

(+) *N*-Substituted benzyl-6-chloro-2,3,4,9-tetrahydrofuro[2,3-*b*]-quinolin-3,4-dione (139-153) 之合成

N-Benzyl-6-chloro-2,3,4,9-tetrahydrofuro[2,3-*b*]-quinolin-3,4-dione (139) 之合成

取化合物 39(2.35g , 0.01mole)懸著於 DMF 30 ml 中 , 加入無水 K_2CO_3 (1.38 g , 0.01 mole)加熱(約 50~60)使之溶解 , 加入 benzyl chloride(11.3g , 0.1mole), 反應 1 小時後加入冰水中 , 以 $CHCl_3$ 萃取 , 取 $CHCl_3$ 層 , 以無水 $MgSO_4$ 乾燥 , 減壓濃縮後 , 收集沉澱物以短程矽膠管柱層析($CHCl_3/EtOH$)沖提 , 再以 MeOH 及 $CHCl_3$ 做再結晶 , 得白色棉絮狀結晶 , 為化合物 139(2.35 g, 72.31 %), mp : 222-225 。光譜數據如下 : MS m/z : 325; IR (KBr) cm^{-1} : 1721.0 ($C_3=O$), 1605.3 ($C_4=O$); UV λ_{max} nm (MeOH) (log ε): 222 (4.79); 1H -NMR ($DMSO-d_6$) δ: 4.95 (2H, s, H-2), 5.57(2H, s, H-10), 7.28-7.37(5H,m, Ar-H) 7.65(1H, d, J=9.0 Hz, H-8), 7.74 (1H, dd, J=9.0Hz, 2.5Hz, H-7), 8.07 (1H, d, J=2.4Hz, H-5); ^{13}C -NMR ($DMSO-d_6$) δ: 46.70(C-10), 76.48 (C-2), 100.69(C-3a), 119.60 (C-8), 125.89 (C-5), 126.87(C-12, C-16), 128.10(C-6), 128.21(C-14), 129.15(C-13, C-15), 129.97 (C-4a), 133.20(C-7), 134.80(C-11), 137.12(C-8a), 170.19(C-9a), 174.81 (C-4), 191.15 (C-3) .

N-*o*-Methylbenzyl-6-chloro-2,3,4,9-tetrahydrofuro[2,3-*b*]quinolin-3,4-dione(140) 之合成

取化合物 39(2.35g , 0.01mole)和 *o*-methylbenzyl chloride (12.7g , 0.1mole)為原料 , 比照化合物 139 的合成法及處理步驟 , 得化合物 140(1.96g , 57.82 %), mp : 278~283 。光譜數據如下: MS m/z : 339; IR (KBr) cm^{-1} : 1728.8 ($C_3=O$), 1605.3 ($C_4=O$); UV λ_{max} nm (MeOH) (log ε): 220 (4.86); 1H -NMR ($DMSO-d_6$) δ: 2.44 (3H, s, $C_{12}-CH_3$), 4.90 (2H, s, H-2), 5.51 (2H, s, H-10), 6.67 (1H, d, J=7.5 Hz, H-13), 7.04-7.26 (3H, m, Ar-H) 7.46(1H, d, J=9.0 Hz, H-8), 7.74(1H, dd, J=9.0 Hz, 2.6 Hz, H-7), 8.13 (1H,d,J=2.6Hz,H-5); ^{13}C -NMR ($DMSO-d_6$) δ: 18.90($C_{12}-CH_3$), 45.24 (C-10), 76.45 (C-2), 100.75 (C-3a), 119.71 (C-8), 124.20 (C-6), 125.91 (C-4a), 126.59 (C-15), 127.67 (C-14), 128.21 (C-13), 130.02 (C-16), 130.71 (C-5), 132.46 (C-7), 133.31 (C-11), 135.38 (C-12), 137.33 (C-8a), 170.29 (C-9a), 174.99 (C-4), 191.05 (C-3) .

