Molecular docking simulation, synthesis and 3D pharmacophore studies of novel 2-substituted-5-nitrobenzimidazole derivatives as anticancer agents targeting VEGFR-2 and c-Met

Hanan Refaat, Heba Mohamed , Kamilia M. Amin, Fadi M. Awadallah

Abstract

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Future University In Egypt (http://www.fue.edu.eg)