Bioorganic & Medicinal Chemistry Letters 20 (2010) 1800

Contents lists available at ScienceDirect

Bioorganic & Medicinal Chemistry Letters

journal homepage: www.elsevier.com/locate/bmcl

Corrigendum

Corrigendum to "Discovery and SAR of cyclic isothioureas as novel NPY Y_1 receptor antagonists" [Bioorg. Med. Chem. Lett. 19 (2009) 6801]

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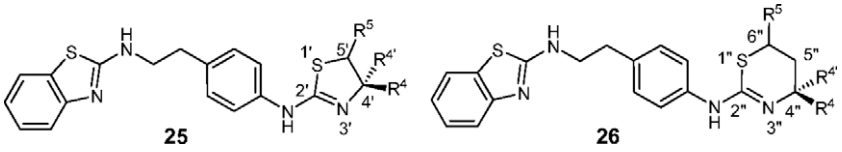
There was an error in Table 2. The corrected table appears below.

Table 2
NPY Y_1 receptor binding affinity of the five- and six-membered cyclic isothiourea analogs with substituted aminothiazole/aminothiazine moieties



Compound	R^4	$R^{4'}$	R^5	K_i[a] (nM)	%Inh[b]
25a	Me	H	H	379	
25b	Et	H	H	153	
25c	*n*-Pr/H[c]		H	482	
25d	*n*-Bu/H[c]		H		69
25e	*i*-Bu	H	H	712	
25f	Ph	H	H	175	
25g	H	Ph	H		0
25h	H	H	Me		61
25i	H	H	OMe	2550	
25j	H	H	Ph		17
26a	Me/H[c]		H	50	
26b	*i*-Pr/H[c]		H	53	
26c	*i*-Bu/H[c]		H	62	
26d	PhEt/H[c]		H	195	
26e	*c*-Hex/H[c]		H	332	
26f	Ph/H[c]		H	55	
26g	*p*-OMePh/H[c]		H	55	
26h	*p*-ClPh/H[c]		H	122	
26i	*m*-ClPh/H[c]		H	107	
26j	*o*-ClPh/H[c]		H	956	
26k	Ph	H	H	30	
26l	H	Ph	H	464	

[a] NPY Y_1 receptor binding affinity K_i values were determined as described in Ref. 18.
[b] Percentage inhibition at 2 μg/mL concentration of testing compounds.
[c] Racemate.

DOI of original article: 10.1016/j.bmcl.2009.09.048
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doi:10.1016/j.bmcl.2010.01.023