N-*m*-Methylbenzyl-6-chloro-2,3,4,9-tetrahydrofuro[2,3-*b*]quinolin-3,4-dione (141) 之合成

取化合物 39(2.35g , 0.01mole)和 *m*-methylbenzyl chloride (12.7g , 0.1mole)為原料 , 比照化合物 139 的合成法及處理步驟 , 得化合物 141(1.98g , 58.41 %), mp:242-245 。光譜數據如下:MS m/z : 339; IR (KBr) cm^{-1} : 1713.3($C_3=O$), 1605.3 ($C_4=O$); UV λ_{max} nm (MeOH) (log ε): 220 (4.84); 1H -NMR ($DMSO-d_6$) δ: 2.25 (3H, s, $C_{13}-CH_3$), 4.95(2H, s, H-2), 5.52(2H, s, H-10), 7.07-7.23 (4H, m, Ar-H), 7.63 (1H, d, J=9.0Hz, H-8), 7.71 (1H, dd, J=9.0Hz, 2.6Hz, H-7), 8.07 (1H, d, J=2.4Hz, H-5); ^{13}C -NMR ($DMSO-d_6$) δ: 21.17 ($C_{13}-CH_3$), 46.72 (C-10), 76.46 (C-2), 100.70

(C-3a), 119.60 (C-8), 123.86 (C-6), 125.88 (C-4a), 127.25 (C-16), 128.21 (C-14), 128.79 (C-15), 129.05 (C-12), 129.96 (C-5), 133.21 (C-7), 134.71 (C-11), 137.17 (C-13), 138.51 (C-8a), 170.22 (C-9a), 174.81 (C-4), 191.17 (C-3).

***N-p*-Methylbenzyl-6-chloro-2,3,4,9-tetrahydrofuro[2,3-*b*]quinolin-3,4-dione(142)之合成**

取化合物 39 (2.35g , 0.01mole) 和 *p*-methylbenzyl chloride (12.7g , 0.1mole) 為原料 , 比照化合物 139 的合成法及處理步驟 , 得化合物 142 (2.62g , 77.29 %), mp:256~259 。光譜數據如下 :MS *m/z*: 339; IR (KBr) cm⁻¹: 1721.0 (C₃=O), 1605.3 (C₄=O); UV λ_{max} nm (MeOH) (log ε): 220 (4.87); ¹H-NMR (DMSO-*d*₆) δ: 2.24 (3H, s, C₁₄-CH₃), 4.95 (2H, s, H-2), 5.52 (2H, s, H-10), 7.14 (2H, d, J=8.2Hz, H-13, H-15), 7.24 (2H, d, J=8.1Hz, H-12, H-16), 7.65 (1H, d, J=9.0Hz, H-8), 7.73 (1H, dd, J=9.0Hz, 1.8Hz, H-7), 8.07 (1H, d, J=2.4Hz, H-5); ¹³C-NMR (DMSO-*d*₆) δ: 20.85 (C₁₄-CH₃), 46.52 (C-10), 76.46 (C-2), 100.66 (C-3a), 119.65 (C-8), 125.87 (C-6), 126.89 (C-13, C-15), 128.22 (C-4), 129.69 (C-12, C-16), 129.95 (C-5), 131.74 (C-11), 133.16 (C-7), 137.11 (C-14), 137.43 (C-8a), 170.18 (C-9a), 174.77 (C-4), 191.13 (C-3).

***N-m*-Methoxybenzyl-6-chloro-2,3,4,9-tetrahydrofuro[2,3-*b*]quinolin-3,4-dione (143) 之合成**

取化合物 39 (2.35g , 0.01mole) 和 *m*-methoxybenzyl chloride (14.3g , 0.1mole) 為原料 , 比照化合物 139 的合成法及處理步驟 , 得化合物 143 (2.35g , 66.20 %), mp:229~232 。光譜數據如下 :MS *m/z*: 355; IR (KBr) cm⁻¹: 1713.3 (C₃=O), 1605.3 (C₄=O); UV λ_{max} nm (MeOH) (log ε): 222 (log =4.70); ¹H-NMR (DMSO-*d*₆) δ: 3.71 (3H, s, C₁₃-OCH₃), 4.95 (2H, s, H-2), 5.53 (2H, s, H-10), 6.83-6.88 (2H, m, H-14, H-16), 6.94 (1H, d, J=1.8Hz, H-12), 7.25 (1H, t, J=8.0Hz, H-15), 7.65 (1H, d, J=9.0Hz, H-8), 7.75 (1H, dd, J=9.0Hz, 2.5Hz, H-7), 8.08 (1H, d, J=2.4Hz, H-5); ¹³C-NMR (DMSO-*d*₆) δ: 46.65 (C-10), 55.32 (C₁₃-OCH₃), 76.48 (C-2), 100.69 (C-3a), 112.96 (C-14), 113.20 (C-8), 118.67 (C-12), 119.60 (C-16), 125.88 (C-6), 128.20 (C-4a), 129.98 (C-15), 130.38 (C-5), 133.21 (C-7), 136.36 (C-11), 137.18 (C-8a), 159.84 (C-13), 170.21 (C-9a), 174.81 (C-4), 191.15 (C-3).

