

Topological Material in the III–V Family: Heteroepitaxial InBi on InAs

Laurent Nicolai^{1, a)}, Ján Minár^{1, b)}, Maria Christine Richter^{2, 3}, Olivier Heckmann^{2, 3}, Jean-Michel Mariot⁴, Uros Djukic², Johan Adell⁵, Mats Leandersson⁵, Janusz Sadowski⁶, Jürgen Braun⁷, Hubert Ebert⁷, Jonathan D. Denlinger⁸, Ivana Vobornik⁹, Jun Fujii⁹, Pavol Šutta¹, Gavin R. Bell¹⁰, Martin Gmitra^{11, 12}, and Karol Hricovini^{2, 3, c)}

¹University of West Bohemia/New Technologies – Research Centre (NTC), Plzeň, Czech Republic

²Laboratoire de Physique des Matériaux et des Surfaces, CY Cergy-Paris Université, Gif-sur-Yvette, France

³LIDYL, Université Paris-Saclay, CEA, CNRS, Gif-sur-Yvette, France

⁴Laboratoire de Chimie Physique–Matière et Rayonnement, Sorbonne Université, CNRS, Paris, France

⁵MAX IV Laboratory, Lund University, Lund, Sweden

⁶Institute of Physics, Polish Academy of Sciences, Warsaw, Poland

⁷Department Chemie, Ludwig-Maximilians-Universität, München, Germany

⁸Advanced Light Source, Lawrence Berkeley National Laboratory, California, USA

⁹TASC Laboratory, Istituto Officina dei Materiali, CNR, Trieste, Italy

¹⁰Department of Physics, University of Warwick, United Kingdom

¹¹Institute of Physics, Pavol Jozef Šafárik University in Košice, Košice, Slovak Republic

¹²Institute of Experimental Physics, Slovak Academy of Sciences, Košice, Slovak Republic

Corresponding author: ^{a)}lnicolai@ntc.zcu.cz

^{b)}jminar@ntc.zcu.cz

^{c)}karol.hricovini@cyu.fr

Abstract, InBi(001) is formed epitaxially on InAs(111)-A by depositing Bi onto an In-rich surface [1]. Angle-resolved photoemission measurements reveal topological electronic surface states, close to the M high symmetry point. This demonstrates a heteroepitaxial system entirely in the III–V family with topological electronic properties. InBi shows coexistence of Bi and In surface terminations, in contradiction with other III–V materials. For the Bi termination, the study gives a consistent physical picture of the topological surface electronic structure of InBi(001) terminated by a Bi bilayer rather than a surface formed by splitting to a Bi monolayer termination. Theoretical calculations based on relativistic density functional theory and the one-step model of photoemission clarify the relationship between the InBi(001) surface termination and the topological surface states, supporting a predominant role of the Bi bilayer termination. Furthermore, a tight-binding model based on this Bi bilayer termination with only Bi–Bi hopping terms, and no Bi–In interaction, gives a deeper insight into the spin texture.

[1] Nicolai *et al.*, Phys. Rev. Res. **6**, 043116 (2024)