Author Correction to: Quantum chemical studies, vibrational analysis, molecular dynamics and docking calculations of some *ent*-kaurane diterpenes from *Annona vepretorum*: a theoretical approach to promising anti-tumor molecules

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In the original published version of the article, some typos are in the text, being in disagreement with the information found in Tables 1, 2, 3 and 4 and in the section “Molecular dynamic studies”.

We would like to inform that instead of reading “C5-C4-C9”, “C11-C13-C14” and “C7-C11” (Geometry and parameters, second paragraph, seventeenth, twenty-fifth and twenty-

eighth lines, respectively), “C2-C6” (Geometry and parameters, third paragraph, third line), “C10-C15” (Geometry and parameters, third paragraph, fifth line), “H51” (Electrostatic potential maps, second paragraph, antepenultimate line) and “1624” (Infrared analysis, second paragraph, ninth line), read C5-C4-C7, C12-C13-C14, C13-C14, C5-C6, C10-C5, H52 and 1642 cm⁻¹ respectively.

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