***N-p*-Methoxybenzyl-6-chloro-2,3,4,9-tetrahydrofuro[2,3-*b*]quinolin-3,4-dione (144) 之合成**

取化合物 39 (2.35g , 0.01mole) 和 *p*-methoxybenzyl chloride (14.3g , 0.1mole) 為原料 , 比照化合物 139 的合成法及處理步驟 , 得化合物 144 (2.75g , 77.46 %), mp:262~265 。光譜數據如下 :MS *m/z*: 355; IR (KBr) cm⁻¹: 1728.8 (C₃=O), 1605.3 (C₄=O); UV λ_{max} nm (MeOH) (log ε): 222 (4.91); ¹H-NMR (DMSO-*d*₆) δ: 3.70 (3H, s, C₁₄-OCH₃), 4.95 (2H, s, H-2), 5.49 (2H, s, H-10), 6.89 (2H, d, J=8.7Hz, H-12, H-16), 7.31 (2H, d, J=8.7 Hz, H-13, H-15), 7.72-7.73 (2H, m, H-7, H-8), 8.07 (1H, d, J=2.0 Hz, H-5); ¹³C-NMR (DMSO-*d*₆) δ: 46.21 (C-10), 55.32 (C₁₄-OCH₃), 76.44

(C-2), 100.68(C-3a), 114.50(C-13, C-15), 119.69(C-8), 125.87(C-6), 126.55 (C-4a), 128.23 (C-11), 128.47 (C-12, C-16), 129.94 (C-5), 133.15 (C-7), 137.06 (C-8a), 159.10 (C-14), 170.17 (C-9a), 174.71 (C-4), 191.16 (C-3).

N-o-Fluorobenzyl-6-chloro-2,3,4,9-tetrahydrofuro[2,3-*b*]quinolin-3,4-dione (145) 之合成

取化合物 **39** (2.35g , 0.01mole) 和 *o*-fluorobenzyl chloride (13.1g , 0.1mole) 為原料 , 比照化合物 **139** 的合成法及處理得化合物 **145** (2.05g , 59.77 %), mp : 247-250 。光譜數據如下 : MS *m/z*: 343; IR (KBr) cm⁻¹: 1721.0 (C₃=O), 1605.3 (C₄=O); UV λ_{max} nm (MeOH) (log ε): 221 (4.82); ¹H-NMR (DMSO-*d*₆) δ: 4.94 (2H, s, H-2), 5.61 (2H, s, H-10), 7.12-7.33 (4H, m, Ar-H), 7.63 (1H, d, J=9.0Hz, H-8), 7.78 (1H, dd, J=9.0Hz, 2.6Hz, H-7), 8.10 (1H, d, J=2.6Hz, H-5); ¹³C-NMR (DMSO-*d*₆) δ: 41.52 (C-10), 76.53 (C-2), 100.80 (C-3a), 115.76 (C-8), 116.17 (C-13), 119.23 (C-6), 121.58 (C-11), 121.86 (C-4a), 125.21 (C-15), 125.98 (C-14), 128.19 (C-5), 128.46 (C-16), 130.27 (C-7), 133.35 (C-8a), 137.11 (C-12), 170.21 (C-9a), 175.01 (C-4), 191.03 (C-3).

