

Supporting Information

Transition Metal-free Regioselective Synthesis of 2-Arylquinolines, Benzo[*h*]quinolines, Pyrido[2,3-*c*]carbazoles and Estimation of Their Photophysical Properties

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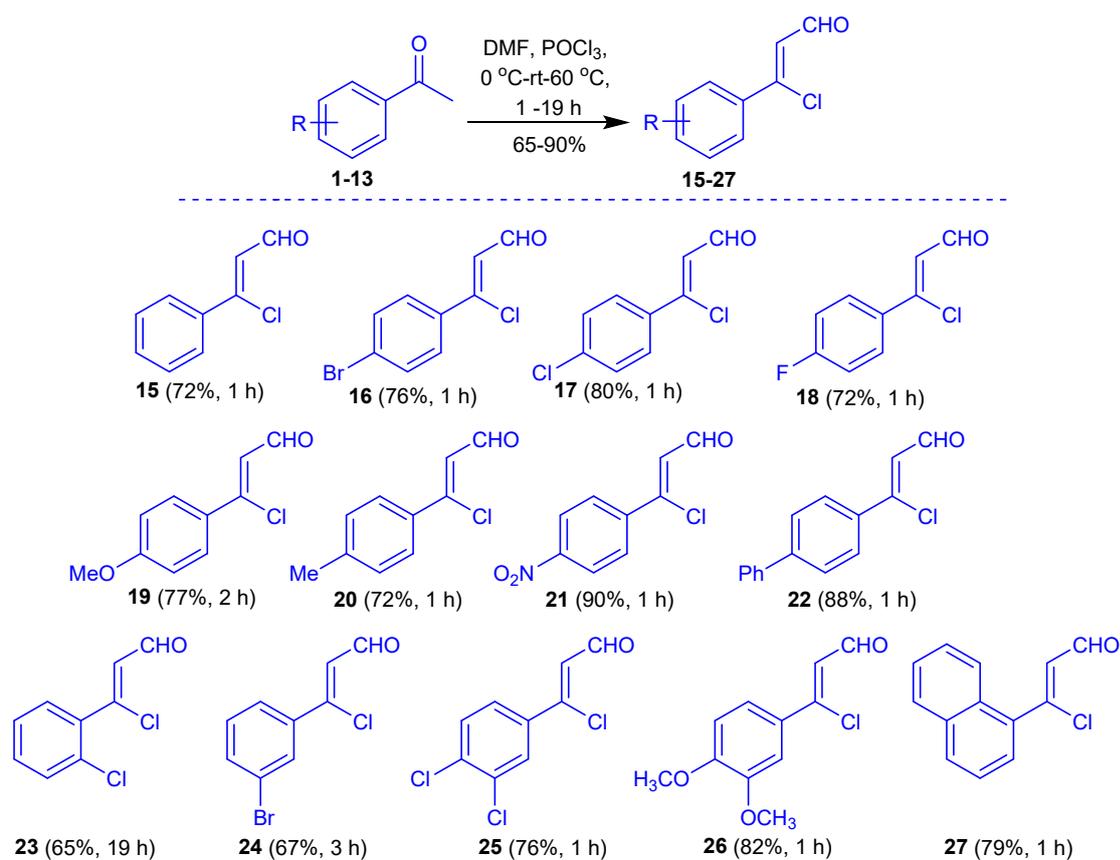
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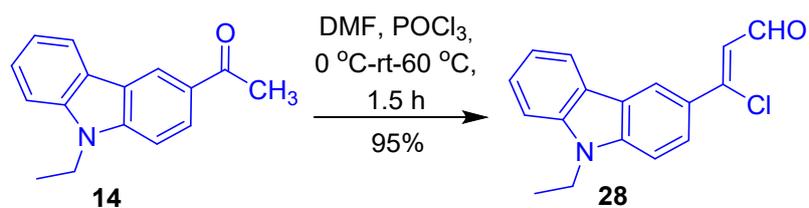
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1. Synthesis of β -Chlorovinyl Aldehydes from various Acetophenones



Scheme S1. Synthesis of β -Chlorovinyl Aldehydes from various Acetophenones



Scheme S2. Synthesis of (Z)-3-chloro-3-(9-ethyl-9H-carbazol-3-yl)acrylaldehyde 28 from 3-acetyl carbazole (14)

2. Experimental data

6,7-Dimethoxy-4-phenylquinoline (15a): Yield = 95% (0.223 g from 0.10 g) as a yellow solid; m.p. 170-172 °C; $R_f = 0.72$ (hexane/EtOAc, 90:10, v/v); IR (Neat): $\nu_{\max} (\text{cm})^{-1} = 1620$ (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 4.03$ (s, 3H, -OCH₃), 4.06 (s, 3H, -OCH₃) 7.06 (s, 1H, ArH), 7.41-7.44 (m, 1H, ArH), 7.50-7.52 (m, 3H, ArH), 7.73 (d, $J = 8.4$ Hz, 1H, ArH), 8.15 (d, $J = 2.4$ Hz, 1H ArH), 8.09-8.11 (m, 2H, ArH) ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 56.2, 56.3, 105.0, 108.5, 117.4, 122.8, 127.4, 128.9, 128.9, 135.0, 140.1, 145.4, 149.8, 152.7, 155.5$ ppm; HRMS (ESI) m/z: calcd. for $\text{C}_{17}\text{H}_{15}\text{NO}_2$ [M + H⁺]: 266.1176, found: 266.1180.

2-(4-Bromophenyl)-6,7-dimethoxyquinoline (16a): Yield = 78% (0.13 g from 0.10 g) as a dark yellow solid; m.p. 192-194 °C; $R_f = 0.25$ (hexane/EtOAc, 95:5, v/v); IR (Neat): $\nu_{\max} (\text{cm})^{-1} = 1615$ (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 4.03$ (s, 3H, -OCH₃), 4.06 (s, 3H, -OCH₃), 7.15 (s, 1H, ArH), 7.57 (s, 1H, ArH), 7.62-7.63 (m, 2H, ArH), 7.78 (d, $J = 8.4$ Hz, 1H, ArH), 7.98-8.01 (m, 2H, ArH), 8.06 (d, $J = 8.4$ Hz, 1H, ArH) ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 56.2, 56.3, 105.0, 108.4, 116.9, 123.0, 123.42, 128.9, 132.0, 135.2, 138.9, 150.0, 152.9, 154.1$ ppm; HRMS (ESI) m/z: calcd. for $\text{C}_{17}\text{H}_{14}\text{BrNO}_2$ [M + H⁺]: 344.0281, found: 344.0284.

2-(4-Chlorophenyl)-6,7-dimethoxyquinoline (17a): Yield = 93% (0.140 g from 0.10 g) as a white solid; m.p. 150-152 °C; $R_f = 0.54$ (hexane/EtOAc, 80:20, v/v); IR (Neat): $\nu_{\max} (\text{cm})^{-1} = 1625$ (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 4.03$ (s, 3H, -OCH₃), 4.06 (s, 3H, -OCH₃), 7.07 (s, 1H, ArH), 7.47 (d, $J = 6.7$ Hz, 3H, ArH), 7.70 (d, $J = 8.4$ Hz, 1H, ArH), 8.06 (d, $J = 8.5$ Hz, 3H, ArH) ppm; ^{13}C NMR (150 MHz, CDCl_3) $\delta = 56.2, 56.3, 105.0, 108.4, 117.0, 122.9, 128.6, 129.1, 135.1, 135.2, 138.5, 145.4, 150.0, 152.8, 154.1$ ppm; HRMS (ESI) m/z: calcd. for $\text{C}_{17}\text{H}_{14}\text{ClNO}_2$ [M+H⁺]: 300.0786; found 300.0790.

2-(4-Fluorophenyl)-6,7-dimethoxyquinoline (18a): Yield = 76% (0.088 g from 0.05 g) as a dark yellow solid; m.p. 122-124 °C; $R_f = 0.67$ (hexane/EtOAc, 80:20, v/v); IR (Neat): $\nu_{\max} (\text{cm})^{-1} = 1617$ (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 4.03$ (s, 3H, -OCH₃), 4.06 (s, 3H, -OCH₃), 7.06 (s, 1H, ArH), 7.19 (t, $J = 8.7$ Hz, 2H, ArH), 7.49 (s, 1H, ArH), 7.68 (d, $J = 8.4$ Hz, 1H, ArH), 8.06 (d, $J = 8.5$ Hz, 1H, ArH), δ 8.09(d, $J = 5.5$ Hz, 1H, ArH), 8.10 (d, $J = 5.5$ Hz, 1H, ArH) .ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 56.2, 56.3, 105.0, 108.3, 115.7, 108.3, 115.7, 115.9$ ($J_{C,F} = 21.4$ Hz), 117.1, 122.8, 129.1, 129.2 ($J_{C,F} = 8.1$ Hz), 135.2, 145.3, 149.9, 152.9, 154.4, 162.8, 164.5 ($J_{C,F} = 246.7$ Hz) ppm; HRMS (ESI) m/z: calcd. For $\text{C}_{17}\text{H}_{14}\text{FNO}_2$ [M + H⁺]: 284.1081found 284.1085.

6,7-Dimethoxy-2-(4-methoxyphenyl)quinoline (19a): Yield = 60% (0.11 g from 0.10 g) as an orange solid; m.p. 224-226 °C; R_f = 0.46 (hexane/EtOAc, 70:30, v/v); IR (Neat): ν_{\max} (cm)⁻¹ = 1595(C=N); ¹H NMR (600 MHz, CDCl₃) δ = 3.86 (s, 3H, -OCH₃), 4.01 (s, 3H, -OCH₃), 4.04 (s, 3H, -OCH₃), 7.02 (t, J = 6.6 Hz, 3H, ArH), 7.46 (s, 1H, ArH), 7.67 (d, J = 8.4 Hz, 1H, ArH), 8.01 (d, J = 8.4 Hz, 1H, ArH), 8.05-8.07 (m, 2H, ArH) ppm; ¹³C NMR (150 MHz, CDCl₃) δ = 55.5, 56.2, 56.3, 105.2, 108.4, 114.3, 117.0, 122.5, 128.5, 132.9, 135.0, 145.4, 149.7, 152.7, 152.2, 160.9 ppm; HRMS (ESI) m/z: calcd. for C₁₈H₁₇NO₃ [M + H⁺]: 296.1281, found: 296.1285.

6,7-Dimethoxy-2-(p-tolyl)quinoline (20a): Yield = 71% (0.075 g from 0.10 g) as a white solid; m.p. 130-132 °C; R_f = 0.63 (hexane/EtOAc, 70:30, v/v); IR (Neat): ν_{\max} (cm)⁻¹ = 1615(C=N); ¹H NMR (600 MHz, CDCl₃) δ = 2.42 (s, 3H, -CH₃), 4.02 (s, 3H, -OCH₃), 4.08 (s, 3H, -OCH₃), 7.26 (s, 1H, ArH), 7.06 (s, 1H, ArH), 7.31 (d, J = 6.6 Hz, 2H, ArH), 7.50 (s, 1H, ArH), 7.70-7.73 (m, 1H, ArH), 8.01-8.04 (m, 2H, ArH) ppm; ¹³C NMR (150 MHz, CDCl₃) δ = 21.4, 56.0, 56.3, 106.1, 108.5, 108.5, 108.5, 117.3, 122.7, 127.3, 129.6, 129.7, 135.0, 137.3, 139.0, 145.3, 149.7, 152.7, 155.5 ppm; HRMS (ESI) m/z: calcd. for C₁₈H₁₇NO₂ [M + H⁺]: 264.1383, found: 264.1387.

6,7-Dimethoxy-2-(4-nitrophenyl)quinoline (21a): Yield = 89% (0.130 g from 0.10 g) as a yellow solid; m.p. 148-150 °C; R_f = 0.28(hexane/EtOAc, 80:20, v/v); IR (neat): ν_{\max} (cm)⁻¹ = 1610 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 4.05 (s, 3H, -OCH₃), 4.08 (s, 3H, -OCH₃), 7.09 (s, 1H, ArH), 7.49 (s, 1H, ArH), 7.78 (d, J = 8.4 Hz, 1H, ArH), 8.12 (d, J = 8.4 Hz, 1H, ArH), 8.30 (d, J = 8.9 Hz, 2H, ArH), 8.35 (d, J = 8.9 Hz, 2H, ArH) ppm; ¹³C NMR (150 MHz, CDCl₃) δ = 56.3, 56.4, 104.9, 108.4, 117.3, 123.6, 124.2, 128.0, 135.4, 145.6, 146.0, 148.1, 150.6, 152.6, 153.2 ppm; HRMS (ESI) m/z: calcd. for C₁₇H₁₄N₂O₄ [M+H⁺]: 311.1026; found 311.1030.

2-([1,1'-Biphenyl]-4-yl)-6,7-dimethoxyquinoline (22a): Yield = 77% (0.109 g from 0.10 g) as a light yellow solid; m.p. 174-176 °C; R_f = 0.37 (hexane/EtOAc, 80:20, v/v); IR (neat): ν_{\max} (cm)⁻¹ = 1635 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 4.04 (s, 3H, -OCH₃), 4.07 (s, 3H, -OCH₃), 7.07 (s, 1H, ArH), 7.37 (t, J = 7.4 Hz, 1H, ArH), 7.46 (d, J = 7.8 Hz, 2H, ArH), 7.52 (s, 1H, ArH), 7.68 (d, J = 7.0 Hz, 2H, ArH), 7.75 (d, J = 8.3 Hz, 2H, ArH), 7.78 (d, J = 8.4 Hz, 1H, ArH), 8.07 (d, J = 8.5 Hz, 1H, ArH), 8.20 (d, J = 8.2 Hz, 2H, ArH) ppm; ¹³C NMR (150 MHz, CDCl₃) δ = 56.2, 56.3, 105.1, 108.5, 117.3, 122.9, 127.2, 127.6, 127.8, 129.0, 135.1, 139.0, 140.8, 141.7, 145.5, 149.8, 152.7, 155.0 ppm; HRMS (ESI) m/z: calcd. for C₂₃H₁₉NO₂ [M+H⁺]: 342.1489; found 342.1493.

6-Phenyl-[1,3]dioxolo[4,5-g]quinoline (15b): Yield = 65% (0.097 g from 0.10 g) as a light yellow solid; m.p. 112-114 °C; R_f = 0.56 (hexane/EtOAc, 90:10, v/v); IR (neat): ν_{\max} (cm)⁻¹ = 1615 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 6.11 (s, 2H, -OCH₂O-), 7.07 (s, 1H, ArH), 7.43 (t, J = 7.3 Hz, 1H, ArH), 7.46 (s, 1H, ArH), 7.51 (t, J = 7.6 Hz, 2H, ArH), 7.72 (d, J = 8.5 Hz, 1H, ArH), 8.02 (d, J = 7.8 Hz, 1H, ArH), 8.10 (d, J = 1.1 Hz, 1H, ArH), 8.11 (d, J = 1.4 Hz, 1H, ArH) ppm; ¹³C NMR (151 MHz, CDCl₃) δ = 101.8, 102.7, 106.3, 117.4, 124.3, 127.4, 128.9, 129.0, 135.7, 139.9, 146.7, 147.9, 151.0, 155.5 ppm; HRMS (ESI) m/z : calcd. for C₁₆H₁₁NO₂ [M + H⁺]: 250.0863, found 250.0867.

6-(4-Bromophenyl)-[1,3]dioxolo[4,5-g]quinoline (16b): Yield = 68% (0.107 g from 0.10 g) as a white solid; m.p. 200-202 °C; R_f = 0.78 (hexane/EtOAc, 95:5, v/v); IR (KBr): ν_{\max} (cm)⁻¹ = 1615 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 6.22 (s, 2H, -OCH₂O), 7.36 (s, 1H, ArH), 7.39 (s, 1H, ArH), 7.70-7.71 (m, 2H, ArH), 7.71 (d, J = 8.5 Hz, 1H, ArH), 8.10-8.16 (d, J = 9.0 Hz, 2H, ArH), 8.24 (d, J = 8.4 Hz, 1H, ArH) ppm; ¹³C NMR (151 MHz, CDCl₃) δ = 1.2, 53.7, 56.1, 56.3, 105.0, 108.4, 117.0, 123.0, 128.6, 129.2, 135.1, 135.3, 138.5, 145.3, 150.0, 152.9, 154.1 ppm; HRMS (ESI) m/z : calcd. for C₁₆H₁₀BrNO₂ [M + H⁺]: 327.9968, found: 327.9972.

8-(4-Chlorophenyl)-[1,3]dioxolo[4,5-g]quinolone (17b): Yield = 41%, (0.07 g from 0.10 g) as a white solid; m.p. 224-226 °C; R_f = 0.63 (hexane/EtOAc, 70:30, v/v); IR (KBr): ν_{\max} (cm)⁻¹ = 1613(C=N); ¹H NMR (600 MHz, CDCl₃) δ = 6.10 (s, 2H, -OCH₂O), 7.04 (s, 1H, ArH), 7.24 (s, 1H, ArH), 7.41 (s, 1H, ArH), 7.45 (d, J = 6.7 Hz, 1H, ArH), 7.66 (d, J = 8.5 Hz, 1H, ArH), 7.99 (d, J = 8.5 Hz, 1H, ArH), 8.05 (d, J = 7.2 Hz, 2H, ArH) ppm; ¹³C NMR (150 MHz, CDCl₃) δ = 0.1, 101.9, 102.7, 106.3, 116.9, 119.9, 124.4, 128.6, 129.1, 135.2, 135.8, 138.3, 146.7, 148.0, 151.1, 154.1 ppm; HRMS (ESI) m/z : calcd. for C₁₆H₁₀ClNO₂ [M + H⁺]: 283.0400, found: 283.0404.

6-(4-Fluorophenyl)-[1,3]dioxolo[4,5-g]quinoline (18b): Yield = 40%, (0.07 g from 0.10 g) as a yellow solid; m.p. 192-194 °C; R_f = 0.85 (hexane/EtOAc, 70:30, v/v); IR (KBr): ν_{\max} (cm)⁻¹ = 1661(C=N); ¹H NMR (600 MHz, CDCl₃) δ = 6.11 (s, 2H, -OCH₂O), 7.66 (d, J = 8.5 Hz, 1H, ArH), 8.01 (d, J = 8.5 Hz, 1H, ArH), 8.08-8.10 (m, 2H, ArH), 7.06 (s, 1H, ArH), 7.16-7.19 (m, 2H, ArH), 7.43 (s, 1H, ArH) ppm; ¹³C NMR (151 MHz, CDCl₃) δ = 101.8, 102.8, 106.2, 115.7, 115.9 ($J_{C,F}$ = 21.5 Hz) 116.0, 117.0, 124.2, 129.2 ($J_{C,F}$ = 8.3 Hz), 136.7, 136.0, 146.8, 147.9, 151.2, 154.5, 163.0, 164.5 ($J_{C,F}$ = 246.9 Hz) ppm; HRMS (ESI) m/z : calcd. for C₁₆H₁₀FNO₂ [M + H⁺]: 268.0768, found: 268.0772.

6-(4-Methoxyphenyl)-[1,3]dioxolo[4,5-g]quinoline (19b): Yield = 88% (0.154 g from 0.10 g) as a yellow solid; m.p. 223-225 °C; $R_f = 0.75$ (hexane/EtOAc, 70:30, v/v); IR (KBr): ν_{\max} (cm)⁻¹ = 1648(C=N); ¹H NMR (600 MHz, CDCl₃) $\delta = 3.86$ (s, 3H, -OCH₃), 6.08 (s, 2H, -OCH₂O), 7.02 (t, $J = 7.5$ Hz, 3H, ArH), 7.41 (s, 1H, ArH), 7.65 (d, $J = 8.5$ Hz, 1H, ArH), 7.96 (d, $J = 8.6$ Hz, 1H, ArH), 8.05 (d, $J = 8.8$ Hz, ArH) ppm; ¹³C NMR (151 MHz, CDCl₃) $\delta = 1.2, 55.5, 101.7, 102.7, 106.2, 114.3, 117.0, 123.8, 128.6, 132.5, 135.5, 146.75, 147.6, 151.0, 155.1, 160.6$ ppm; HRMS (ESI) m/z: calcd. for C₁₇H₁₃NO₃ [M + H⁺]: 280.0968, found: 280.0972.

6-(p-Tolyl)-[1,3]dioxolo[4,5-g]quinoline (20b): Yield = 85% (0.094 g from 0.10 g) as a grey solid; m.p. 180-182 °C; $R_f = 0.41$ (hexane/EtOAc, 95:5, v/v); IR (KBr): ν_{\max} (cm)⁻¹ = 1613 (C=N); ¹H NMR (600 MHz, CDCl₃) $\delta = 2.4$ (s, 3H, -CH₃), 6.1 (s, 2H, -OCH₂O), 7.05 (s, 1H, ArH), 7.26-7.27 (m, 1H, ArH), 7.31 (d, $J = 7.2$ Hz, 2H, ArH), 7.45 (s, 1H, ArH), 7.69-7.71 (m, 1H, ArH), 8.0 (t, $J = 6.6$ Hz, 2H, ArH) ppm; ¹³C NMR (150 MHz, CDCl₃) $\delta = 21.4, 101.7, 102.7, 102.7, 106.3, 106.3, 117.2, 124.1, 127.2, 129.6, 135.6, 137.1, 139.0, 146.7, 147.7, 150.9, 155.4$ ppm; HRMS (ESI) m/z: calcd. for C₁₇H₁₃NO₂ [M + H⁺]: 264.1019, found: 264.1023.

6-([1,1'-Biphenyl]-4-yl)-[1,3]dioxolo[4,5-g]quinoline (22b): Yield = 69% (0.093 g from 0.10 g) as a white solid; m.p. 202-204 °C; $R_f = 0.54$ (hexane/EtOAc, 80:20, v/v); IR (neat): ν_{\max} (cm)⁻¹ = 1600 (C=N); ¹H NMR (600 MHz, CDCl₃) $\delta = 6.12$ (s, 2H, -OCH₂O), 7.08 (s, 1H, ArH), 7.36 – 7.39 (m, 1H, ArH), 7.47 (d, $J = 6.7$ Hz, 3H, ArH), 7.68 (d, $J = 7.7$ Hz, 2H, ArH), 7.75 (d, $J = 8.3$ Hz, 2H, ArH), 7.77 (d, $J = 8.5$ Hz, 1H, ArH), 8.04 (d, $J = 8.4$ Hz, 1H, ArH), 8.20 (d, $J = 8.3$ Hz, 2H, ArH) ppm; ¹³C NMR (150 MHz, CDCl₃) $\delta = 101.8, 102.7, 106.3, 117.3, 124.3, 127.3, 127.6, 127.8, 129.0, 135.7, 138.8, 140.8, 141.8, 147.9, 151.0$ ppm; HRMS (ESI) m/z: calcd. for C₂₂H₁₅NO₂ [M+H⁺]: 326.1176; found 326.1180.

7-Phenyl-2,3-dihydro-[1,4]dioxino[2,3-g]quinoline (15c): Yield = 66% (0.105 g from 0.10 g) as an off white solid; m.p. 146-148 °C; $R_f = 0.71$ (hexane/EtOAc, 80:20, v/v); IR (neat): ν_{\max} (cm)⁻¹ = 1625 (C=N); ¹H NMR (600 MHz, CDCl₃) $\delta = 4.40$ (s, 4H, -OCH₂CH₂O-), 7.22 (s, 1H, ArH), 7.44 (t, $J = 7.3$ Hz, 1H, ArH), 7.51 (t, $J = 7.8$ Hz, 2H, ArH), 7.62 (s, 1H, ArH), 7.71 (d, $J = 8.5$ Hz, 1H, ArH), 8.03 (d, $J = 8.6$ Hz, 1H, ArH), 8.11 (d, $J = 8.3$ Hz, 2H, ArH) ppm; ¹³C NMR (150 MHz, CDCl₃) $\delta = 64.5, 111.7, 114.6, 117.6, 123.3, 127.5, 128.9, 129.1, 135.3, 140.1, 144.3, 144.9, 147.1, 156.0$ ppm; HRMS (ESI) m/z: calcd. for C₁₇H₁₃NO₂ [M+H⁺]: 264.1019; found 264.1023.

(4-Bromophenyl)-2,3-dihydro-[1,4]dioxino[2,3-g]quinoline (16c): Yield = 46% (0.103 g from 0.10 g) as a light yellow solid; m.p. 182-184 °C; $R_f = 0.67$ (hexane/EtOAc, 70:30, v/v); IR (neat): $\nu_{\max} (\text{cm})^{-1} = 1628$ (C=N) ; $^1\text{H NMR}$ (600 MHz, CDCl_3) $\delta = 4.39$ (d, $J = 2.1$ Hz, 4H, - $\text{OCH}_2\text{CH}_2\text{O}$ -), 7.21 (s, 1H, ArH), 7.58 (s, 1H, ArH), 7.62 (d, $J = 8.5$ Hz, 2H, ArH), 7.65 (d, $J = 8.6$ Hz, 1H, ArH), 7.99 (d, $J = 8.6$ Hz, 2H, ArH), 8.02 (d, $J = 8.3$ Hz, 1H, ArH) ppm; $^{13}\text{C NMR}$ (151 MHz, CDCl_3) $\delta = 64.5, 64.6, 111.7, 114.6, 117.1, 122.2, 122.8, 123.4, 123.6, 129.0, 132.0, 135.4, 138.9, 144.5, 144.9, 147.3, 154.6$ ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{17}\text{H}_{12}\text{BrNO}_2$ [$\text{M} + \text{H}$]: 342.0124 found 342.0128.

7-(4-Chlorophenyl)-2,3-dihydro-[1,4]dioxino[2,3-g]quinoline (17c): Yield = 74% (0.110 g from 0.10 g) as an off white solid; m.p. 168-170 °C; $R_f = 0.6$ (hexane/EtOAc, 80:20, v/v); IR (neat): $\nu_{\max} (\text{cm})^{-1} = 1625$ (C=N); $^1\text{H NMR}$ (600 MHz, CDCl_3) $\delta = 4.39$ (d, $J = 2.1$ Hz, 4H, - $\text{OCH}_2\text{CH}_2\text{O}$ -), 7.21 (s, 1H, ArH), 7.47 (d, $J = 8.5$ Hz, 2H, ArH), 7.58 (s, 1H, ArH), 7.66 (d, $J = 8.5$ Hz, 1H, ArH), 8.02 (d, $J = 8.6$ Hz, 1H, ArH), 8.06 (d, $J = 8.5$ Hz, 2H, ArH) ppm; $^{13}\text{C NMR}$ (150 MHz, CDCl_3) $\delta = 64.5, 111.7, 114.6, 117.1, 123.4, 128.7, 129.0, 129.1, 135.2, 135.4, 138.5, 144.5, 144.9, 147.3, 154.6$ ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{17}\text{H}_{12}\text{ClNO}_2$ [$\text{M} + \text{H}^+$]: 298.0629; found 298.0633.

7-(4-Methoxyphenyl)-2,3-dihydro-[1,4]dioxino[2,3-g]quinoline (19c): Yield = 88% (0.041 g from 0.10 g) as a light yellow solid; m.p. 132-134 °C; $R_f = 0.48$ (hexane/EtOAc, 70:30, v/v); IR (neat): $\nu_{\max} (\text{cm})^{-1} = 1624$ (C=N); $^1\text{H NMR}$ (600 MHz, CDCl_3) $\delta = 3.88$ (s, 3H, - OCH_3), 4.40 – 4.38 (m, 2H, - $\text{OCH}_2\text{CH}_2\text{O}$ -), 4.46 – 4.43 (m, 2H, - $\text{OCH}_2\text{CH}_2\text{O}$ -), 7.03 (d, $J = 8.8$ Hz, 2H, ArH), 7.30 (d, $J = 9.1$ Hz, 1H, ArH), 7.66 (d, $J = 9.1$ Hz, 1H, ArH), 7.78 (d, $J = 8.8$ Hz, 1H, ArH), 8.10 (d, $J = 8.8$ Hz, 2H, ArH), 8.39 (d, $J = 8.8$ Hz, 1H, ArH), ppm; $^{13}\text{C NMR}$ (151 MHz, CDCl_3) $\delta = 55.5, 64.5, 111.7, 114.3, 114.5, 117.2, 118.0, 121.8, 122.6, 123.0, 128.7, 129.6, 132.7, 135.2, 144.0, 144.9, 147.0, 155.6, 160.7$ ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{18}\text{H}_{15}\text{NO}_3$ [$\text{M} + \text{H}$]: 294.1125, found 294.1129.

7-(*p*-Tolyl)-2,3-dihydro-[1,4]dioxino[2,3-g]quinoline (20c): Yield = 85% (0.130 g from 0.10 g) as a light brown solid; m.p. 70-72 °C; $R_f = 0.5$ (hexane/EtOAc, 80:20, v/v); IR (neat): $\nu_{\max} (\text{cm})^{-1} = 1620$ (C=N); $^1\text{H NMR}$ (600 MHz, CDCl_3) $\delta = 2.42$ (s, 3H, - CH_3), 4.38 (d, $J = 2.9$ Hz, 4H, - $\text{OCH}_2\text{CH}_2\text{O}$ -), 7.20 (s, 1H, ArH), 7.31 (d, $J = 7.8$ Hz, 2H, ArH), 7.59 (s, 1H, ArH), 7.68 (d, $J = 8.5$ Hz, 1H, ArH), 8.00 (dd, $J = 3.5, 8.3$ Hz, 3H, ArH) ppm; $^{13}\text{C NMR}$ (150 MHz, CDCl_3) $\delta = 21.5,$

64.6, 111.7, 114.6, 117.5, 123.2, 127.4, 129.6, 135.2, 137.3, 139.1, 144.2, 144.9, 147.1, 156.0 ppm; HRMS (ESI) m/z : calcd. for $C_{18}H_{15}NO_2$ $[M+H^+]$: 278.1176; found 278.1180.

5,7-Dimethoxy-2-phenylquinoline (15d): Yield = 43% (0.066 g from 0.10 g) as a yellow oily; R_f = 0.56 (hexane/EtOAc, 90:10, v/v); ; m.p. 112-114 °C; IR (neat): ν_{max} (cm)⁻¹ = 1623 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 3.96 (s, 6H, (-OCH₃)₂), 6.50 (d, J = 2.2 Hz, 1H, ArH), 7.11 (d, J = 2.2 Hz, 1H, ArH), 7.45 (t, J = 7.3 Hz, 1H, ArH), 7.51 (t, J = 7.4 Hz, 2H, ArH), 7.68 (d, J = 8.6 Hz, 1H, ArH), 8.13 (d, J = 7.7 Hz, 2H, ArH), 8.47 (d, J = 8.6 Hz, 1H, ArH) ppm; ¹³C NMR (151 MHz, CDCl₃) δ = 55.7, 55.8, 98.0, 100.1, 115.7, 116.0, 127.6, 128.8, 129.0, 129.3, 129.6, 131.5, 140.0, 150.5, 156.1, 158.1, 161.5 ppm; HRMS (ESI) m/z : calcd. for $C_{17}H_{15}NO_2$ $[M + H^+]$: 266.1176, found 266.1180.

2-(4-Bromophenyl)-5,7-dimethoxyquinoline (16d): Yield = 53%, (0.127 g from 0.10 g) as a white solid; m.p; 140-142 °C; R_f = 0.9 (hexane/EtOAc, 90:10, v/v); IR (KBr): ν_{max} (cm)⁻¹ = 1568 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 3.96 (s, 3H, -OCH₃), 3.98 (s, 3H, -OCH₃), 6.51 (s, 1H, ArH), 7.07 (s, 1H, ArH), 7.63 (t, J = 8.8 Hz, 3H, ArH), 8.01 (d, J = 8.9 Hz, 2H, ArH), 8.48 (d, J = 8.6 Hz, 1H, ArH) ppm; ¹³C NMR (150 MHz, CDCl₃) δ = 54.1, 55.9, 55.9, 98.4, 100.0, 115.7, 115.7, 116.0, 123.8, 129.3, 131.9, 132.0, 138.8, 150.7, 156.1, 156.0, 161.7 ppm; HRMS (ESI) m/z : calcd. for $C_{17}H_{14}BrNO_2$ $[M + H^+]$: 344.0281, found: 344.0285.

2-(4-Chlorophenyl)-5,7-dimethoxyquinoline (17d): Yield = 49% (0.073 g from 0.10 g) as a green solid; m.p. 220-222 °C; R_f = 0.45 (hexane/EtOAc, 90:10, v/v); IR (neat): ν_{max} (cm)⁻¹ = 1616 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 3.98 (s, 3H, -OCH₃), 4.06 (s, 3H, -OCH₃), 6.76 (d, J = 8.4 Hz, 1H, ArH), 6.97 (d, J = 8.5 Hz, 1H, ArH), 7.47 (d, J = 8.6 Hz, 2H, ArH), 7.86 (d, J = 8.8 Hz, 1H, ArH), 8.15 (d, J = 8.7 Hz, 2H, ArH), 8.61 (d, J = 8.7 Hz, 1H, ArH), ppm; ¹³C NMR (151 MHz, CDCl₃) δ = 55.9, 55.9, 56.5, 103.9, 108.0, 118.2, 120.7, 129.0, 129.1, 132.1, 135.5, 138.2, 140.6, 148.9, 149.7, 155.4 ppm; HRMS (ESI) m/z : calcd. for $C_{17}H_{14}ClNO_2$ $[M + H^+]$: 300.0786, found 300.0790.

2-(4-Fluorophenyl)-5,7-dimethoxyquinoline (18d): Yield = 48% (0.095 g from 0.10 g) as an off white solid; m.p. 162-164 °C; R_f = 0.46 (hexane/EtOAc, 90:10, v/v); IR (neat): ν_{max} (cm)⁻¹ = 1599 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 4.16 (s, , 6H, -OCH₃) , 6.69 (s, 1H, ArH), 7.38 (t, J = 8.6 Hz, 2H, ArH), 7.45 (s, 1H, ArH), 7.83 (d, J = 8.6 Hz, 1H, ArH), 8.34 – 8.27 (m, 2H ArH), 8.67 (d, J = 8.6 Hz, 1H, ArH) ppm; ¹³C NMR (151 MHz, CDCl₃) δ = 55.8, 55.9, 98.1, 100.1, 115.7 ($J_{C,F}$ = 5.8 Hz), 115.9, 129.4, 129.5 ($J_{C,F}$ = 8.2 Hz), 131.7, 136.1, 150.5, 156.1, 157.1, 161.7,

163.0, 164.7 ($J_{C,F} = 247.4$ Hz) ppm; HRMS (ESI) m/z : calcd. for $C_{17}H_{14}FNO_2$ [M + H⁺]: 284.1081, Found 284.1085.

5,7-Dimethoxy-2-(*p*-tolyl)quinoline (20d): Yield = 41% (0.075 g from 0.10 g) as a light yellow solid; m.p. 102-104 °C; $R_f = 0.56$ (hexane/EtOAc, 90:10, v/v); IR (neat): ν_{max} (cm)⁻¹ = 1624 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 2.43 (s, 3H, -CH₃), 3.96 (s, 3H, -OCH₃), 3.98 (s, 3H, -OCH₃), 6.49 (d, $J = 2.2$ Hz, 1H, ArH), 7.09 (d, $J = 1.9$ Hz, 1H, ArH), 7.32 (d, $J = 8.4$ Hz, 2H, ArH), 7.67 (d, $J = 8.6$ Hz, 1H, ArH), 8.03 (d, $J = 8.2$ Hz, 2H, ArH), 8.46 (d, $J = 8.6$ Hz, 1H, ArH) ppm; ¹³C NMR (151 MHz, CDCl₃) δ = 21.4, 55.7, 55.9, 97.9, 100.2, 115.6, 115.9, 127.5, 129.6, 131.4, 137.2, 139.2, 150.6, 156.1, 158.1, 161.5 ppm; HRMS (ESI) m/z : calcd. for $C_{18}H_{17}NO_2$ [M + H⁺]: 280.1332, found 280.1336.

5,8-Dimethoxy-2-phenylquinoline (15e): Yield = 51% (0.084 g from 0.10 g) as a brown oily; $R_f = 0.28$ (hexane/EtOAc, 90:10, v/v); ; m.p. 112-114 °C; IR (neat): ν_{max} (cm)⁻¹ = 1620 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 3.98 (s, 3H, -OCH₃), 4.07 (s, 3H, -OCH₃), 6.75 (d, $J = 8.4$ Hz, 1H, ArH), 6.96 (d, $J = 8.5$ Hz, 1H, ArH), 7.44 (d, $J = 7.3$ Hz, 1H, ArH), 7.51 (d, $J = 7.6$ Hz, 2H, ArH), 7.91 (d, $J = 8.4$ Hz, 1H, ArH), 8.20 (d, $J = 7.1$ Hz, 2H, ArH), 8.62 (d, $J = 8.7$ Hz, 1H, ArH) ppm; ¹³C NMR (151 MHz, CDCl₃) δ = 55.9, 56.5, 103.7, 107.9, 118.7, 120.6, 127.5, 127.8, 128.5, 128.9, 129.3, 131.8, 139.8, 140.7, 148.9, 149.8, 156.7 ppm; HRMS (ESI) m/z : calcd. for $C_{17}H_{15}NO_2$ [M + H⁺]: 266.1176, found 266.1180.

2-(4-Bromophenyl)-5,8-dimethoxyquinoline (16e): Yield = 37% (0.078 g from 0.150 g) as a yellow solid; m.p. 134-136 °C; $R_f = 0.56$ hexane/EtOAc, 90:10, v/v); IR (neat): ν_{max} (cm)⁻¹ = 1603 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 3.98 (s, 3H, -OCH₃), 4.06 (s, 3H, -OCH₃), 6.76 (d, $J = 8.4$ Hz, 1H, ArH), 6.97 (d, $J = 8.5$ Hz, 1H ArH), 7.63 (d, $J = 8.5$ Hz, 2H ArH), 7.86 (d, $J = 8.8$ Hz, 1H, ArH), 8.08 (d, $J = 8.5$ Hz, 2H, ArH), δ 8.61 (d, $J = 8.8$ Hz, 1H, ArH) ppm; ¹³C NMR (151 MHz, CDCl₃) δ = 55.9, 56.5, 103.9, 108.0, 118.2, 120.7, 123.9, 129.3, 132.0, 132.1, 138.7, 140.6, 148.9, 149.7, 155.4 ppm; HRMS (ESI) m/z : calcd. for $C_{17}H_{14}BrNO_2$ [M + H⁺]: 344.0281, found 344.0285.

2-(4-Fluorophenyl)-5,8-dimethoxyquinoline (18e): Yield = 30% (0.060 g from 0.10 g) as a brown solid; m.p. =152-154 °C; $R_f = 0.56$ (hexane/EtOAc, 90:10, v/v); IR (neat): ν_{max} (cm)⁻¹ = 1713 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 3.98 (s, 3H, -OCH₃), 4.06 (s, 3H, -OCH₃), 6.75 (d, $J = 8.5$ Hz, 1H ArH), 6.96 (d, $J = 8.4$ Hz, 1H ArH), 7.18 (t, $J = 8.1$ Hz, 2H, ArH), 7.85 (d, $J = 8.8$ Hz, 1H, ArH), 8.21 – 8.16 (m, 2H, ArH), 8.60 (d, $J = 8.8$ Hz, 1H ArH) ppm; ¹³C NMR (151 MHz,

CDCl₃) δ = 55.9, 56.5, 103.7, 107.9, 115.7, 115.8 ($J_{C,F}$ = 21.5 Hz), 118.3, 120.5, 129.6, 129.7 ($J_{C,F}$ = 8.5 Hz), 132.0, 136.0, 140.6, 148.9, 149.7, 155.6, 163.1, 164.7 ($J_{C,F}$ = 247.1 Hz) ppm; HRMS (ESI) m/z : calcd. For C₁₇H₁₄FNO₂ [M + H⁺]: 284.1081, found 284.1085.

6-(4-Bromophenyl)-2,3-dimethylnaphthalene (16f): Yield = 36% (0.10 g from 0.10 g) as a light yellow solid; m.p. 182-184 °C; R_f = 0.75 (hexane/EtOAc, 95:05, v/v); IR (neat): ν_{\max} (cm)⁻¹ = 1628 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 2.45 (s, 3H, -CH₃), 2.48 (s, 3H, -CH₃), 7.55 (s, 1H, ArH), 7.63 (d, J = 8.5 Hz, 2H, ArH), 7.73 (d, J = 8.6 Hz, 1H, ArH), 7.91 (s, 1H, ArH), 8.02 (d, J = 8.5 Hz, 2H, ArH), 8.09 (d, J = 8.5 Hz, 1H, ArH) ppm; ¹³C NMR (151 MHz, CDCl₃) δ = 20.1, 20.5, 20.6, 76.9, 77.1, 77.3, 117.8, 118.0, 123.6, 126.0, 126.8, 129.1, 129.1, 131.9, 132.0, 133.1, 135.9, 136.6, 138.9, 140.2, 147.5, 155.2 ppm; HRMS (ESI) m/z : calcd. for C₁₈H₁₅Br [M + H⁺]: 312.0382, found 312.0386.

2-(2-Chlorophenyl)-6,7-dimethoxyquinoline (23a): Yield = 52% (0.097 g from 0.10 g) as a yellow solid; m.p. 136-138 °C; R_f = 0.2 (hexane/EtOAc, 90:10, v/v); IR (neat): ν_{\max} (cm)⁻¹ = 1624 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 4.04 (s, 6H, (-OCH₃)₂), 7.10 (s, 1H, ArH), 7.34 – 7.41 (m, 2H, ArH), 7.49 – 7.41 (m, 2H, ArH), 7.60 (d, J = 8.3 Hz, 1H, ArH), 7.67 (d, J = 7.5 Hz, 1H, ArH), 8.06 (d, J = 8.3 Hz, 1H, ArH) ppm; ¹³C NMR (151 MHz, CDCl₃) δ = 56.2, 56.3, 105.1, 108.4, 121.1, 122.9, 127.2, 129.7, 130.2, 131.7, 132.5, 134.0, 140.0, 145.1, 150.2, 152.8, 155.3 ppm; HRMS (ESI) m/z : calcd. for C₁₇H₁₄ClNO₂ [M + H⁺]: 300.0786, found 300.0790.

2-(3-Bromophenyl)-6,7-dimethoxyquinoline (24a): Yield = 68% (0.112 g from 0.05 g) as a yellow solid; m.p. 154-156 °C; R_f = 0.32 (hexane/EtOAc, 90:10, v/v); IR (neat): ν_{\max} (cm)⁻¹ = 1620 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 4.03 (s, 3H, -OCH₃), 4.07 (s, 3H, -OCH₃), 7.07 (s, 1H, ArH), 7.37 (t, J = 7.8 Hz, 1H, ArH), 7.49 (s, 1H, ArH), 7.54 – 7.56 (m, 1H, ArH), 7.70 (d, J = 8.5 Hz, 1H, ArH), 8.02 (d, J = 7.9 Hz, 1H, ArH), 8.07 (d, J = 8.4 Hz, 1H, ArH), 8.31 (t, J = 1.8 Hz, 1H, ArH), ppm; ¹³C NMR (151 MHz, CDCl₃) δ = 56.22, 56.34, 105.0, 108.5, 117.1, 123.1, 123.2, 125.8, 130.4, 130.5, 131.9, 135.2, 142.2, 145.4, 150.1, 152.9, 153.8, ppm; HRMS (ESI) m/z : calcd. for C₁₇H₁₄BrNO₂ [M + H⁺]: 344.0281, found 344.0285.

2-(3,4-Dichlorophenyl)-6,7-dimethoxyquinoline (25a): Yield = 86% (0.122 g from 0.10 g) as a white solid; m.p. 176-178 °C; R_f = 0.52 (hexane/EtOAc, 80:20, v/v); IR (neat): ν_{\max} (cm)⁻¹ = 1635 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 4.03 (s, 3H, OCH₃), 4.07 (s, 3H, OCH₃), 7.06 (s, 1H, ArH), 7.47 (s, 1H, ArH), 7.56 (d, J = 9.2 Hz, 1H, ArH), 7.68 (d, J = 7.6 Hz, 1H, ArH), 7.94 (d, J = 8.5 Hz, 1H, ArH), 8.07 (d, J = 8.4 Hz, 1H, ArH), 8.26 (s, 1H, ArH) ppm; ¹³C NMR (150 MHz,

CDCl₃) δ = 56.2, 56.3, 105.0, 108.4, 116.8, 123.2, 126.4, 129.2, 130.8, 133.1, 133.2, 135.3, 140.0, 145.4, 150.2, 152.7, 153.0 ppm; HRMS (ESI) m/z : calcd. for C₁₇H₁₃Cl₂NO₂ [M+H⁺]: 334.0396; found 334.0391.

2-(3,4-Dimethoxyphenyl)-6,7-dimethoxyquinoline (26a): Yield = 34% (0.068 g from 0.10 g) as a light brown solid; m.p. 176-178 °C; R_f = 0.56 (hexane/EtOAc, 90:10, v/v); IR (neat): ν_{\max} (cm)⁻¹ = 1600 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 3.96 (s, 3H, -OCH₃), 4.03 (s, 3H, -OCH₃), 4.05 (s, 3H, -OCH₃), 4.07 (s, 3H, -OCH₃), 6.99 (d, J = 8.3 Hz, 1H, ArH), 7.06 (s, 1H, ArH), 7.51 (s, 1H, ArH), 7.62 (d, J = 8.3 Hz, 1H, ArH), 7.71 (d, J = 8.4 Hz, 1H, ArH), 7.82 (s, 1H, ArH), 8.03 (d, J = 8.4 Hz, 1H, ArH) ppm; ¹³C NMR (151 MHz, CDCl₃) δ = 56.1, 56.3, 105.1, 108.3, 110.0, 111.2, 117.0, 119.9, 122.5, 133.0, 135.0, 145.2, 149.5, 149.6, 150.6, 152.7, 155.0 ppm; HRMS (ESI) m/z : calcd. for C₁₉H₁₉NO₄ [M + H⁺]: 326.1387, found 326.1391.

6,7-Dimethoxy-2-(naphthalen-2-yl)quinoline (27a): Yield = 95% (0.196 g from 0.10 g) as a yellow solid; m.p. 138-140 °C; R_f = 0.68 (hexane/EtOAc, 90:10, v/v); IR (neat): ν_{\max} (cm)⁻¹ = 1618 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 1.52 (t, J = 7.3 Hz, 3H, -CH₃), 4.54 (q, J = 7.3 Hz, 2H, -CH₂), 7.41 (t, J = 7.5 Hz, 1H, ArH), 7.51 (d, J = 8.2 Hz, 2H, ArH), 7.54 (t, J = 7.6 Hz, 1H, ArH), 7.60 (d, J = 8.2 Hz, 1H, ArH), 7.90 (d, J = 9.1 Hz, 1H, ArH), 8.02 (d, J = 8.6 Hz, 1H, ArH), 8.17 (d, J = 8.5 Hz, 2H, ArH), 8.22 (d, J = 9.1 Hz, 1H, ArH), 8.52 (d, J = 7.9 Hz, 1H, ArH), 9.12 (d, J = 8.6 Hz, 1H, ArH) . ppm; ¹³C NMR (151 MHz, CDCl₃) δ = 14.5, 37.9, 109.6, 114.3, 114.5, 118.7, 120.1, 121.9, 123.5, 124.0, 124.7, 128.6, 128.7, 129.0, 131.9, 135.0, 137.2, 138.5, 139.3, 144.8 ppm; HRMS (ESI) m/z : calcd. for C₂₁H₁₇NO₂[M + H⁺]: 316.1332, found 316.1336.

6-(3,4-Dimethoxyphenyl)-[1,3]dioxolo[4,5-g]quinoline (26b): Yield = 69%, (0.136 g from 0.10 g) as a yellow solid; m.p; 160-162 °C; R_f = 0.5 (hexane/EtOAc, 70:30, v/v); IR (KBr): ν_{\max} (cm)⁻¹ = 1595 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 3.95 (d, J = 1.8 Hz, 3H, -OCH₃), 4.03 (d, J = 1.2 Hz, 3H, -OCH₃), 6.10 (d, J = 1.8 Hz, 2H, -OCH₂O-), 6.95-6.98 (m, 1H, ArH), 7.05 (d, J = 1.8 Hz, 1H, ArH), 7.44 (s, 1H, ArH), 7.59-7.61 (m, 1H, ArH), 7.79 (s, 1H, ArH), 7.98 (d, J = 6.6 Hz, 1H, ArH) ppm; ¹³C NMR (151 MHz, CDCl₃) δ = 56.1, 101.9, 102.7, 106.2, 110.5, 111.2, 117.0, 120.0, 123.9, 132.9, 135.7, 146.6, 147.6, 149.6, 150.2, 150.9, 155.0 ppm; HRMS (ESI) m/z : calcd. for C₁₈H₁₅NO₄ [M + H⁺]: 310.1074, found: 310.1079.

6-(Naphthalen-1-yl)-[1,3]dioxolo[4,5-g]quinoline (27b) Yield = 67% (0.147 g from 0.100 g) as a brown solid; m.p 142-144 °C; R_f = 0.55 (hexane/EtOAc, 80:20, v/v); IR (neat): ν_{\max} (cm)⁻¹ = 1613 (C=N); ¹H NMR (600 MHz, CDCl₃) δ = 6.14 (s, 2H, -OCH₂O), 7.15 (s, 1H, ArH), 7.45-

7.47 (m, 1H, ArH), 7.49– 7.52 (m, 2H, ArH), 7.55 (d, $J = 8.3$ Hz, 1H, ArH), 7.57-7.59 (m, 1H, ArH), 7.92(d, $J = 4.3$ Hz, 1H, ArH), 7.93(d, $J = 5.0$ Hz, 1H, ArH), 8.09 (d, $J = 8.3$ Hz, 1H, ArH), 8.12 (d, $J = 8.5$ Hz, 1H, ArH) . ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 101.9, 102.7, 106.3, 121.7, 124.1, 125.5, 125.9, 126.0, 126.6, 127.7, 128.5, 129.0, 131.4, 134.1, 135.2, 138.9, 146.5, 148.1, 151.1, 157.3$ ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{20}\text{H}_{13}\text{NO}_2$ [M + H $^+$]:300.1019, found 300.1023.

7-(3,4-Dimethoxyphenyl)-2,3-dihydro-[1,4]dioxino[2,3-g]quinoline (26c): Yield = 43% (0.103 g from 0.10 g) as a light yellow solid; m.p. 128-130 °C; $R_f = 0.67$ (hexane/EtOAc, 70:30, v/v); IR (neat): ν_{max} (cm^{-1}) = 1618 (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 3.95$ (s, 3H, $-\text{OCH}_3$), 4.04 (s, 3H, $-\text{OCH}_3$), 4.38 (d, $J = 2.6$ Hz, 4H, $-\text{OCH}_2\text{CH}_2\text{O}-$), 6.98 (d, $J = 8.3$ Hz, 1H, ArH), 7.19 (s, 1H, ArH), 7.62 – 7.58 (m, 2H, ArH), 7.67 (d, $J = 8.6$ Hz, 1H, ArH), 7.81 (s, 1H, ArH), 7.98 (d, $J = 8.6$ Hz, 1H, ArH) ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 56.1, 64.5, 110.4, 111.1, 111.7, 114.4, 117.2, 120.0, 123.0, 133.0, 135.2, 144.1, 144.8, 147.1, 149.4, 150.2, 155.5$ ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{19}\text{H}_{17}\text{NO}_4$ [M + H $^+$]: 324.1230, found 324.1234.

2-(3,4-Dimethoxyphenyl)-5,7-dimethoxyquinoline (26d): Yield = 47% (0.068 g from 0.10 g) as a light yellow solid; m.p. 138-140 °C; $R_f = 0.56$ (hexane/EtOAc, 90:10, v/v); IR (neat): ν_{max} (cm^{-1}) = 1619 (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 3.96$ (s, 3H, $-\text{OCH}_3$), 3.97 (s, 3H, $-\text{OCH}_3$), 3.98 (s, 3H, $-\text{OCH}_3$), 4.05 (s, 3H, $-\text{OCH}_3$), 6.49 (d, $J = 2.2$ Hz, 1H, ArH), 6.99 (d, $J = 8.3$ Hz, 1H, ArH), 7.10 (d, $J = 2.3$ Hz, 1H, ArH), 7.64 (d, $J = 2.0$ Hz, 1H, ArH), 7.66 (d, $J = 8.8$ Hz, 1H, ArH), 7.83 (d, $J = 2.1$ Hz, 1H, ArH), 8.45 (d, $J = 8.6$ Hz, 1H, ArH) ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 55.7, 55.8, 56.1, 56.1, 97.9, 100.0, 110.5, 111.2, 115.4, 115.6, 120.3, 131.4, 132.8, 149.4, 150.4, 150.5, 156.1, 157.2, 161.2$ ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{19}\text{H}_{19}\text{NO}_4$ [M + H $^+$]: 326.1387, found 326.1391.

2-(3,4-Dimethoxyphenyl)-5,8-dimethoxyquinoline (26e), Yield = 24% (0.034 g from 0.10 g) as a brown oily; $R_f = 0.56$ (hexane/EtOAc, 90:10, v/v); IR (neat): ν_{max} (cm^{-1}) = 1624 (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 3.94$ (s, 3H, $-\text{OCH}_3$), 3.96 (s, 3H, $-\text{OCH}_3$), 4.04 (s, 3H, $-\text{OCH}_3$), 4.06 (s, 3H, $-\text{OCH}_3$), 6.72 (d, $J = 8.4$ Hz, 1H, ArH), 6.95 (d, $J = 8.4$ Hz, 1H, ArH), 6.97 (d, $J = 8.4$ Hz, 1H, ArH), 7.69 (dd, $J_1 = 8.3$ Hz, $J_2 = 2.1$ Hz, 1H, ArH), 7.86 (d, $J = 8.9$ Hz, 2H, ArH), 8.56 (d, $J = 8.8$ Hz, 1H, ArH) ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 55.8, 56.0, 56.1, 56.2, 103.4, 110.2, 110.7, 111.1, 118.3, 120.3, 120.6, 131.7, 132.7, 140.5, 148.9, 149.4, 149.6, 150.4, 156.3$ ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{19}\text{H}_{19}\text{NO}_4$ [M + H $^+$]: 326.1387, found 326.1391.

2-Phenylbenzo[h]quinoline (15g), Yield = 45% (0.08 g from 0.10 g) as a brown solid; m.p. 64-66 °C; $R_f = 0.69$ (hexane/EtOAc, 95:5, v/v); IR (neat): $\nu_{\max} (\text{cm})^{-1} = 1600$ (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 7.46 - 7.40$ (m, 1H, ArH), 7.57 (t, $J = 7.7$ Hz, 2H, ArH), 7.70 – 7.72 (m, 2H, ArH), 7.75 – 7.77 (m, 1H, ArH), 7.80 (d, $J = 8.7$ Hz, 1H, ArH), 7.92 (d, $J = 7.9$ Hz, 1H, ArH), 8.02 (d, $J = 8.3$ Hz, 1H, ArH), 8.24 (d, $J = 8.3$ Hz, 1H, ArH), 8.36 (d, $J = 7.1$ Hz, 2H, ArH), 9.51 (d, $J = 8.1$ Hz, 1H) ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 119.0, 124.9, 125.2, 125.3, 127.0, 127.3, 127.6, 127.9, 128.3, 128.8, 129.0, 129.3, 131.0, 134.0, 136.7, 139.9, 146.4, 155.7$ ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{19}\text{H}_{13}\text{N}$ [$\text{M} + \text{H}^+$]: 256.1121, found 256.1125.

2-(4-Bromophenyl)benzo[h]quinoline (16g), Yield = 60% (0.08 g from 0.10 g) as a brown solid; m.p. 96-98 °C; $R_f = 0.77$ (hexane/EtOAc, 95:5, v/v); IR (neat): $\nu_{\max} (\text{cm})^{-1} = 1631$ (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 7.69$ (t, $J = 8.4$ Hz, 3H, ArH), 7.72 (d, $J = 6.5$ Hz, 1H, ArH), 7.76 (t, $J = 6.8$ Hz, 1H, ArH), 7.81 (d, $J = 8.8$ Hz, 1H, ArH), 7.92 (d, $J = 7.8$ Hz, 1H, ArH), 7.97 (d, $J = 8.3$ Hz, 1H, ArH), 8.22 (d, $J = 8.5$ Hz, 3H, ArH), 9.45 (d, $J = 8.1$ Hz, 1H, ArH) ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 118.6, 123.8, 124.7, 125.0, 125.8, 125.4, 127.9, 129.0, 129.2, 131.8, 132.0, 132.1, 134.0, 136.7, 136.9, 138.8, 146.4, 154.4$ ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{19}\text{H}_{12}\text{BrN}$ [$\text{M} + \text{H}^+$]: 334.0226, found: 334.0230.

