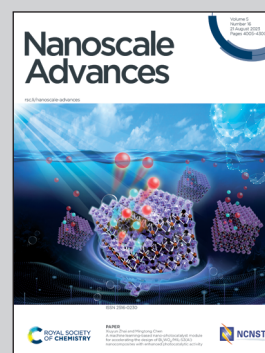


Showcasing research from Professor Samia Subrina's lab,
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Numerical characterization of the electronic and optical
properties of plumbene/hBN heterobilayer using first-principles
study

A detailed investigation of the structural, electronic, and optical
characteristics of a plumbene/hBN heterobilayer has
been conducted using first-principles calculations under the
framework of density functional theory. One of the fascinating
findings of the study is that the composite structure exhibits a
direct bandgap compared to pristine plumbene, which has
an indirect bandgap. In addition, a semiconductor to metallic
transition takes place upon the introduction of biaxial compressive
strain. Optical properties suggest the potential of the proposed
heterobilayer in solar cells and UV photodetectors.

As featured in:



See Samia Subrina *et al.*,
Nanoscale Adv., 2023, 5, 4095.