Scanning tunneling measurements of layers of superconducting 2H-TaSe2: Evidence for a zero-bias anomaly in single layers

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(Received 27 September 2012; revised manuscript received 4 February 2013; published 4 March 2013)

We report a characterization of surfaces of the dichalcogenide TaSe2 using scanning tunneling microscopy and spectroscopy at 150 mK. When the top layer has the 2H structure and the layer immediately below the 1T structure, we find a singular spatial dependence of the tunneling conductance below 1 K, changing from a zero-bias peak on top of Se atoms to a gap in between Se atoms. The zero-bias peak is additionally modulated by the commensurate $3a_0 \times 3a_0$ charge-density wave of 2H-TaSe2. Multilayers of 2H-TaSe2 show a spatially homogeneous superconducting gap with a critical temperature also of 1 K. We discuss possible origins for the peculiar tunneling conductance in single layers.

DOI: 10.1103/PhysRevB.87.094502 PACS number(s): 74.55.+v, 73.63.Bd, 74.45.+c, 74.70.Xa

I. INTRODUCTION

There is a rather general interest arising in transition-metal dichalcogenides because superconductivity, charge-density wave (CDW), quantum criticality, and single-layer physics are found in just a few compounds.1–7 These systems share wave (CDW), quantum criticality, and single-layer physics.

Subsequent analysis showed instead that these patterns are due to the atomic lattice.17,19 The 2H polytype has a unit cell consisting of two TaSe2 units, each built up by two Se triangles separated by Ta atoms. Interlayer bonding is weak through van der Waals forces. The relative arrangement of the Se triangles varies in different polytypes of the same compound. In Fig. 1 we show the structure of the the 2H and the 1T polytypes.

The 2H polytype has a high in-plane resistivity $\Delta T_c \approx 13 K$ and shows a $a$-axis interlayer coupling decreases from $2H-NbSe_2$, $2H-TaS_2$, to $2H-TaSe_2$. Superconductivity in $2H-NbSe_2$ ($T_c = 7 K$) is multiband and $s$-wave, with a strong in-plane anisotropy related to the charge-density wave.1,2,9,10 $2H-TaS_2$ ($T_c = 0.8 K$) shows anisotropic charge-density-wave patterns at the surface. These anisotropic patterns were first proposed to be indicative of a chiral charge-density wave.1,12

Subsequent analysis showed instead that these patterns are due to the atomic lattice.17,19 The relationship between superconductivity and such in-plane anisotropic charge order is yet unclear.

There are few reports about superconductivity in $2H-TaSe_2$, showing superconducting diamagnetic signals and zero resistance at temperatures below 150 mK.8,14,15 Zero-temperature extrapolated critical fields are very low (1.4 mT with the field along the $c$ axis). On the other hand, calculations and angular-resolved photoemission have unveiled the Fermi surface and band structure in bulk.

We measure TaSe$_2$ in a home-built STM arrangement installed in a dilution refrigerator, which is capable of cooling down the microscope to about 100 mK. We can make tunneling conductance curves with a resolution of 20 $\mu$V, which corresponds to an energy resolution above 150 mK. The STM has a sample holder that allows us to change...
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in situ the scanning window. We use a tip of Au, which we clean through repeated indentation on an Au cleaning pad. The samples are 2$H$-TaSe$_2$ crystals grown using iodine vapor transport from stoichiometric prereacted powders. Bulk susceptibility measurements of our samples are shown in the inset of Fig. 3 for two magnetic fields. Susceptibility gives a diamagnetic signal with the same $T_{c}$ as in previous work (150 mK). The temperature dependence of the resistance normalized at ambient temperature is shown in the main panel of Fig. 3 and was measured between 300 and 0.5 K. CDW onset causes a change in the resistance near $T_{CDW} = 122$ K. The residual resistivity ratio of the TaSe$_2$ samples is of 26, implying samples similar to those used to observe quantum oscillations and measure the Fermi surface. We glued the 2$H$-TaSe$_2$ sample using silver epoxy onto the sample holder. We used scotch to remove the upper layers of 2$H$-TaSe$_2$ at ambient conditions, and we optically inspected the result prior to inserting the setup into liquid helium. The scotch cleaving procedure works nicely, but it is more involved than cleaving 2$H$-NbSe$_2$. In 2$H$-NbSe$_2$, it is easy to obtain a clean-looking shiny surface without free-standing sheets. But in our samples of 2$H$-TaSe$_2$, the cleaved surface nearly always consists of big loose sheets which have to be manually removed using tweezers, until we observe a flat and shiny surface. We cooled down the samples 11 times, each time making a new cleave. In five cool downs, we observed a pure 1$T$-TaSe$_2$ surface with the $\sqrt{13} \times \sqrt{13}$ CDW, without any traces of superconductivity. In two cool downs, we found 2$H$-TaSe$_2$.