2-(4-Chlorophenyl)benzo[h]quinoline (17g), Yield = 48% (0.096 g from 0.10 g) as a brown solid; m.p. 100-102 °C; $R_f = 0.76$ (hexane/EtOAc, 95:5, v/v); IR (neat): $\nu_{\max} (\text{cm})^{-1} = 1632$ (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 7.53$ (d, $J = 8.5$ Hz, 2H, ArH), 7.71 (t, $J = 8.8$ Hz, 2H, ArH), 7.76 (t, $J = 7.6$ Hz, 1H, ArH), 7.81 (d, $J = 8.8$ Hz, 1H, ArH), 7.92 (d, $J = 7.9$ Hz, 1H, ArH), 7.97 (d, $J = 8.3$ Hz, 1H, ArH), 8.23 (d, $J = 8.3$ Hz, 1H, ArH), 8.29 (d, $J = 8.5$ Hz, 2H, ArH), 9.46 (d, $J = 8.1$ Hz, 1H, ArH) ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 118.6, 124.8, 125.1, 125.4, 127.1, 127.8, 128.0, 128.4, 128.7, 129.1, 131.8, 134.0, 135.5, 136.7, 136.8, 138.3, 146.4, 154.3$ ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{19}\text{H}_{12}\text{ClN}$ [$\text{M} + \text{H}^+$]: 290.0731, found 290.0736.

2-(4-Fluorophenyl)benzo[h]quinoline (18g), Yield = 50% (0.096 g from 0.10 g) as an off white solid; m.p. 98-100 °C; $R_f = 0.66$ (hexane/EtOAc, 95:5, v/v); IR (neat): $\nu_{\max} (\text{cm})^{-1} = 1623$ (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 7.24$ (d, $J = 8.6$ Hz, 2H, ArH), 7.71 (t, $J = 8.0$ Hz, 2H, ArH), 7.76 (t, $J = 6.8$ Hz, 2H, ArH), 7.80 (d, $J = 8.7$ Hz, 1H, ArH), 7.92 (d, $J = 7.2$ Hz, 1H, ArH), 7.97 (d, $J = 8.3$ Hz, 1H, ArH), 8.23 (d, $J = 8.3$ Hz, 1H, ArH), 8.33 (d, $J = 5.5$ Hz, 1H, ArH), 8.35 (d, $J = 5.5$ Hz, 1H, ArH), 9.46 (d, $J = 8.1$ Hz, 1H, ArH) ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 115.8, 115.9$ ($J_{\text{C,F}} = 21.5$ Hz), 118.6, 124.8, 125.2, 127.0, 127.6, 127.9, 128.4, 129.3, 129.4 ($J_{\text{C,F}} = 8.4$ Hz), 131.8,

134.0, 136.0, 136.8, 146.3, 154.6, 163.1, 164.7 ($J_{C,F} = 247.3$ Hz) ppm; HRMS (ESI) m/z : calcd. for $C_{19}H_{12}FN$ [$M + H^+$]: 274.1027, found 274.1032.

2-(4-Methoxyphenyl)benzo[*h*]quinoline (19g), Yield = 40% (0.080 g from 0.10 g) as off white solid; m.p. 116-118 °C; $R_f = 0.52$ (hexane/EtOAc, 95:5, v/v); IR (neat): ν_{max} (cm)⁻¹ = 1590 (C=N); ¹H NMR (600 MHz, CDCl₃) $\delta = 3.92$ (s, 3H, -OCH₃), 7.09 (d, $J = 8.8$ Hz, 2H, ArH), 7.69 (t, $J = 7.4$ Hz, 2H, ArH), 7.75 (t, $J = 6.8$ Hz, 1H, ArH), 7.78 (d, $J = 8.7$ Hz, 1H, ArH), 7.91 (d, $J = 6.5$ Hz, 1H, ArH), 7.96 (d, $J = 8.4$ Hz, 1H, ArH), 8.20 (d, $J = 8.3$ Hz, 1H, ArH), 8.32 (d, $J = 8.7$ Hz, 2H, ArH), 9.49 (d, $J = 8.1$ Hz, 1H, ArH) ppm; ¹³C NMR (151 MHz, CDCl₃) $\delta = 55.5, 114.3, 118.4, 124.8, 125.3, 126.9, 127.1, 127.9, 128.2, 128.9, 131.9, 132.6, 134.0, 136.5, 146.3, 153.3, 160.9$ ppm; HRMS (ESI) m/z : calcd. for $C_{20}H_{15}NO$ [$M + H^+$]: 286.1226, found 286.1230.

2-([1,1'-Biphenyl]-4-yl)benzo[*h*]quinoline (22g). Yield = 66% (0.090 g from 0.10 g) as a yellow solid; m.p. 138-140 °C; $R_f = 0.71$ (hexane/EtOAc, 80:20, v/v); IR (neat): ν_{max} (cm)⁻¹ = 1635 (C=N); ¹H NMR (600 MHz, CHLOROFORM-*D*) $\delta = 7.41$ (t, $J = 7.4$ Hz, 1H, ArH), 7.50 (t, $J = 7.6$ Hz, 2H, ArH), 7.71 – 7.74 (m, 4H, ArH), 7.77 (d, $J = 8.2$ Hz, 1H, ArH), 7.79 – 7.82 (m, 3H, ArH), 7.93 (d, $J = 7.8$ Hz, 1H, ArH), 8.06 (d, $J = 8.2$ Hz, 1H, ArH), 8.24 (d, $J = 8.2$ Hz, 1H, ArH), 8.44 (d, $J = 8.1$ Hz, 2H, ArH), 9.54 (d, $J = 8.1$ Hz, 1H, ArH) ppm; ¹³C NMR (150 MHz, CHLOROFORM-*D*) $\delta = 118.9, 124.9, 125.2, 125.3, 127.0, 127.3, 127.6, 127.7, 127.9, 128.0, 128.3, 129.0, 132.0, 134.0, 136.7, 138.8, 140.8, 142.1, 146.4, 155.2$ ppm; HRMS (ESI) m/z : calcd. for $C_{25}H_{17}N$ [$M+H^+$]: 332.1434; found 332.1438.

2-(3,4-Dichlorophenyl)benzo[*h*]quinoline (25g). Yield = 56% (0.077 g from 0.10 g) as a light brown solid; m.p. 145-147 °C; $R_f = 0.48$ (hexane/EtOAc, 80:20, v/v); IR (neat): ν_{max} (cm)⁻¹ = 1624 (C=N); ¹H NMR (600 MHz, CDCl₃) $\delta = 7.61$ (d, $J = 8.3$ Hz, 1H, ArH), 7.70 (d, $J = 8.7$ Hz, 1H, ArH), 7.73 (d, $J = 6.6$ Hz, 1H, ArH), 7.76 – 7.80 (m, 1H, ArH), 7.82 (d, $J = 8.7$ Hz, 1H, ArH), 7.91 – 7.94 (m, 1H, ArH), 7.95 (d, 1H, ArH), 8.14 (dd, $J = 8.3, 2.2$ Hz, 1H, ArH), 8.24 (d, $J = 8.3$ Hz, 1H, ArH), 8.46 (d, $J = 2.1$ Hz, 1H, ArH), 9.44 (d, $J = 8.2$ Hz, 1H, ArH) ppm; ¹³C NMR (150 MHz, CDCl₃) $\delta = 118.5, 124.8, 125.1, 125.7, 126.5, 127.3, 128.0, 128.2, 128.6, 129.4, 130.8, 131.7, 133.2, 133.5, 134.1, 136.9, 139.8, 146.4, 153.0$ ppm; HRMS (ESI) m/z : calcd. for $C_{19}H_{11}Cl_2N$ [$M+H^+$]: 324.0341; found 324.0345.

2-(3,4-Dimethoxyphenyl)benzo[*h*]quinoline (26g), Yield = 54% (0.080 g from 0.10 g) as a yellow solid; m.p. 110-112 °C; $R_f = 0.43$ (hexane/EtOAc, 90:10, v/v); IR (neat): ν_{max} (cm)⁻¹ = 1590 (C=N); ¹H NMR (600 MHz, CDCl₃) $\delta = 3.99$ (s, 3H, -OCH₃), 4.11 (s, 3H, OCH₃), 7.04 (d, $J =$

8.3 Hz, 1H, ArH), 7.70 (t, $J = 8.1$ Hz, 2H, ArH), 7.75 (t, $J = 6.8$ Hz, 1H, ArH), 7.78 (d, $J = 8.8$ Hz, 1H, ArH), 7.82 (dd, $J_1 = 8.3$, $J_2 = 2.1$ Hz, 1H, ArH), 7.92 (d, $J = 7.8$ Hz, 1H, ArH), 7.97 (d, $J = 8.3$ Hz, 1H, ArH), 8.07 (s, 1H, ArH), 8.20 (d, $J = 8.3$ Hz, 1H, ArH), 9.46 (d, $J = 8.2$ Hz, 1H) ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 56.2, 110.6, 111.3, 118.6, 120.2, 124.7, 125.0, 125.2, 127.0, 127.2, 127.9, 128.2, 131.9, 132.9, 134.1, 136.6, 146.2, 149.5, 150.5, 155.3$ ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{21}\text{H}_{17}\text{NO}_2$ [$\text{M} + \text{H}^+$]: 316.1332, found 316.1336.

2-(4-Chlorophenyl)-10-ethyl-5 α ,10 α -dihydro-10H-pyrido[2,3-*b*]carbazole (17h), Yield = 33% (0.031 g from 0.05 g) as a yellow solid; m.p. 182-184 °C; $R_f = 0.68$ (hexane/EtOAc, 90:10, v/v); IR (neat): ν_{max} (cm) $^{-1} = 1629$ (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 1.52$ (t, $J = 7.3$ Hz, 3H, $-\text{NCH}_2\text{CH}_3$), 4.54 (q, $J = 7.3$ Hz, 2H, $-\text{NCH}_2\text{CH}_3$), 7.41 (t, $J = 7.5$ Hz, 1H, ArH), 7.51 (d, $J = 8.2$ Hz, 2H, ArH), 7.54 (t, $J = 7.6$ Hz, 1H, ArH), 7.60 (d, $J = 8.2$ Hz, 1H, ArH), 7.90 (d, $J = 9.1$ Hz, 1H, ArH), 8.02 (d, $J = 8.6$ Hz, 1H, ArH), 8.17 (d, $J = 8.5$ Hz, 2H, ArH), 8.22 (d, $J = 9.1$ Hz, 1H, ArH), 8.52 (d, $J = 7.9$ Hz, 1H, ArH), 9.12 (d, $J = 8.6$ Hz, 1H, ArH) . ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 14.5, 37.9, 109.6, 114.3, 114.5, 118.7, 120.1, 121.9, 123.5, 124.0, 124.7, 128.6, 128.7, 129.0, 131.9, 135.0, 137.2, 138.5, 139.3, 144.8$ ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{23}\text{H}_{19}\text{ClN}_2$ [$\text{M} + \text{H}^+$]: 357.1153, found 357.1158.

7-Ethyl-3-(4-fluorophenyl)-7H-pyrido[2,3-*c*]carbazole (18h), Yield = 42% (0.068 g from 0.10 g) as a yellow solid; m.p. 188-190 °C; $R_f = 0.69$ (hexane/EtOAc, 90:10, v/v); IR (neat): ν_{max} (cm) $^{-1} = 1627$ (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 1.53$ (t, $J = 7.3$ Hz, 3H, $-\text{NCH}_2\text{CH}_3$), 4.55 (q, $J = 7.3$ Hz, 2H, $-\text{NCH}_2\text{CH}_3$), 7.23 (t, $J = 8.6$ Hz, 2H, ArH), 7.41 (t, $J = 7.4$ Hz, 1H, ArH), 7.54 (t, $J = 7.6$ Hz, 1H, ArH), 7.60 (d, $J = 8.2$ Hz, 1H, ArH), 7.91 (d, $J = 9.0$ Hz, 1H, ArH), 8.01 (d, $J = 8.6$ Hz, 1H, ArH), 8.23 – 8.20 (m, 3H, ArH), 8.53 (d, $J = 7.9$ Hz, 1H, ArH), 9.12 (d, $J = 8.7$ Hz, 1H, ArH) ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 14.6, 37.9, 109.7, 114.3, 114.5, 115.8, 115.9$ ($J_{\text{C,F}} = 21.4$ Hz), 118.8, 120.1, 121.7, 121.9, 123.5, 123.9, 124.8, 128.6, 129.1, 129.2 ($J_{\text{C,F}} = 8.1$ Hz), 132.0, 136.3, 137.2, 139.3, 144.8, 152.7, 162.8, 164.4 ($J_{\text{C,F}} = 246.5$ Hz) ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{23}\text{H}_{17}\text{FN}_2$ [$\text{M} + \text{H}^+$]: 341.1449, found 341.1445.

3-(3-Bromophenyl)-7-ethyl-7H-pyrido[2,3-*c*]carbazole (24h), Yield = 51% (0.043 g from 0.05 g) as a yellow solid; m.p. 142-144 °C; $R_f = 0.53$ (hexane/EtOAc, 80:20, v/v); IR (neat): ν_{max} (cm) $^{-1} = 1629$ (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 1.52$ (t, $J = 7.3$ Hz, 3H, $-\text{NCH}_2\text{CH}_3$), 4.54 (q, $J = 7.3$ Hz, 2H, $-\text{NCH}_2\text{CH}_3$), 7.39 – 7.42 (m, 2H, ArH), 7.54 (t, $J = 7.6$ Hz, 1H, ArH), 7.59 (dd, $J_1 = 13.7$ Hz, $J_2 = 8.2$ Hz, 2H, ArH), 7.91 (d, $J = 9.1$ Hz, 1H, ArH), 8.01 (d, $J = 8.7$ Hz, 1H, ArH), 8.12

(d, $J = 7.7$ Hz, 1H, ArH), 8.23 (d, $J = 9.1$ Hz, 1H, ArH), 8.42 (s, 1H, ArH), 8.52 (d, $J = 7.9$ Hz, 1H, ArH), 9.12 (d, $J = 8.8$ Hz, 1H, ArH) ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 14.6, 37.9, 109.7, 114.4, 114.5, 118.9, 120.2, 121.9, 123.3, 123.5, 124.2, 124.8, 125.9, 128.7, 130.4, 130.4, 131.8, 132.0, 137.3, 139.3, 142.2, 144.8, 152.0$ ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{23}\text{H}_{17}\text{BrN}_2$ $[\text{M} + \text{H}^+]$: 401.0648, found 401.0644.

3-(6,7-Dimethoxyquinolin-2-yl)-9-ethyl-9H-carbazole (28a), Yield = 62% (0.078 g from 0.05 g) as a green solid; m.p. = 178-180 °C; $R_f = 0.6$ (hexane/EtOAc, 70:30, v/v); IR (neat): ν_{max} (cm) $^{-1} = 1591$ (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 1.47$ (t, $J = 7.3$ Hz, 3H, $-\text{NCH}_2\text{CH}_3$), 4.04 (s, 3H, $-\text{OCH}_3$), 4.09 (s, 3H, $-\text{OCH}_3$), 4.41 (q, $J = 7.3$ Hz, 2H, $-\text{NCH}_2\text{CH}_3$), 7.07 (s, 1H, ArH), 7.28 (d, $J = 7.4$, 1H, ArH), 7.44 (d, $J = 8.2$ Hz, 1H, ArH), 7.50 (dd, $J = 14.5, 7.8$ Hz, 2H, ArH), 7.58 (s, 1H, ArH), 7.88 (d, $J = 8.5$ Hz, 1H, ArH), 8.06 (d, $J = 8.4$ Hz, 1H, ArH), 8.23 (d, $J = 7.7$ Hz, 1H, ArH), 8.28 (d, $J = 8.5$ Hz, 1H, ArH), 8.89 (s, 1H, ArH) ppm; ^{13}C NMR (151 MHz, CHLOROFORM-D) $\delta = 13.9, 37.8, 56.1, 56.3, 105.2, 108.4, 108.7, 117.5, 119.2, 119.6, 120.8, 122.4, 123.5, 123.6, 125.3, 125.9, 131.0, 135.0, 140.6, 140.7, 145.4, 149.4, 152.6, 156.4$ ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{25}\text{H}_{22}\text{N}_2\text{O}_2$ $[\text{M} + \text{H}^+]$: 383.1754, found 383.1759.

6-(9-Ethyl-9H-carbazol-3-yl)-[1,3]dioxolo[4,5-g]quinoline (28b), Yield = 56% (0.065g from 0.10 g) as a yellow solid; m.p. = 178-180 °C; $R_f = 0.69$ (hexane/EtOAc, 90:10, v/v); IR (neat): ν_{max} (cm) $^{-1} = 1628$ (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 1.47$ (t, $J = 7.2$ Hz, 3H, $-\text{NCH}_2\text{CH}_3$), 4.42 (q, $J = 7.3$ Hz, 2H, $-\text{NCH}_2\text{CH}_3$), 6.11 (d, $J = 1.6$ Hz, 2H, $-\text{OCH}_2\text{O}-$), 7.08 (s, 1H, ArH), 7.28 (d, $J = 7.4$ Hz, 1H, ArH), 7.44 (d, $J = 8.2$ Hz, 1H, ArH), 7.52–7.48 (m, 3H, ArH), 7.87 (d, $J = 8.5$ Hz, 1H, ArH), 8.03 (d, $J = 8.3$ Hz, 1H, ArH), 8.22 (d, $J = 7.7$ Hz, 1H, ArH), 8.28 (d, $J = 8.5$ Hz, 1H, ArH), 8.87 (s, 1H, ArH), . ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 14.0, 37.8, 76.9, 77.4, 101.7, 102.8, 106.2, 108.8, 117.5, 119.3, 119.7, 120.9, 123.5, 123.6, 123.7, 125.4, 126.0, 130.9, 135.6, 140.6, 140.7, 140.7, 146.7, 147.4, 150.9, 156.3$ ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{24}\text{H}_{18}\text{N}_2\text{O}_2$ $[\text{M} + \text{H}^+]$: 367.1441, found 367.1445.

2-(9-Ethyl-9H-carbazol-3-yl)benzo[g]quinoline (28g), Yield = 51% (0.065g from 0.10 g) as a dark yellow solid; m.p. 100-102 °C; $R_f = 0.88$ (hexane/EtOAc, 70:30, v/v); IR (neat): ν_{max} (cm) $^{-1} = 1625$ (C=N); ^1H NMR (600 MHz, CDCl_3) $\delta = 1.50$ (t, $J = 7.3$ Hz, 3H, $-\text{NCH}_2\text{CH}_3$), 4.45 (q, $J = 7.3$ Hz, 2H, $-\text{NCH}_2\text{CH}_3$), 7.31 (t, $J = 7.3$ Hz, 1H, ArH), 7.47 (d, $J = 8.2$ Hz, 1H, ArH), 7.53 – 7.51 (m, 1H, ArH), 7.58 (d, $J = 8.5$ Hz, 1H, ArH), 7.72 (s, 1H, ArH), 7.73 (s, 1H, ArH), 7.78 (s,

1H, ArH), 7.81 – 7.79 (m, 1H, ArH), 7.93 (dd, $J = 7.9, 1.3$ Hz, 1H, ArH), 8.16 (d, $J = 8.3$ Hz, 1H, ArH), ArH), 9.07 (d, $J = 1.7$ Hz, 1H, ArH), 9.59 (d, $J = 9.4$ Hz, 1H, ArH), ppm; ^{13}C NMR (151 MHz, CDCl_3) $\delta = 55.9, 55.9, 56.5, 103.9, 108.0, 118.2, 120.7, 129.0, 129.1, 132.1, 135.5, 138.2, 140.6, 148.9, 149.7, 155.4$ ppm; HRMS (ESI) m/z : calcd. for $\text{C}_{27}\text{H}_{20}\text{N}_2$ $[\text{M} + \text{H}^+]$: 373.1699, found 373.1694.

3. ^1H and ^{13}C NMR Spectra of Synthesized Compounds

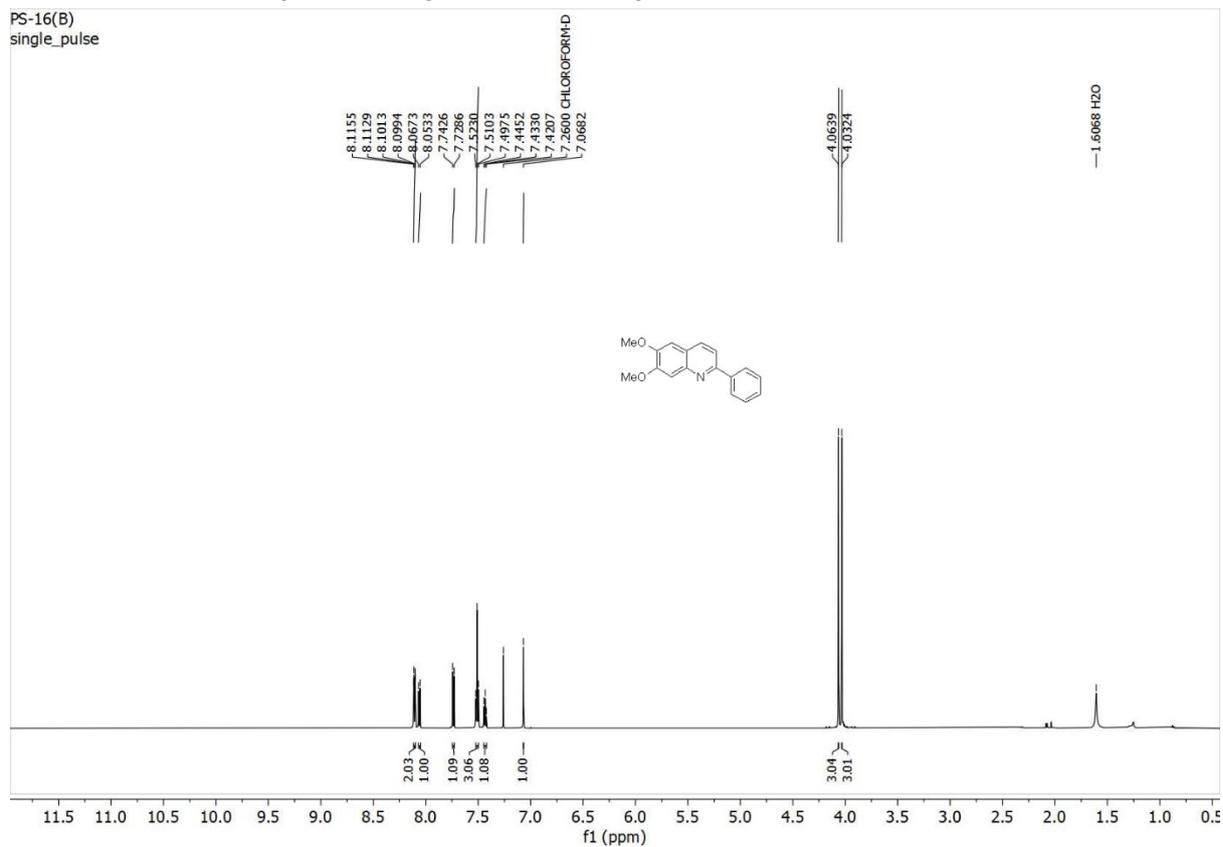


Figure S1. ^1H NMR Spectrum of **15a** in CDCl_3 .

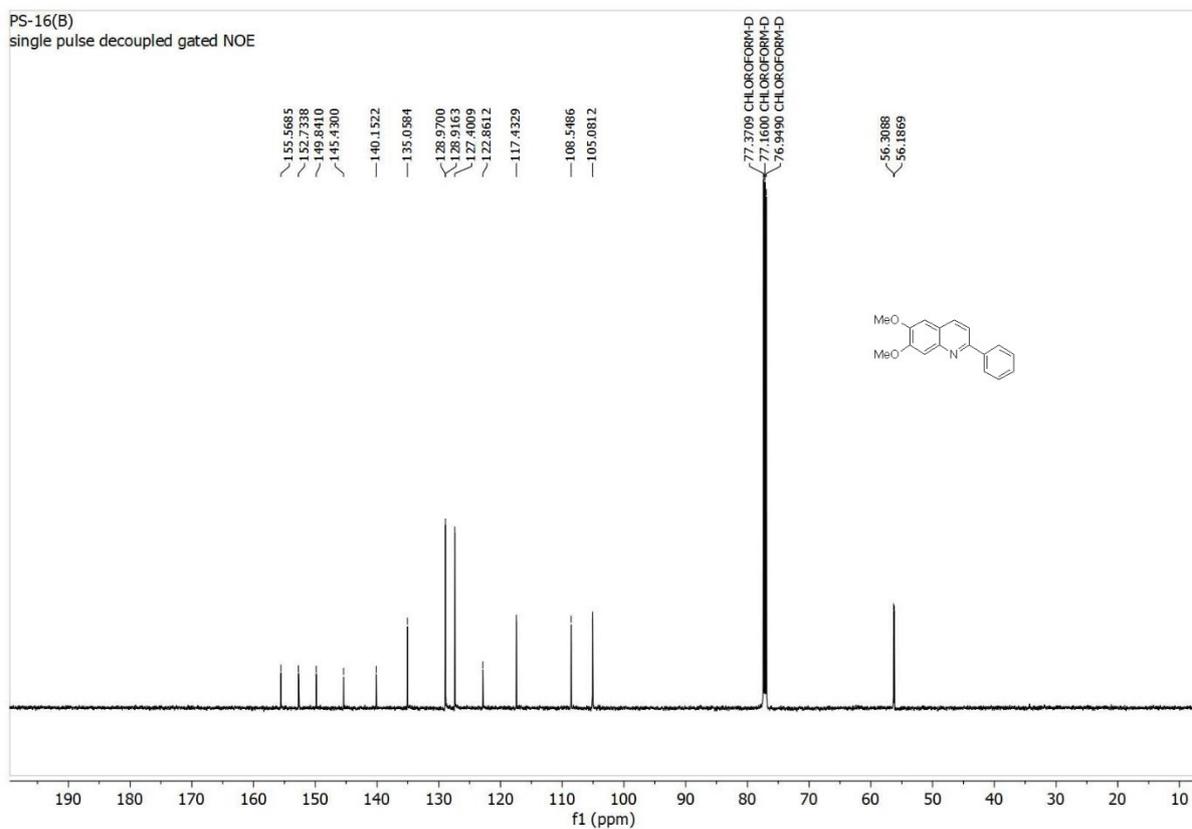


Figure S2. ^{13}C NMR Spectrum of **15a** in CDCl_3 .

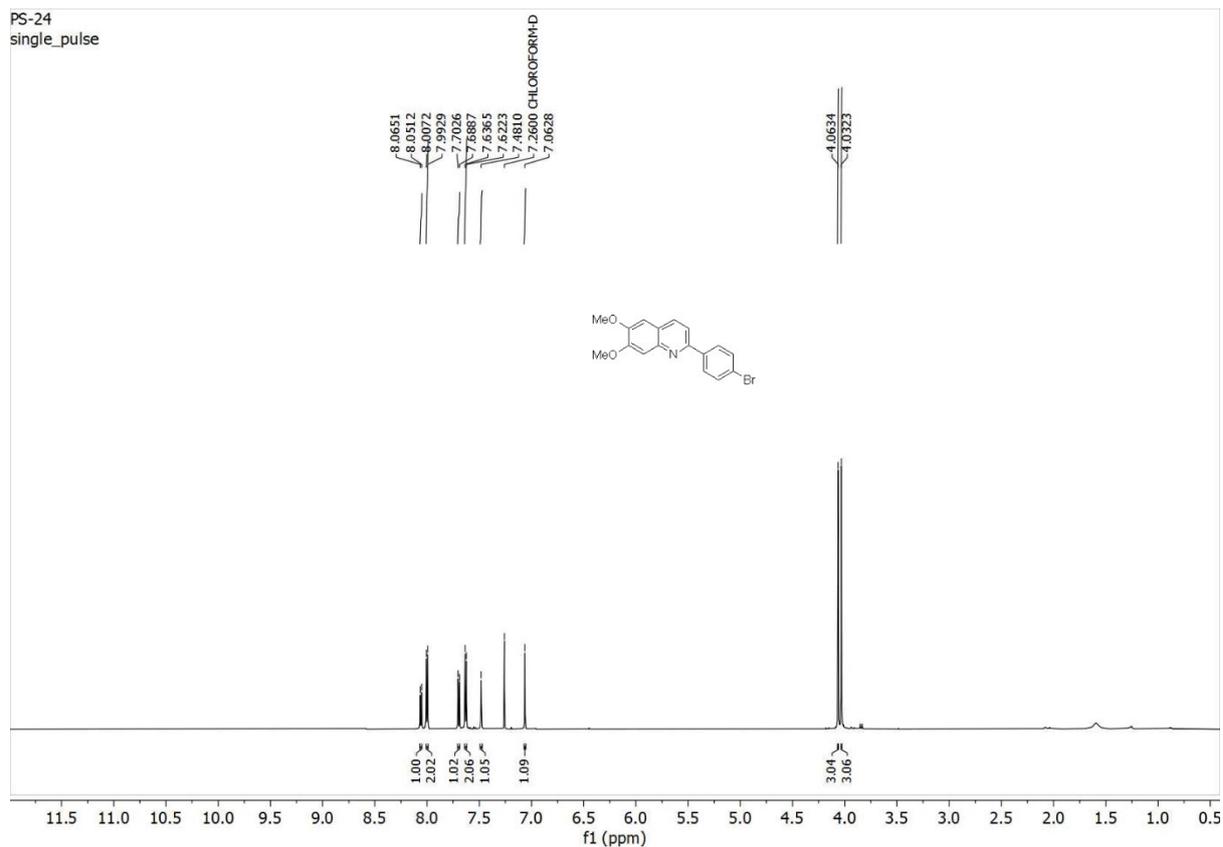


Figure S3. ^1H NMR Spectrum of **16a** in CDCl_3 .

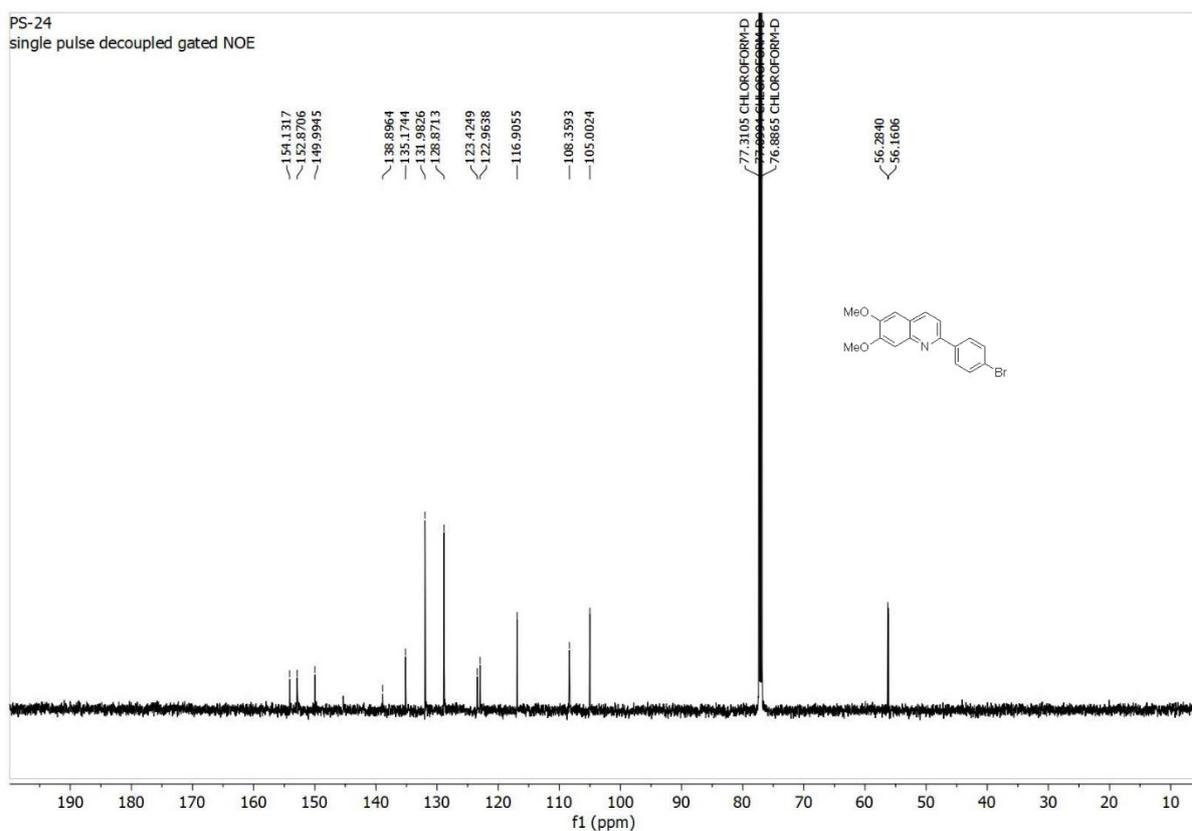


Figure S4. ^{13}C NMR Spectrum of **16a** in CDCl_3 .

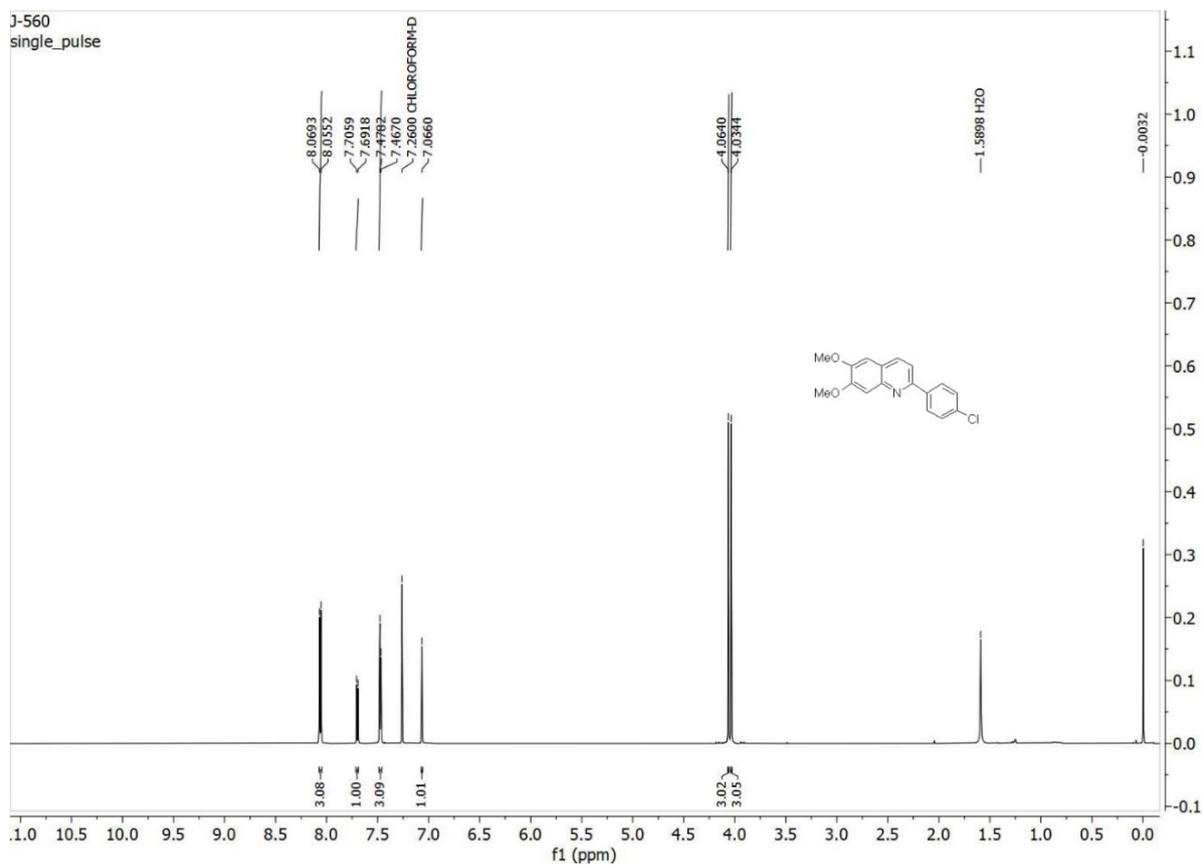


Figure S5. ^1H NMR Spectrum of **17a** in CDCl_3 .

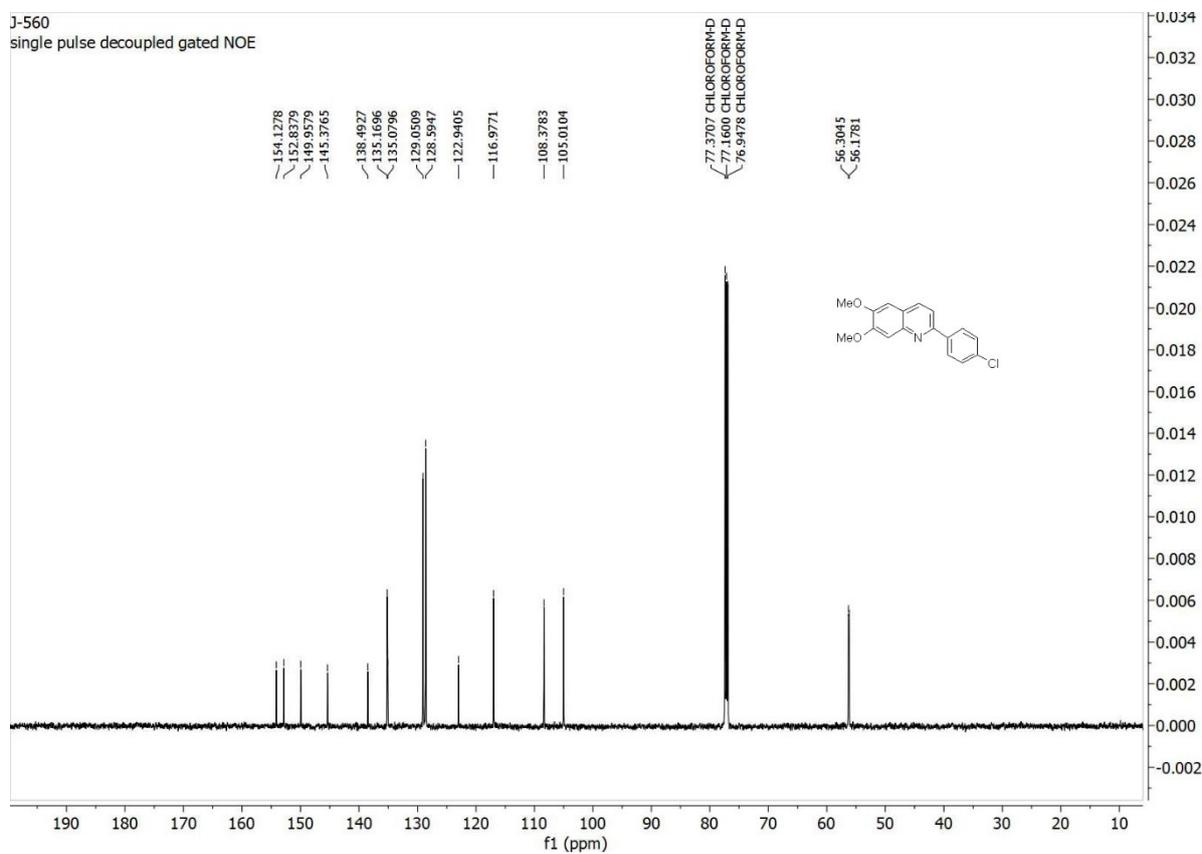


Figure S6. ^{13}C NMR Spectrum of **17a** in CDCl_3 .

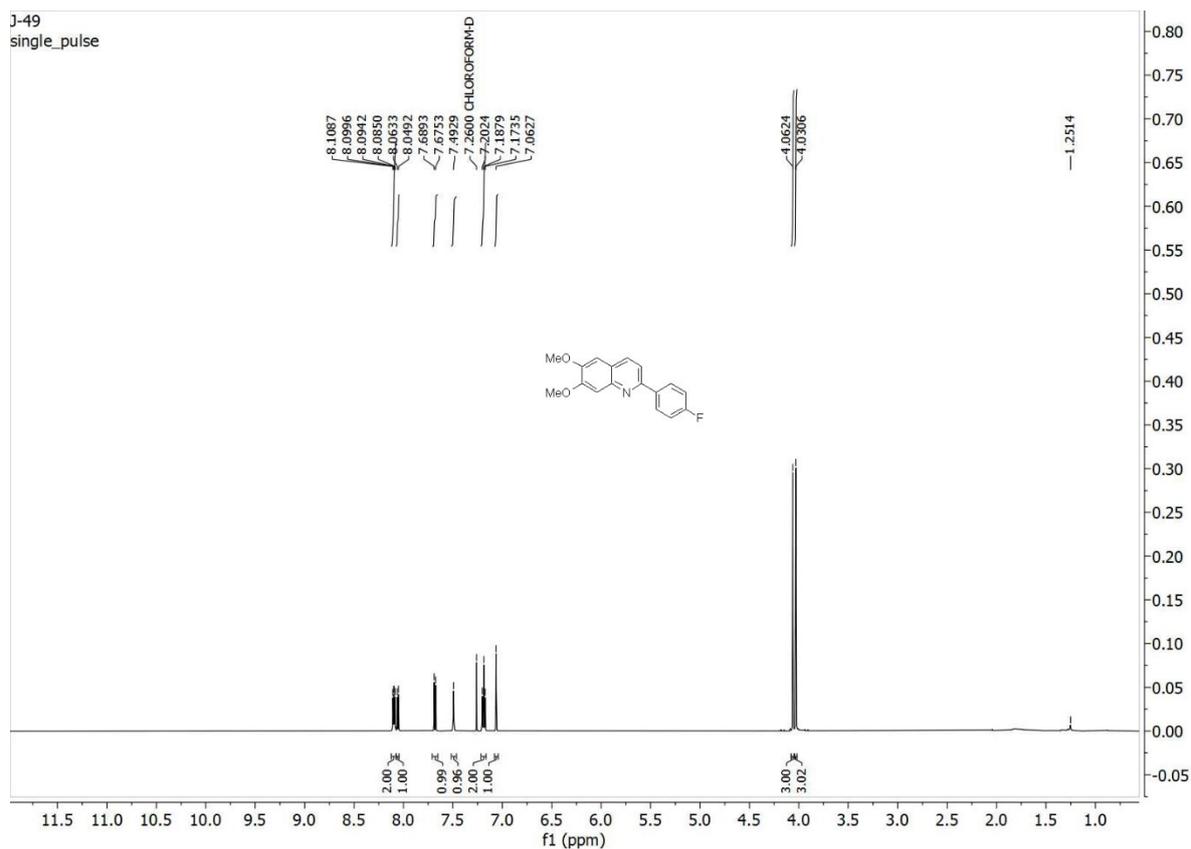


Figure S7. ^1H NMR Spectrum of **18a** in CDCl_3 .

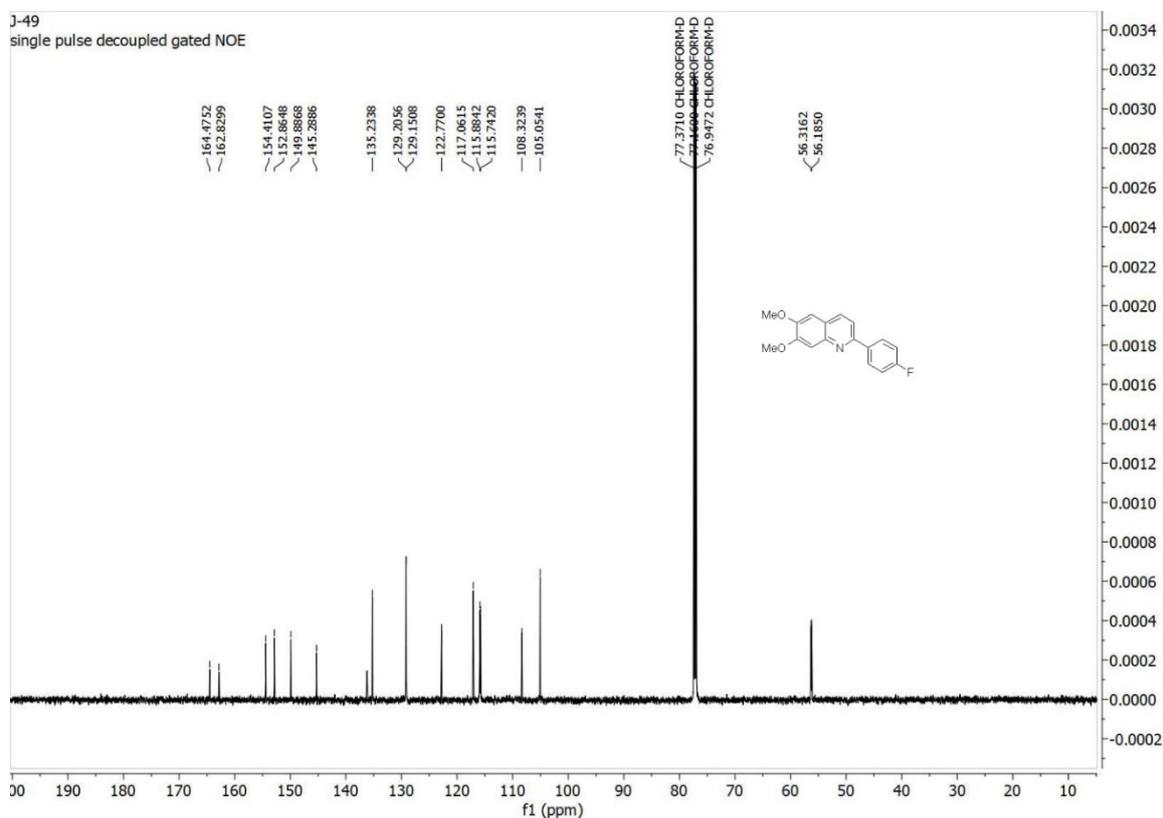


Figure S8. ^{13}C NMR Spectrum of **18a** in CDCl_3 .

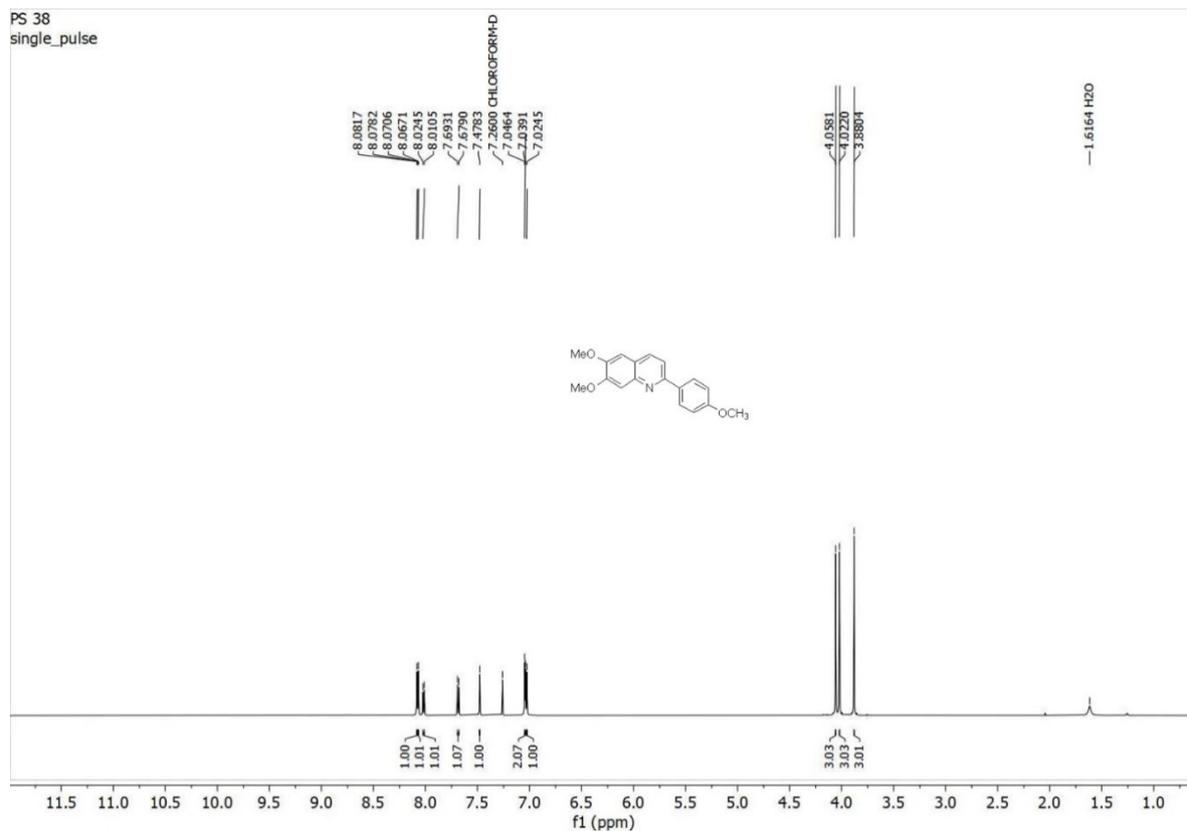


Figure S9. ¹H NMR Spectrum of **19a** in CDCl₃.

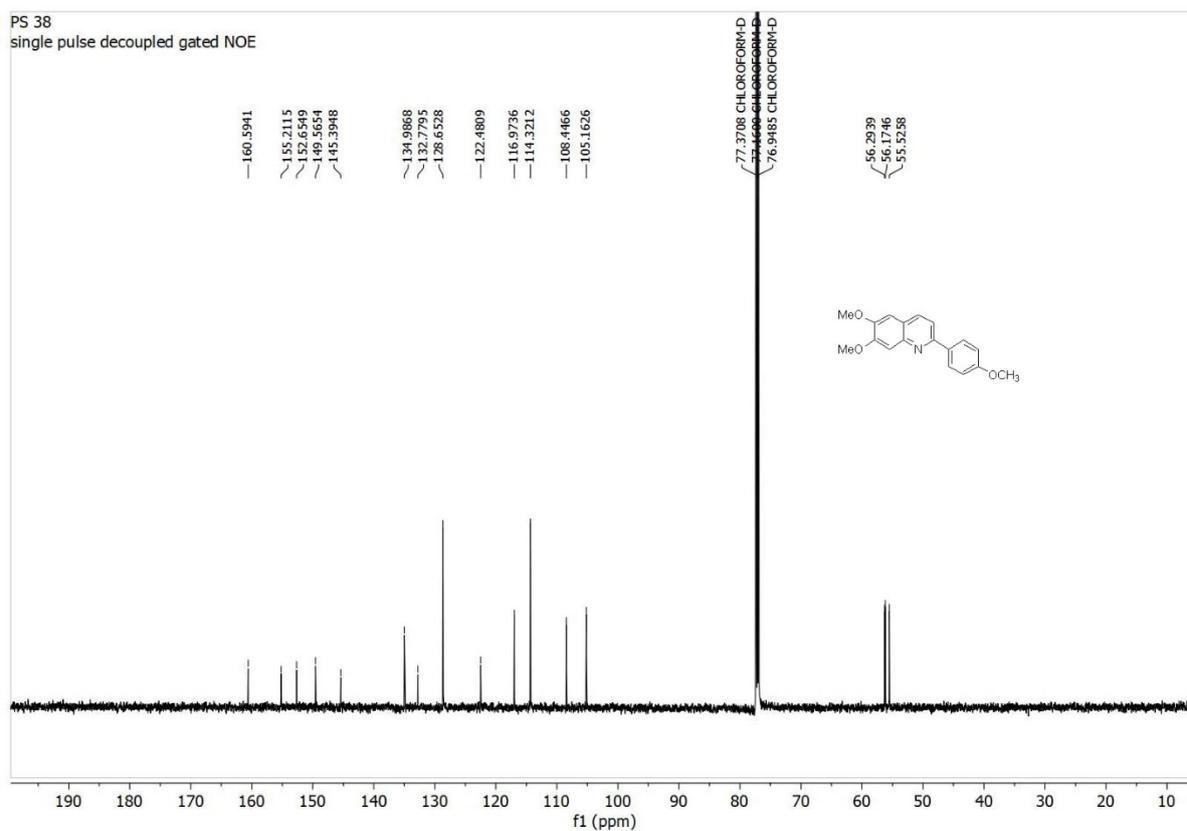


Figure S10. ¹³C NMR Spectrum of **19a** in CDCl₃.

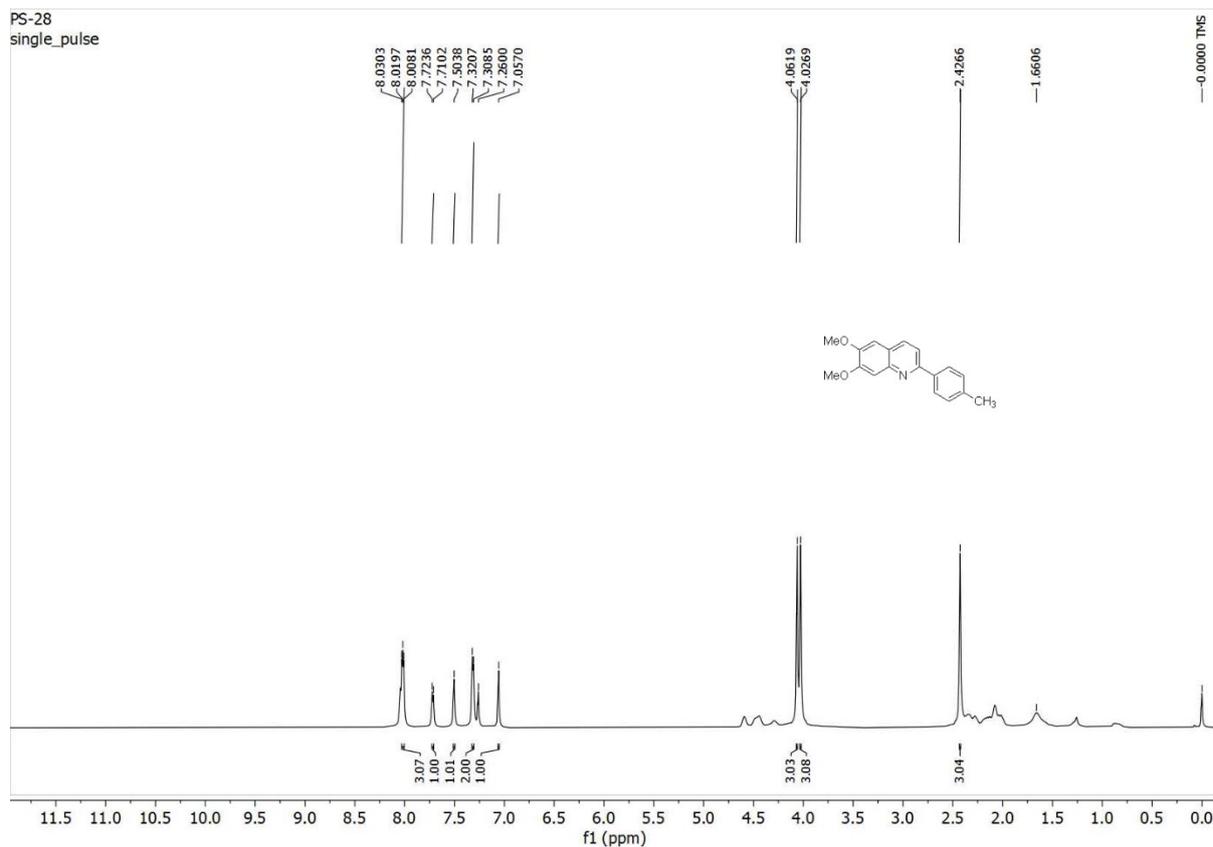


Figure S11. ^1H NMR Spectrum of **20a** in CDCl_3 .

