7% is used in “HSE(7%)”. Again, the main difference is in the Γ -Z path.

is the one along the Γ -Z direction. As one can evince from Fig. 6(a), the position of the valence and conduction bands, that cross the Fermi level in Γ -Z, strongly depends on the on-site repulsion strength included in the GGA+U functional. These bands are shifted down by about 50 meV by the addition of a Hubbard term $U = 3$ eV. This shift increases upon increasing the value of U . As a result, the hole pocket centered at Γ shrinks while the electron pocket centered at Z grows (see Figs. 7(a) and 7(b)). Anyway, the corresponding F values reported in Fig. 8 for the hole pockets at Γ are far too large if compared with the experimental findings. Neither GGA nor GGA+U agree with experimental data. Instead, the use of properly tuned hybrid functionals corrects for this pitfall. While regular HSE opens a fictitious gap in the Γ -Z region, by reducing the exact exchange weight down to 7% the crossing bands acquire an electron character all the way from Γ to Z (Fig. 6(b)). This is reflected in the Fermi surface by the disappearance of the hole pocket at Γ , as seen in Fig. 7(c). In the modified HSE (“HSE(7%)”), the electron pocket becomes a 2D constricted tube that stretches throughout the whole BZ along k_z . The QOs generated by this structure at $\theta = 0$ are given by the extremal cross-sections at Γ (minimum) and Z (maximum). Their values are plotted as histograms in Fig. 8. They nicely correspond to the experimental values of the α and γ orbits, respectively. Therefore, we can identify these two orbits as belonging

to the same Γ -constricted tube, and having Γ -Z as rotational axis. This maps all measured dHvA frequencies into identified pockets, based on a Fermi surface picture provided by our modified HSE.

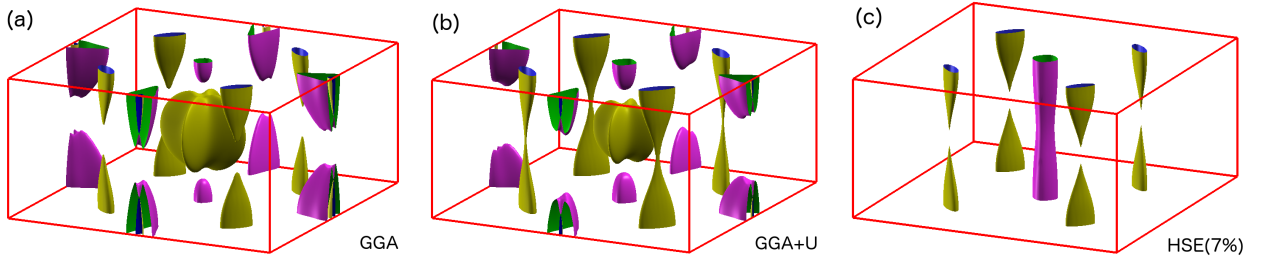


FIG. 7: (color online). Fermi surfaces of BaNiS_2 from *ab initio* calculations with GGA (panel (a)), GGA+U(=3eV) (panel (b)), and modified HSE with 7% of exact exchange (panel (c)). The red lines are the frontiers of the first BZ, drawn in Fig. 1(b) together with the high-symmetry points. The yellow (purple) surfaces indicate Fermi pockets of hole (electron) character.

It is worth noting that in the GGA and GGA+U band structure, there is yet another electron pocket, centered at R (Fig. 6(a)), which is Rashba active⁵. The Rashba splitting gives rise to two nested parabolas, filled by spinors with opposite chirality. One can see the nested shapes in Figs. 7(a) and 7(b). This should yield additional frequencies F , the strongest being as large as $300 T$. This value is not found in the experiment. The improved HSE band structure does not have any pocket at R , because the conduction bands are raised just above the Fermi level, without touching it. This is another feature that highlights the reliability of the modified HSE functional. At variance, the GGA and GGA+U functionals show several drawbacks, and they are unable to reproduce neither quantitatively nor qualitatively the experimental frequency values. This tells us about the importance of non-local screened-exchange interactions in this material, as we will argue in the concluding Sec. IV.

This theoretical analysis allows us to unambiguously assign the S_1 surface to the electron-type quasi-2D tube having Γ -Z as axis with neck and belly centered at Γ (α orbit) and Z (γ orbit), respectively. On the other hand, the 4 S_2 hole-type pockets, corresponding to β , are assigned to 4 conical Fermi surfaces, whose vertical axes are approximately $(-1/4, -1/4, z)$, $(-1/4, 1/4, z)$, $(1/4, -1/4, z)$, and $(1/4, 1/4, z)$, with extremal cross-sections located at $k_z = \pi/c$, and apex at $k_z = 0$.