FIG. 1. (Color online) We show two possible polytypes of TaSe$_2$, the 2$H$ and 1$T$ phases, which consist of two TaSe$_2$ hexagonal prisms rotated with respect to each other, and one trigonal prism. Se atoms are shown as gray spheres and Ta atoms as smaller red spheres. The coordination of Ta is trigonal prismatic in the 2$H$ structure and octahedral in the 1$T$ structure.

FIG. 2. (Color online) We show a schematic representation of scanning tunneling microscopy (STM) images of the atomic arrangements discussed. We highlight the top sheet as 2$H$. It shows a $3a_0 \times 3a_0$ superlattice modulated charge-density wave (CDW). The layer immediately below, highlighted as 1$T$, has a different Moiré pattern with a $\sqrt{13}a_0 \times \sqrt{13}a_0$ CDW. In 2$H$-TaSe$_2$, top and bottom Se atoms are aligned [top figure in (c)]. In 1$T$-TaSe$_2$, the two Se triangles are rotated [bottom figure in (c)]. The middle panels (b) show in more detail the CDW patterns found in each area. There are three inequivalent atomic sites in the pattern, as highlighted by yellow, orange, and white points. The incommensurate CDW $\sqrt{13}a_0 \times \sqrt{13}a_0$ pattern shown in the bottom panel of (b) consists of stars of David, with also three inequivalent sites. The arrangement shown in (a), with one 2$H$-TaSe$_2$ sheet on top of 1$T$-TaSe$_2$, presents the singular superconducting features discussed in the text.
behavior over the whole surface without any trace of superconductivity. This is what we expect for bulk $2H\text{-TaSe}_2$ with the above-mentioned energy resolution, above $T_c$ of 150 mK. In four cool downs, we observed the phenomena described below with mixed hexagonal ($2H$) and trigonal ($1T$) surfaces. In each of them, we studied about ten different 2 $\mu$m × 2 $\mu$m scanning windows, finding the behavior discussed below.

Typically, we scan with a set point at bias voltages of about 2 mV and a tunneling conductance of 1 $\mu$S or below. We did not find a significant dependence of the tunneling curves or the images on the set-point tunneling conductance within an order of magnitude above or below this value. $I-V$ curves are numerically derived, as in previous work.\textsuperscript{10,21} We normalize to the conductance value at 1 mV, and curves are flat between 1 and 2 mV. We did not systematically study curves above this bias voltage value. Images have been made using homemade data acquisition software, and rendered using homemade, WSXM\textsuperscript{25} and MATLAB\textsuperscript{R} programs.

III. RESULTS

The STM topography shows the atomic Se lattice and either the 3$a_0 \times 3a_0$ CDW with the same orientation to the Se lattice and commensurate to it, or Moiré patterns characteristic of a CDW with a periodicity of $\sqrt{3}a_0 \times \sqrt{3}a_0$ rotated by 13.5° with respect to the atomic Se lattice. The first identifies $2H\text{-TaSe}_2$ surfaces, and the second $1T\text{-TaSe}_2$ surfaces (Fig. 2). We find atomically flat $2H\text{-TaSe}_2$ layers, showing immediately below a surface of $1T\text{-TaSe}_2$. Steps between both surfaces are of 1.2 nm, corresponding to two layers of the sandwich Se-Ta-Se. Thus, the upper layers are single-layer crystals of $2H\text{-TaSe}_2$. We also find $1T\text{-TaSe}_2$ or $2H\text{-TaSe}_2$ surfaces over large areas and many different steps. These correspond to multilayer $1T\text{-TaSe}_2$ and $2H\text{-TaSe}_2$, respectively. We have made a thorough spectroscopy and microscopy characterization of all observed surfaces by taking simultaneously topography and tunneling conductance (Figs. 4 and 5). In Fig. 4, we show curves obtained when scanning from the top of a Se atom to an intersite at the center of three Se atoms. On surfaces of $2H\text{-TaSe}_2$ on top of $1T\text{-TaSe}_2$, a clear zero-bias peak is observed on top of the Se atoms [Figs. 4(a) and 5(a)]. The zero-bias peak evolves continuously into a V-shaped conductance at the intersites, following the topography. This is made clear in the Fourier transform of the zero-bias conductance map, which shows Bragg peaks at the positions corresponding to atomic and CDW reciprocal vectors.

