Atomic-Scale Carving of Nanopores into a van der Waals Heterostructure with Slow Highly Charged Ions

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Supporting information

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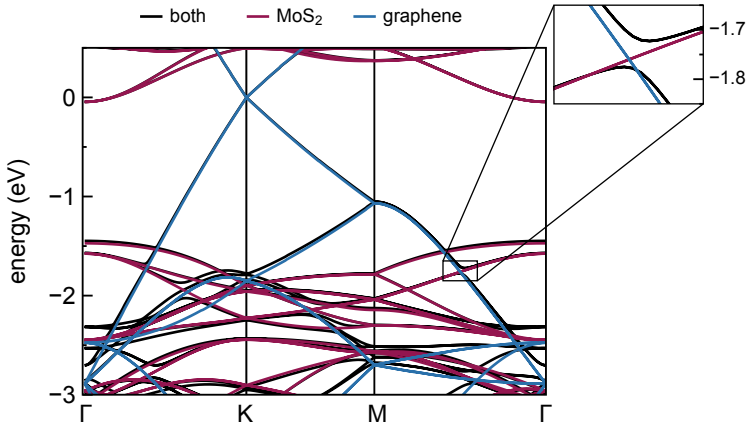

Figure 1: Band structure of a 3×3 MoS_2 + 4×4 graphene bilayer system from density functional theory (see methods). Band structures of the individual, separate layers are shown in red (MoS_2) and blue (graphene). Changes in the bandstructure are minimal, with small avoided crossings (see zoom-in).

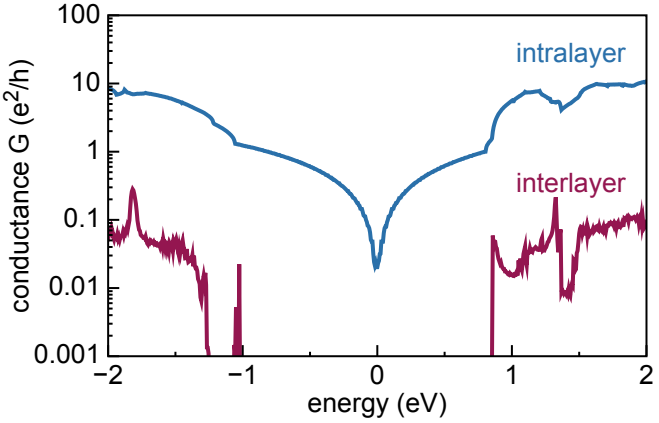

Figure 2: Intralayer conductance of graphene/MoS_2 heterostructure is at least two orders of magnitude larger than the interlayer conductance.