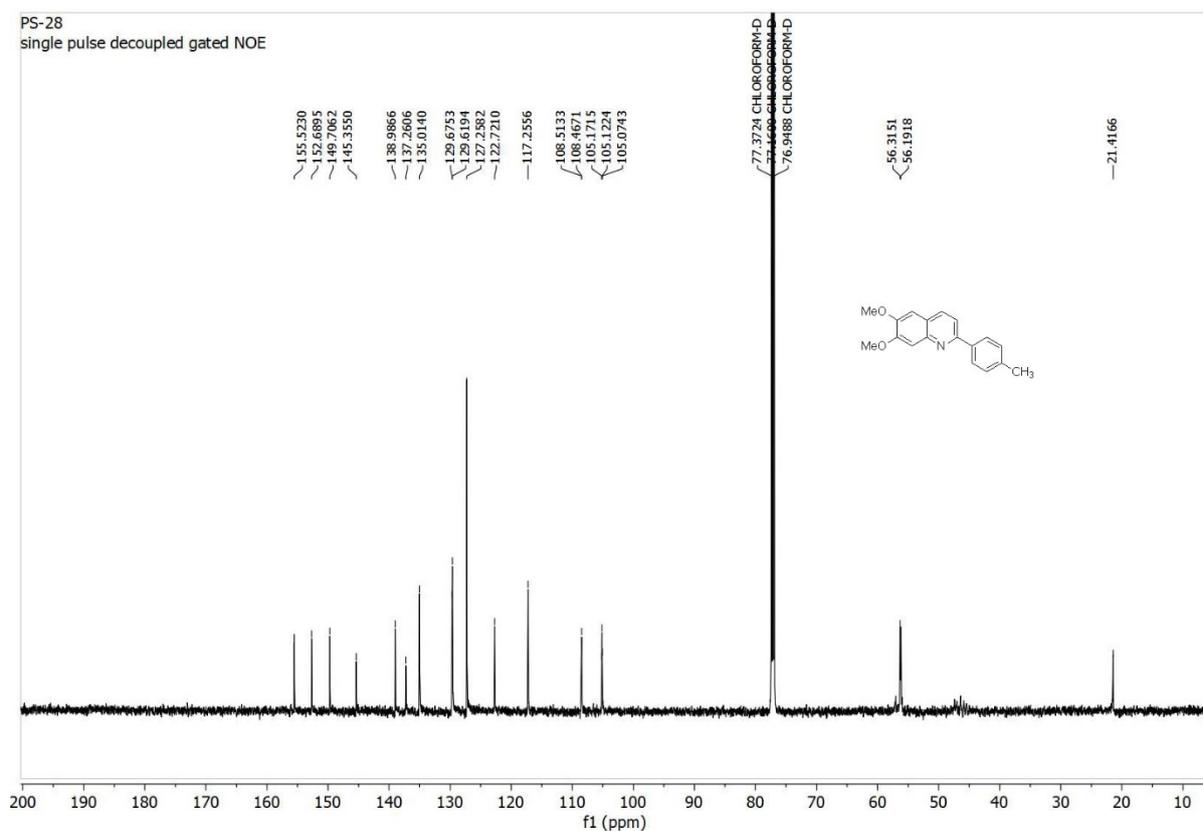


Figure S12. ^{13}C NMR Spectrum of **20a** in CDCl_3 .

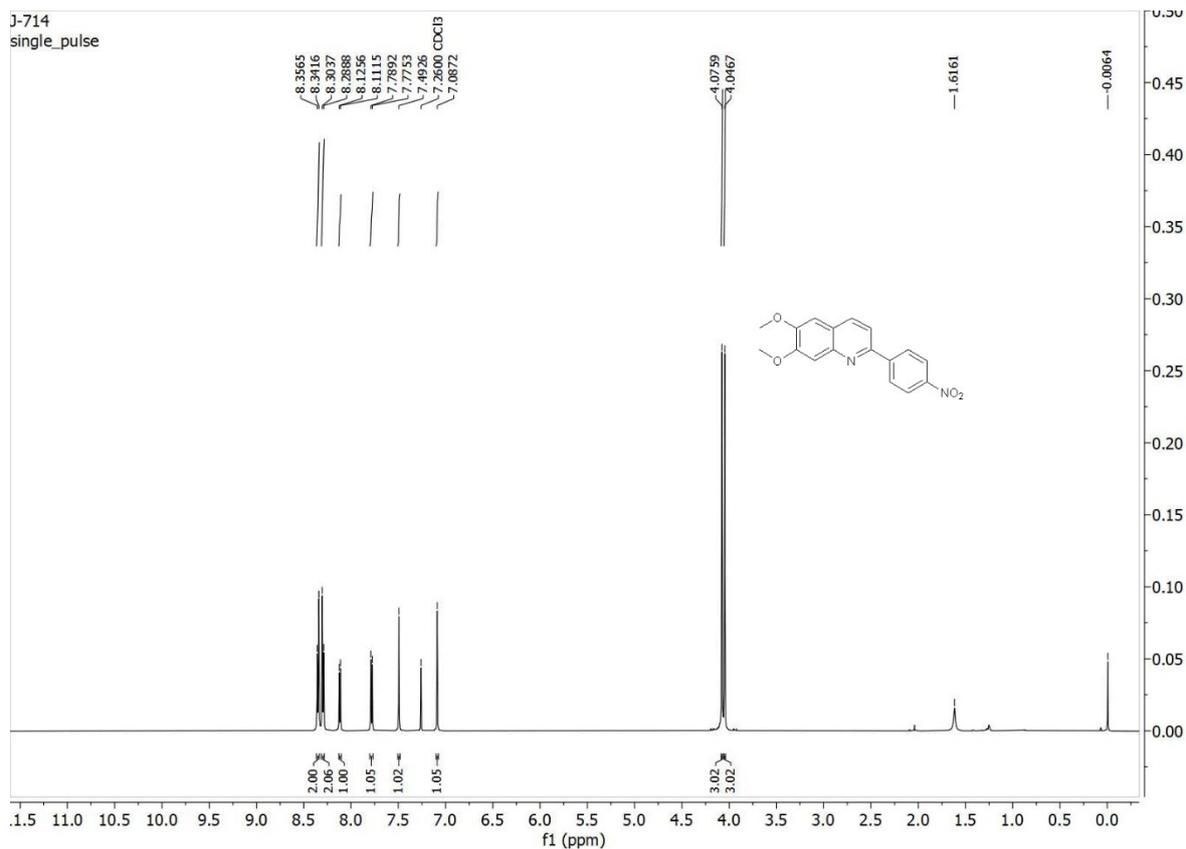


Figure S13. ^1H NMR Spectrum of **21a** in CDCl_3 .

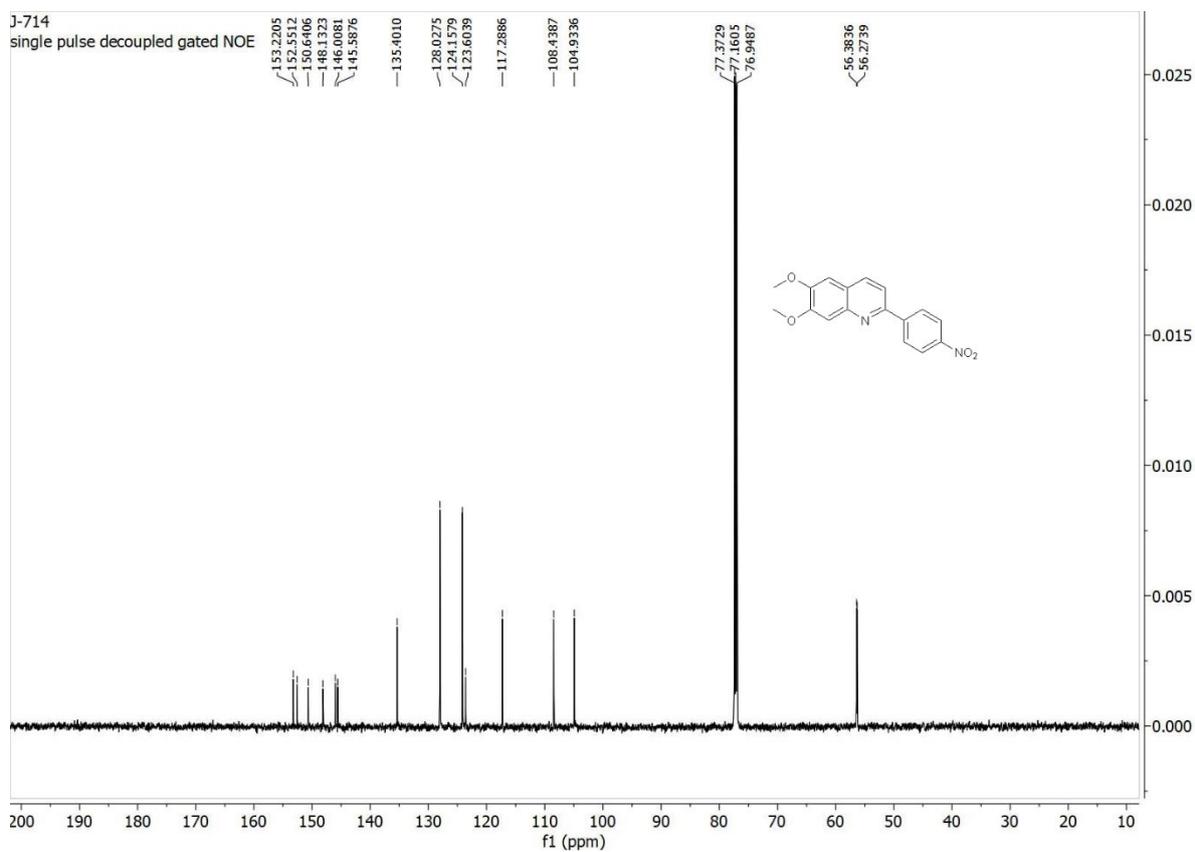


Figure S14. ^{13}C NMR Spectrum of **21a** in CDCl_3 .

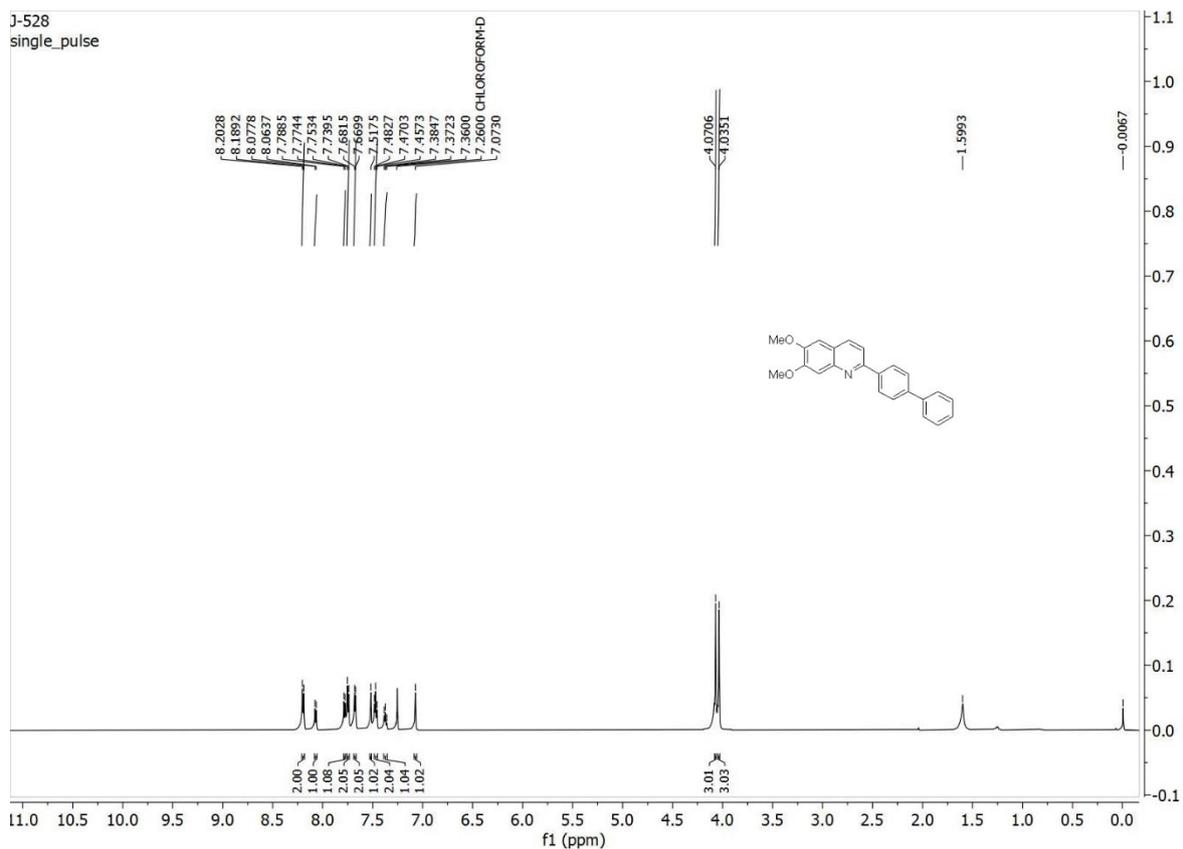


Figure S15. ^1H NMR Spectrum of **22a** in CDCl_3 .

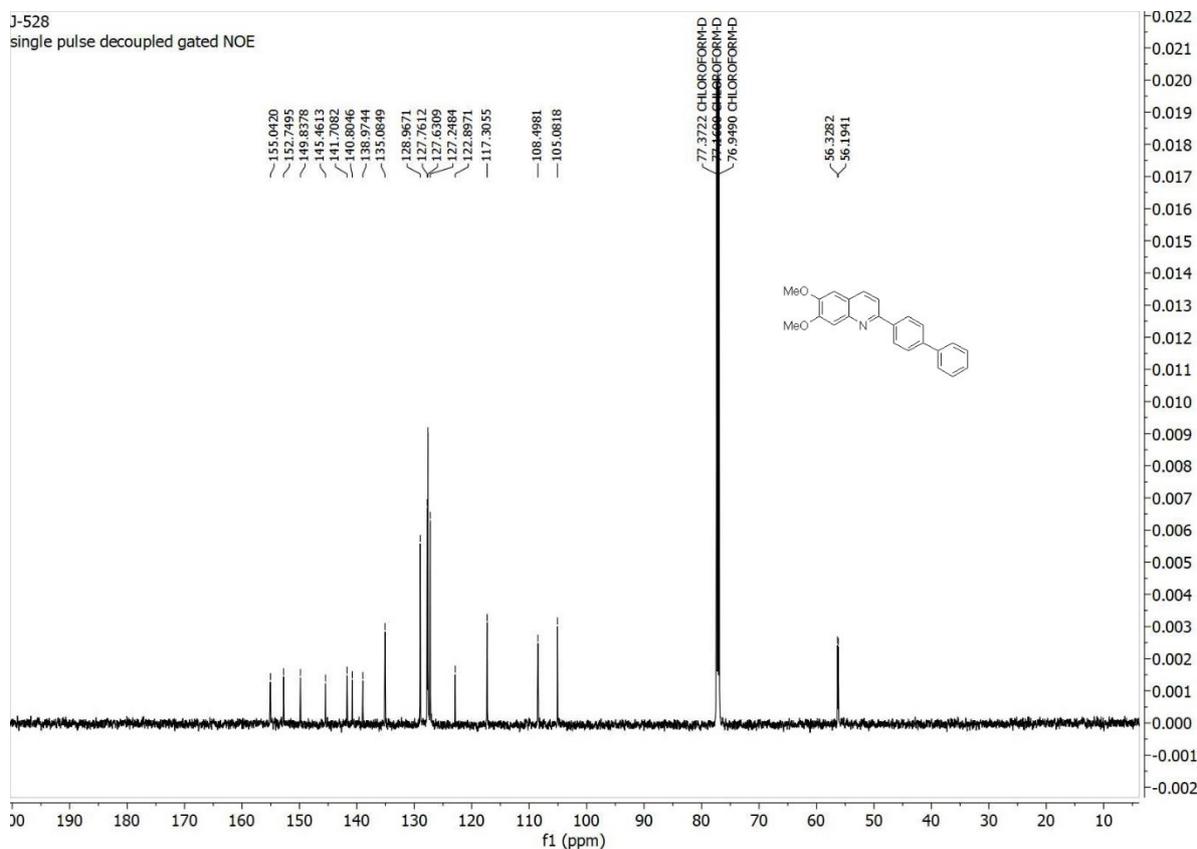


Figure S16. ^{13}C NMR Spectrum of **22a** in CDCl_3 .

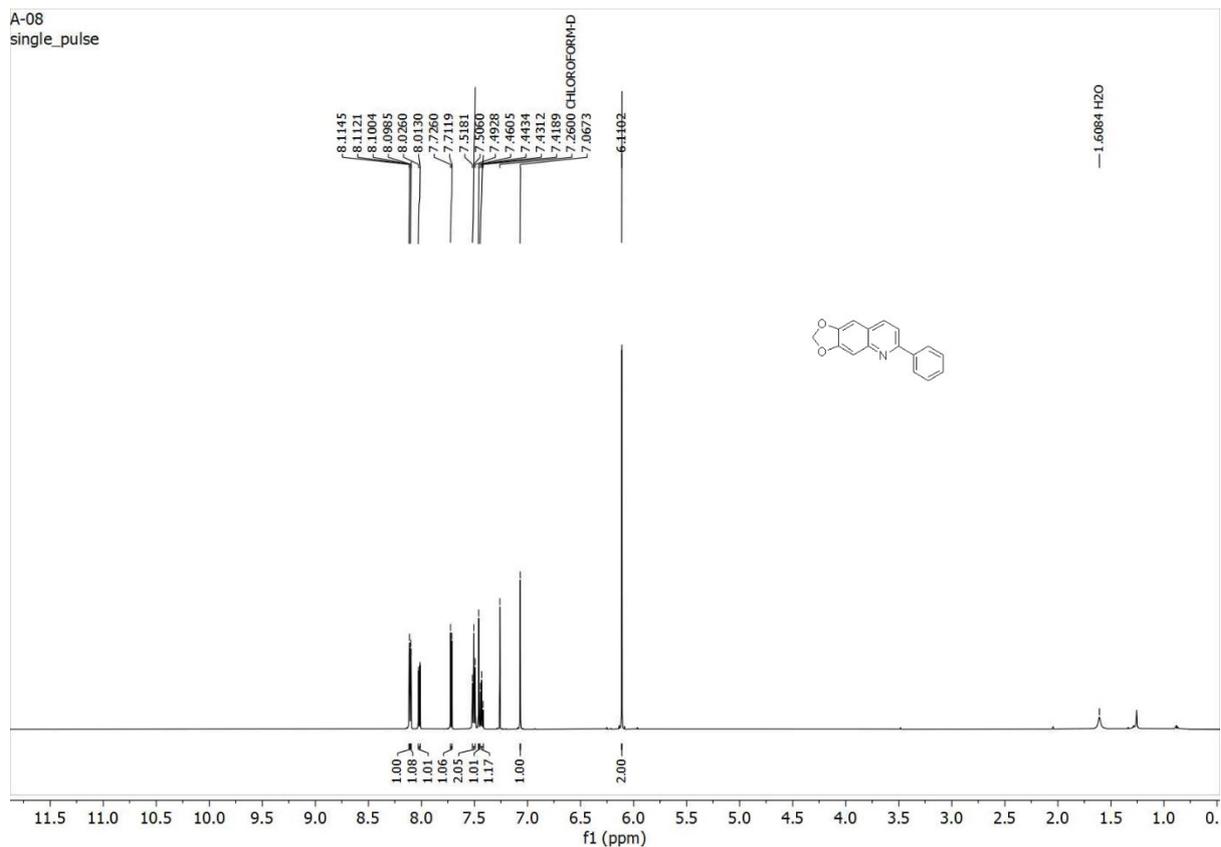


Figure S17. ^1H NMR Spectrum of **15b** in CDCl_3 .

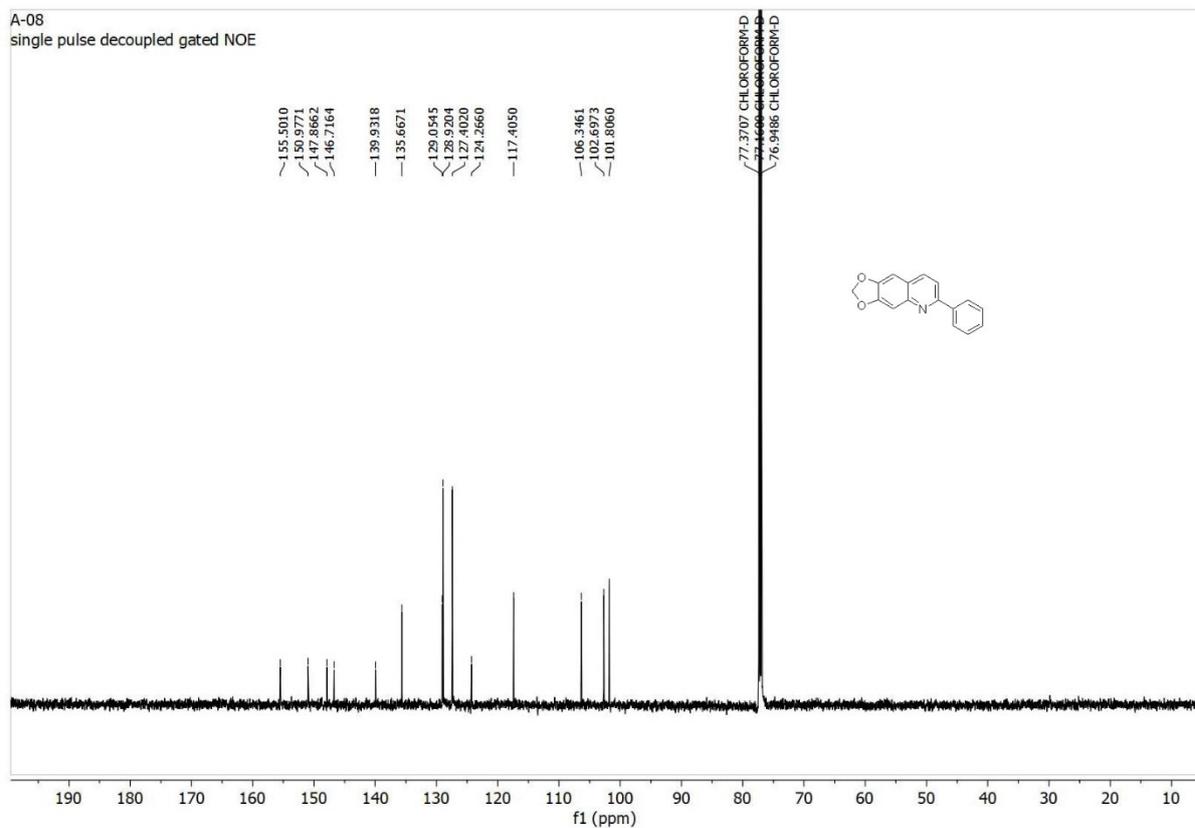


Figure S18. ^{13}C NMR Spectrum of **15b** in CDCl_3 .

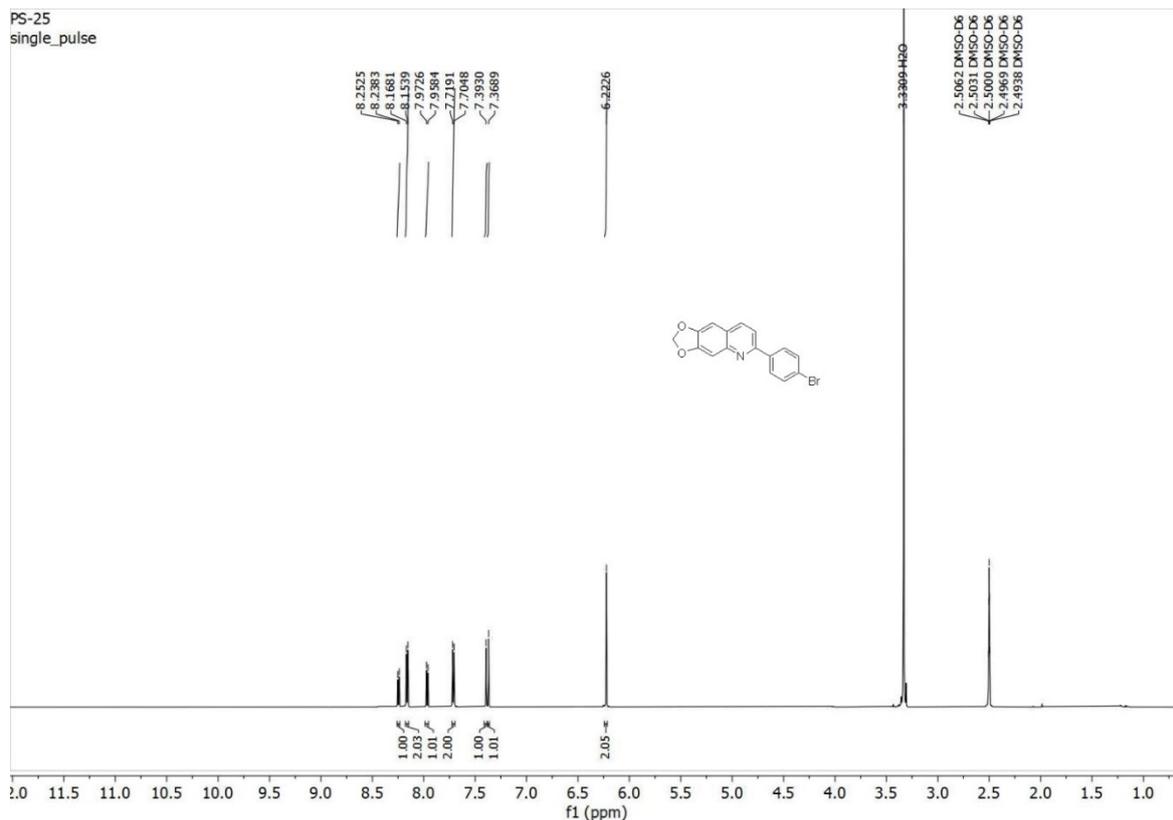


Figure S19. ^1H NMR Spectrum of **16b** in CDCl_3 .

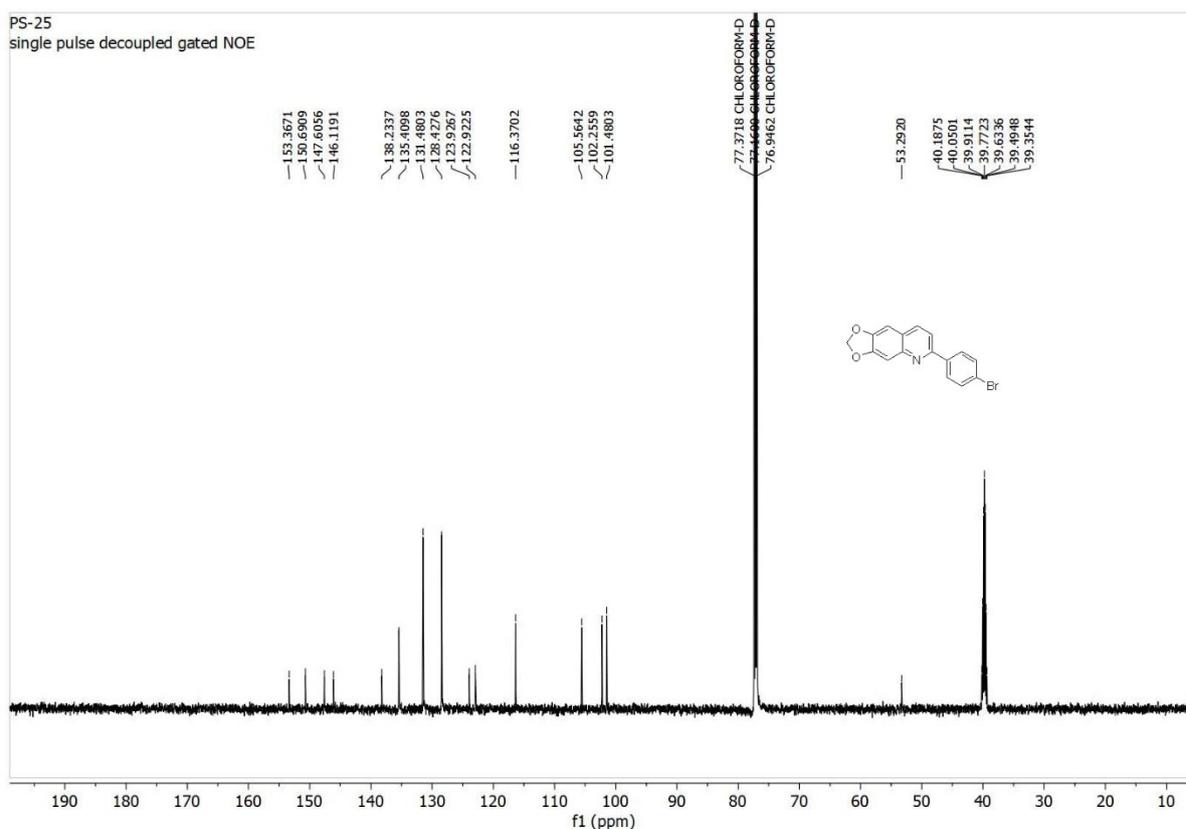


Figure S20. ^{13}C NMR Spectrum of **16b** in CDCl_3 .

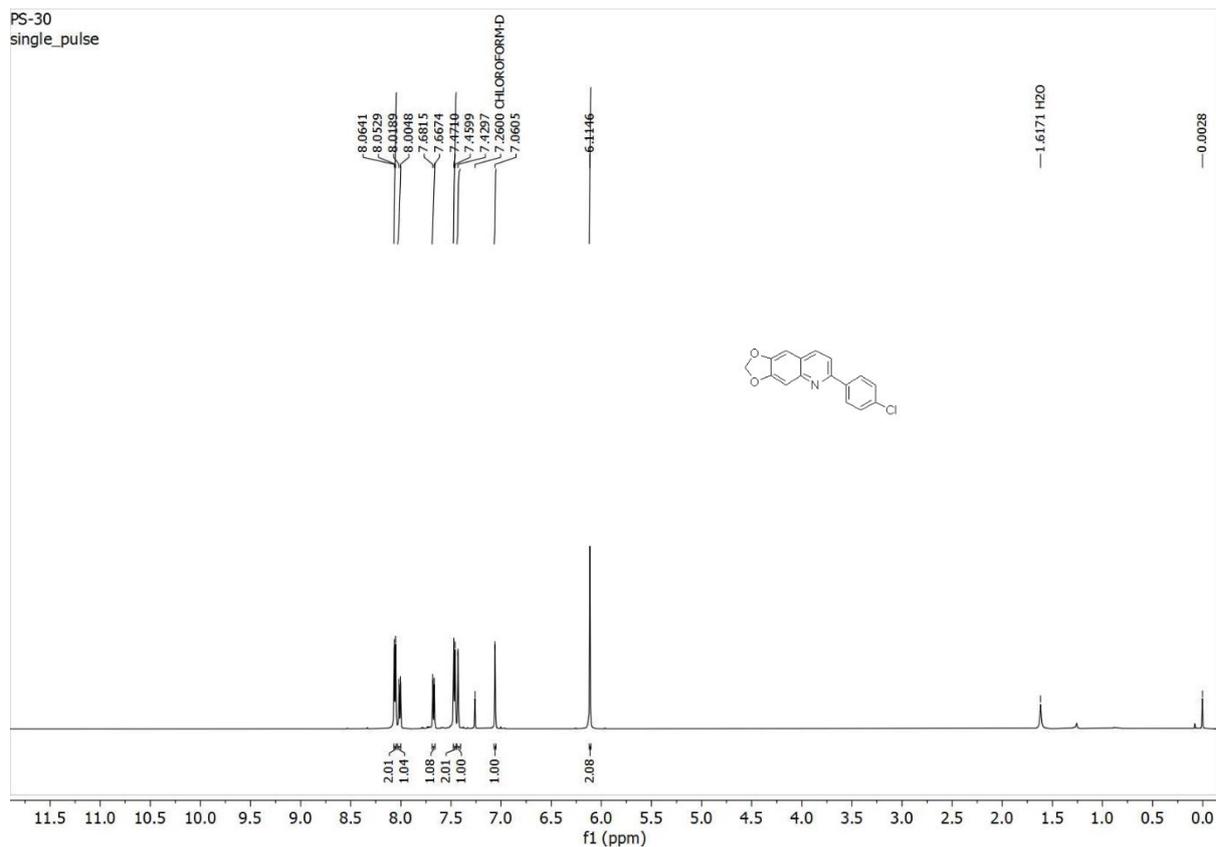


Figure S21. ^1H NMR Spectrum of **17b** in CDCl_3 .

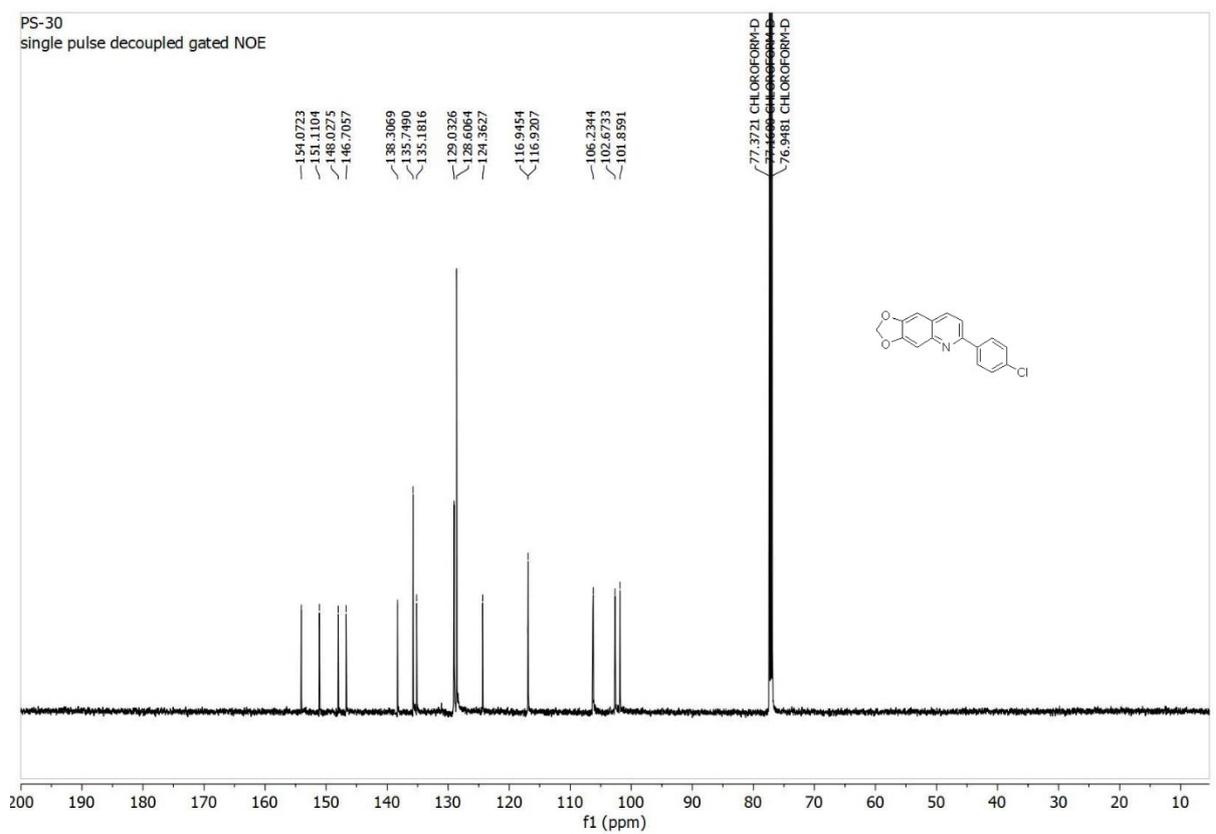


Figure S22. ^{13}C NMR Spectrum of **17b** in CDCl_3 .

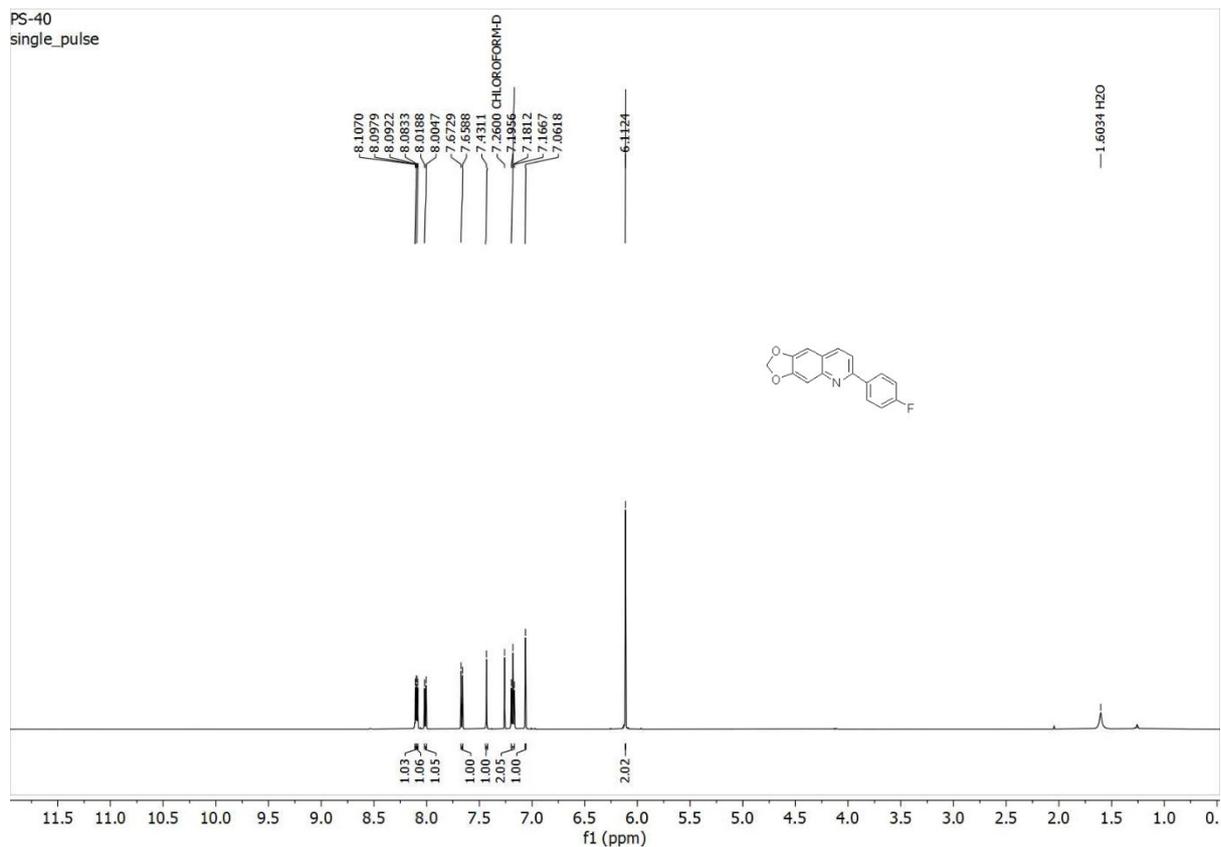


Figure S23. ^1H NMR Spectrum of **18b** in CDCl_3 .

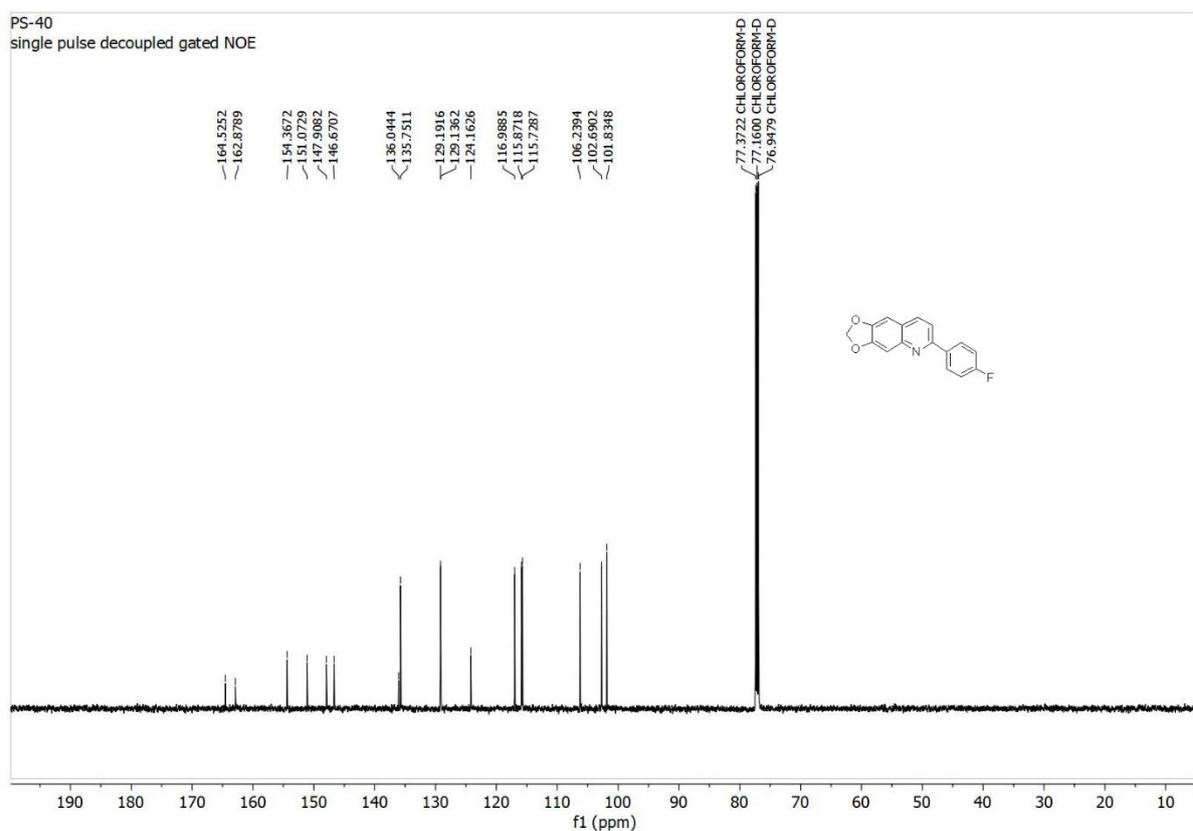


Figure S24. ^{13}C NMR Spectrum of **18b** in CDCl_3 .

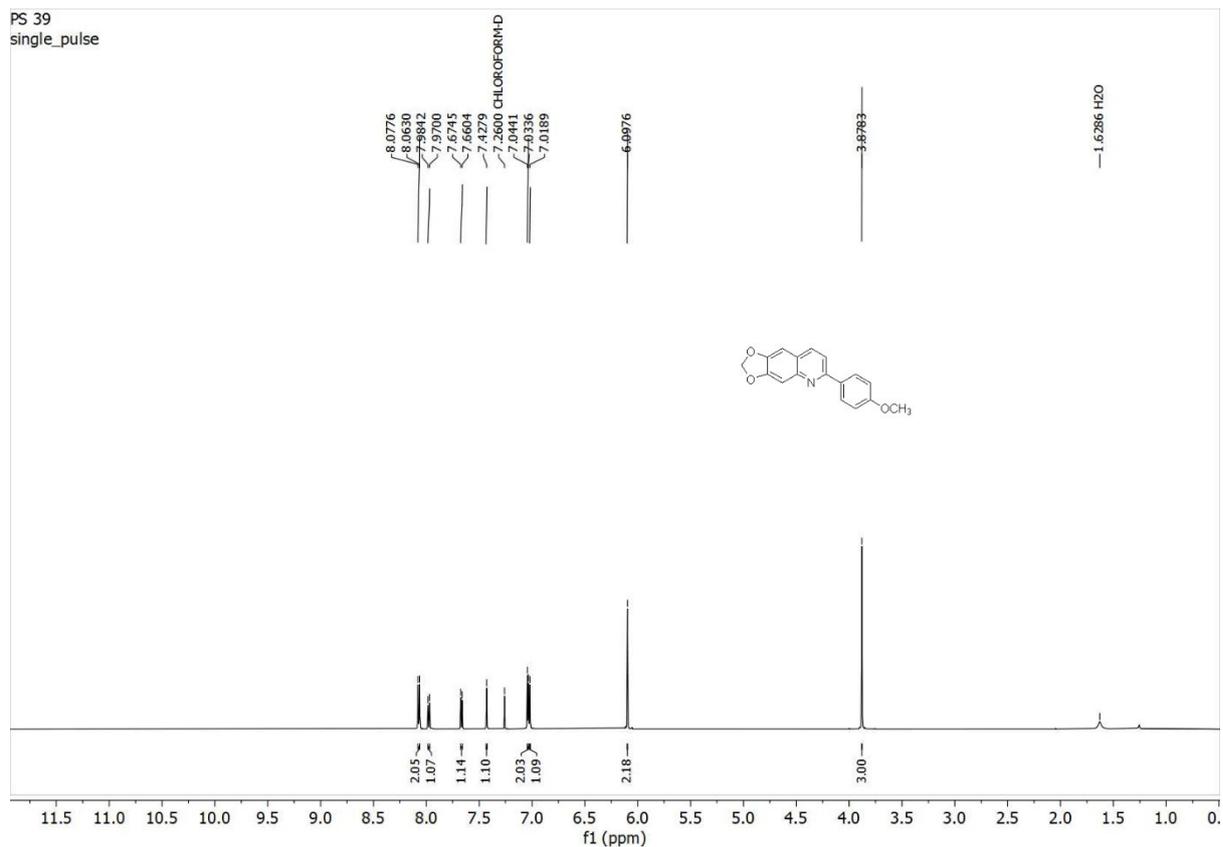


Figure S25. ^1H NMR Spectrum of **19b** in CDCl_3 .

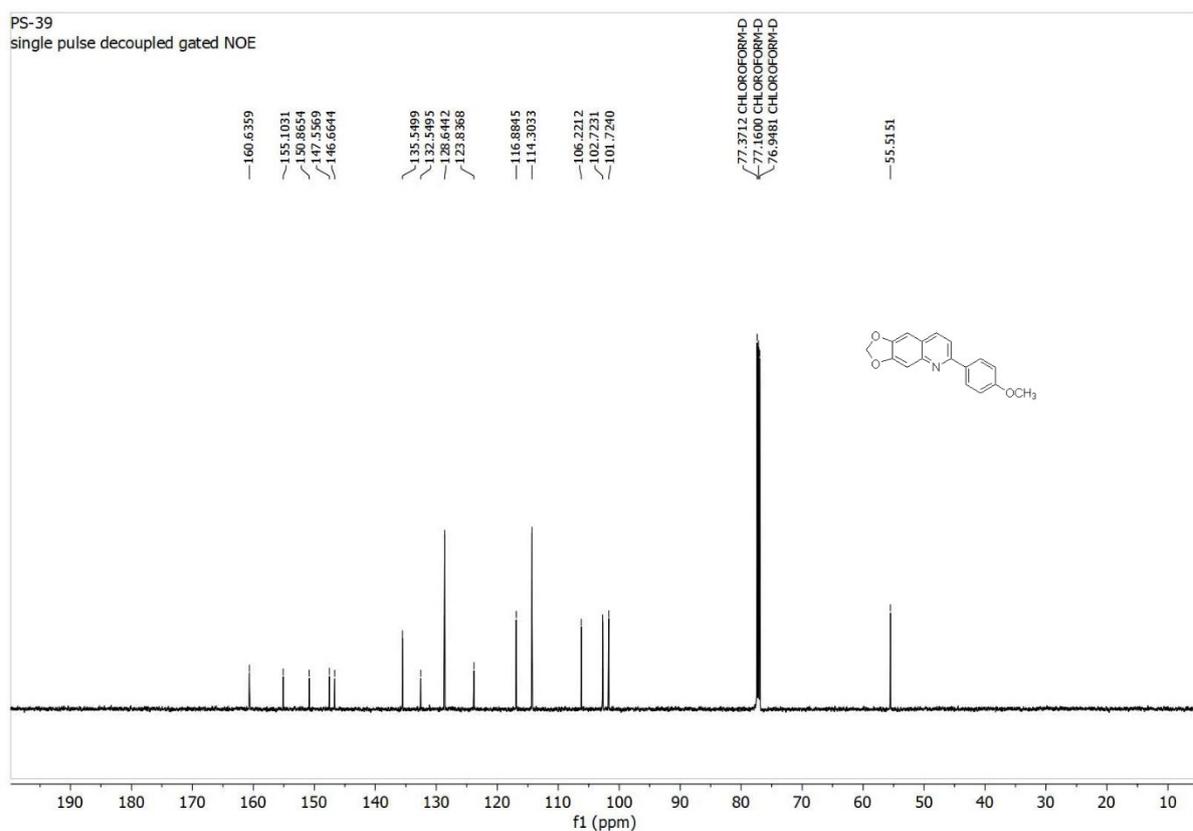


Figure S26. ^{13}C NMR Spectrum of **19b** in CDCl_3 .

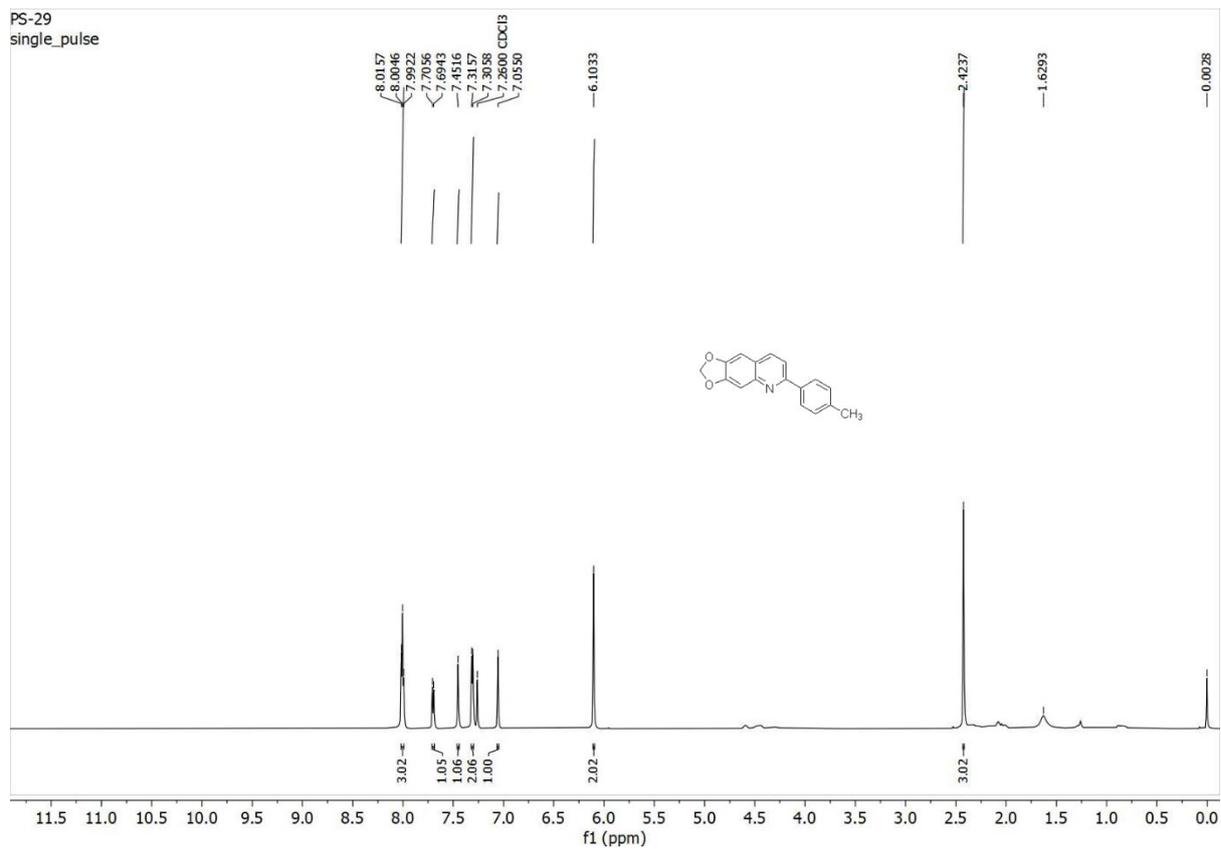


Figure S27. ¹H NMR Spectrum of **20b** in CDCl₃.

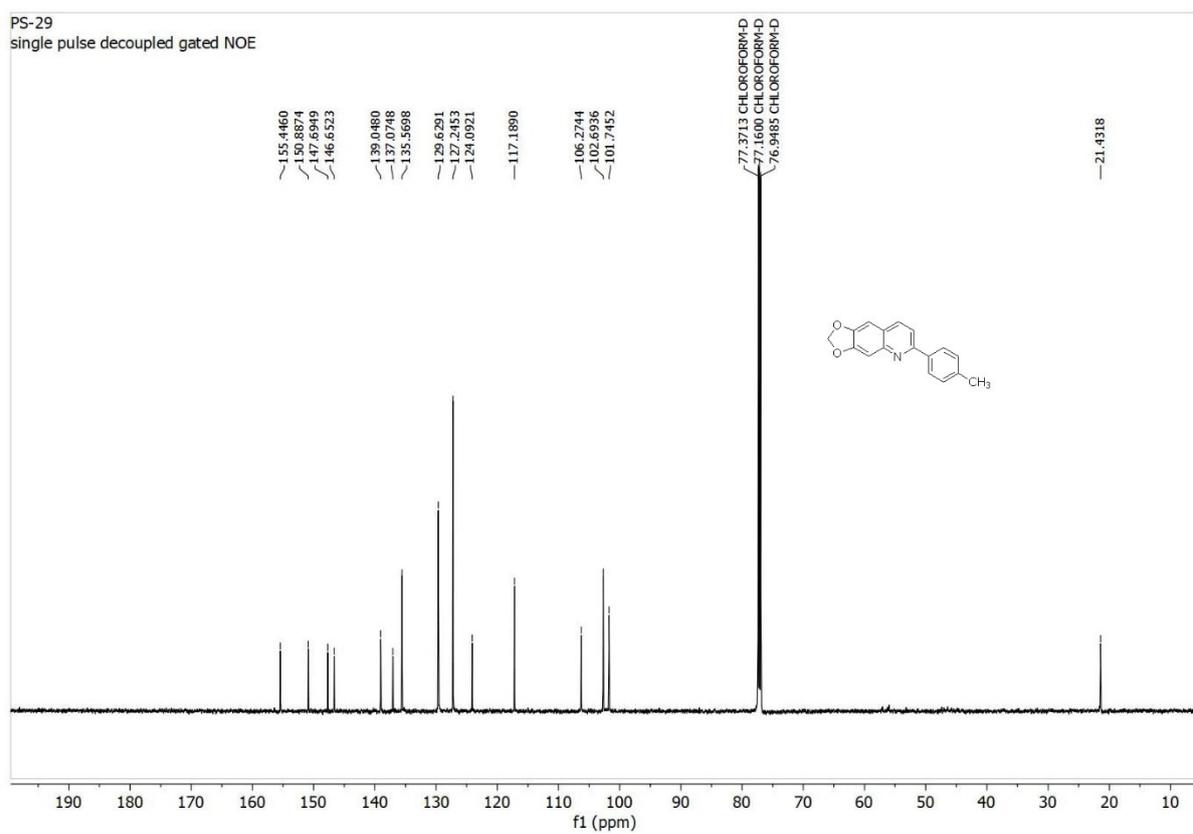


Figure S28. ¹³C NMR Spectrum of **20b** in CDCl₃.