N-m-Fluorobenzyl-6-chloro-2,3,4,9-tetrahydrofuro[2,3-*b*]quinolin-3,4-dione(146) 之合成

取化合物 **39** (2.35g , 0.01mole) 和 *m*-fluorobenzyl chloride (13.1g , 0.1mole) 為原料 , 比照化合物 **139** 的合成法及處理步得化合物 **146** (2.12g , 61.81 %), mp:233~235 。光譜數據如下 : MS *m/z*: 343; IR (KBr) cm⁻¹: 1721.0(C₃=O), 1605.3 (C₄=O); UV λ_{max} nm (MeOH) (log ε): 222 (log =4.74); ¹H-NMR (DMSO-*d*₆) δ: 4.94 (2H, s, H-2), 5.59 (2H, s, H-10), 7.13-7.41 (4H, m, Ar-H), 7.63 (1H, d, J=9.0Hz, H-8), 7.75 (1H, dd, J=9.0Hz, 2.6Hz, H-7), 8.08 (1H, d, J=2.5 Hz, H-5); ¹³C-NMR (DMSO-*d*₆) δ: 46.23(C-10), 76.51(C-2), 100.83(C-3a), 113.79(C-14), 114.52 (C-8), 115.22 (C-12), 119.45 (C-6), 122.91 (C-4a), 125.93 (C-16), 128.25 (C-15), 130.01 (C-5), 131.21 (C-7), 133.24 (C-11), 137.06 (C-8a), 137.73 (C-13), 170.24 (C-9a), 174.87 (C-4), 191.19 (C-3).

N-p-Fluorobenzyl-6-chloro-2,3,4,9-tetrahydrofuro[2,3-*b*]quinolin-3,4-dione (147) 之合成

取化合物 **39** (2.35g , 0.01mole) 和 4-fluorobenzyl chloride (13.1g , 0.1mole) 為原料 , 比照化合物 **139** 的合成法及處理步驟 , 得化合物 **147** (2.71g , 79.00 %), mp:232~234 。光譜數據如下 : MS *m/z*: 353; IR (KBr) cm⁻¹: 1721.0(C₃=O), 1605.3 (C₄=O); UV λ_{max} nm (MeOH) (log ε): 220 (4.90); ¹H-NMR (DMSO-*d*₆) δ: 4.94 (2H, s, H-2), 5.56 (2H, s, H-10), 7.13-7.21 (2H,m,H-13,H-15), 7.40-7.47 (2H,m,H-12,H-16), 7.67 (1H, d, J=9.0Hz, H-8), 7.44 (1H, dd, J=9.0Hz,2.5Hz,H-7), 8.06 (1H,d,J=2.3Hz,H=5); ¹³C-NMR (DMSO-*d*₆) δ: 46.04 (C-10), 76.48 (C-2), 100.76 (C-3a), 115.73 (C-13), 116.16 (C-15), 119.53 (C-8), 125.92 (C-6), 128.25 (C-4a), 129.11 (C-12), 129.28 (C-16), 130.00 (C-5), 130.97 (C-7, C-11), 133.20 (C-8a), 137.02 (C-14), 170.18 (C-9a), 174.80 (C-4), 191.14 (C-3).

*N-o-Chlorobenzyl-6-chloro-2,3,4,9-tetrahydrofuro[2,3-*b*]quinolin-3,4-dione(148)之合成*

取化合物 **39** (2.35g , 0.01mole) 和 *o*-chlorobenzyl chloride (14.7g , 0.1mole) 為原料 , 比照化合物 **139** 的合成法及處理步驟 , 得化合物 **148** (2.22g , 61.84 %) , mp:270-275 。光譜數據如下 :MS *m/z*: 359; IR (KBr) cm⁻¹: 1728.8(C₃=O), 1605.3 (C₄=O); UV λ_{max} nm (MeOH) (log ε): 220 (4.75); ¹H-NMR (DMSO-*d*₆) δ: 4.92 (2H,s,H-2), 5.58 (2H,s,H-10), 7.03 (1H, d, J=1.4Hz, H-16), 7.23-7.36 (2H,m,H-14,H-15), 7.49 (1H,d,J=9.0Hz,H-8), 7.58 (1H,d,J=9.0Hz,H-13), 7.77 (1H,dd, J=9.0Hz,2.6Hz,H-7), 8.13 (1H, d, J=2.6Hz, H-5); ¹³C-NMR (DMSO-*d*₆) δ: 45.13(C-10), 76.57 (C-2), 100.89 (C-3a), 119.30 (C-8), 126.00 (C-6), 127.29 (C-4a), 128.13 (C-14, C-15), 129.84 (C-13), 130.04 (C-5, C-16), 131.66 (C-7), 131.85 (C-12), 133.45 (C-11), 137.23 (C-8a), 170.30 (C-9a), 175.05 (C-4), 191.04 (C-3) .

