

Preparation of chitosan/cellulose composite copper catalyst for green synthesis in the construction of C-Si bond in aqueous phase

Yaoyao Zhang,^a Biao Han,^a Zelang Zhang,^a Xue Zhao,^a Weishuang Li,^a Bojie Li^a and Lei Zhu^{*a,b}

^(a) School of Chemistry and Materials Science, Hubei Key Laboratory of Quality Control of Characteristic Fruits And Vegetables, Hubei Engineering University, Xiaogan, 432000, China)

^(b) Hubei Key Laboratory of Bioinorganic Chemistry & Materia Medica, Huazhong University of Science and Technology, Wuhan 430074, China)

Dedicated to the 80th anniversary celebration of Hubei Engineering University

1. Different raw material ratios of chitosan/cellulose composite microspheres

In order to prepare cross-linked-chitosan-cellulose composite microbeads with good morphology, three different ratios of chitosan and cellulose were prepared according to the procedure of 2.3.1 Preparation of the support material. As a result, while the mass ratios of chitosan and cellulose were 2:1 could obtain better morphology of **CC@Cu**, while 1:1 and 3:1 were unable to form a complete sphere.

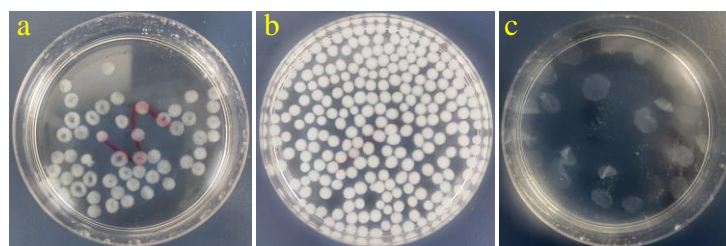
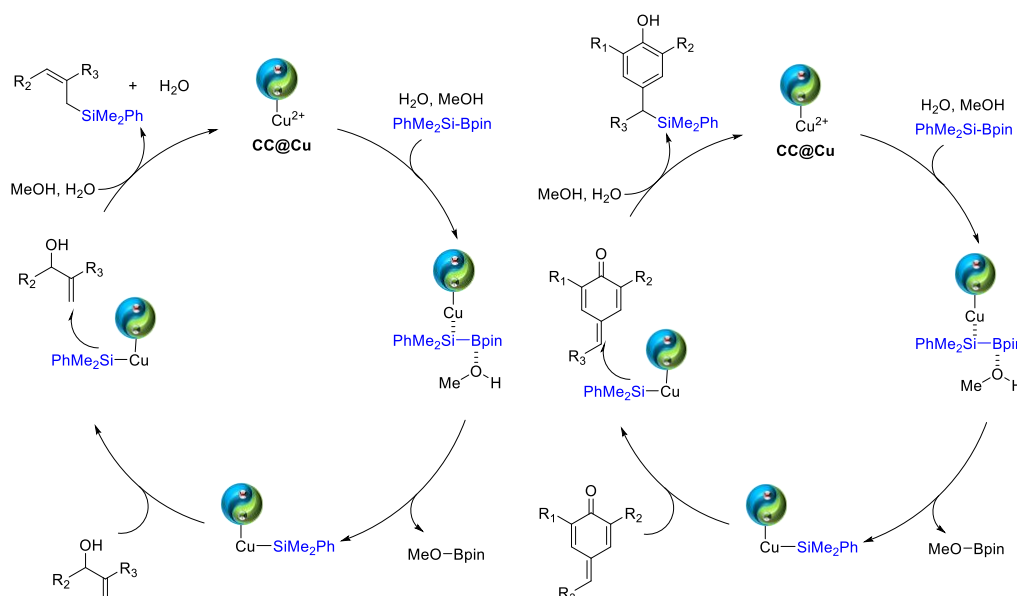


Fig S1. (a)chitosan : cellulose =1:1; (b)chitosan : cellulose =2:1; (c)chitosan : cellulose =3:1;

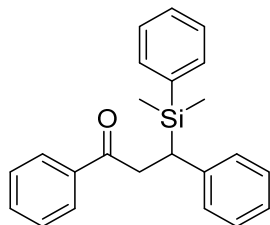
2. Proposed mechanism for the reaction of MBH alcohols and quinones



3. Characterization data for products

Most of adducts are literature-known and obtained characterization data for these compounds is in full agreement with reported data.

3-(Dimethyl(phenyl)silyl)-1,3-diphenylpropan-1-one 3a

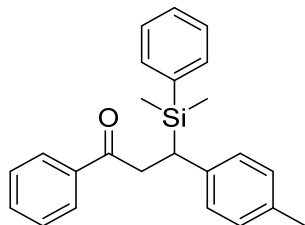


White solid

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.76 – 7.73 (m, 2H), 7.48 – 7.44 (m, 1H), 7.42 – 7.39 (m, 2H), 7.36 – 7.28 (m, 5H), 7.15 – 7.11 (m, 2H), 7.08 – 7.02 (m, 1H), 6.95 – 6.93 (m, 2H), 3.49 (dd, J = 17.1, 10.2 Hz, 1H), 3.19 – 3.05 (m, 2H), 0.26 (s, 3H), 0.20 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 199.1, 142.4, 137.1, 136.9, 134.2, 132.8, 129.4, 128.5, 128.1, 128.0, 127.8, 127.7, 124.8, 39.0, 31.1, – 3.8, – 5.1.

3-(Dimethyl(phenyl)silyl)-1-phenyl-3-(*p*-tolyl)propan-1-one 3b

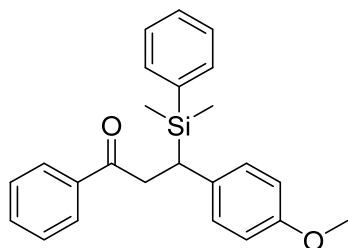


Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.79 (d, J = 7.6 Hz, 2H), 7.52 – 7.46 (m, 3H), 7.40 – 7.33 (m, 5H), 7.00 (d, J = 7.8 Hz, 2H), 6.89 (d, J = 7.7 Hz, 2H), 3.49 (dd, J = 17.0, 10.4 Hz, 1H), 3.21 – 3.04 (m, 2H), 2.27 (s, 3H), 0.30 (s, 3H), 0.23 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 199.2, 139.1, 137.10, 137.05, 134.2, 134.1, 132.8, 129.3, 128.9, 128.5, 128.0, 127.8, 127.6, 39.0, 30.6, 21.0, – 3.7, – 5.2.

3-(Dimethyl(phenyl)silyl)-3-(4-methoxyphenyl)-1-phenylpropan-1-one 3c

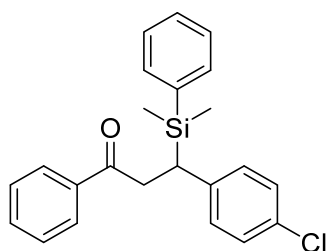


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.80 – 7.78 (m, 2H), 7.52 – 7.45 (m, 3H), 7.41 – 7.33 (m, 5H), 6.93 – 6.88 (m, 2H), 6.75 – 6.72 (m, 2H), 3.75 (s, 3H), 3.46 (dd, J = 17.0, 10.5 Hz, 1H), 3.21 – 3.01 (m, 2H), 0.30 (s, 3H), 0.24 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 199.4, 157.0, 137.1, 137.0, 134.21, 134.19, 132.8, 129.3, 128.6, 128.5, 128.0, 127.8, 113.6, 55.2, 39.2, 30.1, – 3.8, – 5.2.

3-(4-Chlorophenyl)-3-(dimethyl(phenyl)silyl)-1-phenylpropan-1-one 3d

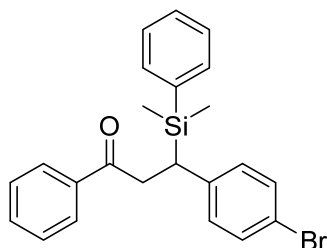


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.80 – 7.78 (m, 2H), 7.54 – 7.50 (m, 1H), 7.45 – 7.34 (m, 7H), 7.15 – 7.13 (m, 2H), 6.90 – 6.88 (m, 2H), 3.48 (dd, J = 17.2, 10.6 Hz, 1H), 3.23 – 3.08 (m, 2H), 0.30 (s, 3H), 0.26 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 198.8, 141.0, 136.9, 136.3, 134.2, 133.0, 130.4, 129.5, 128.9, 128.6, 128.2, 127.94, 127.90, 38.7, 30.7, – 4.0, – 5.2.

3-(4-Bromophenyl)-3-(dimethyl(phenyl)silyl)-1-phenylpropan-1-one 3e



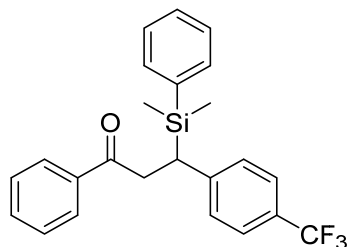
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.79 – 7.77 (m, 2H), 7.54 – 7.49 (m, 1H), 7.44 – 7.33 (m, 7H), 7.28 – 7.26 (m, 2H), 6.85 – 6.81 (m, 2H), 3.46 (dd, J = 17.2, 10.6 Hz, 1H), 3.22 – 3.03 (m, 2H), 0.30 (s, 3H), 0.25 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 198.7, 141.6, 136.9, 136.3, 134.2, 133.0, 131.1, 129.5, 129.3, 128.5, 127.9, 127.9, 118.4, 38.7, 30.8, – 4.0, – 5.2.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 445.0590; found 445.0594.

3-(Dimethyl(phenyl)silyl)-1-phenyl-3-(4-(trifluoromethyl)phenyl)propan-1-one 3f



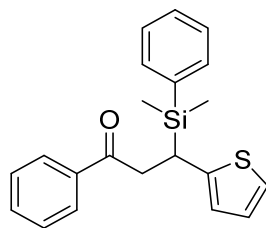
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.82 – 7.80 (m, 2H), 7.55 – 7.51 (m, 1H), 7.44 – 7.34 (m, 9H), 7.08 (d, J = 8 Hz, 2H), 3.56 (dd, J = 17.4, 10.5 Hz, 1H), 3.28 – 3.18 (m, 2H), 0.32 (s, 3H), 0.27 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 198.5, 147.0, 136.8, 136.0, 134.1, 133.1, 129.6, 128.6, 128.0, 127.9, 127.7, 127.1, 126.8, 125.8, 125.10, 125.06, 125.03, 124.99, 123.1, 38.5, 31.5, – 4.0, – 5.2.

HRMS (ESI⁺) calcd for C₂₃H₂₃F₃OSiNa [M+Na]⁺: 435.1364; found 435.1362.

3-(Dimethyl(phenyl)silyl)-1-phenyl-3-(thiophen-2-yl)propan-1-one 3g



Brown oil

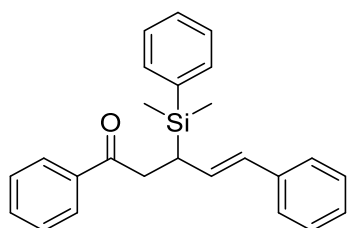
^1H NMR (400 MHz, Chloroform-*d*); δ = 7.81 – 7.79 (m, 2H), 7.53 – 7.48 (m, 3H), 7.41

– 7.33 (m, 5H), 6.97 – 6.96 (m, 1H), 6.85 – 6.83 (m, 1H), 6.60 – 6.59 (m, 1H), 3.48 – 3.41 (m, 2H), 3.20 – 3.12 (m, 1H), 0.39 (s, 3H), 0.31 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 198.6, 146.0, 137.0, 136.6, 134.1, 132.9, 129.5, 128.5, 128.0, 127.9, 126.7, 123.1, 121.5, 40.6, 26.2, – 3.8, – 5.1.

HRMS (ESI⁺) calcd for C₂₁H₂₂OSSiNa [M+Na]⁺: 373.1060; found 373.1053.

(*E*)-3-(dimethyl(phenyl)silyl)-1,5-diphenylpent-4-en-1-one 3h

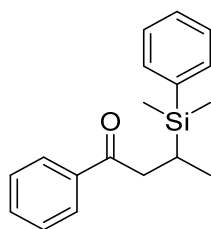


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.62 – 7.59 (m, 2H), 7.35 – 7.29 (m, 3H), 7.22 – 7.16 (m, 5H), 7.06 – 7.00 (m, 4H), 6.96 – 6.92 (m, 1H), 5.99 – 5.88 (m, 2H), 2.94 – 2.81 (m, 2H), 2.49 – 2.43 (m, 1H), 0.18 (s, 3H), 0.17 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 199.5, 138.0, 137.0, 136.8, 134.1, 132.9, 130.7, 129.4, 128.6, 128.4, 128.3, 128.1, 127.9, 126.5, 125.8, 38.2, 29.2, – 4.1, – 5.0.

3-(Dimethyl(phenyl)silyl)-1-phenylbutan-1-one 3i

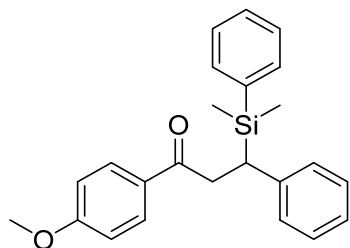


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.83 – 7.81 (m, 2H), 7.55 – 7.51 (m, 3H), 7.43 – 7.37 (m, 5H), 3.02 (dd, J = 15.8, 3.3 Hz, 1H), 2.68 – 2.61 (m, 1H), 1.63 – 1.59 (m, 1H), 0.98 (d, J = 7.3 Hz, 3H), 0.33 (s, 3H), 0.32 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 200.7, 137.6, 137.1, 134.0, 132.8, 129.2, 128.5, 128.1, 127.9, 40.7, 15.9, 14.6, – 4.7, – 5.4.

3-(Dimethyl(phenyl)silyl)-1-(4-methoxyphenyl)-3-phenylpropan-1-one 3j

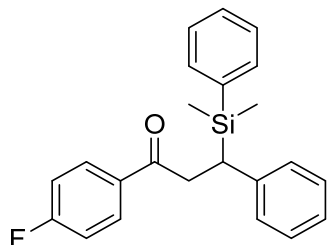


Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.83 – 7.77 (m, 2H), 7.46 – 7.44 (m, 2H), 7.39 – 7.33 (m, 3H), 7.19 – 7.15 (m, 2H), 7.08 – 7.04 (m, 1H), 7.00 – 6.97 (m, 2H), 6.88 – 6.84 (m, 2H), 3.84 (s, 3H), 3.49 (dd, J = 16.6, 10.0 Hz, 1H), 3.18 – 3.09 (m, 2H), 0.30 (s, 3H), 0.24 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 197.6, 163.2, 142.5, 136.9, 134.2, 130.3, 130.2, 129.3, 128.1, 127.8, 127.7, 124.8, 113.6, 55.5, 38.5, 31.2, – 3.8, – 5.1.

3-(Dimethyl(phenyl)silyl)-1-(4-fluorophenyl)-3-phenylpropan-1-one 3k



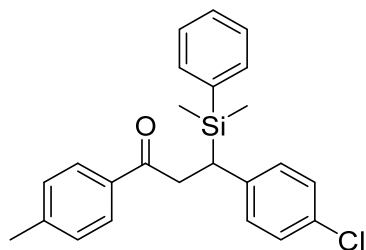
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.82 – 7.79 (m, 2H), 7.46 – 7.33 (m, 5H), 7.21 – 7.17 (m, 2H), 7.10 – 6.98 (m, 5H), 3.49 (dd, J = 17.0, 10.3 Hz, 1H), 3.20 – 3.06 (m, 2H), 0.31 (s, 3H), 0.25 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 197.6, 166.8, 164.3, 142.2, 136.8, 134.2, 133.5, 133.4, 130.6, 130.5, 129.4, 128.2, 127.8, 127.7, 124.9, 115.6, 115.4, 38.9, 31.2, – 3.8, – 5.4.

HRMS (ESI⁺) calcd for C₂₃H₂₃FOSiNa [M+Na]⁺: 385.1401; found 385.1394.

3-(4-Chlorophenyl)-3-(dimethyl(phenyl)silyl)-1-(*p*-tolyl)propan-1-one 3l



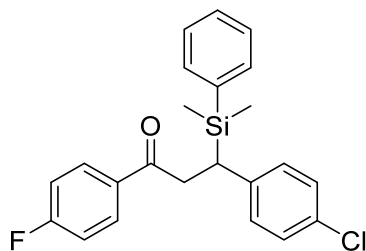
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.70 – 7.68 (m, 2H), 7.44 – 7.33 (m, 5H), 7.20 – 7.18 (m, 2H), 7.14 – 7.10 (m, 2H), 6.90 – 6.86 (m, 2H), 3.44 (dd, J = 16.9, 10.4 Hz, 1H), 3.19 – 3.04 (m, 2H), 2.38 (s, 3H), 0.29 (s, 3H), 0.23 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 198.4, 143.8, 141.1, 136.4, 134.4, 134.2, 130.3, 129.5, 129.2, 128.9, 128.2, 128.1, 127.9, 38.5, 30.8, 21.7, – 4.0, – 5.2.

HRMS (ESI⁺) calcd for C₂₄H₂₅ClOSiNa [M+Na]⁺: 415.1260; found 415.1255.

3-(4-Chlorophenyl)-3-(dimethyl(phenyl)silyl)-1-(4-fluorophenyl)propan-1-one 3m

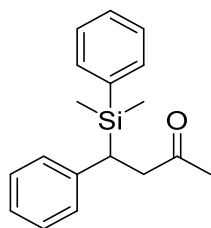


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.82 – 7.78 (m, 2H), 7.44 – 7.33 (m, 5H), 7.15 – 7.12 (m, 2H), 7.08 – 7.03 (m, 2H), 6.95 – 6.86 (m, 2H), 3.42 (dd, J = 17.1, 10.6 Hz, 1H), 3.18 – 3.02 (m, 2H), 0.29 (s, 3H), 0.25 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 197.2, 166.9, 164.3, 140.9, 136.3, 134.1, 133.30, 133.28, 130.6, 130.51, 130.46, 129.5, 128.8, 128.3, 127.9, 115.7, 115.5, 38.7, 30.8, – 4.0, – 5.3.

4-(Dimethyl(phenyl)silyl)-4-phenylbutan-2-one 3n

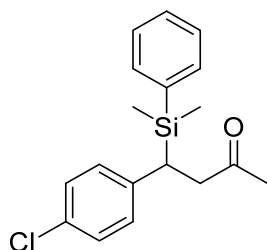


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.42 – 7.32 (m, 5H), 7.22 – 7.17 (m, 2H), 7.11 – 7.06 (m, 1H), 6.95 – 6.93 (m, 2H), 2.96 – 2.87 (m, 2H), 2.68 – 2.59 (m, 1H), 1.95 (s, 3H), 0.24 (s, 3H), 0.22 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 208.3, 142.0, 136.6, 134.2, 129.4, 128.2, 127.8, 127.6, 124.9, 44.0, 31.4, 30.0, – 4.0, – 5.4.

4-(4-Chlorophenyl)-4-(dimethyl(phenyl)silyl)butan-2-one 3o



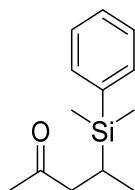
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.42 – 7.32 (m, 5H), 7.17 – 7.13 (m, 2H), 6.86 – 6.82 (m, 2H), 2.90 – 2.82 (m, 2H), 2.87 – 2.60 (m, 1H), 1.96 (s, 3H), 0.23 (s, 3H), 0.22 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 207.7, 140.7, 136.1, 134.1, 130.5, 129.5, 128.8, 128.3, 127.9, 43.8, 30.9, 30.1, – 4.2, – 5.3.

HRMS (ESI⁺) calcd for C₁₈H₂₁ClOSiNa [M+Na]⁺: 339.0944; found 339.0942.

4-(Dimethyl(phenyl)silyl)pentan-2-one 3p

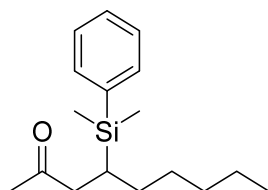


Pale yellow liquid

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.51 – 7.48 (m, 2H), 7.38 – 7.35 (m, 3H), 2.44 – 2.39 (m, 1H), 2.21 – 2.14 (m, 1H), 2.07 (s, 3H), 1.54 – 1.45 (m, 1H), 0.93 (d, J = 7.3 Hz, 3H), 0.28 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 209.5, 137.5, 133.9, 129.1, 127.8, 45.9, 30.0, 15.2, 14.5, – 4.8, – 5.3.

4-(Dimethyl(phenyl)silyl)nonan-2-one 3q

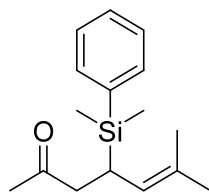


Yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.52 – 7.48 (m, 2H), 7.38 – 7.33 (m, 3H), 2.43 – 2.28 (m, 2H), 2.03 (s, 3H), 1.53 – 1.44 (m, 2H), 1.26 – 1.15 (m, 7H), 0.85 – 0.81 (m, 3H), 0.28 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 209.3, 138.2, 133.9, 129.0, 127.8, 44.6, 32.1, 30.5, 29.9, 28.9, 22.5, 20.3, 14.1, – 3.8, – 4.3.

4-(Dimethyl(phenyl)silyl)-6-methylhept-5-en-2-one 3r

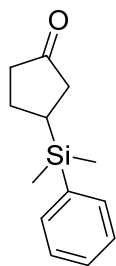


Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.49 – 7.47 (m, 2H), 7.38 – 7.32 (m, 3H), 4.85 – 4.81 (m, 1H), 2.46 – 2.25 (m, 3H), 2.01 (s, 3H), 1.67 (s, 3H), 1.50 (s, 3H), 0.28 (s, 3H), 0.26 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 209.7, 137.2, 134.0, 130.8, 129.2, 127.7, 123.8, 44.4, 29.6, 25.9, 24.9, 18.1, – 4.4, – 5.4.

3-(Dimethyl(phenyl)silyl)cyclopentan-1-one 3s

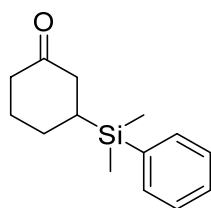


Pale yellow liquid

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.52 – 7.49 (m, 2H), 7.42 – 7.34 (m, 3H), 2.31 – 2.20 (m, 2H), 2.16 – 2.05 (m, 2H), 1.93 – 1.85 (m, 1H), 1.74 – 1.50 (m, 2H), 0.33 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 221.3, 137.0, 133.9, 129.4, 128.0, 40.2, 39.4, 25.0, 24.0, – 4.89, – 4.94.

3-(Dimethyl(phenyl)silyl)cyclohexan-1-one 3t

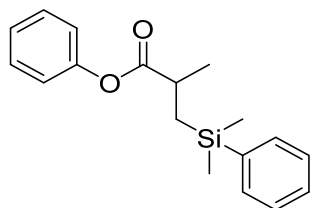


Pale yellow liquid

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.48 – 7.46 (m, 2H), 7.38 – 7.34 (m, 3H), 2.39 – 2.10 (m, 5H), 1.83 – 1.61 (m, 2H), 1.47 – 1.25 (m, 2H), 0.30 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 212.8, 136.6, 133.9, 129.3, 127.9, 42.4, 41.9, 29.8, 27.7, 26.1, – 5.3, – 5.4.

Phenyl 3-(dimethyl(phenyl)silyl)-2-methylpropanoate 3u



Colorless oil

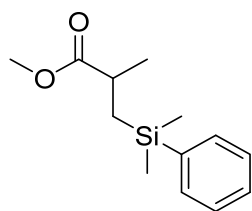
^1H NMR (400 MHz, Chloroform-*d*); δ = 7.59 – 7.56 (m, 2H), 7.41 – 7.35 (m, 5H), 7.24

– 7.20 (m, 1H), 7.01 – 6.98 (m, 2H), 2.82 – 2.76 (m, 1H), 1.50 – 1.44 (m, 1H), 1.32 (d, $J = 12.3$ Hz, 3H), 1.09 – 1.03 (m, 1H), 0.39 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); $\delta = 178.4, 153.2, 141.0, 136.0, 131.7, 131.5, 130.3, 128.02, 123.8, 38.3, 27.4, 23.0, 0.1, 0.0$.

HRMS (ESI⁺) calcd for $\text{C}_{18}\text{H}_{22}\text{O}_2\text{Si}$ $[\text{M}+\text{Na}]^+$: 321.1283; found 321.1281.

Methyl 3-(dimethyl(phenyl)silyl)-2-methylpropanoate **3v**

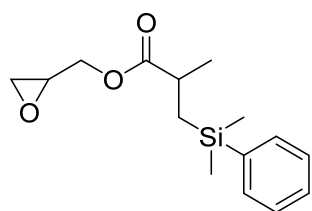


Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); $\delta = 7.51 – 7.49$ (m, 2H), 7.36 – 7.34 (m, 3H), 3.54 (s, 3H), 2.58 – 2.49 (m, 1H), 1.28 – 1.22 (m, 1H), 1.15 (d, $J = 6.8$ Hz, 3H), 0.94 (dd, $J = 7.0, 14.9$ Hz, 1H), 0.29 (d, $J = 2.5$ Hz, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); $\delta = 177.9, 138.7, 133.6, 129.0, 127.8, 51.5, 35.5, 20.9, 20.7, -2.5, -2.6$.

Oxiran-2-ylmethyl 3-(dimethyl(phenyl)silyl)-2-methylpropanoate **3w**



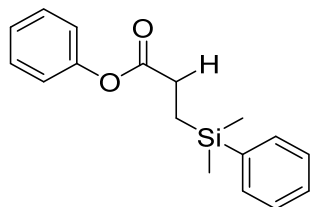
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); $\delta = 7.52 – 7.49$ (m, 2H), 7.37 – 7.34 (m, 3H), 4.29 – 4.20 (m, 1H), 3.80 – 3.71 (m, 1H), 3.16 – 3.11 (m, 1H), 2.83 – 2.81 (m, 1H), 2.61 – 2.55 (m, 2H), 1.33 – 1.27 (m, 1H), 1.17 (d, $J = 7.0$ Hz, 3H), 0.96 – 0.91 (m, 1H), 0.32 – 0.28 (m, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); $\delta = 179.6, 141.0, 136.0, 131.5, 130.2, 67.3, 51.8, 47.1, 38.0, 23.2, 0.0, -0.1$.

HRMS (ESI⁺) calcd for C₁₅H₂₂O₃Si [M+Na]⁺: 301.1226; found 301.1230.

Phenyl 3-(dimethyl(phenyl)silyl)propanoate 3x



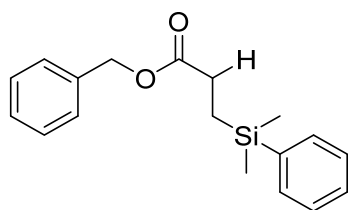
Colorless oil

¹H NMR (400 MHz, Chloroform-*d*); δ = 7.57 – 7.54 (m, 2H), 7.41 – 7.35 (m, 5H), 7.24 – 7.20 (m, 1H), 7.05 – 7.03 (m, 2H), 2.56 – 2.52 (m, 2H), 1.25 – 1.21 (m, 2H), 0.69 (s, 6H).

¹³C NMR (100 MHz, Chloroform-*d*); δ = 173.5, 150.8, 138.0, 133.7, 129.4, 129.3, 128.0, 125.7, 121.5, 29.1, 10.9, – 3.2.

HRMS (ESI⁺) calcd for C₁₇H₂₀O₂Si [M+Na]⁺: 307.1121; found 307.1125.

Benzyl 3-(dimethyl(phenyl)silyl)propanoate 3y

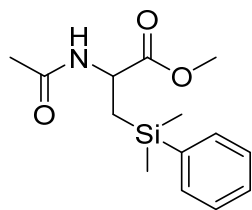


Colorless oil

¹H NMR (400 MHz, Chloroform-*d*); δ = 7.50 – 7.48 (m, 2H), 7.36 – 7.33 (m, 8H), 5.06 (s, 2H), 2.35 – 2.31 (m, 2H), 1.13 – 1.09 (m, 2H), 0.28 (s, 6H).

¹³C NMR (100 MHz, Chloroform-*d*); δ = 174.8, 138.1, 136.0, 133.6, 129.1, 128.6, 128.3, 128.2, 127.9, 66.3, 28.9, 25.0, 10.8, – 3.3.

Methyl 2-acetamido-3-(dimethyl(phenyl)silyl)propanoate 3z

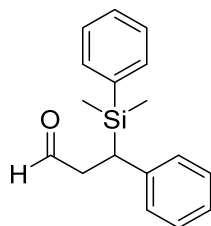


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.50 – 7.48 (m, 2H), 7.39 – 7.36 (m, 3H), 5.64 – 5.62 (m, 1H), 4.64 – 4.58 (m, 1H), 3.57 (s, 3H), 1.72 (m, 3H), 1.42 – 1.36 (m, 1H), 1.26 – 1.20 (m, 1H), 0.35 (s, 3H), 0.32 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 174.0, 169.5, 138.1, 133.5, 129.4, 128.1, 52.2, 49.3, 22.9, 20.2, – 2.8, – 3.2.

3-(Dimethyl(phenyl)silyl)-3-phenylpropanal 3A

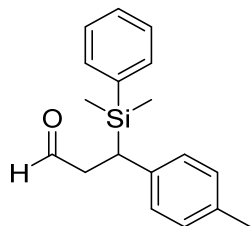


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 9.54 – 9.53 (m, 1H), 7.42 – 7.33 (m, 5H), 7.23 – 7.19 (m, 2H), 7.13 – 7.09 (m, 1H), 6.96 – 6.94 (m, 2H), 2.91 – 2.83 (m, 2H), 2.67 – 2.59 (m, 1H), 0.27 (d, J = 11.4 Hz, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 202.7, 141.1, 136.2, 134.1, 129.5, 128.3, 127.9, 127.7, 125.2, 43.5, 30.1, – 4.2, – 5.5.

3-(Dimethyl(phenyl)silyl)-3-(*p*-tolyl)propanal 3B



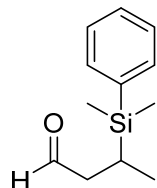
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 9.53 – 9.52 (m, 1H), 7.42 – 7.33 (m, 5H), 7.04

– 7.01 (m, 2H), 6.86 – 6.83 (m, 2H), 2.87 – 2.79 (m, 2H), 2.63 – 2.57 (m, 1H), 2.29 (s, 3H), 0.26 (d, $J = 11.4$ Hz, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); $\delta = 202.9, 137.8, 136.4, 134.6, 134.1, 129.5, 129.1, 127.9, 127.5, 43.5, 29.6, 21.0, -4.1, -5.6$.

3-(Dimethyl(phenyl)silyl)butanal 3C

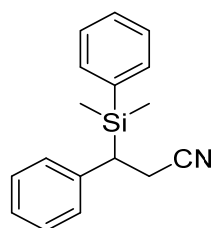


Pale yellow liquid

^1H NMR (400 MHz, Chloroform-*d*); $\delta = 9.67 - 9.66$ (m, 1H), 7.51 – 7.48 (m, 2H), 7.40 – 7.34 (m, 3H), 2.45 – 2.40 (m, 1H), 2.19 – 2.11 (m, 1H), 1.53 – 1.47 (m, 1H), 0.99 (d, $J = 7.4$ Hz, 3H), 0.30 (d, $J = 2.2$ Hz, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); $\delta = 203.5, 137.0, 133.9, 129.3, 127.9, 45.9, 14.6, 13.8, -4.9, -5.4$.

3-(Dimethyl(phenyl)silyl)-3-phenylpropanenitrile 3D

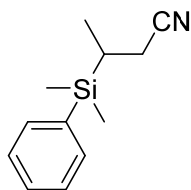


Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); $\delta = 7.43 - 7.37$ (m, 6H), 7.28 – 7.27 (m, 1H), 7.20 – 7.16 (m, 1H), 6.97 – 6.94 (m, 2H), 2.67 – 2.60 (m, 3H), 0.28 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); $\delta = 139.7, 135.2, 134.1, 129.9, 128.6, 128.1, 127.4, 126.0, 119.7, 33.0, 18.9, -4.0, -5.5$.

3-(Dimethyl(phenyl)silyl)butanenitrile 3E

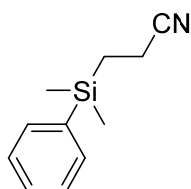


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.49 – 7.47 (m, 2H), 7.42 – 7.37 (m, 3H), 2.42 – 2.37 (m, 1H), 2.11 – 2.04 (m, 1H), 1.27 – 1.23 (m, 1H), 1.17 – 1.15 (m, 3H), 0.35 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 135.9, 133.8, 129.6, 128.1, 119.9, 20.5, 17.3, 14.5, – 4.9, – 5.6.

3-(Dimethyl(phenyl)silyl)propanenitrile 3F

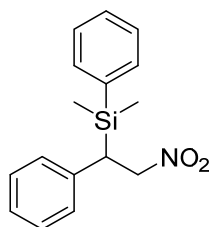


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.50 – 7.47 (m, 2H), 7.41 – 7.37 (m, 3H), 2.29 – 2.25 (m, 2H), 1.17 – 1.13 (m, 2H), 0.36 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 136.6, 133.5, 129.6, 128.2, 121.2, 12.2, 12.0, – 3.5.

Dimethyl(2-nitro-1-phenylethyl)(phenyl)silane 3G

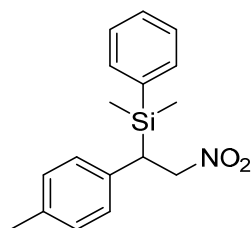


Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.45 – 7.36 (m, 5H), 7.25 – 7.22 (m, 2H), 7.18 – 7.14 (m, 1H), 6.99 – 6.97 (m, 2H), 4.86 (t, J = 13.4 Hz, 1H), 4.52 (dd, J = 13.6, 3.8 Hz, 1H), 3.28 – 3.24 (m, 1H), 0.30 (s, 3H), 0.28 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); $\delta = 137.5, 134.9, 133.9, 130.0, 128.6, 128.2, 127.3, 126.1, 77.1, 36.1, -3.9, -5.4$.

Dimethyl(2-nitro-1-(*p*-tolyl)ethyl)(phenyl)silane 3H

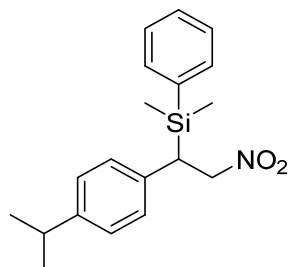


Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); $\delta = 7.46 - 7.37$ (m, 5H), 7.10 – 7.04 (m, 2H), 6.89 – 6.86 (m, 2H), 4.83 (t, $J = 13.5$ Hz, 1H), 4.50 (dd, $J = 13.5, 3.8$ Hz, 1H), 3.26 – 3.20 (m, 1H), 2.29 (m, 3H), 0.30 (s, 3H), 0.28 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); $\delta = 135.6, 135.1, 134.3, 134.0, 130.0, 129.3, 128.2, 127.2, 77.1, 35.7, 21.0, -3.9, -5.5$.

(1-(4-Isopropylphenyl)-2-nitroethyl)dimethyl(phenyl)silane 3I



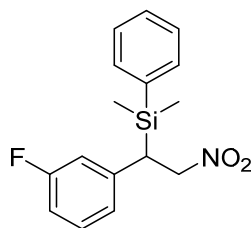
Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); $\delta = 7.45 - 7.36$ (m, 5H), 7.10 – 7.07 (m, 2H), 6.92 – 6.88 (m, 2H), 4.84 (t, $J = 13.5$ Hz, 1H), 4.49 (dd, $J = 13.6, 3.8$ Hz, 1H), 3.27 – 3.21 (m, 1H), 2.88 – 2.81 (m, 1H), 1.22 (d, $J = 7.0$ Hz, 6H), 0.30 (s, 3H), 0.27 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); $\delta = 146.5, 135.2, 134.5, 133.9, 130.0, 128.2, 127.2, 126.6, 77.1, 35.6, 33.6, 24.0, -3.8, -5.5$.

