

Elastic Response of Graphene Nanodomains

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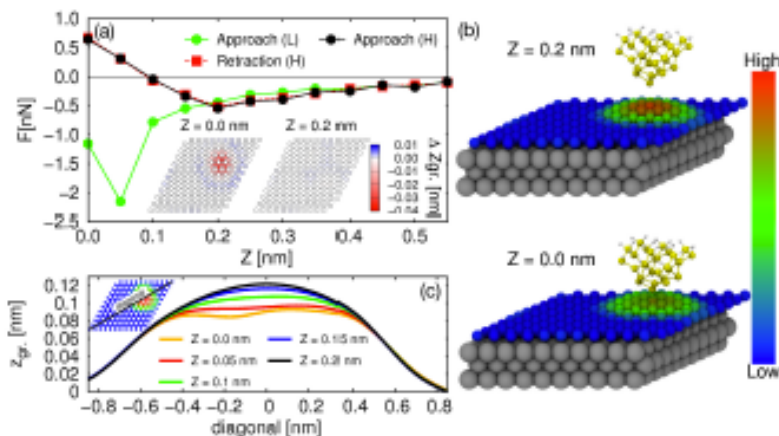
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Due to its exceptional properties graphene has become a topic of utmost importance [1-3]. One of the most remarkable features of a free standing graphene monolayer is its exceptional mechanical strength, which has allowed the development of electromechanical resonators [4,5].

In this study we have focused on the elastic properties of graphene adsorbed on a Ru(0001) surface. When graphene is grown on Ru(0001) a regular arrangement of nanodomains is spontaneously formed. Our experimental measurements [6] show that this periodic structure responds in a perfectly reversible way to perpendicular displacements up to 40% from the equilibrium structure. The elasticity of the nanodomains have been further proven by realistic DFT (density functional theory) calculations, we have estimated a normal stiffness k of about 40N/m [6].

In the figure we show the calculated force versus the distance for the hill site indentation, together with the optimized geometries of the force curve on the hill site of the nanodomain (0.2 nm) and at maximum indentation (0.0 nm). The figure also shows the profiles of the indented graphene hill ($Z \leq 0.2$ nm) along the moiré unit cell diagonal (see inset Fig. c).



References:

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