## Abstract Submitted for the MAR17 Meeting of The American Physical Society

Characterizations of  $ZnO/MoO_3$  superlattices grown by Atomic Layer Deposition Y.S. HONG, Q.Y. CHEN, P.V. WADEKAR, W.C. HSIEH, C.F. CHANG, Department of Physics, National Sun Yat-Sen University, H.C. HUANG, Department of Material Science and Optoelectronics, National Sun Yat-Sen University, C.M. SHIAU, C.H. LEE, Y.P. CHENG, C.Y. DANG, P.C. KUNG, Y.Y. LIANG, S.H. HUANG, Z.Y. WU, C.M. LIN, S.T. YU, L.W. TU, Department of Physics, National Sun Yat-Sen University, N.J. HO, Department of Material Science and Optoelectronics, National Sun Yat-Sen University, H.W. SEO, Department of Physics, Jeju National University, W.K. CHU, Texas Center of Superconductivity and Department of Physics, University of Houston  $- ZnO/MoO_3$  superlattices (SLS) were prepared by ALD on  $Al_2O_3$  substrates at 450K. The growth rates are 0.17 per cycle for MoO<sub>3</sub> and 1.66 for ZnO, according to XRR. The MoO<sub>3</sub> films were found amorphous when deposited separately, while ZnO polycrystalline. However  $MoO_3$  became polycrystalline and ZnO textured grown into SLS, as judged by the electron diffraction patterns. The ZnO thicknesses were fixed at 6 nm per period while  $MoO_3$  varied from 2 to 6 nm. The nanostructures as examined by TEM indeed show expected periodicity consistent with XRR. HRTEM also gave clear interfaces of the SLS with certain regions of imperfection. Thence, we conclude that amorphous  $MoO_3$  would crystallize when grown adjacent to an initial layer of ZnO. PL was employed to investigate the possible variations of their bandgaps when the constituent ZnO and  $MoO_3$  brought together in close proximity of nanoscakes. Electronic band structures according to ab initio calculations will be discussed.

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