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DOI
10.1103/PhysRevMaterials.3.114202

Publication date
2019

Document Version
Final published version

Published in
Physical Review Materials

Citation for published version (APA):
Scattering rate collapse driven by a van Hove singularity in the Dirac semimetal PdTe$_2$

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(Received 14 June 2019; revised manuscript received 25 August 2019; published 18 November 2019)

We present optical measurements of the transition metal dichalcogenide PdTe$_2$. The reflectivity displays an unusual temperature and energy dependence in the far infrared, which we show can only be explained by a collapse of the scattering rate at low temperature, resulting from the vicinity of a van Hove singularity near the Fermi energy. An analysis of the optical conductivity suggests that below 150 K a reduction in the available phase space for scattering takes place, resulting in long-lived quasiparticle excitations. We suggest that this reduction in phase space provides experimental evidence for a van Hove singularity close to the Fermi level. Our data furthermore indicates a very weak electron-phonon coupling. Combined this suggests that the superconducting transition temperature is set by the density of states associated with the van Hove singularity.

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I. INTRODUCTION

The investigation of topological phases of matter has reignited interest and resulted in new insights in the electronic behavior of various relatively simple materials families. One class of materials that is of particular interest for applications is the transition metal dichalcogenides (TMDs). These materials host various forms of topological matter often in combination with a form of spontaneous symmetry breaking. As such, the TMDs offer exciting opportunities to study topological phase transitions, something that was previously only really possible in low temperature quantum Hall systems. One such example is PdTe$_2$, which features both bulk Dirac points [1–3] as well as topological surface states [4] in combination with a transition to a superconducting phase with $T_c = 1.4–2$ K [5–8]. The superconducting state is conventional with an isotropic superconducting gap [8–11]. One peculiarity observed in these experiments is that PdTe$_2$ supports a type-I superconducting state [8,12], but recently some surface effects have been observed that leave open the possibility for an interesting surface superconducting state [13,14]. Superconductivity in PdTe$_2$ is undisputedly of a conventional, electron-phonon coupling origin. Recently, it was pointed out that there is a van Hove singularity close to the Fermi level in PdTe$_2$ [15] and that this could have important consequences for the behavior of the superconducting phase when PdTe$_2$ is intercalated [16] or put under pressure [14].

In this paper we investigate the normal state properties of PdTe$_2$ using optical spectroscopic methods. The most prominent feature in the optical spectra is an optical gap that becomes visible in the low temperature, far-infrared reflectivity with an edge around 37 meV. From an analysis of the optical spectra we identify several features that are consistent with interband transitions between previously observed bulk type-II Dirac states. We use the extended Drude model to investigate the temperature dependent response of the low energy charge carriers and find that as temperature decreases, the optical scattering rate continuously decreases and collapses below our detection limit at temperatures under 60 K. We speculate that this arises from the complete occupation or depletion of a van Hove singularity that sits in close proximity to the Fermi energy. The same analysis demonstrates that the electron-phonon interaction in this material is likely very weak and thus we suggest that the superconducting transition is set by the large density of states associated with the van Hove singularity.

II. METHODS

Single crystals of PdTe$_2$ have been grown using a modified Bridgman method [17] from high quality starting materials. Starting from 99.99% pure Pd and 99.99% pure Te, single crystals were grown in a quartz ampoule at 800 °C. The superconducting characteristics of these crystals, as determined by various experiments, were previously reported in Refs. [8,11,12,14]. Extensive characterization details are reported in Ref. [8] (see also the Supplemental Material Ref. [8]). Energy dispersive x-ray analysis using a scanning electron microscope demonstrates the expected stoichiometry and a homogeneous distribution of the Pd and Te content. Resistivity data shows a residual resistivity ratio, RRR, equal to 30 with a low value of the room temperature resistivity of 23 $\mu$Ωcm. This points to the high quality of these crystals and the very low impurity content. The superconducting transition takes place at $T_c = 1.57$ K. No signs of degradation have been reported for these crystals and oxidation is not expected to play a role. Crystals used in our study were cleaved minutes before sealing the UHV cryostat and, as explained next, all experiments have been repeated several times to confirm reproducibility. These measurements did not show any signs of degradation over the period of the experiments. Reflectivity data was measured in the energy range between 5 meV and 4.35 eV using a Vertex v80 Fourier transform spectrometer. The temperature dependence of the reflectivity was measured using a home build cryostat with a
the RefFit software package developed by A. B. Kuzmenko. The resulting reflectivity data was analyzed using reflectivity spectra are thus obtained between 10 and 300 K in temperature dependent spectrum. We subsequently average all these temperature cycles we obtain a dense sampling of the far-infrared and UV ranges of the spectrum. By combining position stabilized sample holder [18]. Spectra were collected while continuously cooling to base temperature (10 K) and warming up to room temperature. The rate at which we change temperature is 1.3 K/min and we record 1 spectrum every minute. Temperature cycles were repeated up to five times, depending on the frequency range, to ensure reproducibility and improve signal-to-noise ratios, in particular in the far-infrared and UV ranges of the spectrum. By combining these temperature cycles we obtain a dense sampling of the temperature dependent spectrum. We subsequently average all spectra within a 2 K interval. After the temperature cycles on the sample, in situ evaporation of a thin reference layer was used to determine the absolute reflectivity. The temperature cycles were subsequently repeated on this reference layer. The reflectivity spectra are thus obtained between 10 and 300 K in 2 K steps. The resulting reflectivity data was analyzed using the RefFit software package developed by A. B. Kuzmenko [19,20].