HRMS (ESI⁺) calcd for $\text{C}_{19}\text{H}_{25}\text{NO}_2\text{SiNa}$ [$\text{M}+\text{Na}$]⁺: 350.1548; found 350.1547.

(1-(3-Fluorophenyl)-2-nitroethyl)dimethyl(phenyl)silane 3J



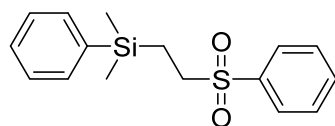
Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.46 – 7.39 (m, 5H), 7.23 – 7.18 (m, 1H), 6.88 – 6.83 (m, 1H), 6.77 – 6.74 (m, 1H), 6.69 – 6.65 (m, 1H), 4.81 (t, J = 13.6 Hz, 1H), 4.52 (dd, J = 13.8, 3.7 Hz, 1H), 3.30 – 3.24 (m, 1H), 0.32 (d, J = 2.0 Hz, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 164.1, 161.7, 140.4, 140.3, 134.3, 133.9, 130.2, 130.1, 130.0, 128.3, 122.93, 122.90, 114.4, 114.1, 113.2, 113.0, 76.6, 36.14, 36.12, – 4.0, – 5.4.

HRMS (ESI⁺) calcd for $\text{C}_{16}\text{H}_{18}\text{FNO}_2\text{SiNa}$ [$\text{M}+\text{Na}$]⁺: 326.0981; found 326.0983.

Dimethyl(phenyl)(2-(phenylsulfonyl)ethyl)silane 3K

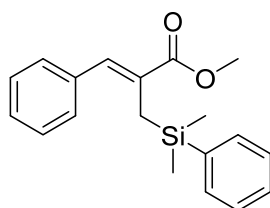


White solid

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.74 – 7.72 (m, 2H), 7.54 – 7.48 (m, 1H), 7.46 – 7.40 (m, 2H), 7.28 – 7.21 (m, 5H), 2.85 – 2.81 (m, 2H), 1.05 – 1.01 (m, 2H), 0.15 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 138.6, 136.5, 133.6, 133.4, 129.6, 129.2, 128.2, 128.1, 52.5, 8.5, – 3.4.

Methyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-3-phenylacrylate 5a

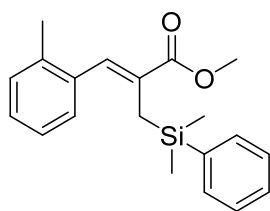


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.56 (s, 1H), 7.49 – 7.46 (m, 2H), 7.37 – 7.27 (m, 8H), 3.71 (s, 3H), 2.41 (s, 2H), 0.27 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 169.3, 138.6, 136.3, 135.7, 133.6, 130.9, 129.09, 129.05, 128.3, 127.8, 127.7, 51.9, 17.4, – 2.6.

Methyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-3-(*o*-tolyl)acrylate 5b



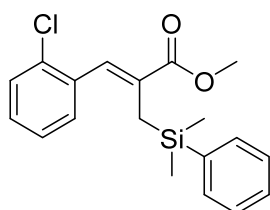
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.50 (s, 1H), 7.32 – 7.29 (m, 2H), 7.25 – 7.18 (m, 3H), 7.12 – 7.09 (m, 2H), 7.06 – 7.00 (m, 2H), 3.64 (s, 3H), 2.18 (s, 2H), 2.14 (s, 3H), 0.13 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 169.1, 138.6, 136.6, 135.6, 135.3, 133.5, 131.4, 130.1, 128.9, 128.4, 127.8, 127.7, 125.5, 51.8, 19.9, 17.0, – 2.6.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 347.1434; found 347.1438.

Methyl (Z)-3-(2-chlorophenyl)-2-((dimethyl(phenyl)silyl)methyl)acrylate 5c



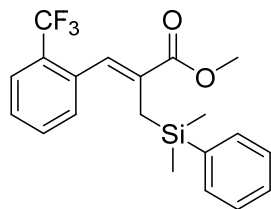
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.56 (s, 1H), 7.42 – 7.27 (m, 6H), 7.22 – 7.11 (m, 3H), 3.71 (s, 3H), 2.26 (s, 2H), 0.20 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 168.7, 138.4, 135.0, 134.0, 133.5, 132.9, 132.8, 129.9, 129.6, 129.01, 128.97, 127.7, 126.4, 52.0, 17.3, – 2.8.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 367.0885; found 367.0892.

Methyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-3-(2-(trifluoromethyl)phenyl)acrylate 5d



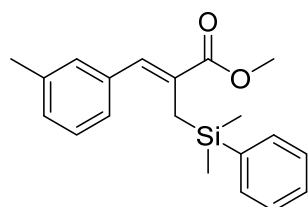
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.69 (s, 1H), 7.67 – 7.64 (m, 1H), 7.42 – 7.28 (m, 7H), 7.20 – 7.17 (m, 1H), 3.73 (s, 3H), 2.21 (s, 2H), 0.22 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 168.6, 138.5, 135.04, 135.02, 133.5, 132.2, 131.4, 130.2, 129.0, 128.7, 128.4, 127.7, 127.6, 126.1, 126.1, 126.01, 125.96, 125.3, 122.6, 52.1, 17.6, – 2.7.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 401.1152; found 401.1155.

Methyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-3-(*m*-tolyl)acrylate 5e

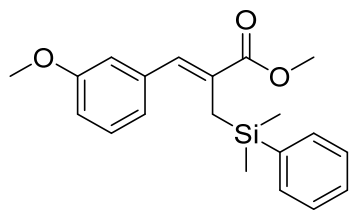


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.51 (s, 1H), 7.47 – 7.45 (m, 2H), 7.36 – 7.29 (m, 3H), 7.23 – 7.19 (m, 1H), 6.89 – 6.87 (m, 1H), 6.82 – 6.79 (m, 2H), 3.73 (s, 3H), 3.68 (s, 3H), 2.41 (s, 2H), 0.26 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 171.7, 162.0, 141.1, 140.2, 138.0, 136.1, 133.7, 131.8, 131.5, 130.2, 124.0, 116.9, 116.1, 57.7, 54.4, 20.0, 0.0.

Methyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-3-(3-methoxyphenyl)acrylate 5f



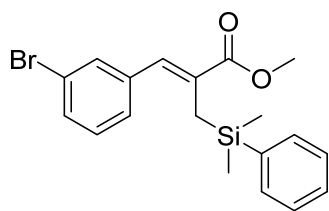
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.52 (s, 1H), 7.49 – 7.46 (m, 2H), 7.39 – 7.29 (m, 3H), 7.20 – 7.16 (m, 1H), 7.09 – 7.05 (m, 3H), 3.70 (s, 3H), 2.39 (s, 2H), 2.25 (s, 3H), 0.27 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 171.9, 141.2, 140.4, 138.6, 138.3, 136.1, 133.1, 132.2, 131.5, 131.1, 130.7, 130.2, 128.8, 54.4, 23.8, 20.0, 0.0.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 363.1380; found 363.1387.

Methyl (Z)-3-(3-bromophenyl)-2-((dimethyl(phenyl)silyl)methyl)acrylate 5g



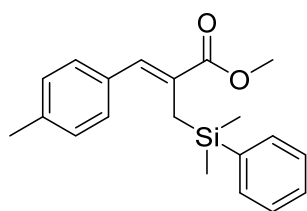
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.46 – 7.42 (m, 3H), 7.39 – 7.29 (m, 5H), 7.15 – 7.11 (m, 2H), 3.70 (s, 3H), 2.53 (s, 2H), 0.27 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 168.9, 138.4, 138.1, 133.8, 133.6, 132.4, 131.7, 130.7, 129.8, 129.2, 127.8, 127.6, 122.4, 52.0, 17.8, – 2.6.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 411.0386; found 411.0386.

Methyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-3-(*p*-tolyl)acrylate 5h

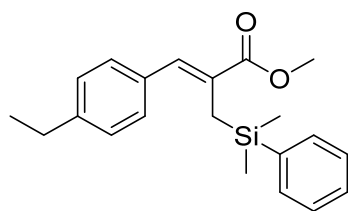


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.53 (s, 1H), 7.49 – 7.47 (m, 2H), 7.38 – 7.31 (m, 3H), 7.20 – 7.18 (m, 2H), 7.10 – 7.08 (m, 2H), 3.68 (s, 3H), 2.41 (s, 2H), 2.34 (s, 3H), 0.27 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 169.5, 138.8, 137.9, 135.7, 133.6, 133.4, 130.0, 129.2, 129.1, 129.0, 127.7, 51.9, 21.3, 17.5, – 2.5.

Methyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-3-(4-ethylphenyl)acrylate **5i**



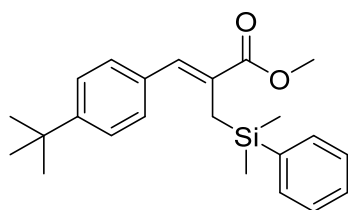
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.54 (s, 1H), 7.51 – 7.47 (m, 2H), 7.38 – 7.31 (m, 3H), 7.24 – 7.22 (m, 2H), 7.14 – 7.12 (m, 2H), 3.70 (s, 3H), 2.68 (dd, J = 15.2, 7.6 Hz, 2H), 2.42 (s, 2H), 1.27 (t, J = 7.6 Hz, 3H), 0.28 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 169.5, 144.2, 138.8, 135.8, 133.61, 133.59, 130.0, 129.3, 129.0, 127.8, 127.7, 51.9, 28.7, 17.5, 15.4, – 2.5.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 361.1594; found 361.1594.

Methyl (Z)-3-(4-(*tert*-butyl)phenyl)-2-((dimethyl(phenyl)silyl)methyl)acrylate **5j**



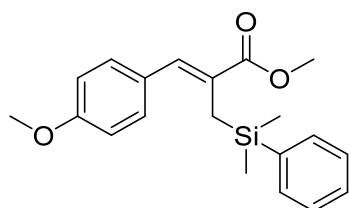
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.50 (s, 1H), 7.46 – 7.44 (m, 2H), 7.32 – 7.26 (m, 5H), 7.23 – 7.20 (m, 2H), 3.65 (s, 3H), 2.39 (s, 2H), 1.29 (s, 9H), 0.25 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 172.0, 153.5, 141.3, 138.1, 136.1, 135.8, 132.5, 131.53, 131.47, 130.2, 127.7, 54.3, 37.1, 33.7, 19.9, 0.0.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 389.1902; found 389.1907.

Methyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-3-(4-methoxyphenyl)acrylate 5k

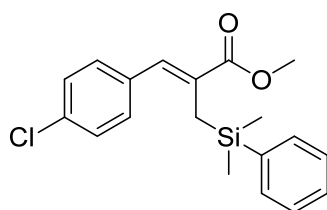


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.45 – 7.42 (m, 3H), 7.30 – 7.27 (m, 3H), 7.21 – 7.19 (m, 2H), 6.76 – 6.74 (m, 2H), 3.75 (s, 3H), 3.62 (s, 3H), 2.35 (s, 2H), 0.21 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 169.6, 159.3, 138.8, 135.4, 133.6, 130.8, 129.0, 128.8, 128.7, 127.8, 113.8, 55.3, 51.9, 17.4, –2.5.

Methyl (Z)-3-(4-chlorophenyl)-2-((dimethyl(phenyl)silyl)methyl)acrylate 5l

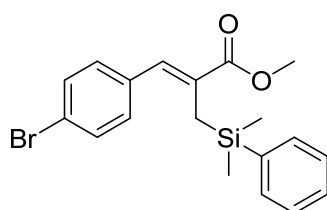


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.46 (s, 1H), 7.45 – 7.42 (m, 2H), 7.38 – 7.30 (m, 3H), 7.24 – 7.20 (m, 2H), 7.16 – 7.12 (m, 2H), 3.70 (s, 3H), 2.35 (s, 2H), 0.26 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 169.1, 138.3, 134.7, 134.2, 133.5, 131.6, 130.3, 129.2, 128.5, 127.8, 52.0, 17.7, – 2.6.

Methyl (Z)-3-(4-bromophenyl)-2-((dimethyl(phenyl)silyl)methyl)acrylate 5m

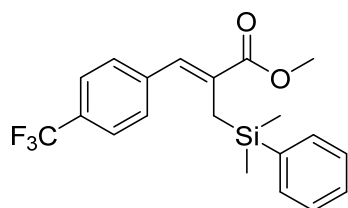


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.44 – 7.42 (m, 3H), 7.39 – 7.30 (m, 5H), 7.08 – 7.06 (m, 2H), 3.70 (s, 3H), 2.35 (s, 2H), 0.26 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 169.1, 138.3, 135.1, 134.2, 133.5, 131.7, 131.5, 130.6, 129.2, 127.8, 121.8, 52.0, 17.7, –2.6.

Methyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-3-(4-(trifluoromethyl)phenyl)acrylate 5n



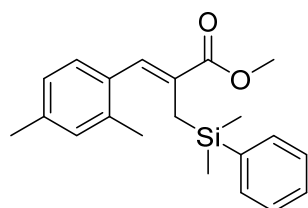
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.48 – 7.46 (m, 3H), 7.39 – 7.36 (m, 2H), 7.32 – 7.24 (m, 5H), 3.70 (s, 3H), 2.34 (s, 2H), 0.24 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 168.8, 139.9, 138.0, 133.8, 133.5, 133.1, 129.6, 129.2, 129.1, 127.8, 125.4, 125.3, 125.23, 125.19, 125.2, 122.7, 52.1, 17.8, –2.7.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 401.1152; found 401.1155.

Methyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-3-(2,4-dimethylphenyl)acrylate 5o



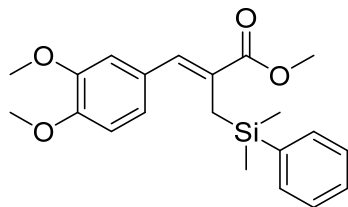
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.56 (s, 1H), 7.41 – 7.38 (m, 2H), 7.36 – 7.28 (m, 3H), 7.03 – 7.00 (m, 2H), 6.92 – 6.89 (m, 1H), 3.70 (s, 3H), 2.32 (s, 3H), 2.26 (s, 2H), 2.19 (s, 3H), 0.21 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 169.2, 138.8, 137.6, 136.6, 135.3, 133.6, 132.6, 130.9, 130.8, 128.9, 128.3, 127.7, 126.2, 51.8, 21.2, 19.9, 17.0, –2.6.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 361.1589; found 361.1594.

Methyl (Z)-3-(3,4-dimethoxyphenyl)-2-((dimethyl(phenyl)silyl)methyl)acrylate 5p



Colorless oil

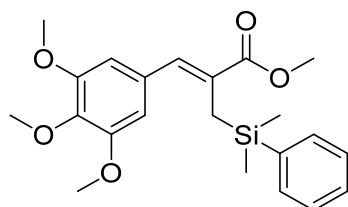
¹H NMR (400 MHz, Chloroform-*d*); δ = 7.50 – 7.46 (m, 3H), 7.36 – 7.29 (m, 3H), 6.93 – 6.90 (m, 1H), 6.82 – 6.76 (m, 2H), 3.88 (s, 3H), 3.72 (s, 3H), 3.69 (s, 3H), 2.44 (s, 2H), 0.28 (s, 6H).

¹³C NMR (100 MHz, Chloroform-*d*); δ = 171.9, 151.3, 151.0, 141.2, 138.0, 136.0, 131.49, 131.46, 131.4, 130.2, 124.9, 114.8, 113.2, 58.3, 58.2, 54.3, 19.9, 0.0.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 393.1491; found 393.1493.

Methyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-3-(3,4,5-trimethoxyphenyl)acrylate 5q

5q



Colorless oil

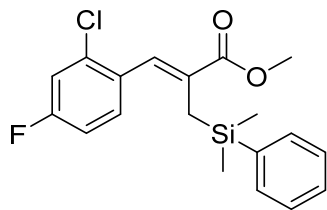
¹H NMR (400 MHz, Chloroform-*d*); δ = 7.47 – 7.44 (m, 3H), 7.33 – 7.28 (m, 3H), 6.50 (s, 2H), 3.84 (s, 3H), 3.70 (s, 3H), 3.69 (s, 6H), 2.45 (s, 2H), 0.29 (s, 6H).

¹³C NMR (100 MHz, Chloroform-*d*); δ = 171.6, 155.3, 141.0, 140.2, 138.0, 135.9, 134.1, 132.6, 131.4, 130.1, 108.8, 63.2, 58.4, 54.3, 19.9, 0.0.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 423.1598; found 423.1598.

Methyl (Z)-3-(2-chloro-4-fluorophenyl)-2-((dimethyl(phenyl)silyl)methyl)acrylate 5r

5r



Colorless oil

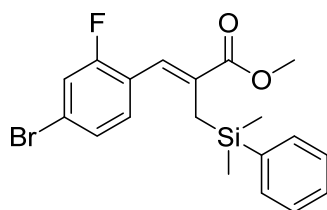
^1H NMR (400 MHz, Chloroform-*d*); δ = 7.48 (s, 1H), 7.41 – 7.28 (m, 5H), 7.12 – 7.06 (m, 2H), 6.85 – 6.80 (m, 1H), 3.71 (s, 3H), 2.23 (s, 2H), 0.22 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 168.6, 163.0, 160.5, 138.2, 134.8, 134.7, 133.5, 133.0, 131.8, 131.1, 131.0, 130.9, 130.8, 129.1, 127.8, 117.1, 116.9, 113.9, 113.7, 52.1, 17.5, – 2.8.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 385.0789; found 385.0797.

Methyl (Z)-3-(4-bromo-2-fluorophenyl)-2-((dimethyl(phenyl)silyl)methyl)acrylate

5s



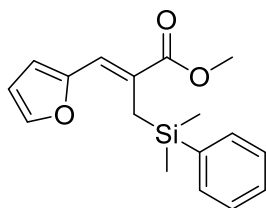
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.40 – 7.37 (m, 3H), 7.34 – 7.28 (m, 3H), 7.20 – 7.17 (m, 1H), 7.12 – 7.09 (m, 1H), 7.02 – 6.96 (m, 1H), 3.71 (s, 3H), 2.27 (s, 2H), 0.24 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 168.5, 161.0, 158.5, 138.1, 133.9, 133.5, 130.73, 130.70, 129.1, 127.8, 127.1, 127.1, 123.4, 123.3, 121.9, 121.8, 119.4, 119.2, 52.1, 18.1, – 2.8.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 429.0288; found 429.0292.

Methyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-3-(furan-2-yl)acrylate 5t



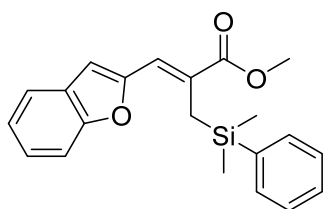
Brown oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.54 – 7.52 (m, 2H), 7.42 – 7.41 (m, 1H), 7.34 – 7.33 (m, 3H), 7.28 (s, 1H), 6.47 – 6.42 (m, 2H), 3.64 (s, 3H), 2.60 (s, 2H), 0.25 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 169.1, 152.3, 143.2, 139.0, 133.6, 128.9, 127.6, 127.2, 122.3, 114.2, 111.8, 51.9, 18.5, – 3.0.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 323.1070; found 323.1074.

Methyl (Z)-3-(benzofuran-2-yl)-2-((dimethyl(phenyl)silyl)methyl)acrylate 5u

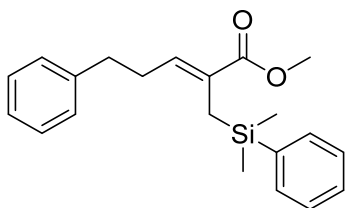


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.44 – 7.39 (m, 3H), 7.25 – 7.22 (m, 2H), 7.20 – 7.14 (m, 4H), 7.11 (m, 1H), 6.64 (s, 1H), 3.55 (s, 3H), 2.64 (s, 2H), 0.15 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 168.8, 155.1, 154.0, 138.8, 133.6, 130.9, 129.0, 128.1, 127.7, 125.5, 123.2, 122.2, 121.3, 111.2, 110.5, 52.1, 19.3, -2.9.

Methyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-5-phenylpent-2-enoate 5v



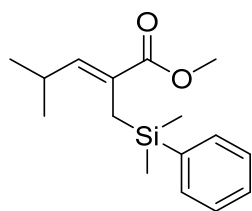
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.48 – 7.46 (m, 2H), 7.33 – 7.29 (m, 3H), 7.24

– 7.21 (m, 2H), 7.18 – 7.14 (m, 1H), 7.01 – 7.05 (m, 2H), 6.63 (t, $J = 7.2$ Hz, 1H), 3.59 (s, 3H), 2.60 (t, $J = 7.8$ Hz, 2H), 2.23 (dd, $J = 15.6, 7.6$ Hz, 2H), 1.98 (s, 2H), 0.25 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); $\delta = 168.6, 141.3, 138.7, 138.2, 133.6, 129.8, 129.1, 128.4, 128.3, 127.8, 127.7, 126.0, 51.6, 34.8, 30.8, 17.2, -2.8$.

Methyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-4-methylpent-2-enoate 5w



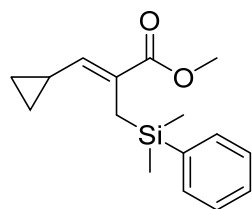
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); $\delta = 7.52 - 7.50$ (m, 2H), 7.35 – 7.34 (m, 3H), 6.41 (d, $J = 10.2$ Hz, 1H), 3.61 (s, 3H), 2.41 – 2.30 (m, 1H), 2.01 (s, 2H), 0.88 (d, $J = 6.6$ Hz, 6H), 0.28 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); $\delta = 169.0, 146.1, 138.8, 133.6, 129.0, 127.7, 126.8, 51.5, 28.1, 21.9, 16.7, -2.9$.

HRMS (ESI⁺) calcd for $\text{C}_{23}\text{H}_{23}\text{BrOSiNa}$ [$\text{M}+\text{Na}$]⁺: 299.1431; found 299.1438.

Methyl (Z)-3-cyclopropyl-2-((dimethyl(phenyl)silyl)methyl)acrylate 5x



Colorless oil

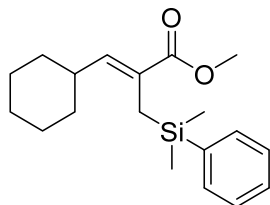
^1H NMR (400 MHz, Chloroform-*d*); $\delta = 7.55 - 7.53$ (m, 2H), 7.36 – 7.34 (m, 3H), 6.04 (d, $J = 10.6$ Hz, 1H), 3.60 (s, 3H), 2.14 (s, 2H), 1.40 – 1.31 (m, 1H), 0.81 – 0.79 (m, 2H), 0.53 – 0.49 (m, 2H), 0.32 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); $\delta = 168.5, 144.5, 139.0, 133.7, 129.0, 127.7,$

126.7, 51.5, 16.7, 12.3, 8.5, – 2.9.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 297.1279; found 297.1281.

Methyl (Z)-3-cyclohexyl-2-((dimethyl(phenyl)silyl)methyl)acrylate 5y



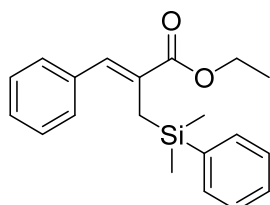
Colorless oil

¹H NMR (400 MHz, Chloroform-*d*); δ = 7.53 – 7.50 (m, 2H), 7.36 – 7.33 (m, 3H), 6.42 (d, *J* = 10.1 Hz, 1H), 3.62 (s, 3H), 2.05 – 1.93 (m, 3H), 1.66 – 1.60 (m, 4H), 1.28 – 1.23 (m, 1H), 1.16 – 0.96 (m, 5H), 0.28 (s, 6H).

¹³C NMR (100 MHz, Chloroform-*d*); δ = 169.1, 144.6, 138.8, 133.6, 129.0, 127.7, 127.0, 51.5, 38.0, 31.9, 25.9, 25.6, 16.8, – 2.9.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 339.1750; found 339.1751.

Ethyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-3-phenylacrylate 5z



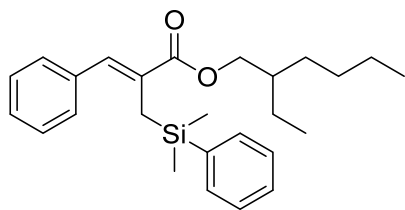
Colorless oil

¹H NMR (400 MHz, Chloroform-*d*); δ = 7.41 (s, 1H), 7.33 – 7.31 (m, 2H), 7.21 – 7.18 (m, 3H), 7.15 – 7.10 (m, 5H), 4.05 (dd, *J* = 14.2, 7.1 Hz, 2H), 2.26 (s, 2H), 1.17 – 1.12 (m, 3H), 0.12 (s, 6H).

¹³C NMR (100 MHz, Chloroform-*d*); δ = 171.3, 141.2, 138.9, 137.9, 136.1, 133.7, 131.6, 131.5, 130.8, 130.2, 63.4, 27.5, 19.8, 16.7, 0.0.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 347.1440; found 347.1438.

2-Ethylhexyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-3-phenylacrylate 5A



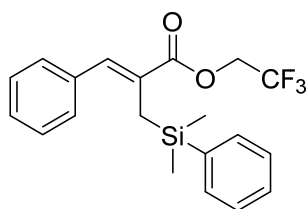
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.43 (s, 1H), 7.37 – 7.35 (m, 2H), 7.24 – 7.13 (m, 8H), 3.95 – 3.93 (m, 2H), 2.31 (s, 2H), 1.54 – 1.51 (m, 1H), 1.32 – 1.20 (m, 8H), 0.84 – 0.80 (m, 6H), 0.15 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 169.0, 138.7, 136.4, 135.2, 133.6, 131.4, 129.11, 129.06, 128.3, 127.8, 67.4, 38.8, 30.6, 29.0, 23.9, 23.1, 17.4, 14.2, 11.1, – 2.5.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 431.2382; found 431.2377.

2,2,2-Trifluoroethyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-3-phenylacrylate 5B



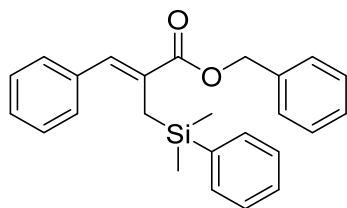
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.61 (s, 1H), 7.47 – 7.44 (m, 2H), 7.38 – 7.29 (m, 8H), 4.47 (dd, *J* = 17.2, 8.5 Hz, 2H), 2.43 (s, 2H), 0.26 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 167.1, 138.2, 137.6, 135.7, 133.6, 129.4, 129.22, 129.19, 128.4, 128.3, 127.8, 124.5, 121.8, 61.3, 60.9, 60.6, 60.2, 17.5, – 2.7.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 401.1150; found 401.1155.

Benzyl (Z)-2-((dimethyl(phenyl)silyl)methyl)-3-phenylacrylate 5C



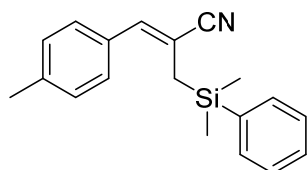
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.56 (s, 1H), 7.41 – 7.32 (m, 7H), 7.31 – 7.26 (m, 3H), 7.24 – 7.20 (m, 5H), 5.11 (s, 2H), 2.39 (s, 2H), 0.18 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 168.7, 138.6, 136.3, 136.1, 135.9, 133.6, 131.0, 129.14, 129.08, 128.6, 128.36, 128.35, 128.2, 127.9, 127.8, 66.7, 17.4, – 2.5.

HRMS (ESI⁺) calcd for C₂₃H₂₃BrOSiNa [M+Na]⁺: 409.1593; found 409.1594.

(Z)-2-((dimethyl(phenyl)silyl)methyl)-3-(*p*-tolyl)acrylonitrile 5D



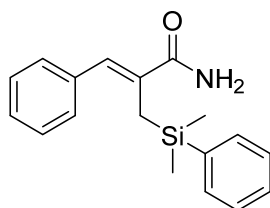
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.54 – 7.48 (m, 4H), 7.41 – 7.38 (m, 3H), 7.19 – 7.04 (m, 2H), 6.54 (s, 1H), 2.36 (s, 3H), 2.06 (s, 2H), 0.46 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 141.8, 139.6, 136.9, 133.7, 131.5, 129.6, 129.4, 128.1, 128.0, 119.9, 107.0, 26.2, 21.4, – 3.4

HRMS (ESI⁺) calcd for C₁₉H₂₂NSi [M+H]⁺: 292.1522; found 292.1523.

(Z)-2-((dimethyl(phenyl)silyl)methyl)-3-phenylacrylamide 5E



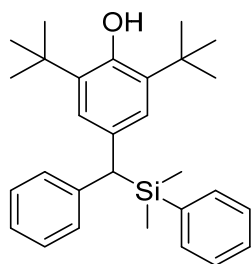
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.41 – 7.38 (m, 2H), 7.29 – 7.11 (m, 8H), 7.00 (s, 1H), 5.56 (s, 2H), 2.23 (s, 2H), 0.23 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 175.1, 140.8, 138.6, 137.7, 136.1, 133.1, 131.8, 131.3, 130.9, 130.4, 130.1, 20.5, 0.0.

HRMS (ESI⁺) calcd for C₁₈H₂₁NOSiNa [M+Na]⁺: 318.1290; found 318.1291.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(phenyl)methyl)phenol 7a

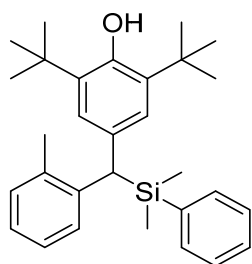


Yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.35 – 7.28 (m, 2H), 7.25 – 7.19 (m, 5H), 7.16 – 7.08 (m, 3H), 6.86 (s, 2H), 4.93 (s, 1H), 3.62 (s, 1H), 1.33 (s, 18H), 0.27 (s, 3H), 0.25 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 151.3, 143.1, 137.9, 135.2, 134.5, 132.3, 129.0, 128.9, 128.2, 127.5, 125.4, 125.0, 45.5, 34.3, 30.4, – 3.0, – 3.4.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(*o*-tolyl)methyl)phenol 7b

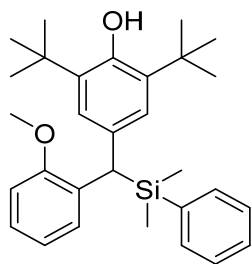


Yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.38 – 7.31 (m, 1H), 7.30 – 7.28 (m, 5H), 7.15 – 7.03 (m, 3H), 6.82 (s, 2H), 4.93 (s, 1H), 3.84 (s, 1H), 2.27 (s, 3H), 1.33 (s, 18H), 0.33 (s, 3H), 0.29 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 151.1, 141.5, 138.4, 136.6, 135.1, 134.5, 131.9, 130.5, 129.4, 128.9, 127.5, 125.5, 125.4, 125.2, 40.1, 34.3, 30.3, 20.6, – 2.5, – 3.5.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(2-methoxyphenyl)methyl)phenol 7c

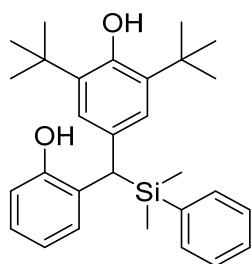


Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.36 – 7.32 (m, 1H), 7.31 – 7.28 (m, 4H), 7.24 – 7.21 (m, 1H), 7.15 – 7.11 (m, 1H), 6.92 (s, 2H), 6.89 – 6.81 (m, 2H), 4.93 (s, 1H), 4.20 (s, 1H), 3.72 (s, 3H), 1.36 (s, 18H), 0.28 (s, 3H), 0.25 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 156.7, 151.1, 139.1, 135.0, 134.4, 132.5, 131.7, 130.2, 128.7, 127.3, 126.0, 125.7, 120.2, 110.4, 55.1, 36.6, 34.3, 30.4, – 2.8, – 3.4.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(2-hydroxyphenyl)methyl)phenol 7d



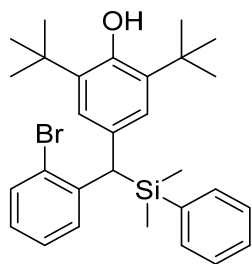
Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.39 – 7.35 (m, 1H), 7.34 – 7.28 (m, 5H), 7.10 – 7.06 (m, 1H), 6.90 – 6.86 (m, 1H), 6.84 (s, 2H), 6.80 – 6.77 (m, 1H), 5.00 (s, 1H), 4.81 (s, 1H), 3.95 (s, 1H), 1.33 (s, 18H), 0.38 (s, 3H), 0.35 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 153.8, 151.6, 138.2, 135.5, 134.4, 131.1, 130.6, 129.2, 129.0, 127.6, 126.7, 125.5, 120.5, 116.1, 37.0, 34.3, 30.3, – 2.6, – 3.5.

HRMS (ESI⁺) calcd for C₂₉H₃₉O₂Si [M+H]⁺: 447.1719; found 447.2716.

4-((2-Bromophenyl)(dimethyl(phenyl)silyl)methyl)-2,6-di-*tert*-butylphenol 7e

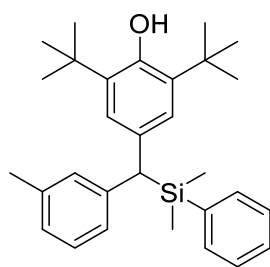


Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.55 – 7.53 (m, 1H), 7.37 – 7.27 (m, 6H), 7.20 – 7.16 (m, 1H), 6.99 – 6.95 (m, 1H), 6.87 (s, 2H), 4.95 (s, 1H), 4.35 (s, 1H), 1.32 (s, 18H), 0.32 (s, 3H), 0.26 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 151.4, 142.6, 137.7, 135.2, 134.6, 133.0, 131.3, 130.4, 129.1, 127.6, 127.0, 126.6, 126.1, 125.6, 43.0, 34.3, 30.3, – 2.7, – 3.8.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(*m*-tolyl)methyl)phenol 7f

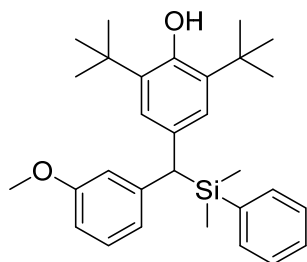


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.33 – 7.29 (m, 1H), 7.25 – 7.20 (m, 4H), 7.10 – 7.08 (m, 1H), 6.95 – 6.89 (m, 3H), 6.82 (s, 2H), 4.89 (s, 1H), 3.53 (s, 1H), 2.23 (s, 3H), 1.30 (s, 18H), 0.24 (s, 3H), 0.21 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 151.3, 142.8, 138.0, 137.5, 135.2, 134.5, 132.3, 130.0, 128.9, 127.9, 127.4, 125.8, 125.7, 125.3, 45.4, 34.3, 30.3, 21.6, – 3.1, – 3.5.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(3-methoxyphenyl)methyl)phenol 7g

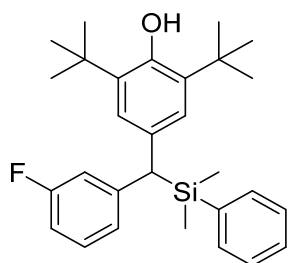


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.39 – 7.34 (m, 1H), 7.33 – 7.28 (m, 4H), 7.18 – 7.14 (m, 1H), 6.91 (s, 2H), 6.80 – 6.78 (m, 1H), 6.72 – 6.69 (m, 2H), 4.98 (s, 1H), 3.70 (s, 3H), 3.64 (s, 1H), 1.38 (s, 18H), 0.31 (s, 3H), 0.30 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 159.3, 151.4, 144.6, 137.9, 135.3, 134.6, 132.1, 129.02, 129.00, 127.5, 125.4, 121.4, 114.1, 111.0, 55.0, 45.7, 34.4, 30.4, – 3.1, – 3.3.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(3-fluorophenyl)methyl)phenol 7h



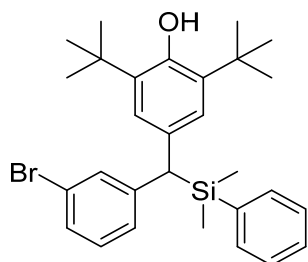
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.33 – 7.29 (m, 1H), 7.25 – 7.19 (m, 4H), 7.15 – 7.10 (m, 1H), 6.88 – 6.85 (m, 1H), 6.81 – 6.74 (m, 4H), 4.94 (s, 1H), 3.58 (s, 1H), 1.30 (s, 18H), 0.24 (s, 3H), 0.22 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 161.5, 151.5, 145.9, 145.8, 137.4, 135.4, 134.4, 131.6, 129.4, 129.3, 129.1, 127.6, 125.5, 124.5, 124.4, 115.6, 115.4, 111.9, 111.7, 45.5, 45.4, 34.3, 30.3, – 3.2, – 3.4.