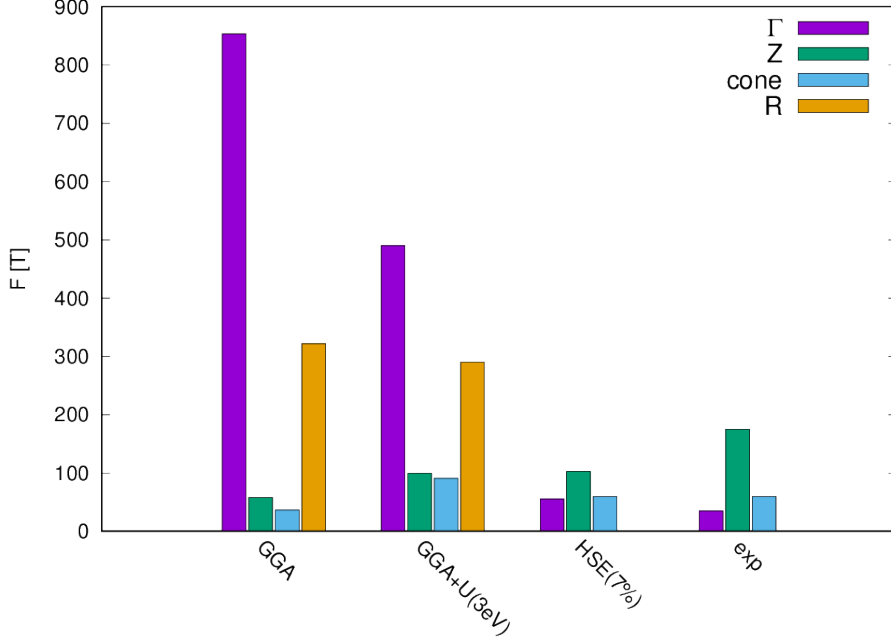


FIG. 8: (color online). Computed QOs frequencies for $\theta = 0$ by employing the same density functionals that give the Fermi surfaces plotted in Fig. 7. Those frequencies correspond to the extremal cross-sections containing the k -points reported in the caption. For the modified HSE with 7% of exact exchange (“HSE(7%)”), there is no pocket around R, at variance with the GGA and GGA+U cases.

IV. CONCLUSIONS

Shubnikov-de Haas and de Haas-van Alphen oscillations are reported for the first time in the undoped BaNiS₂ compensated semimetal. The estimated mean free path of 400 Å indicates a low concentration of defects in the measured single crystals. We detected three independent main frequencies corresponding to small cyclotron orbits, covering less than 1% of the FBZ cross-section area. Angle dependences of these frequencies suggest the existence of two types of pockets, which is confirmed by DFT electronic structure calculations when a modified HSE hybrid functional with 7% of exact exchange is used.

The first type (S_1) is a constricted tube of electron-like character and having the Γ -Z direction as rotational symmetry axis. By symmetry there is just one pocket of this type. The remaining four hole-like pockets belong to the second type (S_2). They are oriented in the z direction, and centered in the proximity of one of the $(-1/4, -1/4, z)$, $(-1/4, 1/4, z)$, $(1/4, -1/4, z)$, and $(1/4, 1/4, z)$ axes. For this second type, the measured

frequency indicates a conventional ellipsoidal shape due to parabolic bands. According to band structure calculations the unprobed part of these pockets has a conic-like dispersion and should result in high-mobility carriers. This could be verified experimentally by the determination of the angle dependence of the effective mass.

At first sight, the present results may be in contradiction with our model of the magnetotransport data including three types of charge carriers⁸: minority holes with a high mobility and majority holes and electrons. However, the latter model is based on the Drude law of parabolic bands, which supposes homogeneous effective masses and scattering times of charge carriers, and is not adapted for linear dispersions characteristic of the present hole pockets. It turns out that the recourse to two discrete and homogeneous populations of charge carriers is sufficient to simulate the properties of the conic-like and ellipsoid-like parts of the S_2 pockets, hosting hole carriers with different mobilities, very high for the conic-like part, lower in the ellipsoid sector.

By carrying out *ab initio* electronic structure calculations, we showed that the use of a modified hybrid functional is necessary to explain the experimental values. This modified functional has a reduced ratio of exact exchange, if compared with the standard HSE functional. We found that 7% of the exact exchange yields a good quantitative agreement with the QOs results. The agreement is lost, even qualitatively, in the case we use a regular GGA functional. Adding the local Hubbard repulsion to the GGA in the GGA+U framework does not improve the GGA picture. This indicates the importance of non-local screened-exchange interactions in the system. Therefore, in the BaNiS₂ semimetal, the main role of the electronic correlation is non-local, rather than Hubbard-like.

The role played by the non-local screened exchange can be relevant to other transition metal compounds, such as the iron selenide. Like BaNiS₂, FeSe is a multiband compensated semimetal, although its correlation strength is known to be much larger than the one in BaNiS₂^{13,15,16,38}. Nevertheless, there are features that are not explained by a local self energy only¹⁸. The present study suggests that the non-local screened-exchange interactions cannot be neglected in 3d semimetals, particularly when the screening is poor, as a consequence of *d*-electron correlation and charge-carrier compensation. To model these systems, advanced *ab initio* methods are necessary, which include non-local correlations, such as hybrid functionals or many-body methods with non-local self energy, such as GW, GW+DMFT, and their variants^{39,40}. Combining QOs measurements with various types of first-principles cal-

culations turns out to be very important to identify the main renormaliation mechanisms of the low-energy electronic structure.

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