III. RESULTS

In Fig. 1 the reflectivity for a few selected temperatures is shown. The far-infrared reflectivity of PdTe$_2$ displays characteristic signatures of a good metal: At room temperature the reflectivity follows a Hagen-Rubens behavior from which we extract a DC resistivity of approximately 14 $\mu$Ωcm, which is a little bit lower than what is observed in transport experiments [14,17,21]. As temperature decreases several features appear in the reflectivity. The most prominent of these is a steplike feature at 37 meV is indicated by S1, while a second feature 17 meV is indicated by S2. Both features become more prominent with decreasing temperature. The inset shows the reflectivity over the full range.

FIG. 1. Reflectivity for selected temperatures. A steplike feature that develops at approximately 37 meV is indicated by S1, while a second feature 17 meV is indicated by S2. Both features become more prominent with decreasing temperature. The inset shows the reflectivity over the full range.

From the measured reflectivity one can obtain the optical conductivity using two methods: Kramers-Kronig transformation or through modeling. Although both methods should work equally well, modeling provides physically correct and flexible extrapolations outside the measured experimental range. We use a minimal Drude-Lorentz model with six oscillators at all temperatures. The model consists of a single Drude term, followed by a Lorentz oscillator with a central frequency that varies from 150 cm$^{-1}$ (at 10 K) to 400 cm$^{-1}$ (300 K) (i.e., from 18–49 meV). Next we need two midinfrared oscillators, which are more or less temperature independent and have central frequencies of 3600 and 4300 cm$^{-1}$ (0.45 and 0.53 eV). The visible part of the spectrum can be modelled with a strong optical transition at 1.4 eV and a weaker one around 4 eV. The real part of the dielectric function extrapolates to a value $\varepsilon_{\infty} = 3.5$.

Based on the Drude-Lorentz modeling we apply a variational dielectric function [20] that effectively carries out the Kramers-Kronig transformation. The resulting optical conductivity is displayed in Fig. 2. As expected from the Drude-Lorentz model, we observe a midinfrared band centered around 0.5 eV and a strong transition around 1.5 eV. The midinfrared band has a significant fine structure: We observe transitions at 0.27 eV, 0.42 eV, 0.52 eV, 0.65 eV, and 0.83 eV. To determine the origin of these transitions, we turn to band-structure calculations [1,15,21,23]. These indicate that the bands near the Fermi level consist predominantly of Te 5p orbitals. In particular, Ref. [1] shows that along the $\Gamma$–$A$ direction a band crossing takes place, resulting in a type-II bulk Dirac point at binding energy around $-0.65$ eV. At the same time surface states are observed with binding energies of $-1.1$ eV and $-1.7$ eV. To determine whether the optical transitions observed here arise from these bands, the optical selection rules need to be evaluated. Although the bands around the Fermi level are all derived from $p$ orbitals, the character of the hybridized bands is altered as a
gap around 0.03 eV at low temperature. The optical transitions features prominently in the visible range (1.5 eV). The midinfrared band centered around 0.5 eV. A second interband transition serves a midinfrared band centered around 0.5 eV. A second interband transition features prominently in the visible range (1.5 eV). The optical transitions features prominently in the visible range (1.5 eV). The dispersion. We note that in a recent calculation [15] it was observed that there is a van Hove singularity (vHS) at an energy just below the Fermi level. In such a case, it is completely occupied at low temperature and a partial depletion of the vHS leads to a strong increase of the scattering rate.

