## Theoretical Study on Tuning Band Gap and Electronic Properties of Atomically Thin Nanostructured ${\rm MoS_2/Metal}$ Cluster Heterostructures

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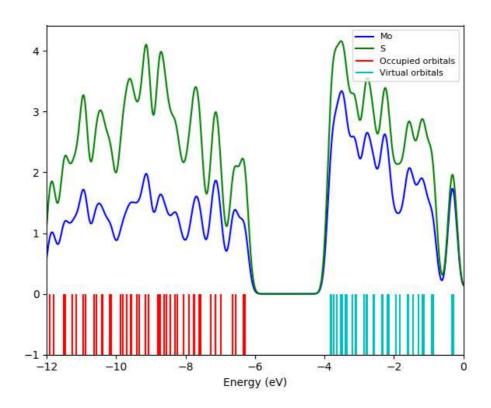


Figure 1: Density of states (DOS) of  $MoS_2$ .

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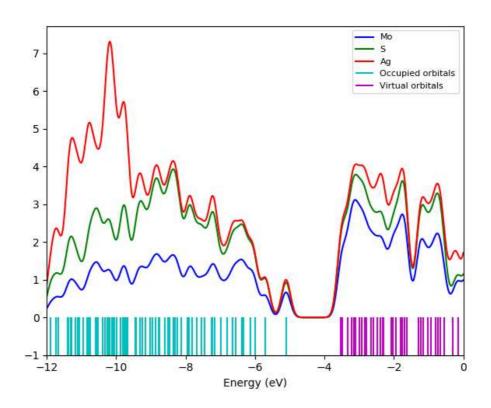


Figure 2: Density of states (DOS) of Silver/MoS $_2$ .

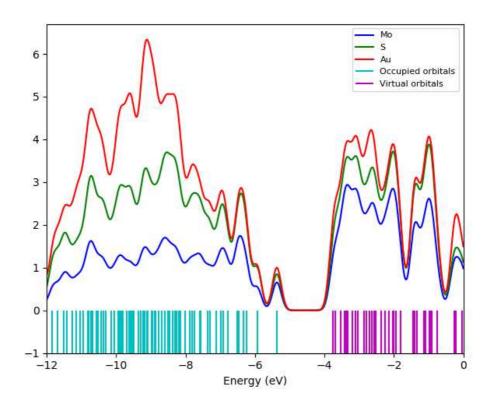


Figure 3: Density of states (DOS) of  $Gold/MoS_2$ .