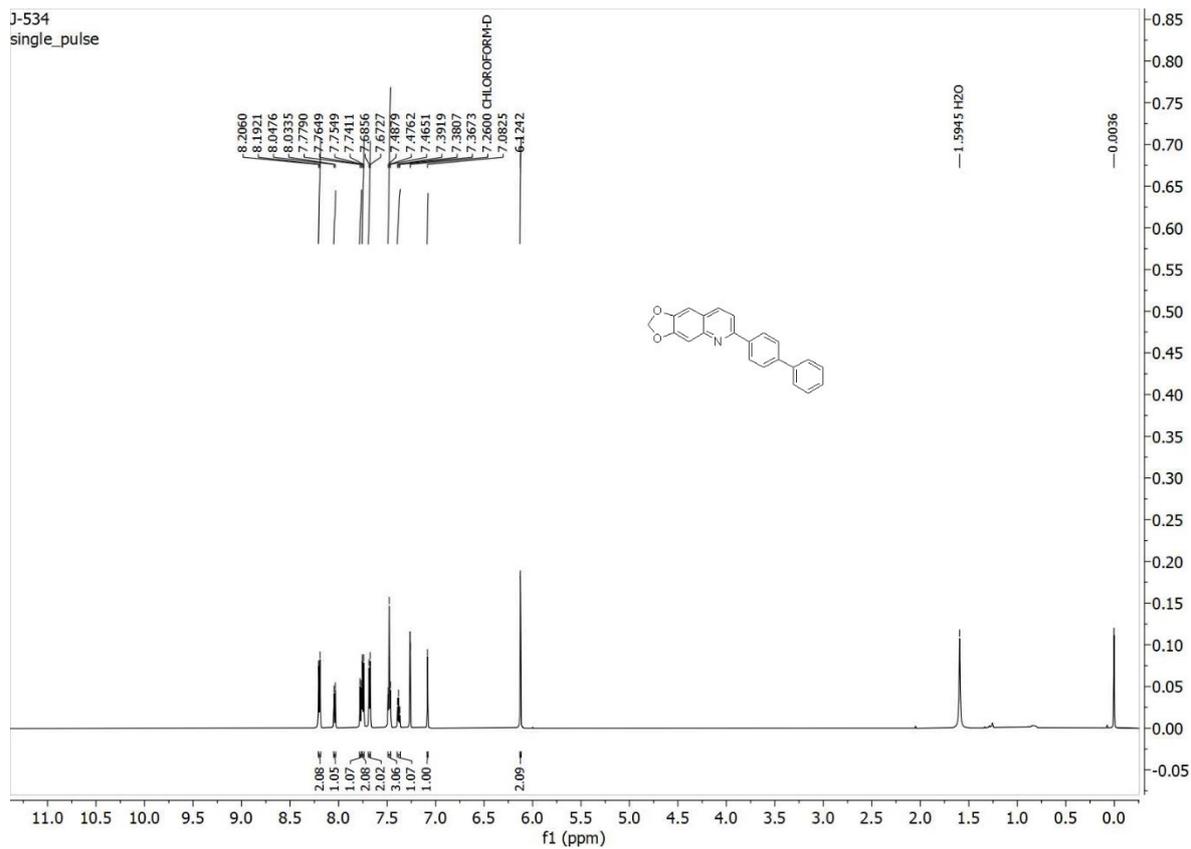


Figure S29. ^1H NMR Spectrum of **22b** in CDCl_3 .

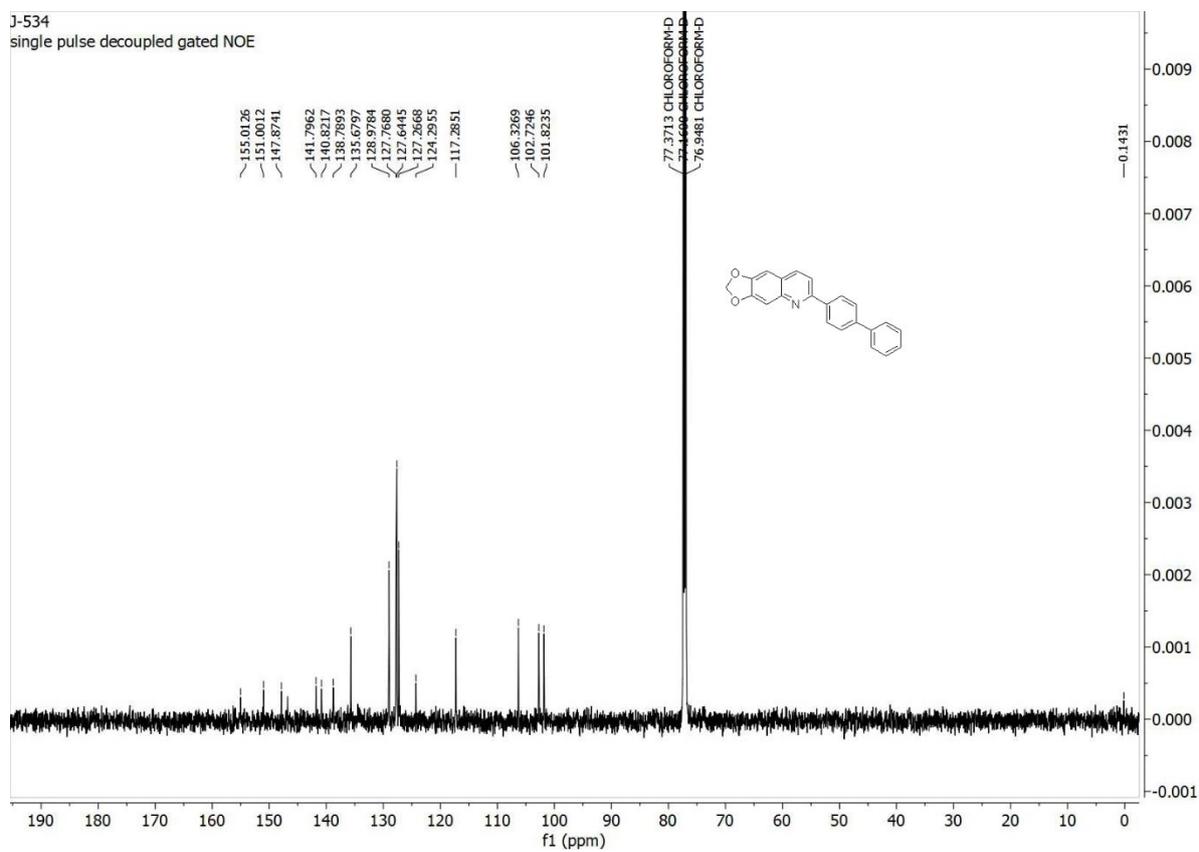


Figure S30. ^{13}C NMR Spectrum of **22b** in CDCl_3 .

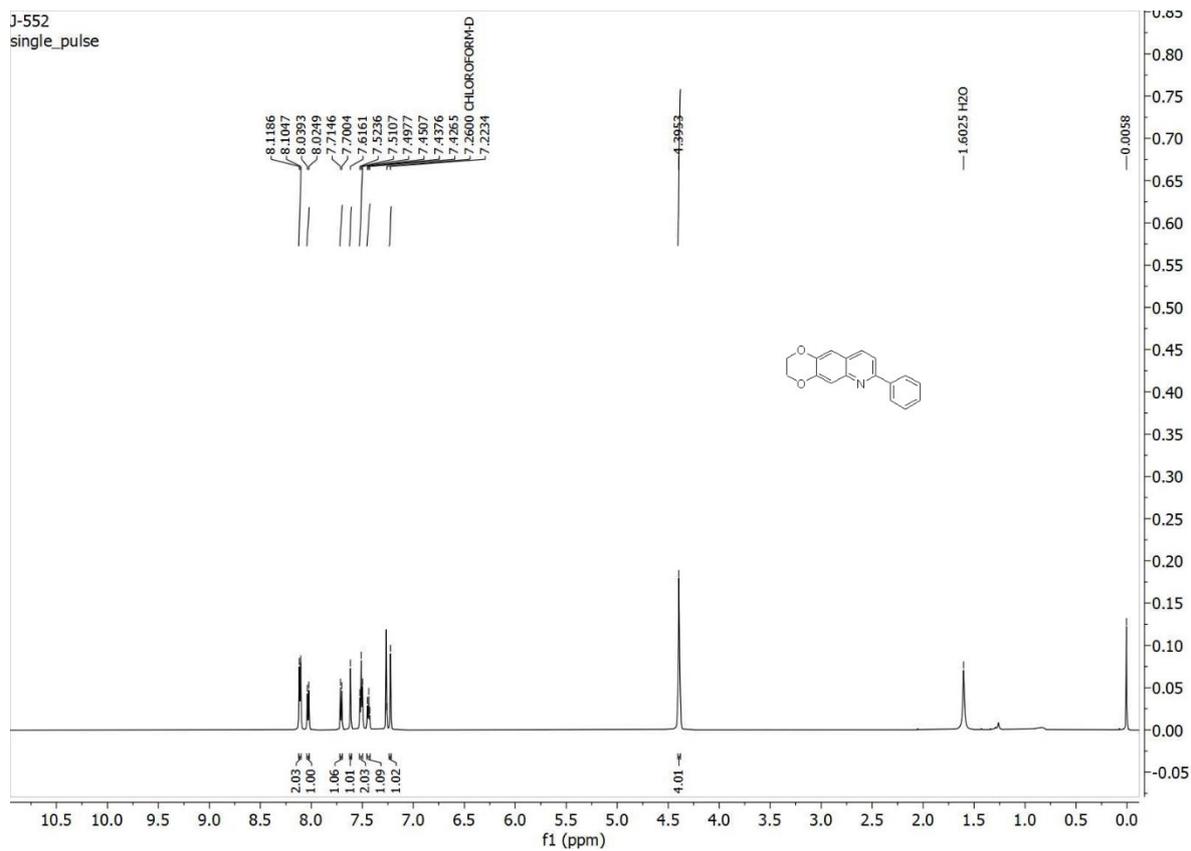


Figure S31. ^1H NMR Spectrum of **15c** in CDCl_3 .

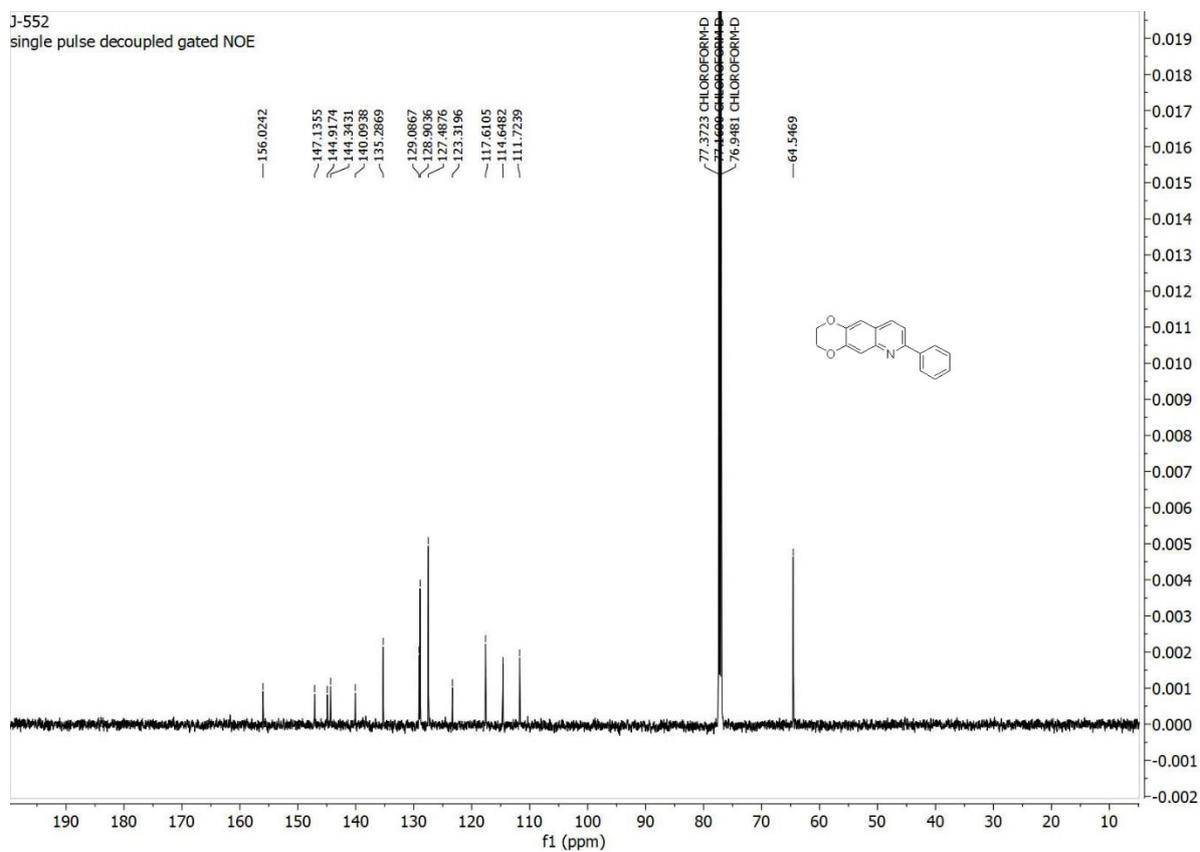


Figure S32. ^{13}C NMR Spectrum of **15c** in CDCl_3 .

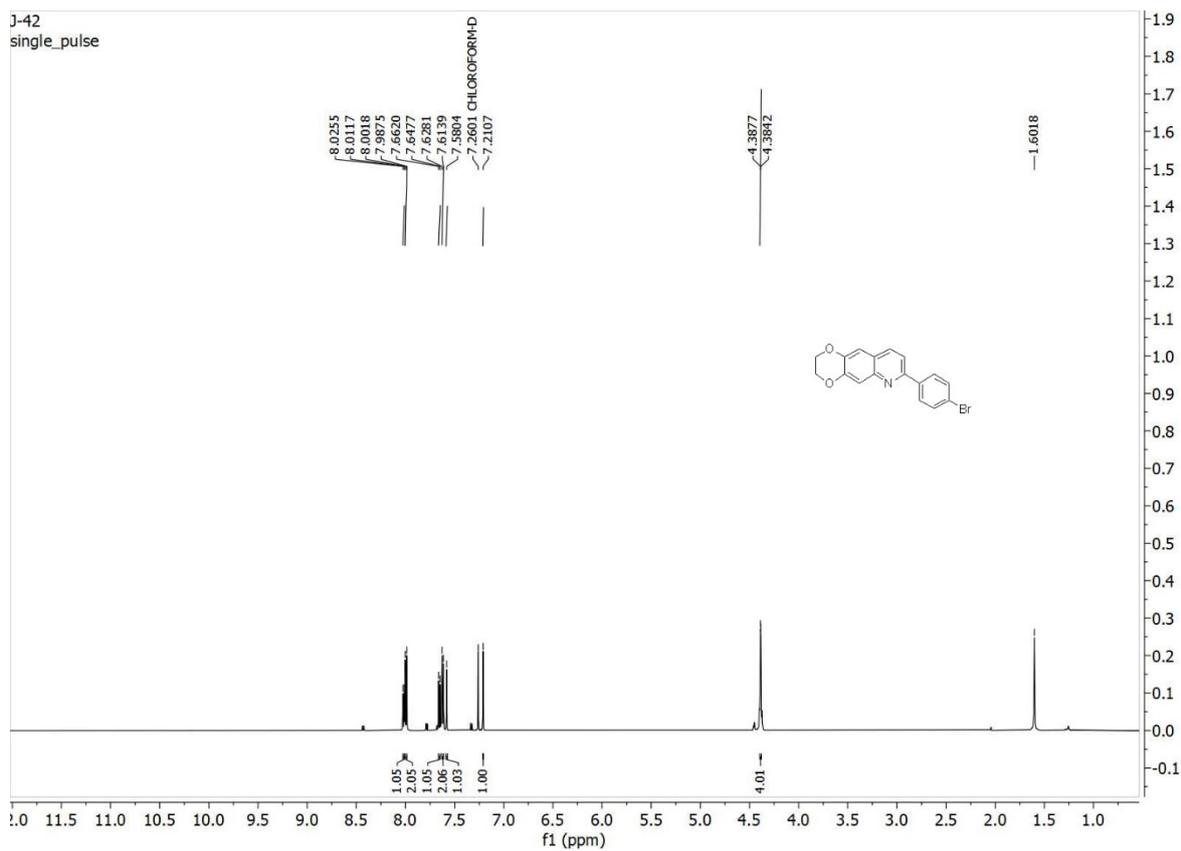


Figure S33. ^1H NMR Spectrum of **16c** in CDCl_3 .

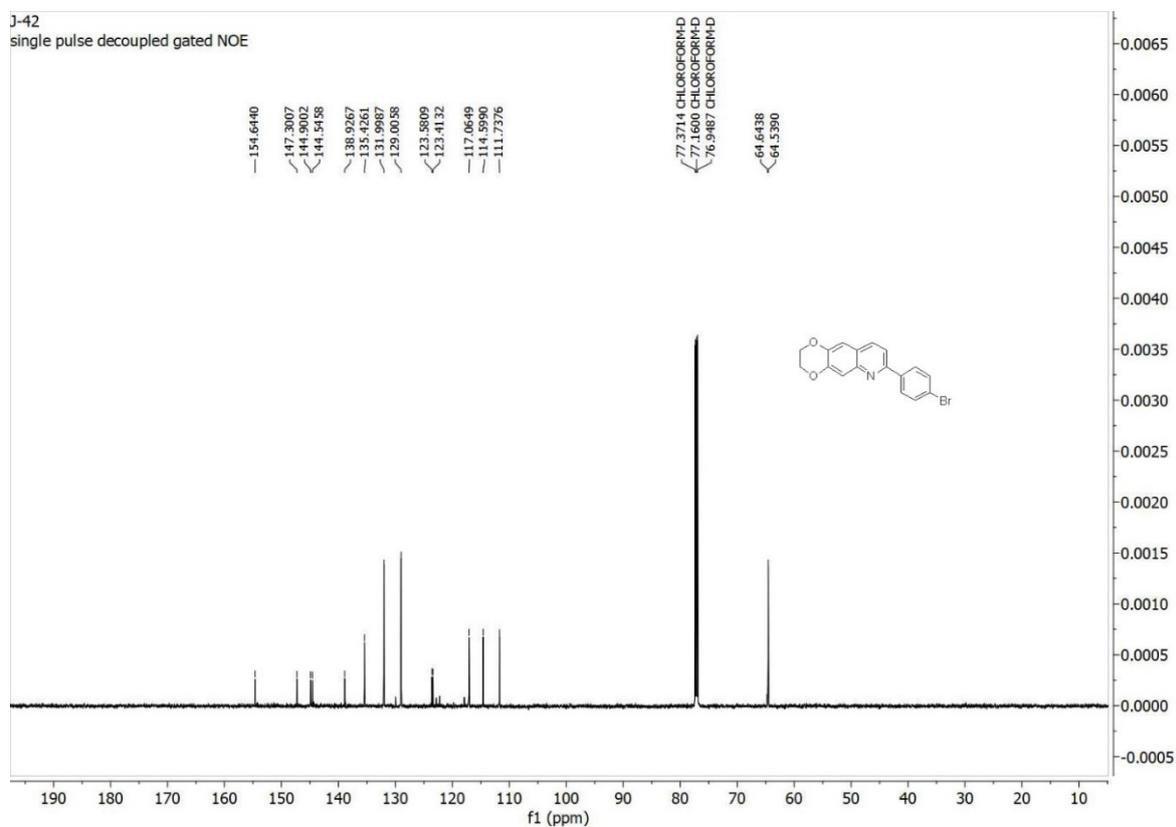


Figure S34. ^{13}C NMR Spectrum of **16c** in CDCl_3 .

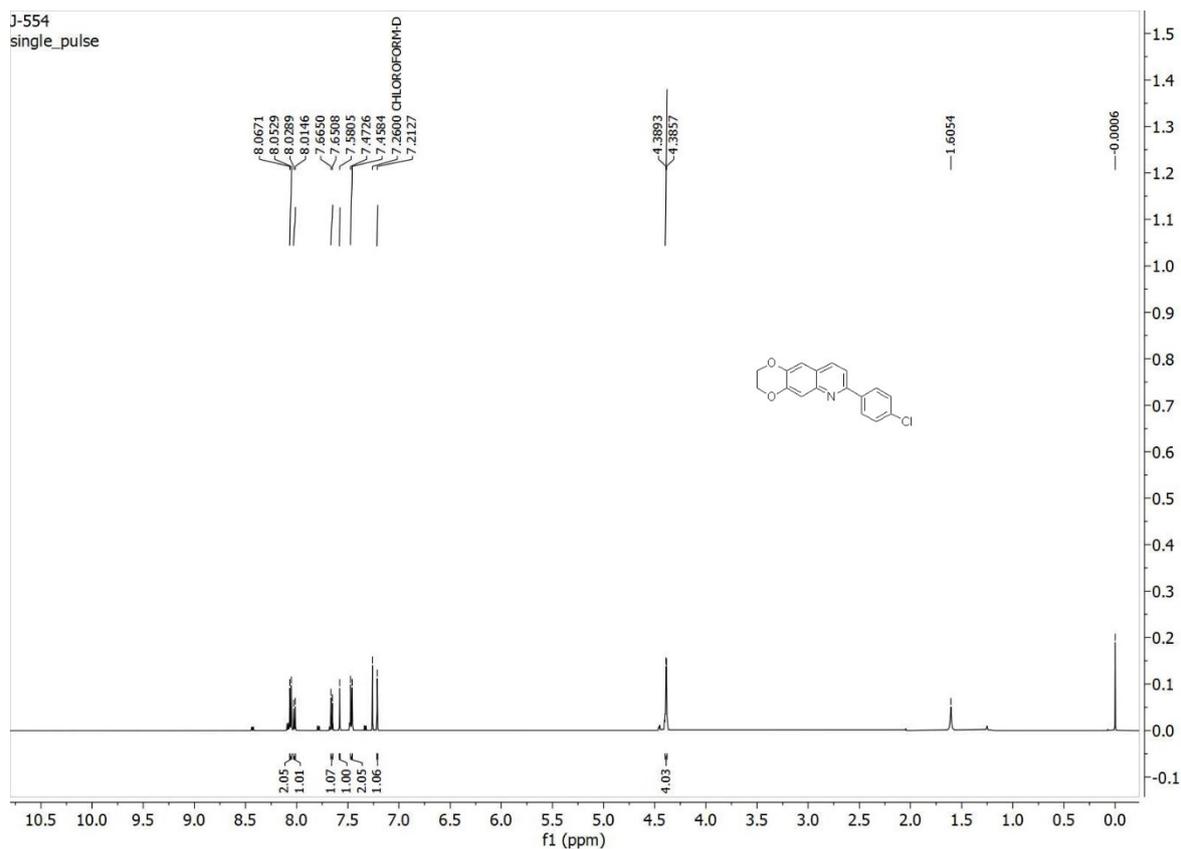


Figure S35. ^1H NMR Spectrum of **17c** in CDCl_3 .

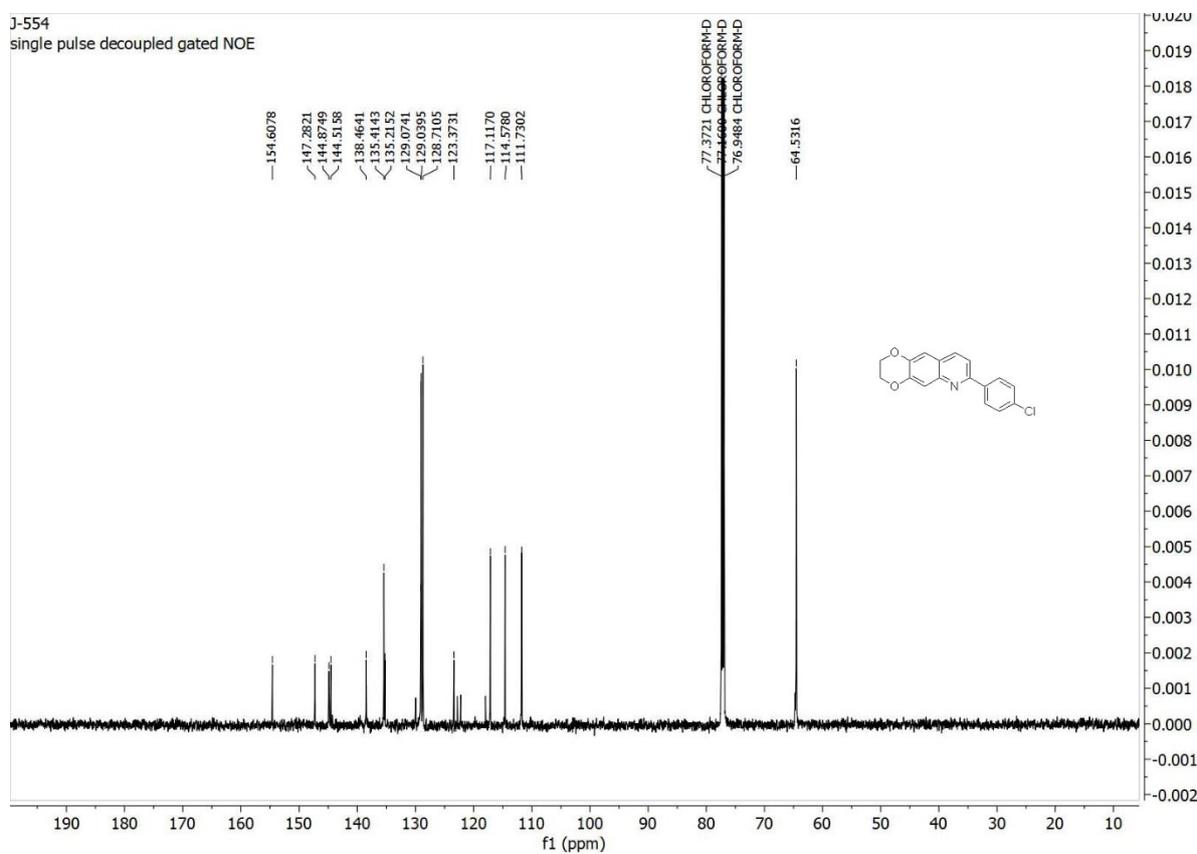


Figure S36. ^{13}C NMR Spectrum of **17c** in CDCl_3 .

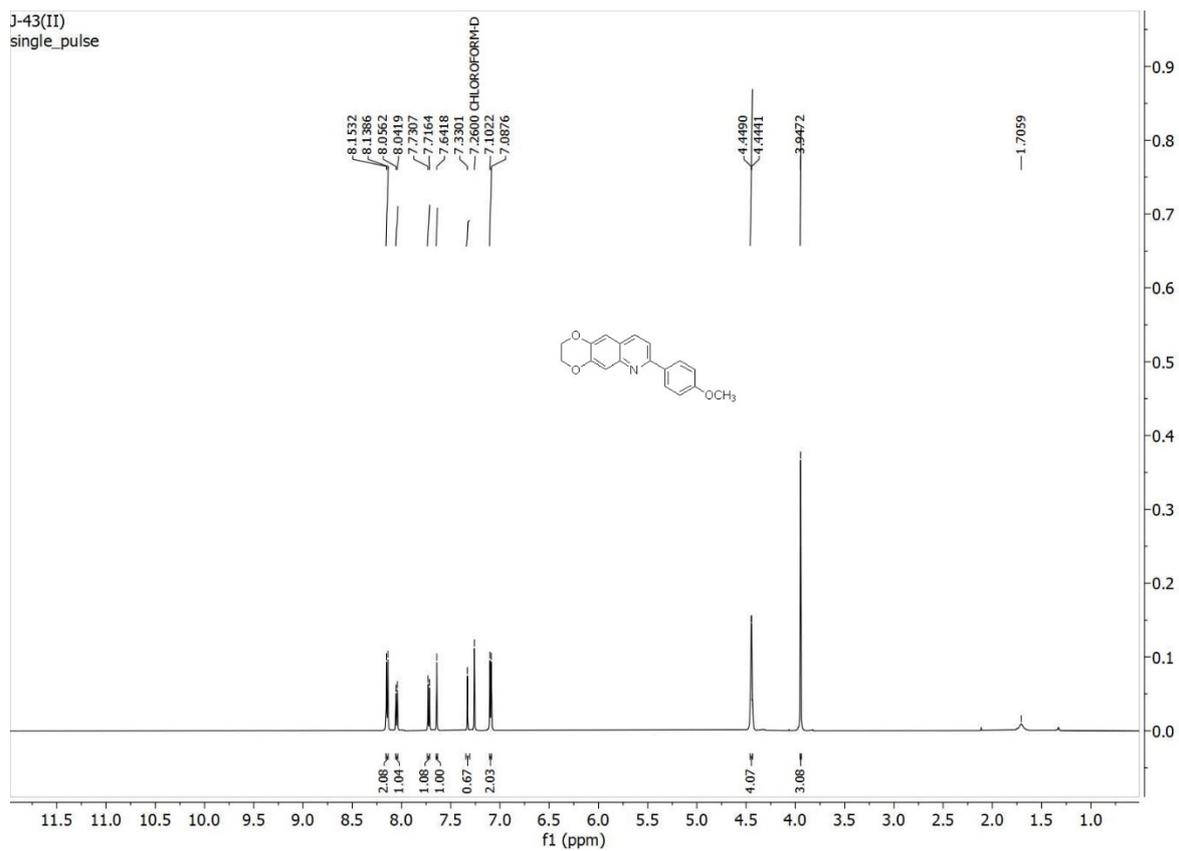


Figure S37. ^1H NMR Spectrum of **19c** in CDCl_3 .

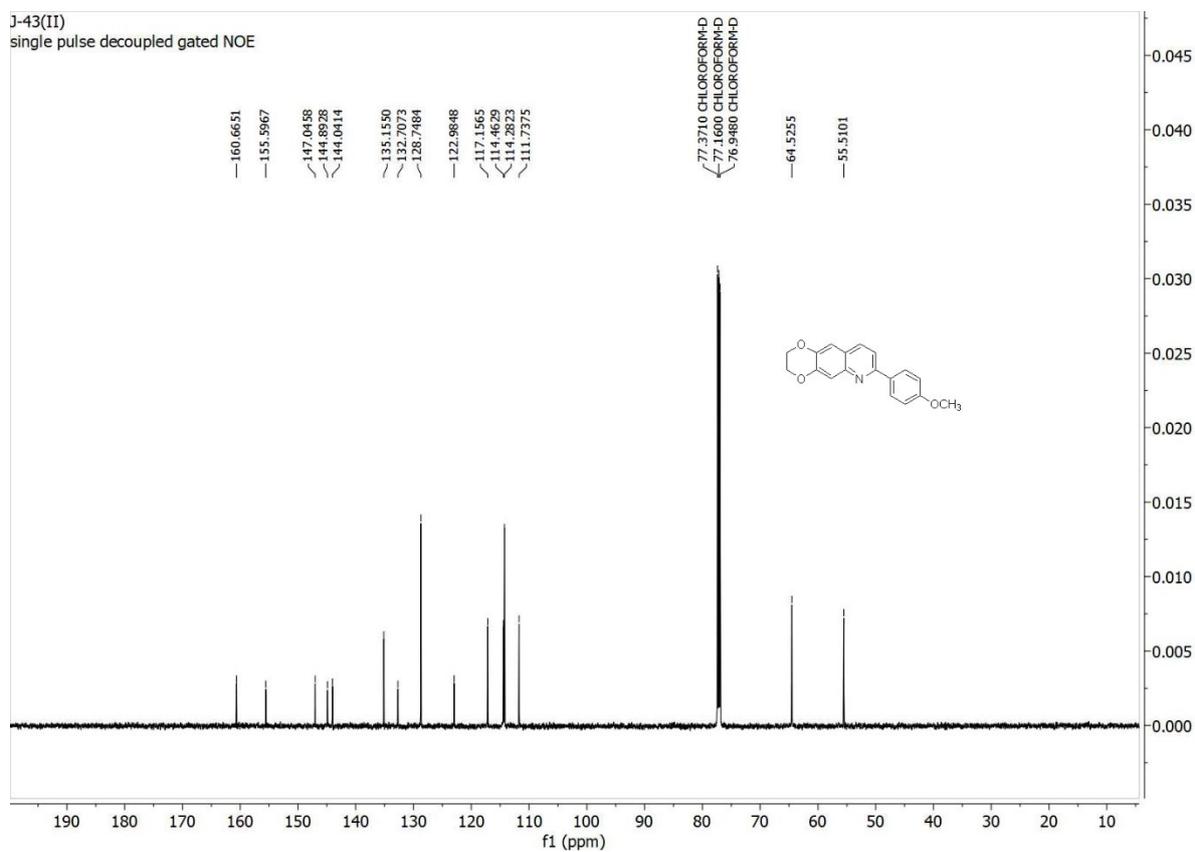


Figure S38. ^{13}C NMR Spectrum of **19c** in CDCl_3 .

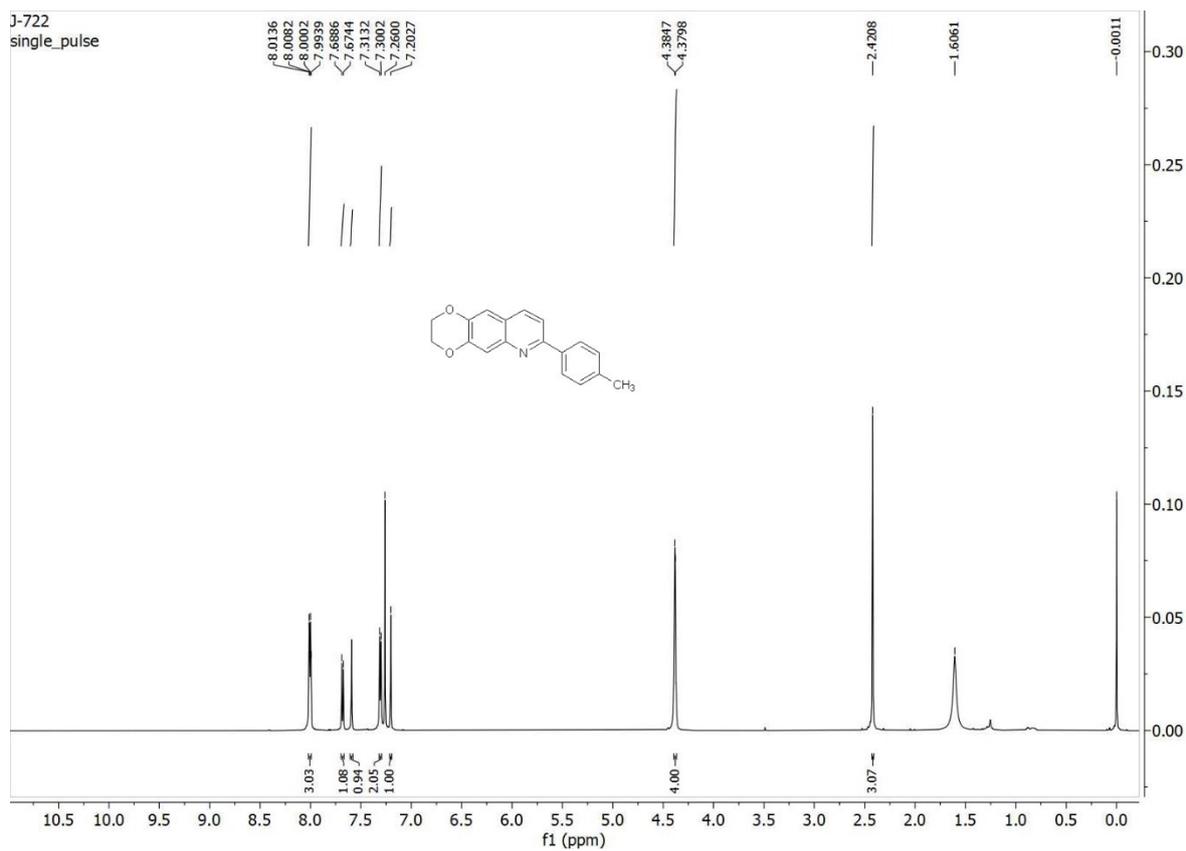


Figure S39. ^1H NMR Spectrum of **20c** in CDCl_3 .

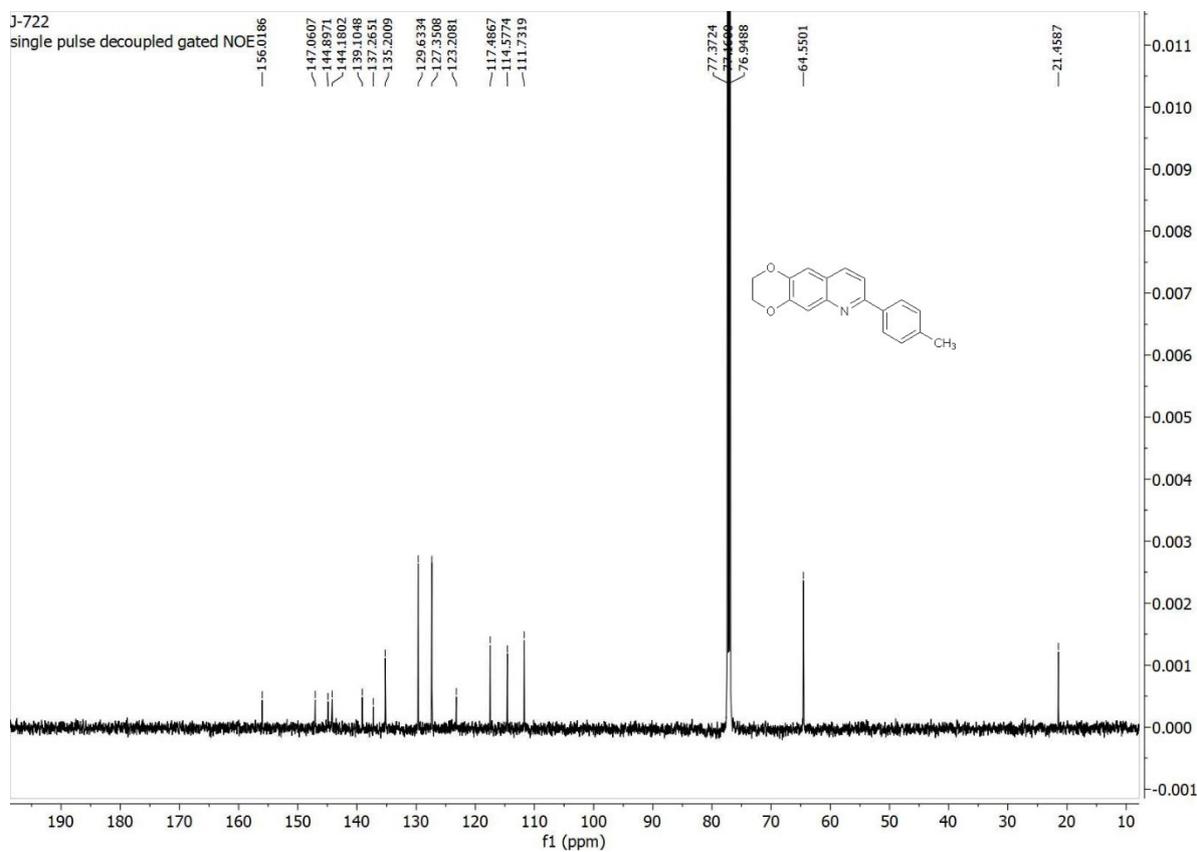


Figure S40. ^{13}C NMR Spectrum of **20c** in CDCl_3 .

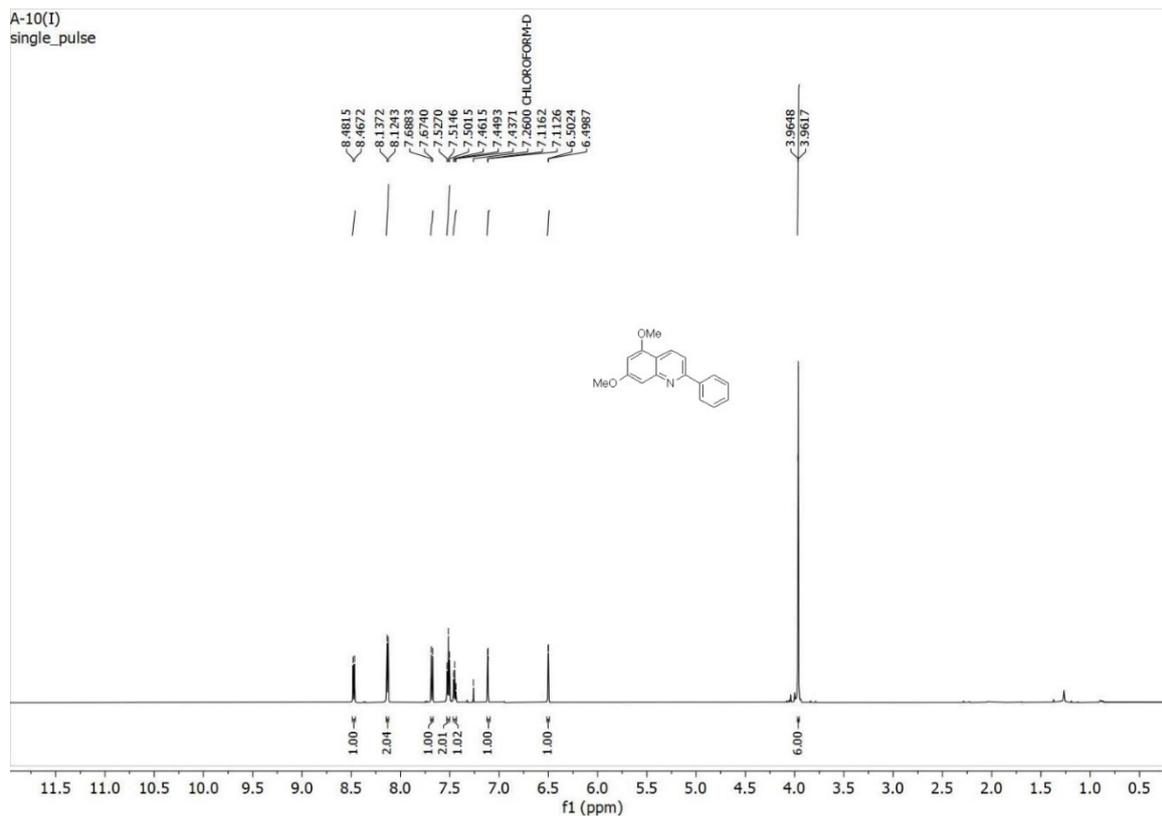


Figure S41. ^1H NMR Spectrum of **15d** in CDCl_3 .

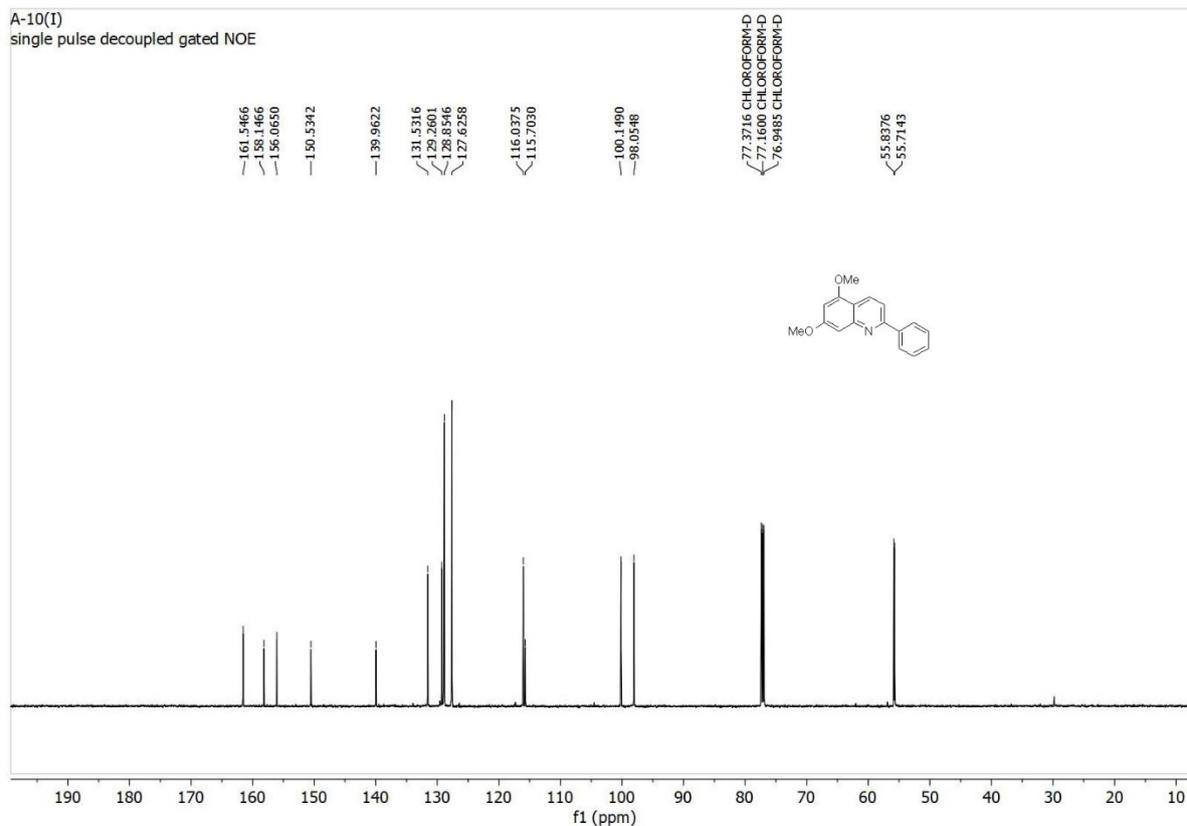


Figure S42. ^{13}C NMR Spectrum of **15d** in CDCl_3 .

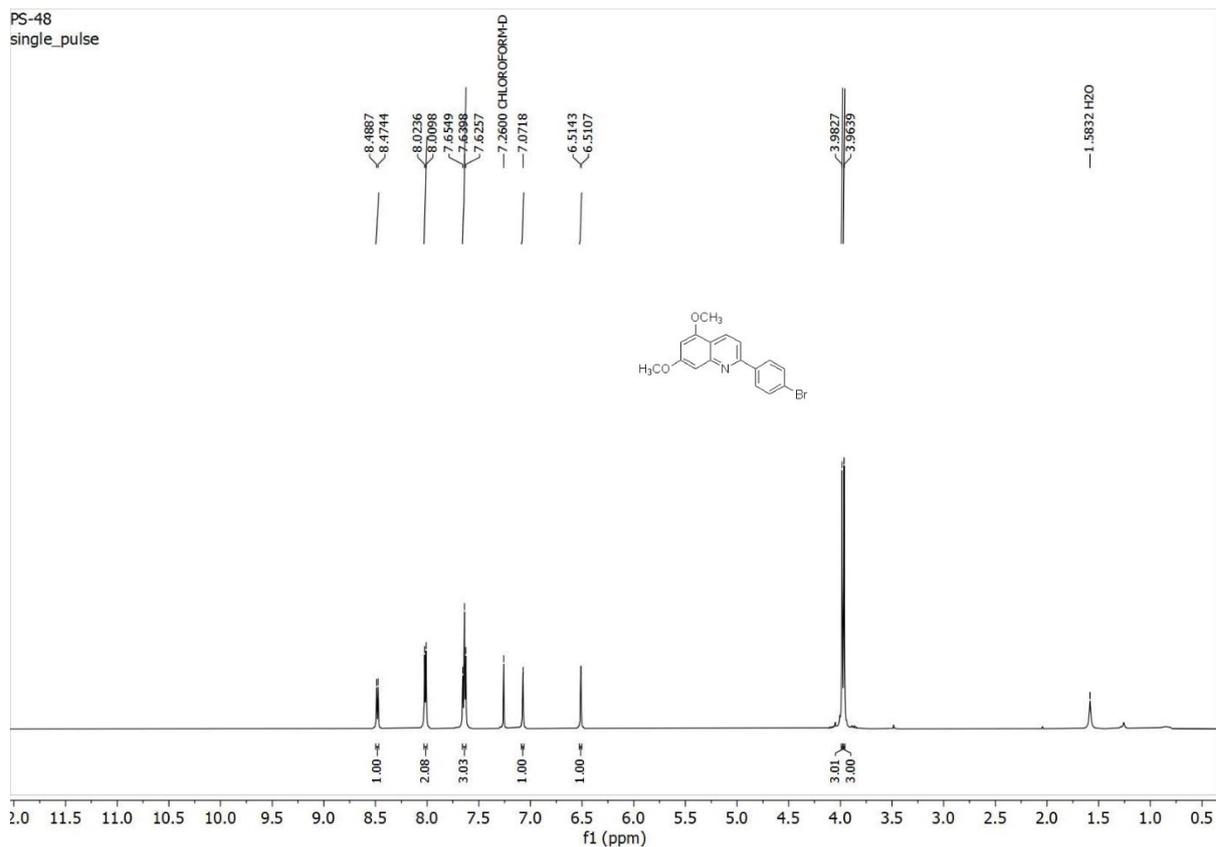


Figure S43. ^1H NMR Spectrum of **16d** in CDCl_3 .

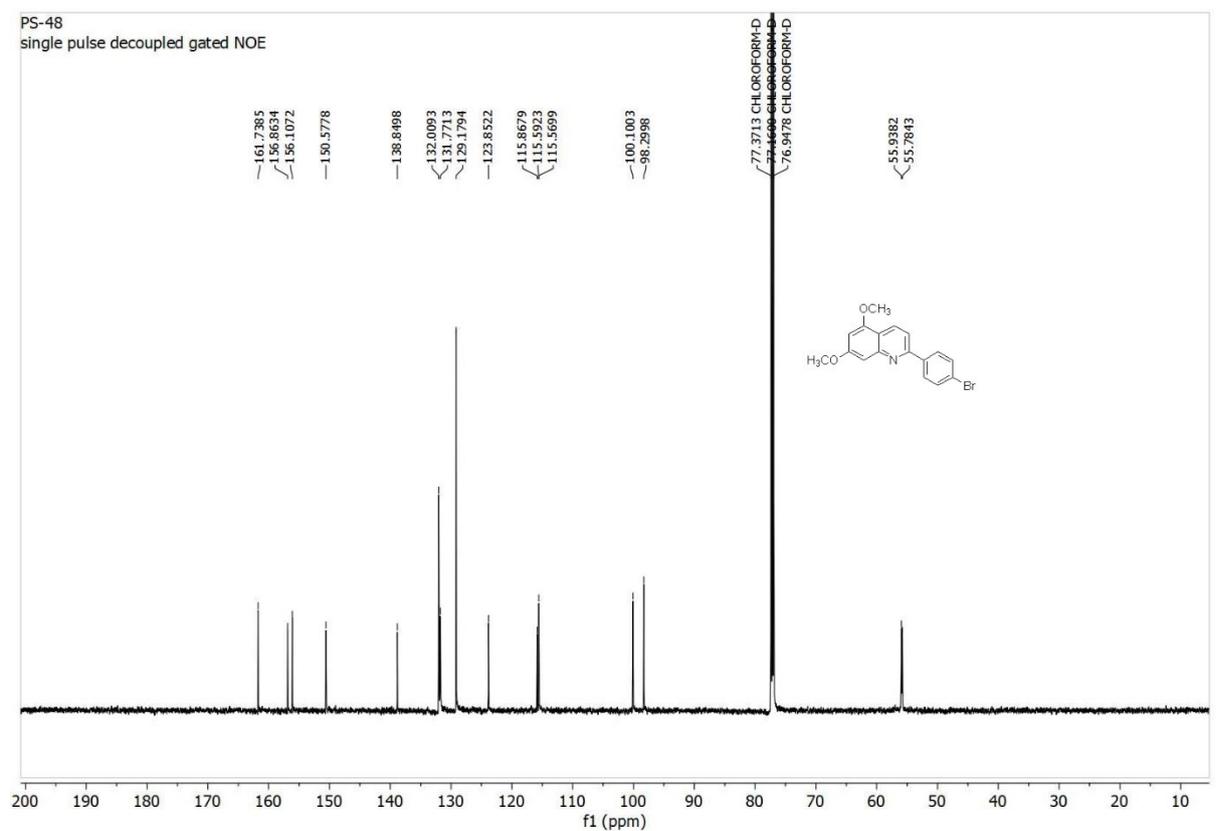


Figure S44. ^{13}C NMR Spectrum of **16d** in CDCl_3 .

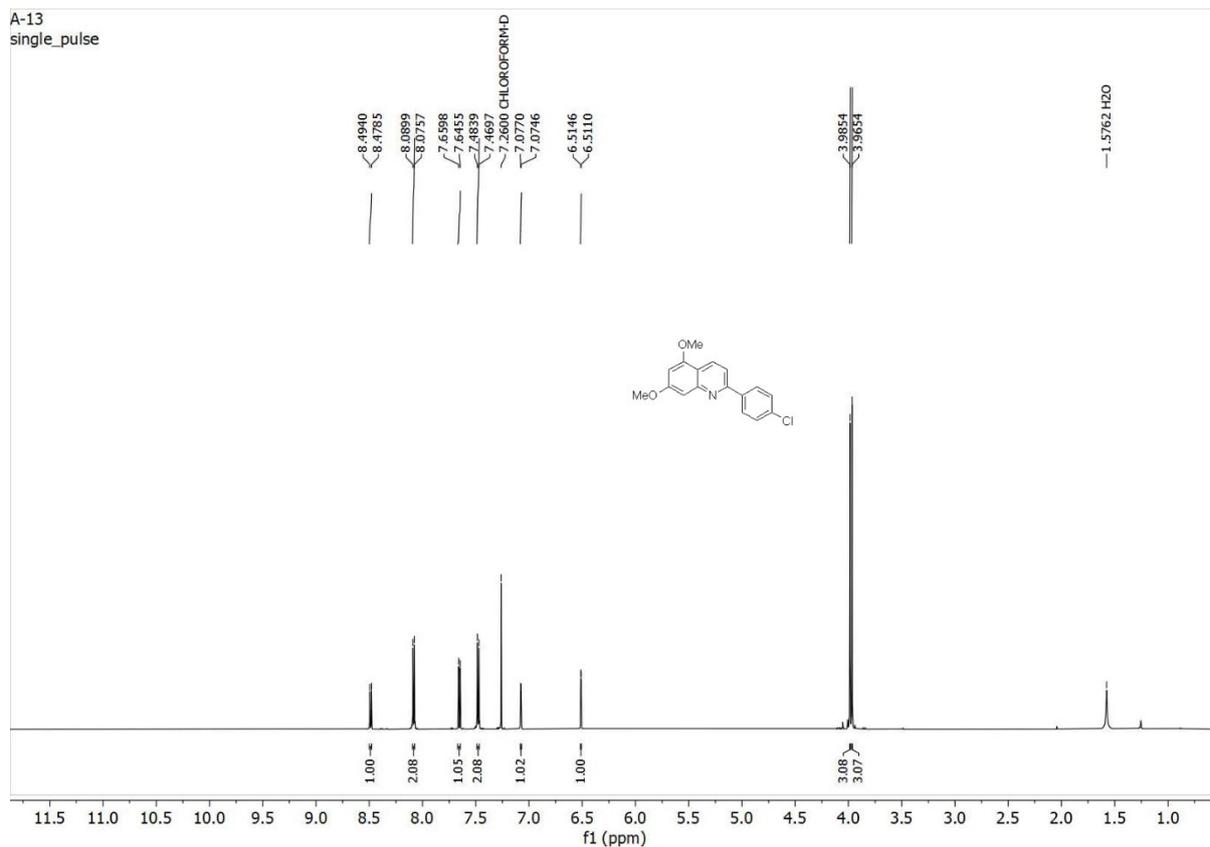


Figure S45. ¹H NMR Spectrum of **17d** in CDCl₃.