HRMS (ESI⁺) calcd for C₂₉H₃₈FOSi [M+H]⁺: 449.2676; found 449.2672.

4-((3-Bromophenyl)(dimethyl(phenyl)silyl)methyl)-2,6-di-*tert*-butylphenol 3i



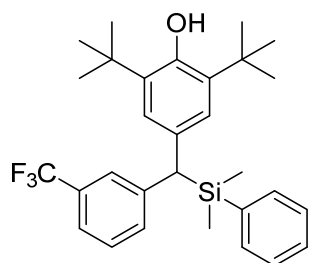
Yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.33 – 7.29 (m, 1H), 7.25 – 7.24 (m, 1H), 7.22 – 7.17 (m, 5H), 7.05 – 6.99 (m, 2H), 6.78 (s, 2H), 4.88 (s, 1H), 3.56 (s, 1H), 1.23 (s, 18H), 0.16 (s, 3H), 0.15 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 151.5, 145.5, 137.2, 135.4, 134.5, 131.9, 131.5, 129.6, 129.2, 128.0, 127.6, 127.2, 125.4, 122.2, 45.3, 34.3, 30.3, – 3.2, – 3.5.

HRMS (ESI⁺) calcd for C₂₉H₃₈BrOSi [M+H]⁺: 509.1875; found 509.1871.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(3-(trifluoromethyl)phenyl)methyl)phenol 3j

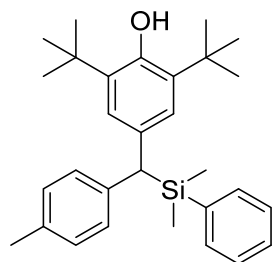


Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.27 – 7.22 (m, 3H), 7.20 – 7.15 (m, 4H), 7.12 – 7.09 (m, 2H), 6.74 (s, 2H), 4.93 (s, 1H), 3.52 (s, 1H), 1.29 (s, 18H), 0.23 (s, 3H), 0.21 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 151.6, 144.0, 137.0, 135.5, 134.5, 132.0, 131.3, 129.3, 128.5, 127.6, 125.5, 125.4, 125.3, 121.79, 121.75, 45.6, 34.4, 30.3, – 3.4, – 3.5.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(*p*-tolyl)methyl)phenol 7k



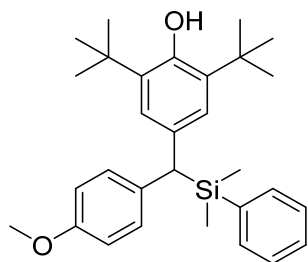
Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.33 – 7.28 (m, 1H), 7.25 – 7.21 (m, 4H), 7.03

– 6.99 (m, 4H), 6.81 (s, 2H), 4.89 (s, 1H), 3.54 (s, 1H), 2.26 (s, 3H), 1.30 (s, 18H), 0.23 (s, 3H), 0.21 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 151.2, 139.9, 138.1, 135.1, 134.5, 134.3, 132.5, 128.9, 128.84, 128.79, 127.4, 125.3, 45.0, 34.3, 30.3, 21.0, – 3.0, – 3.4.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(4-methoxyphenyl)methyl)phenol 7l



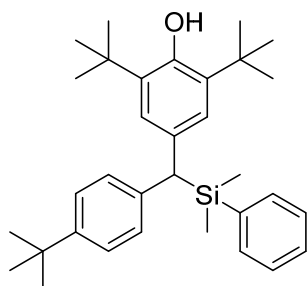
Yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.35 – 7.31 (m, 1H), 7.28 – 7.23 (m, 4H), 7.08 – 7.05 (m, 2H), 6.83 (s, 2H), 6.79 – 6.76 (m, 2H), 4.92 (s, 1H), 3.77 (s, 3H), 3.56 (s, 1H), 1.33 (s, 18H), 0.26 (s, 3H), 0.24 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 157.1, 151.2, 138.1, 135.2, 135.2, 134.5, 132.7, 129.9, 128.9, 127.5, 125.2, 113.6, 55.2, 44.3, 34.3, 30.3, – 3.0, – 3.4.

2,6-Di-*tert*-butyl-4-((4-(*tert*-butyl)phenyl)(dimethyl(phenyl)silyl)methyl)phenol 7m

7m



Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.35 – 7.29 (m, 1H), 7.25 – 7.20 (m, 6H), 7.08 – 7.06 (m, 2H), 6.84 (s, 2H), 4.90 (s, 1H), 3.55 (s, 1H), 1.32 (s, 18H), 1.27 (s, 9H), 0.26 (s, 3H), 0.23 (s, 3H).

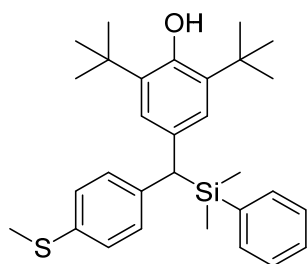
^{13}C NMR (100 MHz, Chloroform-*d*); δ = 151.3, 147.6, 139.8, 138.1, 135.1, 134.5,

132.5, 128.9, 128.4, 127.4, 125.4, 125.0, 45.2, 34.30, 34.28, 31.5, 30.4, – 3.1, – 3.3.

HRMS (ESI⁺) calcd for C₃₃H₄₇OSi [M+H]⁺: 487.3396; found 487.3396.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(4-(methylthio)phenyl)methyl)phenol

7n



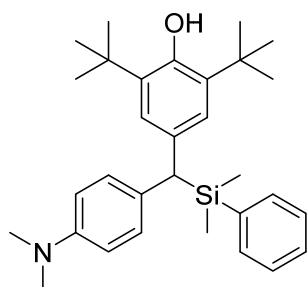
Pale yellow oil

¹H NMR (400 MHz, Chloroform-*d*); δ = 7.38 – 7.29 (m, 5H), 7.19 – 7.10 (m, 4H), 6.88 (s, 2H), 4.99 (s, 1H), 3.62 (s, 1H), 2.49 (s, 3H), 1.38 (s, 18H), 0.31 (s, 6H).

¹³C NMR (100 MHz, Chloroform-*d*); δ = 151.4, 140.3, 137.8, 135.3, 134.5, 134.1, 132.1, 129.4, 129.0, 127.5, 126.9, 125.4, 45.0, 34.3, 30.3, 16.3, – 3.1, – 3.4.

HRMS (ESI⁺) calcd for C₃₀H₄₁OSSi [M+H]⁺: 477.2647; found 477.2652.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(4-(dimethylamino)phenyl)methyl)phenol 7o



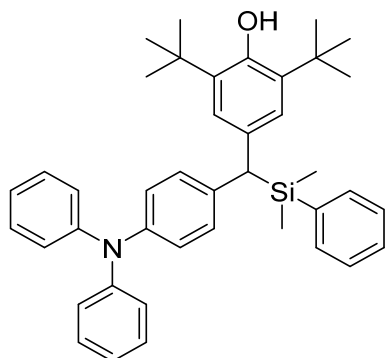
Orange oil

¹H NMR (400 MHz, Chloroform-*d*); δ = 7.34 – 7.31 (m, 1H), 7.26 – 7.25 (m, 4H), 7.05 – 7.01 (m, 2H), 6.82 (s, 2H), 6.67 – 6.63 (m, 2H), 4.89 (s, 1H), 3.50 (s, 1H), 2.89 (s, 6H), 1.31 (s, 18H), 0.25 (s, 3H), 0.23 (s, 3H).

¹³C NMR (100 MHz, Chloroform-*d*); δ = 151.1, 148.4, 138.4, 135.0, 134.6, 133.0, 131.4, 129.6, 128.8, 127.4, 125.2, 113.0, 44.1, 41.0, 34.3, 30.3, – 2.9, – 3.4.

HRMS (ESI⁺) calcd for C₃₁H₄₄NOSi [M+H]⁺: 474.3192; found 474.3189.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(4-(diphenylamino)phenyl)methyl)phenol 7p



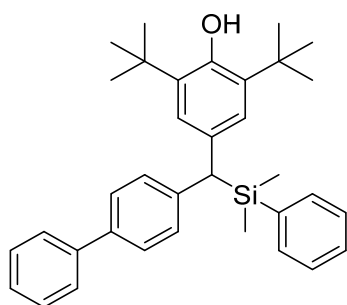
Orange oil

¹H NMR (400 MHz, Chloroform-*d*); δ = 7.35 – 7.29 (m, 2H), 7.25 – 7.19 (m, 7H), 7.06 – 7.02 (m, 6H), 6.99 – 6.94 (m, 4H), 6.86 (s, 2H), 4.95 (s, 1H), 3.57 (s, 1H), 1.34 (s, 18H), 0.29 (s, 3H), 0.26 (s, 3H).

¹³C NMR (100 MHz, Chloroform-*d*); δ = 151.3, 148.0, 137.9, 137.8, 135.2, 134.5, 132.4, 129.6, 129.1, 128.9, 127.4, 125.3, 124.7, 123.6, 122.1, 44.9, 34.3, 30.3, – 3.2, – 3.4.

HRMS (ESI⁺) calcd for C₄₁H₄₈NOSi [M+H]⁺: 598.3505; found 598.3508.

4-([1,1'-Biphenyl]-4-yl(dimethyl(phenyl)silyl)methyl)-2,6-di-*tert*-butylphenol 7q



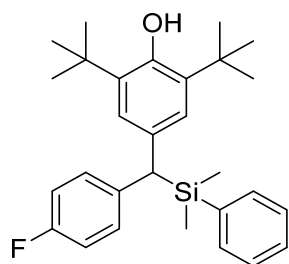
Yellow oil

¹H NMR (400 MHz, Chloroform-*d*); δ = 7.54 – 7.51 (m, 2H), 7.42 – 7.33 (m, 4H), 7.31 – 7.27 (m, 1H), 7.25 – 7.13 (m, 7H), 6.83 (s, 2H), 4.89 (s, 1H), 3.60 (s, 1H), 1.28 (s, 18H), 0.24 (s, 3H), 0.22 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 151.4, 142.3, 141.1, 137.8, 137.6, 135.3, 134.6, 132.2, 129.2, 129.0, 128.7, 127.5, 126.9, 126.9, 126.8, 125.5, 45.3, 34.3, 30.4, – 3.1, – 3.3.

HRMS (ESI⁺) calcd for C₃₅H₄₂OSiNa [M+Na]⁺: 529.2903; found 529.2904.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(4-fluorophenyl)methyl)phenol 7r



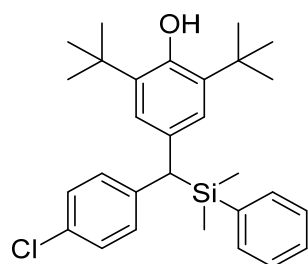
Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.30 – 7.26 (m, 1H), 7.23 – 7.16 (m, 4H), 7.04 – 7.00 (m, 2H), 6.86 – 6.82 (m, 2H), 6.77 (s, 2H), 4.90 (s, 1H), 3.54 (s, 1H), 1.28 (s, 18H), 0.20 (s, 3H), 0.19 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 159.6, 151.4, 138.8, 138.7, 137.6, 135.4, 134.5, 132.2, 130.2, 130.1, 129.1, 127.6, 125.3, 115.0, 114.8, 44.5, 34.4, 30.3, – 3.1, – 3.4.

HRMS (ESI⁺) calcd for C₂₉H₃₈FOSi [M+H]⁺: 449.2676; found 449.2674.

2,6-Di-*tert*-butyl-4-((4-chlorophenyl)(dimethyl(phenyl)silyl)methyl)phenol 7s



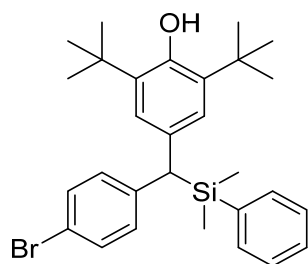
Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.26 – 7.22 (m, 1H), 7.19 – 7.11 (m, 4H), 7.08 – 7.05 (m, 2H), 6.94 – 6.92 (m, 2H), 6.72 (s, 2H), 4.56 (s, 1H), 3.48 (s, 1H), 1.23 (s, 18H), 0.15 (s, 3H), 0.14 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 151.5, 141.7, 137.4, 135.4, 134.5, 131.8, 130.6, 130.1, 129.1, 128.2, 127.6, 125.4, 44.9, 34.3, 30.3, - 3.2, - 3.4.

HRMS (ESI⁺) calcd for C₂₉H₃₈ClOSi [M+H]⁺: 465.2380; found 465.2379.

4-((4-Bromophenyl)(dimethyl(phenyl)silyl)methyl)-2,6-di-*tert*-butylphenol 7t



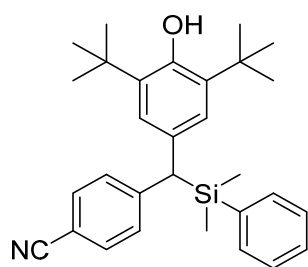
Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.36 – 7.32 (m, 2H), 7.31 – 7.21 (m, 5H), 6.99 – 6.97 (m, 2H), 6.82 (s, 2H), 4.96 (s, 1H), 3.87 (s, 1H), 1.32 (s, 18H), 0.25 (s, 3H), 0.24 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 151.5, 142.2, 137.4, 135.4, 134.5, 131.7, 131.1, 130.5, 129.1, 127.6, 125.4, 118.6, 44.9, 34.3, 30.3, - 3.2, - 3.4.

4-((3,5-Di-*tert*-butyl-4-hydroxyphenyl)(dimethyl(phenyl)silyl)methyl)benzonitrile 7u

7u

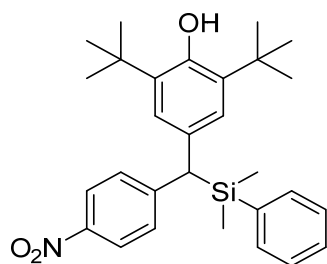


Yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.78 – 7.76 (m, 2H), 7.68 – 7.64 (m, 1H), 7.58 – 7.55 (m, 2H), 7.52 – 7.50 (m, 2H), 7.47 – 7.45 (m, 2H), 7.15 (s, 2H), 5.33 (s, 1H), 4.00 (s, 1H), 1.64 (s, 18H), 0.57 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 151.8, 149.3, 136.7, 135.7, 134.4, 131.9, 130.8, 129.4, 129.1, 127.7, 125.7, 119.4, 108.4, 46.4, 34.4, 30.3, - 3.2, - 3.5.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(4-nitrophenyl)methyl)phenol 7v



Pale yellow oil

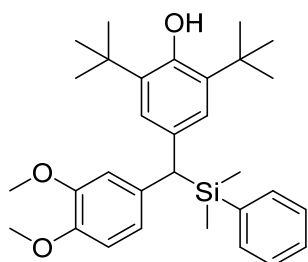
^1H NMR (400 MHz, Chloroform-*d*); δ = 8.07 – 8.04 (m, 2H), 7.39 – 7.35 (m, 1H), 7.31 – 7.27 (m, 2H), 7.23 – 7.18 (m, 4H), 6.89 (s, 2H), 5.04 (s, 1H), 3.77 (s, 1H), 1.35 (s, 18H), 0.29 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 151.9, 151.7, 145.4, 136.6, 135.8, 134.4, 129.5, 128.9, 127.7, 125.7, 123.4, 46.3, 34.4, 30.3, – 3.2, – 3.5.

HRMS (ESI⁺) calcd for C₂₉H₃₈NO₃Si [M+H]⁺: 476.2621; found 476.2622.

2,6-Di-*tert*-butyl-4-((3,4-dimethoxyphenyl)(dimethyl(phenyl)silyl)methyl)phenol 7w

7w



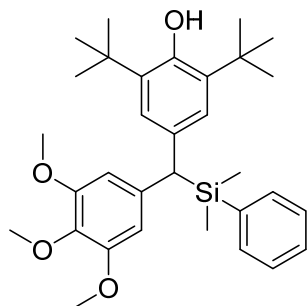
Pale yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.32 – 7.29 (m, 1H), 7.25 – 7.24 (m, 1H), 7.22 – 7.20 (m, 3H), 6.83 (s, 2H), 6.72 – 6.65 (m, 2H), 6.56 – 6.55 (m, 1H), 4.91 (s, 1H), 3.80 (s, 3H), 3.62 (s, 3H), 3.50 (s, 1H), 1.30 (s, 18H), 0.21 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 151.3, 148.2, 146.4, 137.9, 135.5, 135.2, 134.7, 132.4, 129.0, 127.5, 125.1, 120.7, 112.2, 110.9, 55.8, 55.4, 44.9, 34.3, 30.3, – 3.1, – 3.5.

HRMS (ESI⁺) calcd for C₃₁H₄₃O₃Si [M+H]⁺: 491.2981; found 491.2976.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(3,4,5-trimethoxyphenyl)methyl)phenol 7x



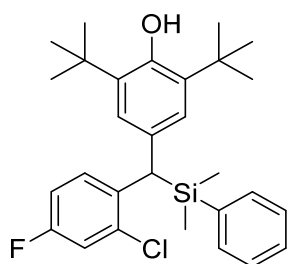
Yellow oil

¹H NMR (400 MHz, Chloroform-*d*); δ = 7.29 – 7.21 (m, 2H), 7.19 – 7.16 (m, 3H), 6.82 (s, 2H), 6.25 (m, 2H), 4.89 (s, 1H), 3.71 (s, 3H), 3.60 (s, 6H), 3.44 (s, 1H), 1.28 (s, 18H), 0.19 (s, 3H), 0.18 (s, 3H).

¹³C NMR (100 MHz, Chloroform-*d*); δ = 152.6, 151.4, 138.6, 137.7, 135.3, 134.8, 132.0, 129.1, 127.5, 125.1, 105.7, 60.9, 55.7, 46.0, 34.4, 30.3, – 3.2, – 3.4.

HRMS (ESI⁺) calcd for C₃₂H₄₅O₄Si [M+H]⁺: 521.3087; found 521.3087.

2,6-Di-*tert*-butyl-4-((2-chloro-4-fluorophenyl)(dimethyl(phenyl)silyl)methyl)phenol 7y



Colorless oil

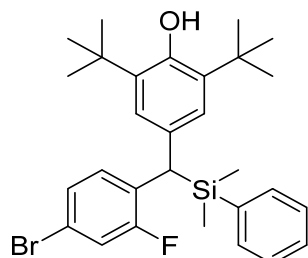
¹H NMR (400 MHz, Chloroform-*d*); δ = 7.43 – 7.40 (m, 1H), 7.38 – 7.35 (m, 4H), 7.30 – 7.26 (m, 1H), 7.18 – 7.15 (m, 1H), 6.96 – 6.91 (m, 1H), 6.89 (s, 2H), 5.04 (s, 1H), 4.33 (s, 1H), 1.39 (s, 18H), 0.37 (s, 3H), 0.33 (s, 3H).

¹³C NMR (100 MHz, Chloroform-*d*); δ = 161.3, 158.9, 151.5, 137.4, 136.9, 136.9, 135.3, 134.8, 134.5, 131.2, 131.1, 131.0, 129.2, 127.6, 125.5, 116.9, 116.7, 113.7, 113.5,

39.3, 34.3, 30.3, - 2.8, - 3.7.

HRMS (ESI⁺) calcd for C₂₉H₃₇FCIOSi [M+H]⁺: 483.2286; found 483.2289.

4-((4-Bromo-2-fluorophenyl)(dimethyl(phenyl)silyl)methyl)-2,6-di-*tert*-butylphenol 7z



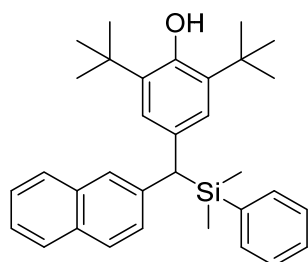
Yellow oil

¹H NMR (400 MHz, Chloroform-*d*); δ = 7.37 – 7.32 (m, 1H), 7.29 – 7.28 (m, 4H), 7.15 – 7.11 (m, 3H), 6.92 (s, 2H), 4.94 (s, 1H), 4.39 (s, 1H), 1.35 (s, 18H), 0.21 (s, 6H).

¹³C NMR (100 MHz, Chloroform-*d*); δ = 154.4, 151.2, 138.1, 138.0, 135.2, 134.4, 132.0, 131.2, 128.9, 127.4, 126.1, 125.1, 123.8, 118.8, 36.7, 34.3, 30.3, 26.8, 24.3, - 3.0, - 4.0.

HRMS (ESI⁺) calcd for C₂₉H₃₇FBrOSi [M+H]⁺: 527.1781; found 527.1777.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(naphthalen-2-yl)methyl)phenol 7A

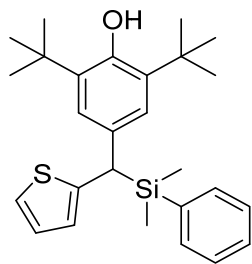


Yellow oil

¹H NMR (400 MHz, Chloroform-*d*); δ = 7.80 – 7.78 (m, 1H), 7.73 – 7.69 (m, 2H), 7.60 (s, 1H), 7.45 – 7.36 (m, 3H), 7.35 – 7.29 (m, 5H), 6.95 (s, 2H), 4.97 (s, 1H), 3.82 (s, 1H), 1.36 (s, 18H), 0.32 (s, 3H), 0.30 (s, 3H).

¹³C NMR (100 MHz, Chloroform-*d*); δ = 151.4, 140.7, 137.9, 135.3, 134.6, 133.6, 132.2, 131.5, 129.1, 128.3, 127.54, 127.52, 127.48, 126.8, 125.7, 125.5, 124.8, 45.6, 34.3, 30.3, - 3.0, - 3.3.

2,6-Di-*tert*-butyl-4-((dimethyl(phenyl)silyl)(thiophen-2-yl)methyl)phenol 7B

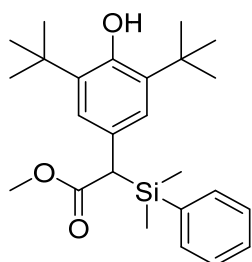


Yellow oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.22 – 7.18 (m, 1H), 7.13 – 7.08 (m, 4H), 6.90 – 6.89 (m, 1H), 6.76 -6.74 (m, 1H), 6.70 (s, 2H), 6.60 (s, 1H), 4.81 (s, 1H), 3.74 (s, 1H), 1.20 (s, 18H), 0.15 (s, 3H), 0.14 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 151.4, 146.0, 137.0, 135.2, 134.5, 131.5, 129.1, 127.4, 126.6, 124.8, 124.1, 122.1, 40.2, 34.3, 30.3, – 3.7, – 4.1.

Methyl 2-(3,5-di-*tert*-butyl-4-hydroxyphenyl)-2-(dimethyl(phenyl)silyl)acetate 7C

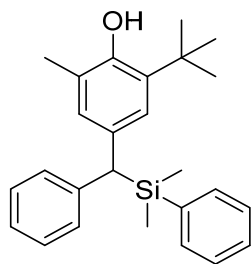


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.38 – 7.33 (m, 1H), 7.32 – 7.28 (m, 4H), 6.93 (s, 2H), 4.99 (s, 1H), 3.59 (s, 3H), 3.49 (s, 1H), 1.34 (s, 18H), 0.34 (s, 3H), 0.33 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 173.7, 151.7, 135.8, 135.2, 134.1, 129.4, 127.6, 125.8, 125.0, 51.3, 45.3, 34.2, 30.3, – 4.3, – 4.4.

2-(*tert*-Butyl)-4-((dimethyl(phenyl)silyl)(phenyl)methyl)-6-methylphenol 7D

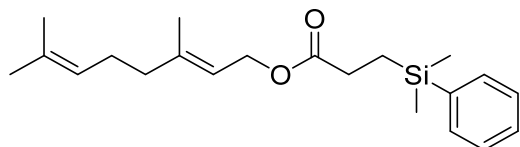


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.37 – 7.31 (m, 1H), 7.30 – 7.28 (m, 4H), 7.23 – 7.19 (m, 2H), 7.14 – 7.09 (m, 3H), 6.85 (s, 1H), 6.75 (s, 1H), 4.54 (s, 1H), 3.62 (s, 1H), 2.13 (s, 3H), 1.31 (s, 9H), 0.30 (s, 3H), 0.28 (s, 3H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 150.3, 143.0, 137.9, 135.0, 134.6, 133.2, 129.1, 129.0, 128.8, 128.2, 127.5, 126.0, 125.0, 122.7, 45.0, 34.6, 29.8, 16.2, – 3.1, – 3.2.

(*E*)-3,7-dimethylocta-2,6-dien-1-yl 3-(dimethyl(phenyl)-14-sulfanyl)propanoate 9a



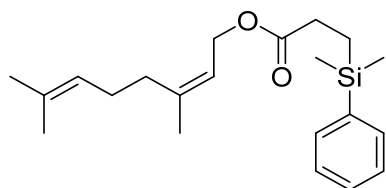
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.51 – 7.49 (m, 2H), 7.37 – 7.36 (m, 3H), 5.34 – 5.30 (m, 1H), 5.10 – 5.07 (m, 1H), 4.57 – 4.55 (m, 2H), 2.31 – 2.27 (m, 2H), 2.09 – 2.04 (m, 4H), 1.70 – 1.68 (m, 6H), 1.60 (s, 3H), 1.12 – 1.08 (m, 2H), 0.29 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 175.0, 142.2, 138.2, 133.6, 131.8, 129.1, 127.9, 123.8, 118.3, 61.4, 39.6, 28.9, 26.3, 25.7, 17.7, 16.5, 10.9, – 3.3.

HRMS (ESI⁺) calcd for C₂₁H₃₃O₂Si [M+H]⁺: 345.2250; found 345.2248.

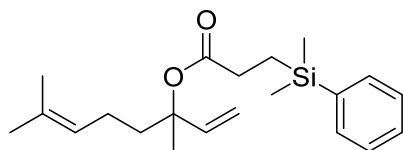
(*Z*)-3,7-dimethylocta-2,6-dien-1-yl 3-(dimethyl(phenyl)silyl)propanoate 9b



Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.51 – 7.49 (m, 2H), 7.37 – 7.33 (m, 3H), 5.35 – 5.32 (m, 1H), 5.11 – 5.07 (m, 1H), 4.54 – 4.52 (m, 2H), 2.30 – 2.26 (m, 2H), 2.21 – 2.06 (m, 4H), 1.76 (s, 3H), 1.68 (s, 3H), 1.60 (s, 3H), 1.12 – 1.07 (m, 2H), 0.29 (s, 6H).
 ^{13}C NMR (100 MHz, Chloroform-*d*); δ = 174.9, 142.6, 138.2, 133.6, 132.2, 129.1, 127.9, 123.6, 119.2, 61.1, 32.2, 28.9, 26.7, 25.7, 23.6, 17.7, 10.9, – 3.3.
HRMS (ESI⁺) calcd for C₂₁H₃₂O₂SiNa [M+Na]⁺: 367.2069; found 367.2064.

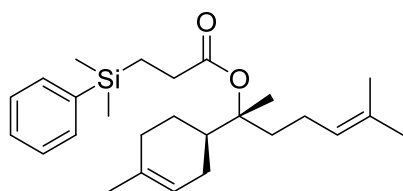
3,7-Dimethylocta-1,6-dien-3-yl 3-(dimethyl(phenyl)silyl)propanoate 9c



Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.52 – 7.49 (m, 2H), 7.37 – 7.26 (m, 3H), 5.98 – 5.91 (m, 1H), 5.16 – 5.06 (m, 2H), 2.25 – 2.21 (m, 2H), 1.99 – 1.93 (m, 2H), 1.87 – 1.72 (m, 2H), 1.67 (s, 3H), 1.58 (s, 3H), 1.52 (s, 3H), 1.09 – 1.04 (m, 2H), 0.28 (s, 6H).
 ^{13}C NMR (100 MHz, Chloroform-*d*); δ = 173.7, 141.9, 138.3, 133.6, 131.8, 129.1, 127.9, 123.9, 113.1, 82.7, 39.8, 29.8, 25.7, 23.6, 22.4, 17.6, 10.8, – 3.2.
HRMS (ESI⁺) calcd for C₂₁H₃₂O₂SiNa [M+Na]⁺: 367.2069; found 367.2073.

(*R*)-6-methyl-2-((*R*)-4-methylcyclohex-3-en-1-yl)hept-5-en-2-yl (dimethyl(phenyl)silyl)propanoate 9d



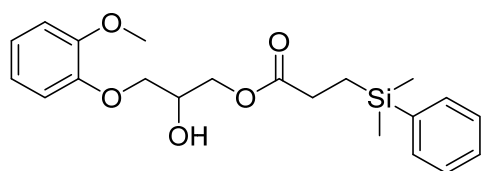
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.52 – 7.49 (m, 2H), 7.37 – 7.26 (m, 3H), 5.35 – 5.31 (m, 1H), 5.11 – 5.08 (m, 2H), 4.57 – 4.52 (m, 2H), 2.30 – 2.25 (m, 2H), 2.12 – 2.03 (m, 7H), 1.76 – 1.58 (m, 11H), 1.43 – 1.38 (m, 1H), 1.27 – 1.23 (m, 1H), 1.11 – 1.07 (m, 2H), 0.29 (s, 6H).
 ^{13}C NMR (100 MHz, Chloroform-*d*); δ = 174.9, 142.5, 138.2, 135.9, 133.6, 131.6,

129.1, 127.9, 124.3, 124.2, 119.2, 61.1, 32.4, 32.0, 28.9, 26.6, 26.4, 25.8, 23.6, 23.4, 17.7, 10.9, – 3.3.

HRMS (ESI⁺) calcd for C₂₆H₄₁O₂Si [M+H]⁺: 413.2876; found 413.2873.

2-Hydroxy-3-(2-methoxyphenoxy)propyl 3-(dimethyl(phenyl)silyl)propanoate 9e



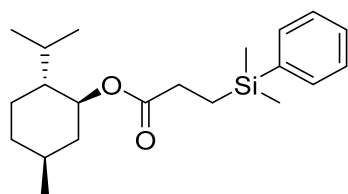
Colorless oil

¹H NMR (400 MHz, Chloroform-*d*); δ = 7.50 – 7.48 (m, 2H), 7.37 – 7.35 (m, 3H), 7.01 – 6.88 (m, 4H), 4.25 – 4.18 (m, 3H), 4.09 – 3.97 (m, 2H), 3.85 (s, 3H), 2.34 – 3.00 (m, 2H), 1.11 – 1.07 (m, 2H), 0.29 (s, 6H).

¹³C NMR (100 MHz, Chloroform-*d*); δ = 175.0, 150.0, 147.9, 138.0, 133.6, 129.2, 127.9, 122.6, 121.1, 115.8, 112.0, 71.5, 68.5, 65.2, 55.8, 28.7, 10.8, – 3.3.

HRMS (ESI⁺) calcd for C₂₁H₂₉O₅Si [M+H]⁺: 389.1784; found 389.1783.

(1*S*,2*R*,5*S*)-2-isopropyl-5-methylcyclohexyl 3-(dimethyl(phenyl)silyl)propanoate 9f



Colorless oil

¹H NMR (400 MHz, Chloroform-*d*); δ = 7.52 – 7.49 (m, 2H), 7.38 – 7.35 (m, 3H), 4.69 – 4.62 (m, 1H), 2.27 – 2.22 (m, 2H), 1.97 – 1.80 (m, 2H), 1.69 – 1.30 (m, 7H), 1.60 (s, 2H), 1.10 – 1.05 (m, 2H), 0.89 – 0.87 (m, 6H), 0.75 (d, *J* = 7 Hz, 3H), 0.29 (s, 6H).

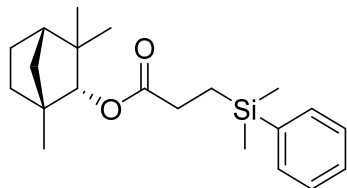
¹³C NMR (100 MHz, Chloroform-*d*); δ = 174.6, 138.3, 133.6, 129.1, 127.9, 74.0, 47.0, 40.9, 34.3, 31.4, 29.2, 26.2, 23.4, 22.1, 20.8, 16.3, 10.9, – 3.24, – 3.27.

HRMS (ESI⁺) calcd for C₂₁H₃₄O₂SiNa [M+Na]⁺: 369.2226; found 369.2227.

(1R,2R,4S)-1,3,3-trimethylbicyclo[2.2.1]heptan-2-yl

3-

(dimethyl(phenyl)silyl)propanoate 9g



Colorless oil

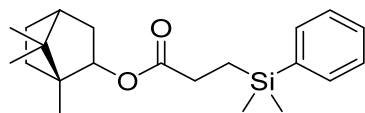
^1H NMR (400 MHz, Chloroform-*d*); δ = 7.54 – 7.50 (m, 2H), 7.40 – 7.32 (m, 3H), 4.35 (s, 1H), 2.32 – 2.28 (m, 2H), 1.75 – 1.55 (m, 5H), 1.48 – 1.40 (m, 1H), 1.19 – 1.13 (m, 3H), 1.10 (s, 3H), 1.02 (s, 3H), 0.75 (s, 3H), 0.30 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 175.5, 138.3, 133.6, 129.1, 127.9, 86.0, 48.34, 48.25, 41.4, 39.5, 29.7, 28.9, 26.6, 25.9, 20.2, 19.5, 11.0, – 3.23, – 3.24.

HRMS (ESI⁺) calcd for C₂₁H₃₃O₂Si [M+H]⁺: 345.2250; found 345.2250.

(1S)-1,7,7-trimethylbicyclo[2.2.1]heptan-2-yl 3-(dimethyl(phenyl)silyl)propanoate

9h



Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.51 – 7.49 (m, 2H), 7.37 – 7.35 (m, 3H), 4.65 – 4.62 (m, 1H), 2.26 – 2.22 (m, 2H), 1.81 – 1.50 (m, 6H), 1.10 – 1.06 (m, 3H), 0.96 (s, 3H), 0.86 – 0.82 (m, 6H), 0.29 (s, 6H).

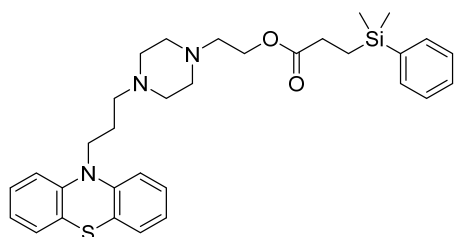
^{13}C NMR (100 MHz, Chloroform-*d*); δ = 174.5, 138.3, 133.6, 129.1, 127.9, 80.9, 48.7, 46.9, 45.0, 38.8, 33.8, 29.2, 27.1, 20.2, 20.0, 11.5, 10.8, – 3.2, – 3.3.

HRMS (ESI⁺) calcd for C₂₁H₃₃O₂Si [M+H]⁺: 345.2250; found 345.2247.