In the $1T\text{-TaSe}_2$ layers [Figs. 4(b) and 5(b)], we find featureless flat tunneling conductance curves. When we find $2H\text{-TaSe}_2$ [Figs. 4(c) and 5(c)] over different layers, clear superconducting tunneling features are observed with quasiparticle peaks located somewhat below 200 $\mu$V. These features do not change as a function of the position over the surface, as shown by the zero-bias conductance maps and their Fourier transform in Fig. 5(c).

In Fig. 6, we show representative examples of single layers of $2H\text{-TaSe}_2$ imaged over larger areas (top panels), together with four zoom-ups of topography (middle panels) and zero-bias conductance maps (bottom panels) at the $2H$ single layers and at the $1T$ underlayer. Steps between single-layer

FIG. 4. (Color online) Tunneling conductance vs bias voltage in single-layer $2H\text{-TaSe}_2$ on top of $1T\text{-TaSe}_2$ (a), as in Fig. 2, in $1T\text{-TaSe}_2$ (b) and in multilayer $2H\text{-TaSe}_2$ (c) taken at 0.15 K. The black line on top of the yellow curve in c is a fit to the superconducting gap as described in Sec. IV. Different tunneling conductance curves are taken when the tip moves from a Se atom (black) to an intersite (yellow). Insets show the path on atomic size topography images involving a hexagon of Se atoms. Following other paths of the sixfold symmetry leads to qualitatively the same features.

$2H\text{-TaSe}_2$ correspond to the size of one unit cell of $2H\text{-TaSe}_2$, around 1.2 nm. Typically, the lateral size of the single-layer sheets is between 50 and 300 nm, and the boundary with the underlying $1T\text{-TaSe}_2$ layer is sharp.

More detailed real-space imaging of the $2H\text{-TaSe}_2$ single layer is shown in Fig. 7. The atomic and CDW modulations are observed simultaneously on the zero-bias conductance map and the topography. In particular, the Fourier transform image of the zero-bias conductance map shows the atomic Se lattice and the CDW. In real space, we observe that the highest value for the zero-bias quasiparticle peak [blue curve in Fig. 7(b)] coincides with the brightest Se atom due to the CDW [white points in Fig. 7(a)]. On the other hand, the V-shaped dip [yellow curves in Figs. 4(a) and 5(a) and in Fig. 7(b)]
between Se atoms is homogeneous. Thus, there is a very strong modulation of the zero-bias peak at the Se atoms with the position related to the CDW order. The V-shape of the intersites shows no CDW modulation, and persists until bias voltages up to approximately 0.7 mV.

Accordingly, the conductance maps at bias voltages different from zero show a smooth evolution of the Moiré patterns presented in Fig. 5(a). The contrast related to the Se lattice and the CDW is shown in Fig. 8, where we plot the Fourier amplitude of the lattice and CDW peaks observed in the Fourier transform of the zero-bias conductance map in Fig. 5(a) for different bias voltages. The CDW peaks are maintained up to about 150 μV, where they start to decrease. Above 300 μV, all the atomic positions present similar conductance, and the only contrast between them and the intersites is observed in the conductance maps. This gives the sixfold modulation of the hexagonal lattice until roughly 0.7 mV, where all features in the conductance maps disappear. Thus, as shown in Fig. 7, the zero-bias conductance peak has a small energy scale of roughly 150 μV and is linked to the CDW, whereas the V-shaped dip at Se atoms is broad and survives up to higher bias voltages.

