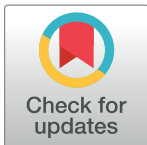


CORRECTION

# Correction: Unveiling a novel transient druggable pocket in BACE-1 through molecular simulations: Conformational analysis and binding mode of multisite inhibitors

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There is an error in the Funding section. The correct Funding statement is as follows: This work was supported by the Ministerio de Economía y Competitividad (SAF2014-57094-R), Generalitat de Catalunya (GC; 2014SGR1189, 2014SGR52), and ICREA Academia (FJL). The Consorci de Serveis Universitaris de Catalunya (CSUC; FJL) is acknowledged for providing computational resources. This work used the ARCHER UK National Supercomputing Service (<http://www.archer.ac.uk>), via the HECBioSim consortium (EPSRC Grant EP/L000253/1), and the Spanish Barcelona Supercomputer Center (<https://www.bsc.es>; project BCV-2017-2-0012). A fellowship from GC to ODP is gratefully acknowledged. The funders had no role in study design, data collection and analysis, decision to publish, or preparation of the manuscript.



## Reference

1. Di Pietro O, Juárez-Jiménez J, Muñoz-Torrero D, Laughton CA, Luque FJ (2017) Unveiling a novel transient druggable pocket in BACE-1 through molecular simulations: Conformational analysis and binding mode of multisite inhibitors. PLoS ONE 12(5): e0177683. <https://doi.org/10.1371/journal.pone.0177683> PMID: 28505196

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