2-(4-(3-(10H-phenothiazin-10-yl)propyl)piperazin-1-yl)ethyl

3-

(dimethyl(phenyl)silyl)propanoate 9i



Colorless oil

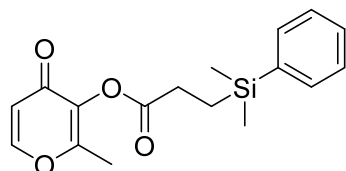
^1H NMR (400 MHz, Chloroform-*d*); δ = 7.51 – 7.48 (m, 2H), 7.39 – 7.35 (m, 4H), 7.17 – 7.11 (m, 2H), 7.02 – 6.84 (m, 5H), 4.15 (t, J = 6.0 Hz, 2H), 3.90 (t, J = 6.8 Hz, 2H), 2.59 – 2.44 (m, 12H), 2.30 – 2.25 (m, 2H), 1.95 – 1.92 (m, 2H), 1.10 – 1.05 (m, 2H), 0.29 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 174.8, 146.4, 144.4, 138.1, 133.5, 133.1, 133.0, 129.5, 129.1, 127.82, 127.80, 127.44, 127.35, 124.7, 123.4, 122.8, 122.2, 115.8, 115.7, 61.7, 56.5, 55.4, 53.2, 53.1, 45.3, 28.8, 24.1, 10.7, – 3.4.

HRMS (ESI⁺) calcd for C₃₂H₄₂N₃O₂SSi [M+H]⁺: 560.2767; found 560.2769.

2-methyl-4-oxo-4*H*-pyran-3-yl 3-(dimethyl(phenyl)silyl)propanoate

9j



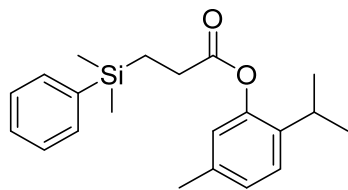
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.66 – 7.64 (m, 1H), 7.53 – 7.50 (m, 2H), 7.38 – 7.37 (m, 3H), 6.40 (d, J = 5.8 Hz, 1H), 2.60 – 2.55 (m, 2H), 2.21 (s, 3H), 1.24 – 1.19 (m, 2H), 0.33 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 172.1, 171.7, 159.0, 154.1, 138.7, 137.9, 133.6, 129.2, 128.0, 116.9, 28.4, 15.0, 10.9, – 3.3.

HRMS (ESI⁺) calcd for C₁₈H₂₅O₃Si [M+H]⁺: 317.1573; found 317.1574.

2-Isopropyl-5-methylphenyl 3-(dimethyl(phenyl)silyl)propanoate **9k**



Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.60 – 7.54 (m, 2H), 7.41 – 7.33 (m, 3H), 7.19 – 7.17 (m, 1H), 7.02 – 7.00 (m, 1H), 6.74 (s, 1H), 2.99 – 2.88 (m, 1H), 2.57 – 2.52 (m, 2H), 1.30 (s, 3H), 1.25 – 1.23 (m, 2H), 1.18 (d, J = 6.9 Hz, 6H), 0.36 (s, 6H).

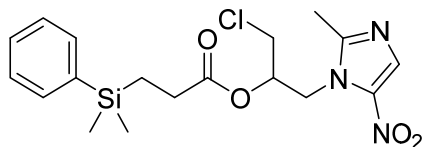
^{13}C NMR (100 MHz, Chloroform-*d*); δ = 173.7, 148.0, 138.0, 137.0, 136.5, 133.6, 129.2, 128.0, 127.1, 126.4, 122.7, 29.0, 27.0, 23.1, 20.9, 11.0, – 3.2.

HRMS (ESI⁺) calcd for C₂₁H₂₈O₂SiNa [M+Na]⁺: 363.1756; found 363.1755.

1-Chloro-3-(2-methyl-5-nitro-1*H*-imidazol-1-yl)propan-2-yl

3-

(dimethyl(phenyl)silyl)propanoate 9l



Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.94 (s, 1H), 7.48 – 7.45 (m, 2H), 7.38 – 7.36 (m, 3H), 5.36 – 5.33 (m, 1H), 4.74 – 4.70 (m, 1H), 4.49 – 4.43 (m, 1H), 3.80 – 3.71 (m, 2H), 2.48 (s, 3H), 2.26 – 2.09 (m, 2H), 1.03 – 0.88 (m, 2H), 0.27 (s, 6H).

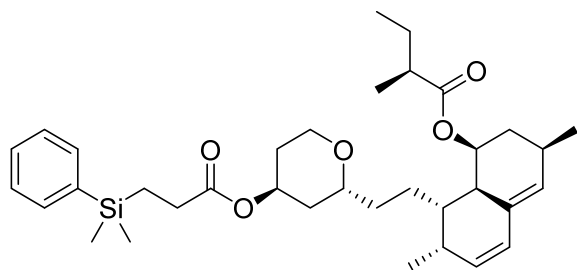
^{13}C NMR (100 MHz, Chloroform-*d*); δ = 173.5, 151.0, 138.6, 137.6, 133.5, 133.2, 129.3, 128.0, 70.7, 47.1, 43.4, 28.6, 14.4, 10.7, – 3.37, – 3.43.

HRMS (ESI⁺) calcd for C₁₈H₂₅ClN₃O₄Si [M+H]⁺: 410.1303; found 410.1307.

(1*S*,3*R*,7*S*,8*S*,8*aR*)-8-(2-((2*R*,4*S*)-4-((3-

(dimethyl(phenyl)silyl)propanoyl)oxy)tetrahydro-2*H*-pyran-2-yl)ethyl)-3,7-

dimethyl-1,2,3,7,8,8*a*-hexahydronaphthalen-1-yl (*S*)-2-methylbutanoate 9m



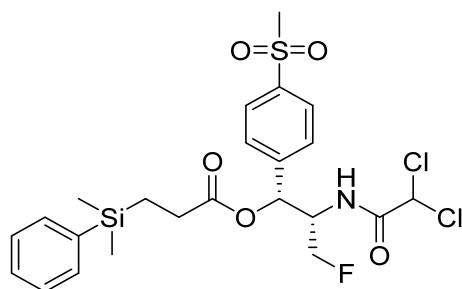
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.61 – 7.59 (m, 1H), 7.50 – 7.48 (m, 2H), 7.39 – 7.35 (m, 5H), 5.99 – 5.97 (m, 1H), 5.79 – 5.75 (m, 1H), 5.53 – 5.51 (m, 1H), 5.36 – 5.34 (m, 1H), 3.93 – 3.89 (m, 1H), 2.46 – 2.19 (m, 6H), 2.01 – 1.87 (m, 2H), 1.80 – 1.74 (m, 2H), 1.69 – 1.58 (m, 5H), 1.45 – 1.33 (m, 3H), 1.11 – 1.06 (m, 6H), 0.87 – 0.83 (m, 6H), 0.41 (s, 3H), 0.34 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 176.6, 173.9, 135.6, 133.9, 133.1, 133.0, 131.6, 129.7, 129.6, 128.3, 128.1, 127.9, 78.7, 67.8, 41.4, 37.2, 36.7, 32.7, 32.6, 30.7, 29.9, 28.6, 27.5, 26.9, 24.8, 22.8, 16.3, 15.1, 13.9, 11.7, – 5.4, – 5.5.

HRMS (ESI⁺) calcd for C₃₅H₅₂O₅SiNa [M+Na]⁺: 603.3482; found 603.3485.

(1*R*,2*S*)-2-(2,2-dichloroacetamido)-3-fluoro-1-(4-(methylsulfonyl)phenyl)propyl 3-(dimethyl(phenyl)silyl)propanoate 9n



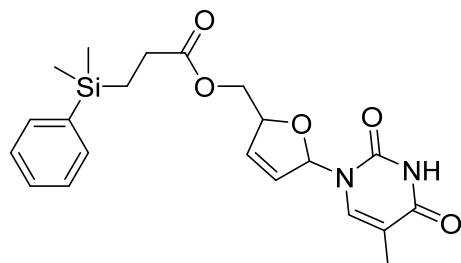
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); δ = 7.98 – 7.96 (m, 2H), 7.60 – 7.58 (m, 2H), 7.48 – 7.34 (m, 5H), 6.91 – 6.87 (m, 1H), 6.11 (d, J = 7.4 Hz, 1H), 5.88 (s, 1H), 4.53 – 4.18 (m, 3H), 3.06 (s, 3H), 2.36 – 2.31 (m, 2H), 1.07 (t, J = 8.7 Hz, 2H), 0.28 (s, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); δ = 173.9, 164.2, 142.4, 141.2, 137.6, 133.5, 129.3, 128.1, 128.1, 128.0, 82.4, 80.7, 72.6, 66.0, 54.2, 54.0, 44.5, 28.8, 10.7, – 3.35, – 3.37.

HRMS (ESI⁺) calcd for C₂₃H₂₈Cl₂FNO₅SSiNa [M+Na]⁺: 570.0716; found 570.0711.

(5-(5-Methyl-2,4-dioxo-3,4-dihydropyrimidin-1(2H)-yl)-2,5-dihydrofuran-2-yl)methyl 3-(dimethyl(phenyl)silyl)propanoate 9o



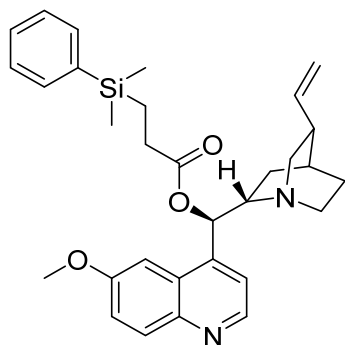
Colorless oil

¹H NMR (400 MHz, Chloroform-*d*); δ = 9.23 (s, 1H), 7.48 – 7.46 (m, 2H), 7.36 – 7.33 (m, 3H), 7.15 (s, 1H), 7.00 – 6.98 (m, 1H), 6.23 – 6.21 (m, 1H), 5.88 – 5.85 (m, 1H), 5.00 – 4.99 (m, 1H), 4.38 – 4.33 (m, 1H), 4.18 – 4.15 (m, 1H), 2.30 – 2.26 (m, 2H), 1.80 (s, 3H), 1.11 – 1.07 (m, 2H), 0.28 (s, 6H).

¹³C NMR (100 MHz, Chloroform-*d*); δ = 174.5, 163.8, 150.8, 137.7, 135.4, 133.5, 133.1, 129.3, 128.0, 127.4, 111.2, 89.7, 84.3, 64.8, 28.7, 12.6, 10.7, – 3.27, – 3.30.

HRMS (ESI⁺) calcd for C₂₁H₂₆N₂O₅SiNa [M+Na]⁺: 437.1509; found 437.1513.

(*R*)-(6-methoxyquinolin-4-yl)((1*S*,2*S*,4*S*,5*R*)-5-vinylquinuclidin-2-yl)methyl 3-(dimethyl(phenyl)silyl)propanoate 9p



Colorless oil

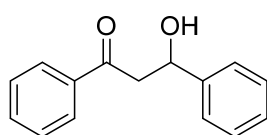
¹H NMR (400 MHz, Chloroform-*d*); δ = 8.73 (d, *J* = 4.6 Hz, 1H), 8.02 (d, *J* = 9.2 Hz, 1H), 7.48 – 7.44 (m, 3H), 7.38 – 7.26 (m, 5H), 6.48 (s, 1H), 5.87 – 5.79 (m, 1H), 5.04 – 4.99 (m, 2H), 3.94 (s, 3H), 3.38 – 3.33 (m, 1H), 3.07 – 3.01 (m, 2H), 2.65 – 2.58 (m,

2H), 2.35 – 2.26 (m, 3H), 1.89 – 1.84 (m, 2H), 1.54 – 1.47 (m, 3H), 1.08 – 1.04 (m, 2H), 0.26 (d, $J = 4.3$ Hz, 6H).

^{13}C NMR (100 MHz, Chloroform-*d*); $\delta = 174.0, 157.9, 147.5, 144.8, 143.6, 141.7, 137.8, 133.5, 131.8, 129.2, 127.9, 127.1, 121.9, 119.0, 114.6, 101.5, 59.1, 56.5, 55.7, 42.4, 29.0, 27.5, 10.7, -3.30, -3.32$.

HRMS (ESI⁺) calcd for $\text{C}_{31}\text{H}_{39}\text{N}_2\text{O}_3\text{Si}$ $[\text{M}+\text{H}]^+$: 515.2730; found 515.2733.

3-Hydroxy-1,3-diphenylpropan-1-one 10

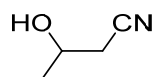


Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); $\delta = 7.97 - 7.95$ (m, 2H), 7.61 – 7.57 (m, 1H), 7.49 – 7.43 (m, 4H), 7.41 – 7.37 (m, 2H), 7.33 – 7.29 (m, 1H), 5.37 – 5.34 (m, 1H), 3.71 (br, 1H), 3.39 – 3.37 (m, 2H).

^{13}C NMR (100 MHz, Chloroform-*d*); $\delta = 200.2, 143.0, 136.5, 133.7, 128.8, 128.6, 128.2, 127.7, 125.8, 70.0, 47.5$.

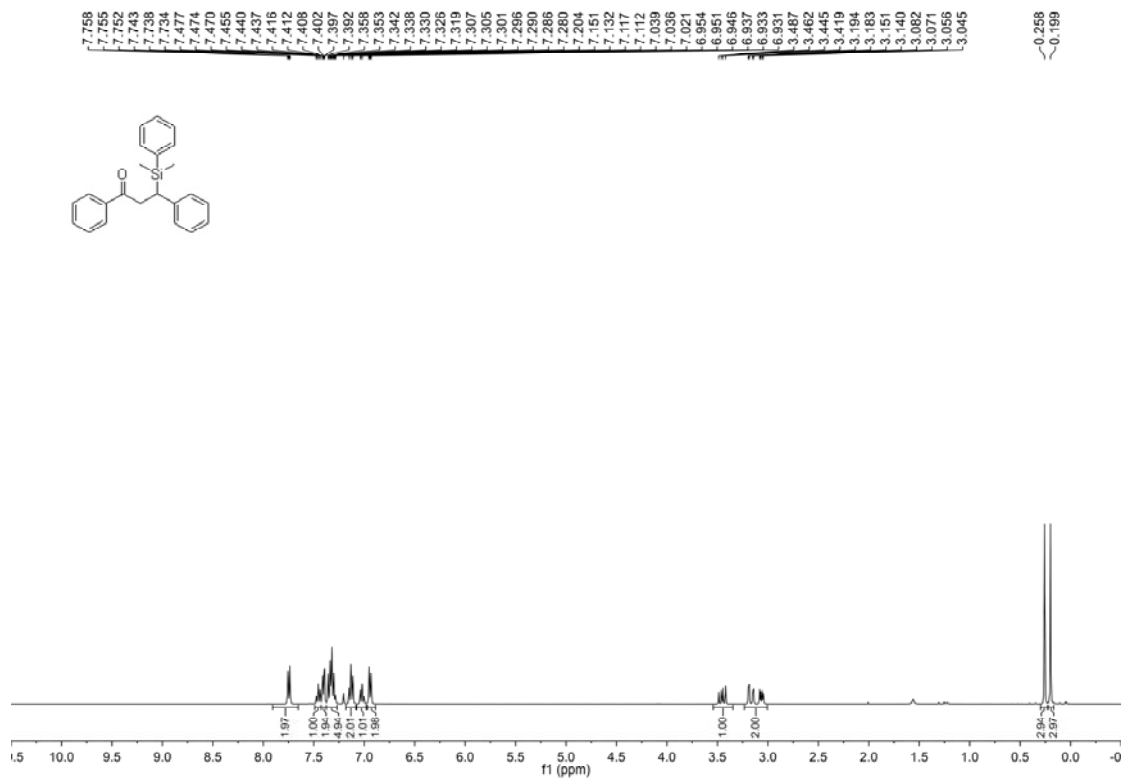
3-Hydroxybutanenitrile 11



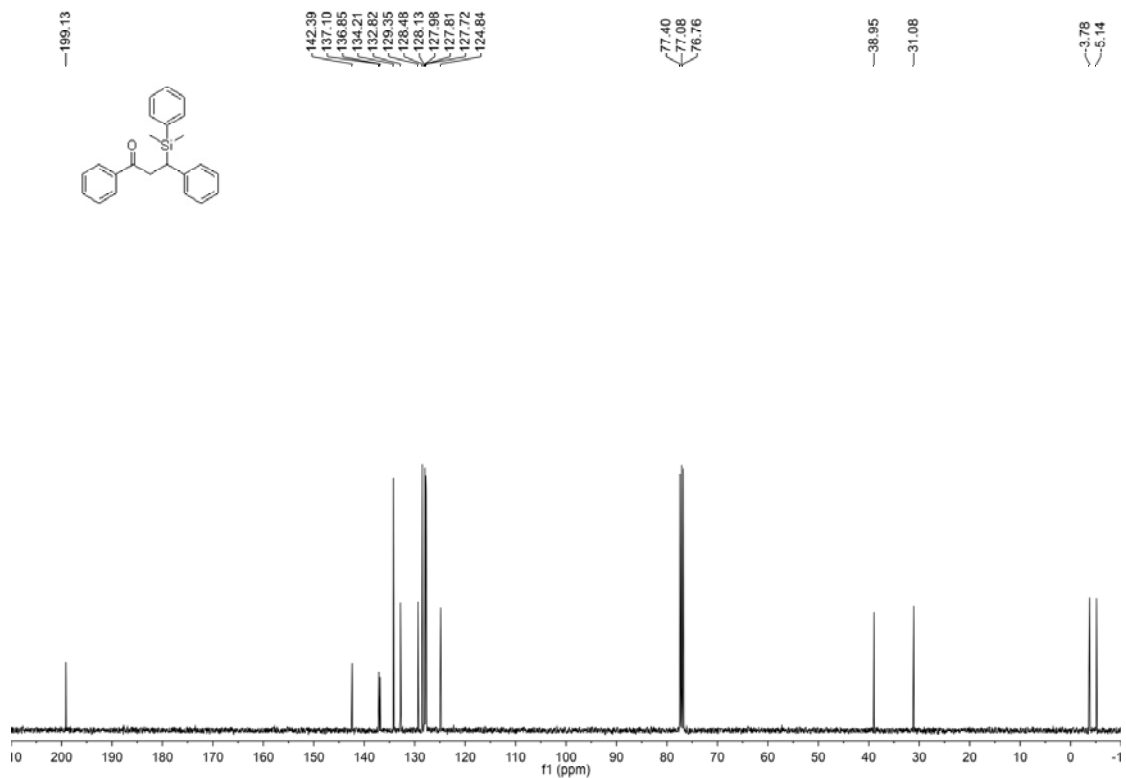
Colorless oil

^1H NMR (400 MHz, Chloroform-*d*); $\delta = 4.19 - 4.16$ (m, 1H), 2.57 – 2.48 (m, 2H), 1.94 (br, 1H), 1.38 (d, $J = 4.2$ Hz, 3H).

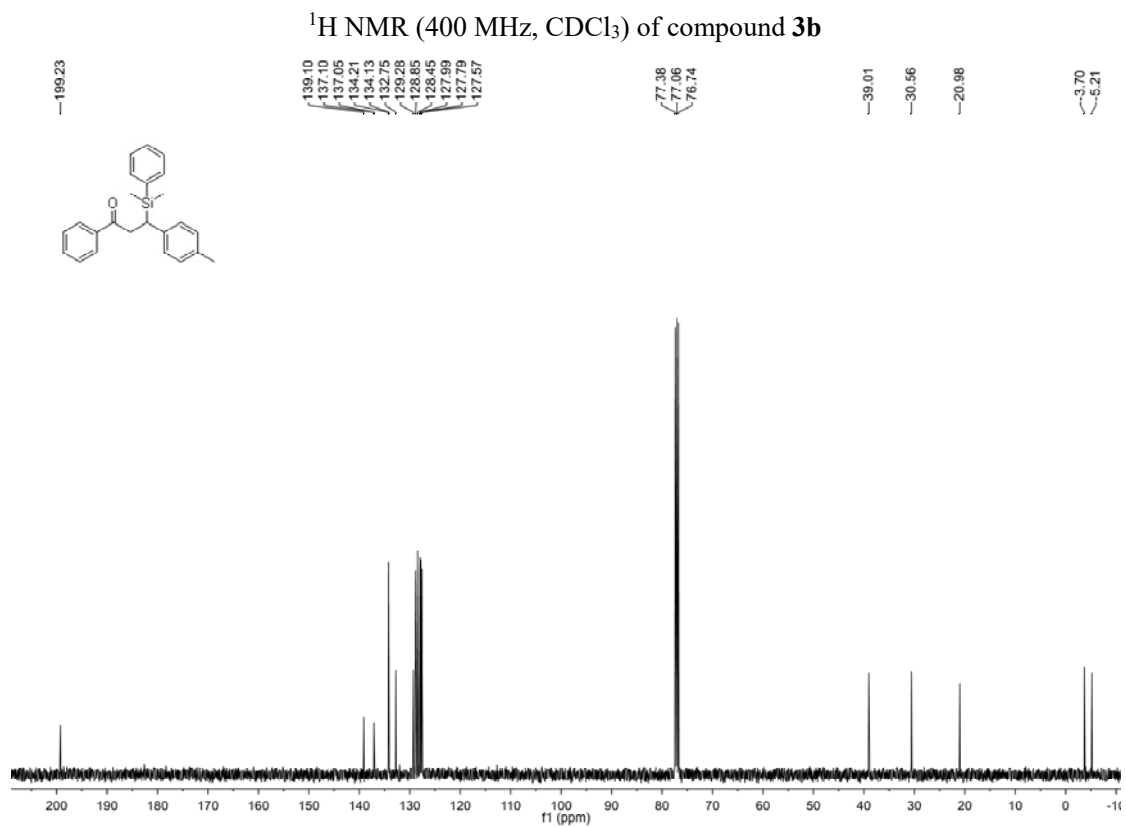
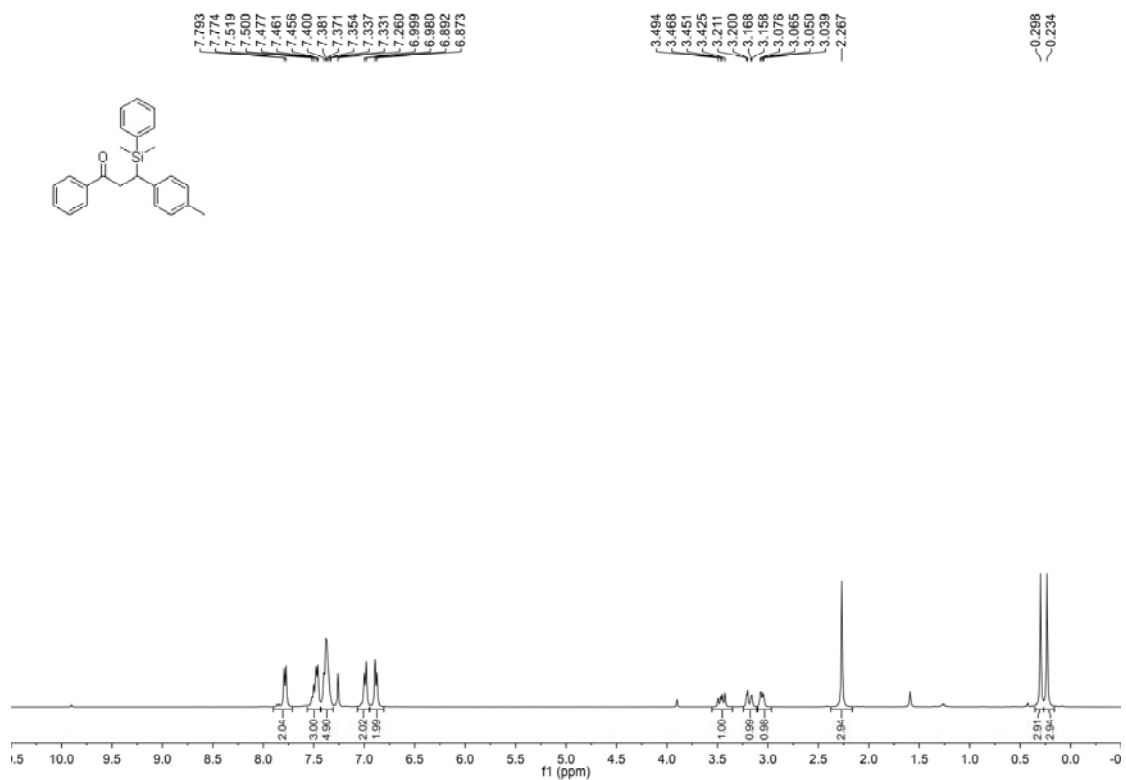
^{13}C NMR (100 MHz, Chloroform-*d*); $\delta = 99.9, 64.2, 27.5, 22.8$.

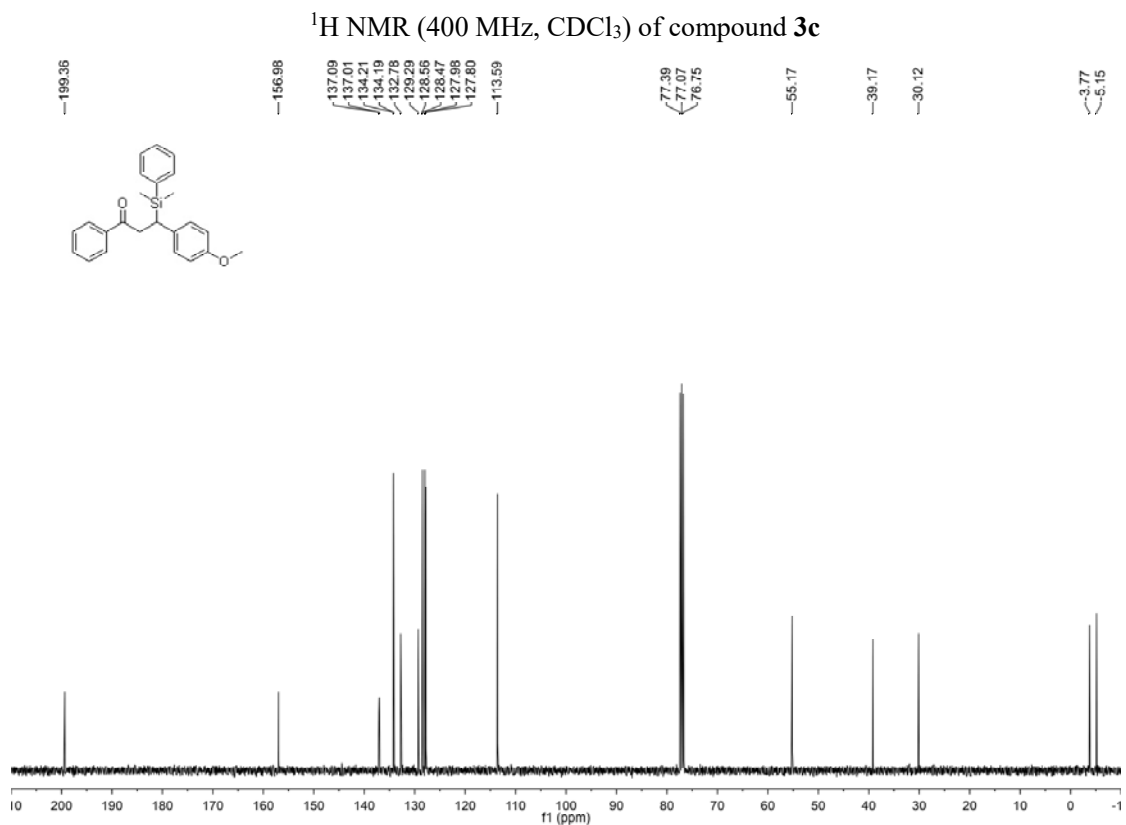
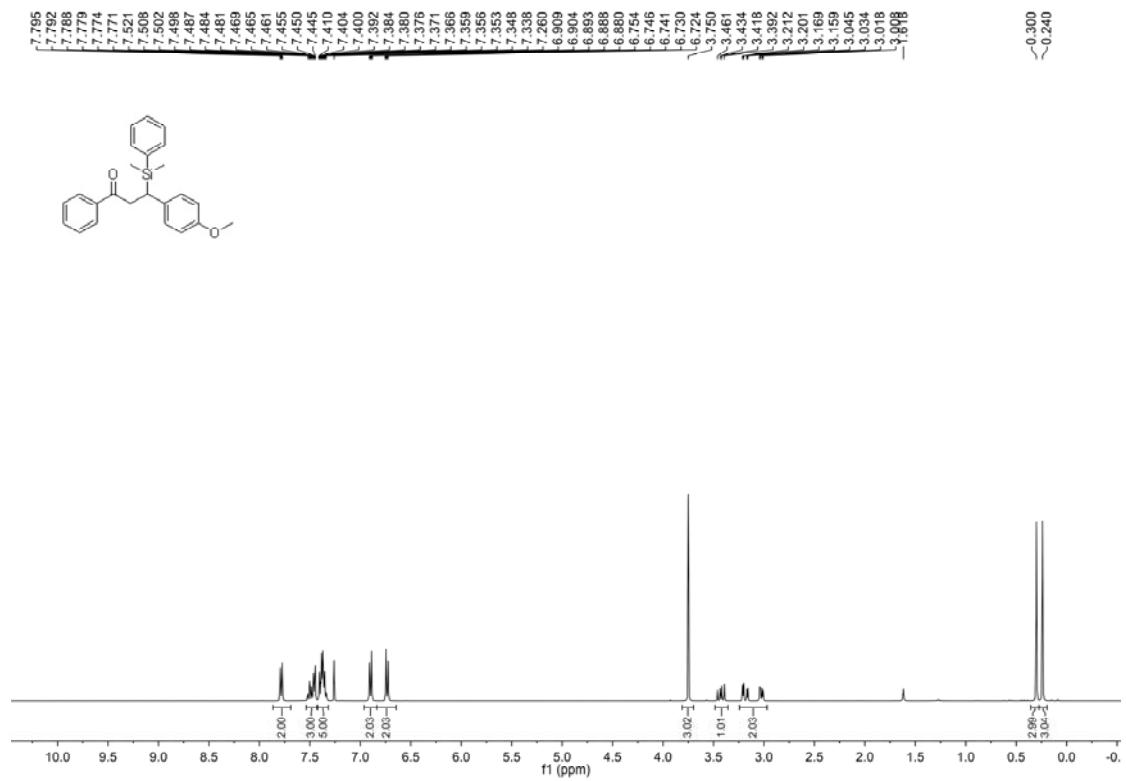


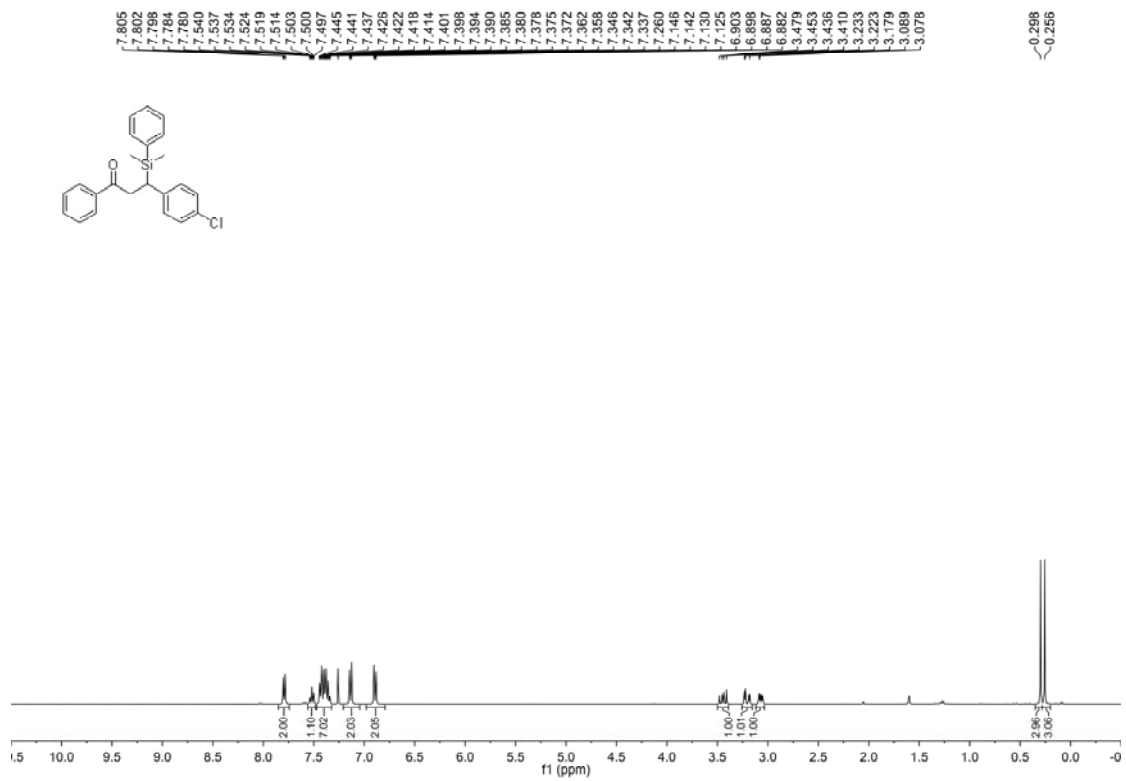
^1H NMR (400 MHz, CDCl_3) of compound **3a**



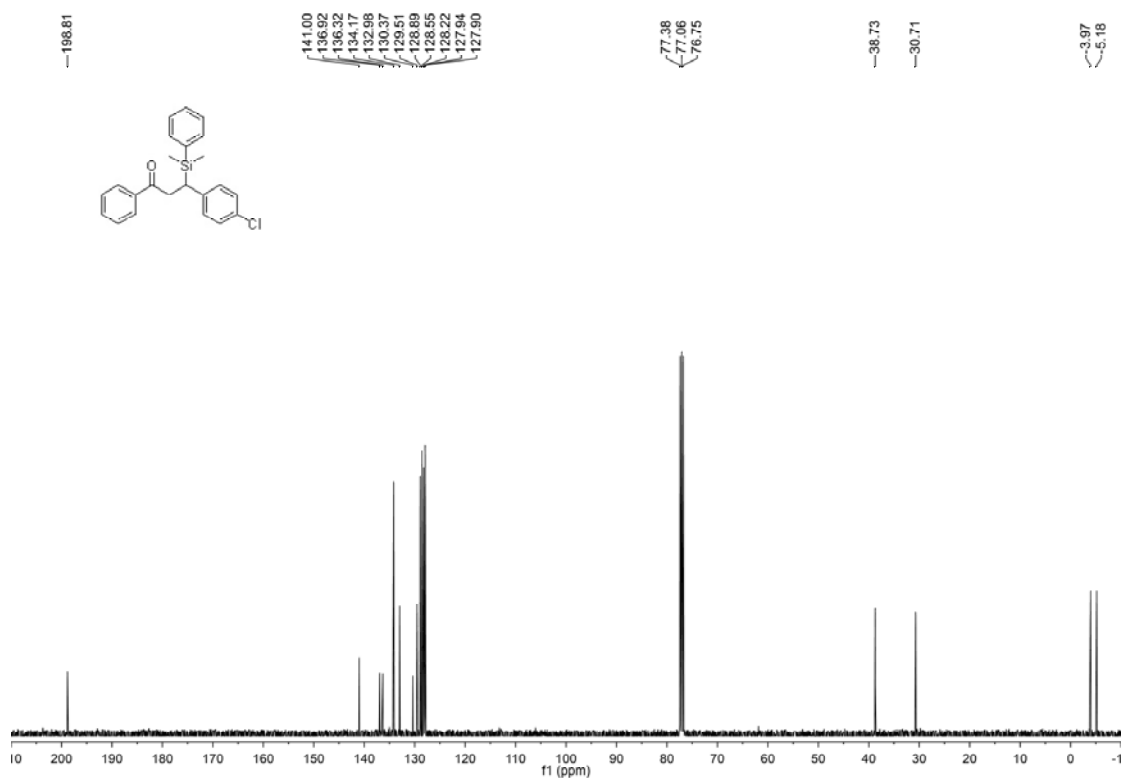
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **3a**