*N-m-Chlorobenzyl-6-chloro-2,3,4,9-tetrahydrofuro-[2,3-*b*]quinolin-3,4-dione (149) 之合成*

取化合物 **39** (2.35g , 0.01mole) 和 *m*-chlorobenzyl chloride (14.7g , 0.1mole) 為原料 , 比照化合物 **139** 的合成法及處理步驟 , 得化合物 **149** (2.35g , 65.46%) , mp:245~248 。光譜數據如下 :MS *m/z*: 359; IR (KBr) cm⁻¹: 1721.0(C₃=O), 1605.3 (C₄=O); UV λ_{max} nm (MeOH) (log ε): 221 (4.67); ¹H-NMR (DMSO-*d*₆) δ: 4.94 (2H, s, H-2), 5.58 (2H, s, H-10), 7.29-7.38 (3H, m, H-14, H-15, H-16), 7.51 (1H, s, H12), 7.63 (1H, d, J=9.0Hz, H-8), 7.76 (1H, dd, J=9.0Hz, 2.6Hz, H=7), 8.09 (1H,d, J=2.6Hz,H-5); ¹³C-NMR (DMSO-*d*₆) δ: 46.15 (C-10), 76.52 (C-2), 100.87 (C-3a), 119.46 (C-8), 125.46 (C-6), 125.96 (C-4a), 126.86 (C-16), 128.17 (C-12, C-14), 130.04 (C-15), 131.03 (C-5), 133.28 (C-13), 133.83 (C-7), 137.06 (C-11), 137.40 (C-8a), 170.26 (C-9a), 174.91 (C-4) 191.21 (C-3) .

*N-p-Chlorobenzyl-6-chloro-2,3,4,9-tetrahydrofuro[2,3-*b*]quinolin-3,4-dione(150) 之合成*

取化合物 **39** (2.35g , 0.01mole) 和 *p*-chlorobenzyl chloride (14.7g , 0.1mole) 為原料 , 比照化合物 **139** 的合成法及處理步驟 , 得化合物 **150** (2.85g , 79.39%) , mp:253-255 。光譜數據如下 :MS *m/z*: 359; IR (KBr) cm⁻¹: 1721.0(C₃=O), 1605.3 (C₄=O); UV λ_{max} nm (MeOH) (log ε): 222(4.87); ¹H-NMR (DMSO-*d*₆) δ: 4.94 (2H, s, H-2), 5.57 (2H, s, H-10), 7.39 (4H, s, Ar-H), 7.64 (1H, d, J=9.0Hz, H-8), 7.75 (1H, dd, J=9.0Hz, 2.6Hz,H=7), 8.08 (1H, d, J=2.4Hz, H-5); ¹³C-NMR (DMSO-*d*₆) δ: 46.09 (C-10), 76.50 (C-2), 100.77 (C-3a), 119.51 (C-8), 125.94 (C₆), 128.24 (C-4a), 128.90 (C-13, C-15), 129.08 (C-12, C-16), 130.03 (C-5), 132.76 (C-14), 133.25 (C-7), 133.88 (C-11), 137.03 (C-8a), 170.21 (C-9a), 174.84 (C-4), 191.14 (C-3) .

*N-o-Nitrobenzyl-6-chloro-2,3,4,9-tetrahydrofuro[2,3-*b*]quinolin-3,4-dione (151) 之合成*