We investigate the enhancement of the scattering rate using the extended Drude model [24]. Figure 3 displays both the real and imaginary part of the optical scattering rate, where the latter is commonly referred to as the mass enhancement factor. Figure 3(a) shows the frequency dependent optical scattering rate for a few selected temperatures. The inset of panel 3(b) displays the temperature dependence of the optical scattering rate at 25 meV. Starting at low temperature, we observe a frequency independent scattering rate (below 30 meV), almost below our detection limit, 1/τ(0, T = 10 K) ≈ 1 ± 0.5 meV. Furthermore, the scattering rate is temperature independent up to 50 K as the inset in panel 3(b) shows. This suggests that the scattering at low temperature and energy is completely dominated by impurity scattering. As temperature increases, the scattering rate acquires a frequency dependence that slowly turns linear over an extended energy range as clearly seen in the data above 200 K. The temperature dependence of the optical scattering rate [inset Fig. 3(b)] shows a marked increase as temperature increases above 50 K. This increase in the scattering rate occurs when temperature becomes comparable to the energy of the lowest mode frequencies in the Eliashberg function α²F(ω) [15] and we therefore suggest that at this temperature electron-phonon scattering becomes relevant. At even higher temperatures (∼150 K), a strong temperature dependence sets in and the scattering rate increases more or less linearly. The temperature where this takes place is indeed around the temperature where we expect the vHS to become partially occupied and the phase space for scattering to be significantly enhanced, thus confirming the first of the expected optical signatures of the presence of the vHS. The second feature, a redistribution of spectral weight, is also observed in our data. This is clearly visible from Fig. 2 where at low temperature spectral weight is piling up below the energy scale set by the vHS. At low temperature the integrated spectral weight shows a plateau at low energies (40 meV) before further increasing as interband transitions start to contribute. The integrated spectral weight

![Graph](image_url)
The observation of a frequency and temperature independent optical scattering rate at low temperature leaves an interesting open question about the appearance of superconductivity below 1.5 K [5,6,8]. Since the mechanism of superconductivity in PdTe$_2$ is expected to be electron-phonon coupling driven, one would expect a signature of this interaction in spectroscopic experiments. In the superconducting state, such signatures are often visible as 'peak-dip-hump' structures, most famously observed in Giaever's first tunneling experiments [25]. Similar features have also been observed in the far-infrared spectra of superconductors [26] and these are often taken as important clues to the mechanism underlying the superconducting state. Our observations of a featureless and temperature independent scattering rate suggest that the electron-phonon coupling is very weak below 50 K. Using the standard BCS result for $T_c = 1.14\hbar\omega_0\exp(-1/\Delta_0)$, our experimental data indeed imply an important role for the vHS in driving the superconducting instability as suggested in Ref. [15]. The highest phonon frequency to be expected in PdTe$_2$ is of order 22 meV [15]. One would expect some change in the frequency dependence of the scattering rate around this energy. Even if the dominant electron-phonon coupling would be below our experimental energy range, we should detect a corresponding frequency dependence of the optical scattering rate in our experimental window. A final important indication for the absence of significant electron-phonon interaction is given by the mass enhancement factor, displayed in Fig. 3(b). There appears to be some mass enhancement at room temperature, but as the temperature decreases the mass enhancement decreases as well and reaches a minimum of approximately $m^*/m \approx 1.1$ at $T = 10$ K.

It is interesting to consider the role of the vHS in other dichalcogenide systems, such as PtTe$_2$ and PtSe$_2$. Band-structure calculations have been performed for PtTe$_2$ in Refs. [21] and [15]. Both calculations indeed show that the vHS appears at higher energy, e.g., further away from the Fermi level, which would result in a reduction of $N(0)$. If there is an important role for the vHS in determining $T_c$, PtTe$_2$ should have a lower $T_c$ and is indeed not superconducting. Nonsuperconducting PtSe$_2$ has been studied in Ref. [27], where a vHS appears around 0.5 eV above the Fermi level. These authors find that upon electron doping, their calculations predict a similar $T_c$ as for PdTe$_2$ ($T_c = 2.15$ K). The authors indicate that the superconducting transition is driven by a softening of the acoustic phonon branch, but we note that for the case where the doping is changed by 0.5 electron/unit cell, the vHS comes precipitously close to the Fermi level.

To summarize, we have presented optical spectra of the superconductor PdTe$_2$. The optical data features several mid-infrared and visible range optical transitions, some of which may be related to optical transitions within the bulk Dirac cone. In the far infrared we observe a sharp plasma edgelike feature at low temperature. A strong reduction of the optical scattering rate and a reshuffling of spectral weight provide evidence that this edge is the optical signature of a van Hove singularity some 30 meV above the Fermi level. This observation combined with a frequency independent optical scattering rate provides experimental support for a superconducting transition driven by a large density of states associated with a van Hove singularity.
ACKNOWLEDGMENTS

We would like to thank A. de Visser for useful comments during the preparation of this paper. This work is part of the research programme Strange Metals (Grant No. 16METL01) of the former Foundation for Fundamental Research on Matter (FOM), which is financially supported by the Netherlands Organisation for Scientific Research (NWO).