Additions and Corrections

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Xiang-Qun Xie,* Xiu-Wen Han, Jian-Zhong Chen, Michael Eissenstat, and Alexandros MakriyannisHigh-Resolution NMR and Computer Modeling Studies of the Cannabimimetic Aminoalkylindole Prototype WIN-55212-2

Page 4026. Figure 5 is incorrect. The correct Figure 5 appears below.

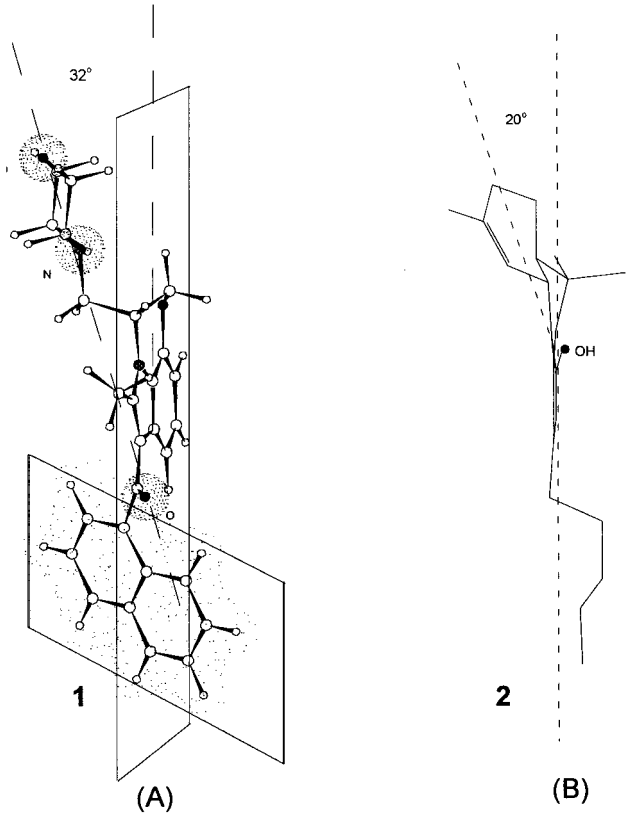

Figure 5. (A) Graphical representation of the favored conformations of **1**, in which the atoms highlighted with van der Waals electron surfaces are potential hydrogen-bonding centers, while the aromatic rings constitute the hydrophobic components. (B) Conformation of **2** from X-ray crystallographic data.

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