取化合物 **39** (2.35g , 0.01mole) 和 *o*-nitrobenzyl chloride (15.8g , 0.1mole) 為原料 , 比照化合物 **139** 的合成法及處理步驟 , 得化合物 **151**(2.11g , 57.03%), mp:280-285 。光譜數據如下:MS *m/z*: 370; IR (KBr) cm⁻¹: 1721.0(C₃=O), 1605.3 (C₄=O); UV λ_{max} nm (MeOH) (log ε): 221(4.81); ¹H-NMR (DMSO-*d*₆) δ: 4.89 (2H, s, H-2) , 5.93 (2H, s, H-10) , 7.11 (1H, m, H-16), 7.58-7.63 (2H, m, H-14, H-15) , 7.67(2H, d, J=1.3Hz, H=7, H-8), 8.12(1H,s,H-5),8.27(1H,m,H-13); ¹³C-NMR (DMSO-*d*₆) δ: 45.31(C-10) , 76.55(C-2) , 100.97(C-3a) , 119.77(C-8) , 125.92(C-6) , 126.02 (C-13) , 127.19 (C-4a) , 128.25 (C-14) , 129.34 (C-16) , 130.18 (C-5) , 130.25 (C-11) , 133.26 (C-7) , 135.03 (C-15) , 137.26 (C-8a) , 147.32 (C-12) , 170.44 (C-9a) , 175.25 (C-4) , 191.04 (C-3) .

*N-m-Nitrobenzyl-6-chloro-2,3,4,9-tetrahydrofuro[2,3-*b*]quinolin-3,4-dione (152) 之合成*

取化合物 **39** (2.35g , 0.01mole) 和 *m*-nitrobenzyl chloride (15.8g , 0.1mole) 為原料 , 比照化合物 **139** 的合成法及處理步驟 , 得化合物 **152**(2.32g , 62.70 %) , mp:294-298 。光譜數據如下:MS *m/z*: 370; IR (KBr) cm⁻¹: 1713.3(C₃=O), 1605.3 (C₄=O); UV λ_{max} nm (MeOH) (log ε): 219 (log =4.71); ¹H-NMR (DMSO-*d*₆) δ: 4.94 (2H, s, H-2) , 5.74 (2H, s, H-10) , 7.61 (1H, m, H-16), 7.71~7.75 (3H, m, H-7, H-8, H-15), 8.11 (1H, d, J=2.3Hz, H-5), 8.15 (1H, d, J=1.1Hz, H-14) , 8.31 (1H, d, J=7.1Hz, H-12); ¹³C-NMR (DMSO-*d*₆) δ: 46.04(C-1O), 76.56(C-2) , 100.92(C-3a) , 119.43 (C-8) , 122.16 (C-14) , 123.07 (C-6) , 126.02 (C-4a) , 128.30 (C-12) , 130.11 (C-15) , 130.74 (C-5) , 133.21 (C-7) , 133.34 (C-16) , 137.03 (C-11) , 137.23 (C-8a) , 148.30 (C-13) , 170.26 (C-9a) , 175.02 (C-4) , 191.14 (C-3) .

*N-p-Nitrobenzyl-6-chloro-2,3,4,9-tetrahydrofuro[2,3-*b*]quinolin-3,4-dione (153) 之合成*

取化合物 **39** (2.35g , 0.01mole) 和 *p*-nitrobenzyl chloride (15.8g , 0.1mole) 為原料 , 比照化合物 **139** 的合成法及處理步驟得化合物 **153** (2.78g , 75.14 %) , mp : >300 。光譜數據如下:MS *m/z*: 370; IR (KBr) cm⁻¹: 1721.0(C₃=O), 1605.3 (C₄=O); UV λ_{max} nm (MeOH) (log ε): 219(4.86); ¹H-NMR (DMSO-*d*₆) δ: 4.94 (2H, s, H-2) , 5.74 (2H, s, H-10) , 7.60-7.64 (3H, m, H-8, H-12, H-16), 7.75 (1H, dd, J=9.0Hz, 2.6Hz, H-7) , 8.11(1H,d,J=2.5Hz,H-5),8.18(2H,d,J=8.7Hz,H-13,H-15); ¹³C-NMR (DMSO-*d*₆) δ: 46.29 (C-10) , 76.56 (C-2) , 100.87 (C-3a) , 119.38 (C-8) , 124.15 (C-13, C-15), 126.03(C-6), 128.14(C-12, C-16), 128.28(C-4a), 130.14(C-5), 133.32 (C-7) , 137.04 (C-11) , 142.60 (C-8a) , 147.35 (C-14) , 170.26 (C-9a) , 174.97 (C-4), 191.08 (C-3) .