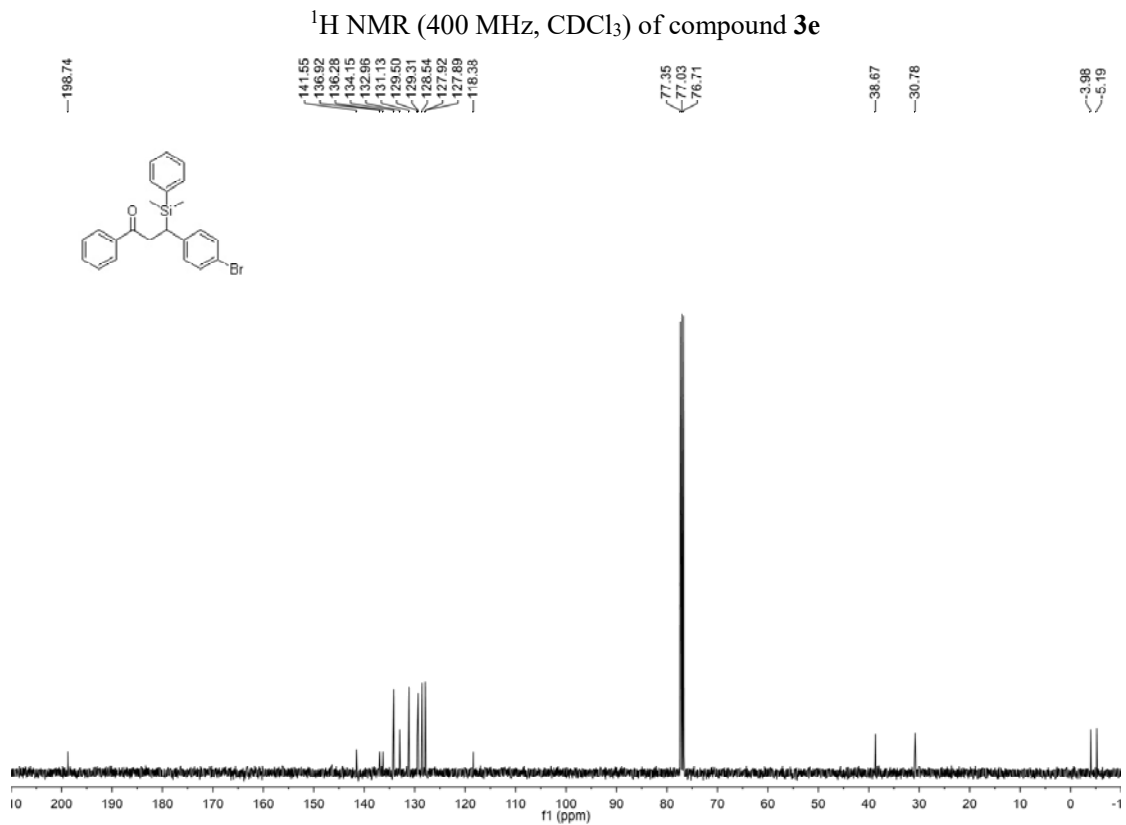
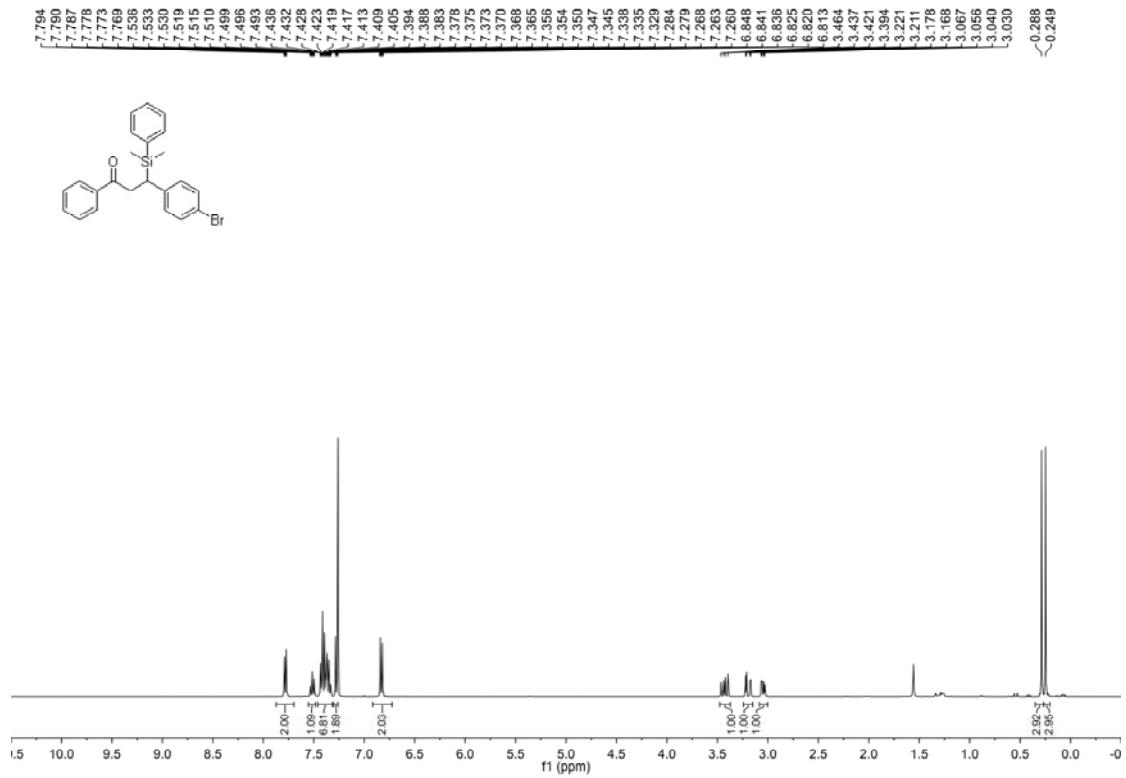


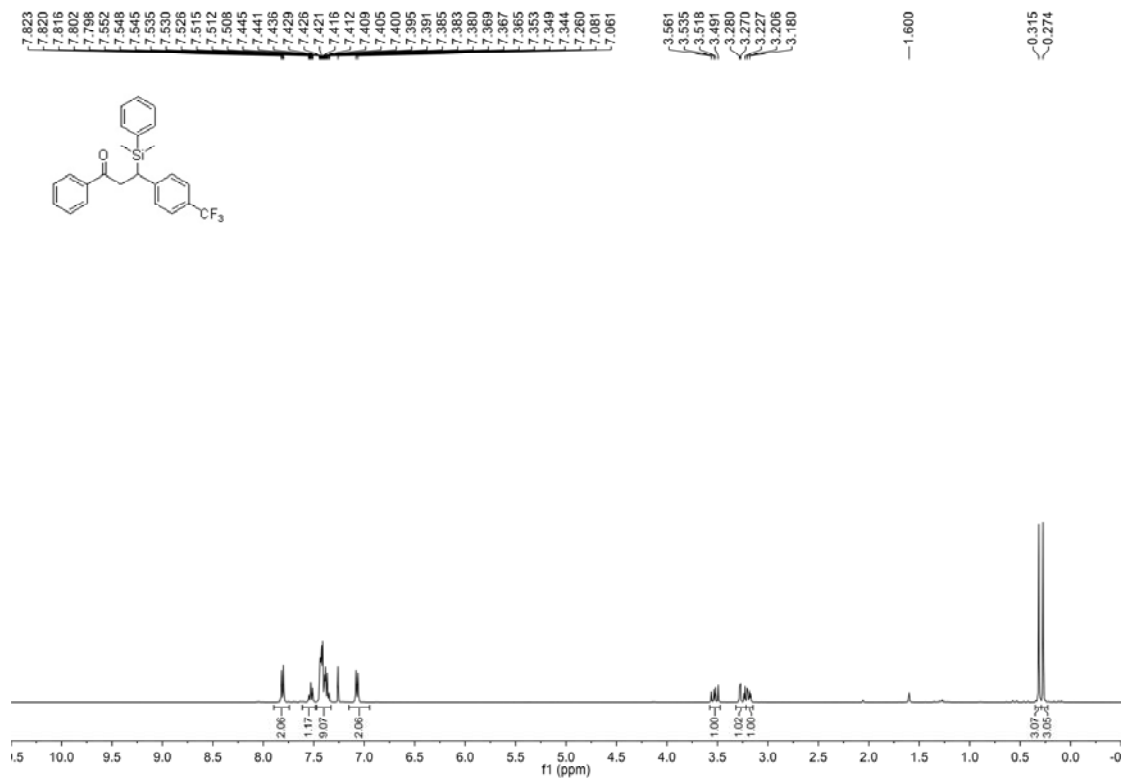


^1H NMR (400 MHz, CDCl_3) of compound **3d**

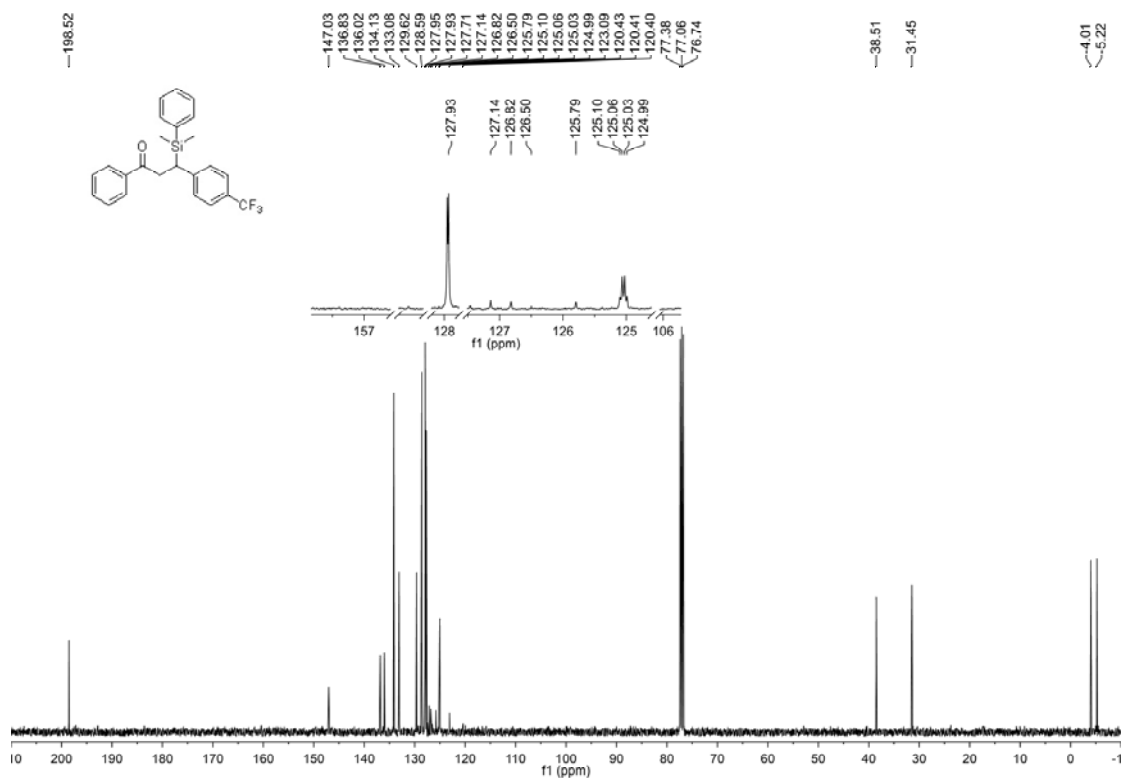


$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **3d**

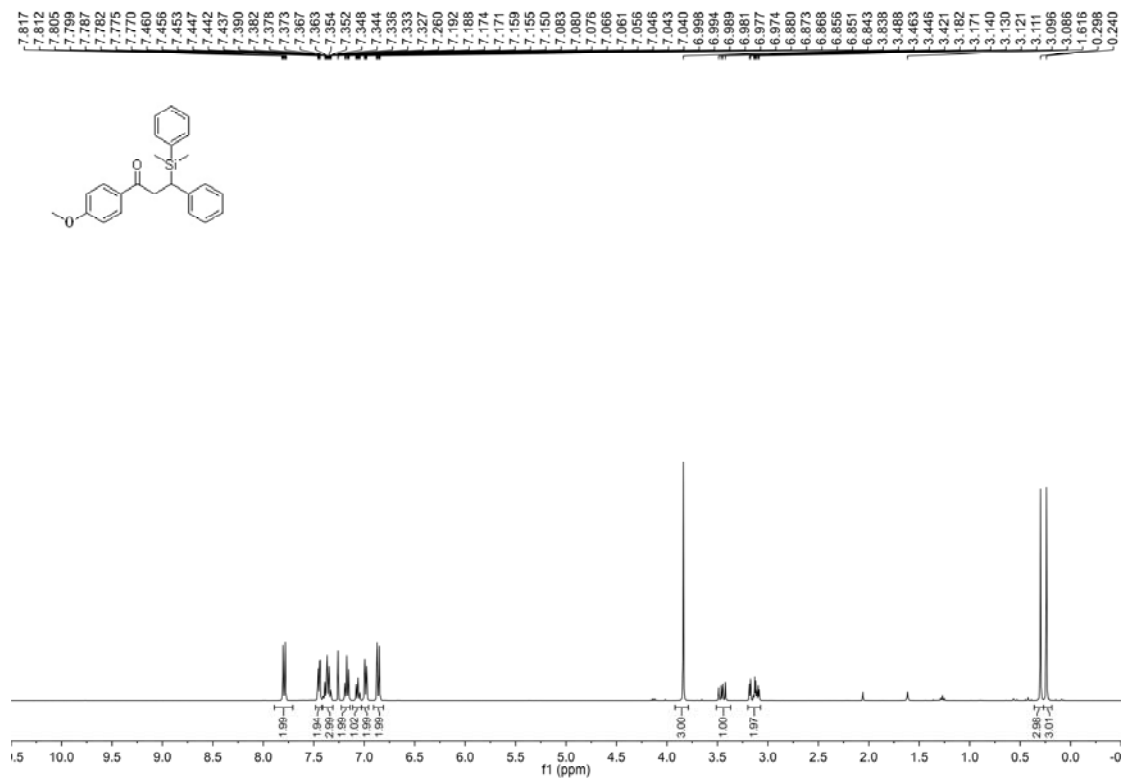




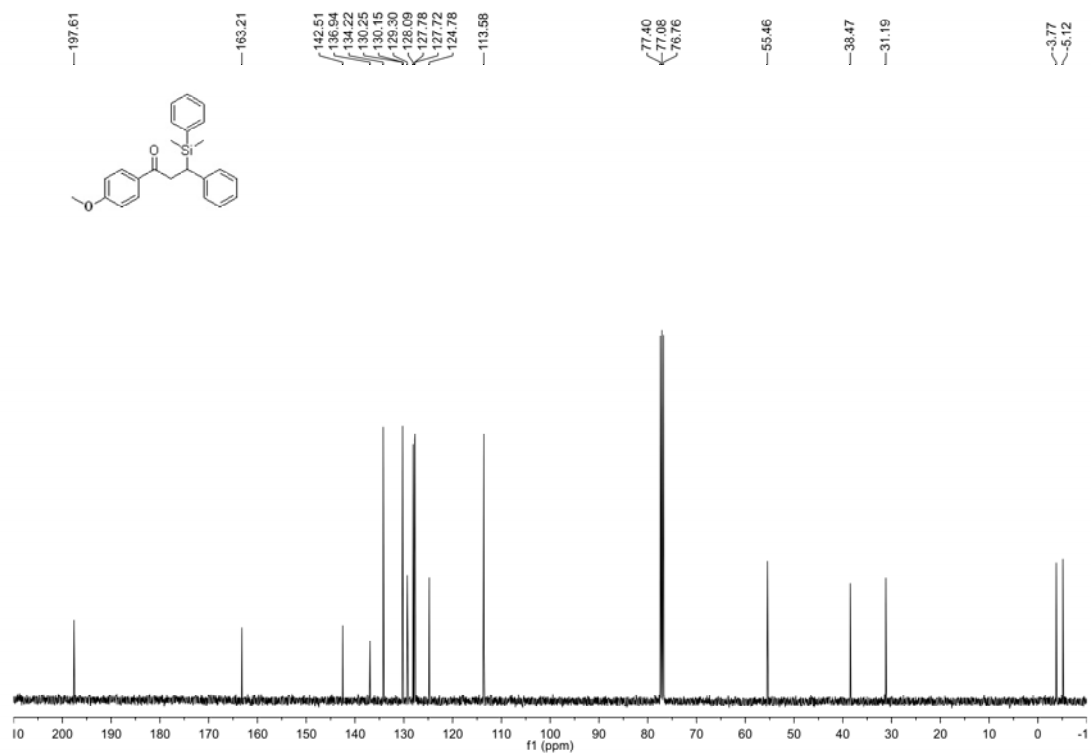
¹H NMR (400 MHz, CDCl₃) of compound **3f**



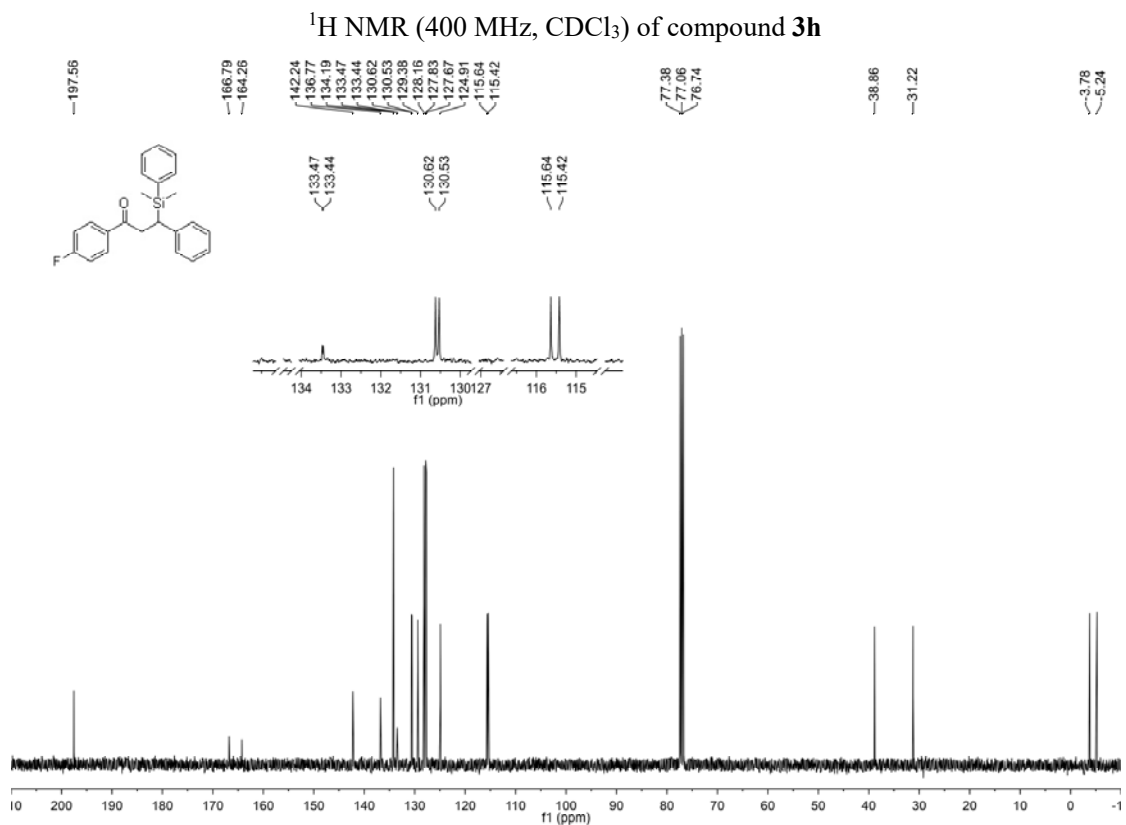
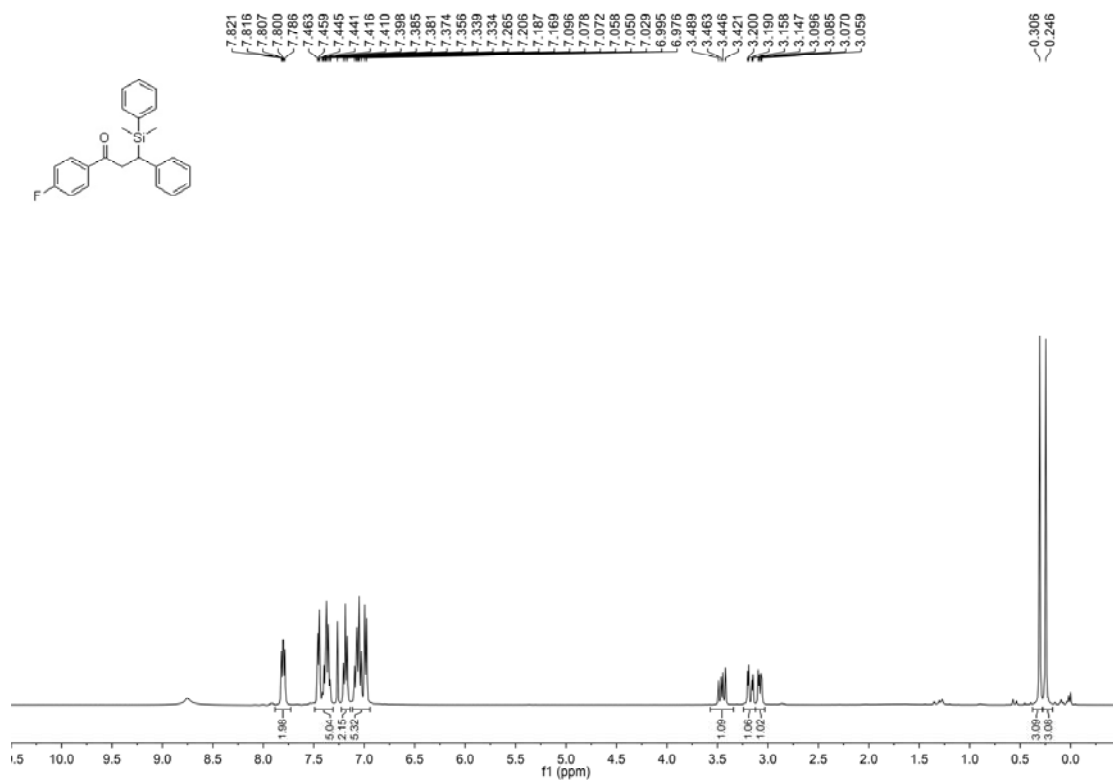
¹³C {¹H} NMR (100 MHz, CDCl₃) of compound **3f**

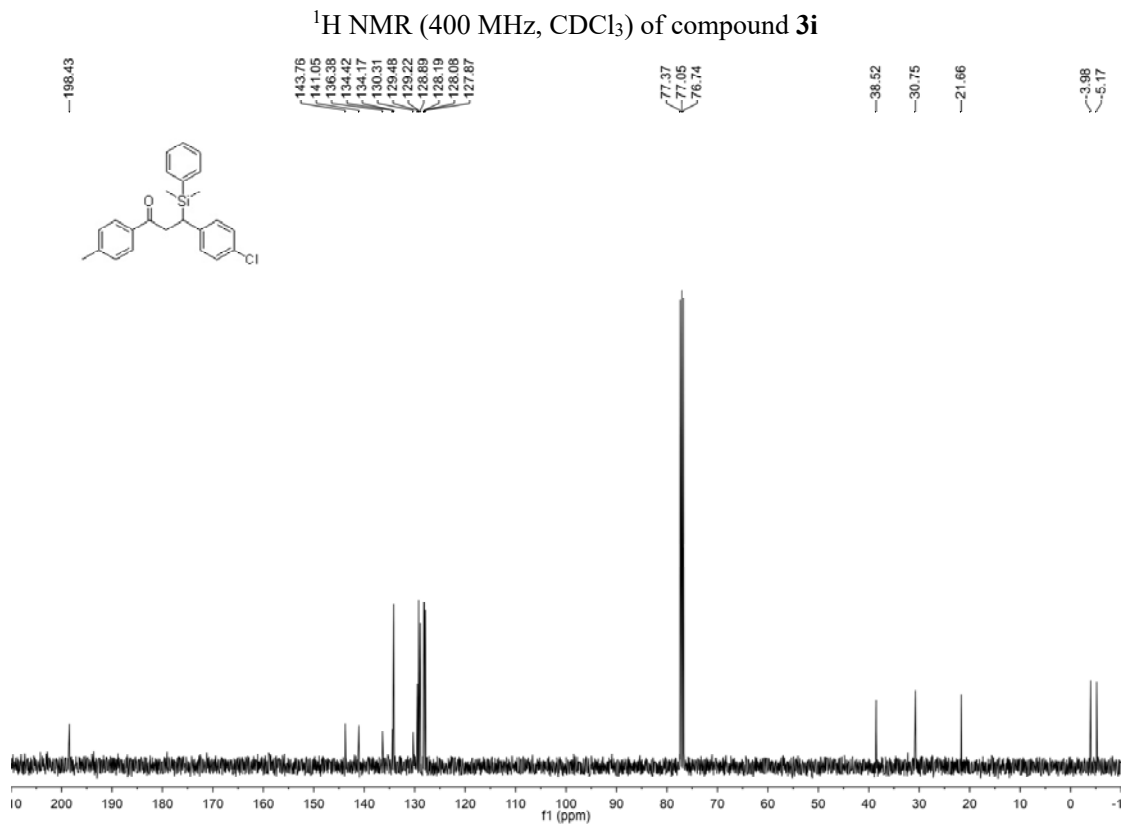
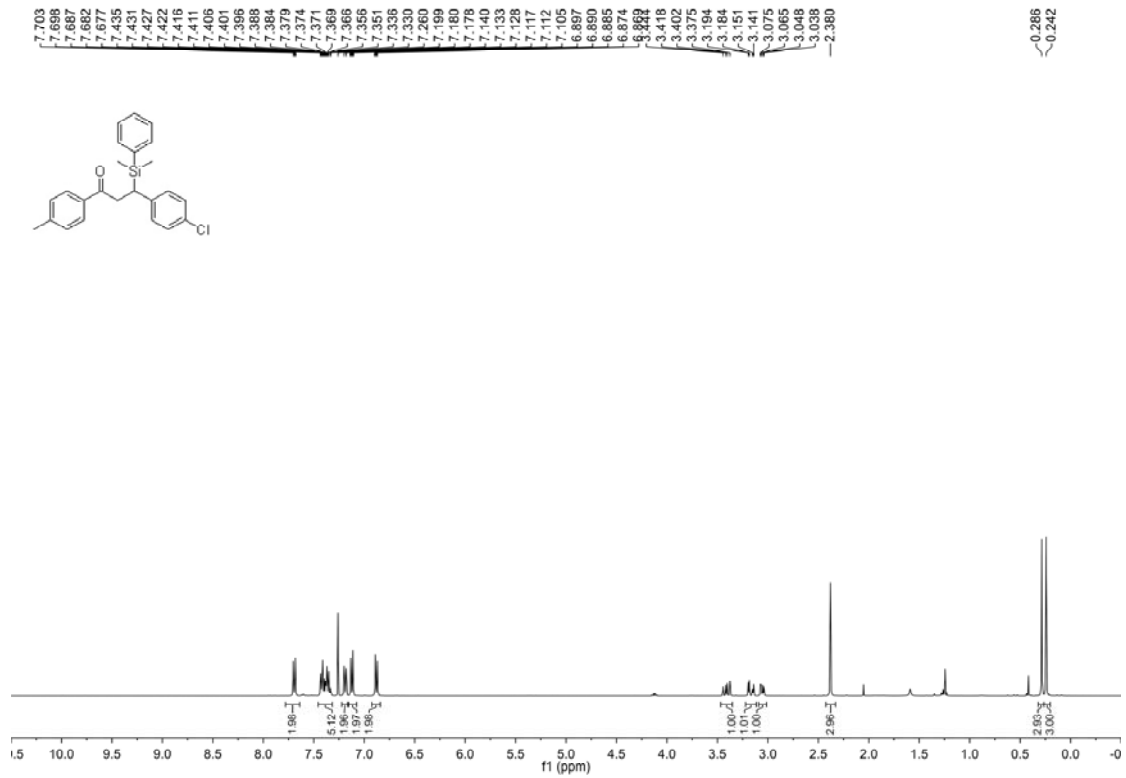


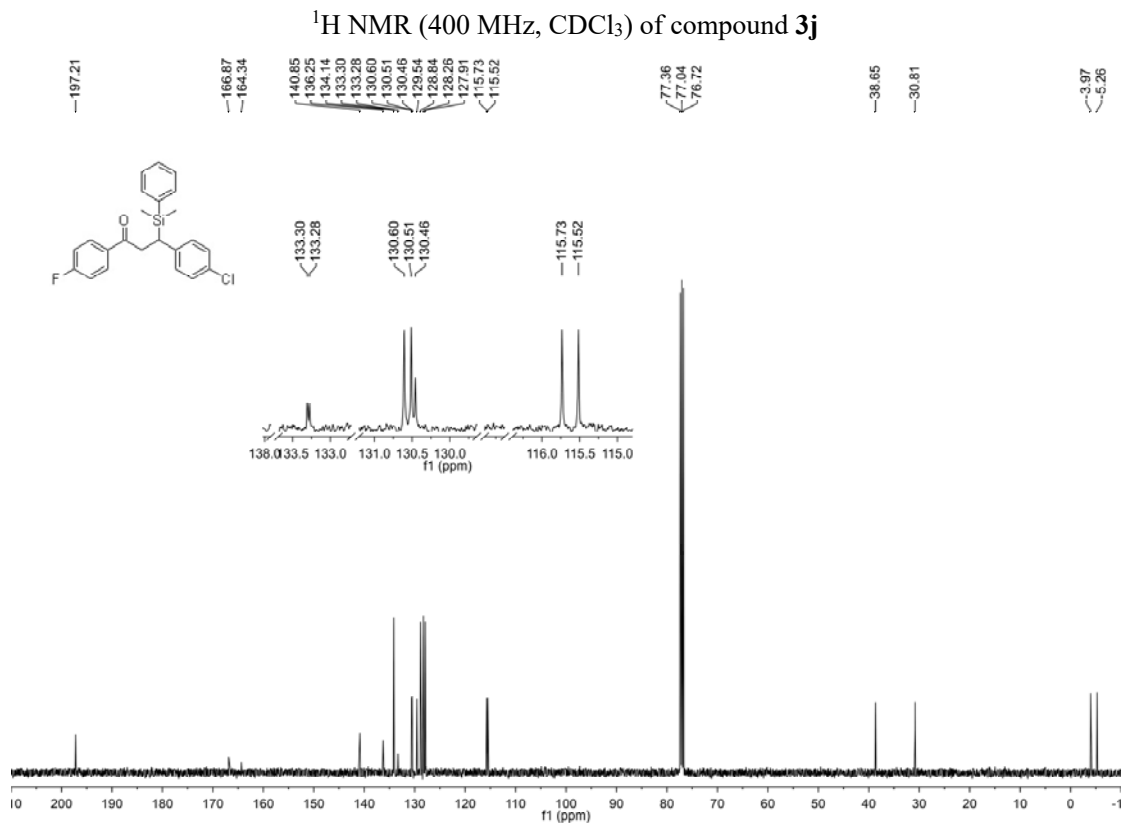
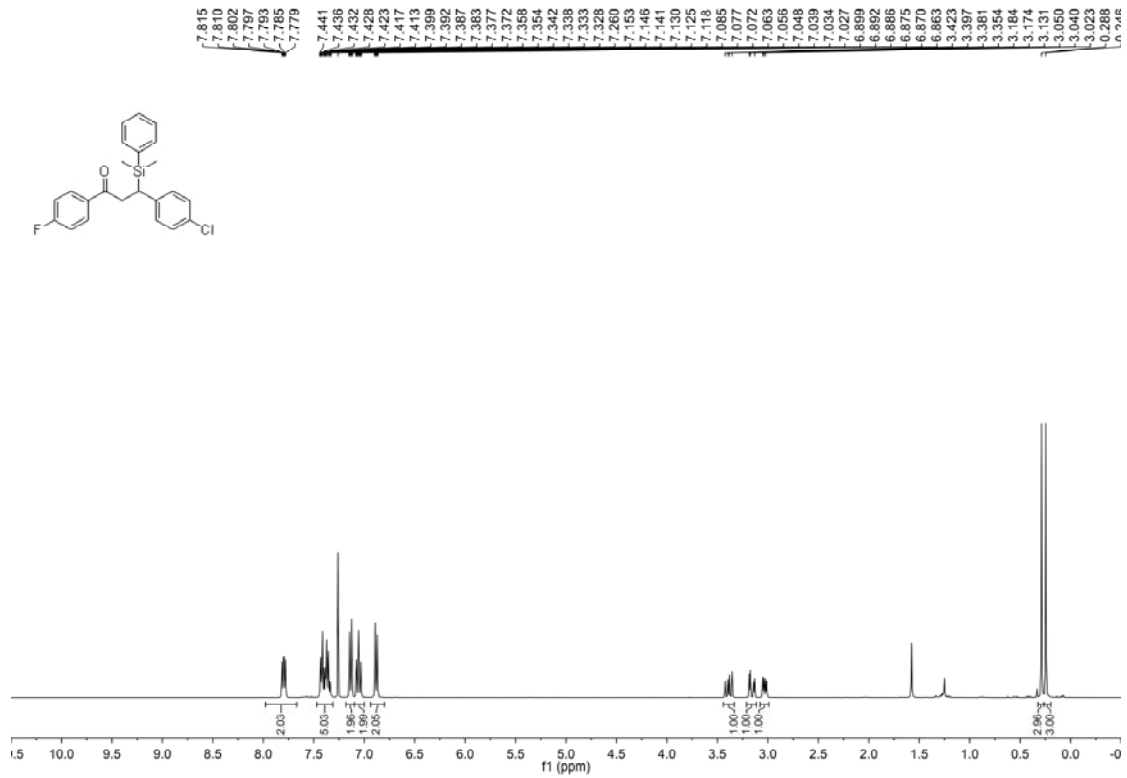
^1H NMR (400 MHz, CDCl_3) of compound **3g**

