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Erratum to "*Multilevel B-Spline Repulsive Energy in Nanomodeling of Graphenes*" [Journal of Surface Engineered Materials and Advanced Technology Vol. 4 No. 2 (April 2014) 75-86]

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In addition, please remove **Figure 5** because it has little relevance with the proposed method. Discard also the corresponding description on page 84: "As a next test, we consider the complex band structures for using the DFT and SE computations whose results are respectively displayed in **Figure 5(a)**, **Figure 5(b)** for the graphene with chirality (1,0). The plots depict band lines which are not shown as continuous curves but as sets of sampling points. The points which are purely real and explicitly complex are depicted in red and green respectively".