When we increase the temperature or the magnetic field, we observe that these features disappear from the tunneling conductance curves, which become flat above approximately 1 K (insets of Fig. 9). Above fields of 100 millitesla, the peak and dip disappear into flat tunneling conductance curves. The weak-coupling superconducting gap equation \( \Delta = 1.76k_B T_c \) gives, for \( T_c = 1 \) K, \( \Delta = 150 \) μeV, which coincides with the width of the zero bias peak in the single hexagonal layer (Fig. 7) and the size of the superconducting gap in multilayers [Fig. 4(c)].

Similar temperature dependences are observed for the superconducting gap measured in \( 2H-TaSe_2 \) multilayers (Fig. 9). Note that the superconducting gap observed in multilayers is strongly smeared, with a high number of states at zero bias, contrasting well developed gap structures found in the other \( 2H \) transition-metal dichalcogenides.

Magnetic fields of 10 mT, applied perpendicular to the sample, lead to flat tunneling conductance curves both in single layers and in multilayers. Such fields are at the lower limit of our coil and magnet power supply system, and thus we did not follow the temperature dependence of the upper critical field nor study spatial gap dependences under field.
FIG. 6. (Color online) Large topographic images in the top panels and zoom-ups of different areas in the middle panels. The bottom panels are conductance maps taken at zero bias in the same areas as the small size topographic images shown in the middle panels, showing similar color scales as in Fig. 5. We observe a 2H-TaSe₂ single layer on top of 1T-TaSe₂. The hexagonal layer shows the features in the conductance discussed in the top panels of Figs. 4 and 5, and the trigonal layer is featureless. Images here are unfiltered and have been taken at 0.15 K. The height profile marked by a blue line in the top right panel is shown in the top right inset.

IV. DISCUSSION

Our sample shows, in the bulk, clearly 2H-TaSe₂ features (Fig. 3). However, on the surface we can find 1T-TaSe₂ CDW on some regions. Thus, the surface properties can significantly change with respect to the bulk. Tunneling conductance maps are featureless in 1T-TaSe₂ surfaces, which is in itself not very surprising, and shows that this polytype has no noticeable physics at energies of a mV or below. On the other hand, 2H-TaSe₂ surfaces are found with a critical temperature of 1 K. This is at odds with the superconducting transition found in the bulk (0.15 K). The critical temperature of these materials is easily enhanced through pressure or strain.18,28–30 The increase of $T_c$ highlights surface strains or surface-induced slight modifications in the electron-phonon coupling in some areas. Such modifications are probably more difficult to observe in other dichalcogenides such as NbSe₂ or NbS₂ where the bulk $T_c$ is higher and closer to the maximum $T_c$ obtained in these materials under pressure, which lies around 9 K.30,31 In our TaSe₂ samples, we did not observe any feature in the resistivity around 1 K. This means that the layers showing the superconducting gap of Fig. 4 (bottom panel) represent a very small volume fraction of the sample. The strongly broadened BCS features observed in Fig. 4 (bottom panel) also show that superconductivity is not that of a typical bulk and clean BCS $s$-wave superconductor.

At present, there are no clear-cut data of in-plane and out-of-plane coherence lengths of the bulk 2H-TaSe₂ superconductor with $T_c = 0.15$ K. Values of $\xi_\parallel = 500$ nm (in-plane) and $\xi_\perp = 200$ nm (out-of-plane) have been obtained from the zero-temperature extrapolation of the out-of-plane and in-plane critical fields (using $H_{c,\perp} = \Phi_0/2\pi\mu_0\xi_\perp^2$ and $H_{c,\parallel} = \Phi_0/2\pi\mu_0\xi_\parallel^2$).15 However, resistivity measurements show an anisotropy of nearly three orders of magnitude (700), which is clearly at odds with the far smaller anisotropy found in the critical field measurements.32 Fermi surface measurements also point to strongly two-dimensional bands, so that the out-of-plane coherence length $\xi_\perp$ should be probably far below the value from critical field measurements.8,16 Regarding the in-plane coherence length $\xi_\parallel$, it is striking that the values given are much larger than
FIG. 7. (Color online) Topography (top panel) and tunneling conductance map at zero bias (bottom panel) of a hexagonal layer on top of a trigonal layer. Images, taken at 0.15 K, are filtered for clarity. The blue, dark, and light green curves of (b) are taken all three on top of Se atoms, corresponding to the color code shown in the bottom panel of (a). The yellow curve is taken in between Se atoms. The curves on top of the Se atoms all show a zero-bias conductance peak, whose height is modulated as a function of the position. Blue curves are located at the Se position, which also shows the highest peak, whose height is modulated as a function of the position. The blue, dark, and light green curves of (b) are taken all three on top of a trigonal layer. Images, taken at 0.15 K, are filtered for clarity. The blue, dark, and light green curves of (b) are taken all three on top of Se atoms, corresponding to the color code shown in the bottom panel of (a). Dark green and light green curves are located at the other Se positions with different charge modulations. Yellow curve is independent of the position in the CDW modulation.