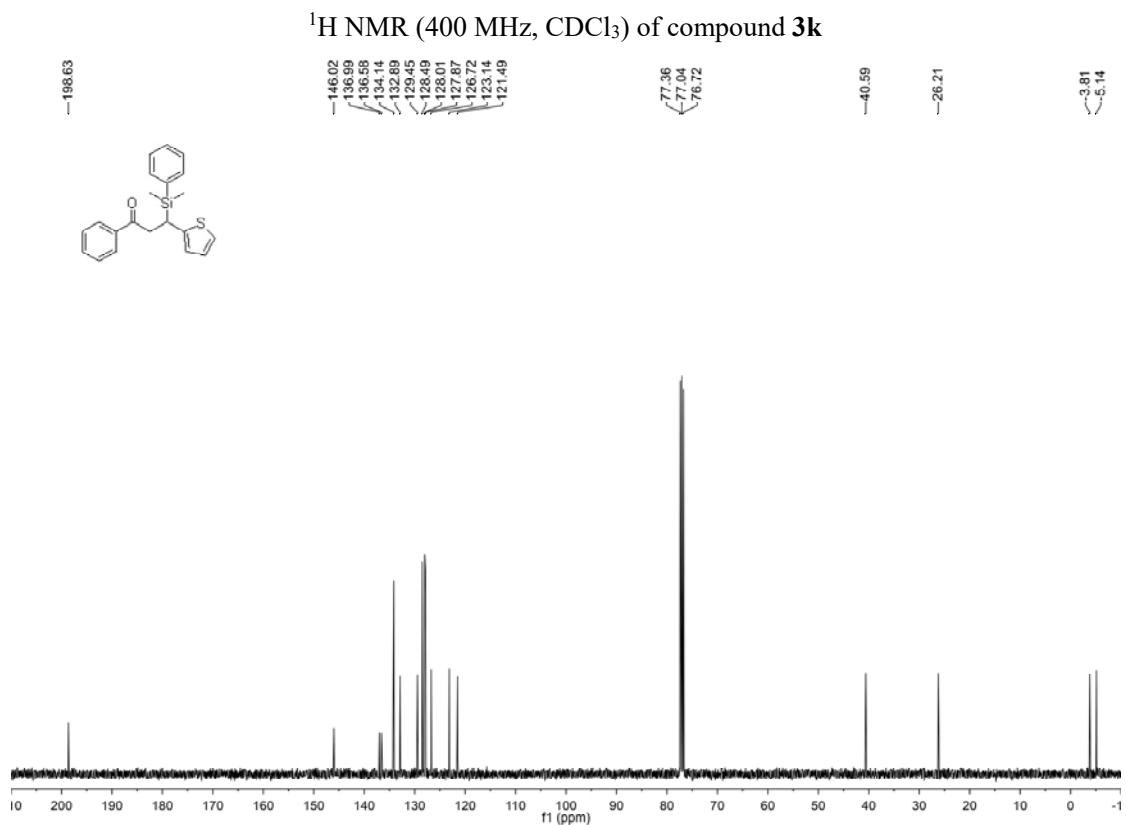
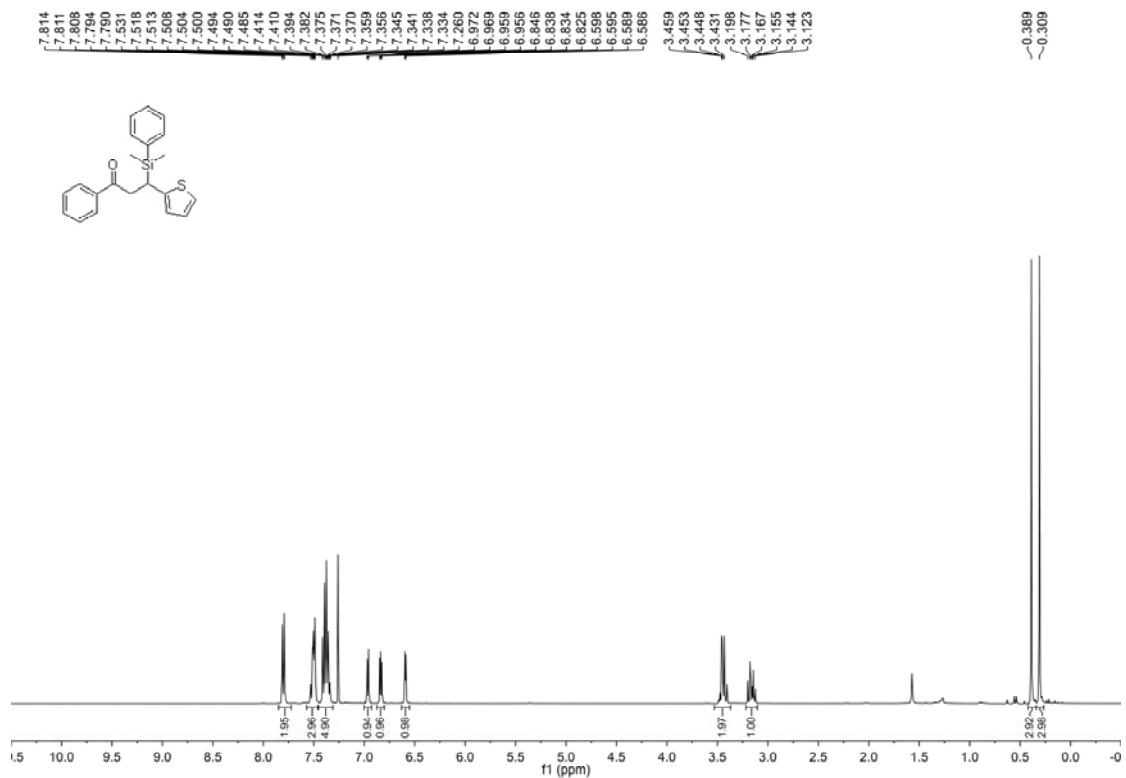


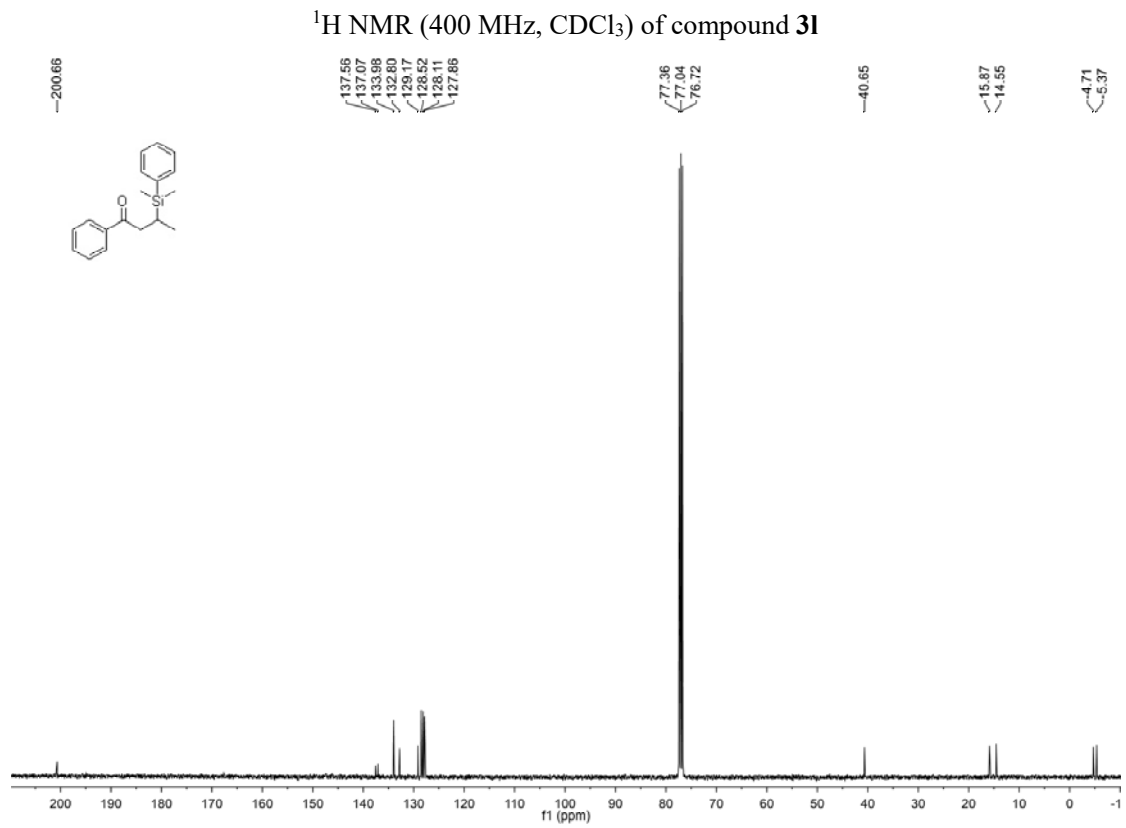
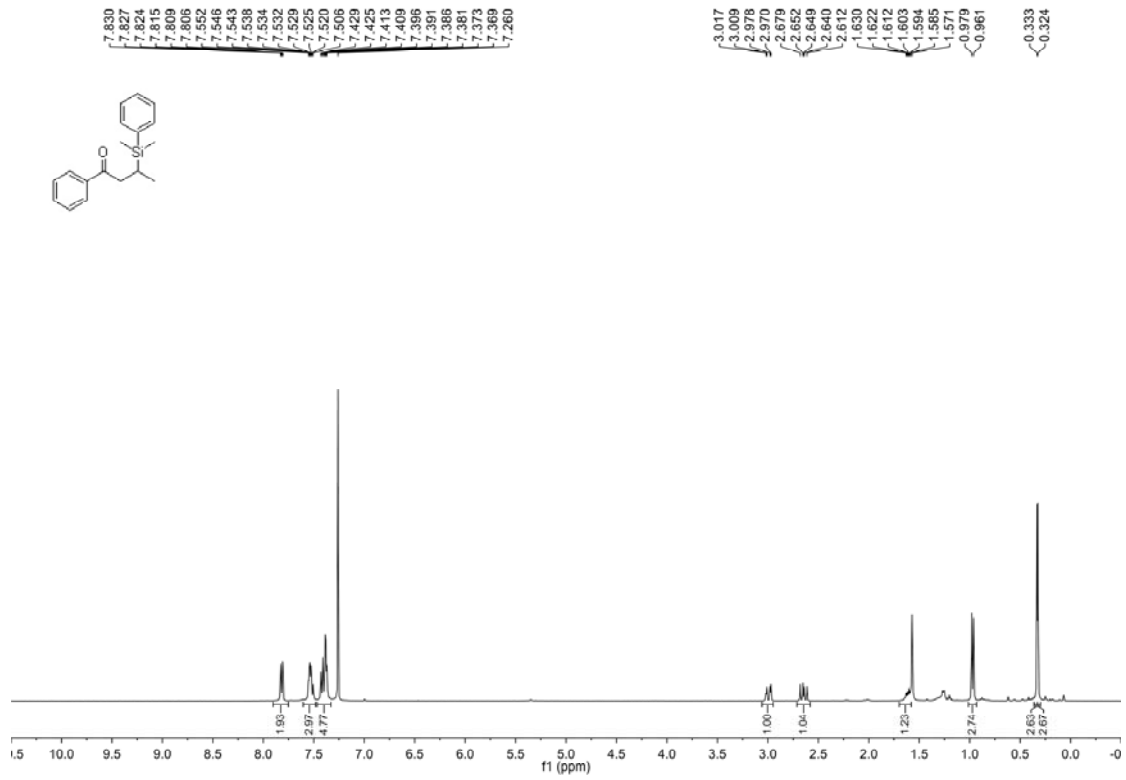
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **3g**



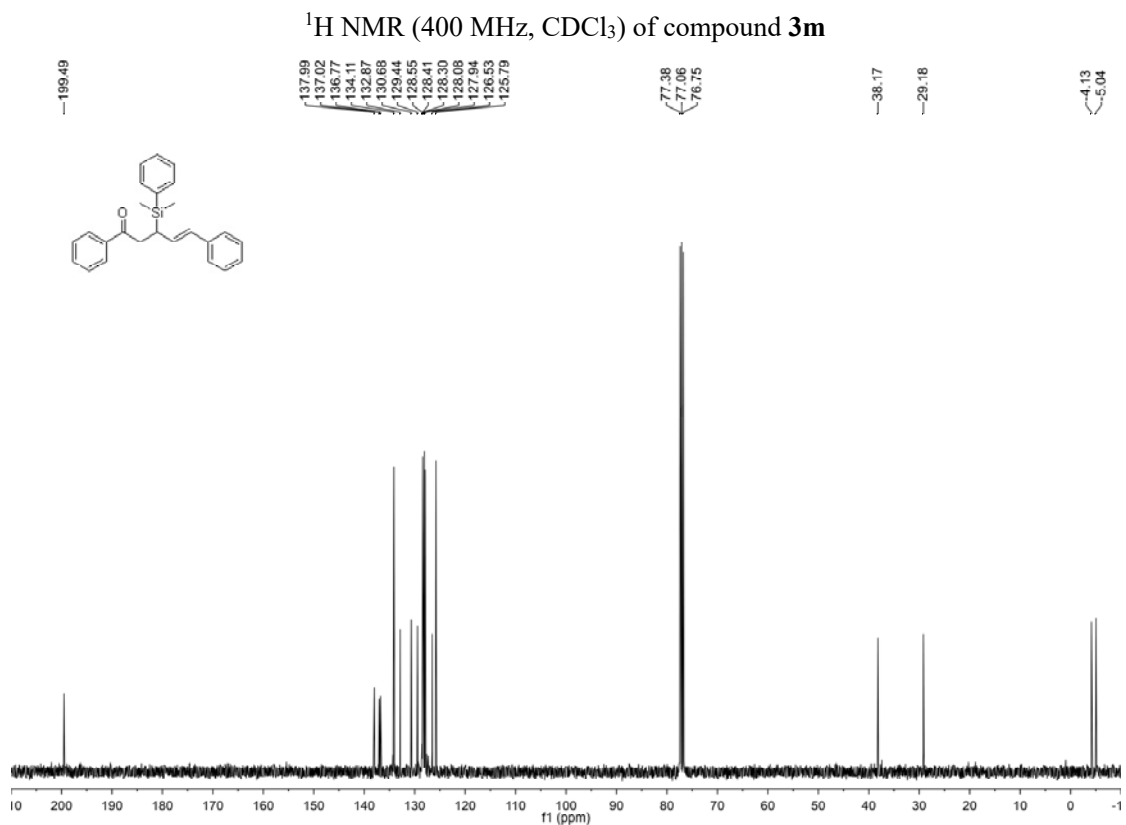
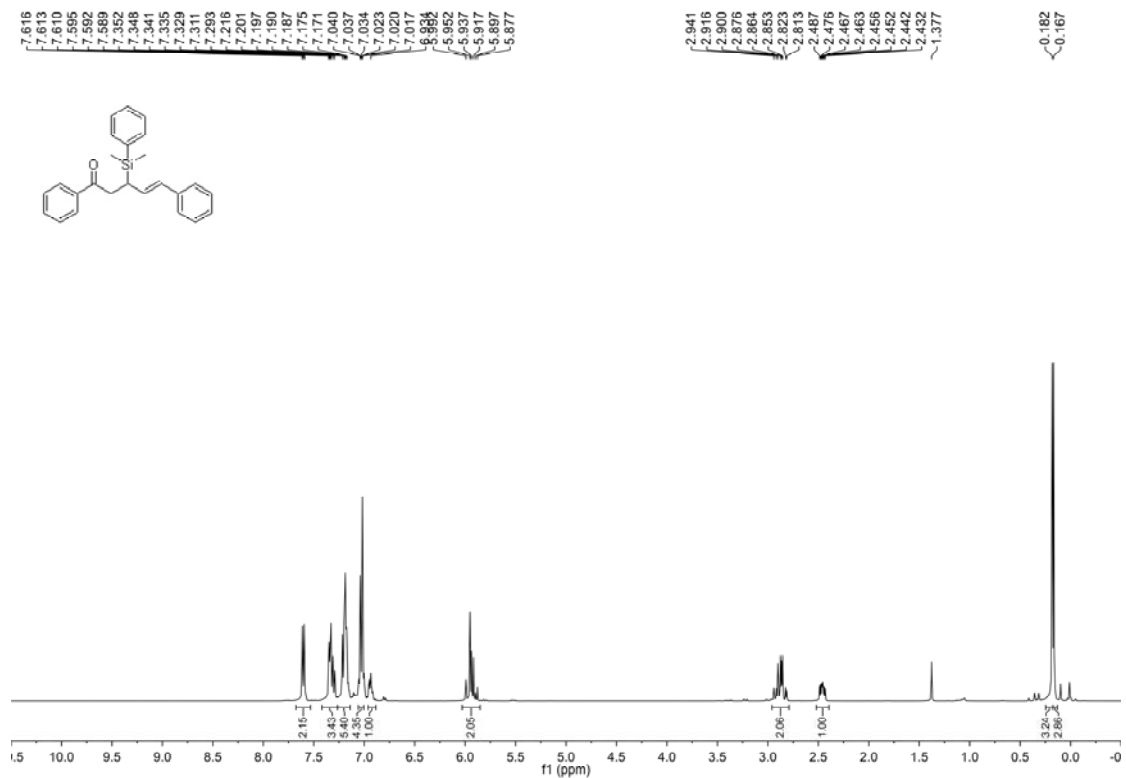


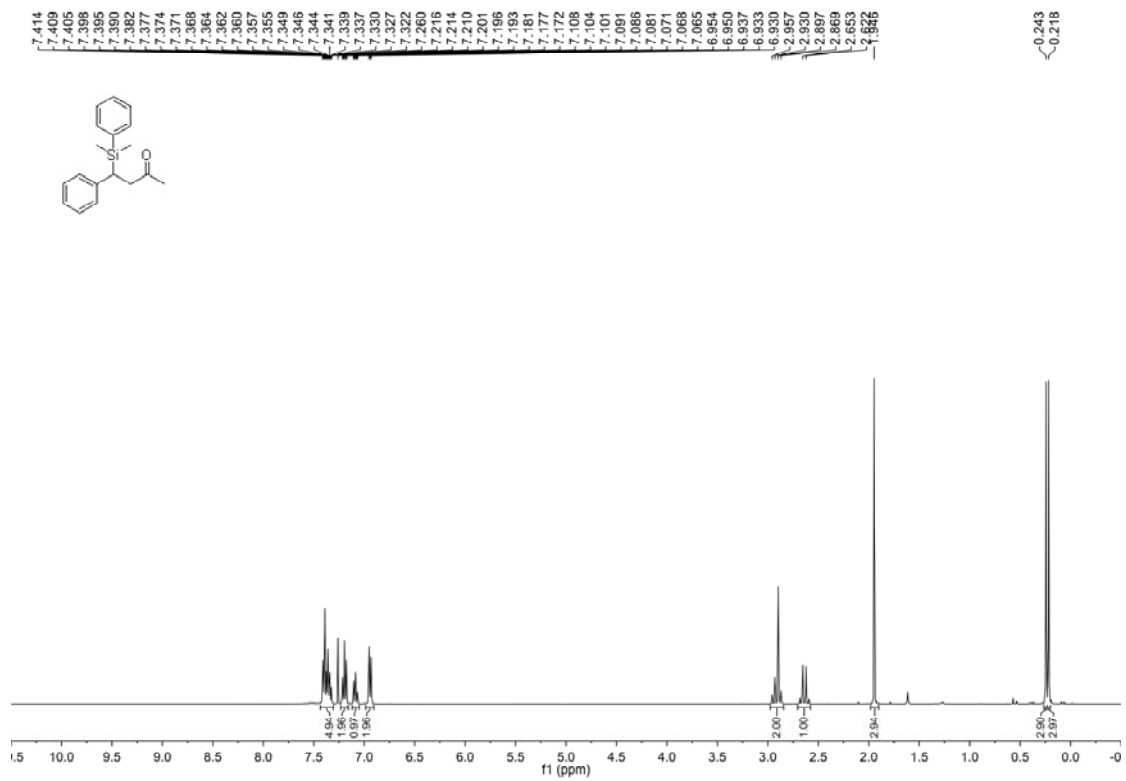




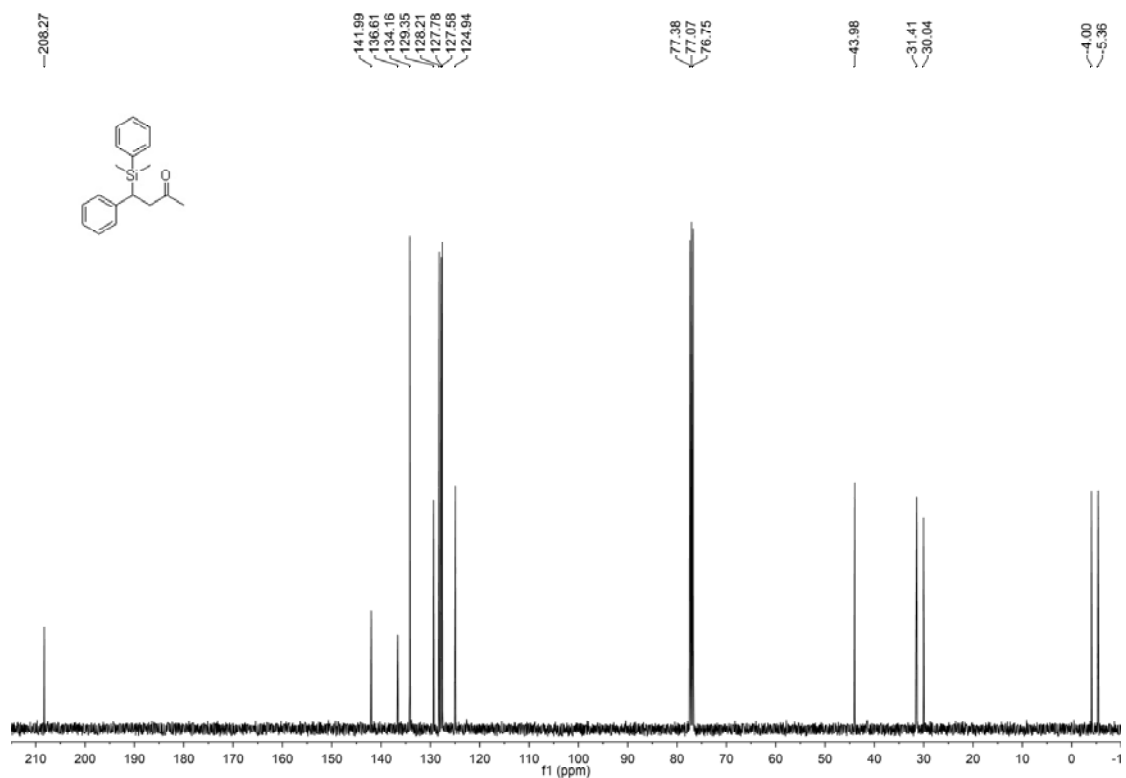


$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **31**

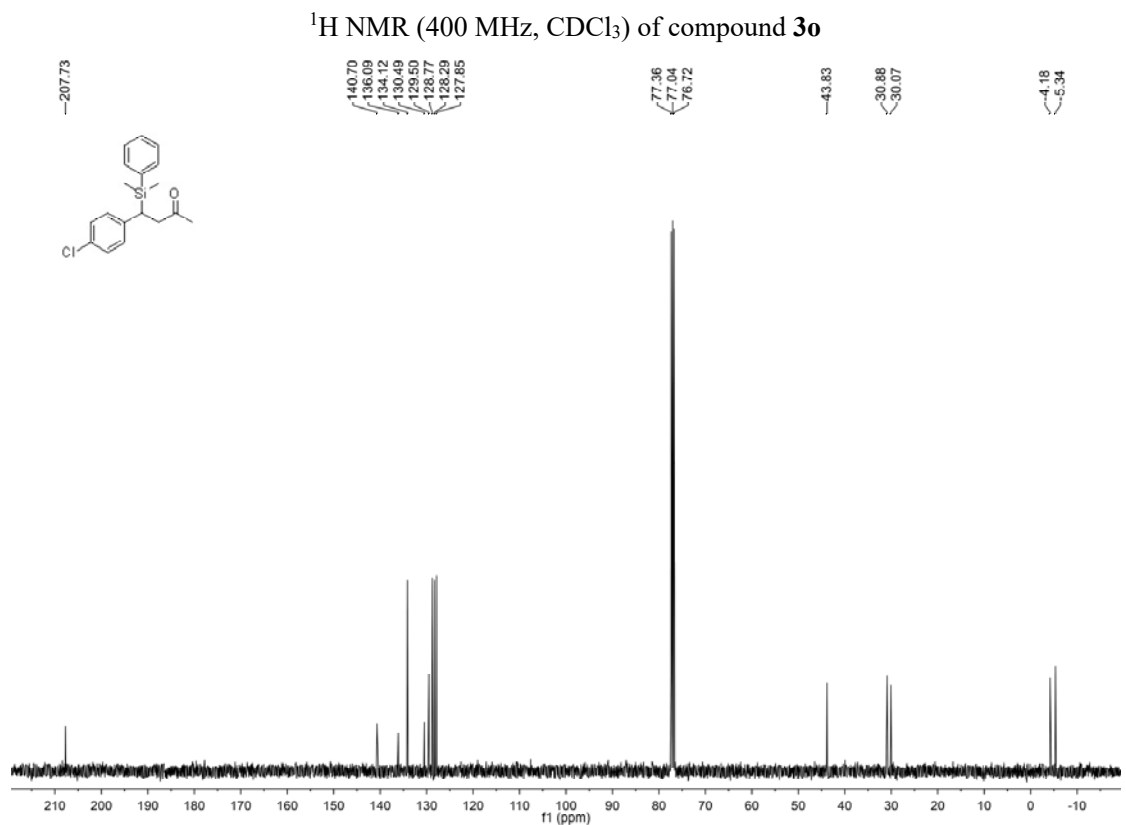
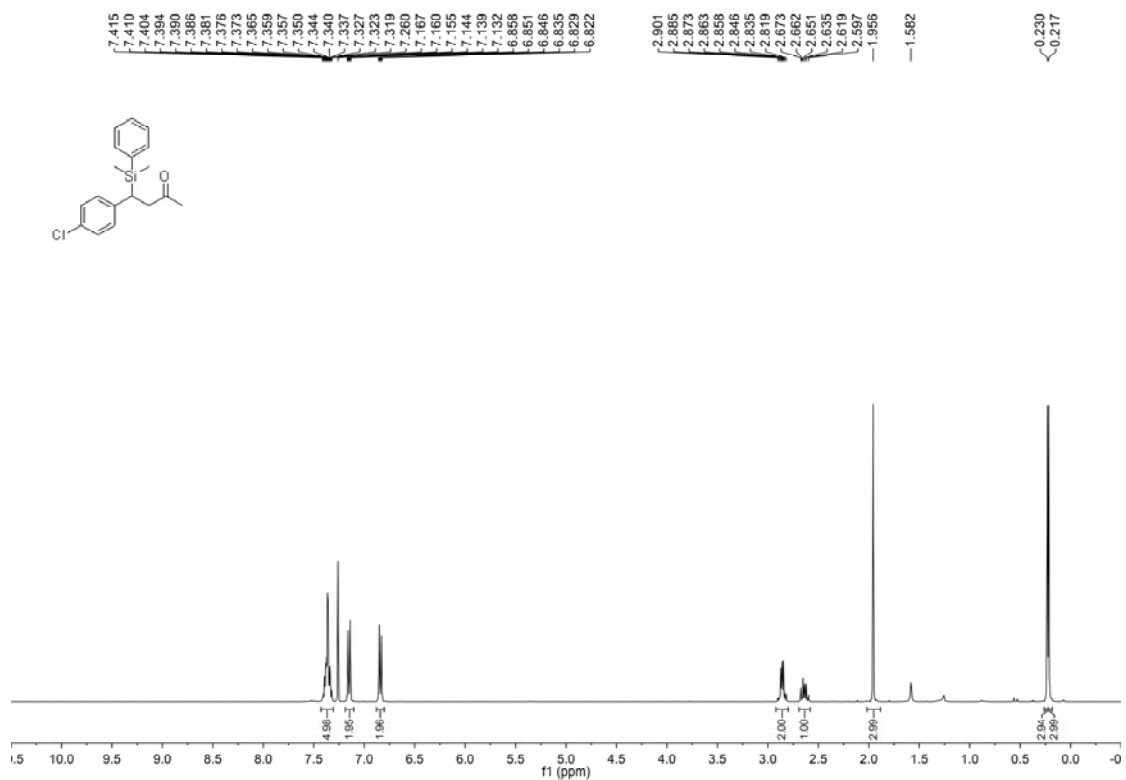




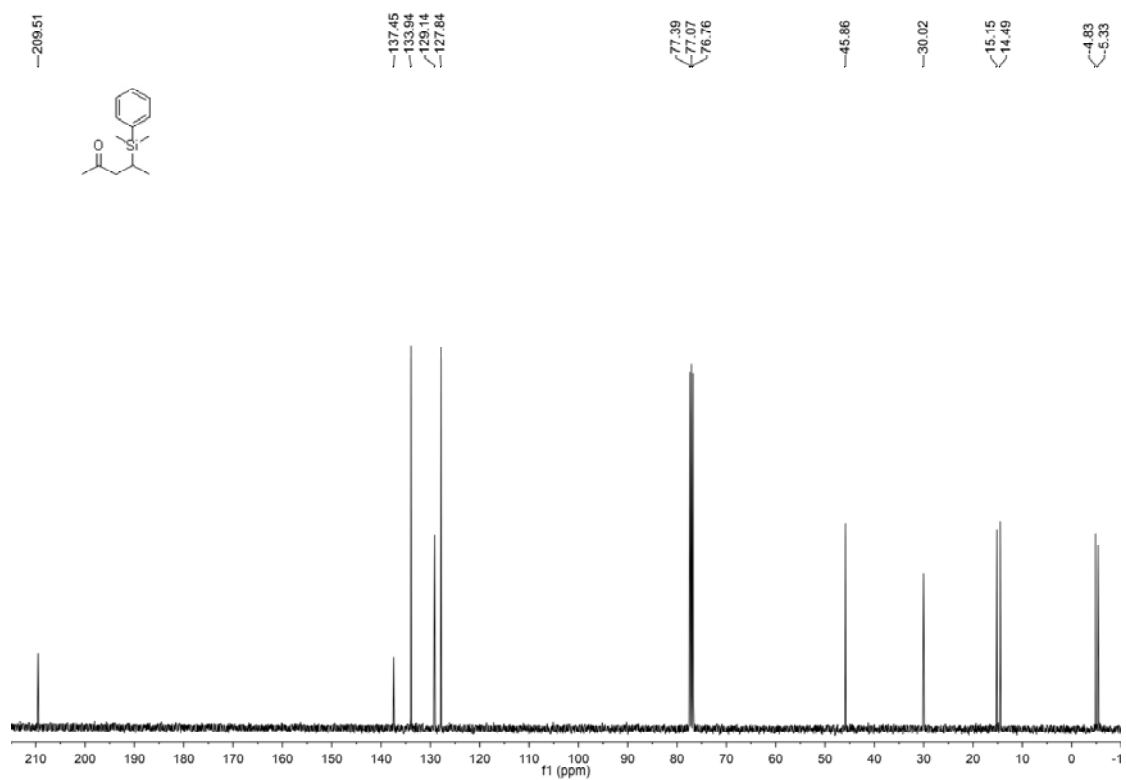
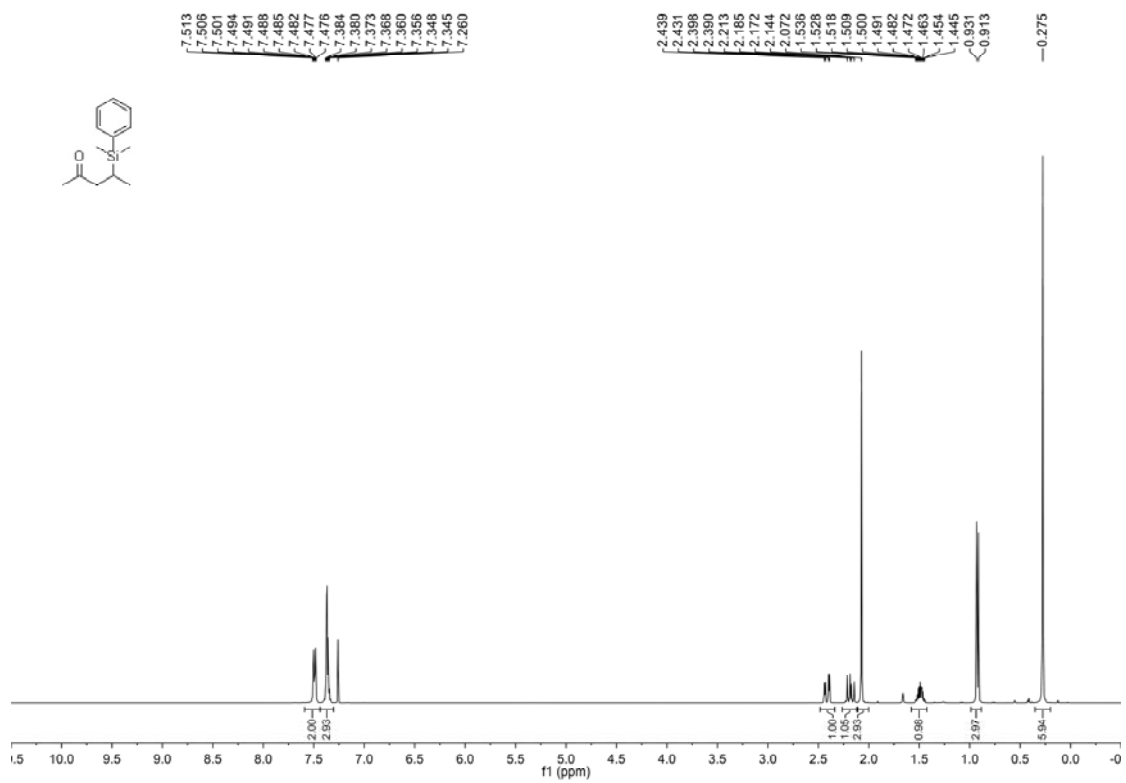
^1H NMR (400 MHz, CDCl_3) of compound **3n**

