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

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Correction for 'Exploiting butyrylcholinesterase inhibitors through a combined 3-D pharmacophore modeling, QSAR, molecular docking, and molecular dynamics investigation' by Sunil Kumar *et al.*, *RSC Adv.*, 2023, **13**, 9513–9529, <https://doi.org/10.1039/D3RA00526G>.

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.

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