

Electric field modulation of electronic properties in type-II phosphorene/PbI₂ van der Waals heterojunction

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Supporting information:

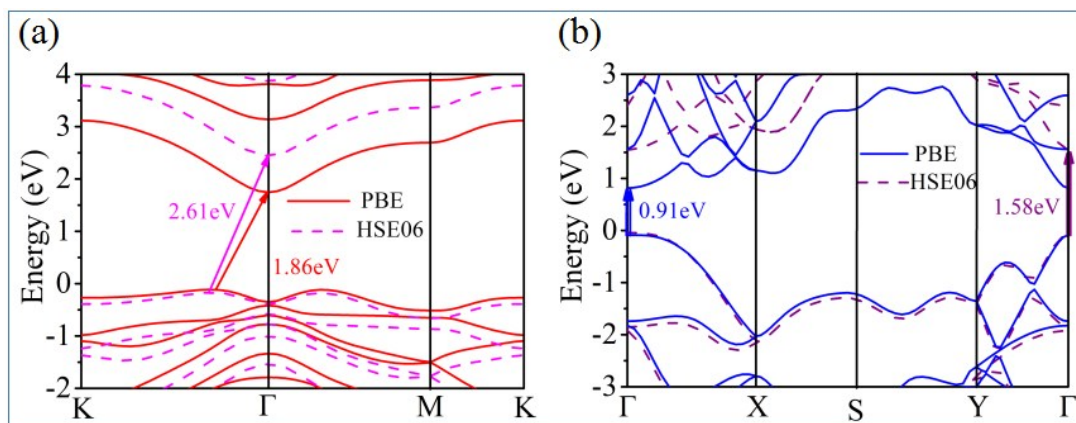


Fig. S1 The band structures of PbI₂ (a) and phosphorene (b) monolayer. The solid and dash line are the band structure calculated by the PBE and HSE06.

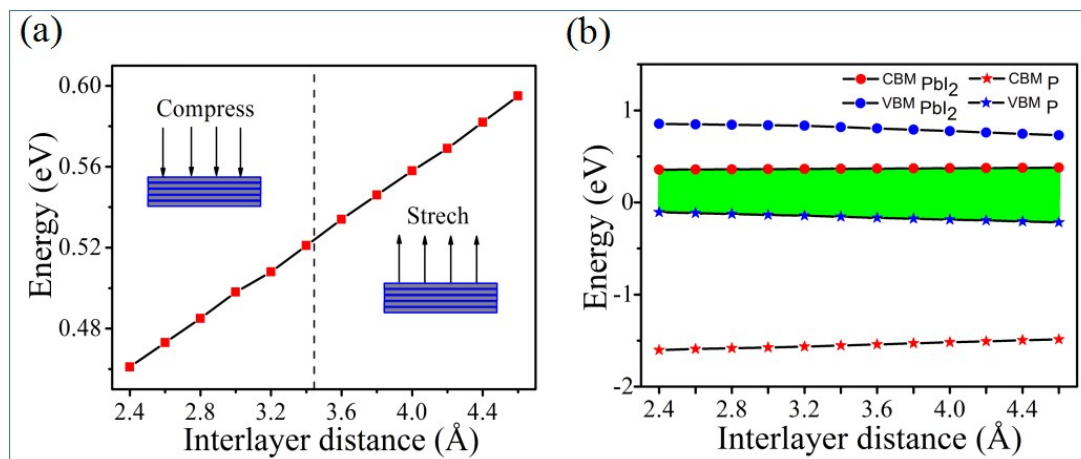


Fig. S2 Variation of band gap (a) and band edge (b) of the P/PbI₂ vdW heterojunction under different interlayer distance.