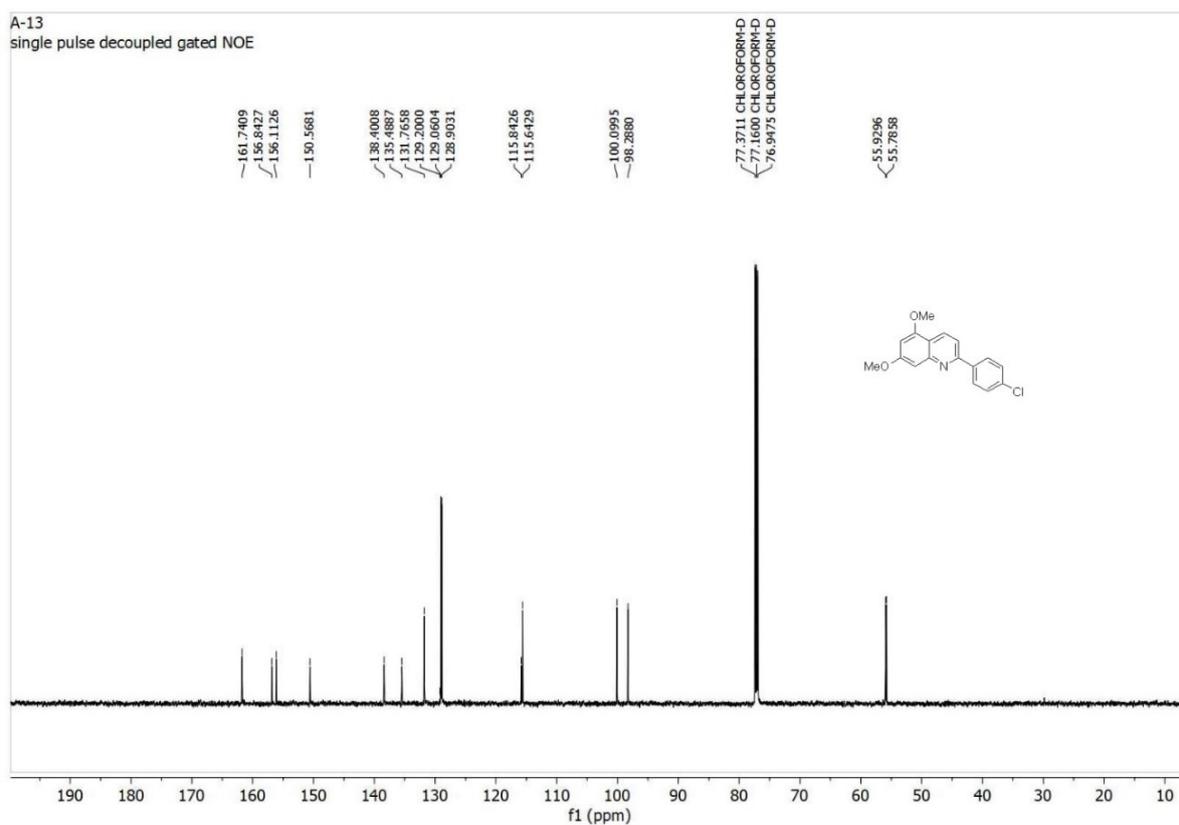


Figure S46. ¹³C NMR Spectrum of **17d** in CDCl₃.

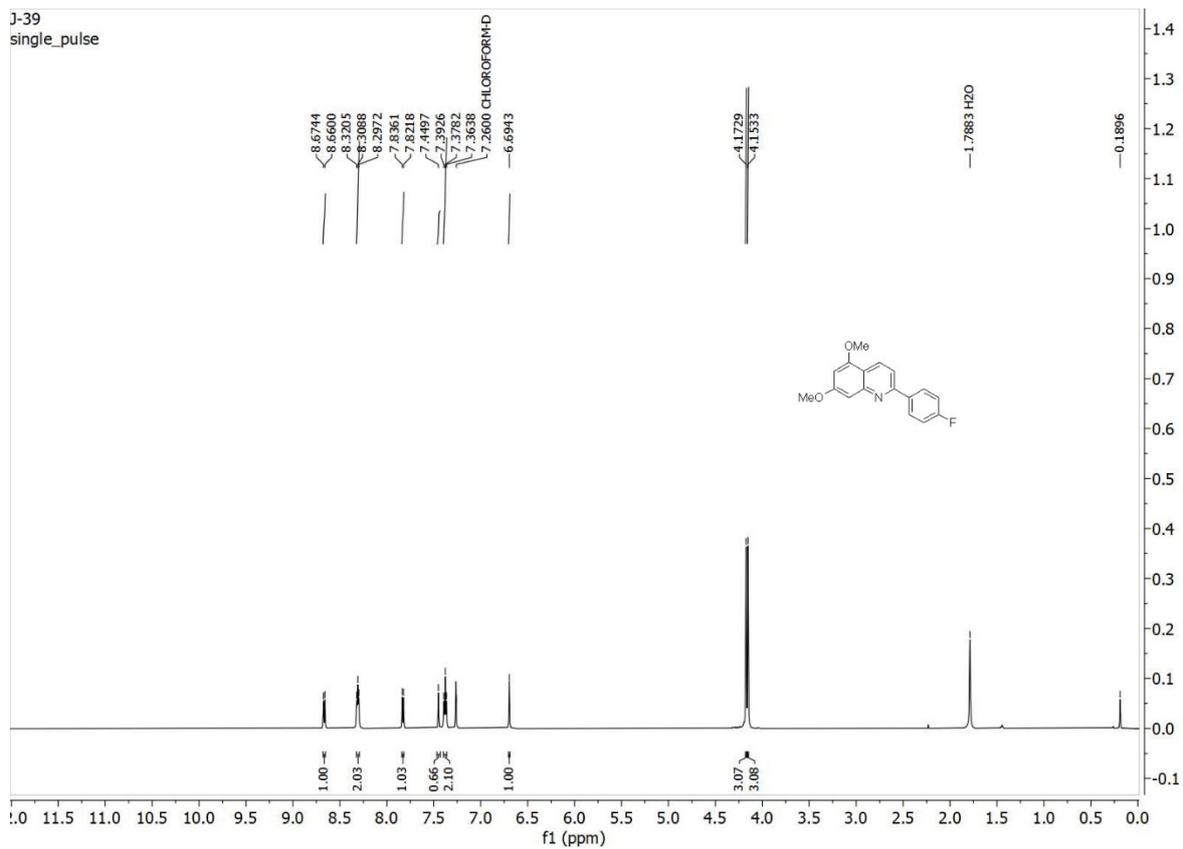


Figure S47. ¹H NMR Spectrum of **18d** in CDCl₃.

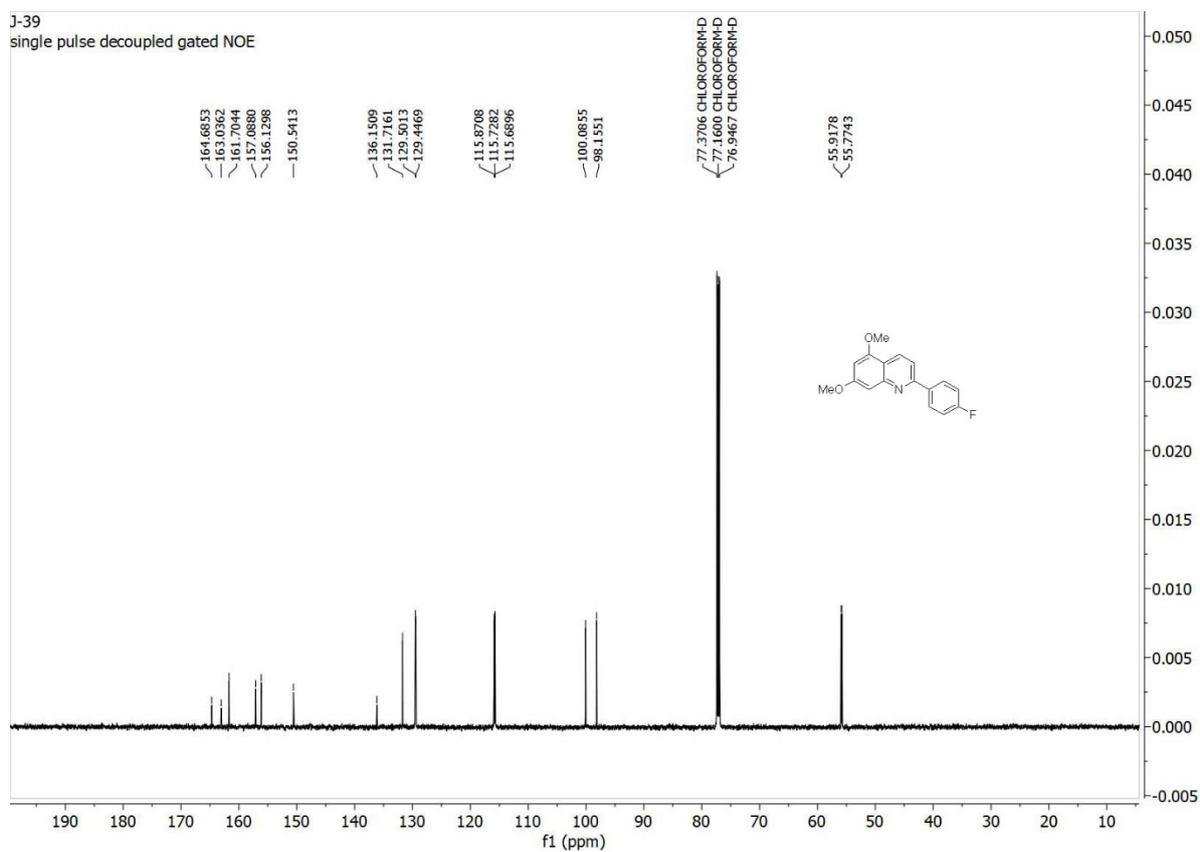


Figure S48. ¹³C NMR Spectrum of **18d** in CDCl₃.

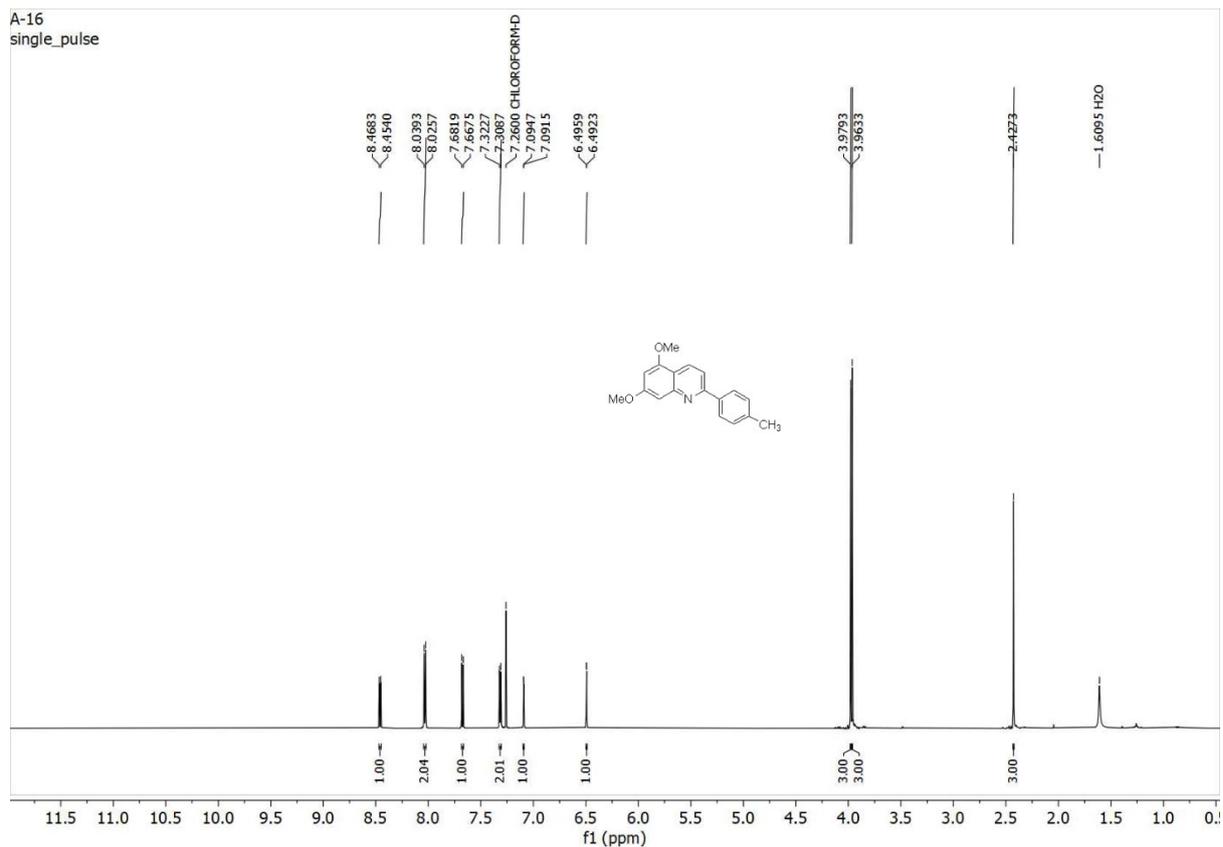


Figure S49. ¹H NMR Spectrum of **20d** in CDCl₃.

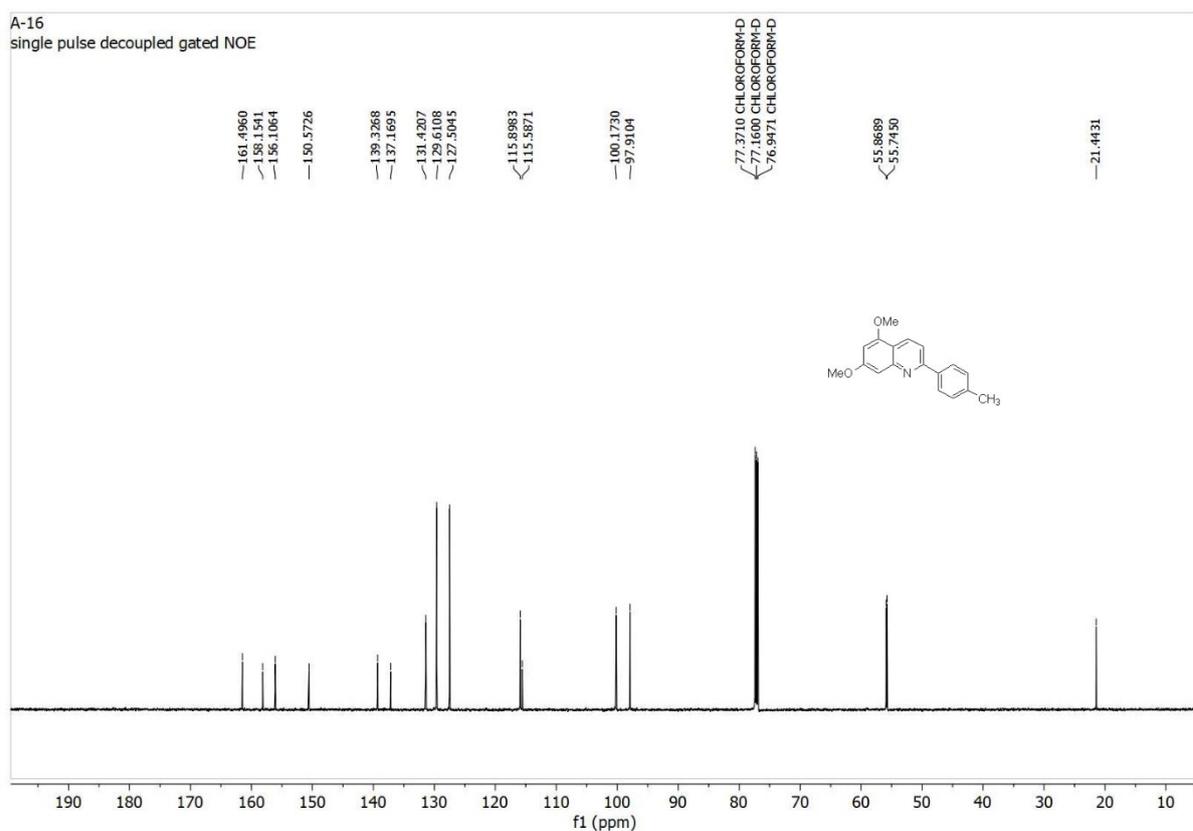


Figure S50. ¹³C NMR Spectrum of **20d** in CDCl₃.

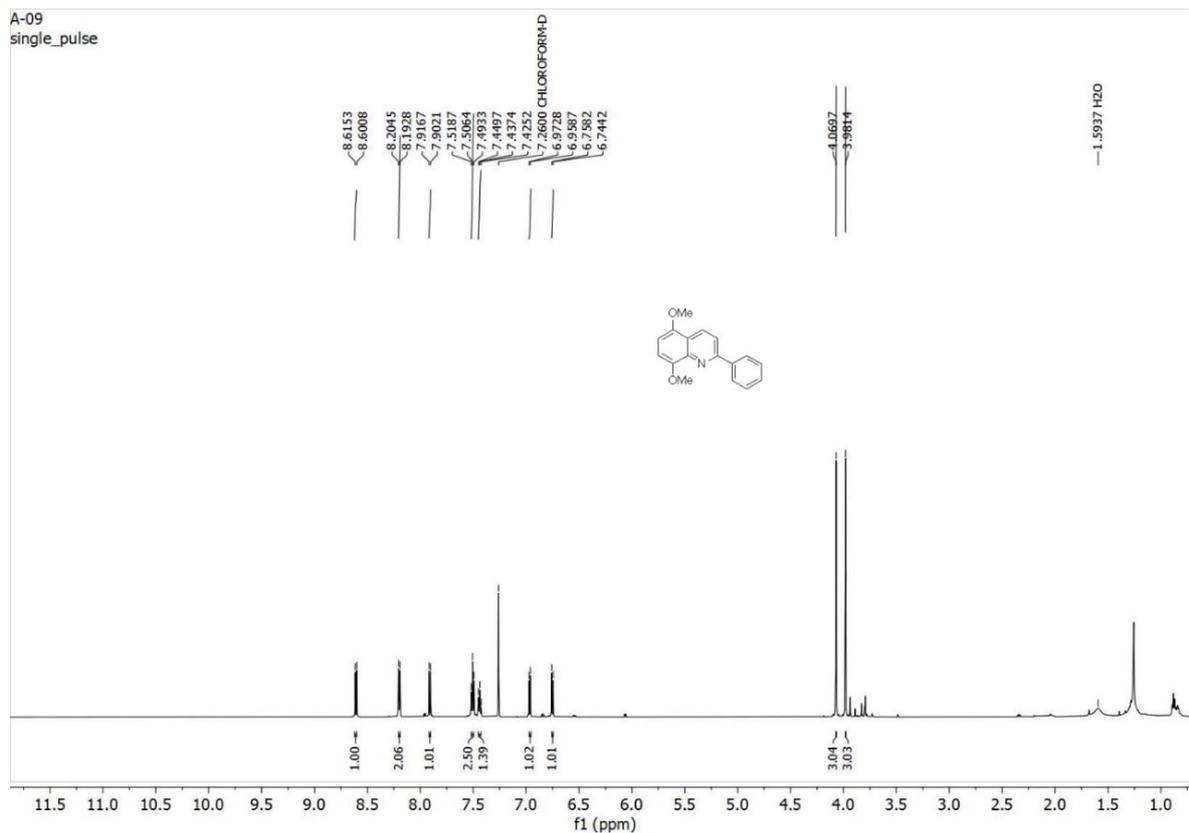


Figure S51. ¹H NMR Spectrum of **15e** in CDCl₃.

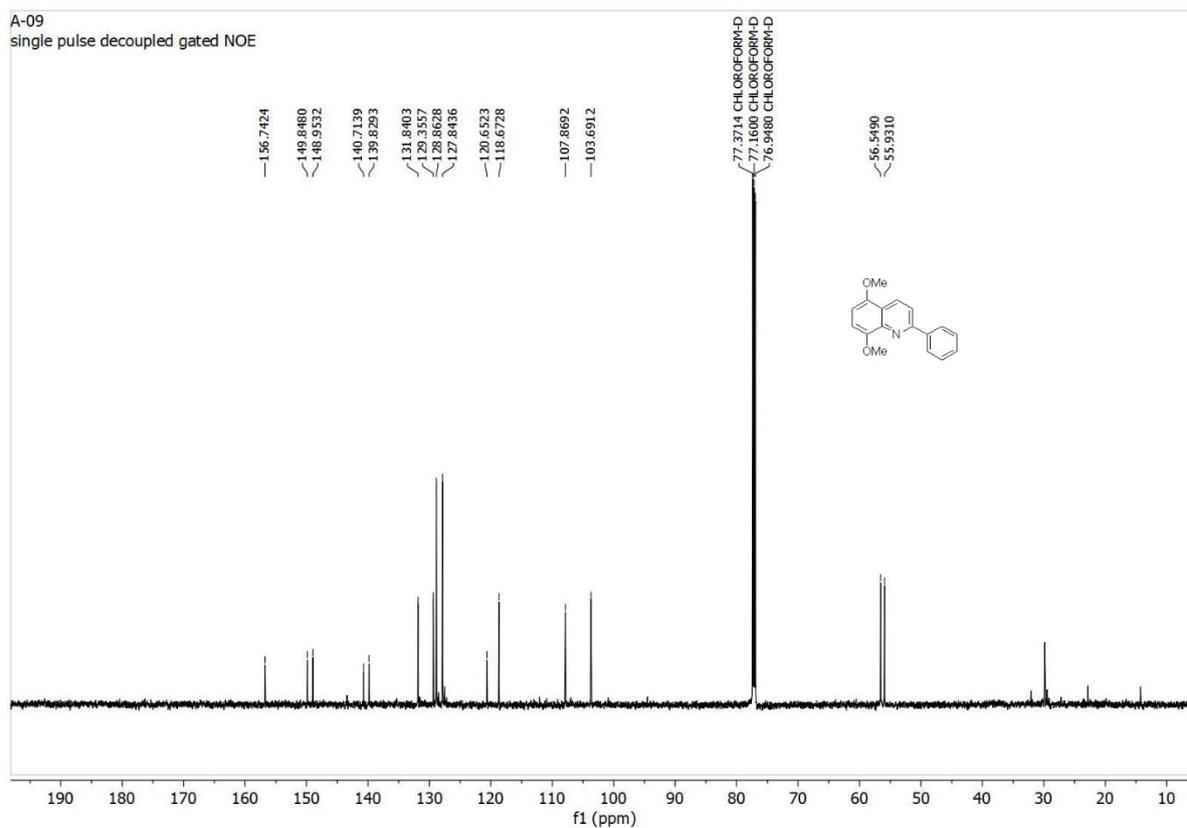


Figure S52. ¹³C NMR Spectrum of **15e** in CDCl₃.

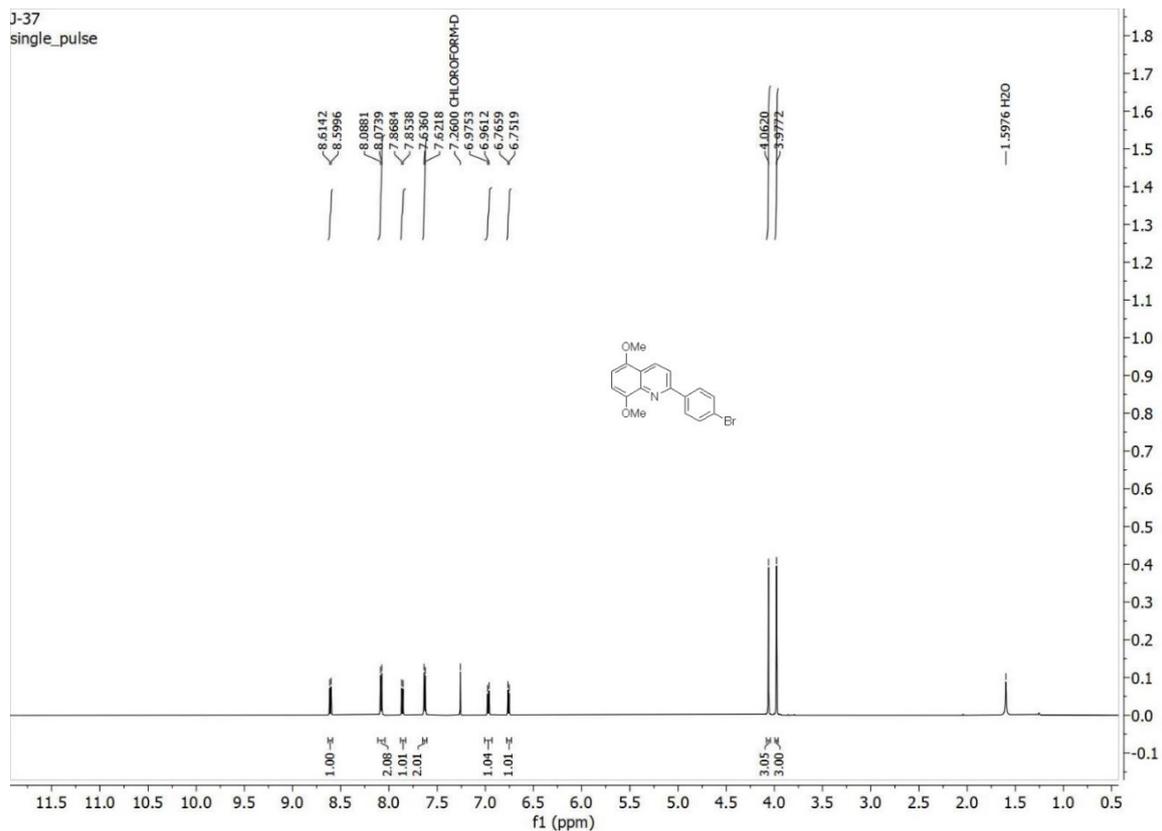


Figure S53. ^1H NMR Spectrum of **16e** in CDCl_3 .

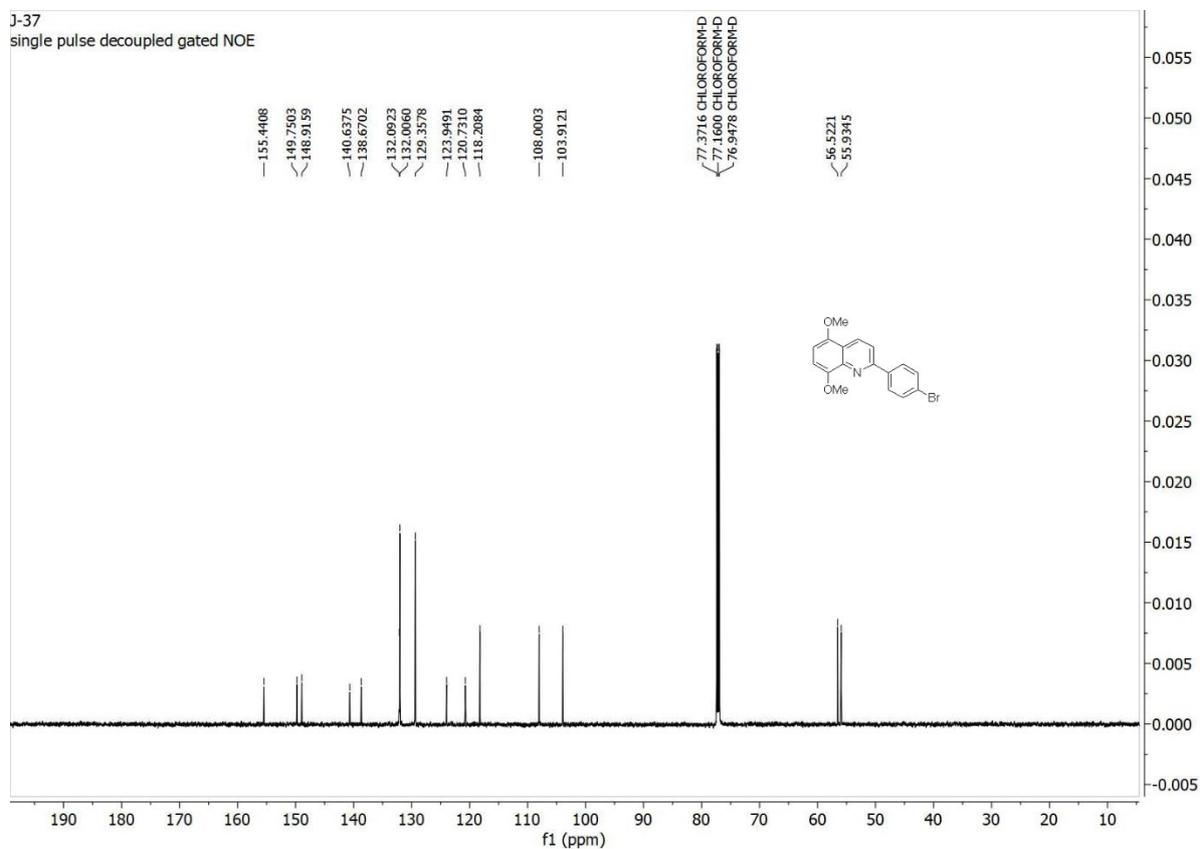


Figure S54. ^{13}C NMR Spectrum of **16e** in CDCl_3 .

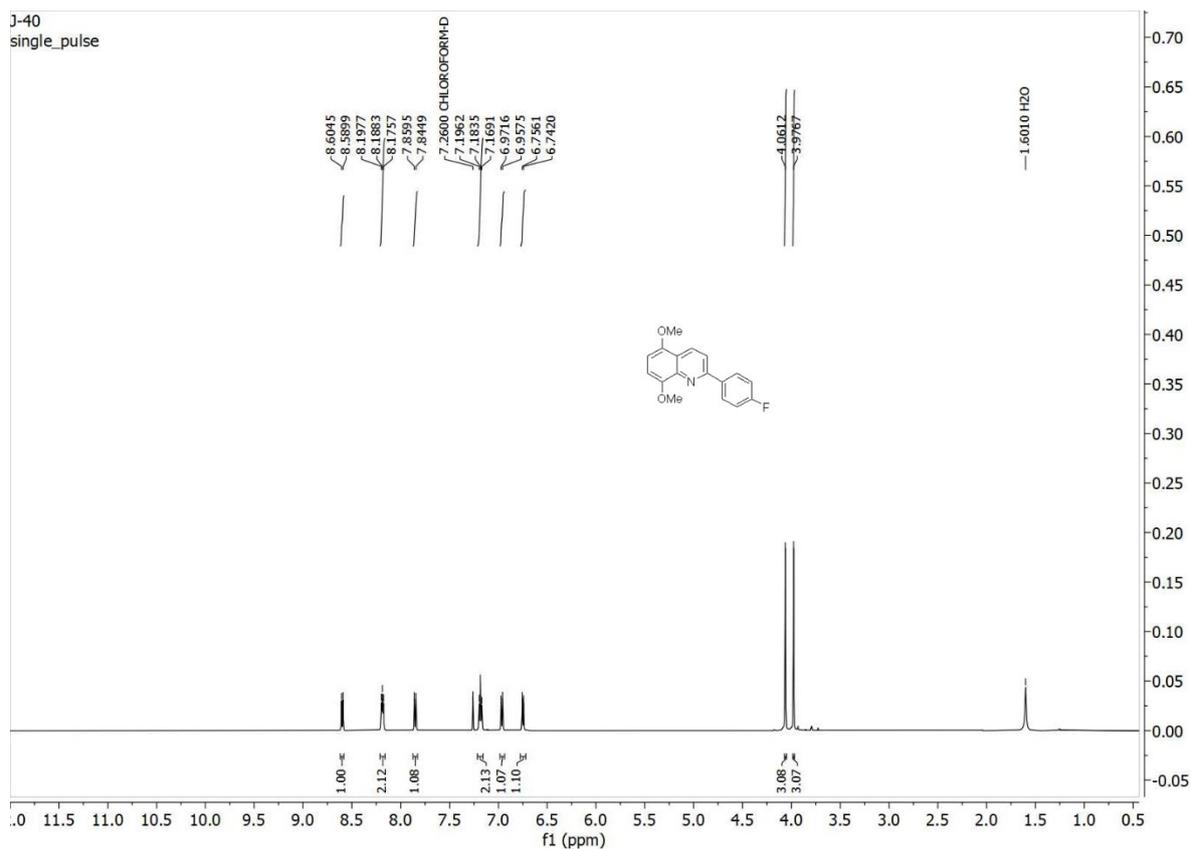


Figure S55. ^1H NMR Spectrum of **18e** in CDCl_3 .

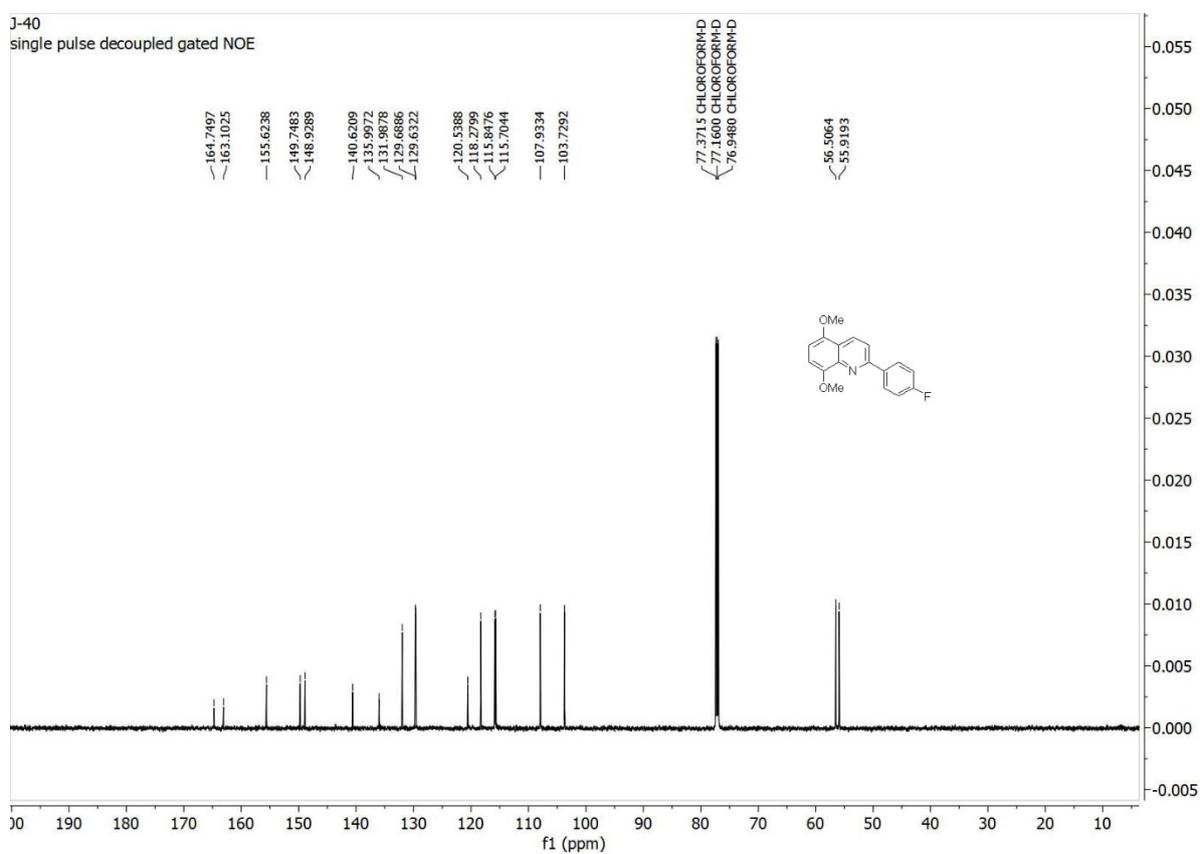


Figure S56. ^{13}C NMR Spectrum of **18e** in CDCl_3 .

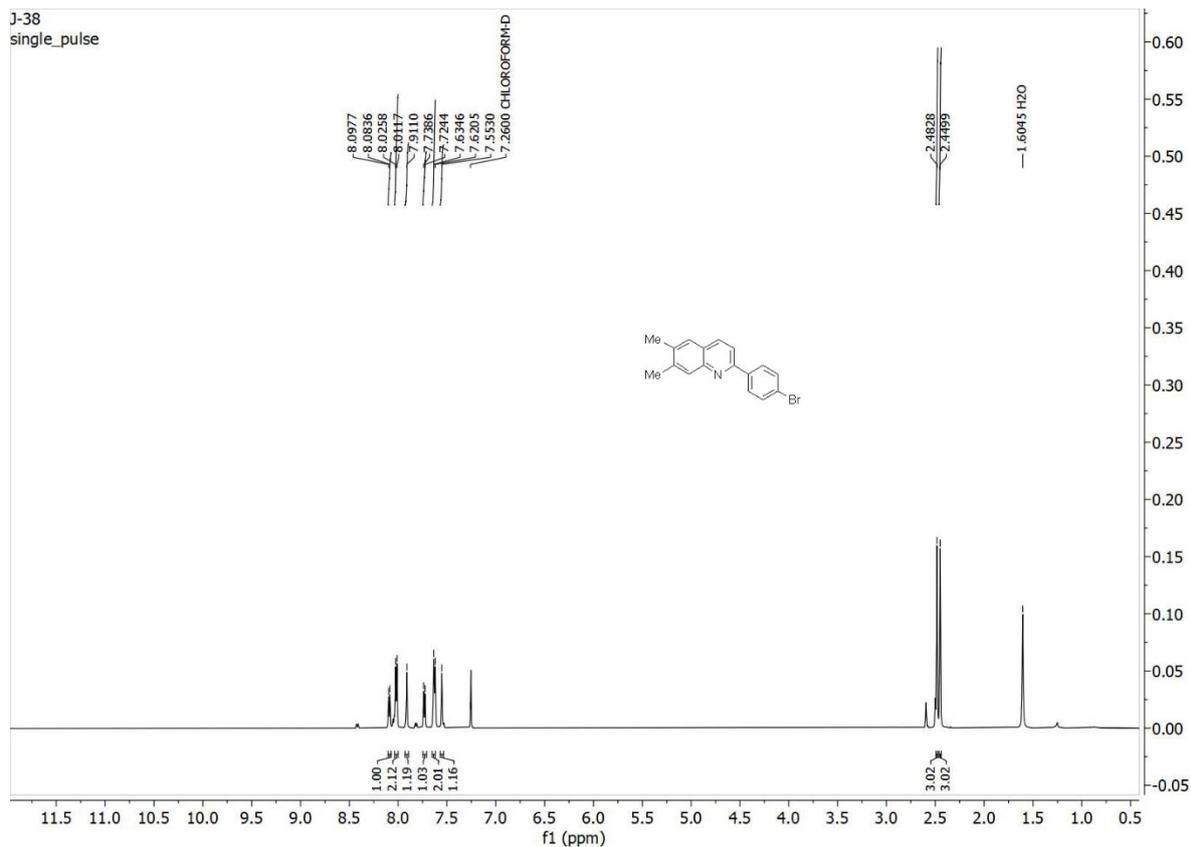


Figure S57. ^1H NMR Spectrum of **16f** in CDCl_3 .

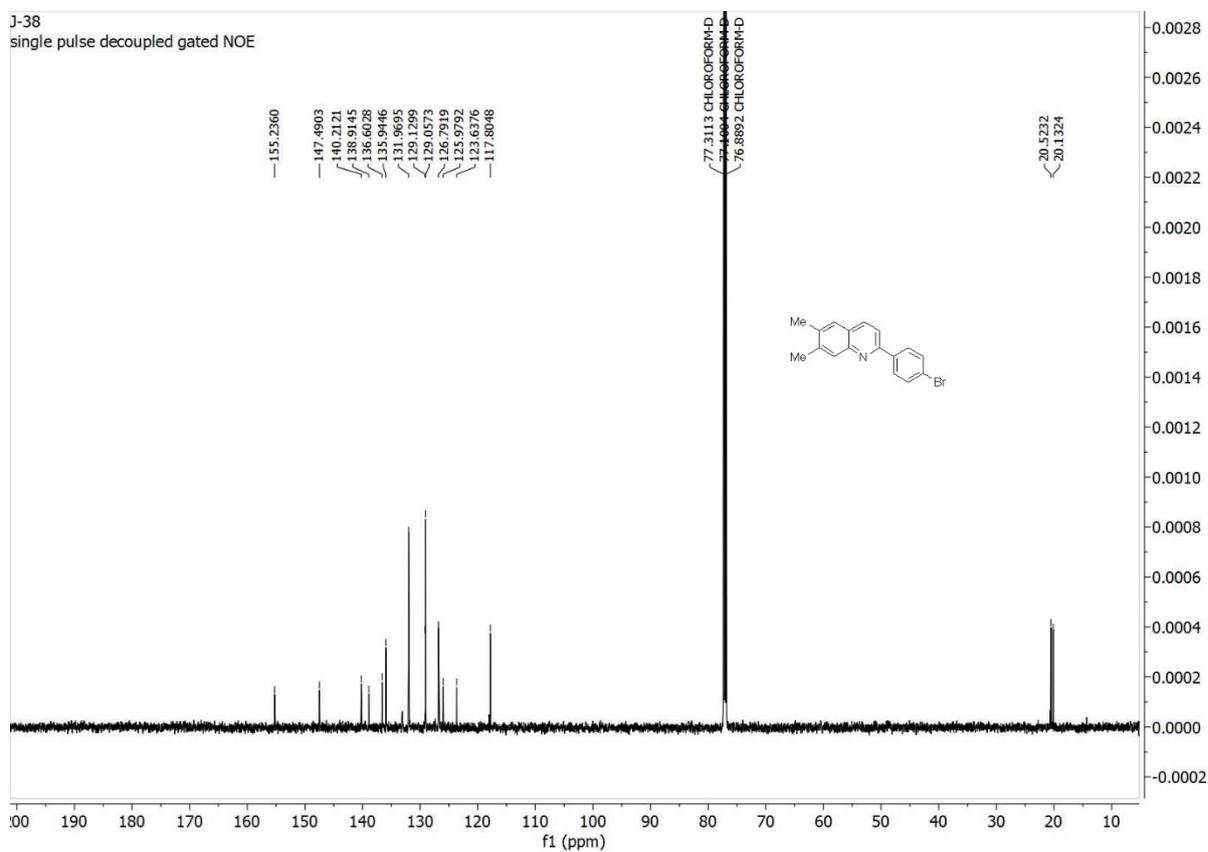


Figure S58. ^{13}C NMR Spectrum of **16f** in CDCl_3 .

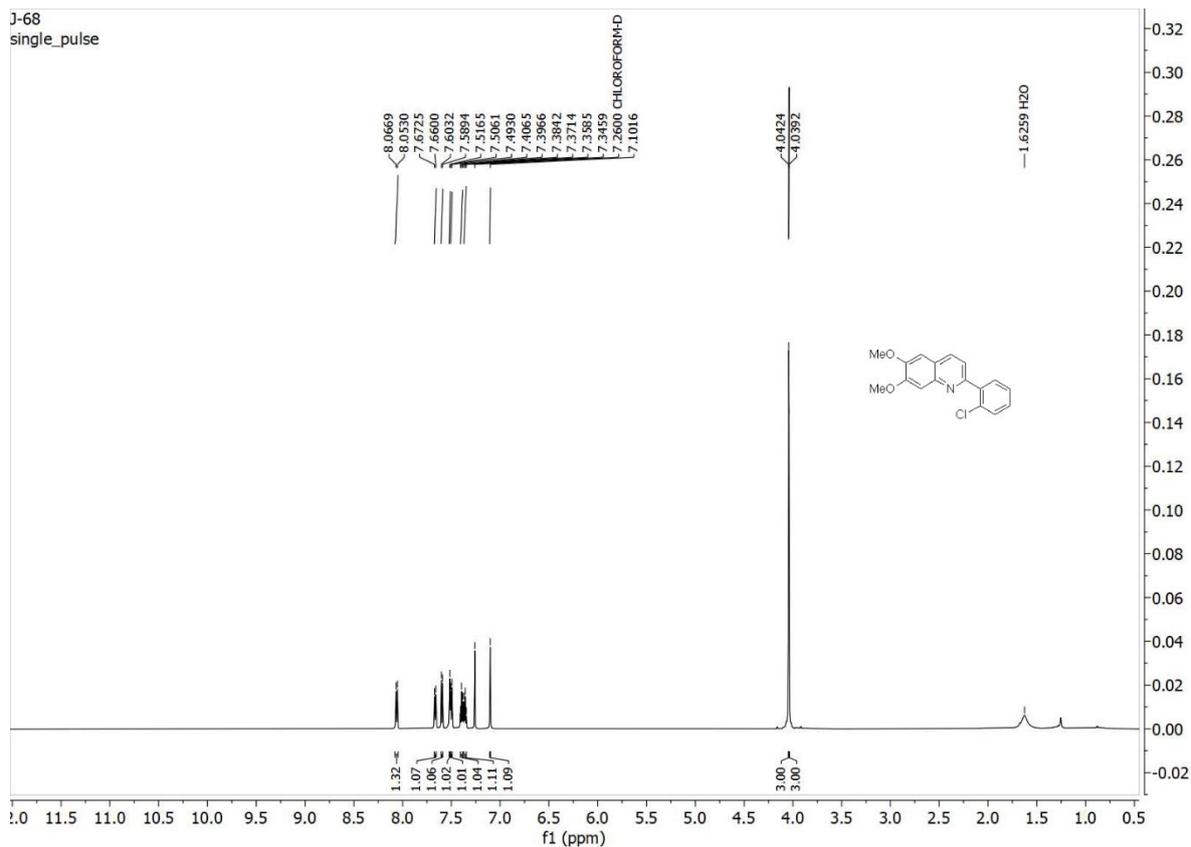


Figure S59. ¹H NMR Spectrum of **23a** in CDCl₃.

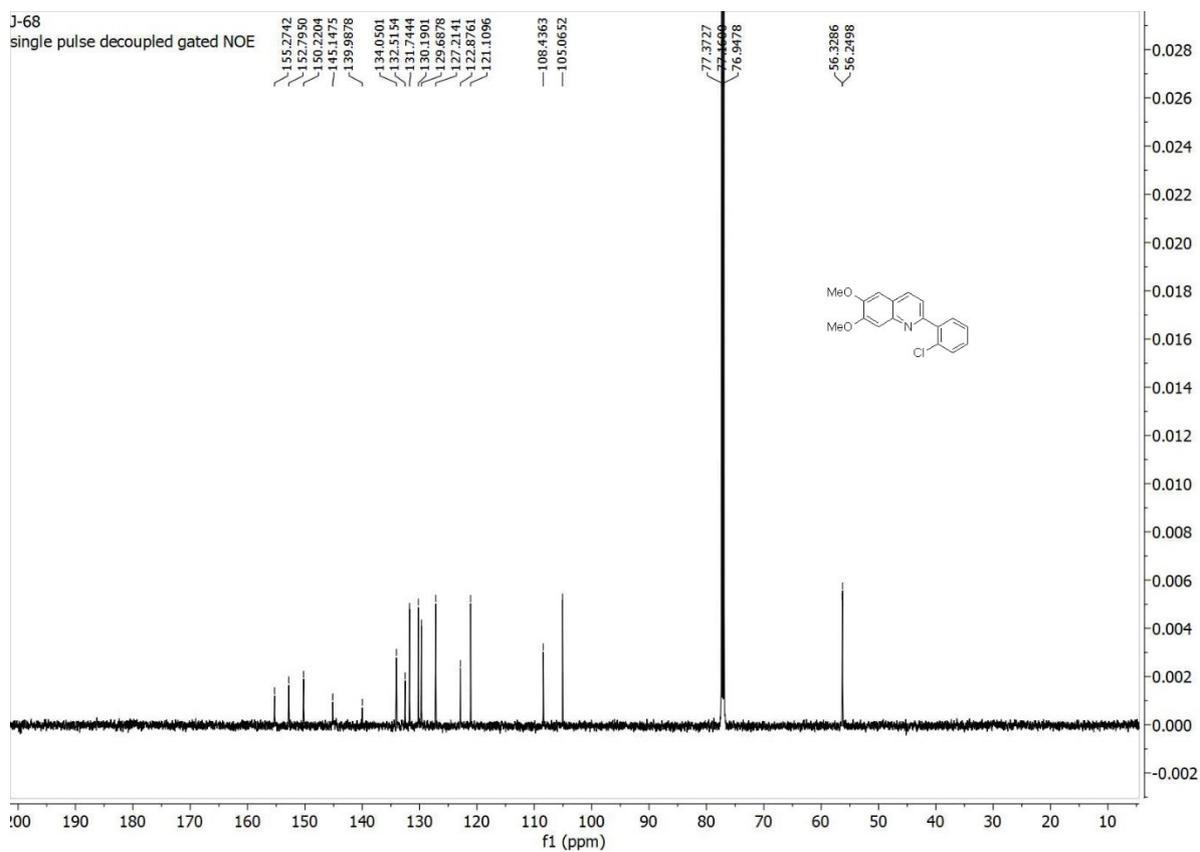


Figure S60. ¹³C NMR Spectrum of **23a** in CDCl₃.

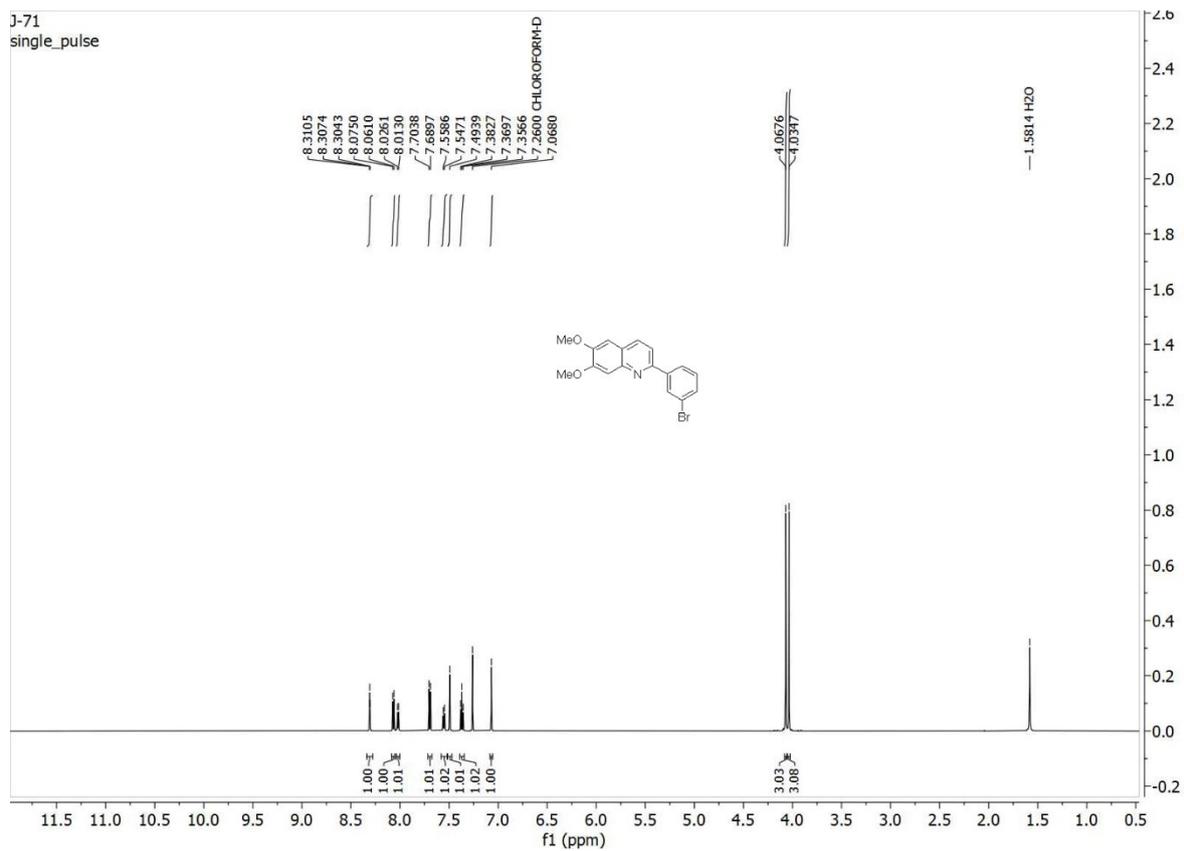


Figure S61. ^1H NMR Spectrum of **24a** in CDCl_3 .

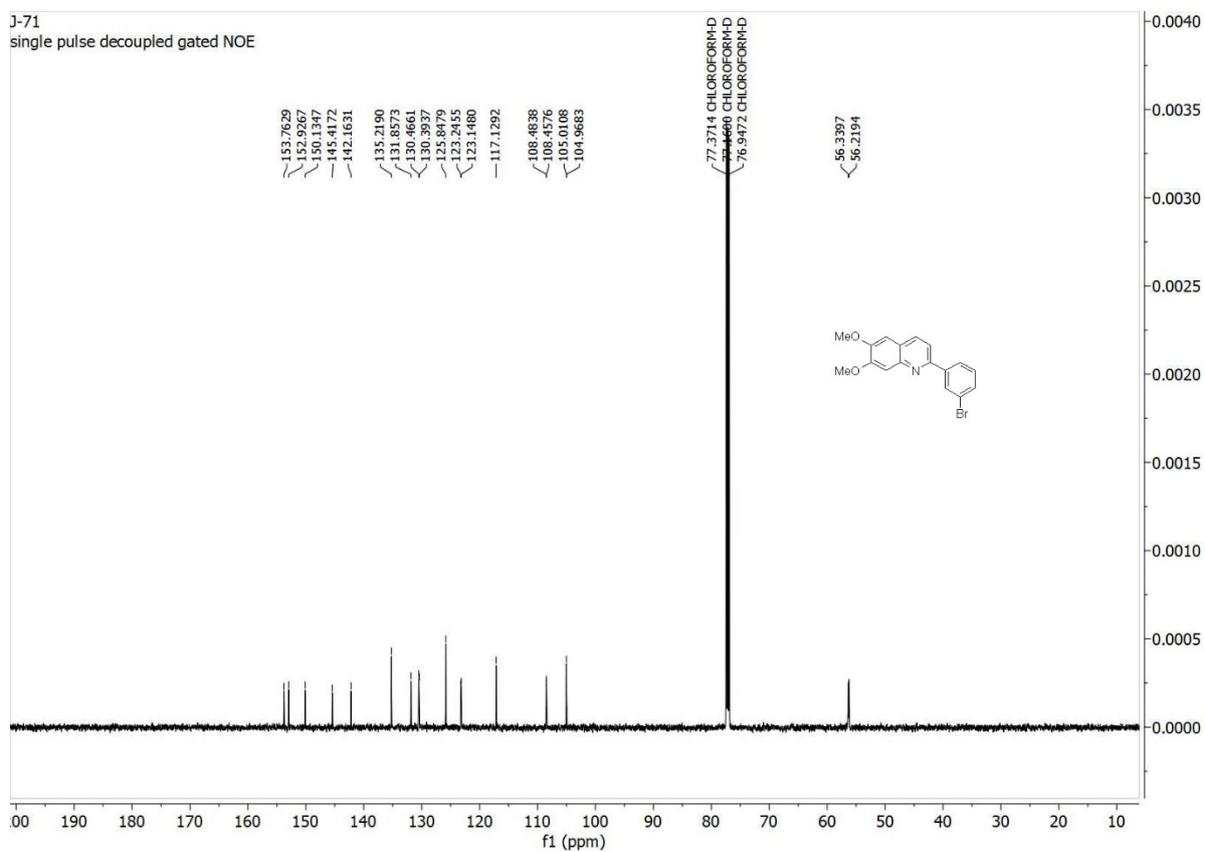


Figure S62. ^{13}C NMR Spectrum of **24a** in CDCl_3 .

