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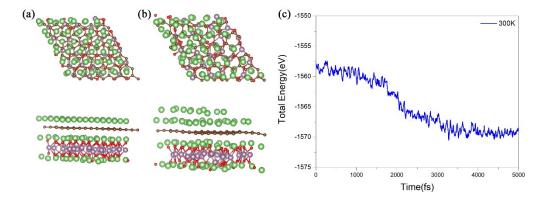


Figure S1. (a) and (b) The structures of MoO_2/G nanocomposite with three layers Li atoms adsorption before and after 5ps AIMD simulations at 300K. (c) The total energy shifts with time of AIMD simulation at 300 K.