FIG. 8. The Fourier amplitude of the first three Bragg peaks in single layers of 2H-TaSe$_2$ on top of 1T-TaSe$_2$ along the direction of the sixfold modulation of highest contrast in the tunneling conductance maps. Note the use of log-scale to highlight the noise background. The central Bragg peak has been removed. The two peaks at the lower reciprocal lengths give the CDW pattern, and the third peak is due to the atomic Se lattice. When increasing the bias voltage, the sixfold modulation remains for the CDW peaks until it disappears above 300 μV. The largest part of this modulation comes from the local variations in the zero-bias peak amplitude. The modulation at distances of the atomic lattice (third Bragg peak) remains until 600 μV. This is due to the modulation of the V-shaped dip at intersites.

those reported in 2H-NbSe$_2$, of 10 nm. With such high values, the search for a vortex lattice requires very large scanning ranges, above the size of the flat 2H-TaSe$_2$ areas observed in our experiment. Using the value we find here for the superconducting gap Δ, we can make a simple estimation of the superconducting coherence length of multilayers of 2H-TaSe$_2$ with $T_c$ at 1 K and find $\xi \parallel \approx \hbar v_F / 2 \Delta$ and $v_F = 4.8 \times 10^7$ m/s from Ref. 33) $\xi \parallel \approx 500$ nm, of similar order to that of the values discussed in the bulk.8,14,15

The zero-bias conductance peak found on single layers of 2H-TaSe$_2$ on top of 1T-TaSe$_2$ is probably our most intriguing result. We can compare (Fig. 9) the temperature evolution of the zero-bias conductance peak in single layers with the evolution of the superconducting gap in multilayers. We can fit the tunneling conductance curves in multilayer 2H-TaSe$_2$ to $s$-wave BCS theory using a gap of $\Delta = 150$ μeV and a broadening lifetime parameter$^{31}$ of $\Gamma = 55$ μeV (black line shown in the bottom panel of Fig. 4). The temperature evolution of the gap parameter $\Delta(T)$ is shown by the open black points in Fig. 9. $\Delta(T)$ is below the nearly parabolic dependence expected within BCS theory, and it clearly disappears at 1 K. On the other hand, the zero-bias conductance peak in single-layer 2H-TaSe$_2$ follows well a Gaussian shape with a width of $\sigma = 0.1$ mV. When increasing temperature, the width of the peak increases above temperature-induced broadening. The width of the peak in single layers inversely scales with the decrease of the gap in multilayers. This shows that the origin of the zero-bias conductance peak in single layers is related to superconductivity. It also shows that the destruction of superconducting correlations by temperature roughly follows $\Delta(T)$ in single layers, and the zero-bias peak is correspondingly broadened by the temperature-induced gap decrease.