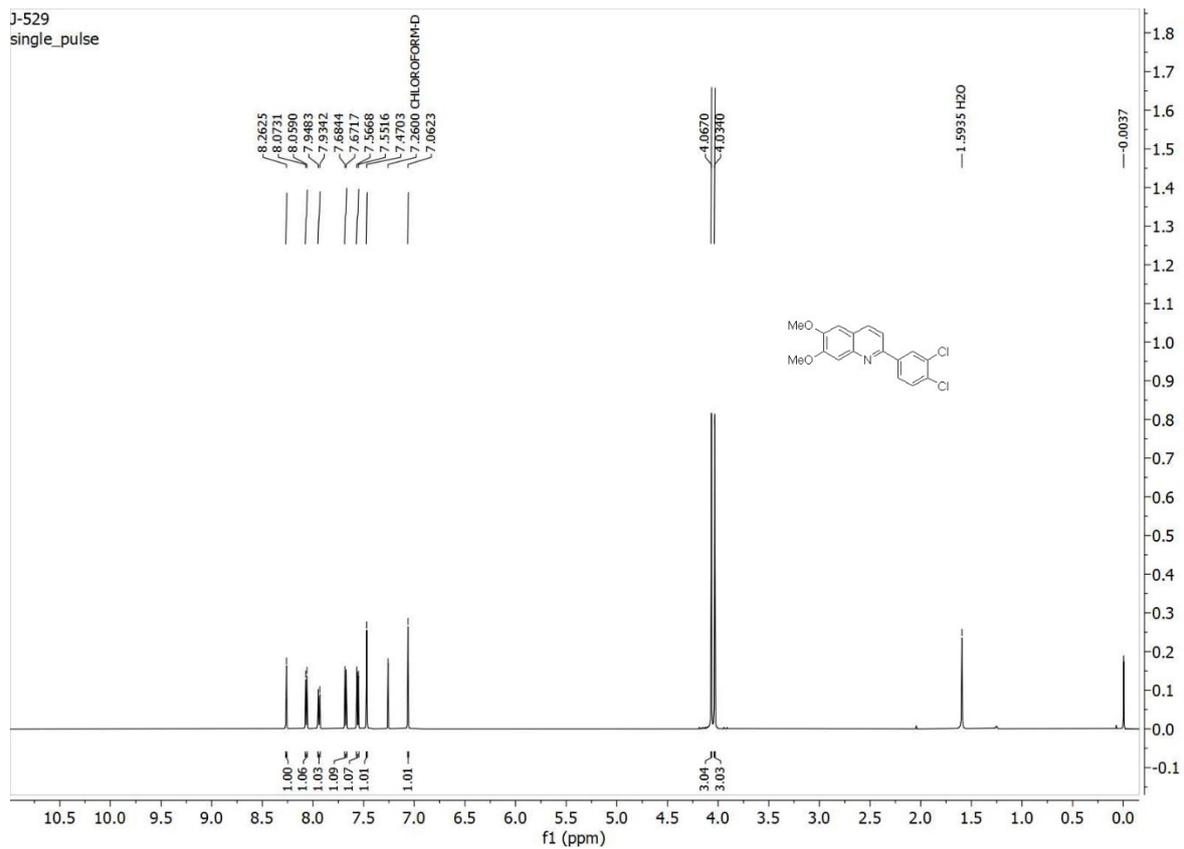


Figure S63. ^1H NMR Spectrum of 25a in CDCl_3 .

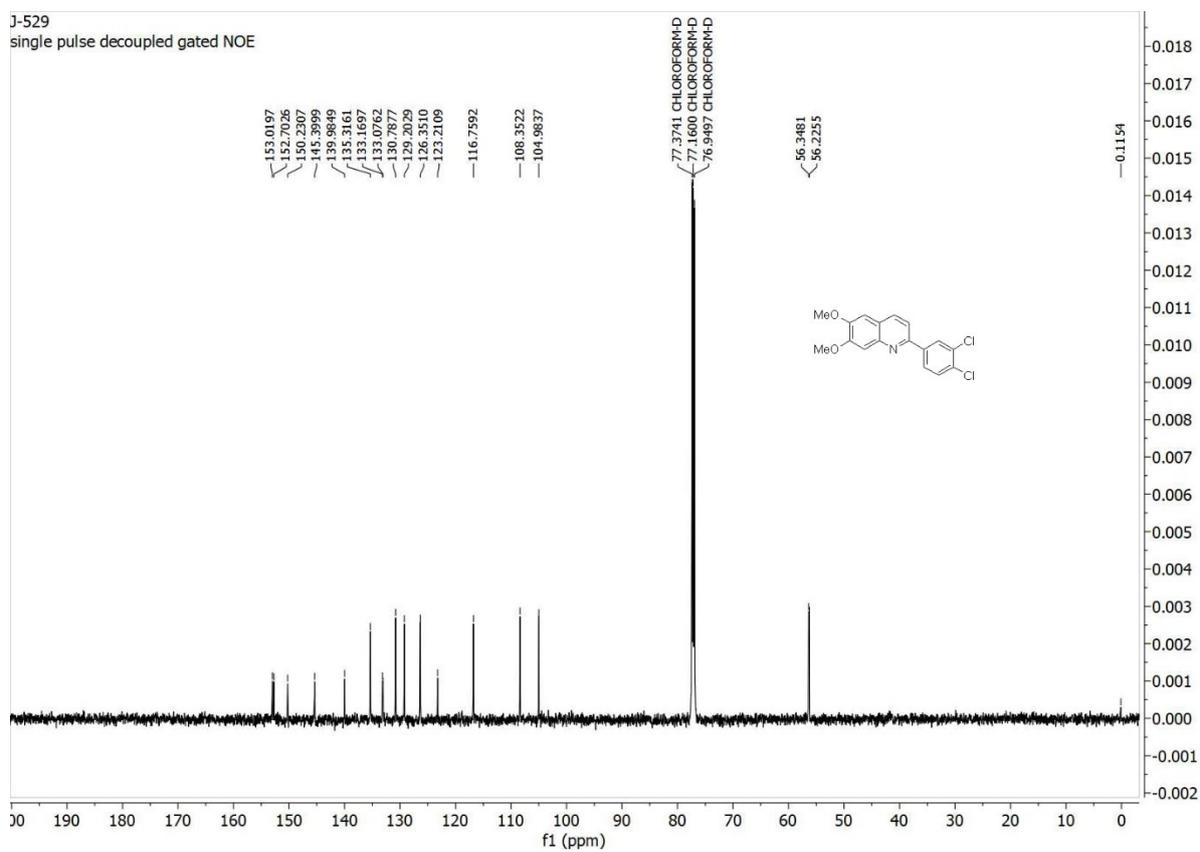


Figure S64. ^{13}C NMR Spectrum of 25a in CDCl_3 .

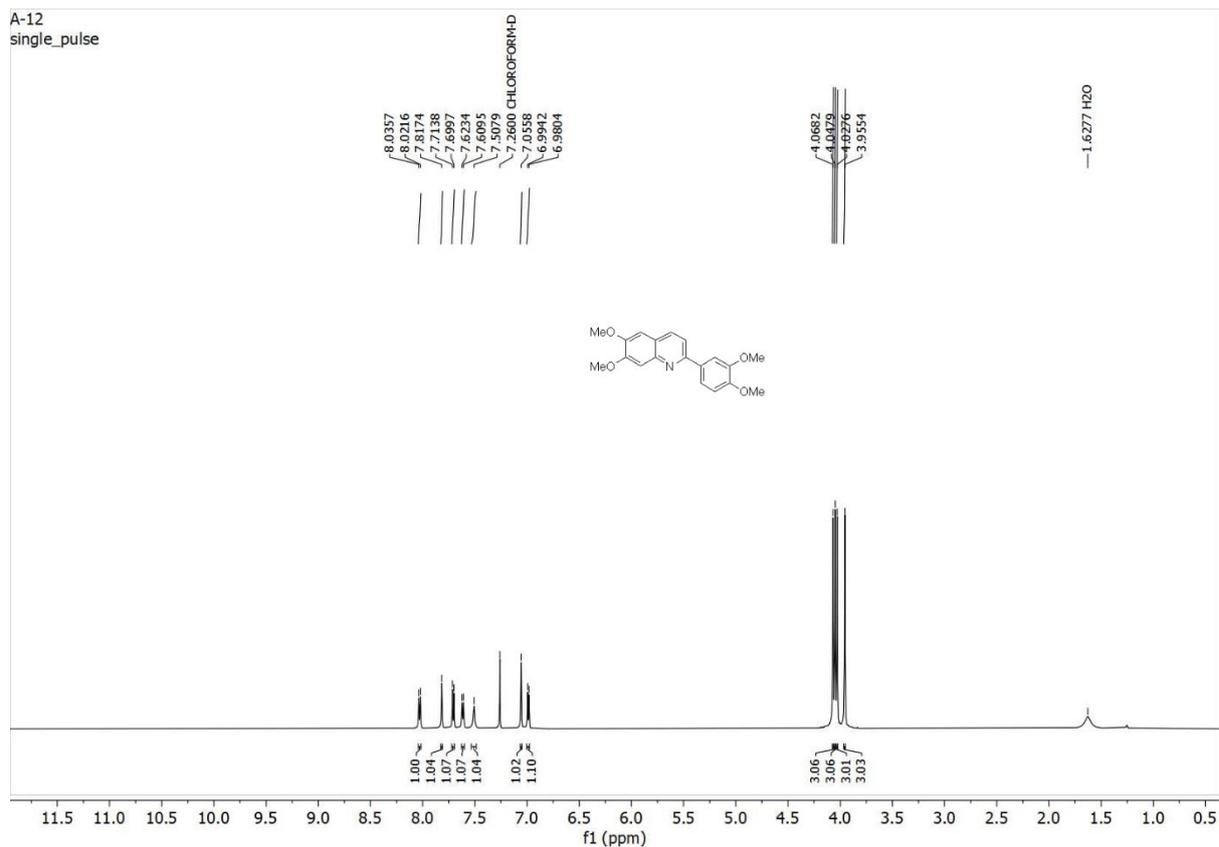


Figure S65. ^1H NMR Spectrum of **26a** in CDCl_3 .

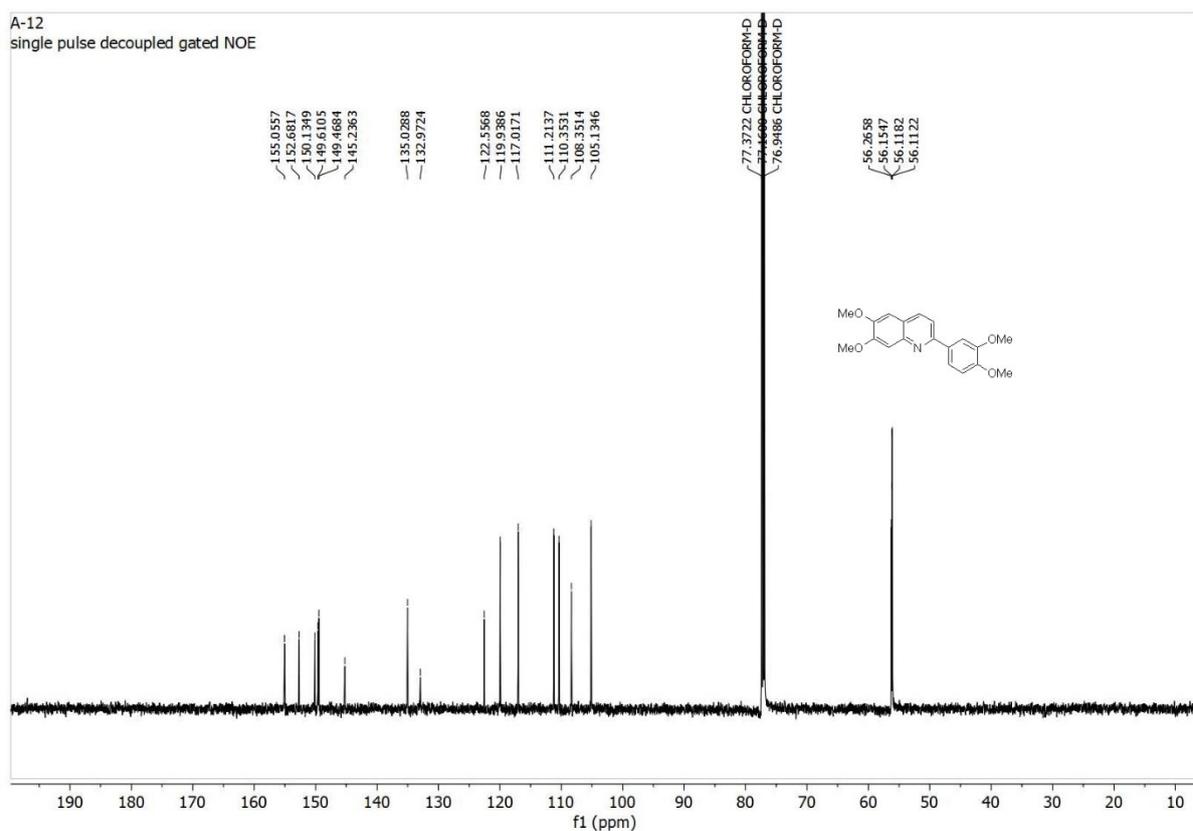


Figure S66. ^{13}C NMR Spectrum of **26a** in CDCl_3 .

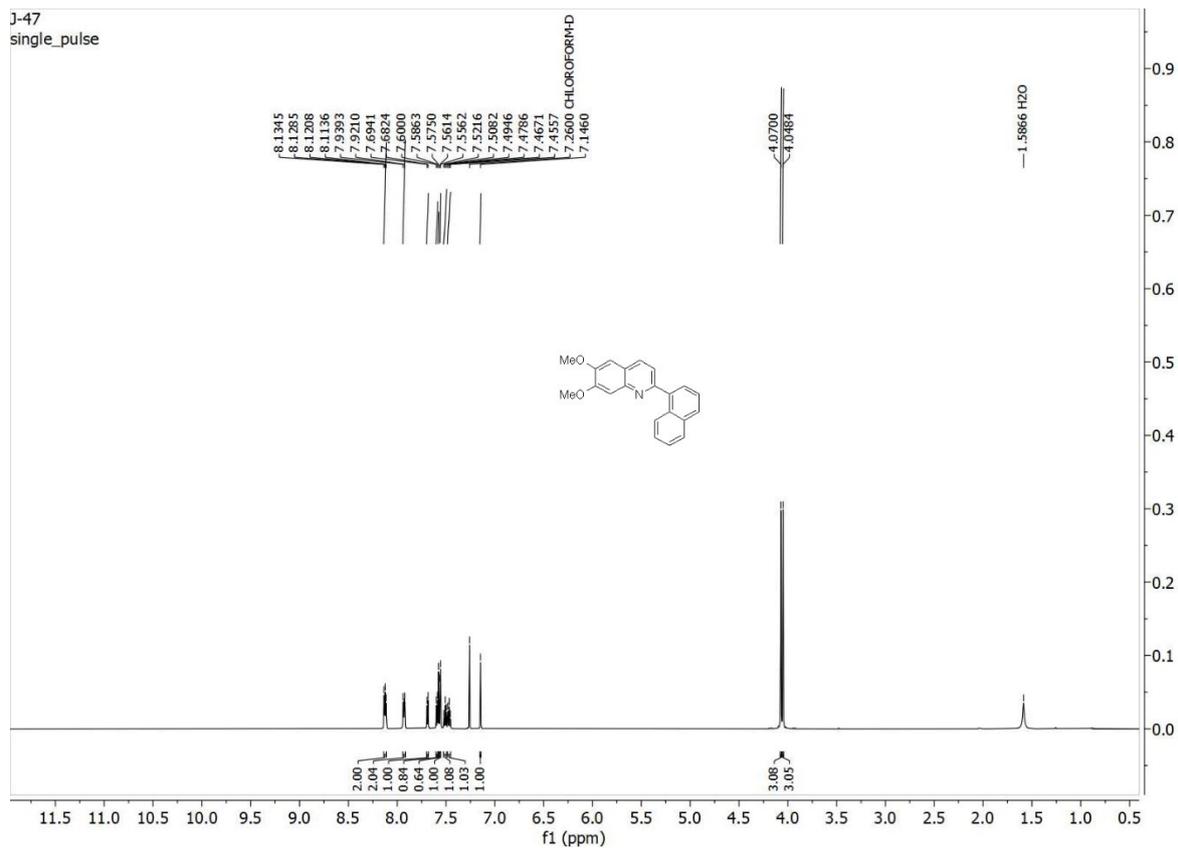


Figure S67. ¹H NMR Spectrum of **27a** in CDCl₃.

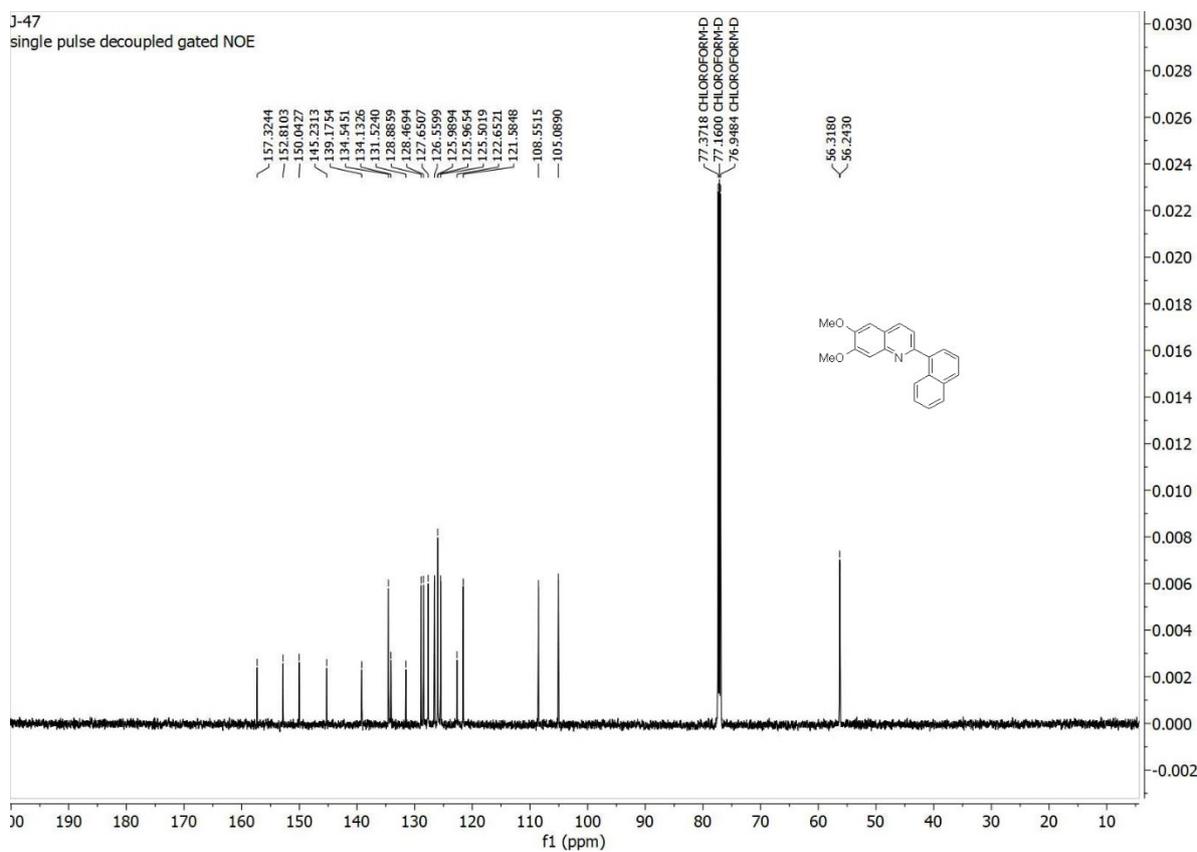


Figure S68. ¹³C NMR Spectrum of **27a** in CDCl₃.

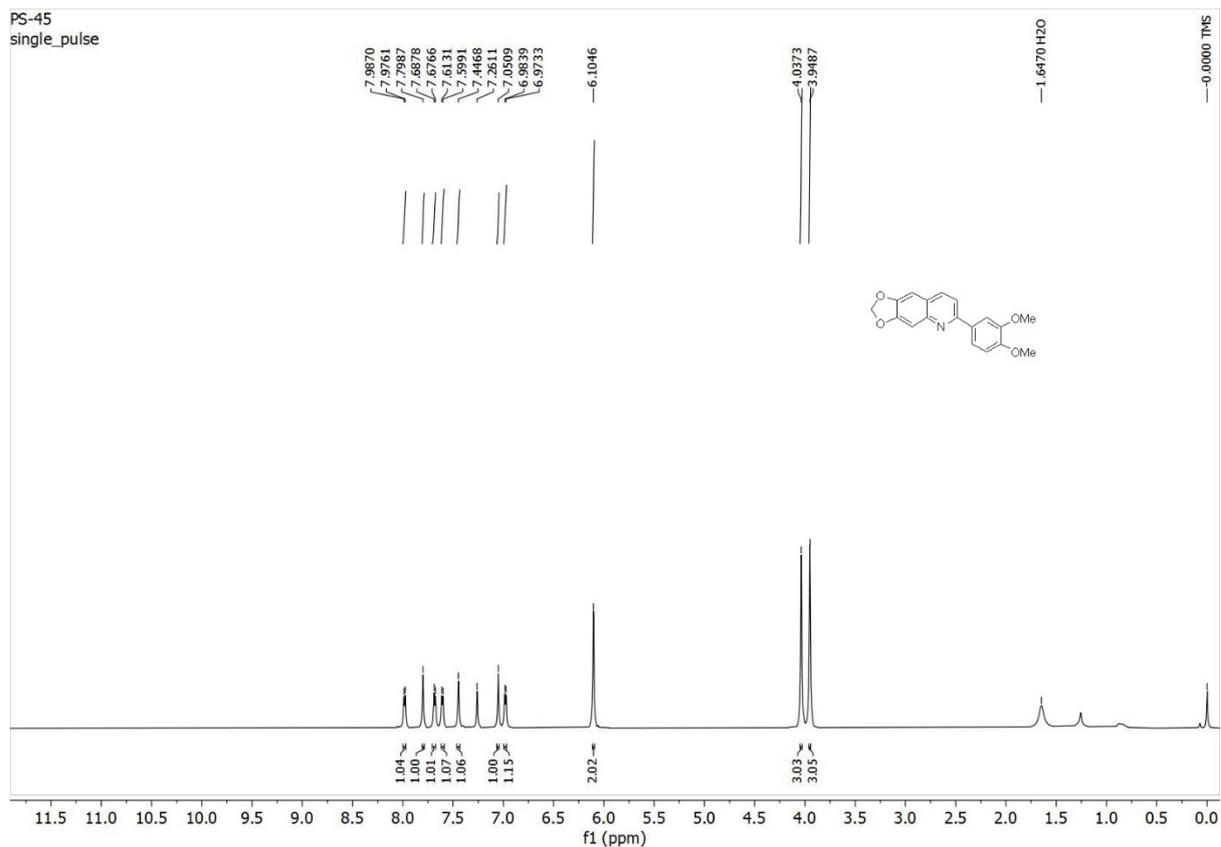


Figure S69. ^1H NMR Spectrum of **26b** in CDCl_3 .

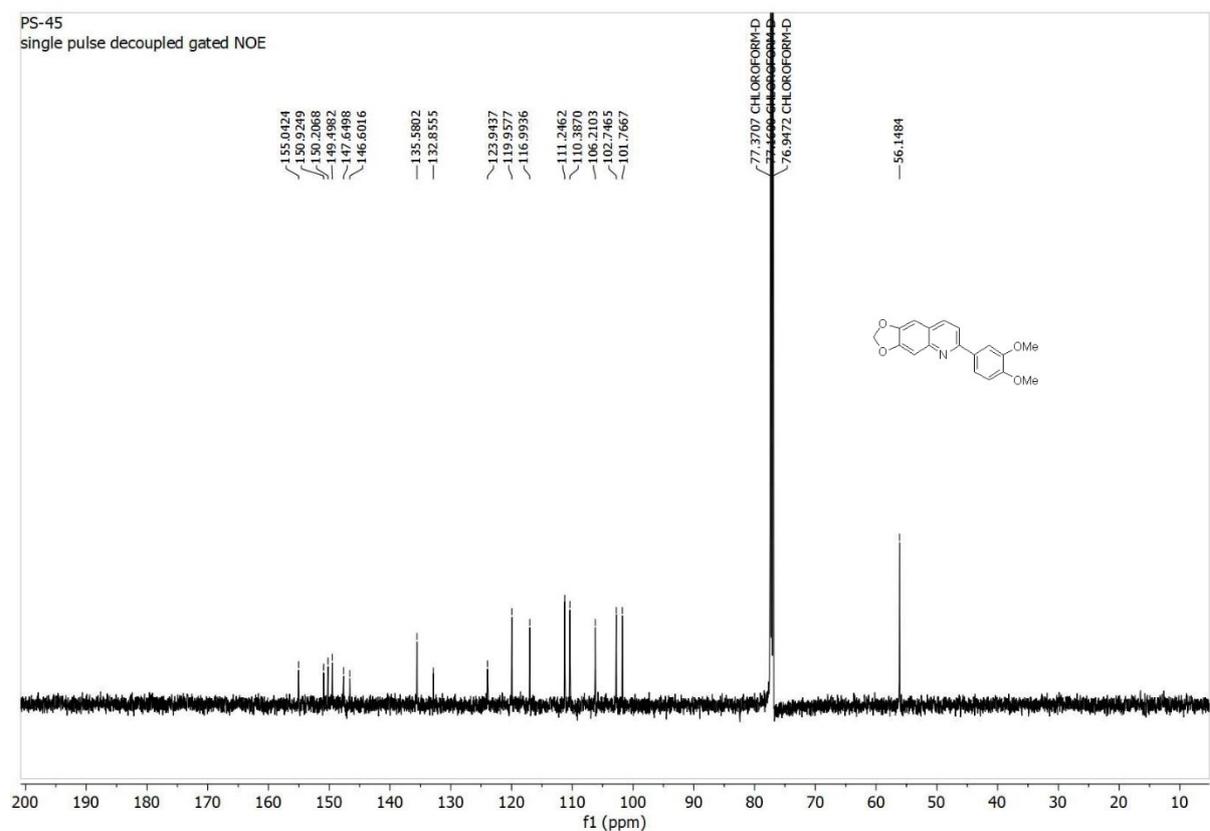


Figure S70. ^{13}C NMR Spectrum of **26b** in CDCl_3 .

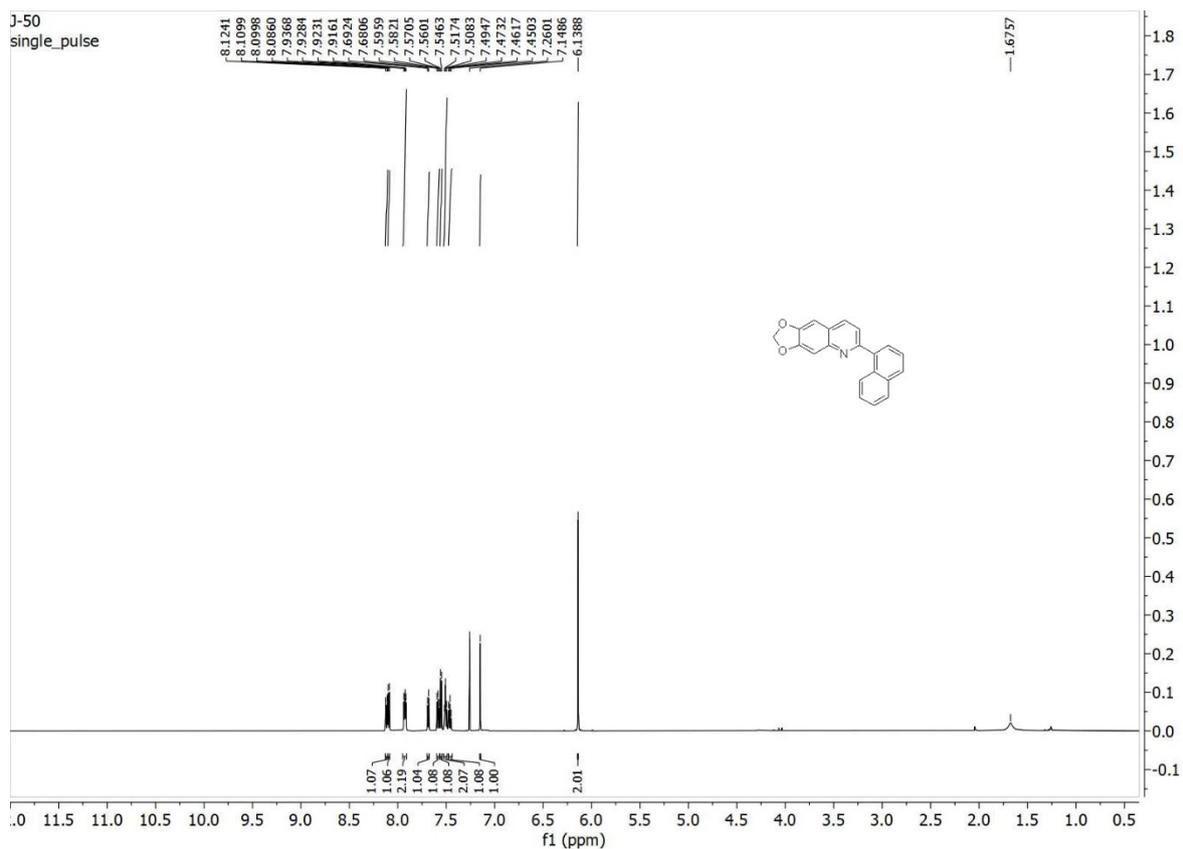


Figure S71. ^1H NMR Spectrum of **27b** in CDCl_3 .

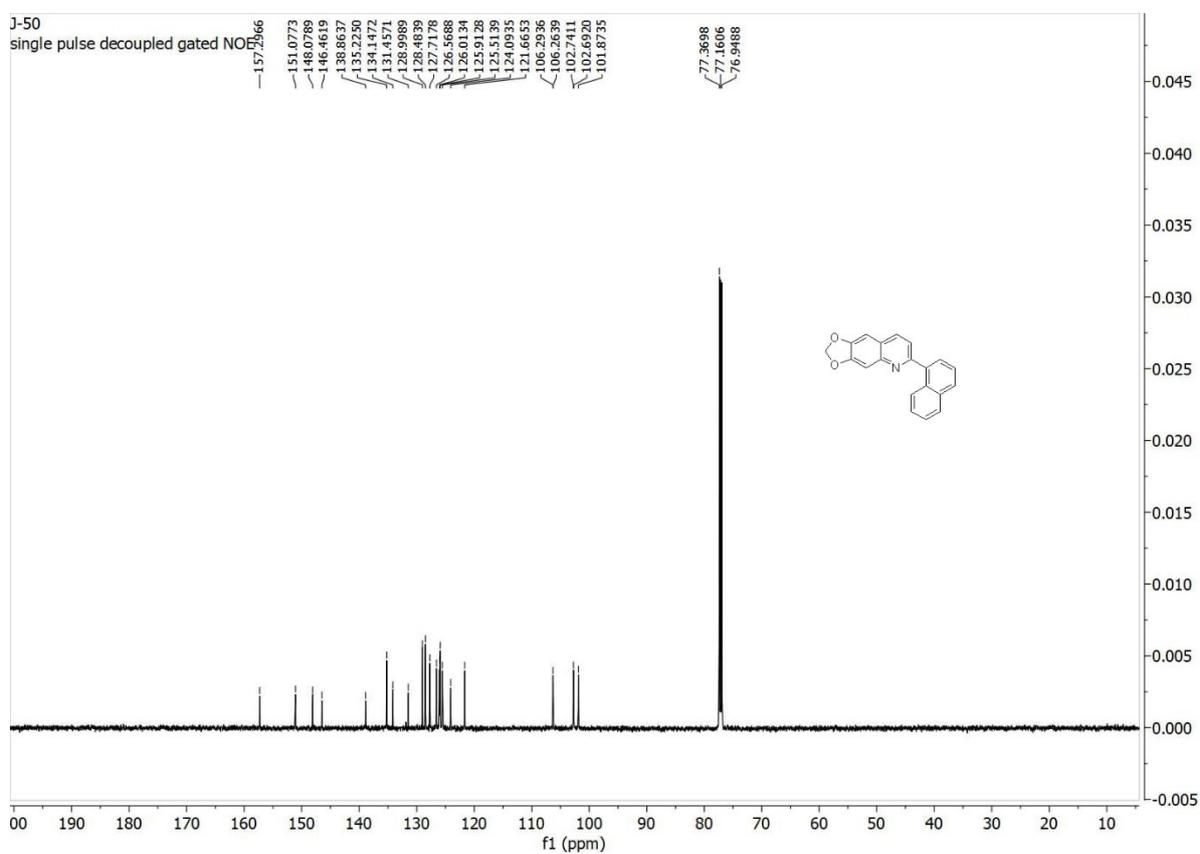


Figure S72. ^{13}C NMR Spectrum of **27b** in CDCl_3 .

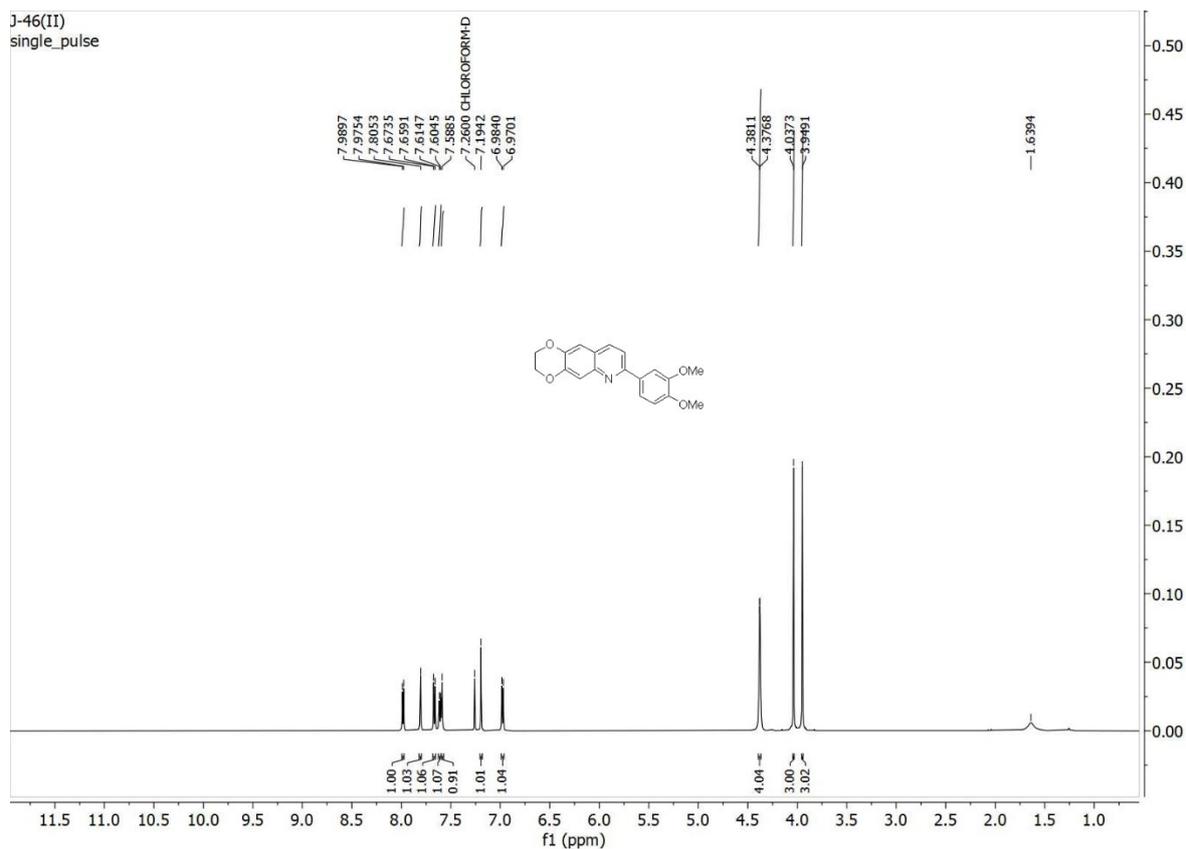


Figure S73. ^1H NMR Spectrum of **26c** in CDCl_3 .

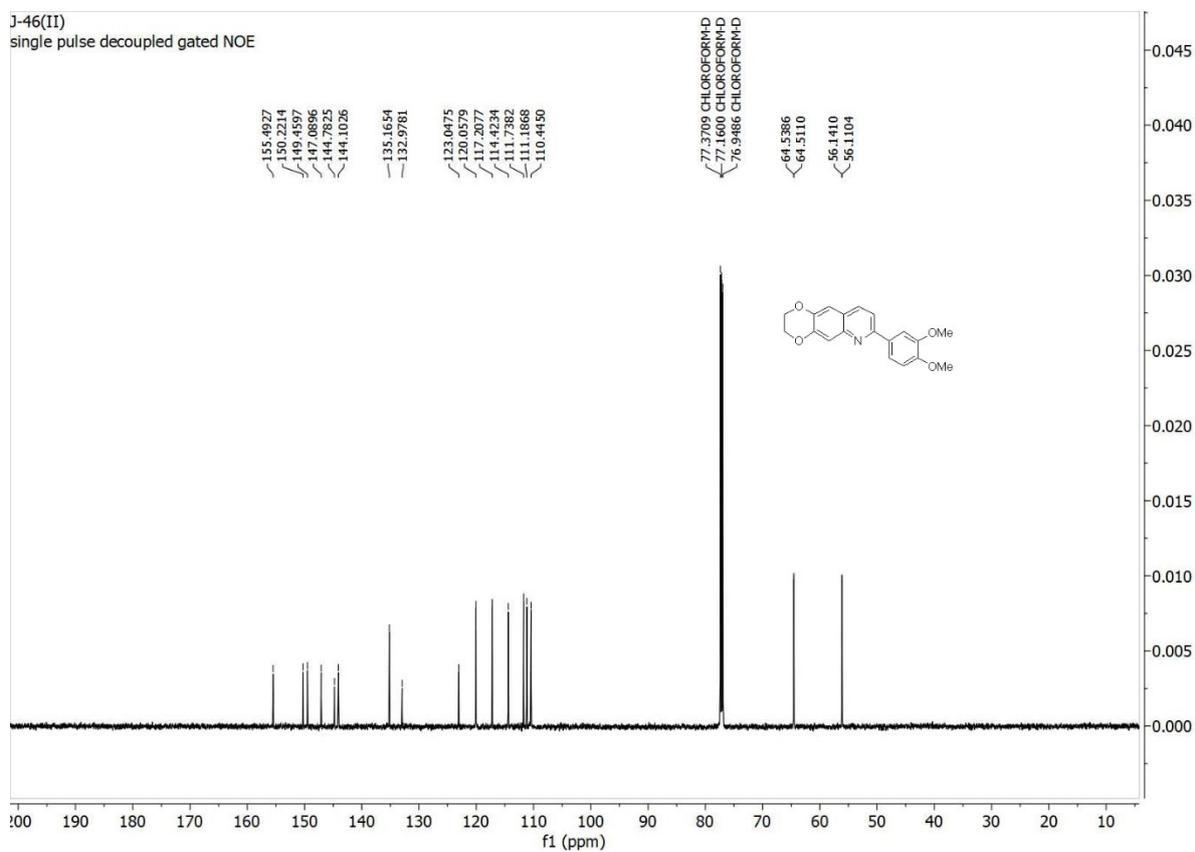


Figure S74. ^{13}C NMR Spectrum of **26c** in CDCl_3 .

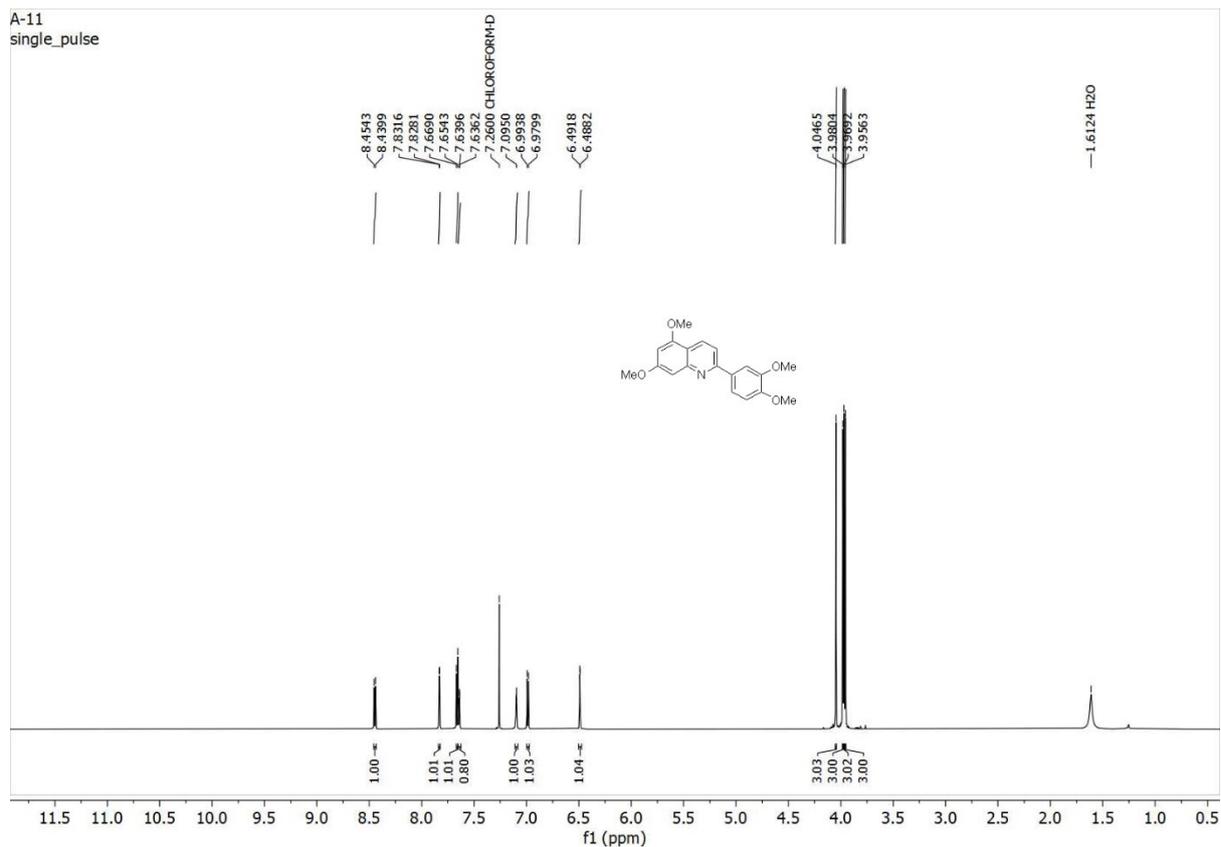


Figure S75. ^1H NMR Spectrum of **26d** in CDCl_3 .

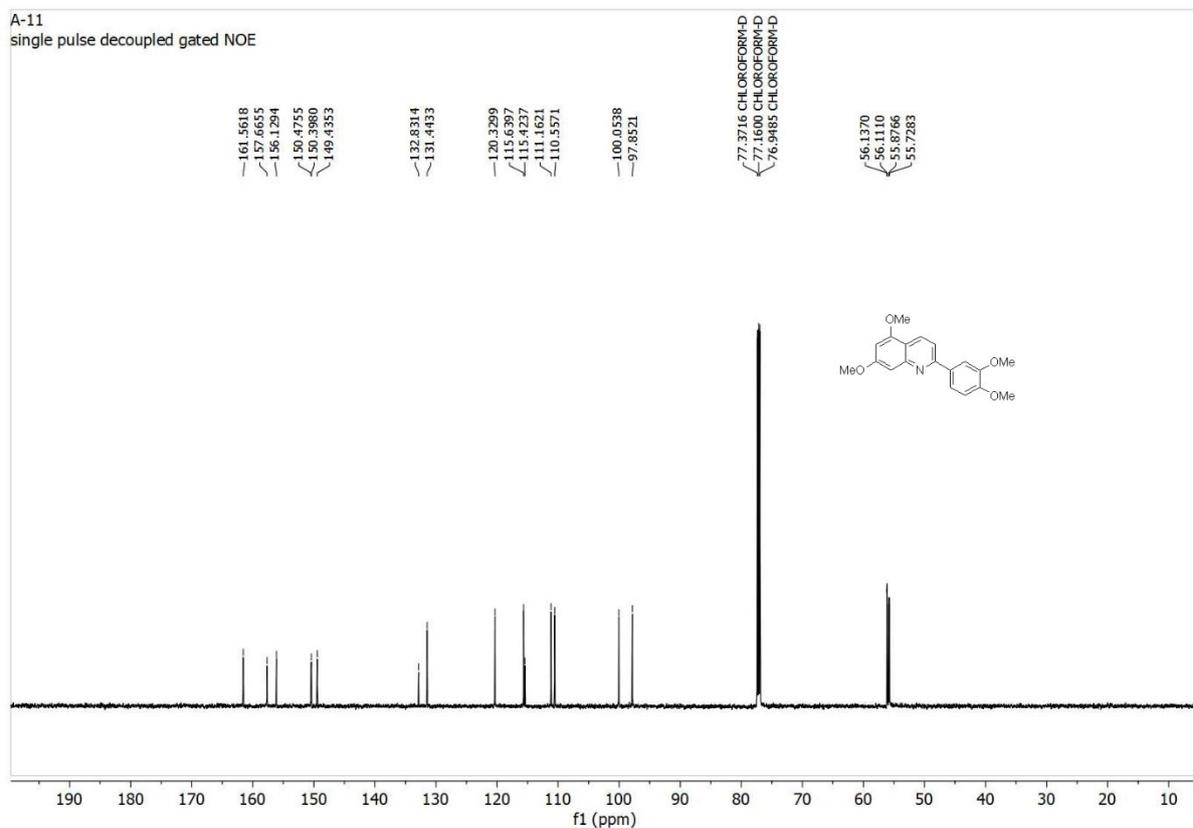


Figure S76. ^{13}C NMR Spectrum of **26d** in CDCl_3 .

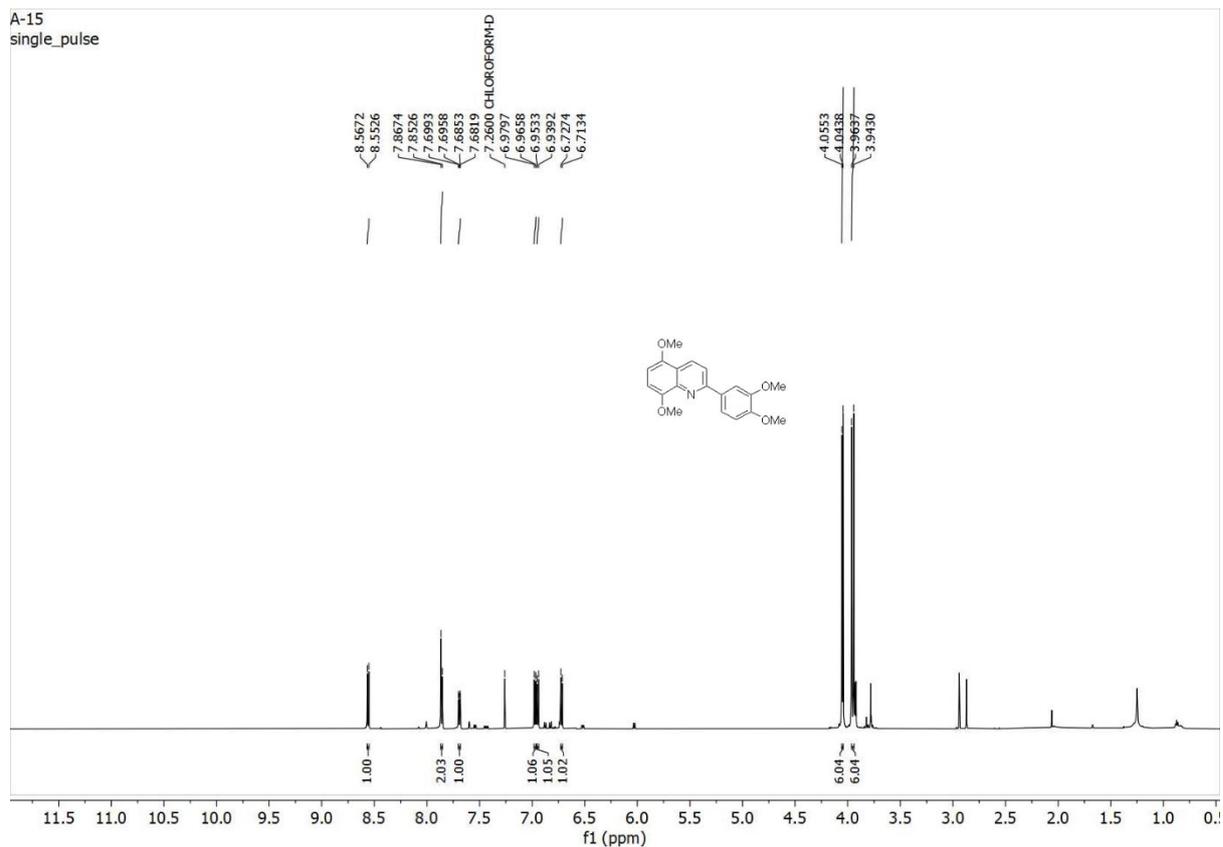


Figure S77. ^1H NMR Spectrum of **26e** in CDCl_3 .

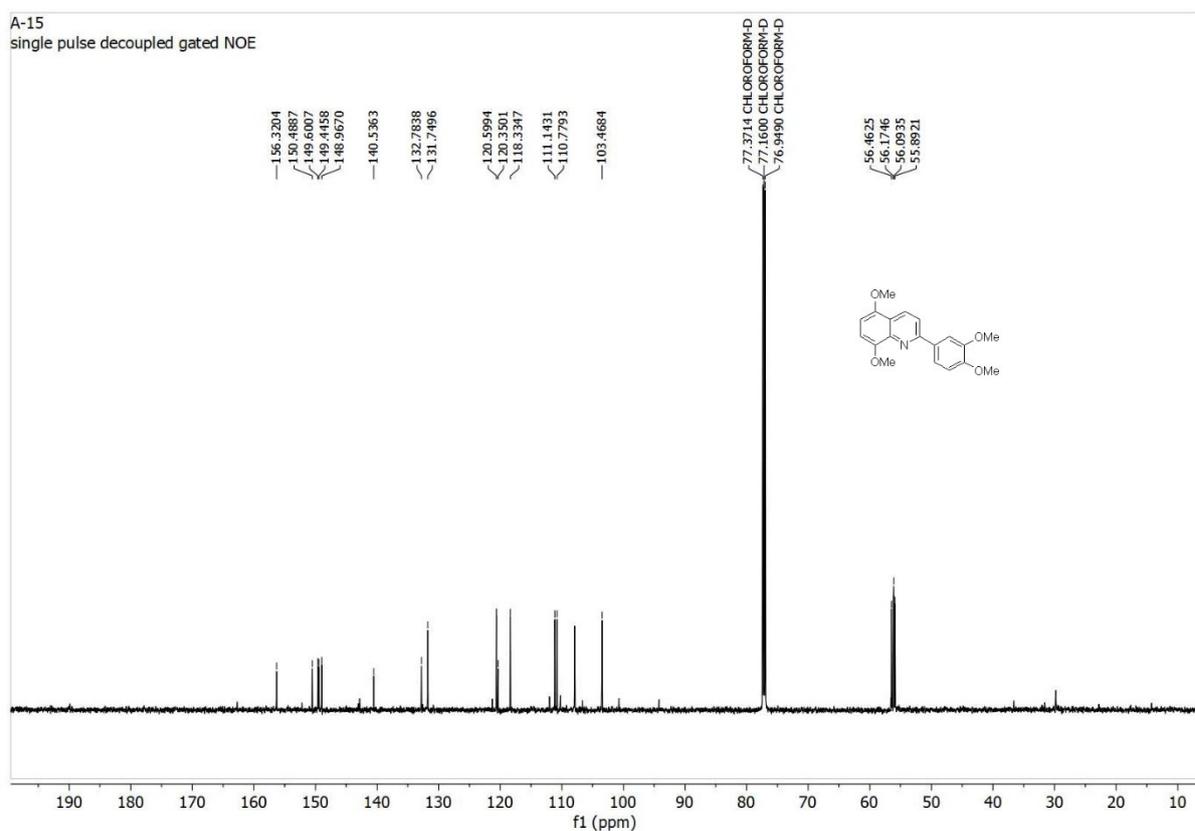


Figure S78. ^{13}C NMR Spectrum of **26e** in CDCl_3 .

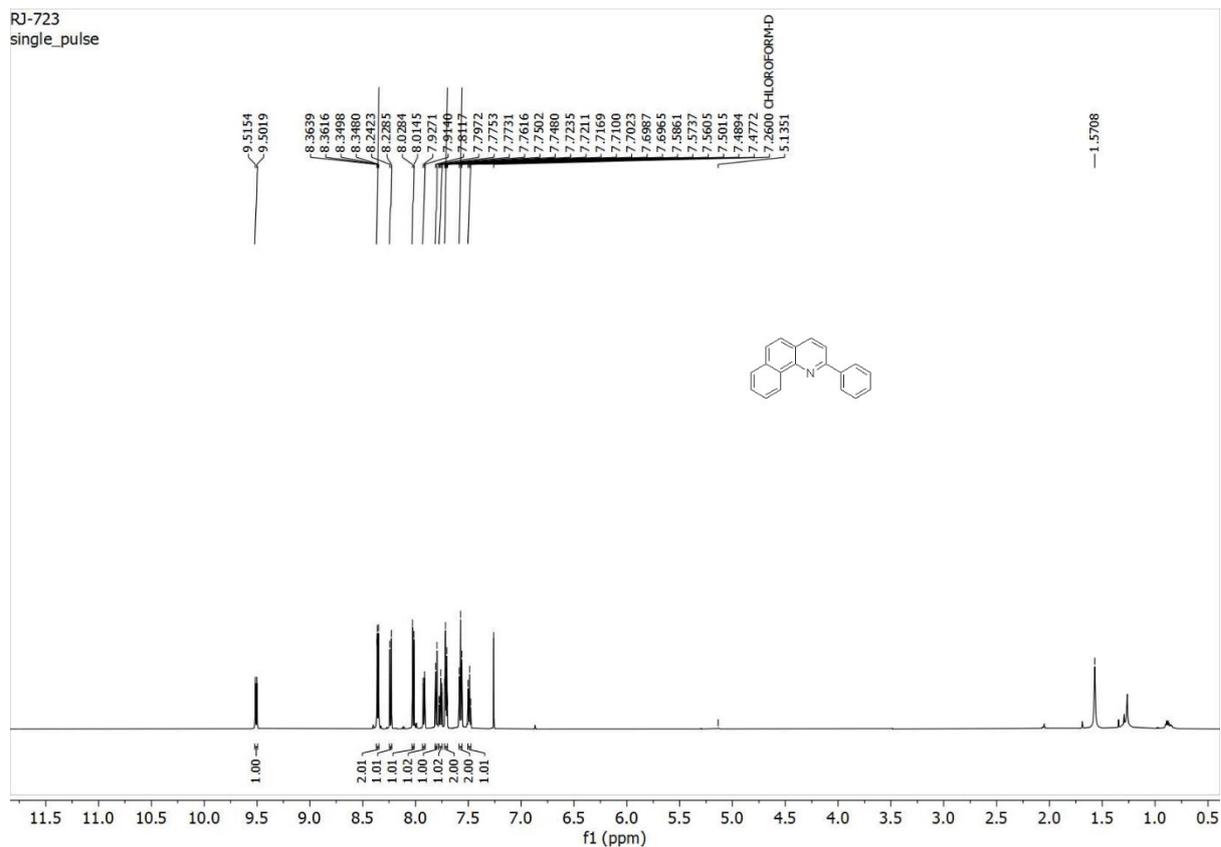


Figure S79. ^1H NMR Spectrum of **15g** in CDCl_3 .