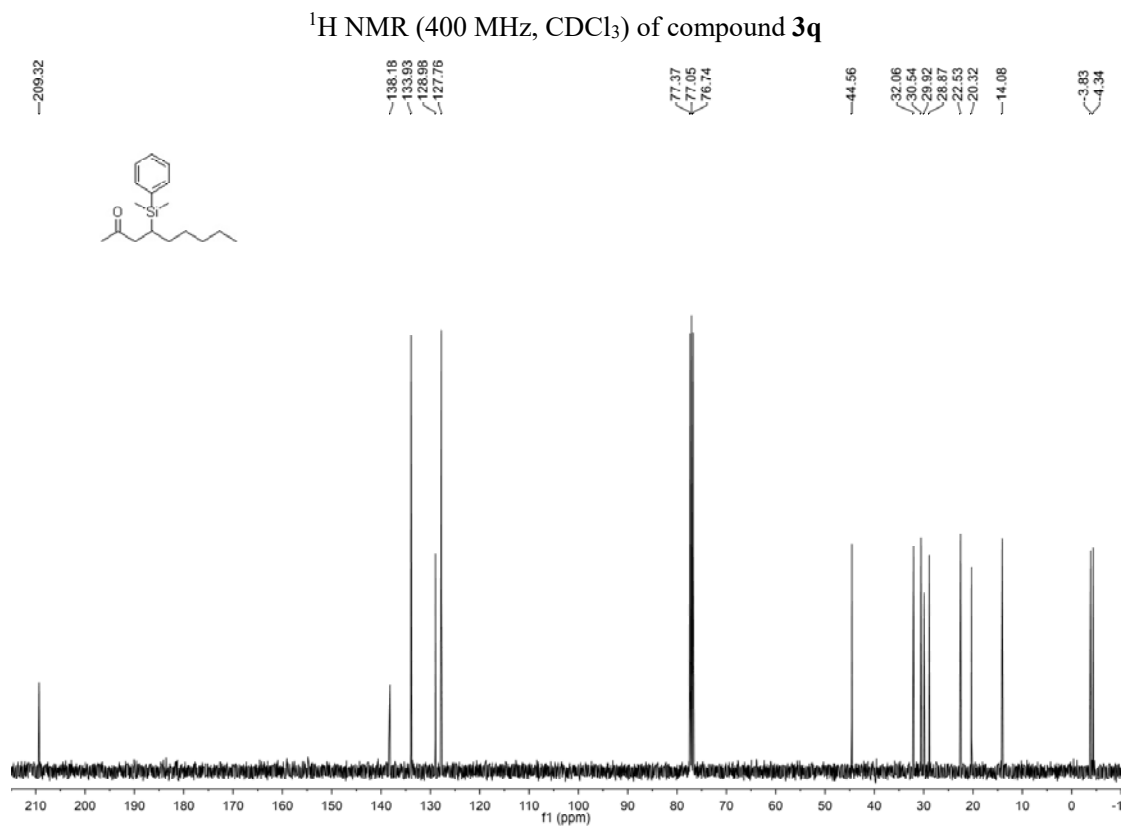
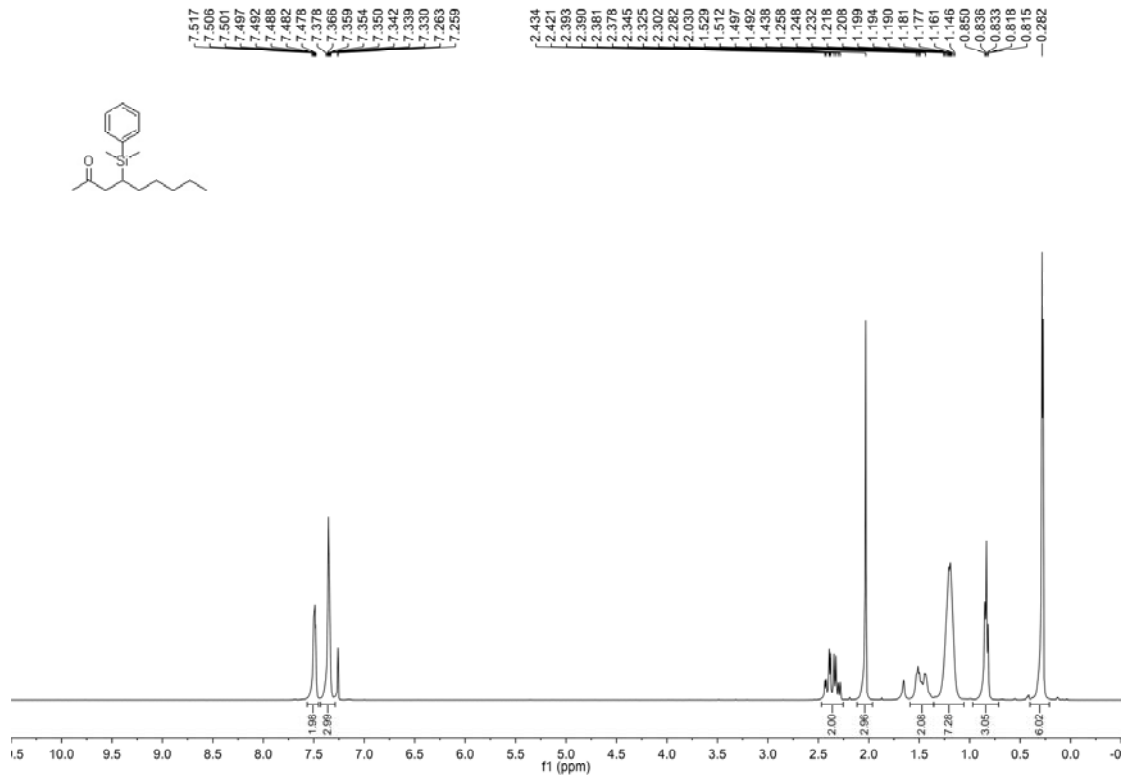


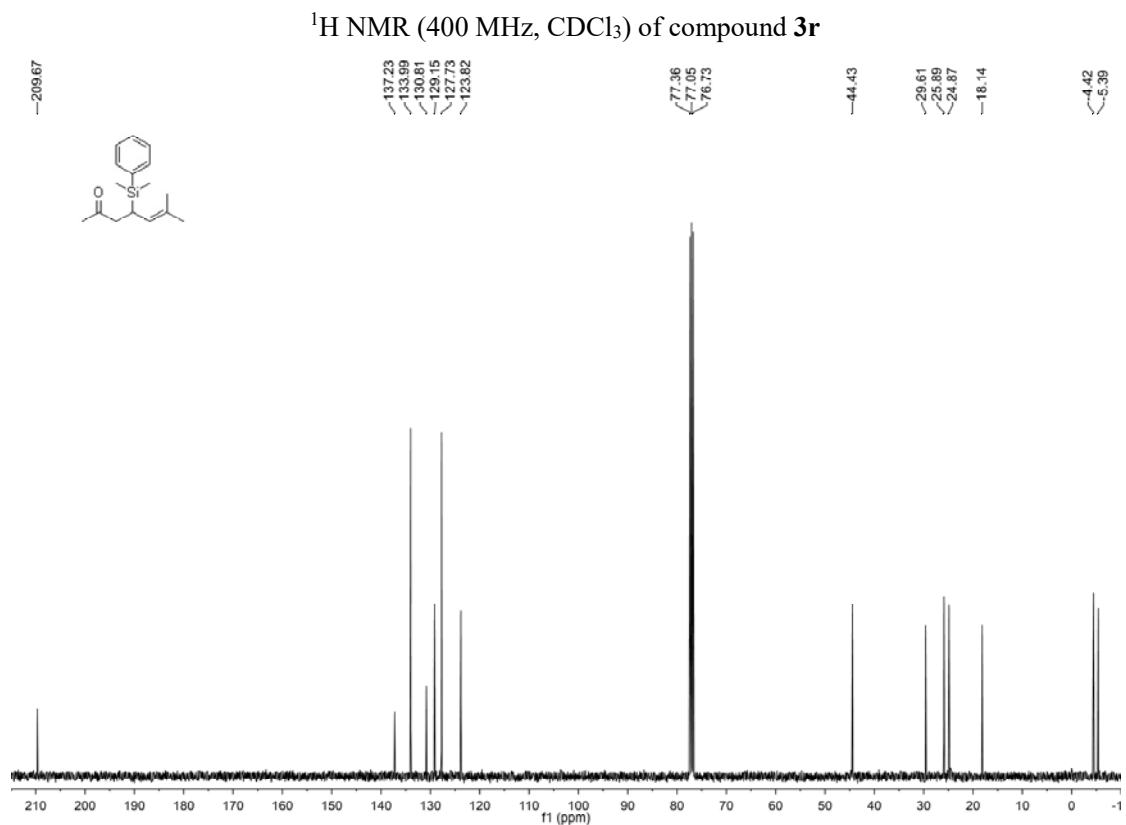
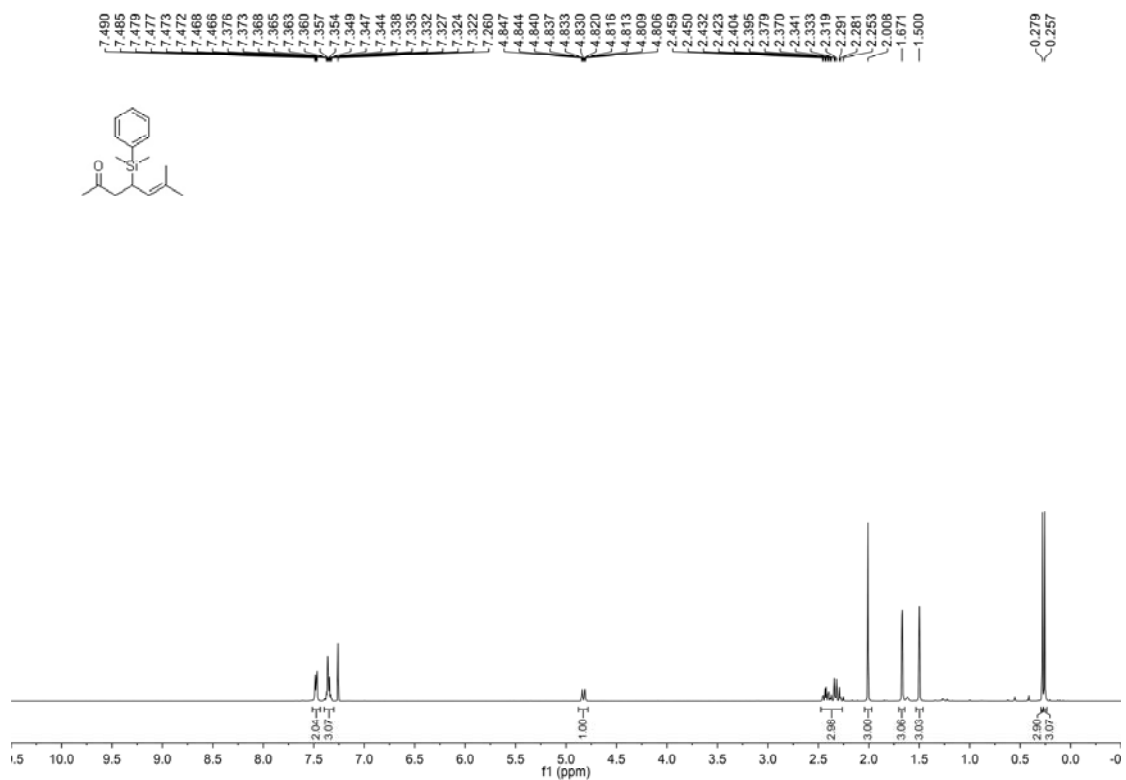
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **3n**

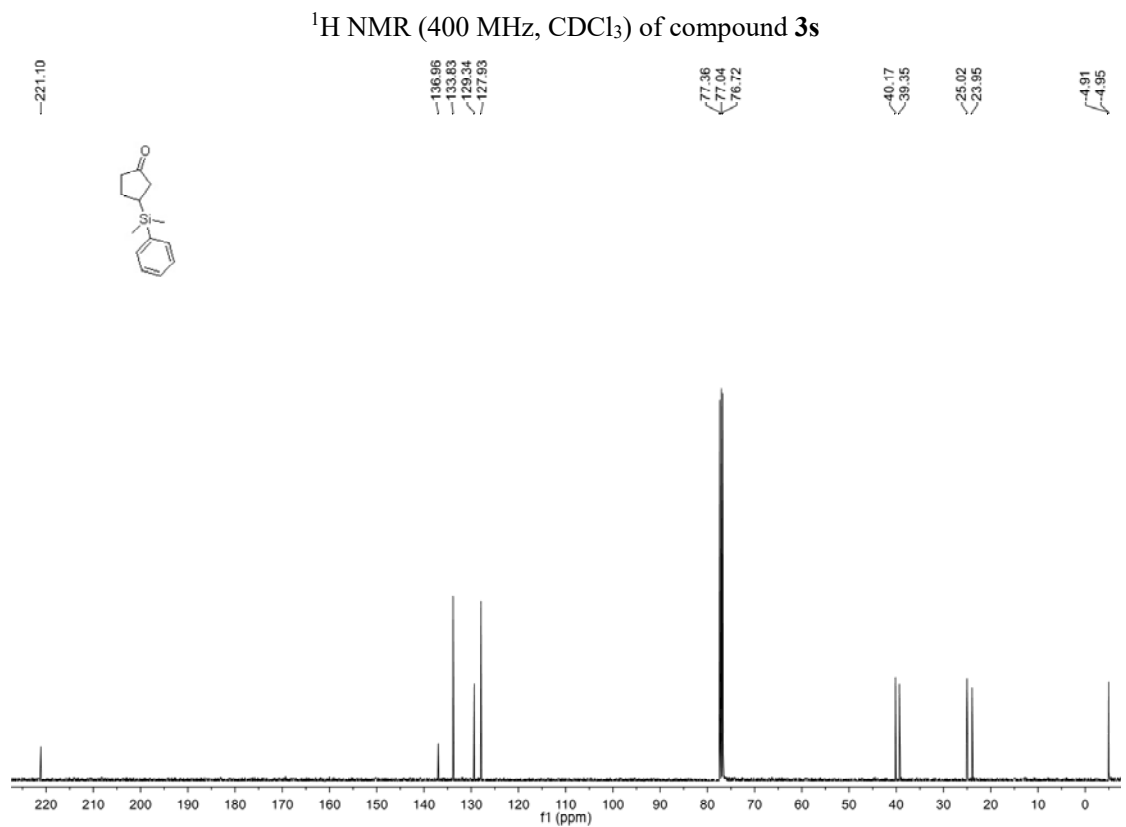
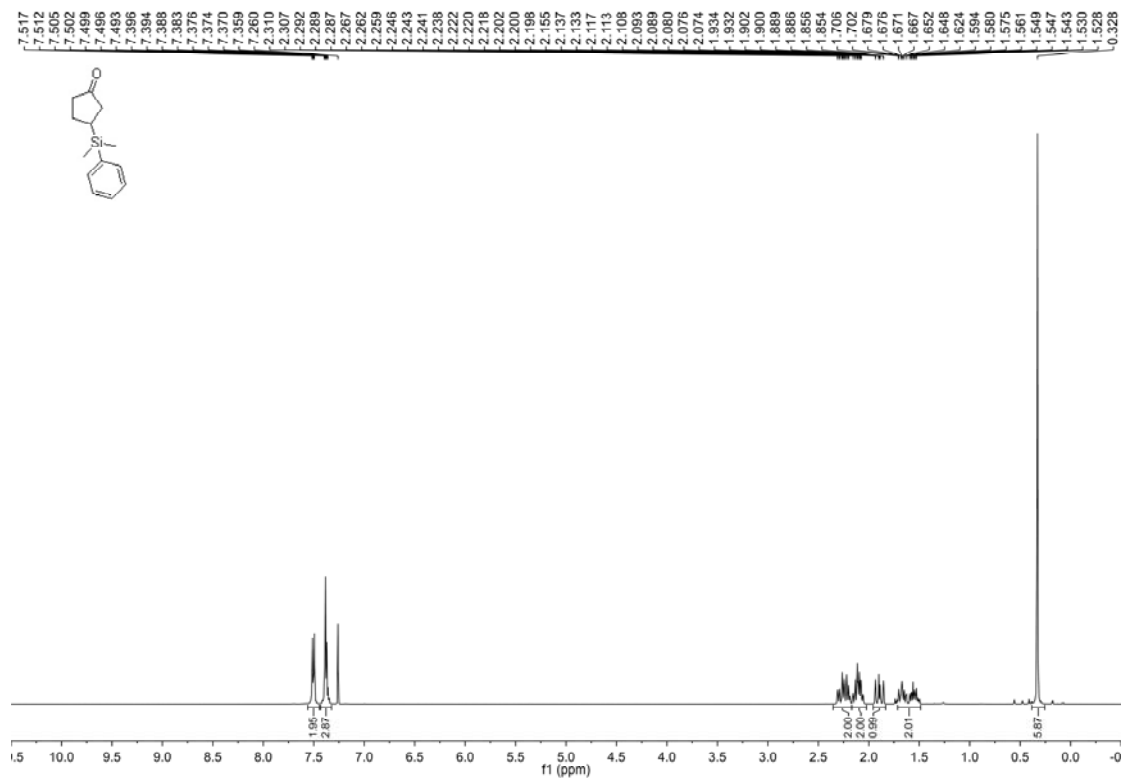


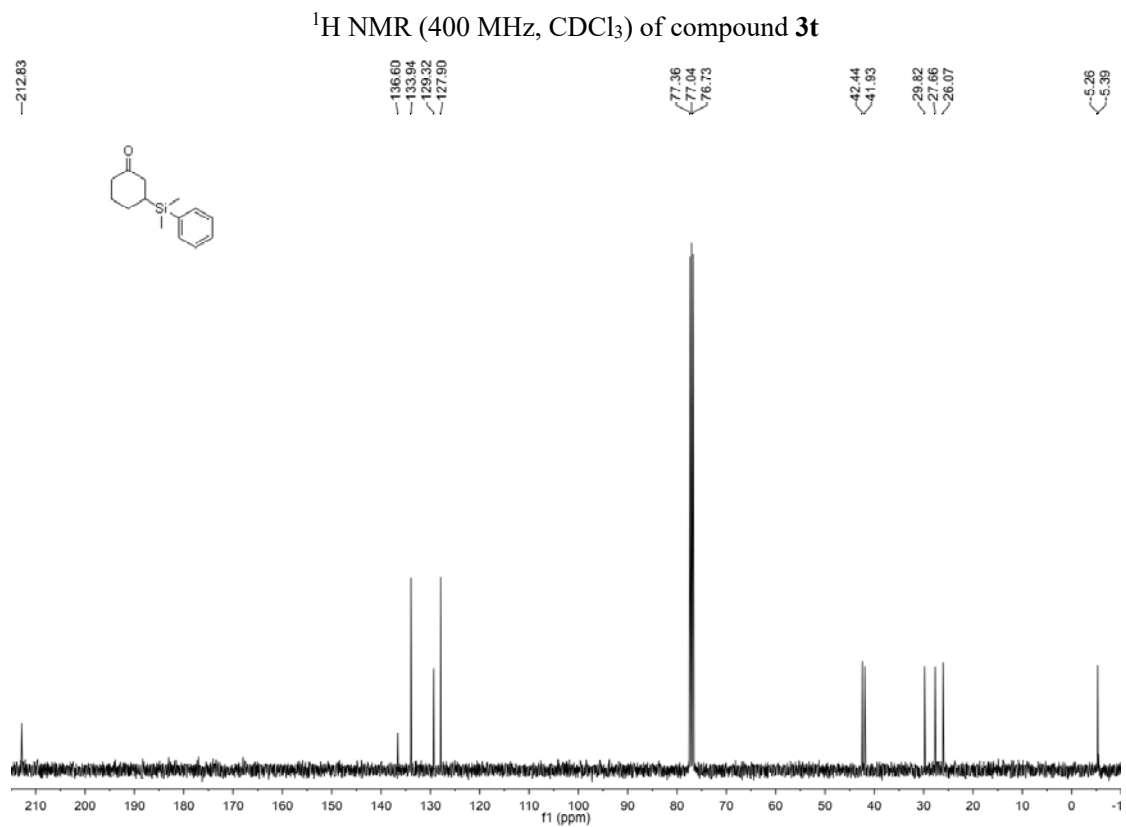
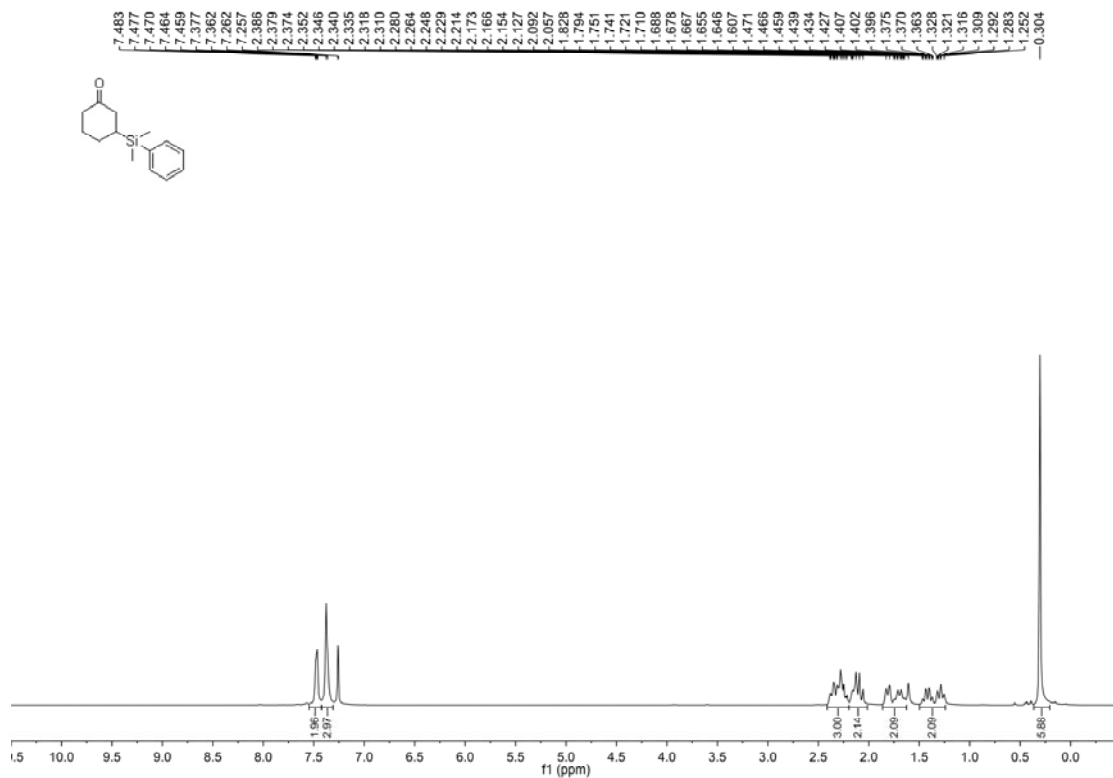
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **3o**

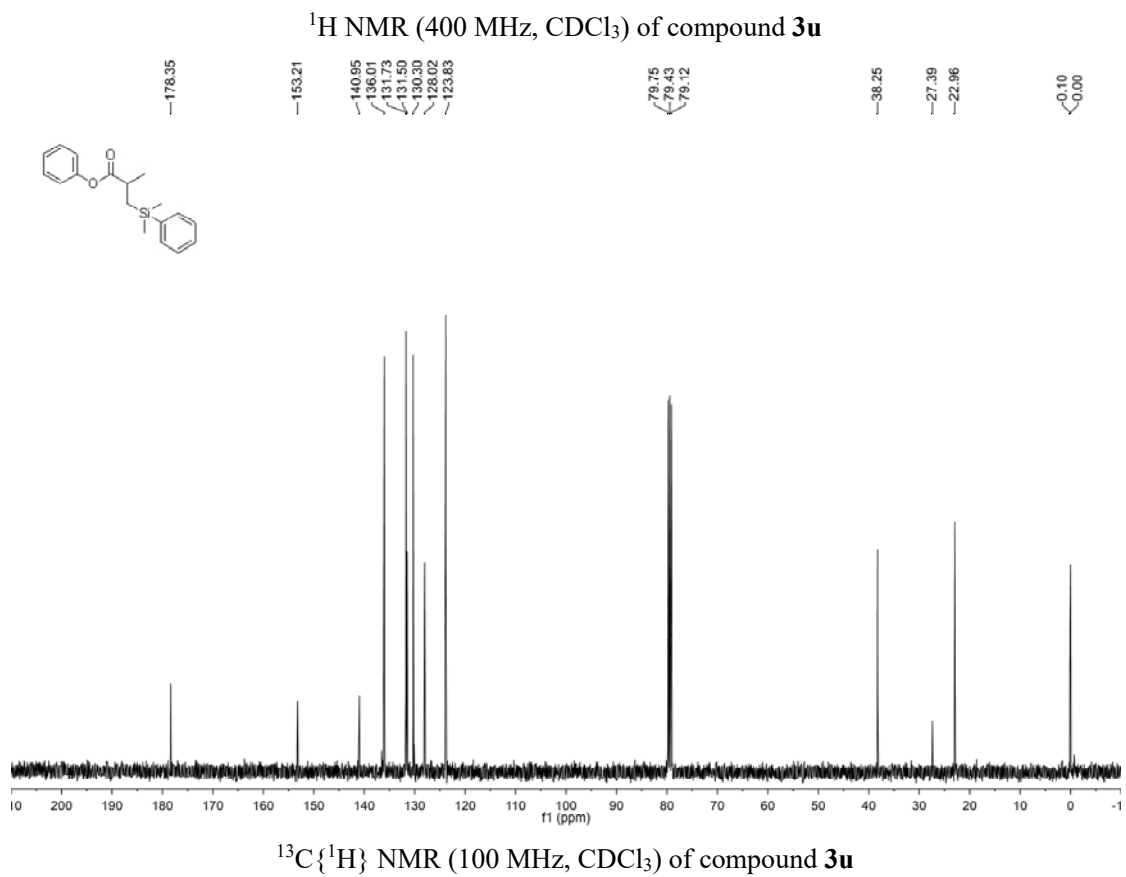
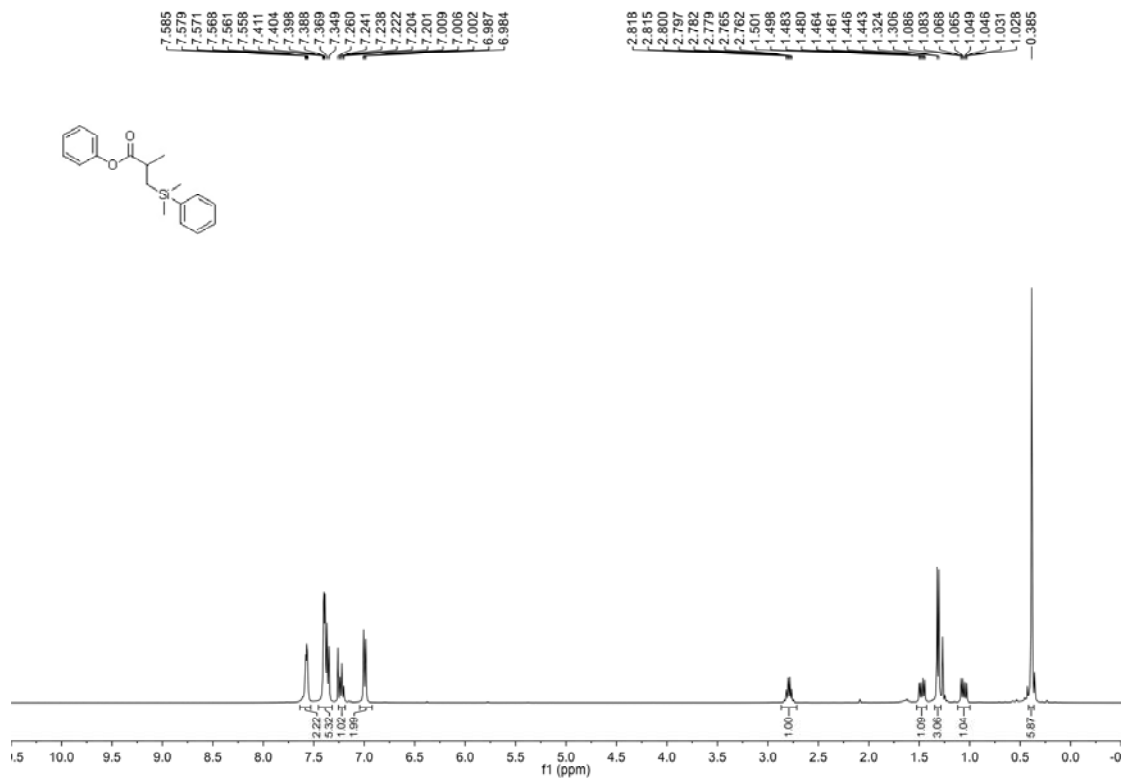


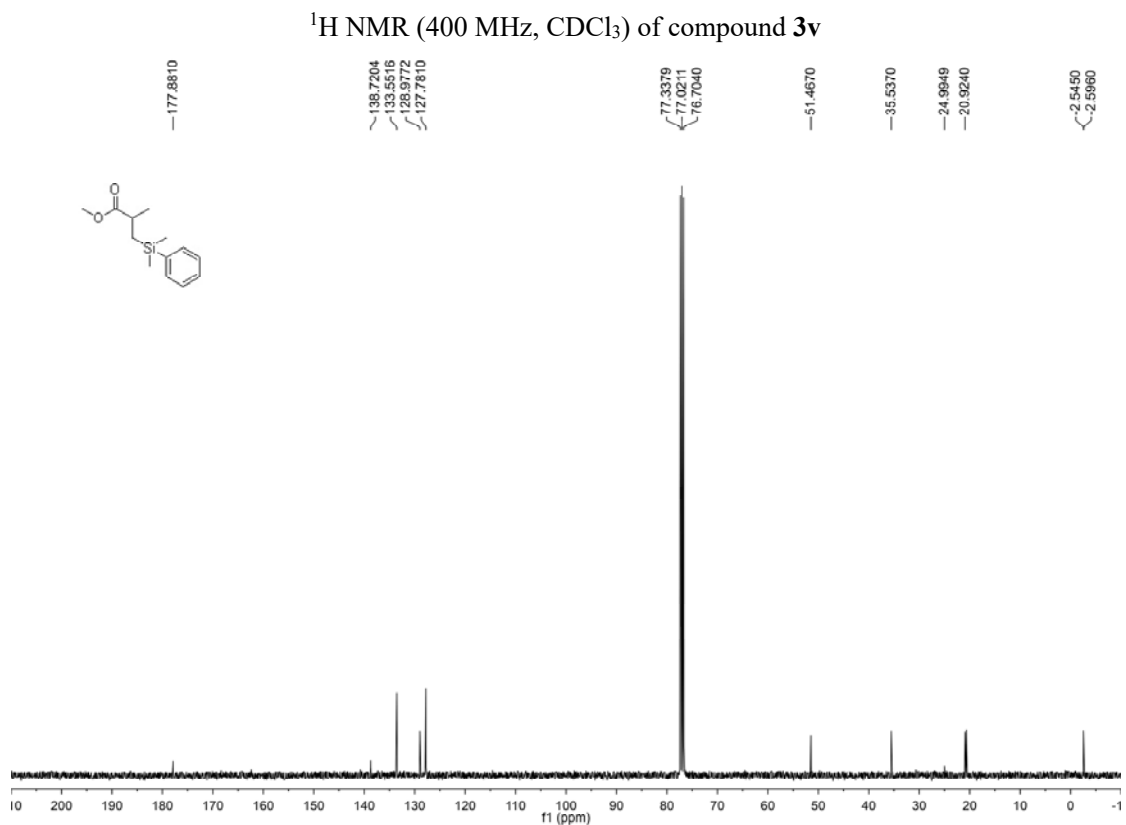
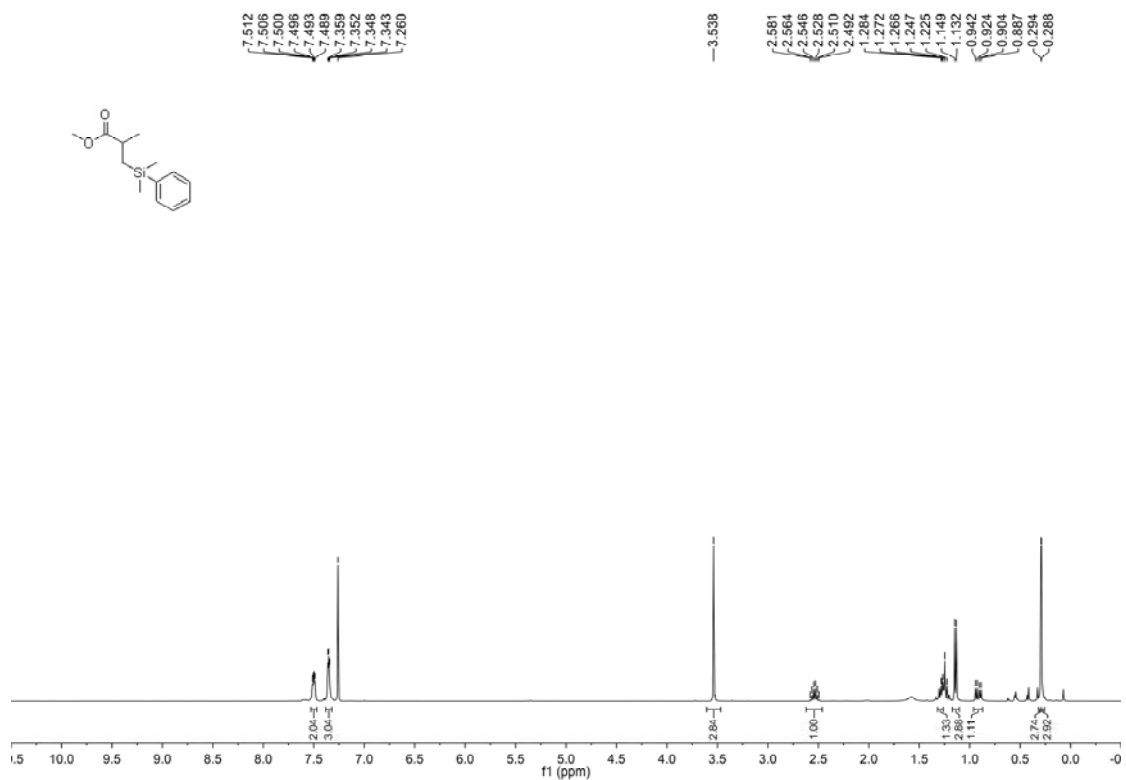




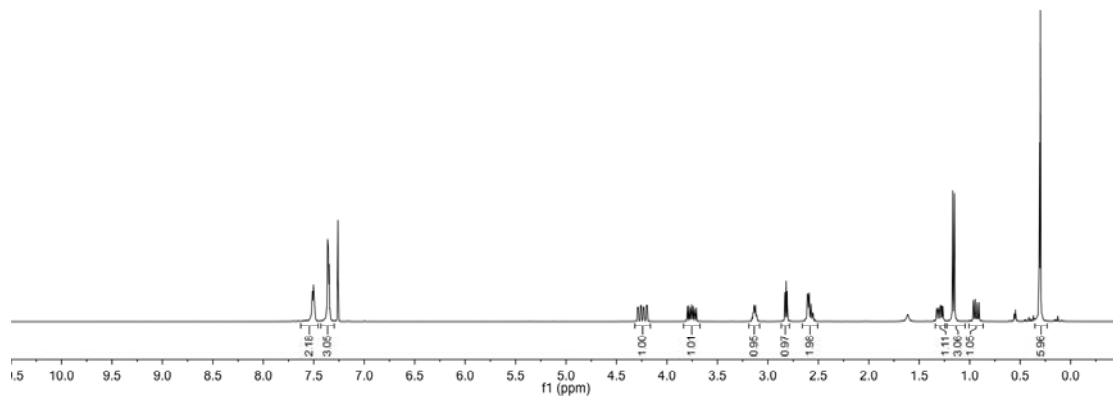
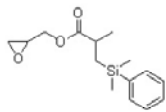






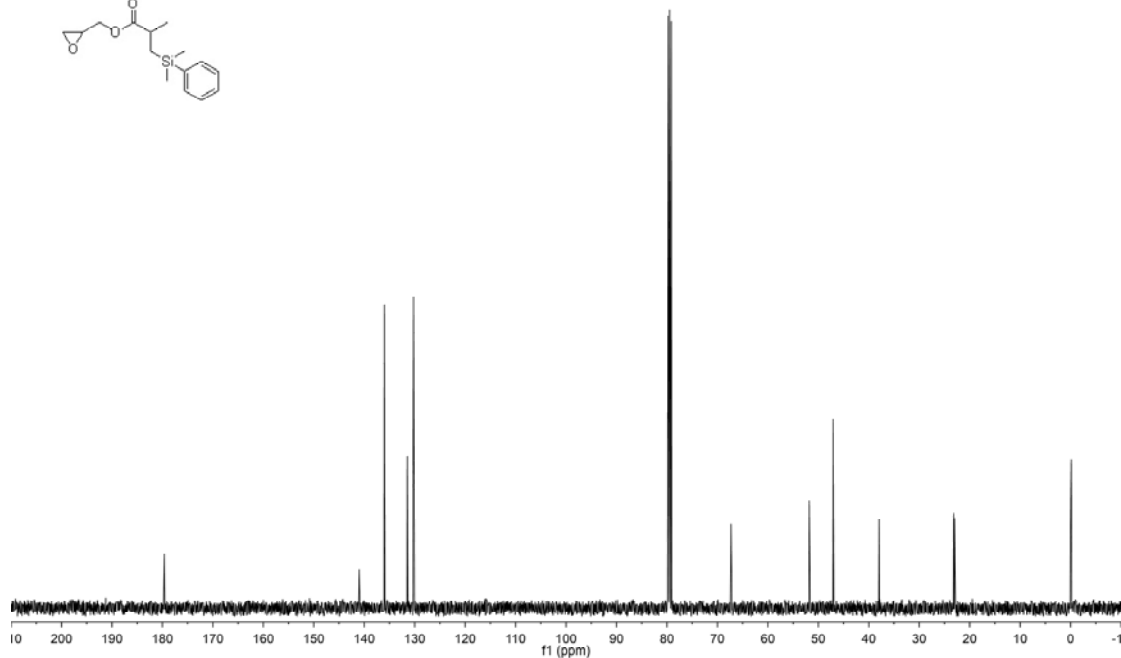
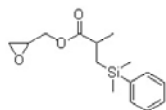