To discuss the atomic spatial dependence of the zero-bias conductance peak, it is useful to mention the known features of the Fermi surface of 2H-TaSe$_2$ (Fig. 10; see Ref. 16). At room temperature, it shows holelike sections centered at the $\Gamma$ and $K$ points, and electronlike dogbone-formed sheets around the $M$ point, coming from two different bands derived from Ta d electrons.16,35 One is for the $\Gamma$- and $K$-centered sheets, which has a saddle point in between, and the other is for the $M$-centered sheet. Angular-resolved photoemission shows that small pseudogaps open on the $K$ pocket at high temperatures, precluding CDW order. At low temperatures, below the CDW transition, the Brillouin zone becomes three times smaller, and real band gaps appear. The $\Gamma$ pocket remains intact, and the $K$ pocket is destroyed, due to nesting features in the $\Gamma$ and $M$ sheets, with significant softening of $\Sigma_1$ phonons.28 The $M$ dogbone sheet also suffers great changes. The low-temperature Fermi surface consists of circular pockets at the $\Gamma$ point of the new Brillouin zone and rounded triangles at the new $K$ points. The $M$ dogbone breaks up into parts. One is located around the $\Gamma$ and the other one around the $K$ point, both of which
are new due to the $3 \times 3$ CDW state. The strong spin-orbit coupling of the Ta $5d$ levels has a pronounced influence on the Fermi surface topology and could be at the origin of the gap on the dogbone sheets. There exists one part of the Fermi surface which remains untouched by the CDW order, and is strongly two-dimensional, namely the big circle around $\Gamma$. In single-layer $2H$-TaSe$_2$, the CDW has the same structure as in the bulk, and thus the Fermi surface, at least the part involved in charge order, is likely to have the same features, too.

Atomic-sized changes in the superconducting tunneling conductance have been often observed previously in systems with an anisotropic gap structure, such as $2H$-NbSe$_2$ or the high-$T_c$ cuprates. They result from the anisotropic interaction between the tip and the sample. The tunneling conductance probes the density of states of different parts of the band structure. The variations in the height of the zero-bias peak with the CDW [Fig. 7(b)] show the involvement of charge order in shaping the zero voltage anomaly. Thus, the peak is related to the bands where the CDW forms, i.e., the rounded triangles and the pockets stemming from the dogbone sheet.

Isolated single-layer crystals of dichalcogenide materials have been obtained previously on different substrates through repeated exfoliation. In spite of extensive searches, no clear experimental evidence of superconductivity has been found in them. On the other hand, in situ grown single-layer surfaces, mostly of Pb, are creating a rich playground, demonstrating that superconductivity can form in atomically thin crystals. In all cases, the substrate and the related interface plays a fundamental role, for instance the covalent bonding to a Si(111) substrate in submonolayers of Pb, or the interaction with the SrTiO$_3$ substrate in FeSe. In any case, the superconducting critical temperature $T_c$ decreases when achieving ultimate thickness, and the superconducting tunneling conductance shows $s$-wave BCS like gap features. Theory proposes significant electron-electron interactions in doped graphene layers or in single-layer MoS$_2$, which give $d$-wave or superconductivity changing sign in different Fermi surface sheets.

Thus, both the observed strong critical temperature increase close to the surface and the atomic size variation of the tunneling conductance between a zero-bias peak and a V-shaped dip are reported here. We can speculate as to the various possible origins of this behavior.

Regarding the zero-bias conductance peak, zero-energy resonances have been found previously in a number of systems, possibly receiving the most attention in superconductors and in low-dimensional structures with charging effects. Often, the zero-bias peaks demonstrate the presence of a flat band.

In superconductors, resonances close to the Fermi level are found, e.g., at the core of magnetic vortices because of multiple Andreev reflection. The first such state is located very close to zero energy (at $\Delta^2 / E_F$, which is generally small) and has been seen in the STM experiments as a zero-bias peak. At zero magnetic field, in-gap bound states also arise at magnetic impurities in $s$-wave superconductors. Their energy location depends on the relative amplitude of electron and hole impurity wave functions, which is governed by the exchange interaction between the localized magnetic moments of the impurity and the Cooper pairs, or by the scattering phase shift. When scattering is resonant, the impurity bound state occurs exactly at the Fermi level, and the impurity spin is screened by an in-gap state oppositely polarized, in a similar way to Kondo screening in a normal metal through singlet formation. Magnetic properties have been found in gated MoS$_2$, but in TaSe$_2$, there has been until now no evidence for magnetic interactions. Thus, it seems difficult to discuss here bound states formed through magnetic scattering.

On the other hand, zero-energy resonant bound states located exactly at the Fermi level arise in reduced symmetry superconductors, such as $d$-wave or more complex superconductors, when some kind of scattering leads to a sign change or a phase slip of the underlying wave function. A zero voltage...
conductance peak can appear thus close to impurities, at a surface or close to crystal boundaries. An intriguing possibility is that some sort of unconventional superconductivity appears within the $2H$-$\text{TaSe}_2$ sheets, either in the form of a $d$-wave order parameter or of sign changing superconductivity between different sheets. Such a possibility would imply that the superconducting properties change from conventional $s$-wave in multilayers to unconventional reduced symmetry superconductivity when decreasing the thickness of the sample down to a single layer.

