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## Correction: Exploiting butyrylcholinesterase inhibitors through a combined 3-D pharmacophore modeling, QSAR, molecular docking, and molecular dynamics investigation

 Sunil Kumar,<sup>a</sup> Amritha Manoharan,<sup>a</sup> Jayalakshmi J,<sup>a</sup> Mohamed A. Abdelgawad,<sup>bc</sup> Wael A. Mahdi,<sup>d</sup> Sultan Alshehri,<sup>d</sup> Mohammed M. Ghoneim,<sup>ef</sup> Leena K. Pappachen,<sup>a</sup> Subin Mary Zachariah,<sup>a</sup> T. P. Aneesh<sup>\*a</sup> and Bijo Mathew<sup>\*a</sup>

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The authors regret that the email addresses were associated with the incorrect institution. The corrected association is as shown above.

The Royal Society of Chemistry apologises for these errors and any consequent inconvenience to authors and readers.

<sup>a</sup>Department of Pharmaceutical Chemistry, Amrita School of Pharmacy, Amrita Vishwa Vidyapeetham, AIMS Health Sciences Campus, Kochi, 682 041, India. E-mail: [bijomathew@aims.amrita.edu](mailto:bijomathew@aims.amrita.edu); [bijovilaventgu@gmail.com](mailto:bijovilaventgu@gmail.com); [aneeshtp@aims.amrita.edu](mailto:aneeshtp@aims.amrita.edu)

<sup>b</sup>Department of Pharmaceutical Chemistry, College of Pharmacy, Jouf University, Sakaka, 72341, Saudi Arabia

<sup>c</sup>Department of Pharmaceutical Organic Chemistry, Faculty of Pharmacy, Beni-Suef University, Beni-Suef, Egypt

<sup>d</sup>Department of Pharmaceutics, College of Pharmacy, King Saud University, Riyadh, 11451, Saudi Arabia

<sup>e</sup>Department of Pharmacy Practice, College of Pharmacy, AlMaarefa University, Ad Diriyah, 13713, Saudi Arabia

<sup>f</sup>Pharmacognosy and Medicinal Plants Department, Faculty of Pharmacy, Al-Azhar University, Cairo, 11884, Egypt