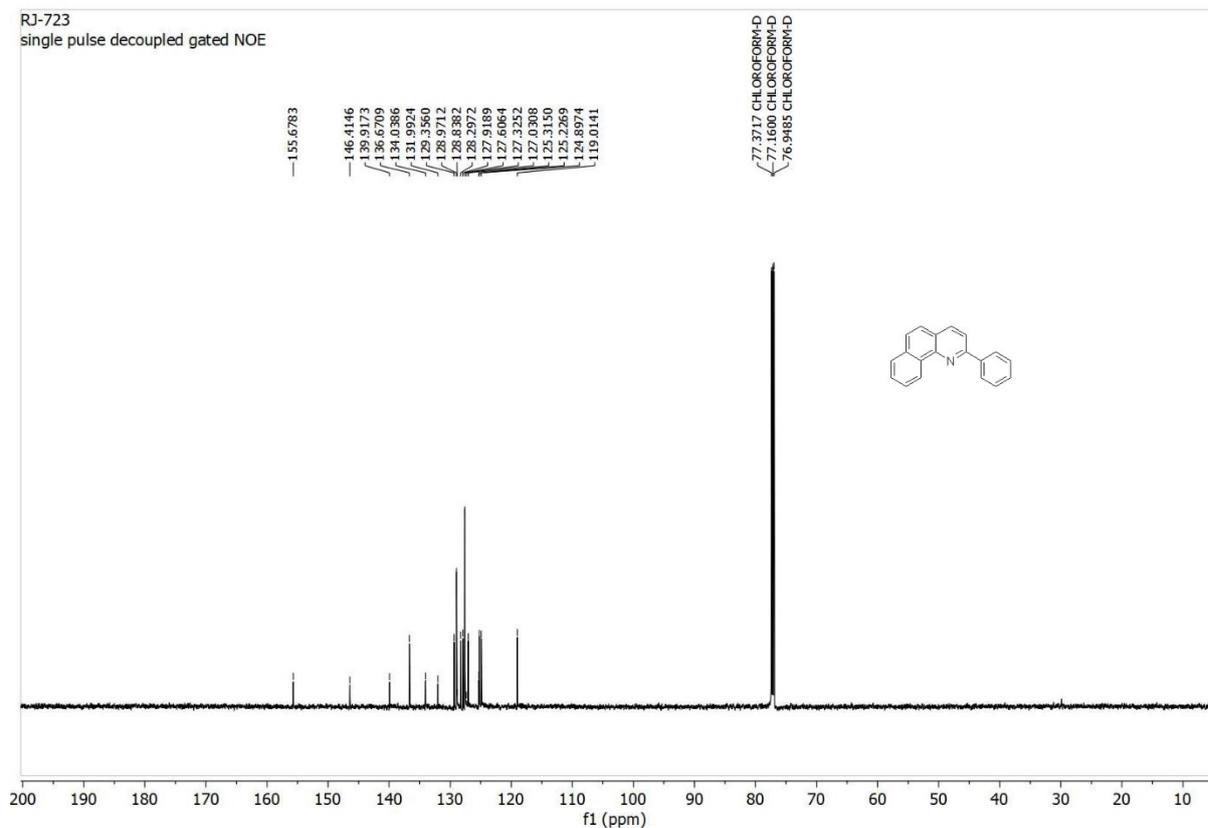


Figure S80. ^{13}C NMR Spectrum of **15g** in CDCl_3 .

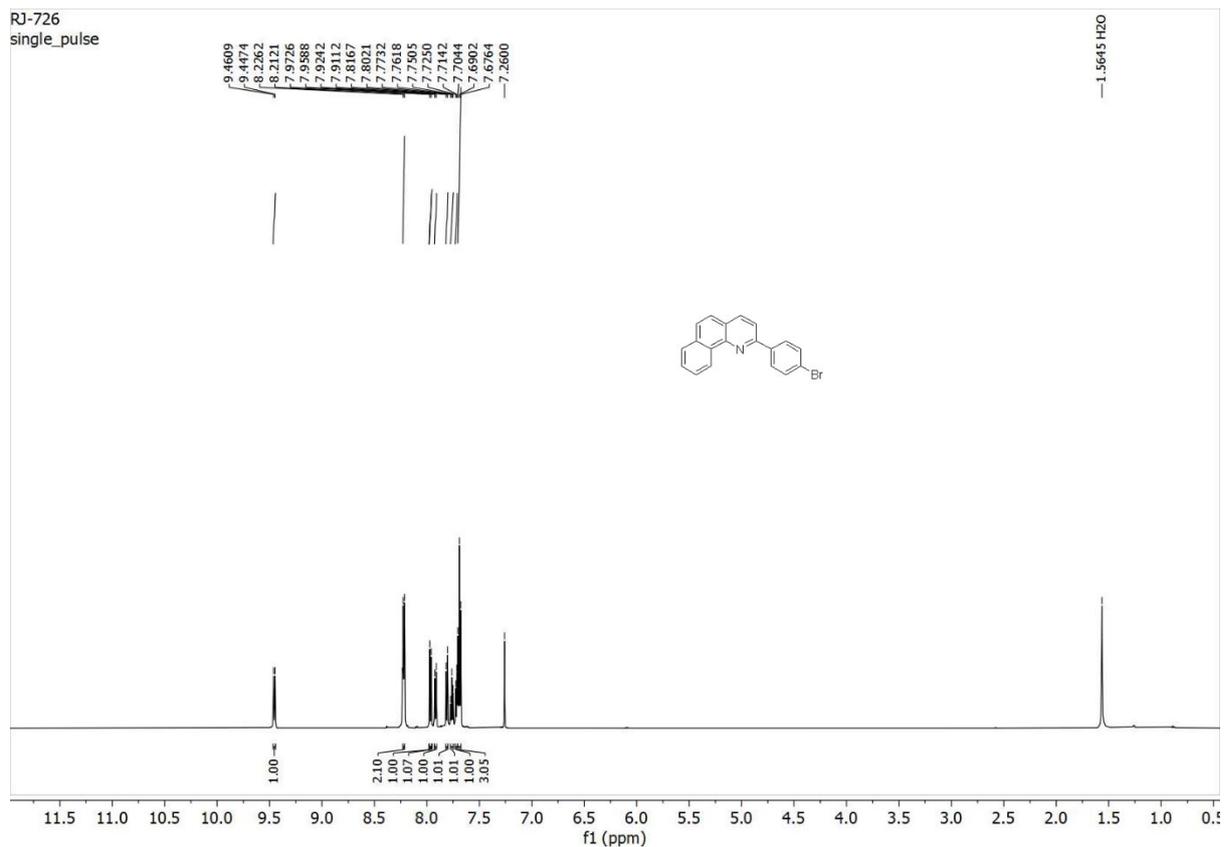


Figure S81. ^1H NMR Spectrum of **16g** in CDCl_3 .

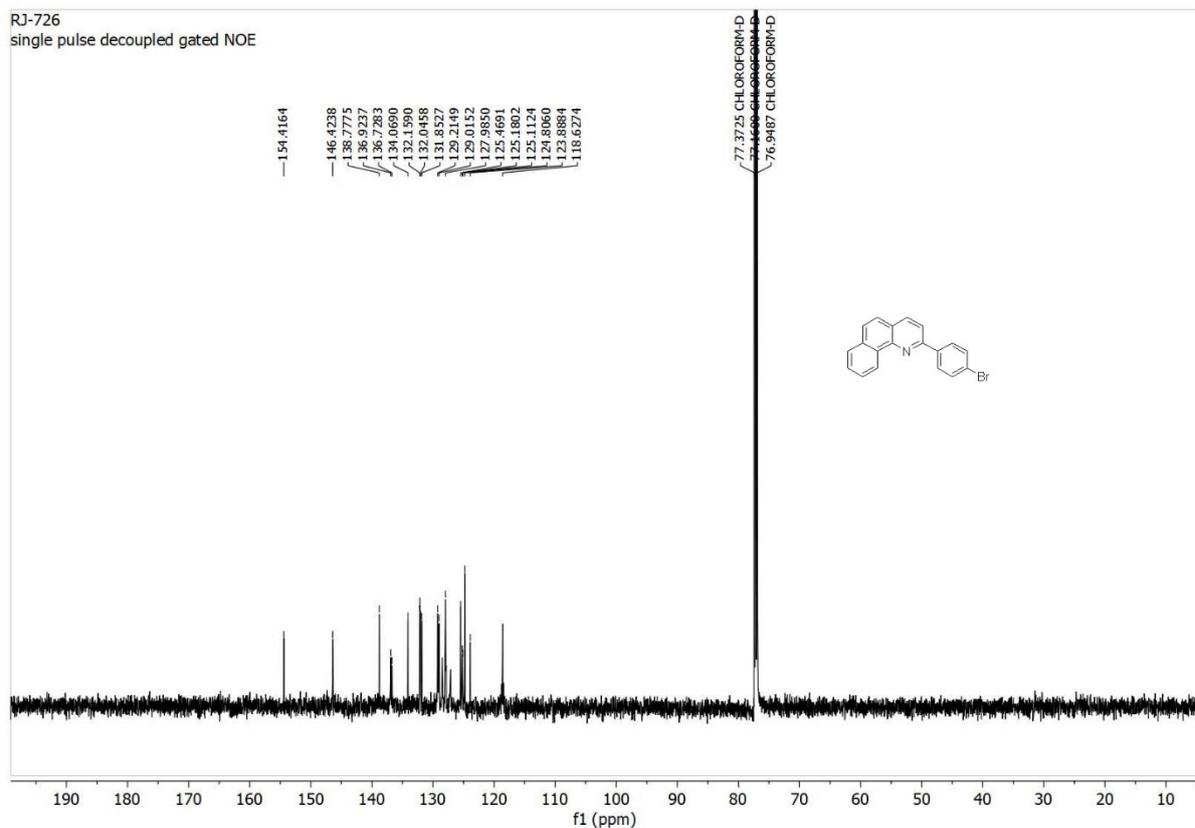


Figure S82. ^{13}C NMR Spectrum of **16g** in CDCl_3 .

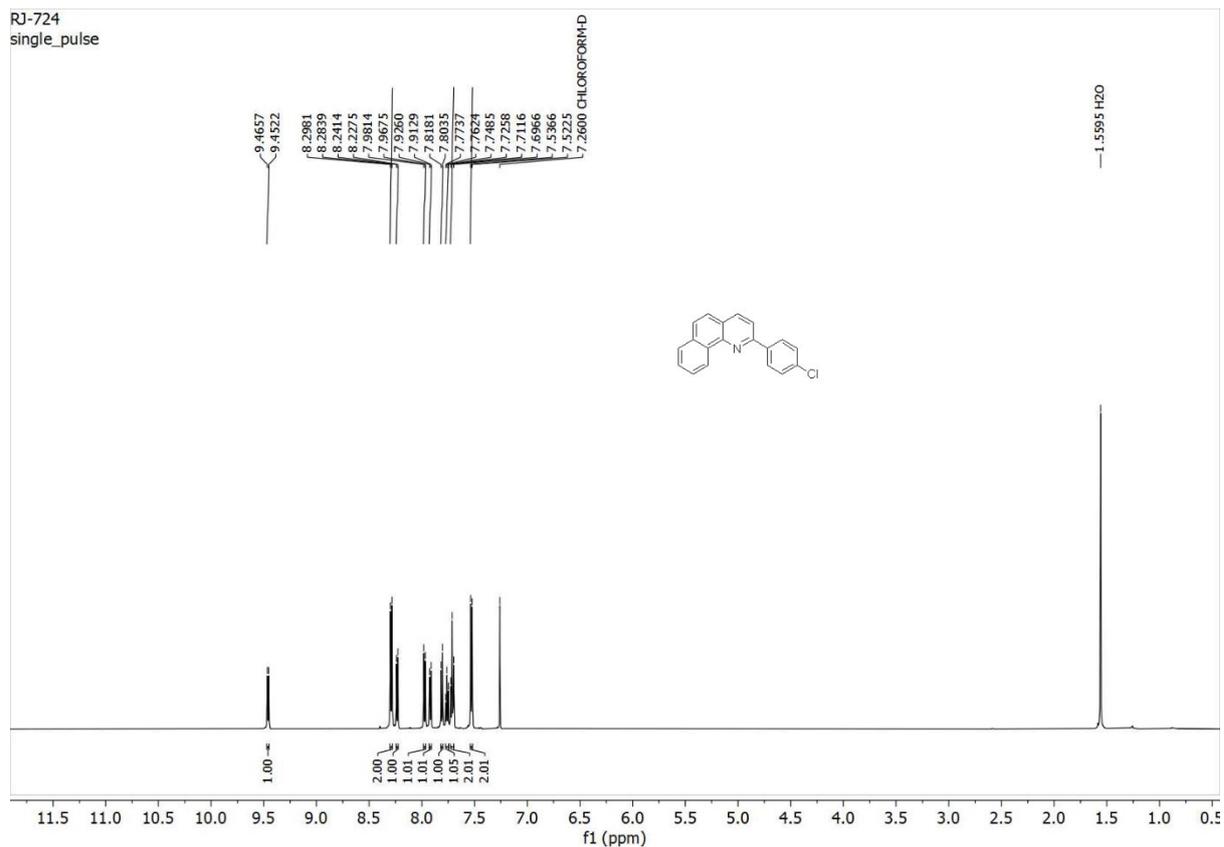


Figure S83. ¹H NMR Spectrum of **17g** in CDCl₃.

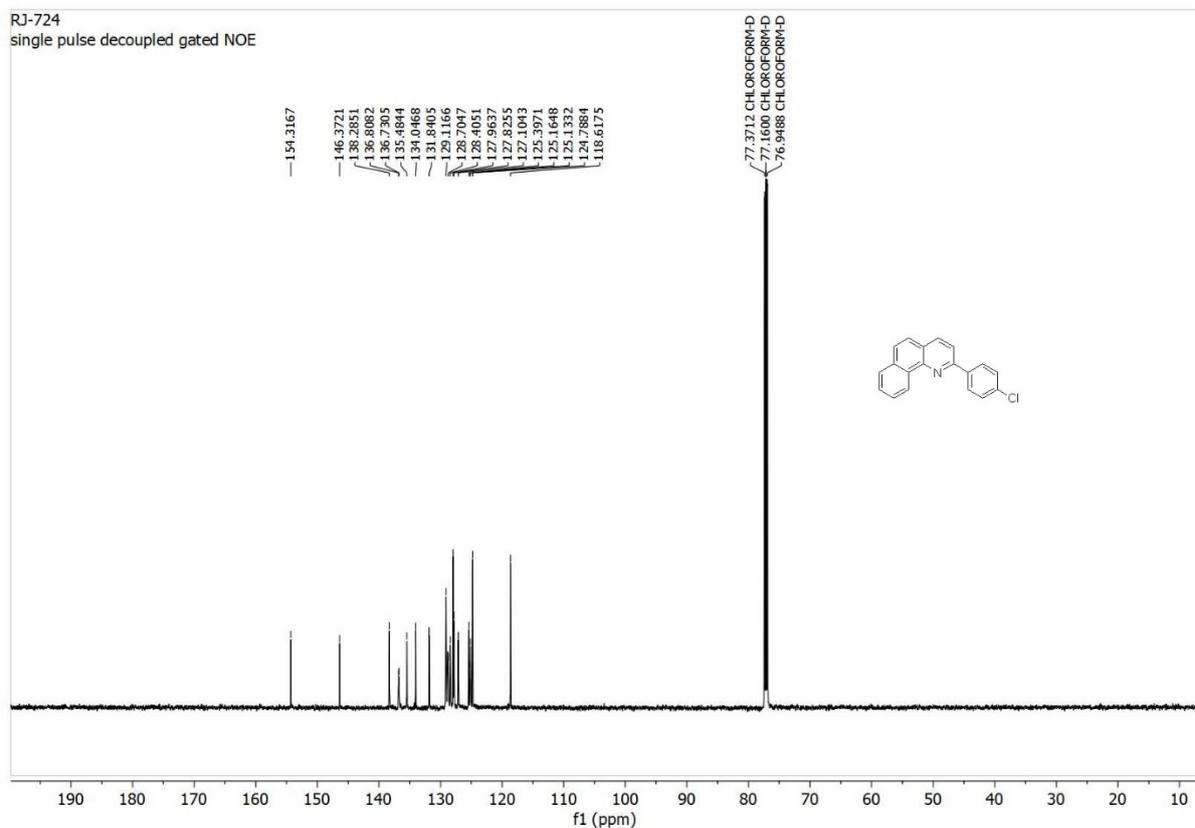


Figure S84. ¹³C NMR Spectrum of **17g** in CDCl₃.

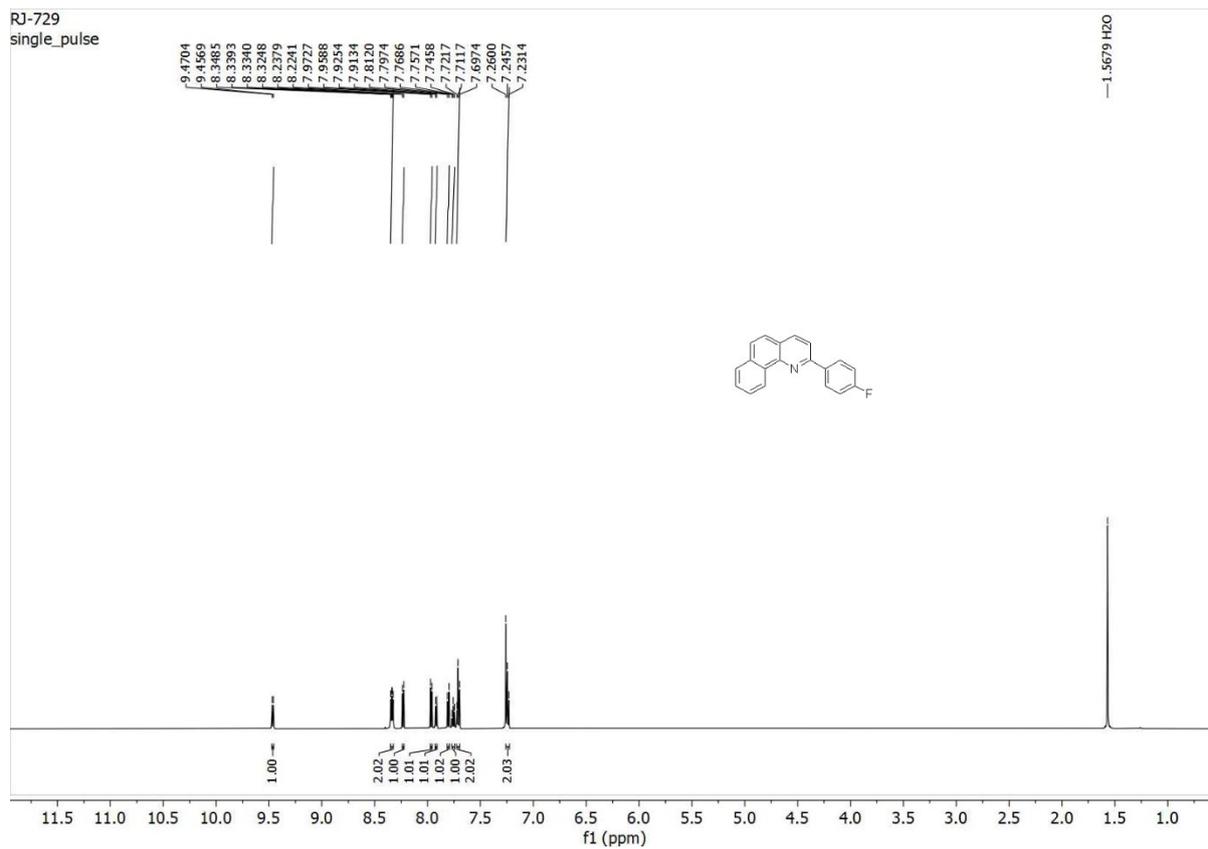


Figure S85. ^1H NMR Spectrum of **18g** in CDCl_3

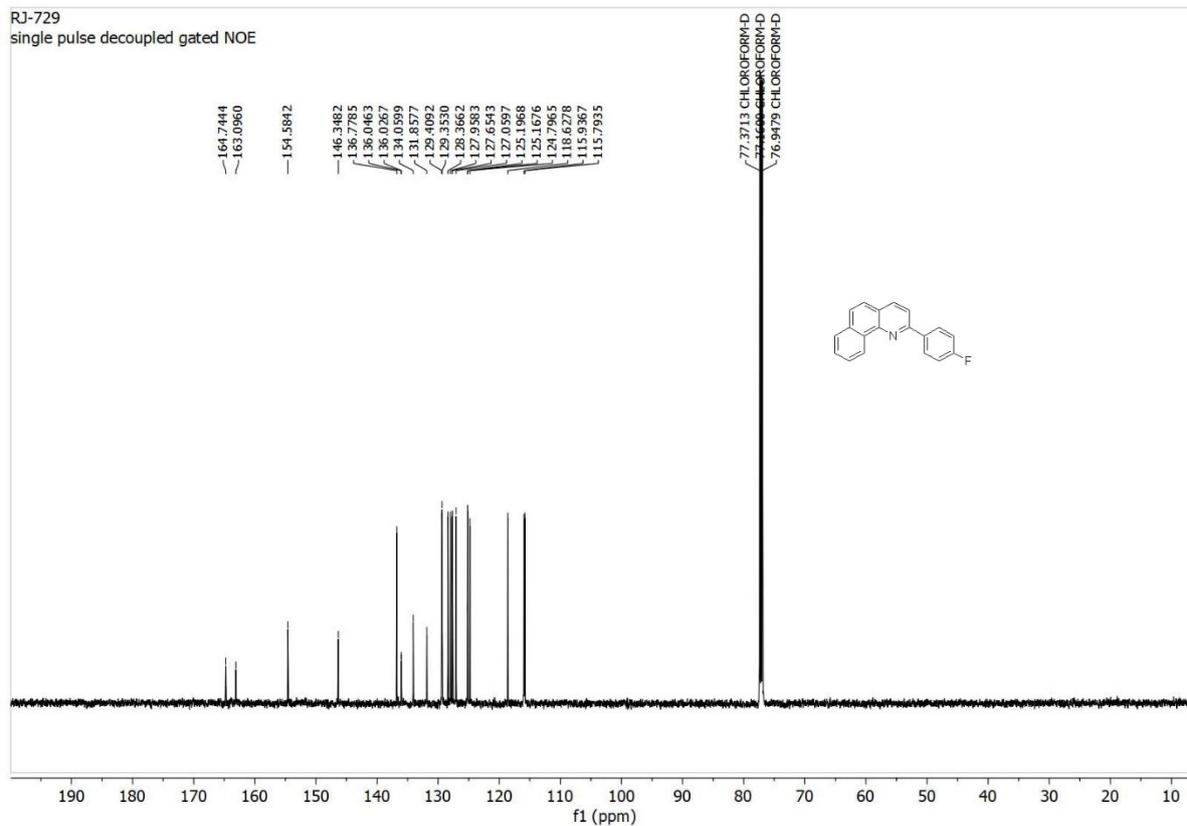


Figure S86. ^{13}C NMR Spectrum of **18g** in CDCl_3 .

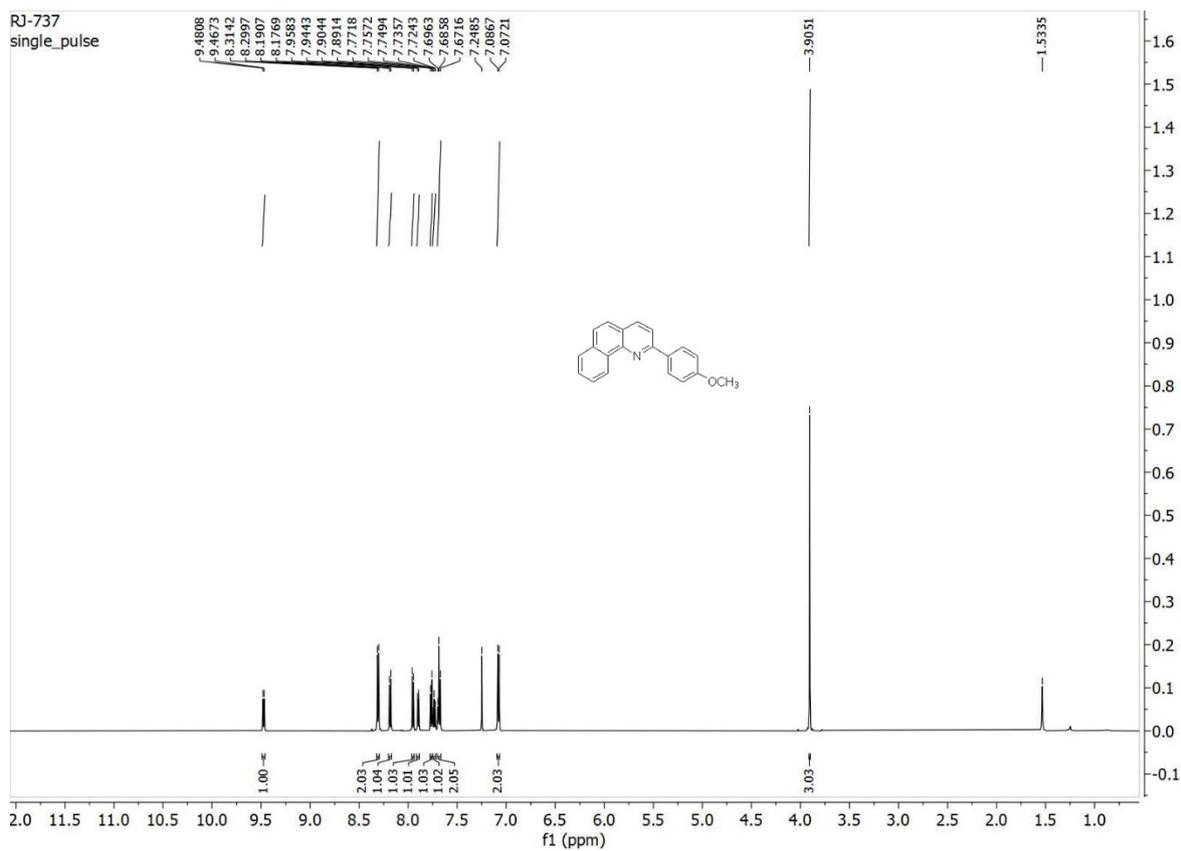


Figure S87. ^{13}C NMR Spectrum of **19g** in CDCl_3 .

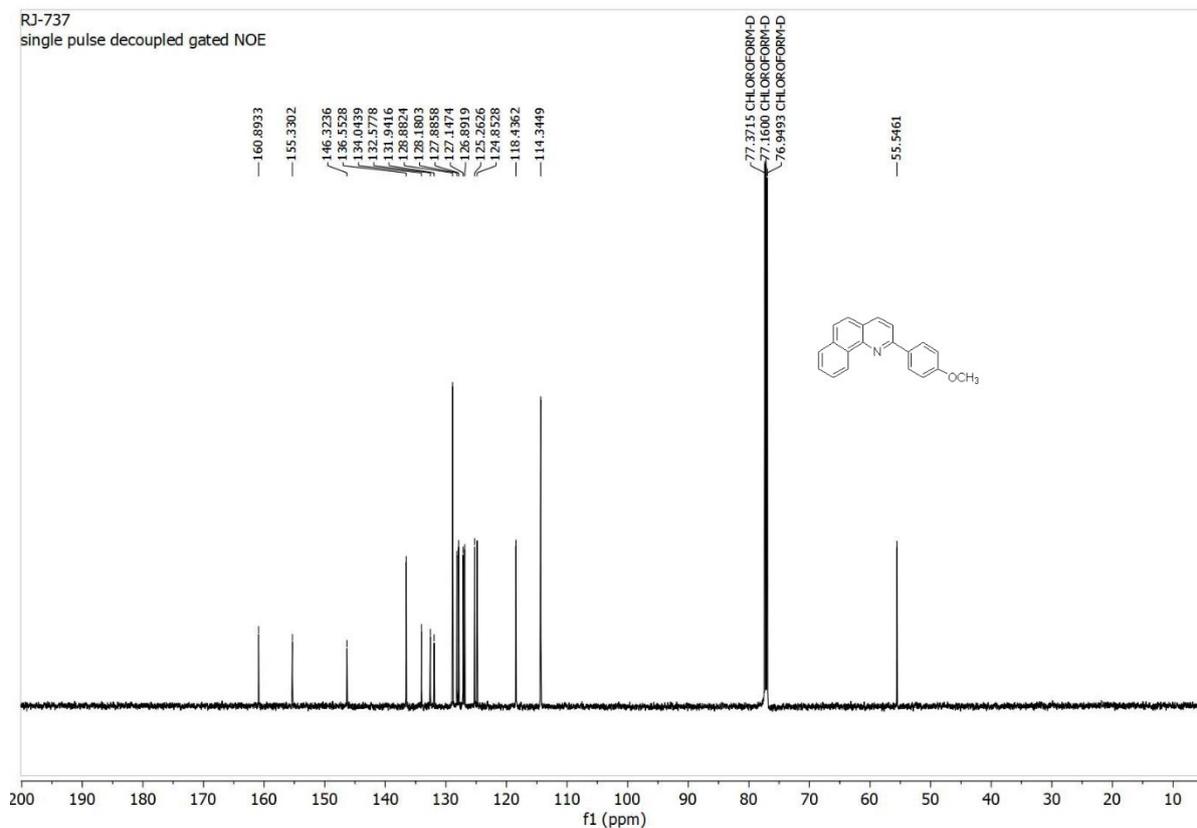


Figure S88. ^{13}C NMR Spectrum of **19g** in CDCl_3 .

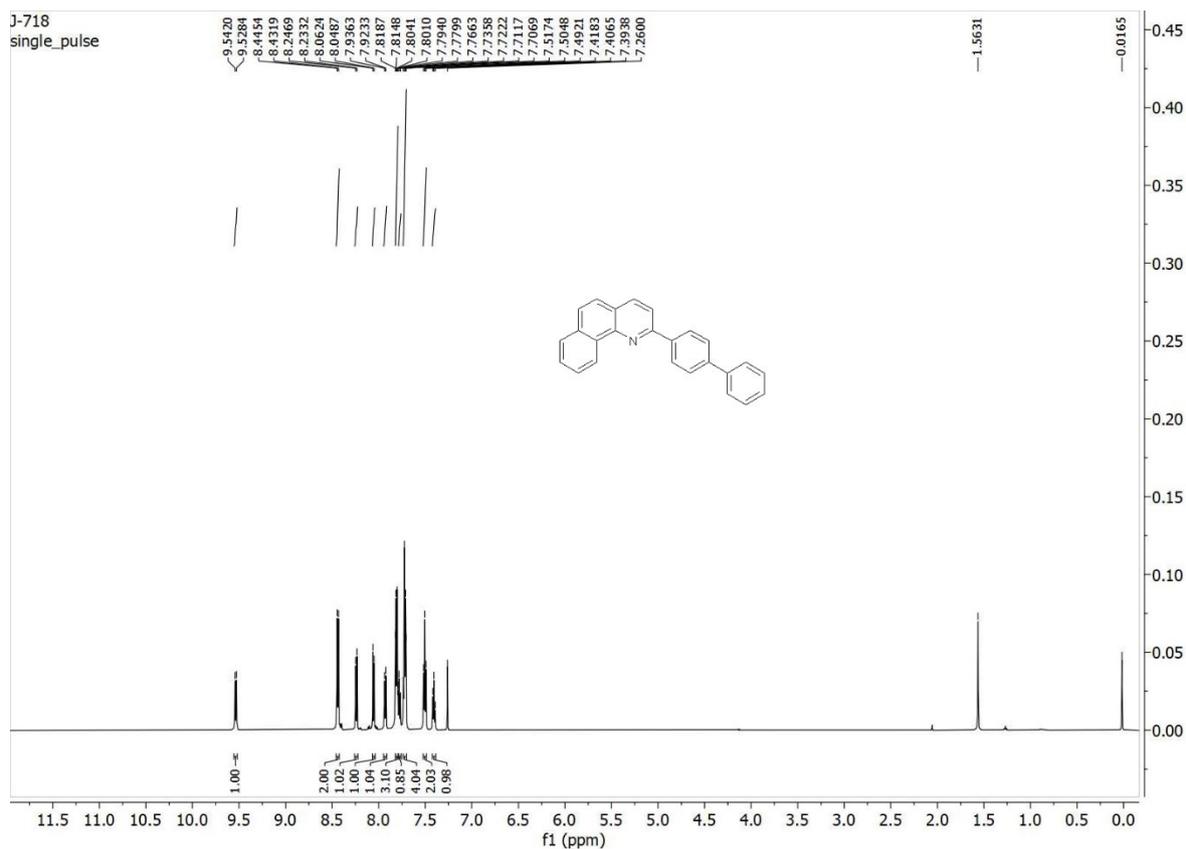


Figure S89. ¹H NMR Spectrum of **22g** in CDCl₃.

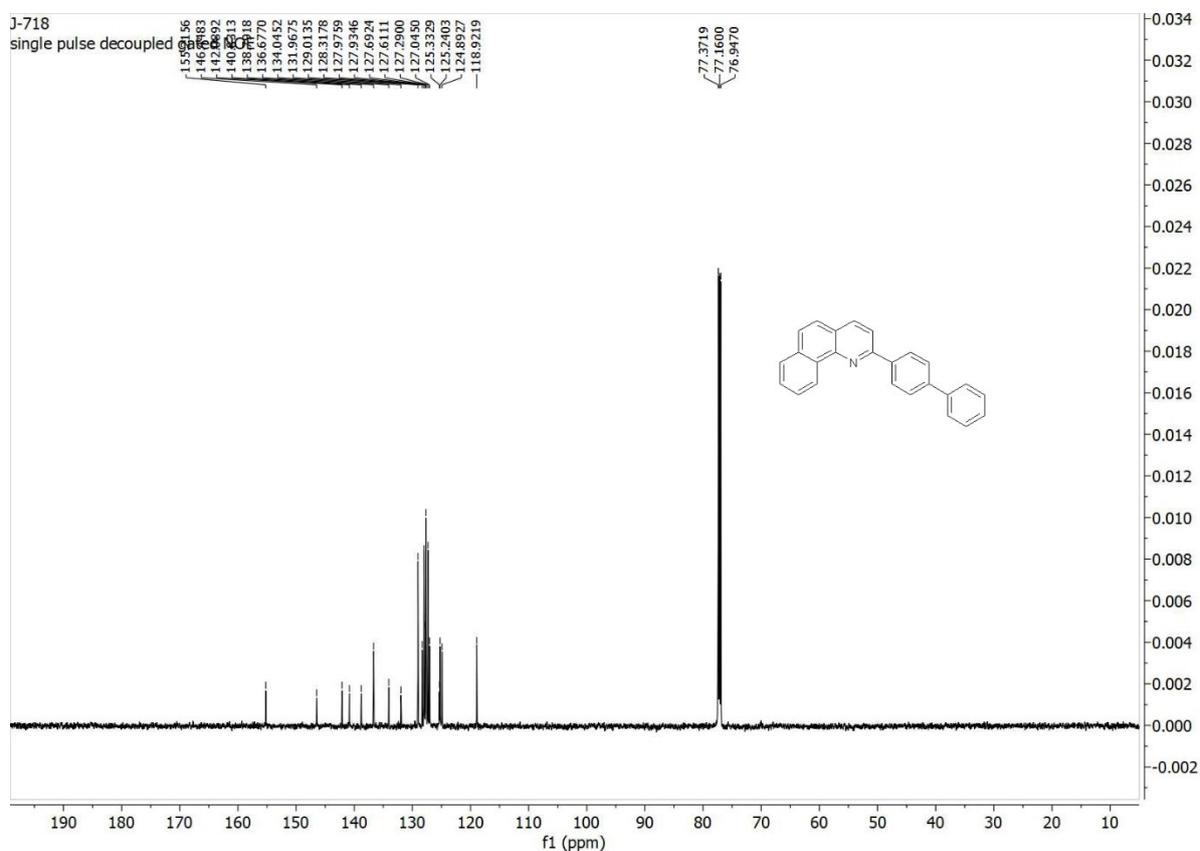


Figure S90. ¹³C NMR Spectrum of **22g** in CDCl₃.

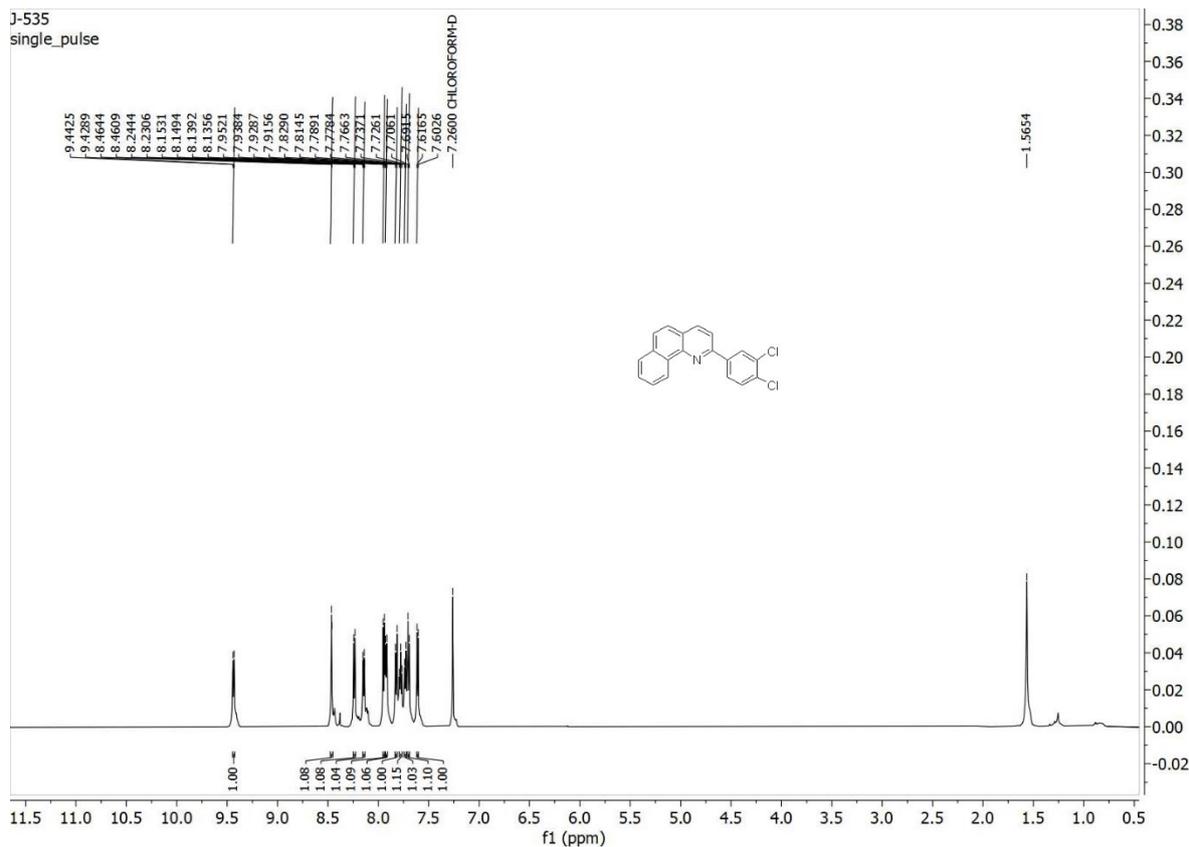


Figure S91. ¹H NMR Spectrum of 25g in CDCl₃.

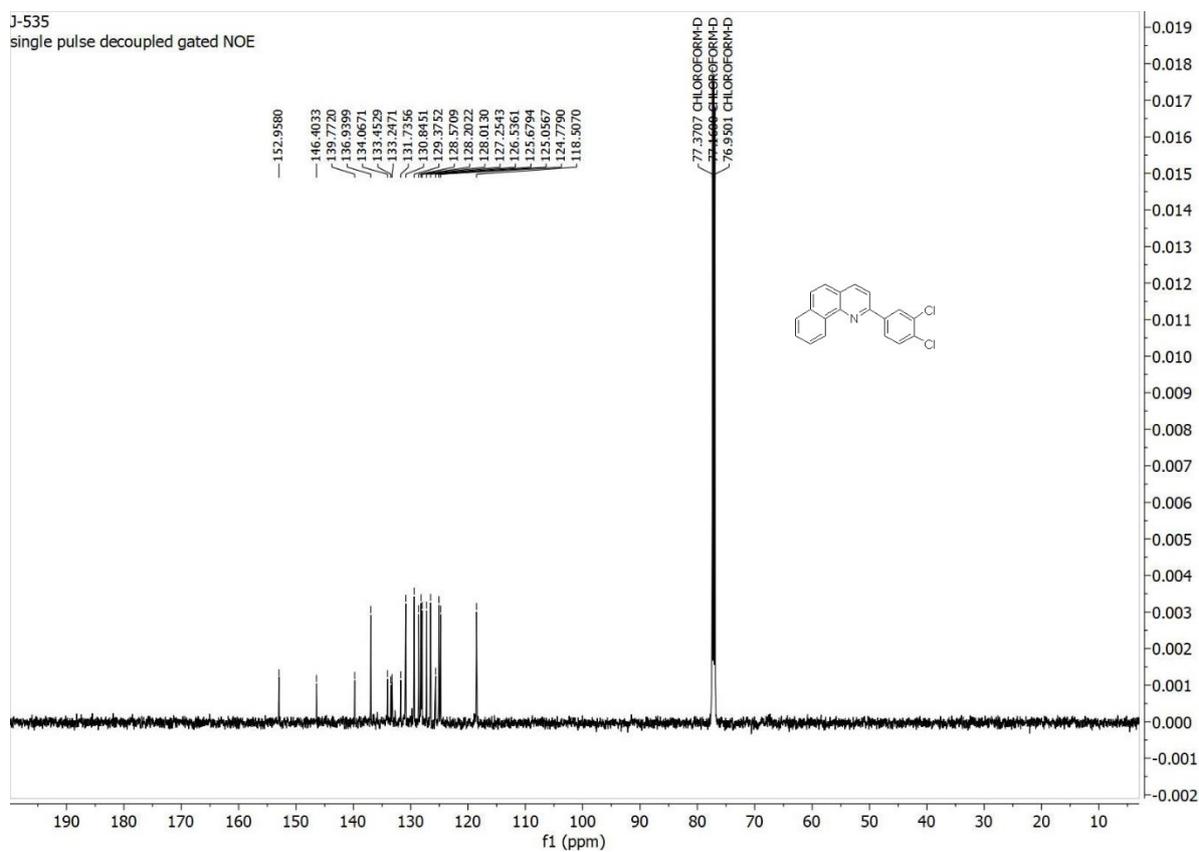


Figure S92. ¹³C NMR Spectrum of 25g in CDCl₃.

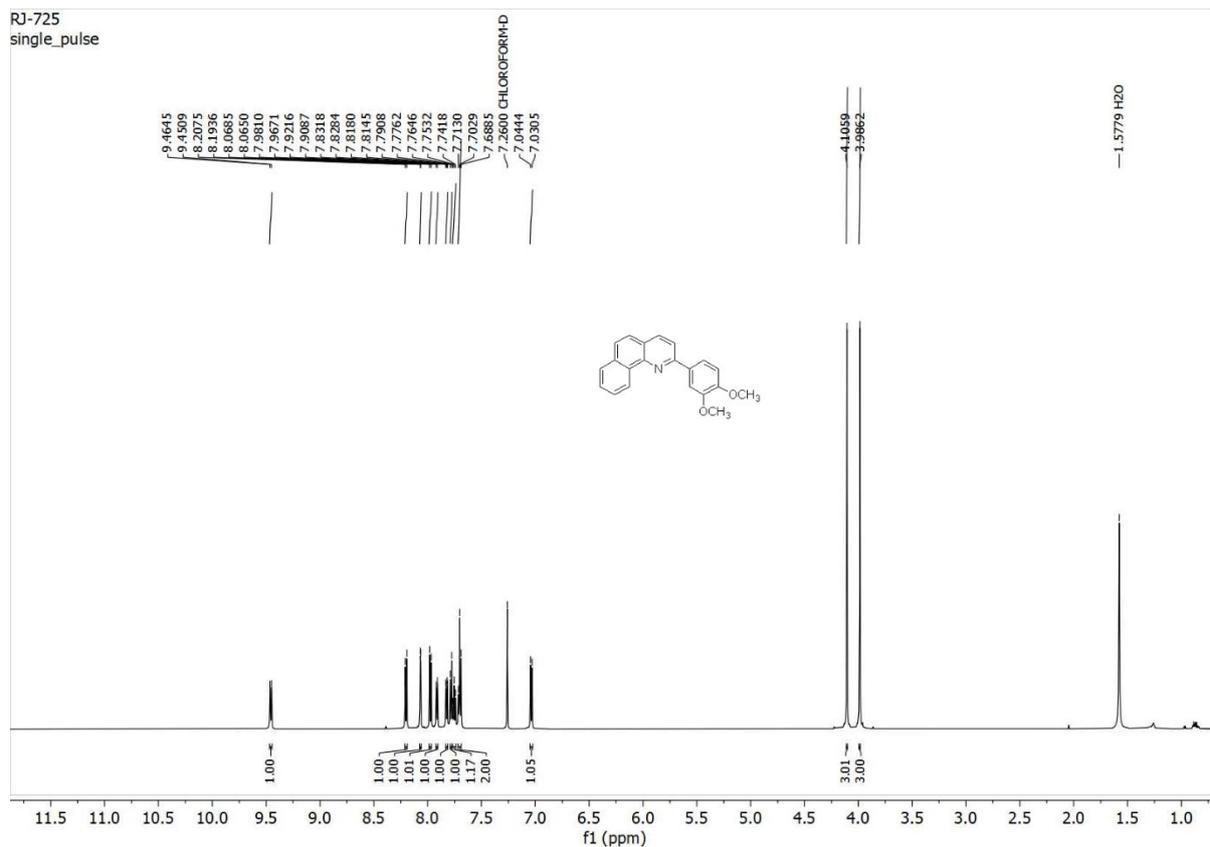


Figure S93. ¹H NMR Spectrum of 26g in CDCl₃.

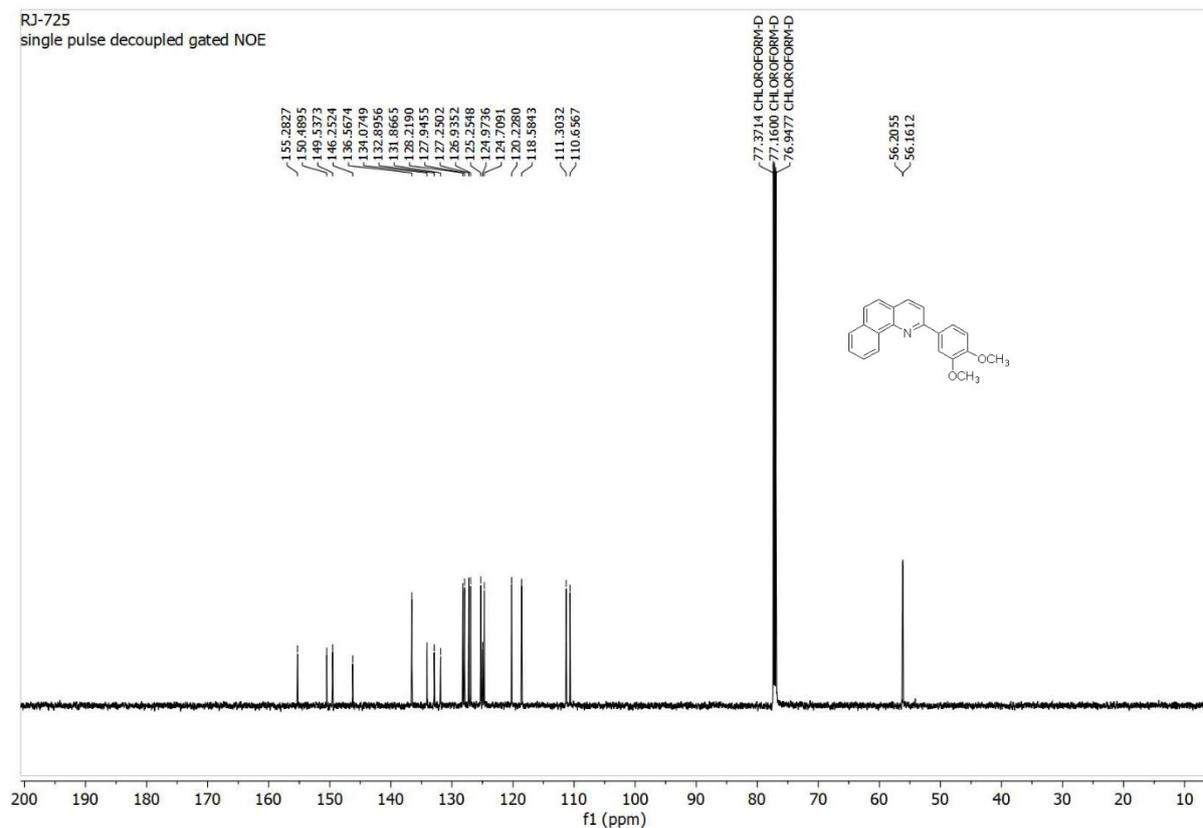


Figure S94. ¹³C NMR Spectrum of 26g in CDCl₃.

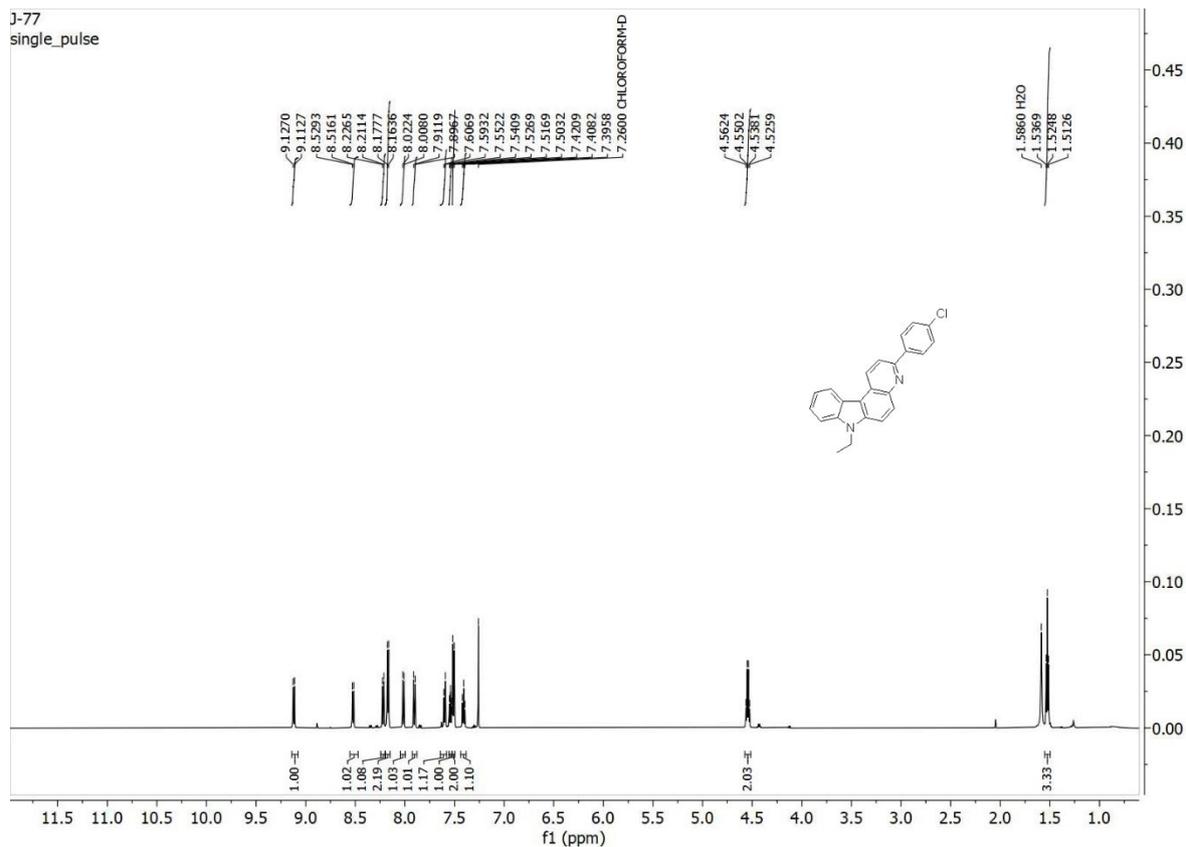


Figure S95. ^1H NMR Spectrum of **17h** in CDCl_3 .

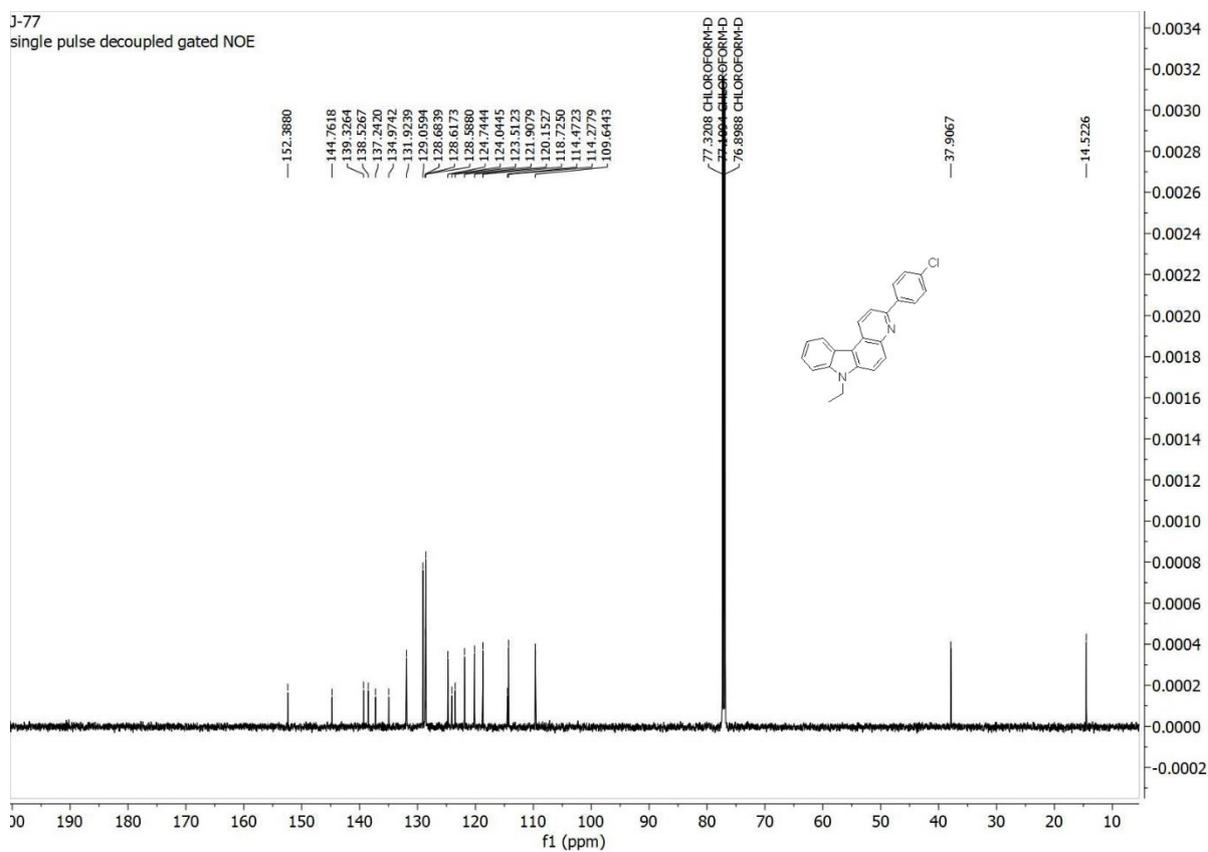


Figure S96. ^{13}C NMR Spectrum of **17h** in CDCl_3 .