Resonant scattering gives sharp states with vanishing energy width. In the experiment, the observed peak is often broadened. Peak widths of several tenths of mV, being mostly smaller than the gap value, are found in many cases. For instance, the peak observed at the vortex core in $2H$-$\text{NbSe}_2$ is about $1/3$ of the gap value. The zero-bias conductance peak we observe here has a strong broadening, of the same order as the gap value. Broadening of zero-bias conductance peaks in superconductors has been related to impurities, random disorder, or complex gap variations over the Fermi surface. We do not find evidence for impurities or random disorder. But gap size changes in different parts of the Fermi surface are likely to appear in the involved Fermi surface of $2H$-$\text{TaSe}_2$.

Other zero-bias conductance peaks have been discussed in low-dimensional structures, quantum dots, and graphene. Bound states are formed by confinement or at interfaces, and, under appropriate conditions, these can lead to a zero-energy state. For instance, edge states in graphite ribbons have been predicted to show flat bands near the edges. Electrostatic gating of graphene has been proposed to lead to confinement-induced sharp peaks in the density of states. The combination of Coulomb blockade and quantum dots can exhibit zero-bias anomalies, and interface bound states have been predicted at the interface between graphene and superconductors. The role of interface and charging effects seems difficult to discuss in our single layers of $2H$-$\text{TaSe}_2$ with the available data. The very recent discovery of superconductivity in gate-tuned MoS$_2$ devices shows that superconductivity can arise at the interface between a dichalcogenide, which is semiconducting in the bulk, and a superconductor.

The V-shaped conductance dip [yellow curves in Fig. 4(a)] can be related to decreasing density of states due to superconducting correlations, in particular the decrease below $150 \mu V$. Such curves are closely odd with conventional BCS expressions, and highlight a rather peculiar density of states. They do not show any charge-order related modulations, and could be thus reflecting behavior of the parts of the Fermi surface which are not affected by charge order.

Let us remark that the crossing between regions with widely different tunneling conductance occurs very sharply, just at atomic size at the step between both layers. This implies that there is significant electronic decoupling between the topmost $2H$-$\text{TaSe}_2$ and the substrate $1T$-$\text{TaSe}_2$ layers. The $2H$-$\text{TaSe}_2$ showing the zero-bias peak acts like a separate single layer weakly coupled to its substrate. This, together with the multiband properties of the Fermi surface, could be the origin of peculiar electronic features. Separated layers may include, in particular, enough electron-electron repulsion to establish sign changing superconductivity.

In summary, we have found highly anomalous tunneling conductance features on the surface of $2H$-$\text{TaSe}_2$. We observe a significant increase of the critical temperature close to the surface, and a zero-bias peak in single-layer crystals of $2H$-$\text{TaSe}_2$ when they lie on top of a surface of $1T$-$\text{TaSe}_2$. The zero-bias peak is modulated by a charge-density wave, and coexists with a V-shaped conductance dip. We do not fully understand the microscopic origin for the zero-bias peak, but we have shown that it disappears at the same temperature at which superconducting correlations disappear in multilayered $2H$-$\text{TaSe}_2$. We briefly discuss different possibilities to explain the peak, including the presence of a resonant state at the Fermi level. Single-layer properties of these materials have unexpected low-energy features which are totally different from the bulk.

**ACKNOWLEDGMENTS**

We acknowledge discussions with F. Guinea, V. Vinokur, T. Baturina, A. I. Buzdin, and P. Monceau. We also acknowledge advice and discussions about the dichalcogenides with J. L. Vicent. The Laboratorio de Bajas Temperaturas is associated with the ICM of the CSIC. This work was supported by the EU (ERC Advanced Grant SPINMOL and COST MP-1201), the Spanish MINECO (Consolider-Ingenio in Molecular Nanoscience, CSD2007-00010 and projects FIS2011-23488, MAT2011-25046, MAT2011-22785 and ACI-2009-0905, co-financed by FEDER), by the Comunidad de Madrid (program Nanobiomagnet) and the Generalitat Valenciana (Programs Prometeo and ISIC-NANO).