Pergamon

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Corrigendum

Corrigendum to "Synthesis and Preliminary Biological Evaluation of [^{3}H]-MRE 3008-F20: the First High Affinity Radioligand Antagonist for the Human A_3 Adenosine Receptors"†

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Unfortunately, the structures in Figure 1 and Scheme 1 on page 210 are incorrect. The correct Figure 1 and Scheme 1 are as follows.

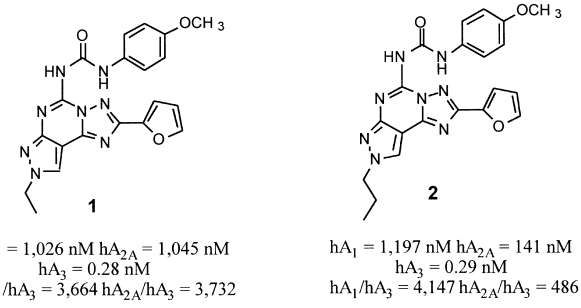
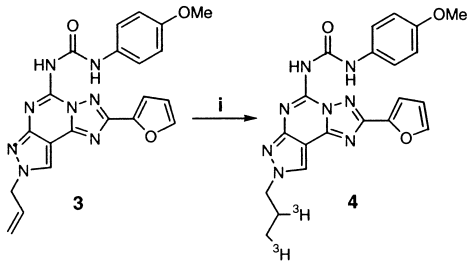


Figure 1. Structures and binding affinities of human A_3 adenosine receptor antagonists.



Scheme 1. Reagents: i: Tritium gas, 10% Pd-C, EtOAc:EtOH 2:1.

†PII of original article S0960-894X(99)00674-5.

PII: S0960-894X(00)00239-0