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Erratum

Erratum to "Novel Potent 5-HT_3 Receptor Ligands Based on the Pyrrolidone Structure. Effects of the Quaternization of the Basic Nitrogen on the Interaction with 5-HT_3 Receptor" [Bioorg. Med. Chem. 10 (2002) 2681]†

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The publishers would like to apologise for an error on page 2681 of the above manuscript. In the second column, line 9, the sentence should read:
"By means of the integration of a classical medicinal chemistry approach [based on extensive qualitative structure–affinity relationship (SAFIR) studies] with a computational approach (receptor mapping, homology modelling and docking simulation) we obtained a picture of the complexity of the ligand–receptor interaction.[5]"

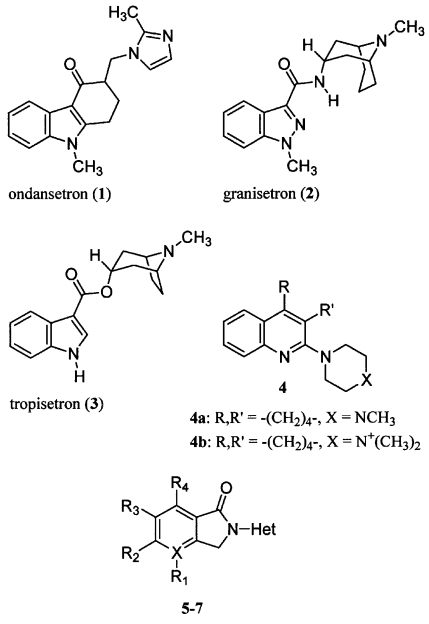

Chart 1.

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