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Erratum

Erratum to "Synthesis, biological activity, and X-ray crystal structural analysis of diaryl ether inhibitors of malarial enoyl acyl carrier protein reductase. Part 1: 4′-Substituted triclosan derivatives" [Bioorg. Med. Chem. Lett. 15 (2005) 5247–5252]

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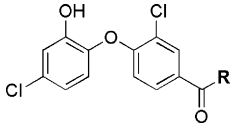
The publisher regrets that the graphics (chemical structures) for Tables 2 and 3 were reversed. The corrected tables appear here.

Table 2. Inhibitory properties of selected 4′-amide derivatives



Compound	R	EC_{50}[a] 3D7/Dd2 (μM)	PfENR[a] IC_{50} (nM)
35	NH_2	96.3 ± 11.7/100 ± 16	120 ± 35
37	N(H)Me	77.8 ± 25.3/100 ± 22	310 ± 66
38	N(H)Bn	7.5 ± 2.8/20.6 ± 3.6	3000 ± 800
39	NMe_2	68.4 ± 13.2/95.7 ± 11.6	460 ± 110
40	*N*-Pyrrolidine	20.7 ± 6.0/42.6 ± 12.8	810 ± 300
41	*N*-Piperidine	11.7 ± 2.2/18.0 ± 3.0	680 ± 120
42	*N*-Morpholine	63.5 ± 14.1/94.5 ± 10.0	200 ± 10

[a] Values reported as means ± standard error.

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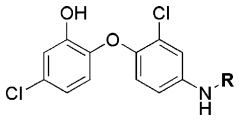
doi:10.1016/j.bmcl.2005.12.054

Table 3. Inhibitory properties of selected 4′-aniline derivatives



Compound	R	EC_{50}[a] 3D7/Dd2 (μM)	PfENR[a] IC_{50} (nM)
30	H	120 ± 21/140 ± 24	360 ± 92
43	Ac	52.2 ± 17.3/50.9 ± 19.2	57 ± 24
44	Bz	57.2 ± 14.7/83.9 ± 15.0	170 ± 16
45	$C(O)NH_2$	110 ± 24/120 ± 22	160 ± 45
46	C(O)*N*-Morpholine	66.8 ± 14.1/86.4 ± 14.9	250 ± 49
47	SO_2Ph	55.1 ± 10.2/77.5 ± 8.0	370 ± 100
48	SO_2(1-Naphthyl)	34.3 ± 5.6/43.7 ± 12.2	590 ± 88
49	SO_2(2-Naphthyl)	23.5 ± 3.2/33.9 ± 9.6	2500 ± 900
50	SO_2CF_3	>120/>120	140 ± 34
51	CH_2Ph	35.8 ± 5.3/47.5 ± 12.2	140 ± 69
52	CH_2(2-CNPh)	26.7 ± 2.9/45.4 ± 6.5	560 ± 120
53	CH_2(2-HOPh)	4.5 ± 0.5/5.6 ± 1.3	320 ± 59

[a] Values reported as means ± standard error.