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Corrigendum

Corrigendum to "Structure-Based De Novo Design of Non-nucleoside Adenosine Deaminase Inhibitors" [Bioorg. Med. Chem. Lett. 13 (2003) 1115]☆

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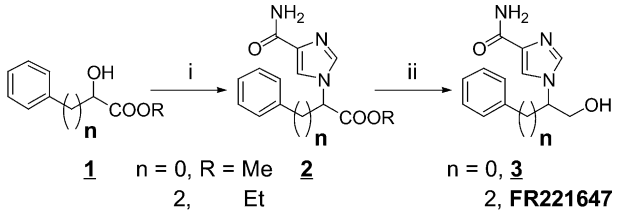
The authors regret that there were errors contained in the above article. The structure of compounds **1** and **2** in Scheme 2 were incorrect. The correct structures are shown below:



The K_i value of EHNA in Table 3 was incorrect. The correct K_i value of EHNA is 37 nM.

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