

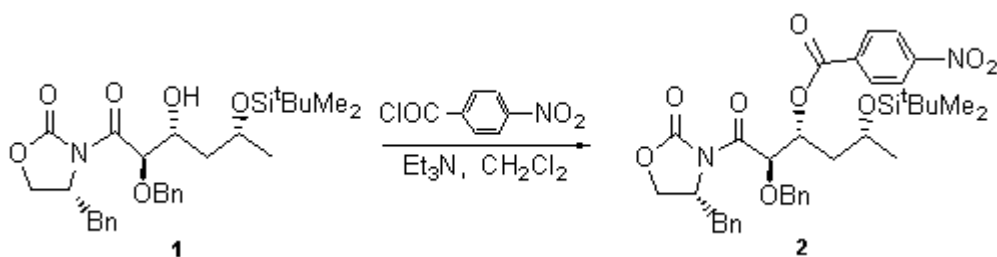
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3-[5-(*tert*-Butyldimethylsilyloxy)-3-*p*-nitrobenzoyl-1-oxo-2-(phenylmethoxy)hexyl]-4-(phenylmethyl)-2-oxazolidinone

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A mixture of alcohol **1** (121 mg, 0.23 mmol) [1], triethylamine (48 l, 0.34 mmol) and *p*-nitrobenzoyl chloride (51 mg, 0.28 mmol) in dichloromethane (2 ml) was stirred for 2 h. at 30°C. The reaction mixture was poured into saturated aqueous sodium hydrogen carbonate (5 ml), extracted into ethyl acetate (2 x 10 ml), washed with water (2 x 5 ml) and dried over sodium sulfate. Removal of the solvent under reduced pressure and purification of the residue by flash chromatography, using light petroleum-ethyl acetate (8:2) as eluent afforded the title compound **2** (129 mg, 83%) as a colourless oil.

$[\alpha]_D -41.34$ (c 1.024, CHCl_3).

IR (cm^{-1} , neat): 1783s, 1729s, 1714s, 1386m, 1103m.

^1H NMR (400 MHz, CDCl_3): 0.01, 0.02 (6H, s, SiMe_2), 0.85 (9H, s, Bu^t), 1.21 (3H, d, $J_{6',5'} 6.0$ Hz, $\text{H6}'$), 2.10-2.14 (2H, m, $\text{H4}'$), 2.62 (1H, dd, $J_{\text{gem}} 13.5$ and $J 9.5$ Hz, CHCH^APh), 3.15 (1H, dd, $J_{\text{gem}} 13.5$ and $J 3.3$ Hz, CHCH^BPh), 3.76 (1H, dd, $J_{\text{gem}} 8.8$ and $J_{5A,4} 8.8$ Hz, H5^A), 3.95-4.00 (1H, m, $\text{H5}'$), 4.06 (1H, dd, $J_{\text{gem}} 8.8$ and $J_{5B,4} 2.5$ Hz, H5^B), 4.47-4.51 (1H, m, H4), 4.70 (2H, s, OCH_2Ph), 5.55-5.60 (2H, m, $\text{H2}'$, $\text{H3}'$), 7.15-7.42 (10H, m, Ph), 8.18 (2H, d, $J 8.8$ Hz, PhNO_2), 8.27 (2H, d, $J 8.8$ Hz, PhNO_2).

^{13}C NMR (100 MHz, CDCl_3): -4.1, -3.9 (CH_3 , SiMe_2), 18.8 (quat., CMe_3), 24.0 (CH_3 , $\text{C6}'$), 25.5 (CH_3 , CMe_3), 38.3 (CH_2 , CHCH_2Ph), 41.5 (CH_2 , $\text{C4}'$), 56.2 (CH , C4), 66.3 (CH_2 , C5), 67.1 (CH , $\text{C5}'$), 73.5 (CH , $\text{C3}'$), 74.2 (CH_2 , OCH_2Ph), 78.5 (CH , $\text{C2}'$), 124.2, 124.3, 128.1, 128.9, 129.2, 129.6, 130.1, 131.3, 131.5 [CH , 3 x Ph (last 5 peaks coincidental)], 135.5 (quat., CHCH_2Ph), 135.7 (quat., $\text{OC}=\text{OC}$), 137.6 (quat., OCH_2Ph), 151.4 (quat., CNO_2), 153.6 (quat., C2) 164.0 (quat., $\text{OC}=\text{O}$), 170.7 (quat., $\text{C1}'$).

CI-MS: (FAB, NBA matrix) 677 (MH^+ , 2%), 619 (4), 569 ($\text{MH}^+ - \text{HOCH}_2\text{Ph}$, 1), 545 ($\text{MH}^+ - \text{HOSiMe}_2\text{Bu}^t$, 4), 224 (5), 178 ($\text{C}_{10}\text{H}_{12}\text{NO}_2$, 6), 159 ($\text{C}_8\text{H}_{19}\text{OSi}$, 9), 150 (15), 136 (14), 91 (CH_2Ph , 100), 73 (27).

Anal. calc. for $\text{C}_{36}\text{H}_{44}\text{N}_2\text{O}_9\text{Si}$ MH^+ (CI, NH_3), 677.2901; found MH^+ , 677.2894.

Reference

- Brimble, M. A.; Park, J. S. O. *J. Chem. Soc. Perkin Trans. I* **2000**, 697-709.

Sample availability: available from the authors.

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