

Visualization of grain boundaries in 2D Transition Metal Telluride

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Abstract

2D transition-metal dichalcogenides and their van der Waals heterostructures recently have drawn increasing attention owing to their promising properties. Particularly, tellurium-based compounds such as MoTe_2 and WTe_2 exhibit semiconducting and semi-metallic characteristics, which are associated with their distinct polymorphic crystal structures [1]. These various lattice phases make MoTe_2 a favourable candidate for both fundamental studies and device applications. However, the energy difference between the 2H and 1T' phases of MoTe_2 is small, as well as 1T'- MoTe_2 is thermodynamically stable [2] make it challenging for phase-selective synthesis. In addition, the location of grain boundary and lattice orientation is required when dealing with non-exfoliated samples, as the mechanical and electrical properties governed by lattice orientation and strongly influenced by presence of grain boundary. Herein, we report a controlled synthesis of highly crystalline 1T'- and 2H- MoTe_2 crystals by APCVD technique. Furthermore, we investigated the existence of grain boundaries using a simple tool called 'polarized optical microscopy', which then further supported by angle resolved Raman scattering and SHG [3]. Polycrystalline 1T'- MoTe_2 are normally Mo-terminated and shows anisotropic optical absorption leading to the clear visualization of lattice domains. On the other hand, 2H- MoTe_2 lattice grains does not exhibits any difference under polarized light. Grain boundary visualization

in 2D tellurides using this simple tool paves the way for the realization of grain boundary free real devices with desired lattice orientation.

References

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Figures

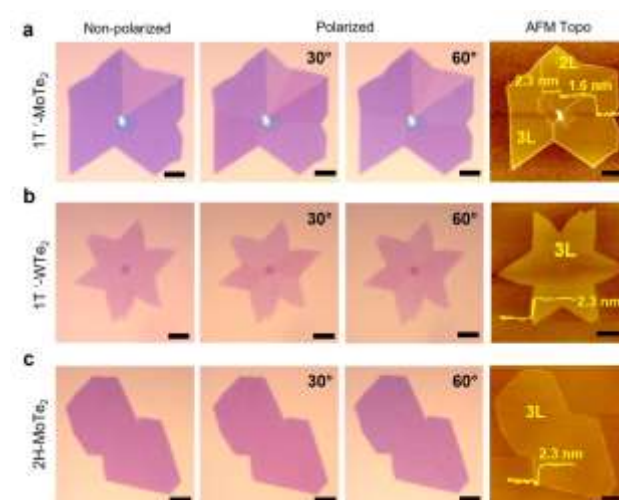


Figure 1: Non-polarized, polarized optical images with different rotational angles and AFM topography of (a) 1T'- MoTe_2 , (b) 1T'- WTe_2 and (c) 2H- MoTe_2 respectively. Scale bar, 10 μm .