Additions and Corrections

1995, Volume 38

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 coworkers [*J. Med. Chem.* **1995**, *38*, 3652−3659].

Table 2. In Vitro Ligand Binding Results (Corrected)

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

JM970116E

1996, Volume 39

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-Deoxyapiolactone.

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

JM9701435