Correction to: Evaluation of affinity of bioactive isolates from various coffee extracts through binding with PPAR- γ with the use of isothermal titration calorimetry and docking simulation to prevent antidiabetic effects

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In the original publication of the article, under the section ‘Molecular modelling’, the sentence that read as “The chemical structures of CGAs and caffeine molecules were built up and fully optimized (B3LYP/6-31+G dip) [29, 30]” should read as “The chemical structures of CGAs and caffeine molecules were built up manually using Gasteiger scheme [29, 30]”

The reference [29] was published incorrectly. The corrected reference [29] is given below:

Reference

29. Budryn G, Grzelczyk J, Pérez-Sánchez H. Binding of red clover isoflavones to actin as a potential mechanism of anti-metastatic activity restricting the migration of cancer cells. *Molecules*. 2018;23:2471.

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