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Corrigendum

# Corrigendum to "Synthesis and in vitro pharmacological studies of new C(4) modified salvinorin A analogues" [Bioorg. Med. Chem. Lett. 16 (2006) 5498–5502]

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The authors regret that this article contains two errors:
(1) On page 5500, left-hand column, in Scheme 6, "Synthesis of the isostere (**23**) of salvinorin acid" should be changed to "Synthesis of the tetrazole (**23**)".
(2) On page 5502, the NMR data for compound **3** were incorrectly given, and those reported data belong to 8-*epi* **3**. The revised $^{1}H$ and $^{13}C$ NMR data of **3** should be revised as follows: $^{1}H$ NMR ($CDCl_3$, 300 MHz) $\delta$ 9.91 (1H, d, $J = 1.2$ Hz), 7.43 (1H, d, $J = 0.9$ Hz), 7.41 (1H, t, $J = 1.8$ Hz), 6.38 (1H, dd, $J = 0.9$ and 1.8 Hz), 5.54 (1H, dd, $J = 5.4$ and 11.7 Hz), 5.17 (1H, dd, $J = 7.2$ and 11.7 Hz), 2.73 (1H, dd, $J = 3.3$ and 12.9 Hz), 2.51 (1H, dd, $J = 5.1$ and 13.5 Hz), 2.39 (1H, m), 2.10–2.30 (4H, m,), 2.13 (3H, s), 1.55–1.90 (4H, m), 1.46 (3H, s), 1.07 (3H, s); $^{13}C$ NMR ($CDCl_3$, 75 MHz) $\delta$ 202.2, 200.6, 170.9, 170.0, 143.8, 139.4, 125.1, 108.3, 75.2, 72.0, 63.6, 59.7, 51.2, 43.2, 42.5, 38.0, 35.4, 28.3, 20.6, 17.9, 17.8, 15.1.

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