Supporting Information

Moiré potential, lattice corrugation, and band gap variation in a twist-free $MoTe_2/MoS_2$ heterobilayer

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Fig. S1. Top and side views of the atomic structure of a $MoTe_2/MoS_2$ bilayer with (7×7)/(8×8) [A], (8×8)/(9×9) [B], and (10×10)/(11×11) [C] stackings. Lattice corrugation in the MoS_2 layer is clearly seen.

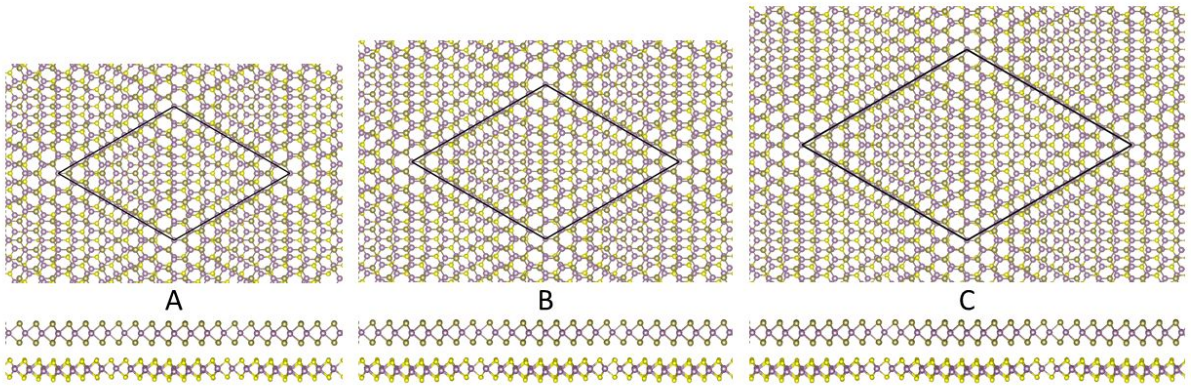

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Fig. S2. Density of states of a $MoTe_2/MoS_2$ bilayer. The local DOS projected on the Mo atoms at or near *AA*, *AB*, and *BA* sites with three-fold rotational symmetry are drawn in bold lines.