

98.5(C-21), 167.5(C-22), 101.0(C-1), 74.8(C-2), 78.4(C-3), 71.8(C-4), 78.7(C-5), 63.0(C-6)。以上MS、¹H-NMR和¹³C-NMR数据与文献报道一致^[3],故确定化合物为斯垂特萨米碱。

化合物 : 黄色针状结晶(甲醇), ESFMS (*m/z*): 545 [M + H]⁺。¹H-NMR (500 MHz, CD₃OD) : 2.81 (1H, m, H-5a), 3.14 (1H, m, H-5b), 2.80 (2H, m, H-6), 7.47 (1H, d, *J* = 7.9 Hz, H-9), 6.99 (1H, t, *J* = 7.2 Hz, H-10), 7.09 (1H, t, *J* = 7.2 Hz, H-11), 7.32 (1H, d, *J* = 7.9 Hz, H-12), 2.07 (2H, m, H-14), 3.30 ~ 3.33 (1H, m, H-15), 7.56 (1H, s, H-17), 3.02 (1H, m, H-18a), 3.52 (1H, d, *J* = 10.8 Hz, H-18b), 4.92 (1H, d, *J* = 7.4 Hz, H-19), 1.74 (1H, m, H-20), 5.84 (1H, d, *J* = 9.3 Hz, H-21), 3.64 (3H, s, H-OCH₃), 4.78 (1H, d, *J* = 11.0 Hz, H-1), 3.33 ~ 3.42 (4H, m, H-2, 3, 4, 5), 3.62 (1H, m, H-6a), 3.85 (1H, d, *J* = 2.1 Hz, H-6b)。¹³C-NMR (125 MHz, CD₃OD) : 133.6(C-2), 93.3(C-3), 54.1(C-5), 23.2(C-6), 111.8(C-7), 127.3(C-8), 120.2(C-9), 120.5(C-10), 123.7(C-11), 112.9(C-12), 138.9(C-13), 43.4(C-14), 27.1(C-15), 111.8(C-16), 154.7(C-17), 59.8(C-18), 74.8(C-19), 41.4(C-20), 98.0(C-21), 169.3(C-22), 101.8(C-1), 75.0(C-2), 78.2(C-3), 71.9(C-4), 78.6(C-5), 63.1(C-6), 52.0(C-OCH₃)。以上MS、¹H-NMR和¹³C-NMR数据与文献报道一致^[4],故确定化合物为卡丹宾碱。

化合物 : 黄色粉末, ESFMS (*m/z*): 547 [M + H]⁺。¹H-NMR (500 MHz, CD₃OD) : 7.53 (1H, s, H-17), 7.36 (1H, d, *J* = 7.8 Hz, H-9), 7.28 (1H, d, *J* = 8.1 Hz, H-12), 7.03 (1H, t, *J* = 7.9, 8.1 Hz, H-11), 6.96 (1H, t, *J* = 7.8, 8.0 Hz, H-10), 5.54 (1H, d, *J* = 9.1 Hz, H-21), 4.76 (1H, d,

J = 8.0 Hz, H-1), 4.20 (1H, m, H-19), 3.81 (1H, m, H-6a), 3.80 (1H, m, H-3), 3.78 (3H, s, H-OCH₃), 3.61 (1H, m, H-6b), 3.23 (1H, m, H-18a), 3.03 (1H, m, H-15), 2.96 (1H, m, H-18b), 2.32 (1H, d, *J* = 14.5 Hz, H-14a), 2.00 (1H, m, H-20), 1.76 (1H, m, H-14b)。¹³C-NMR (125 MHz, CD₃OD) : 136.0(C-2), 66.7(C-3), 51.3(C-5), 23.5(C-6), 111.5(C-7), 128.5(C-8), 118.9(C-9), 122.4(C-10), 120.1(C-11), 112.4(C-12), 138.7(C-13), 37.5(C-14), 34.3(C-15), 112.4(C-16), 153.9(C-17), 59.6(C-18), 64.5(C-19), 44.8(C-20), 96.8(C-21), 169.3(C-22), 101.3(C-1), 74.9(C-2), 78.2(C-3), 71.4(C-4), 78.6(C-5), 63.1(C-6), 52.4(C-OCH₃)。以上MS、¹H-NMR和¹³C-NMR数据与文献报道一致^[4],故确定化合物为二氢卡丹宾碱。

化合物 ~ 的理化性质和波谱数据与文献对照, 分别鉴定为缝籽嗪甲醚^[5]、柯诺辛^[6]、去氢钩藤碱^[7]、异去氢钩藤碱^[7]、钩藤碱^[6]和异钩藤碱^[7]。

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从2008年第1期开始本刊所刊用文章文后的参考文献使用原语种撰写,按照国家标准《文后参考文献著录规则》(GB/T 7714-2005)书写。具体参考文献书写示范例见本刊2009年第40卷第1期上刊登的“《中草药》杂志2009年投稿须知”。