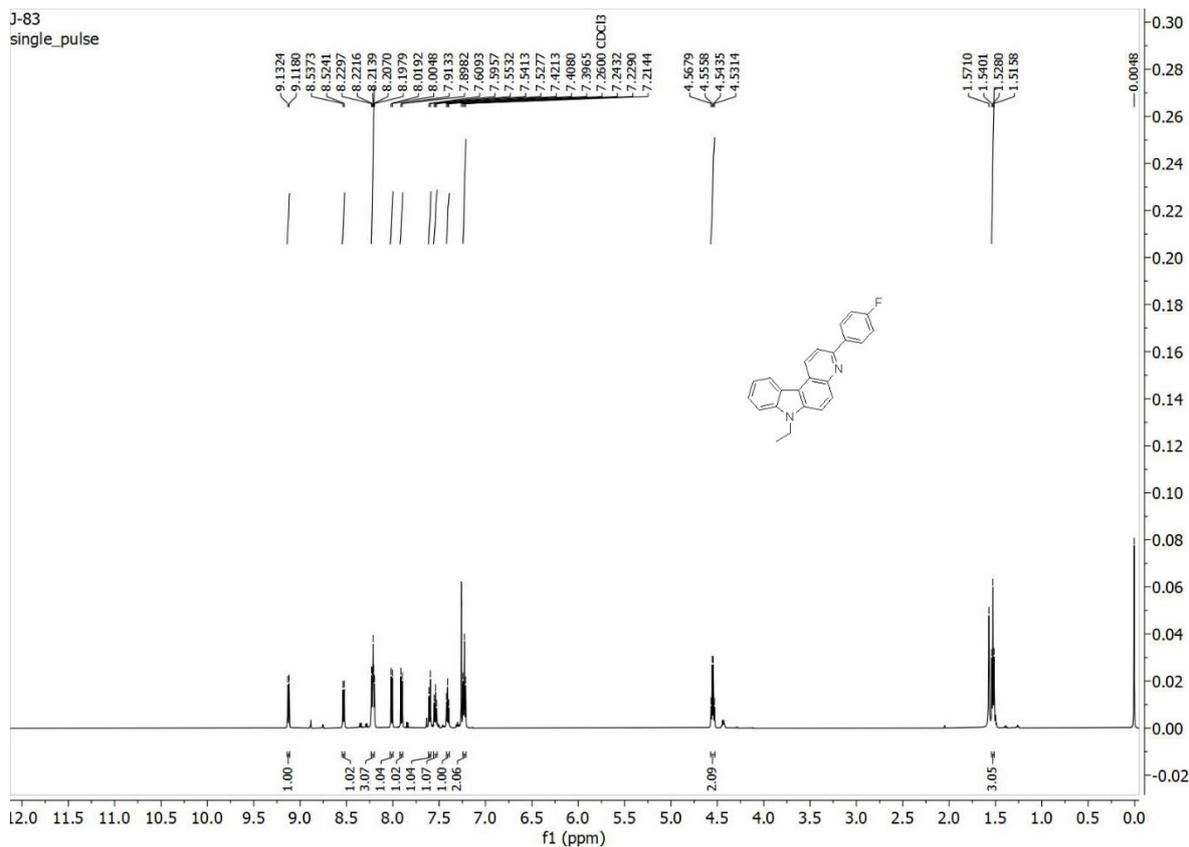


Figure S97. ^1H NMR Spectrum of **18h** in CDCl_3 .

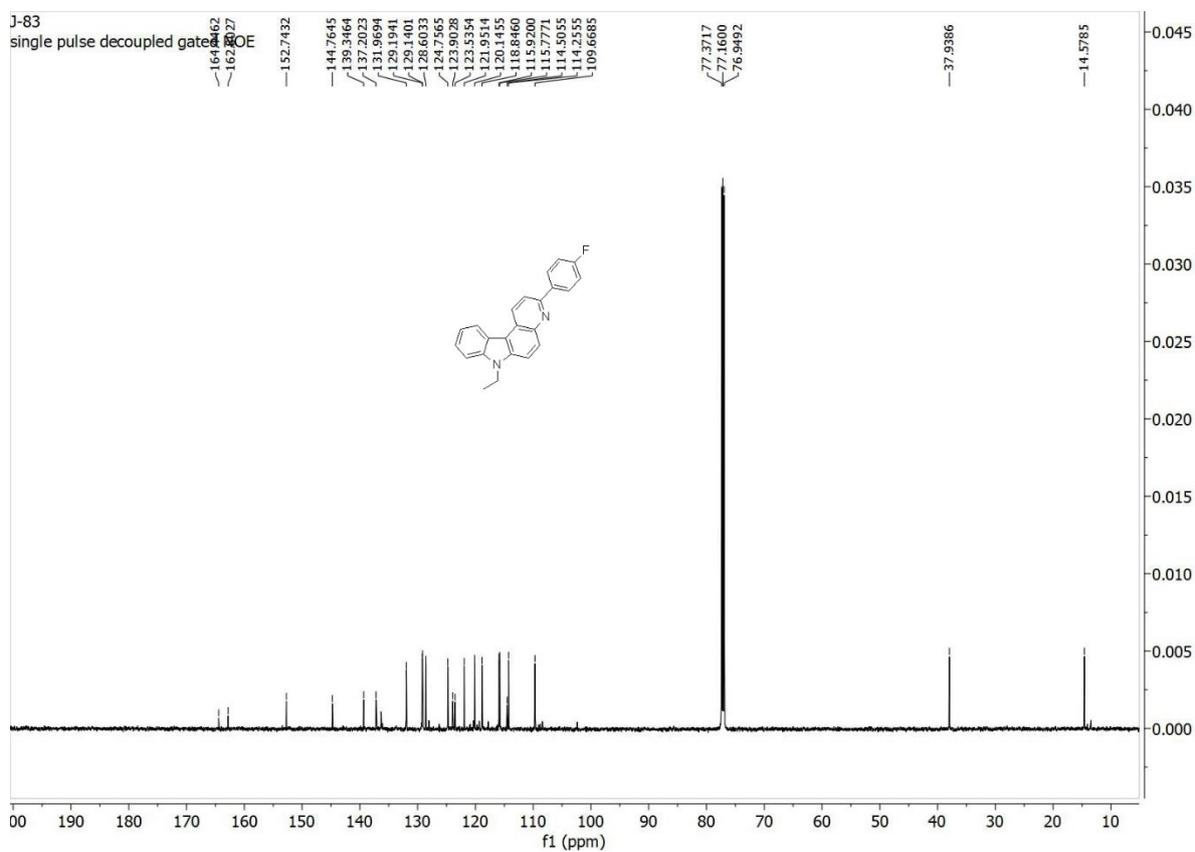


Figure S98. ^{13}C NMR Spectrum of **18h** in CDCl_3 .

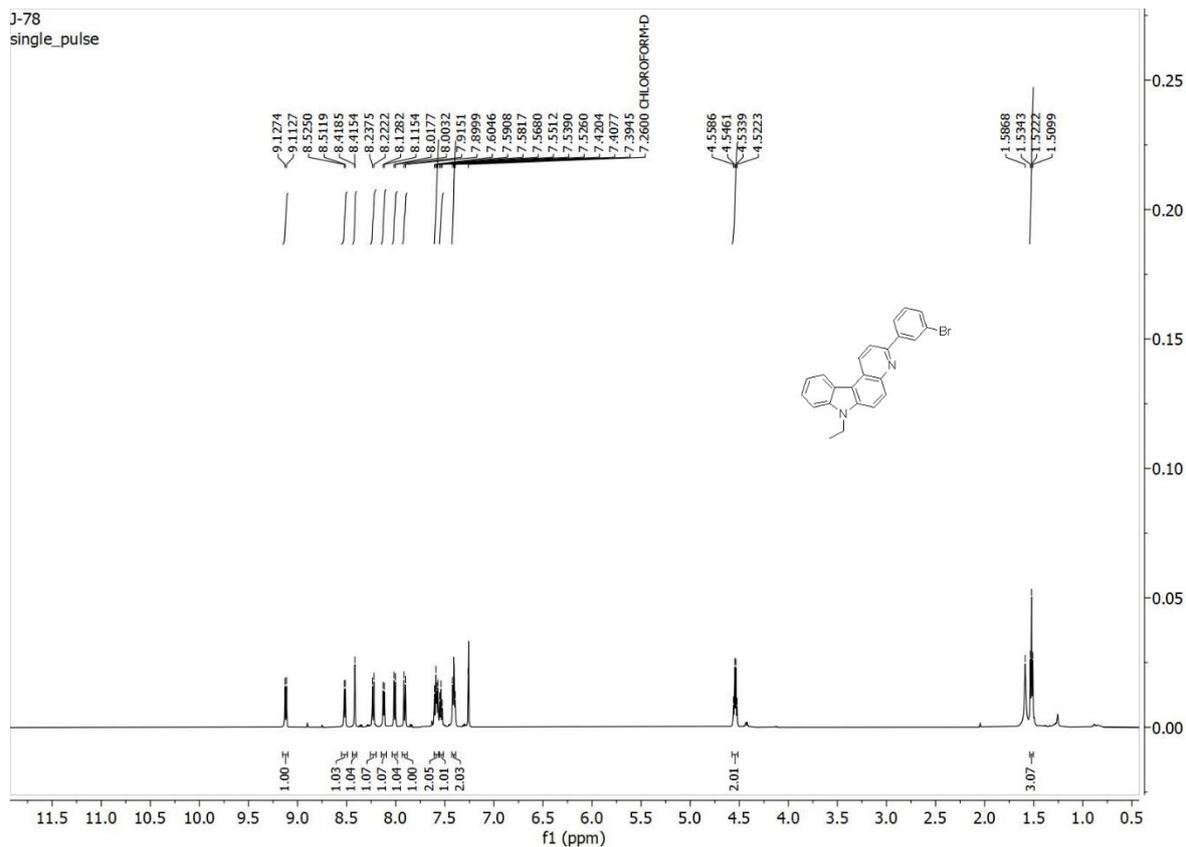


Figure S99. ^1H NMR Spectrum of **24h** in CDCl_3 .

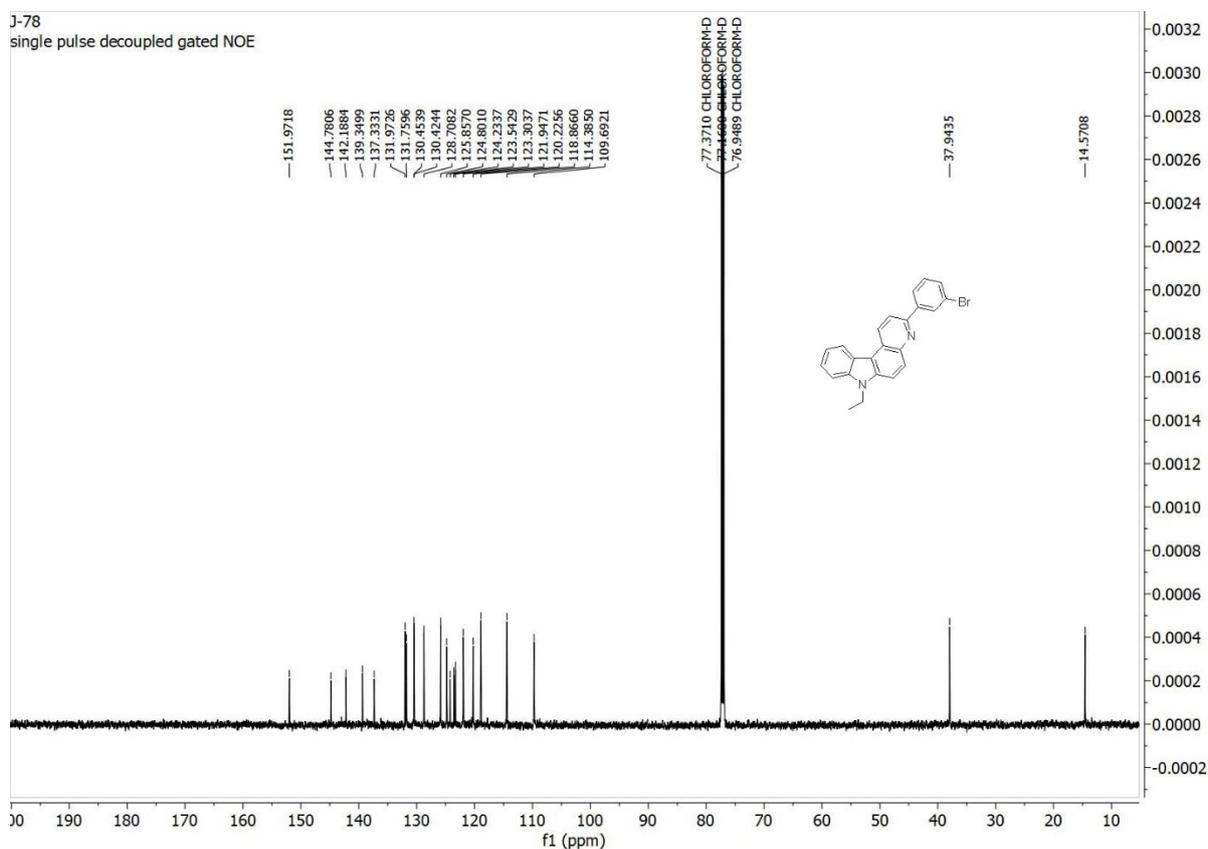


Figure S100. ^{13}C NMR Spectrum of **24h** in CDCl_3 .

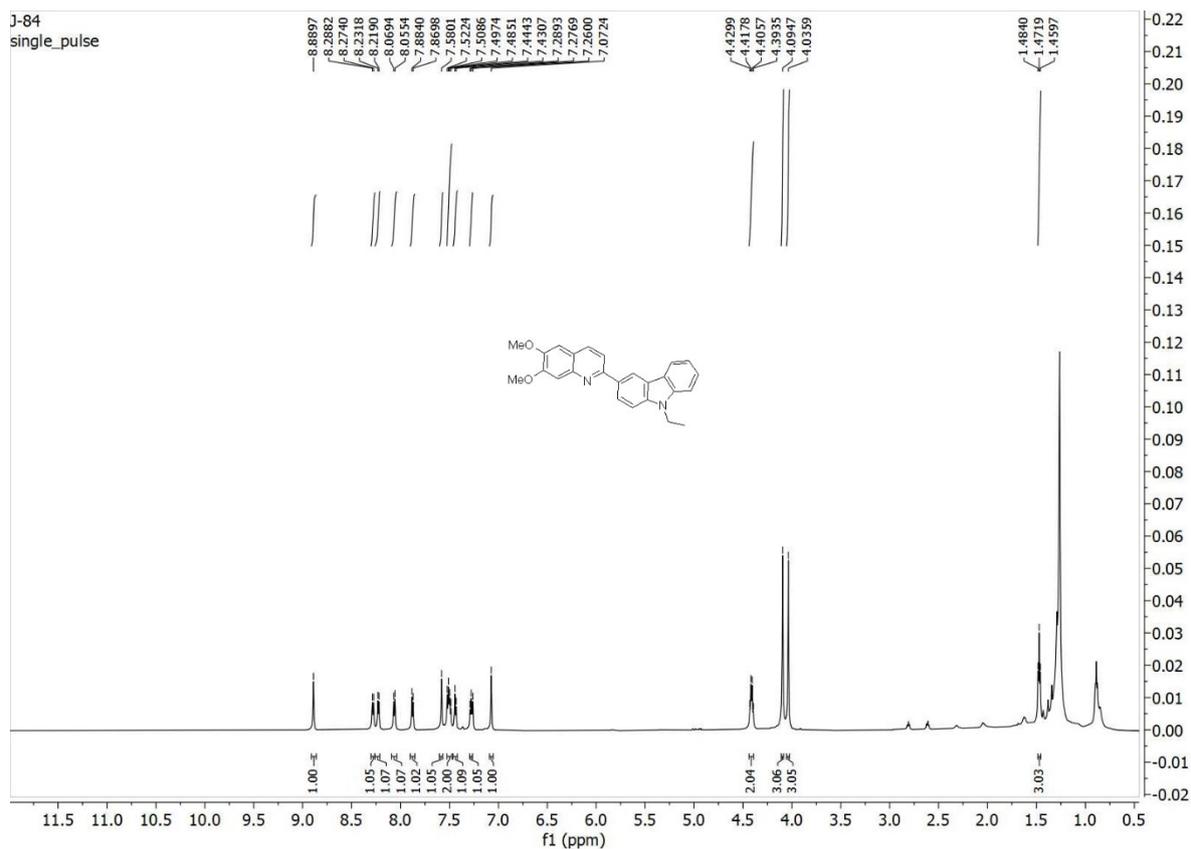


Figure S101. ^1H NMR Spectrum of **28a** in CDCl_3 .

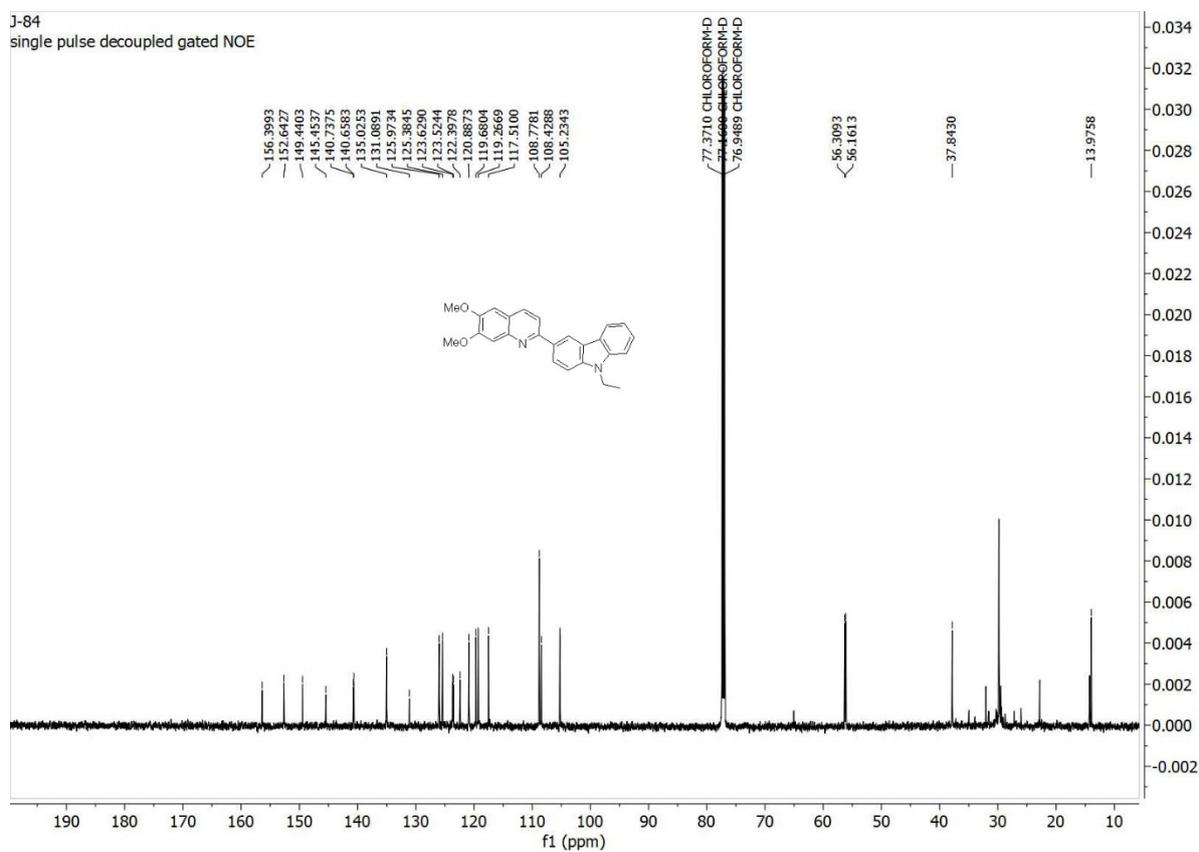


Figure S102. ^{13}C NMR Spectrum of **28a** in CDCl_3 .

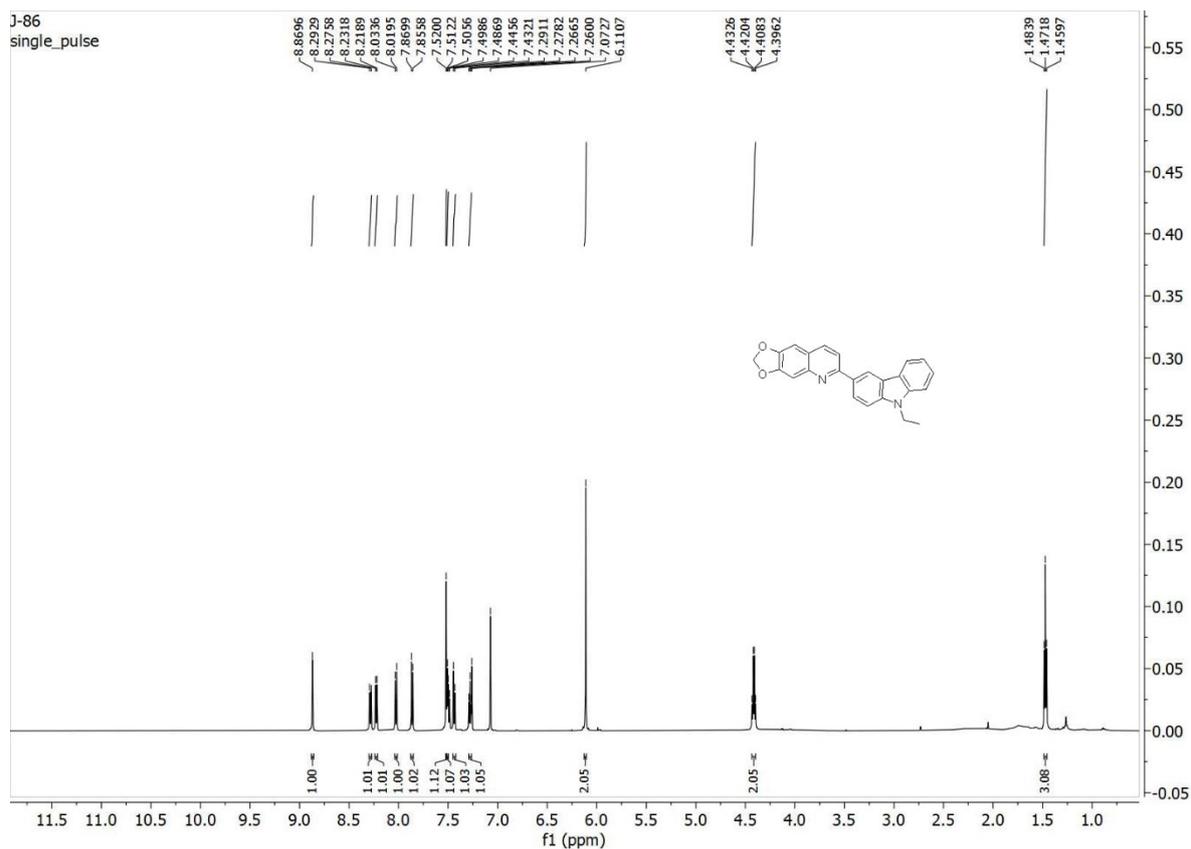


Figure S103. ^1H NMR Spectrum of **28b** in CDCl_3 .

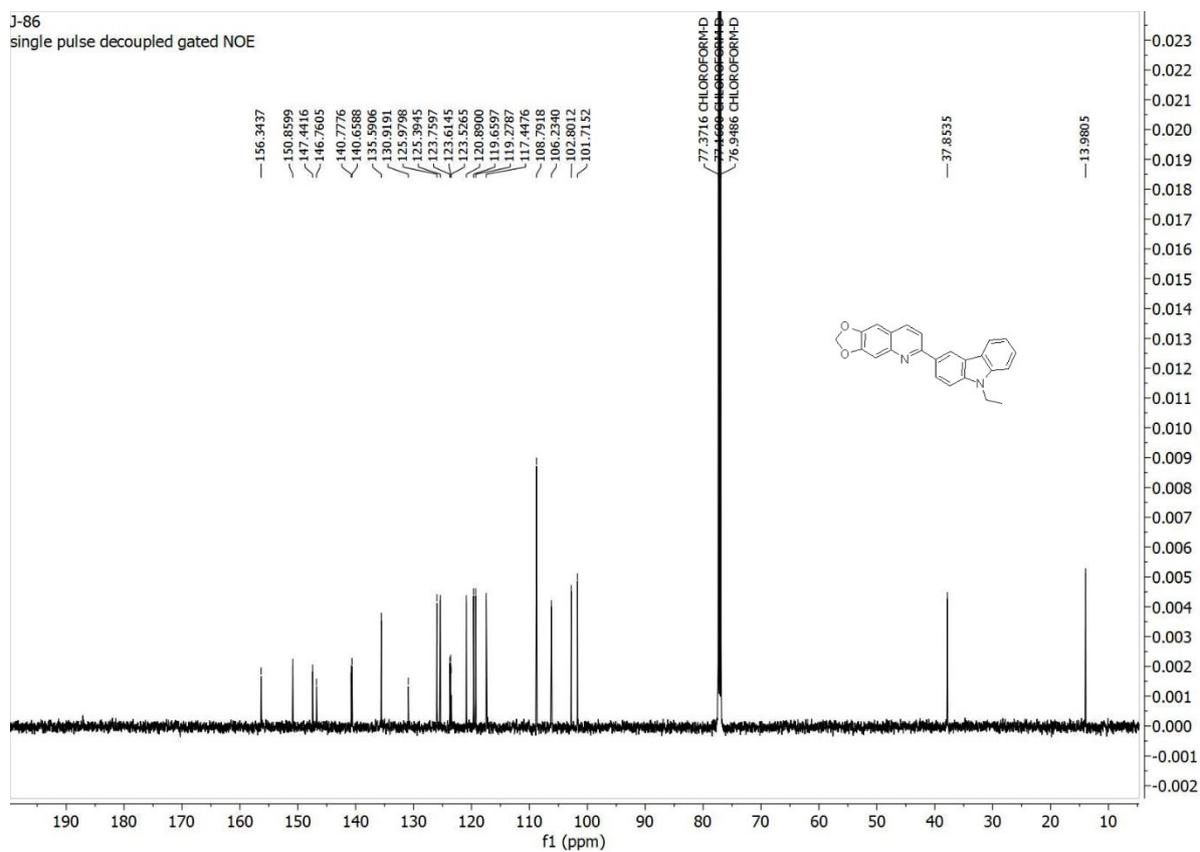


Figure S104. ^{13}C NMR Spectrum of **28b** in CDCl_3 .

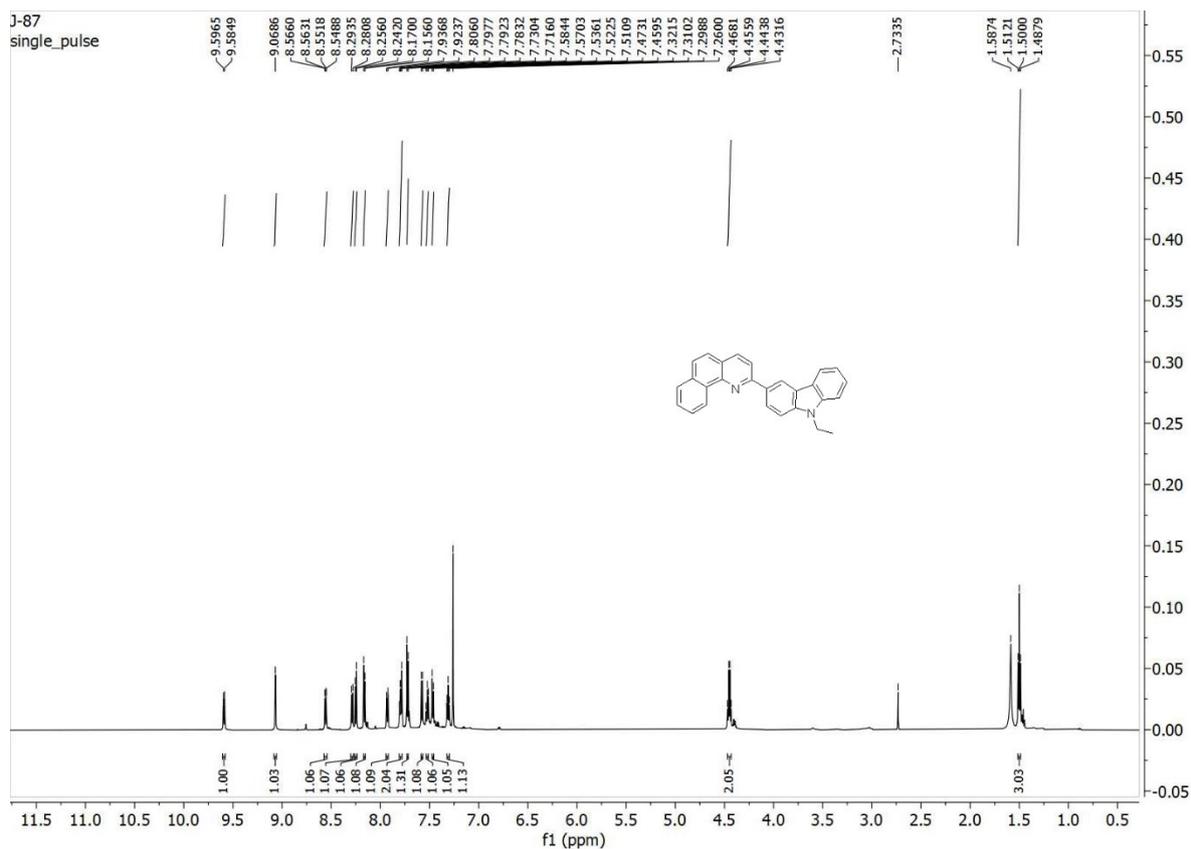


Figure S105. ^1H NMR Spectrum of **28g** in CDCl_3 .

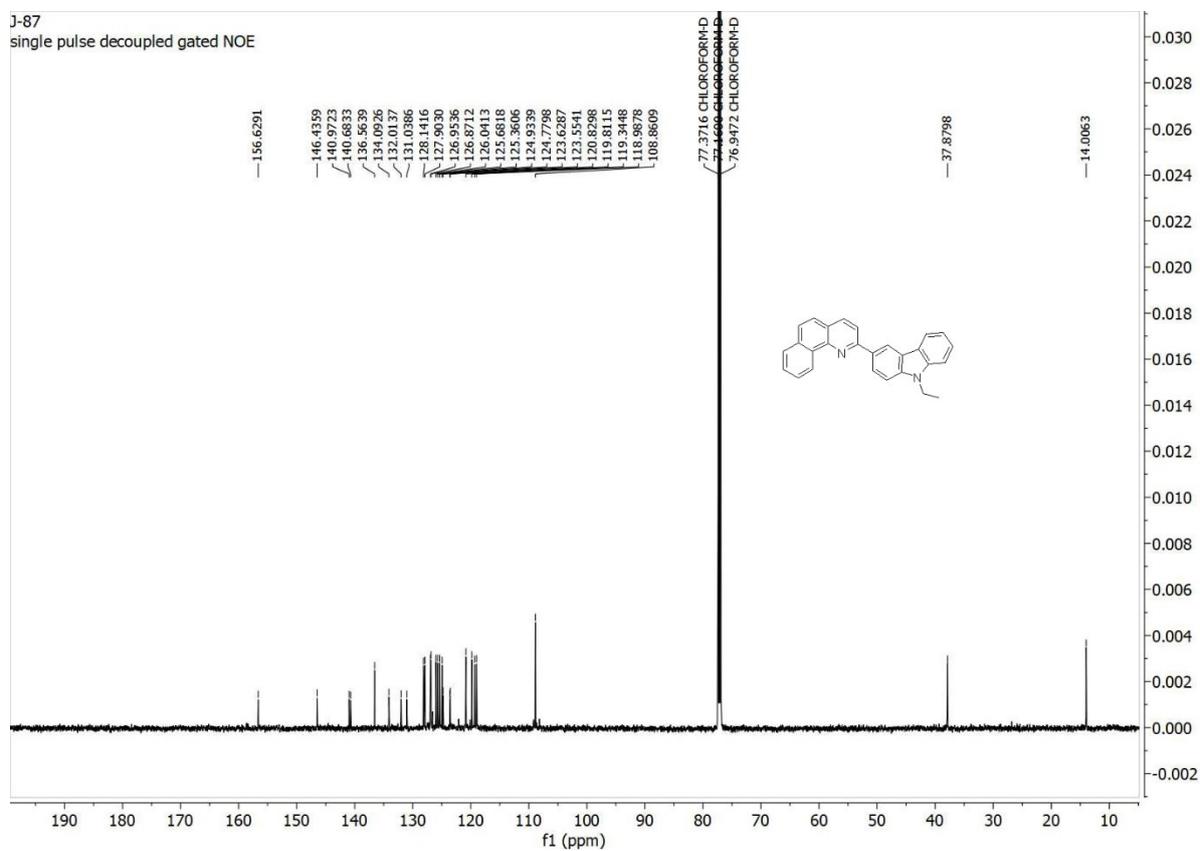


Figure S106. ^{13}C NMR Spectrum of **28g** in CDCl_3 .

4. Photophysical data of synthesised compounds

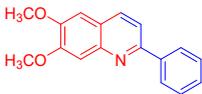
The fluorescent quantum yield (Φ_F) was measured relative to quinine sulfate ($\Phi_F = 0.546$) (0.1 M H₂SO₄ at 350 nm excitation) as a reference compound. For the measurement of UV-Vis absorption and fluorescence emission of samples, stock solution (1.0 mM) was prepared in CHCl₃ and diluted to final concentration (2.0 μ M) using CHCl₃. These quantum yields (QY) were calculated by using the equation as follows:

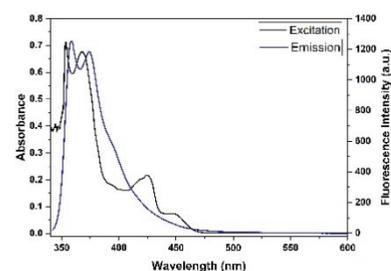
$$\Phi_S = \Phi_R \times \frac{I_S}{I_R} \times \frac{A_R}{A_S} \times \frac{\eta_S^2}{\eta_R^2}$$

R – Reference; *S* – Sample

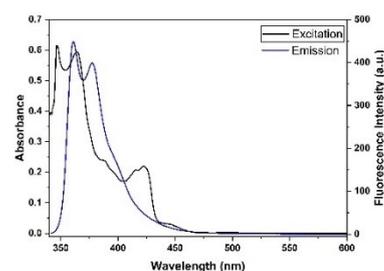
Where Φ is the quantum yields, η is the refractive index of the solvent, *I* is the integrated fluorescence intensity and *A* is the absorbance.

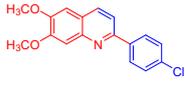
Figure S107. Photophysical properties and graphical data of quinoline-2-carboxylate derivatives:

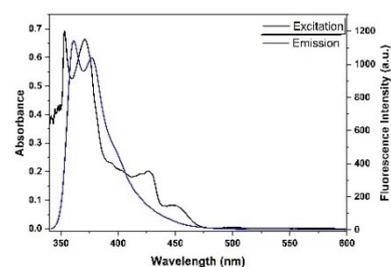
15a 	UV-Vis		Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Inten	sity	
	330	359	1253		

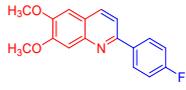


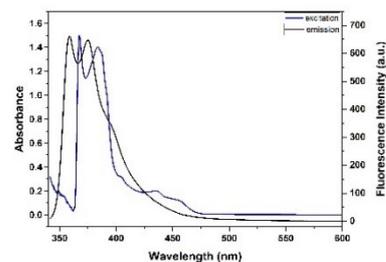
16a 	UV-Vis		Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Inten	sity	
	330	361	448		

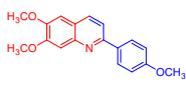


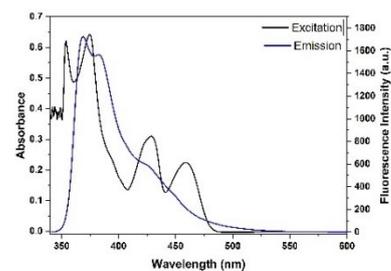
17a 	UV-Vis		Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Inten	sity	
	330	361	1142		



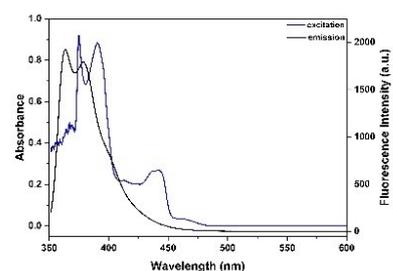
18a	UV-Vis	Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Intensity	
	330	359	662	0.08

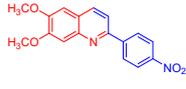


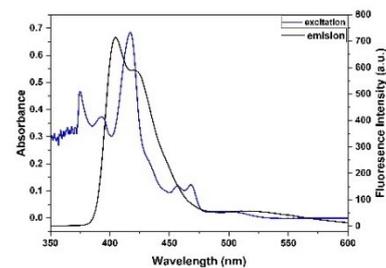
19a	UV-Vis	Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Intensity	
	330	368	1719	0.24

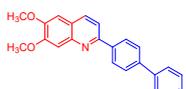


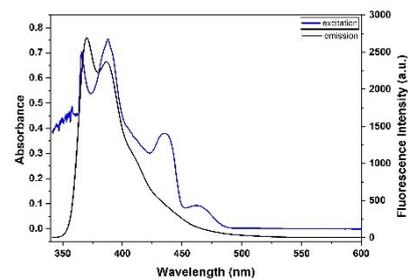
20a	UV-Vis	Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Intensity	
	330	369	2026	0.22

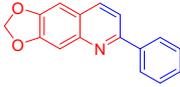


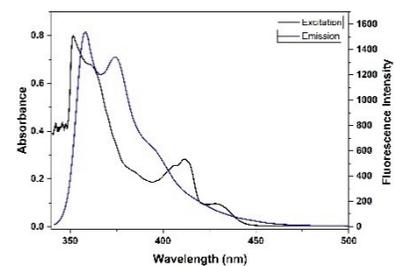
21a	UV-Vis	Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Intensity	
	330	406	713	0.25

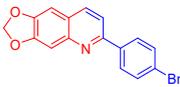


22a	UV-Vis	Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Intensity	
	330	370	2689	0.26



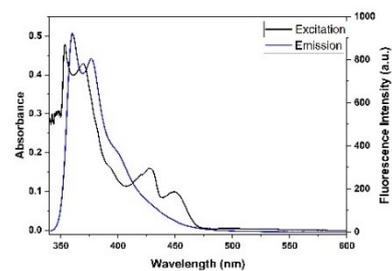
15b	UV-Vis	Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Intensity	
	330	358	1540	0.18

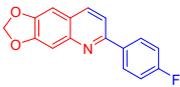


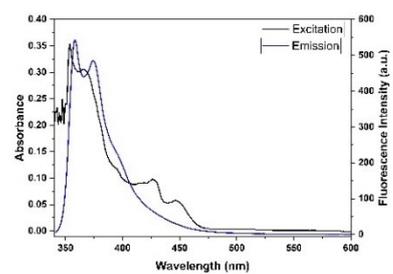
16b	UV-Vis	Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Intensity	
	330	361	448	0.06



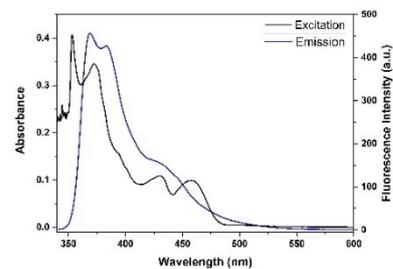
17b	UV-Vis	Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Intensity	
	330	360	922	0.20

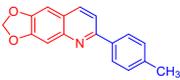


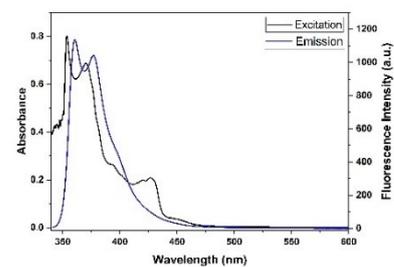
18b	UV-Vis	Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Intensity	
	330	358	542	0.19

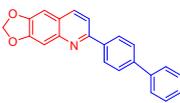


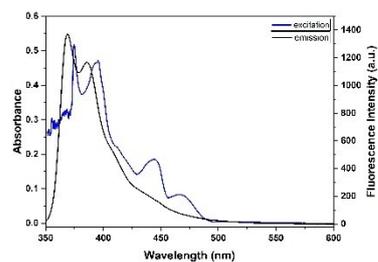
19b	UV-Vis	Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Intensity	
	330	369	457	0.19

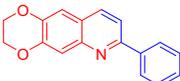


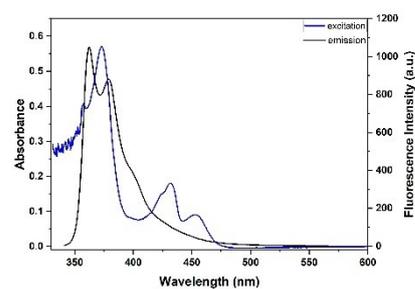
20b	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	361	1134	0.18	

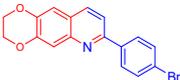


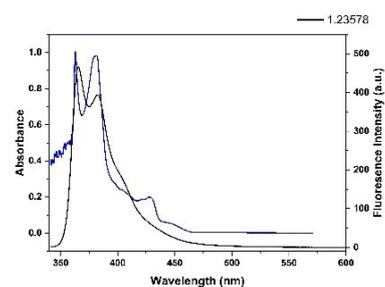
22b	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	369	1370	0.26	

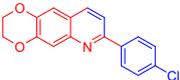


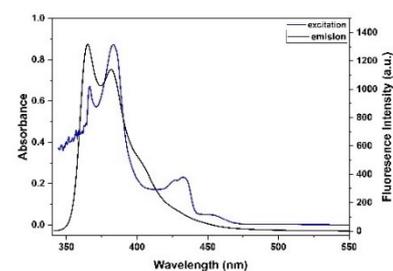
15c	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	361	1046	0.18	



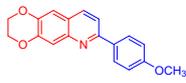
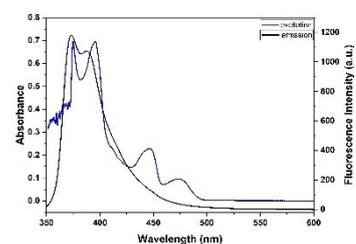
16c	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	365	465	0.08	



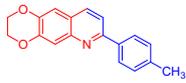
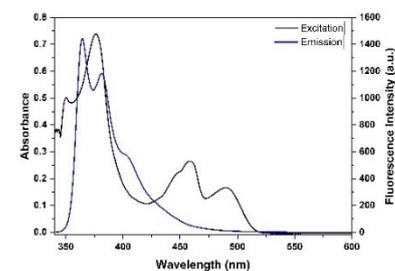
17c	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	365	1318	0.18	



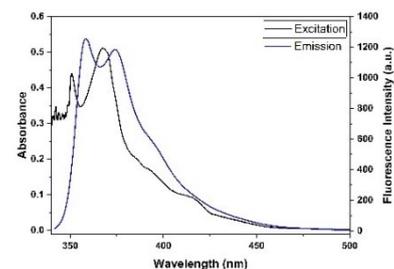
19c	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten	sity	
	330	373	1179		

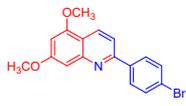
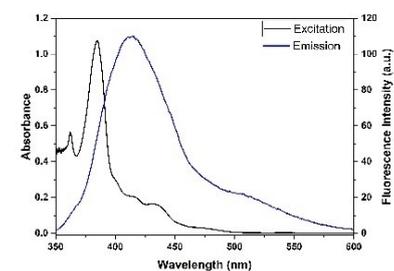
20c	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten	sity	
	330	364	1431		

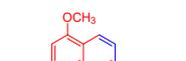
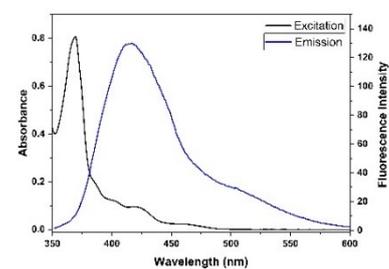
15d	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten	sity	
	330	360	1250		

16d	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten	sity	
	330	414	109		

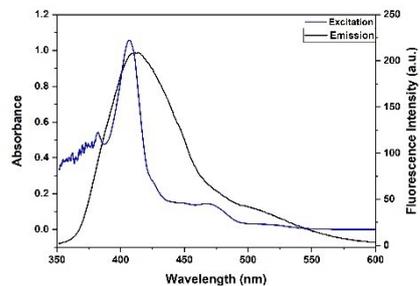



17d	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten	sity	

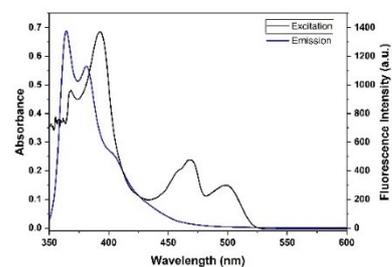



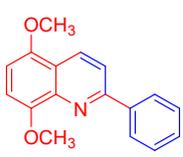
	330	416	129	0.09
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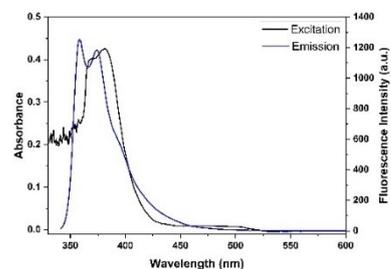
18d 	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	412	208	0.09	

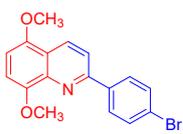


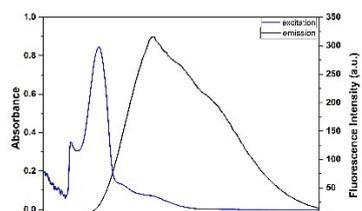
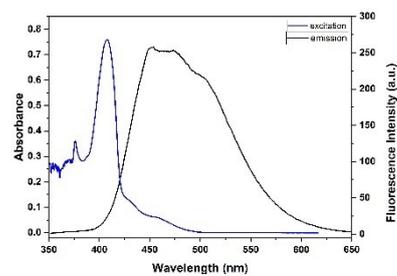
20d 	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	364	1373	0.19	



15e 	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	358	1251	0.20	

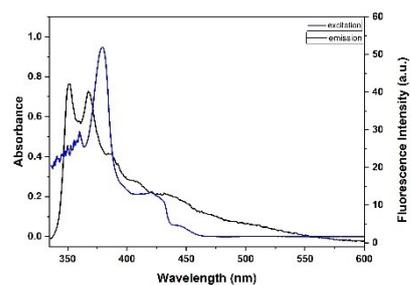


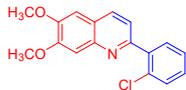
16e 	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	453	257	0.06	

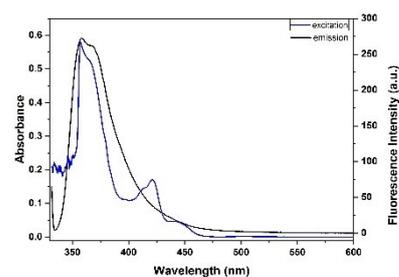


18e	UV-Vis	Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Intensity	
	330	450	316	0.33

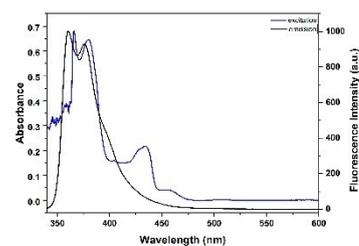
16f	UV-Vis	Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Intensity	
	330	351	42	0.01

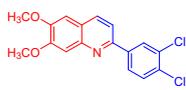


23a	UV-Vis	Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Intensity	
	330	358	272	0.06

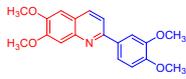


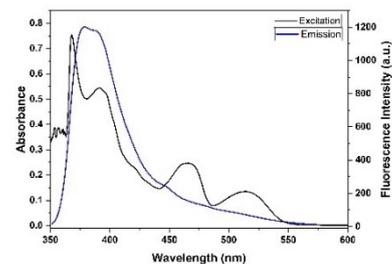
24a	UV-Vis	Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Intensity	
	330	360	1008	0.16

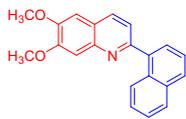


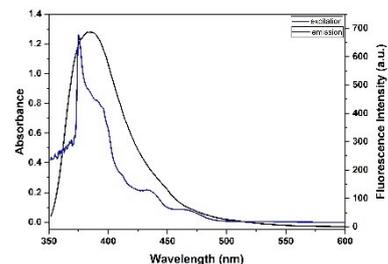
25a	UV-Vis	Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Intensity	
	330	363	1916	0.25

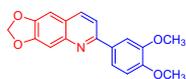


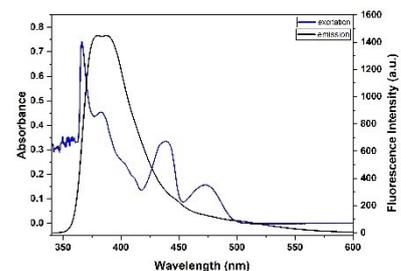
26a	UV-Vis	Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Intensity	
	330	379	1200	0.23

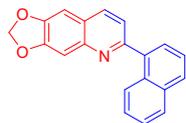


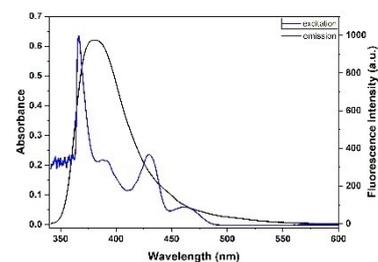
27a	UV-Vis	Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Intensity	
	330	385	686	0.15

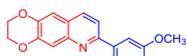


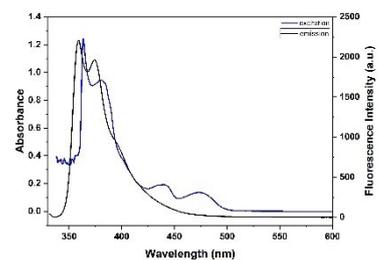
26b	UV-Vis	Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Intensity	
	330	386	1448	0.20



27b	UV-Vis	Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Intensity	
	330	381	974	0.19

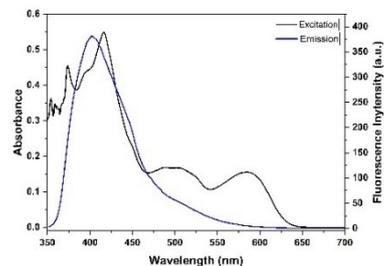


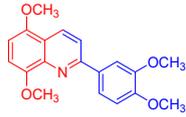
26c	UV-Vis	Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Intensity	
				

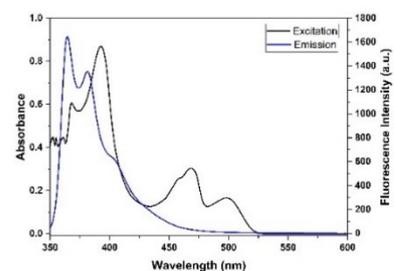


	330	358	2206	0.22
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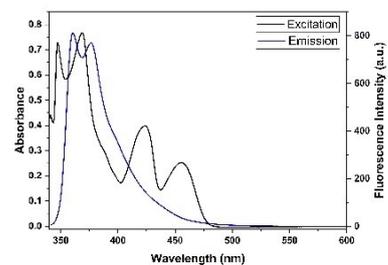
26d 	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	402	382	0.15	

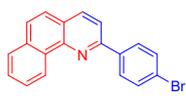


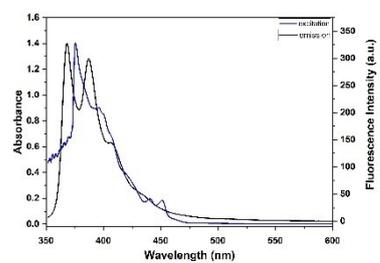
26e 	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	364	1635	0.17	



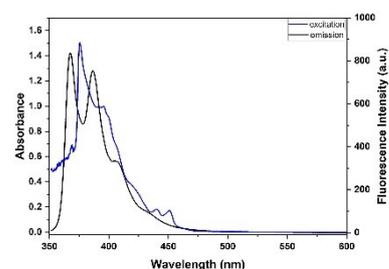
15g 	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	365	832	0.16	



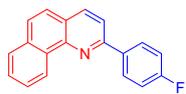
16g 	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	367	326	0.05	

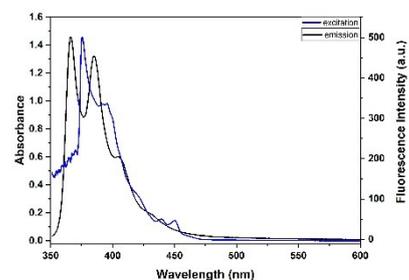


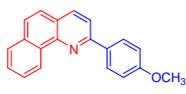
17g	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		

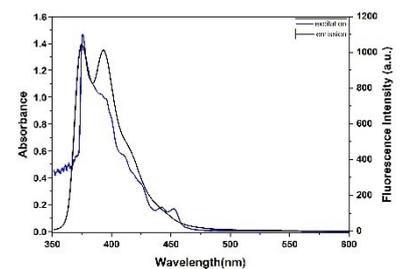


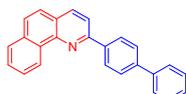
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity	
	330	367	833	0.14

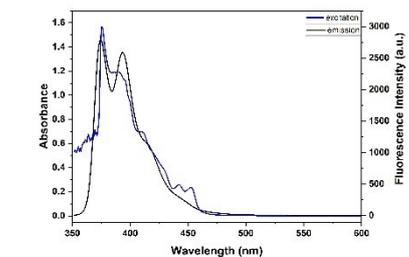
18g 	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	365	505	0.10	

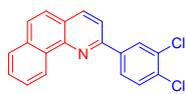


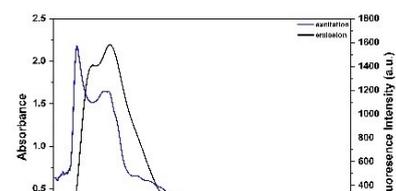
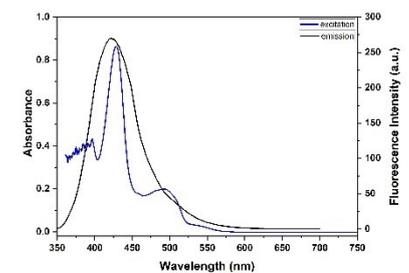
19g 	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	374	1043	0.21	

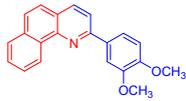


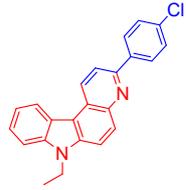
22g 	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	374	2784	0.34	

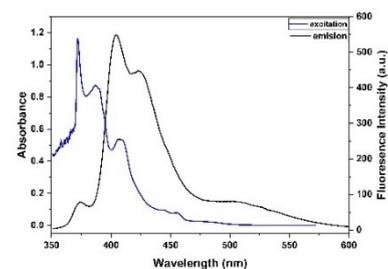


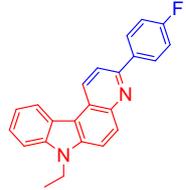
25g 	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten sity		
	330	422	270	0.09	

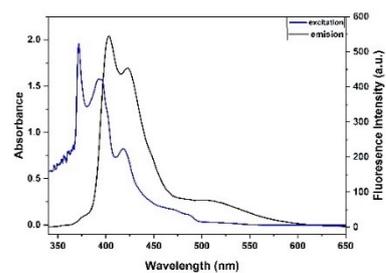


26g 	UV-Vis	Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Inten sity	
	330	399	1582	0.20

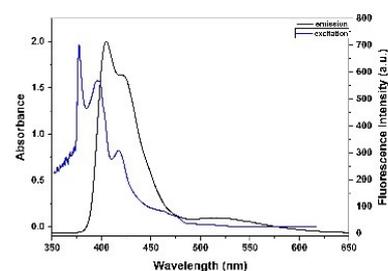
17h 	UV-Vis	Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Inten sity	
	330	404	547	0.28



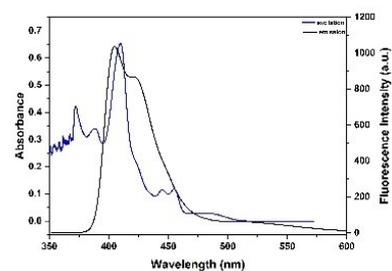
18h 	UV-Vis	Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Inten sity	
	330	403	545	0.14



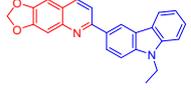
24h 	UV-Vis	Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Inten sity	
	330	406	713	0.25

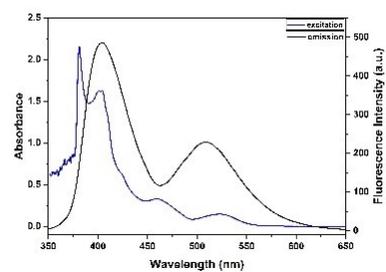


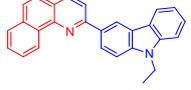
28a	UV-Vis	Fluorescence		Φ_F
	λ_{Ex} (nm)	λ_{Em} (nm)	Inten sity	



		(nm)	sity	
	330	405	1031	0.36

28b 	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten	sity	
	330	404	485	0.11	



28g 	UV-Vis		Fluorescence		Φ_F
	λ_{EX} (nm)	λ_{Em} (nm)	Inten	sity	
	330	409	1716	0.27	

