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Supporting Infomation

Tunable Electronic Properties of Nitrogen and Sulfur Doped Graphene: Density Functional Theory Approach

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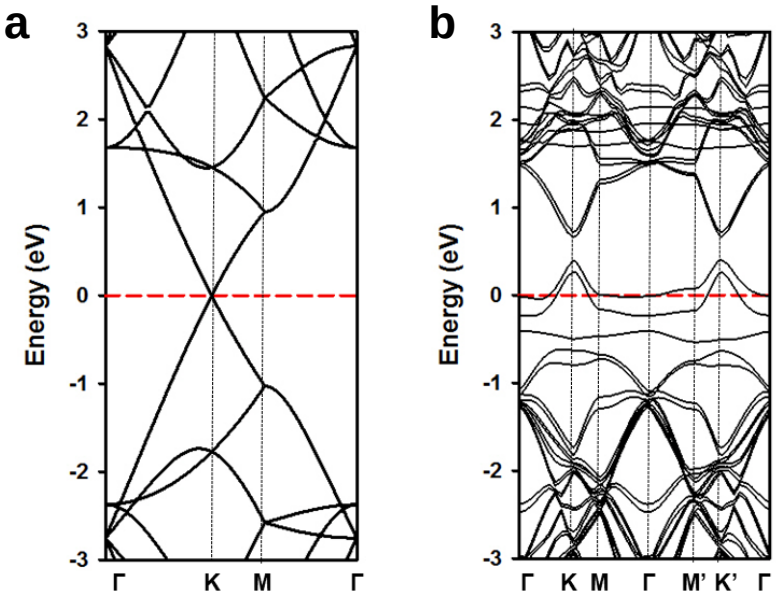

Figure S1. Calculated band structures of the (**a**) pristine graphene and (**b**) graphene with mono-vacancy.