

Abstract Submitted  
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**Single surface state at a strongly correlated electronic structure in  $\text{Na}_2\text{IrO}_3$ : a candidate d-electron topological insulator** QIANG WANG, University of Colorado at Boulder, Los Alamos National Laboratory, YUE CAO, JUSTIN WAUGH, University of Colorado at Boulder, TONGFEI QI, OLEKSANDR KORNETA, GANG CAO, University of Kentucky, DANIEL DESSAU, University of Colorado at Boulder — We have performed angle-resolved photoemission spectroscopy (ARPES) on  $\text{Na}_2\text{IrO}_3$ , a 5d transition metal oxide (TMO) which is a strong insulator with a honeycomb lattice structure and has been theoretically proposed as a candidate for a new class of topological insulators (TIs). The near  $E_F$  electronic structure of  $\text{Na}_2\text{IrO}_3$  was carefully mapped, which shows an overall agreement to the first-principle calculations with spin-orbit (SO) coupling and electron correlation (U), though certain discrepancy remains. Specifically, we found an extra electron-like pocket near the Fermi level with Dirac-cone-like dispersion around  $\Gamma$  point. The further photon energy dependent studies show no  $k_z$ -dispersion of this electron-like pocket, and the metal deposition studies show a great enhancement and sharpening of this feature. These results confirm its surface state nature and suggest a possible single topological surface state at the Brillouin zone (BZ) center in  $\text{Na}_2\text{IrO}_3$ .

Qiang Wang  
Los Alamos National Laboratory

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