Direct Electron-Diffraction Evidence of Charge-Density-Wave Formation in NbSe$_3$


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Direct Electron-Diffraction Evidence of Charge-Density-Wave Formation in NbSe₃


Department of Physics, Hokkaido University, Sapporo 060, Japan

and

K. Yamaya and Y. Abe

Department of Nuclear Engineering, Hokkaido University, Sapporo 060, Japan

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In NbSe₃ one-dimensional diffuse scattering was observed above 140 K and the satellite spots at a lower temperature. They correspond to the Kohn anomaly and the superlattice coupled with charge-density wave, respectively. The superlattice was found to be incommensurate, with the period a' = a, b' ≈ 4.1b, and c' = 1.

Recently, the transition-metal trichalcogenides MX₃ (M: Nb, Ta; X: S, Se) have been the subjects of several investigations. The needle-shaped materials have the chain structure of the transition-metal atoms surrounded by the chalcogen atoms along the needle axis. The distance between metal atoms along the chain is much shorter than the interchain distances. Therefore, the interactions between chains seem to be weak and one-dimensional behavior is expected in MX₃.

One of the MX₃ family, NbSe₃, is characterized by two cusps in the temperature variation of the electrical resistivity along the b axis (needle axis) at 59 and 143 K. These anomalies have been explained by assuming the formation of a charge-density wave (CDW). Especially the suppression of the two peaks in the resistivity by strong electrical fields is explained as Zener breakdown across the very small gaps at the Fermi surface induced by the formation of CDW. In this Letter we present the first direct evidence of CDW formation, by electron diffraction above 130 K.

NbSe₃ was synthesized by direct reaction of Nb and Se without carriers. The morphology and the resistivity anomalies of our crystals were the same as those reported by Haen et al. The residual resistivity ratio was ~70. It was found from x-ray diffraction studies with both the precession camera and the counter method, that the (0k0) reflections were missing when k = 2n + 1. This extinction rule indicates that the space group of NbSe₃ is either P2₁ or P2₁/m, which have higher symmetry than Pm by Meerschaut and Rouxel. It seems that the presence of the 2₁ screw axis along the needle axis is a common feature of MX₃. The electron diffraction of NbSe₃ was performed with the acceleration voltage of 100 kV. The sample holder was kept at 130 K. But the temperature of the individual crystal may be slightly higher as a result of irradiation by the electron beam, though a heat link to the sample holder would release some of the heat.

Figure 1 shows the b*-c* plane of the reciprocal space above 140 K. The anomaly in diffuse scattering can be clearly seen to be one-dimensional. The position of this one-dimensional anomaly along the b* axis is (0.244 ± 0.004)b*, which is a little smaller than the commensurate value of 0.25b*. This incommensurate streak

FIG. 1. Electron diffraction pattern of NbSe₃ above 140 K, showing one-dimensional anomaly in diffuse scattering (marked by arrows).

which yields satisfactory lattice-vibrational spectra predicts negative (unphysical) values for the heat of formation of vacancies in metals such as Al.
corresponds to the Kohn anomaly. In Fig. 2, which shows the $a^*b^*$ plane below 140K, the satellite spots are exactly at the position of the Kohn anomaly. From a similar observation on the $b^*c^*$ plane below 140 K, it was found that the satellite spots have no $a^*$ and $c^*$ components, at least at this temperature. These observations establish that an incommensurate superlattice with $a^* = a$, $b^* = 4.1b$, and $c^* = c$ occurs in NbSe$_3$. The incommensurability itself is a strong support of the CDW mechanism. Thus our electron diffraction experiments confirmed the existence of CDW in NbSe$_3$.

We reported that TaS$_3$ undergoes the Peierls transition at 220 K. Comparing the two materials, the following points are worth noting:

(i) The low-temperature phase of TaS$_3$ has the commensurate superperiod of $c^* = 4c$ within the accuracy of 1%, while it is incommensurate in NbSe$_3$. (Note that the chain axis is the $b$ axis in NbSe$_3$ and the $c$ axis in TaS$_3$.)

(ii) From the nearly equal periods of the one-dimensional modulation, it is expected that the Fermi surfaces of these materials have common features in the normal state, e.g., nearly the same relative position in the Brillouin zone.

(iii) TaS$_3$ is semiconducting below the transition temperature, while NbSe$_3$ remains metallic. The Fermi surface of NbSe$_3$ must be nonplanar. From the resistivity change below 140 K, we estimate that half of the Fermi surface vanishes at the transition.

(iv) In TaS$_3$, the coexistence of the diffuse lines at 0.25$c^*$ with satellites was observed, while in NbSe$_3$, the diffuse lines are absent below the transition as shown in Fig. 2. The diffuse streak connecting the parent spots perpendicular to the chain axis was often observed in both materials. It was not determined whether the diffuse streak is due to the soft phonon mode or the poor crystallization.

(v) In TaS$_3$, the satellite spots also have components perpendicular to the chain axis, while in NbSe$_3$, only the component along the chain axis was observed.

We could not extend the electron diffraction study below 130 K because of the limitation of our cooling apparatus. It must be very interesting to determine the nature of the 59-K anomaly.

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FIG. 1. Electron diffraction pattern of NbSe$_3$ above 140 K, showing one-dimensional anomaly in diffuse scattering (marked by arrows).
FIG. 2. Electron diffraction pattern of NbSe$_3$ below 140 K, showing the satellite spots (marked by arrows).