## Abstract Submitted for the MAR14 Meeting of The American Physical Society

SPE-LEEM Studies on the Surface and Electronic Structure of 2-D Transition Metal Dichalcogenides (Part II)<sup>1</sup> WENCAN JIN, PO-CHUN YEH, NADER ZAKI, DATONG ZHANG, Columbia Univ, JERZY SADOWSKI, ABDULLAH AL-MAHBOOB, Brookhaven National Laboratory, AREND VAN DE ZANDE, Energy Frontier Research Center, Columbia Univ, DANIEL CH-ENET, JERRY DADAP, IRVING HERMAN, Columbia Univ, PETER SUTTER, Brookhaven National Laboratory, JAMES HONE, RICHARD OSGOOD, Columbia Univ — In this work, we studied the surface and electronic structure of monolayer and few-layer exfoliated MoS<sub>2</sub> and WSe<sub>2</sub>, as well as chemical-vapor-deposition (CVD) grown MoS<sub>2</sub>, using Spectroscopic Photoemission and Low Energy Electron Microscope (SPE-LEEM). LEEM measurements reveal that, unlike exfoliated MoS<sub>2</sub>, CVD-grown  $MoS_2$  exhibits grain-boundary alterations due to surface strain. However, LEEM and micro-probe low energy electron diffraction show that the quality of  $\text{CVD-grown MoS}_2$  is comparable to that of exfoliated  $\text{MoS}_2$ . Micrometer-scale angleresolved photoemission spectroscopy (ARPES) measurement on exfoliated  $MoS_2$  and WSe<sub>2</sub> single-crystals provides direct evidence for the shifting of the valence band maximum from  $\Gamma$  to K, when the layer number is thinned down to one, as predicted by density functional theory. Our measurements of the k-space resolved electronic structure allow for further comparison with other theoretical predictions and with transport measurements.

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