

Additions and Corrections

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George A. Brine,* Peter A. Stark, Young Liu, F. Ivy Carroll, P. Singh, Heng Xu, Richard B. Rothman: Enantiomers of Diastereoisomeric *cis-N*-[1-(2-Hydroxy-2-phenylethyl)-3-methyl-4-piperidyl]-*N*-phenylpropanamides: Synthesis, X-ray Analysis, and Biological Activities.

Resynthesis and reanalysis of the four stereoisomers revealed that isomers (2*S*,3*R*,4*S*)-**1a** and (2*R*,3*S*,4*R*)-**1c** had been reversed in the binding data reported in Table 2. A corrected Table 2 (abbreviated to stereoisomer binding data) is given below. All other reported biological data are correct. Our binding data are now consistent with data later reported by Wang and co-workers [*J. Med. Chem.* **1995**, *38*, 3652–3659].

Table 2. In Vitro Ligand Binding Results (Corrected)

compound	K_i (nM \pm SD) [B \pm SD]			μ/δ	μ/κ
	μ	δ	κ		
(2 <i>S</i> ,3 <i>R</i> ,4 <i>S</i>)- 1a (HCl salt)	0.005 \pm 0.002 [1.11 \pm 0.05]	84.06 \pm 10.74 [0.96 \pm 0.10]	41.7 \pm 1.4 [1.03 \pm 0.03]	16812	8340
(2 <i>R</i> ,3 <i>R</i> ,4 <i>S</i>)- 1b (oxalate salt)	0.013 \pm 0.002 [0.94 \pm 0.28]	103.42 \pm 13.20 [0.77 \pm 0.08]	122.2 \pm 7.3 [1.05 \pm 0.06]	7955	9400
(2 <i>R</i> ,3 <i>S</i> ,4 <i>R</i>)- 1c (HCl salt)	47.7 \pm 7.21 [0.98 \pm 0.05]	> 1.5 μ M	> 0.5 μ M	ND	ND
(2 <i>S</i> ,3 <i>S</i> ,4 <i>R</i>)- 1d (oxalate salt)	16.47 \pm 1.07 [0.92 \pm 0.02]	> 3 μ M	> 0.5 μ M	ND	ND

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Jeewoo Lee, Nancy E. Lewin, Peter Acs, Peter M. Blumberg, and Victor E. Marquez*: Conformationally Constrained Analogues of Diacylglycerol. 13.¹ Protein Kinase C Ligands Based on Templates Derived from 2,3-Dideoxy-L-erythro(*threo*)-hexono-1,4-lactone and 2-Deoxyapilactone.

Page 4915. In Scheme 8, compound **28** should go to target **6** (not **4**) and compound **34** should go to target **4** (not **6**).

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