Accepted Manuscript

Rational design of carbamate-based dual binding site and central AChE inhibitors by a “biooxidisable” prodrug approach: Synthesis, *in vitro* evaluation and docking studies

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PII: S0223-5234(18)30488-4

DOI: 10.1016/j.ejmech.2018.05.057

Reference: EJMECH 10468

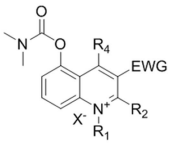
To appear in: *European Journal of Medicinal Chemistry*

Received Date: 16 February 2018

Revised Date: 10 April 2018

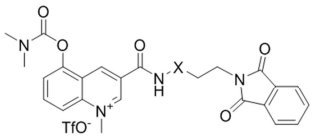
Accepted Date: 31 May 2018

Please cite this article as: M.-L. Țînțaș, V. Gembus, F. Alix, Anaï. Barré, Gaë. Coadou, L. Truong, M. Sebban, C. Papamicaël, H. Oulyadi, V. Levacher, Rational design of carbamate-based dual binding site and central AChE inhibitors by a “biooxidisable” prodrug approach: Synthesis, *in vitro* evaluation and docking studies, *European Journal of Medicinal Chemistry* (2018), doi: 10.1016/j.ejmech.2018.05.057.



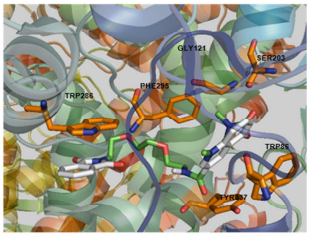
*h*AChE inhibitors **2a-p** binding on active site

SAR analysis



4 X = $-(CH_2CH_2O)_2-$

IC_{50} *h*AChE : 6 nM



- Synthesis and SAR evaluation of compounds **2a-p**
- *In vitro* evaluation of **4** and its prodrug **3** against *h*AChE activity
- Molecular docking simulation of **4**