

Semiconductor and topological phases in lateral heterostructures constructed from germanene and AsSb monolayers

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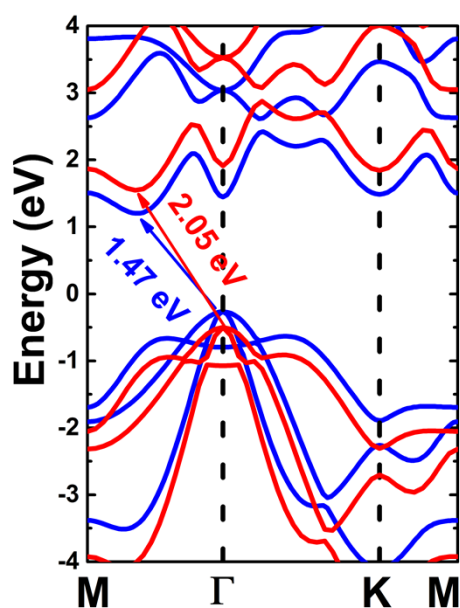


Figure S1: Electronic band structure of AsSb monolayer calculated with PBE (blue curve) and HSE06 (red curve) functionals.

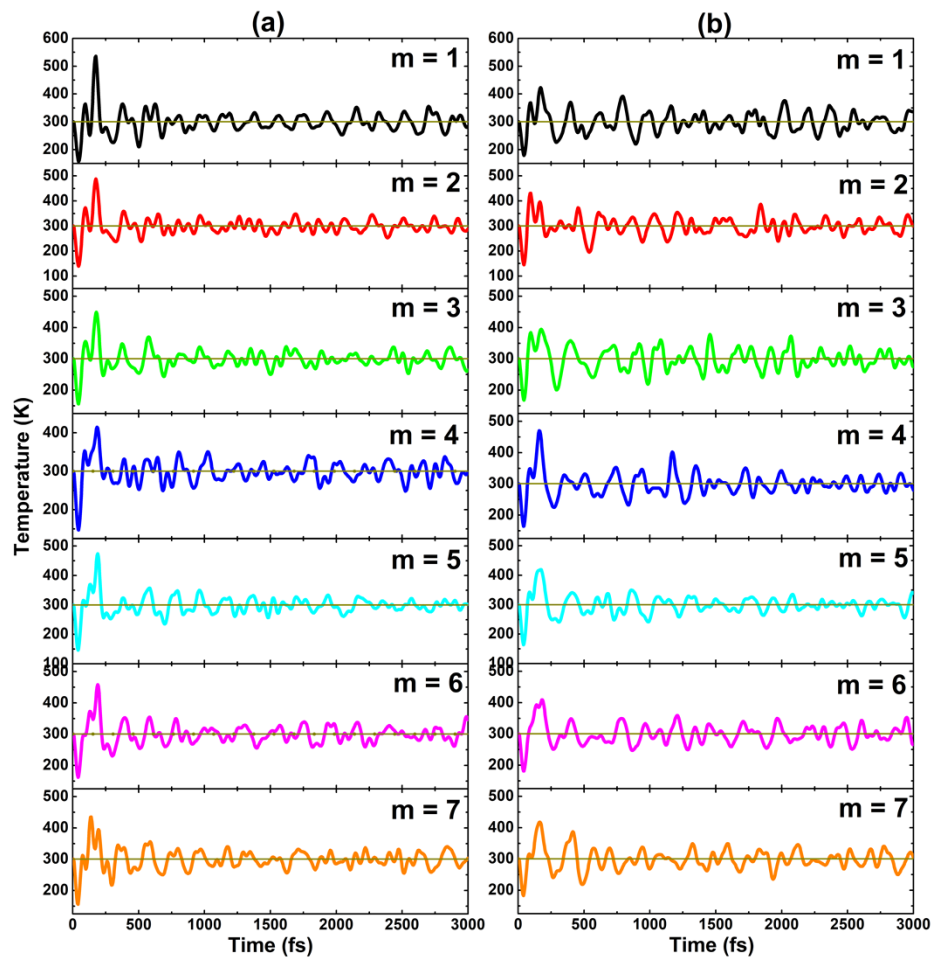


Figure S2: AIMD simulations at 300 K of (a) Armchair- and (b) Zigzag-interline $(\text{Ge})_m\text{-(AsSb)}_{8-m}$ lateral heterostructures.

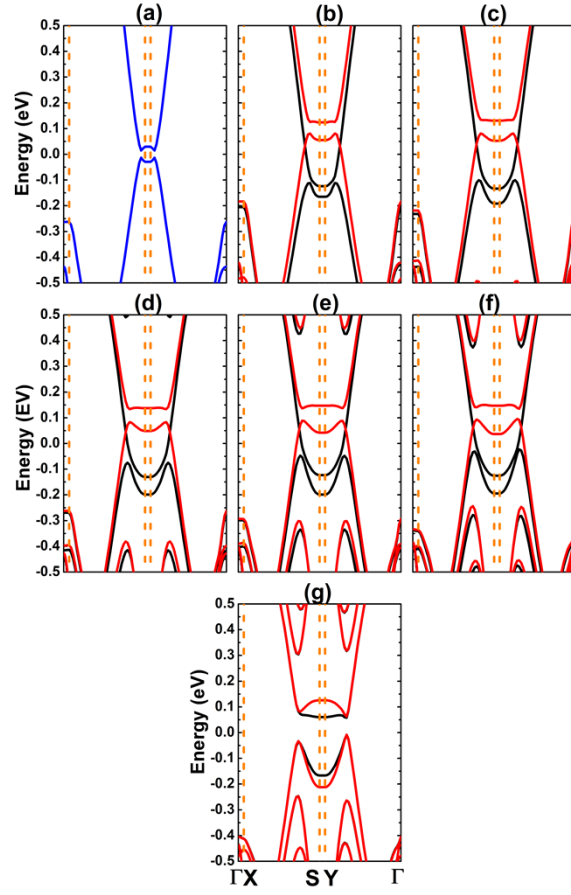


Figure S3: Electronic band structure (Blue curve: non-spin polarization; Black curve: Spin-up; Red curve: Spin-down) of zigzag-interline $(\text{Ge})_m\text{-(AsSb)}_{8-m}$ lateral heterostructures with (a) $m = 1$, (b) $m = 2$, (c) $m = 3$, (d) $m = 4$, (e) $m = 5$, (f) $m = 6$, and (g) $m = 7$.