7.519
7.512
7.507
7.506
7.503
7.499
7.495
7.372
7.363
7.356
7.352
7.347
7.341
7.261
4.282
4.284
4.281
4.234
4.227
4.204
4.196
3.799
3.783
3.768
3.755
3.752
3.740
3.725
3.709
3.139
3.132
3.129
3.122
2.831
2.820
2.819
2.808
2.613
2.609
2.605
2.601
2.595
2.594
2.584
2.580
2.583
2.573
2.570
1.329
1.321
1.310
1.292
1.284
1.273
1.265
1.168
1.150
0.963
0.962
0.944
0.926
0.925
0.907
0.319
0.310
0.306
0.289
0.281

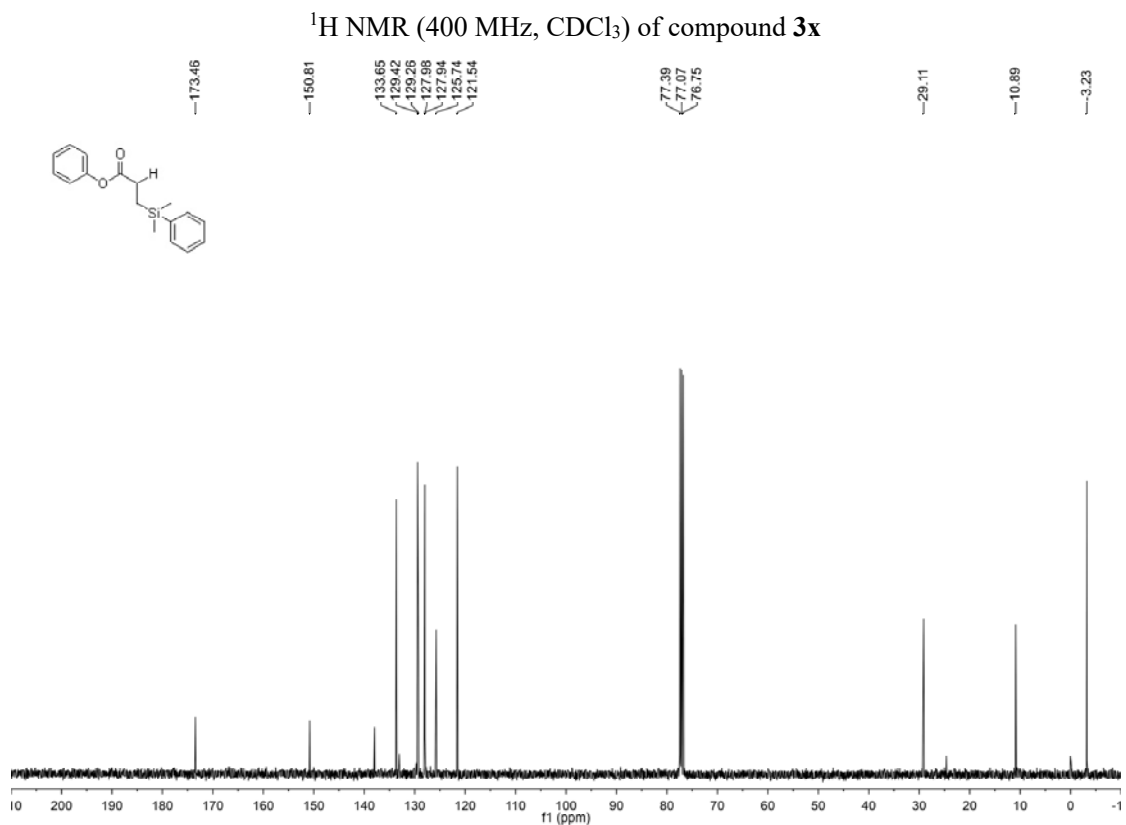
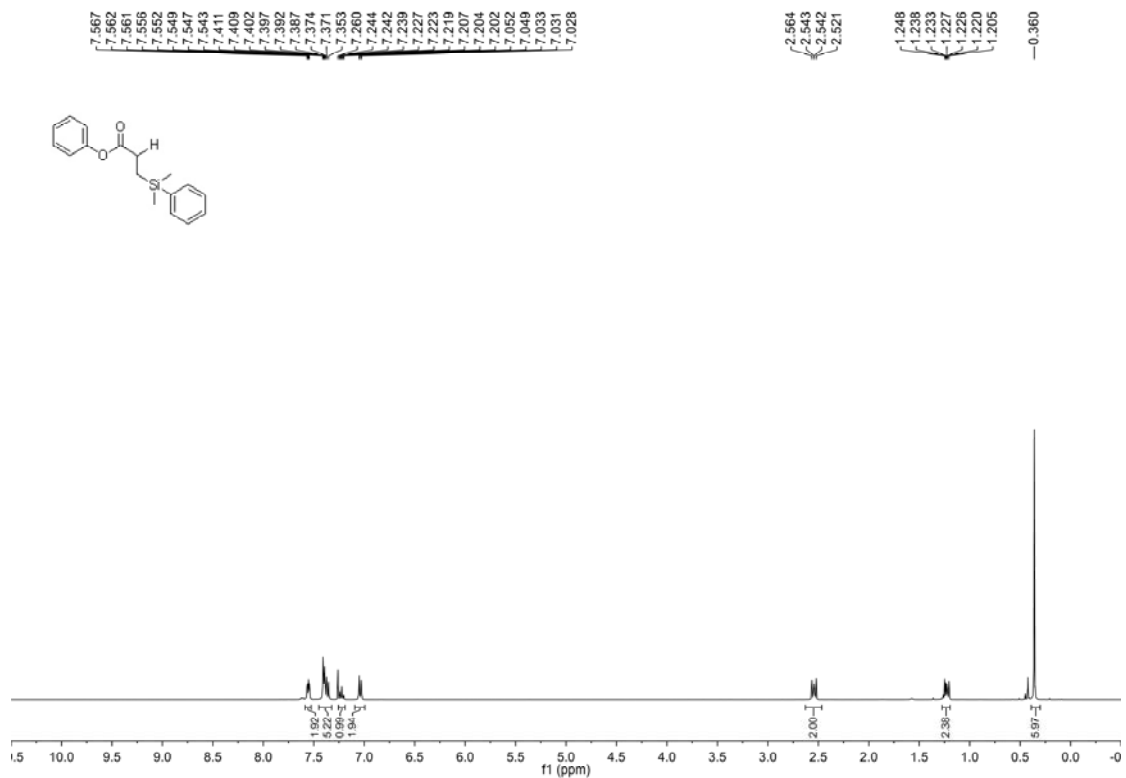


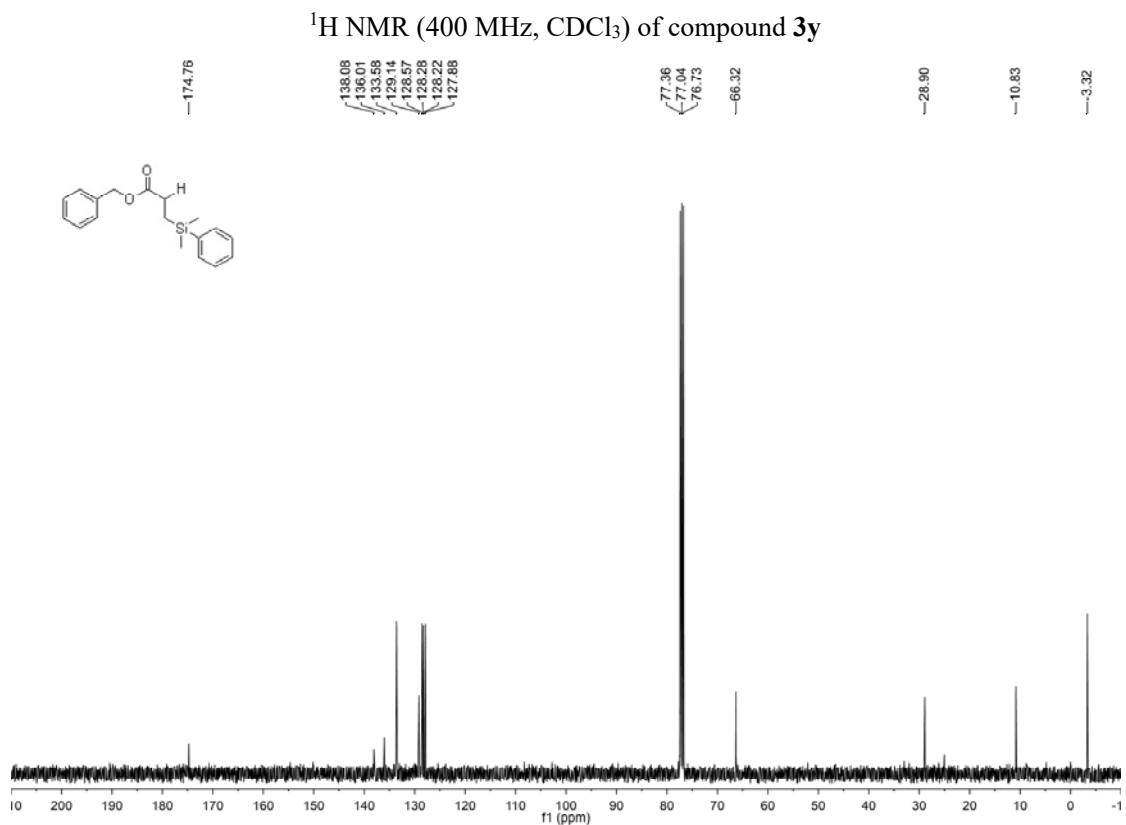
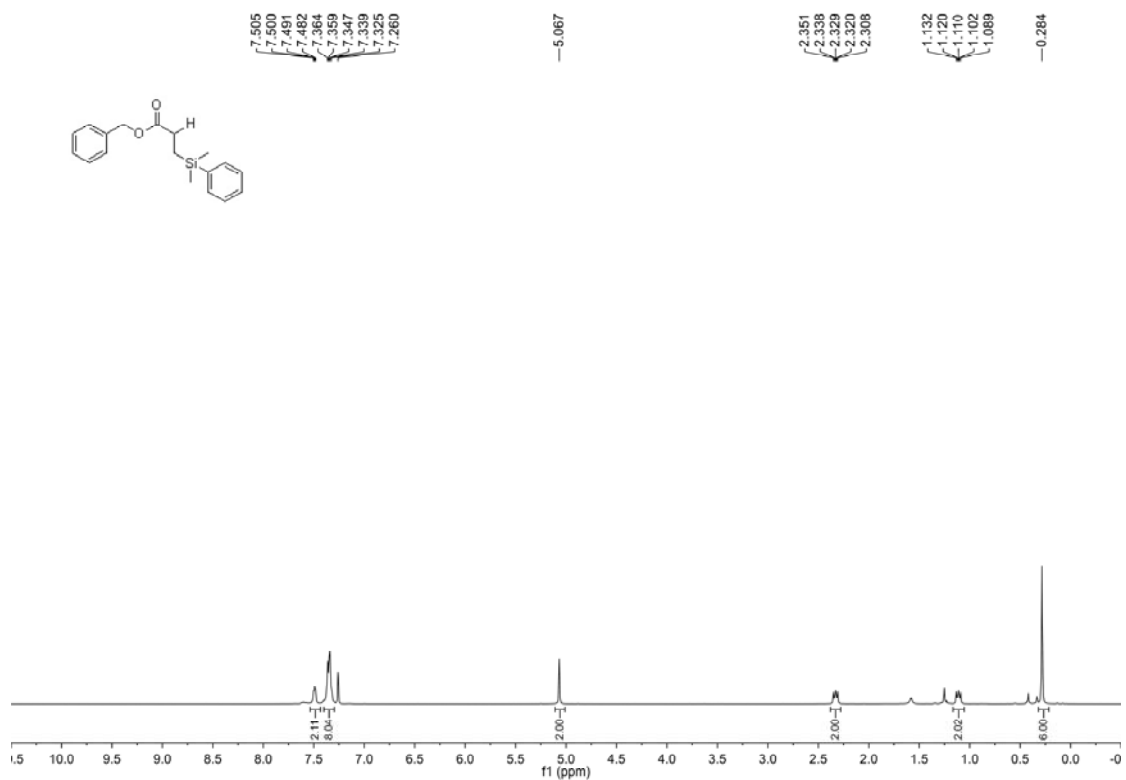
^1H NMR (400 MHz, CDCl_3) of compound **3w**

-179.62
141.00
136.01
131.46
130.24
79.79
79.47
79.15
-67.30
-51.79
-47.06
-37.96
-23.17
0.00
-0.13

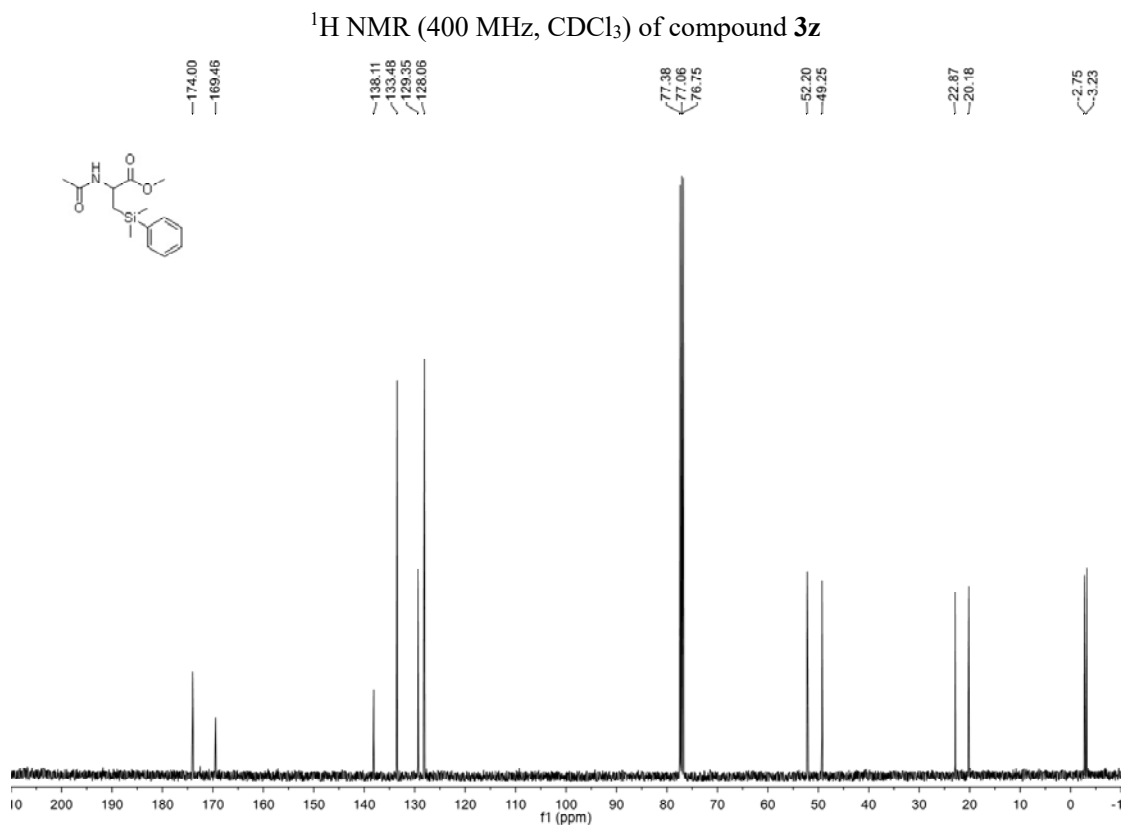
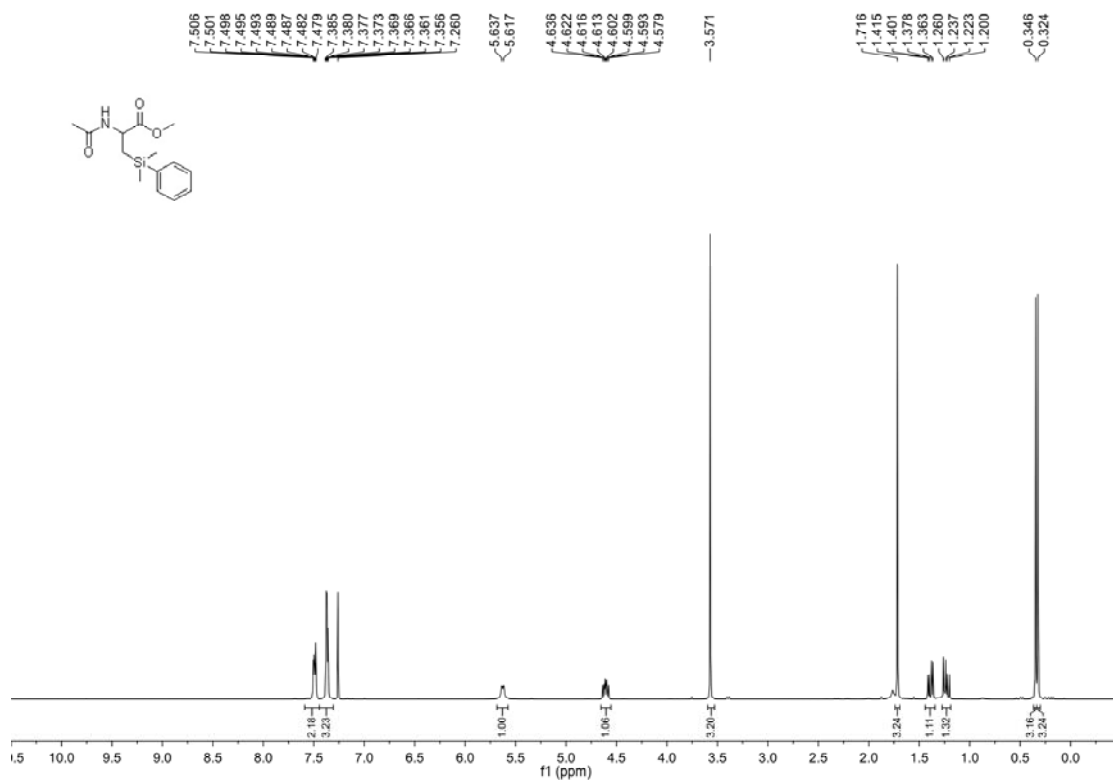


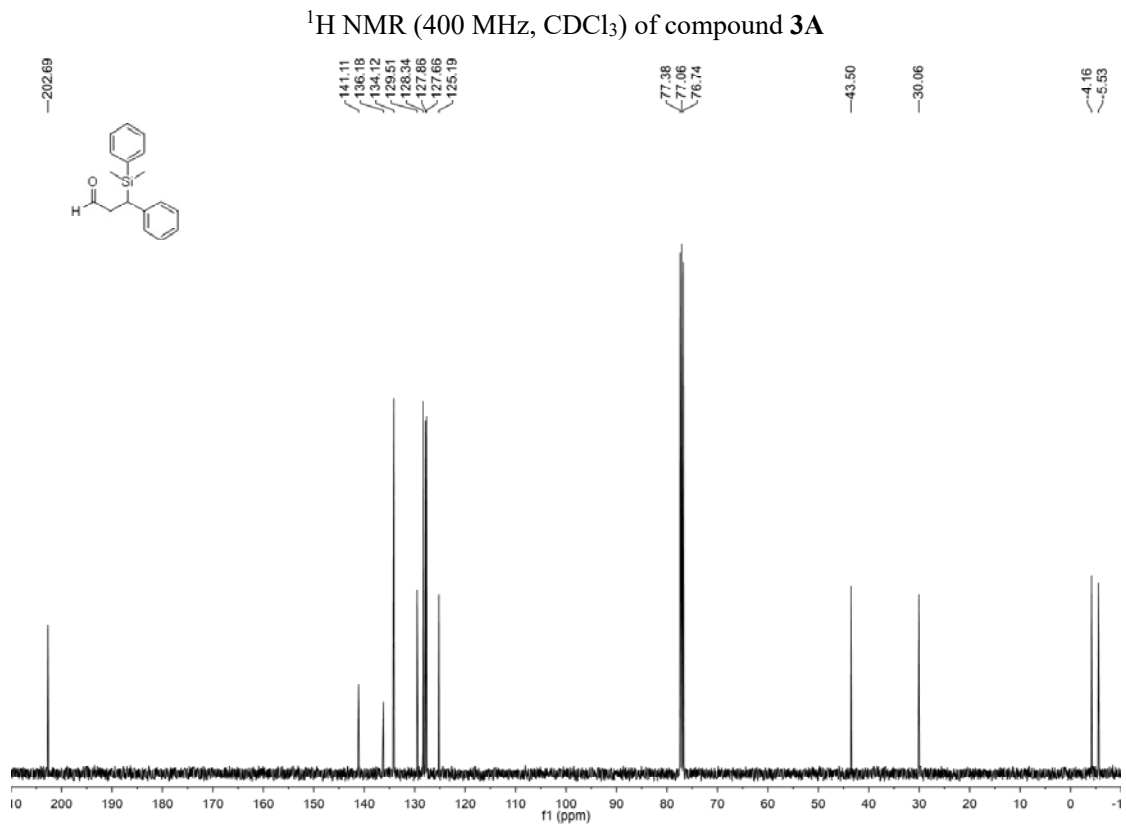
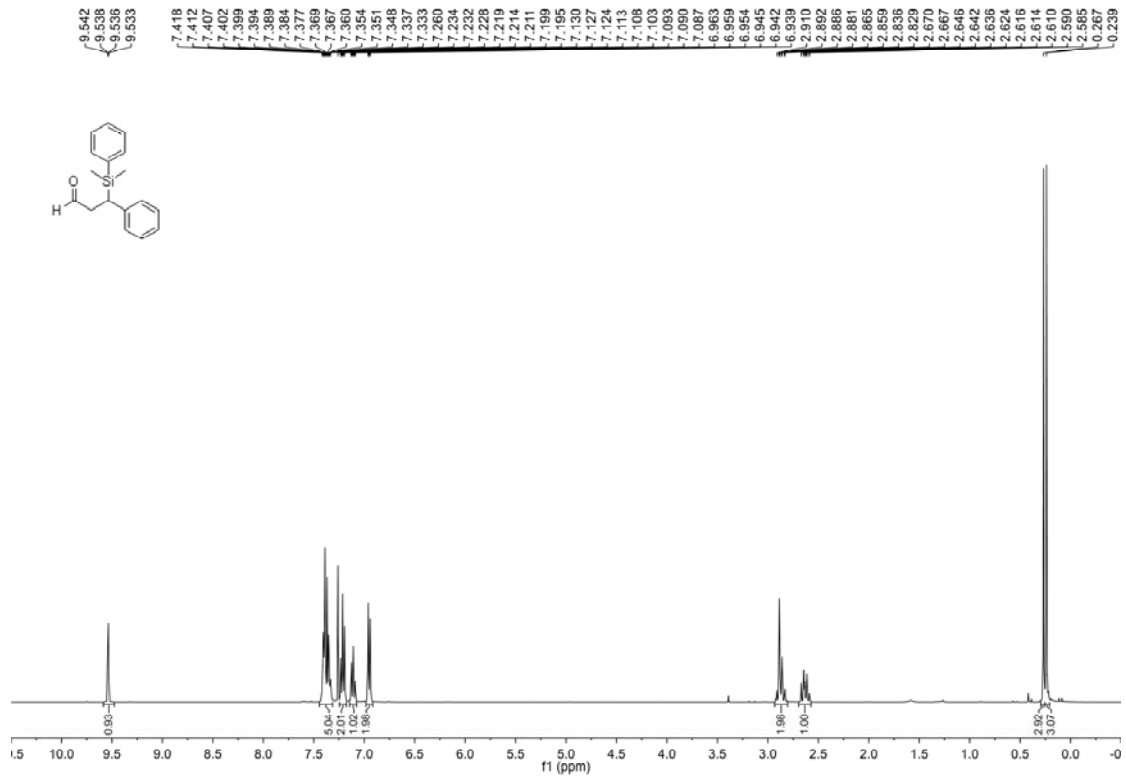
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **3w**

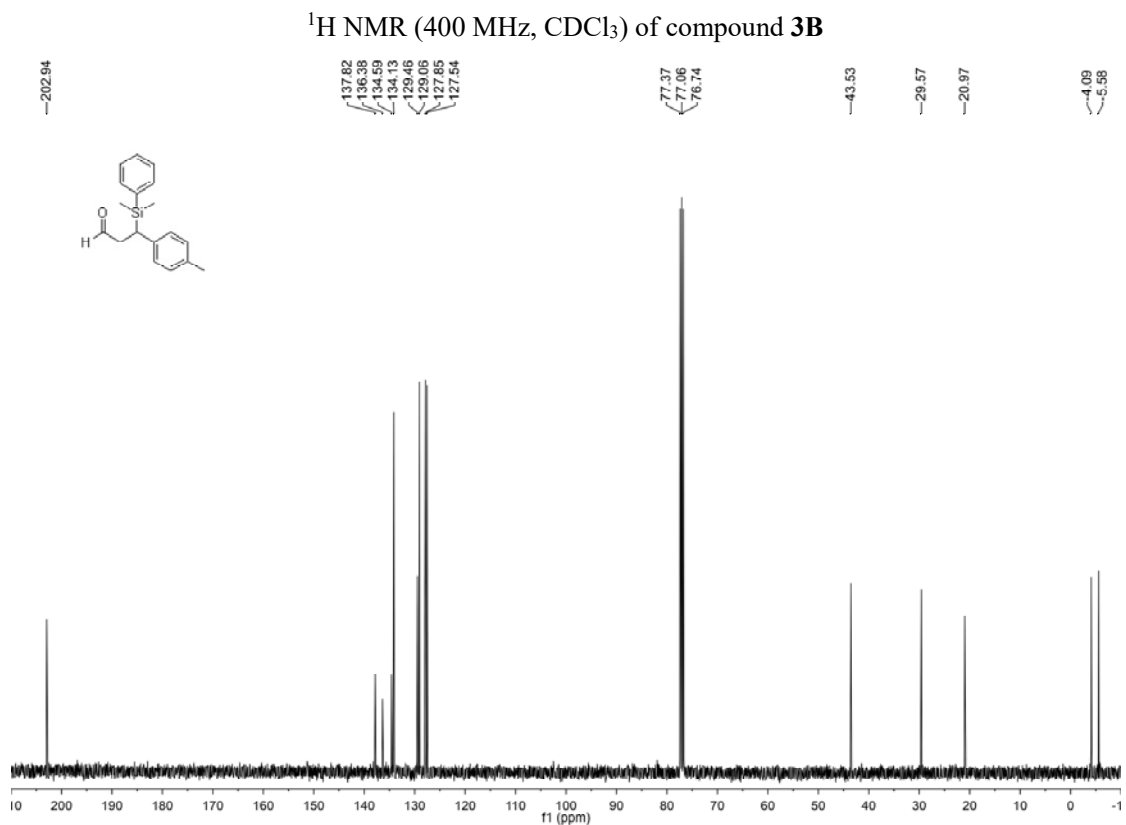
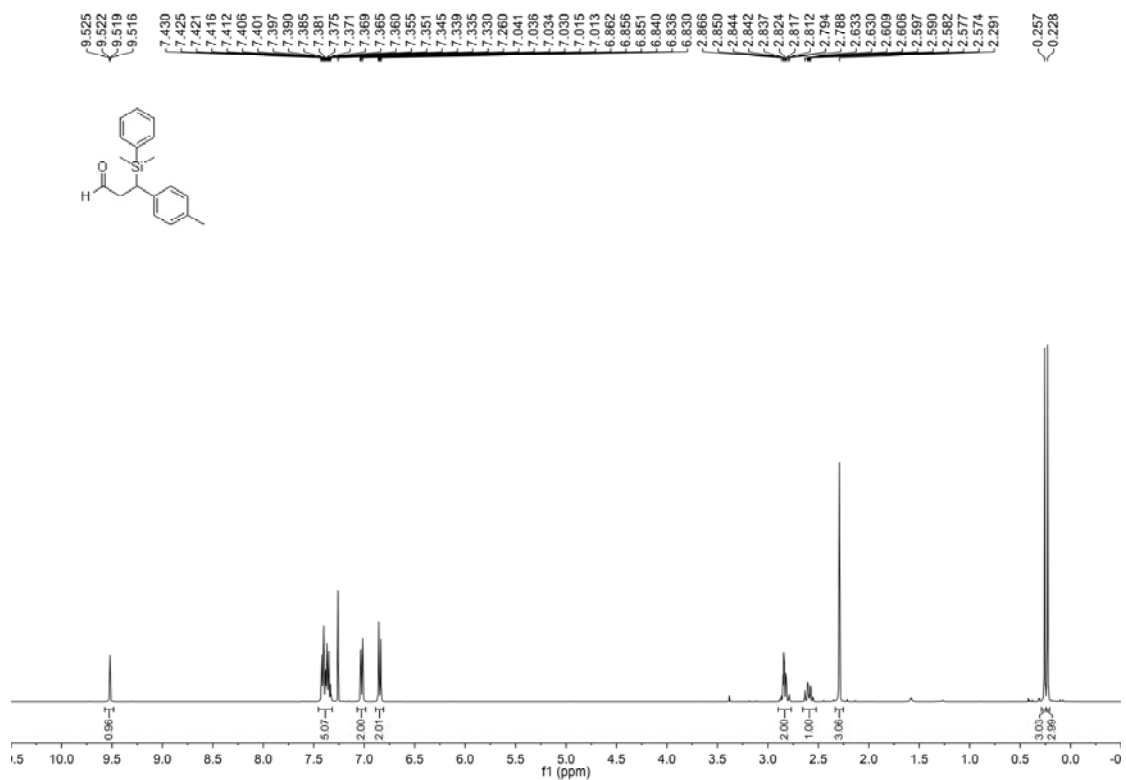


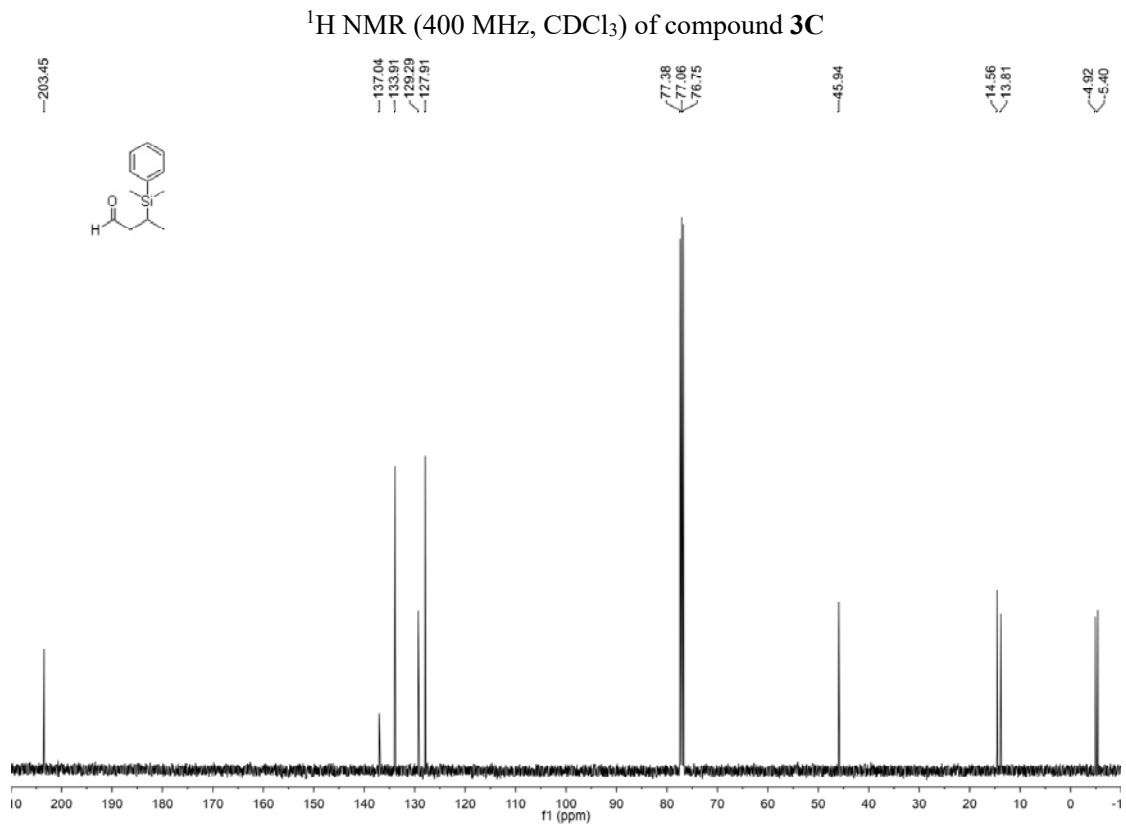
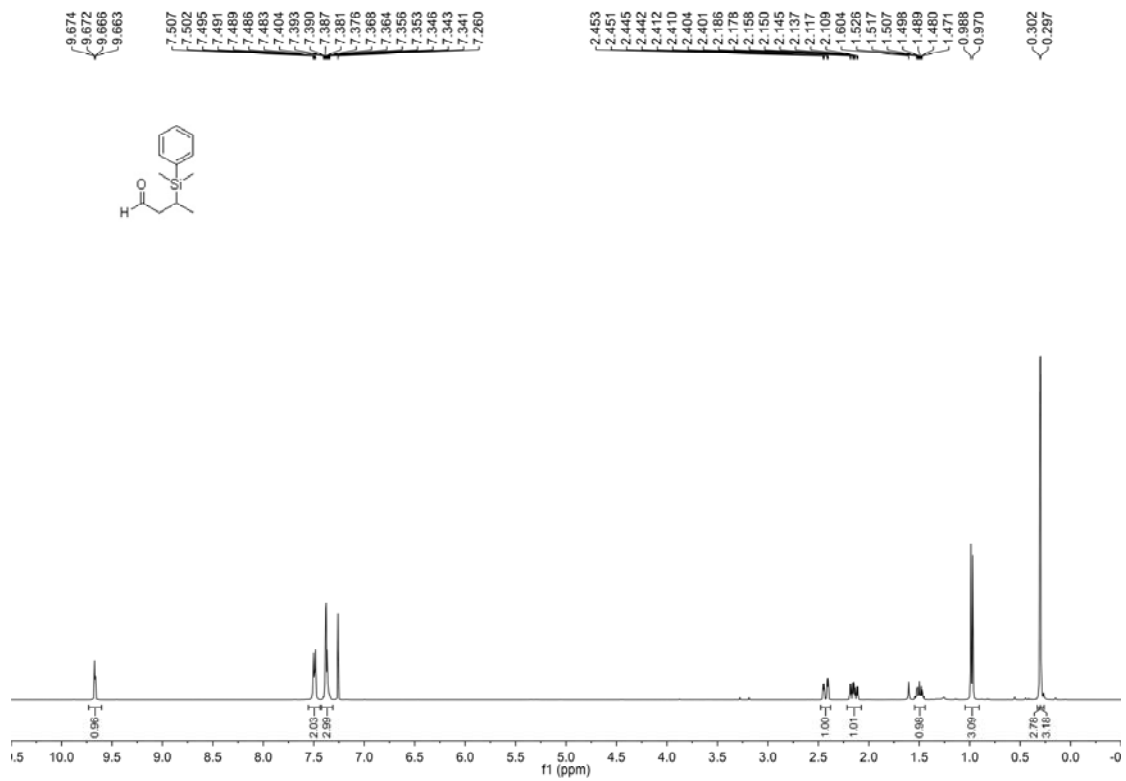


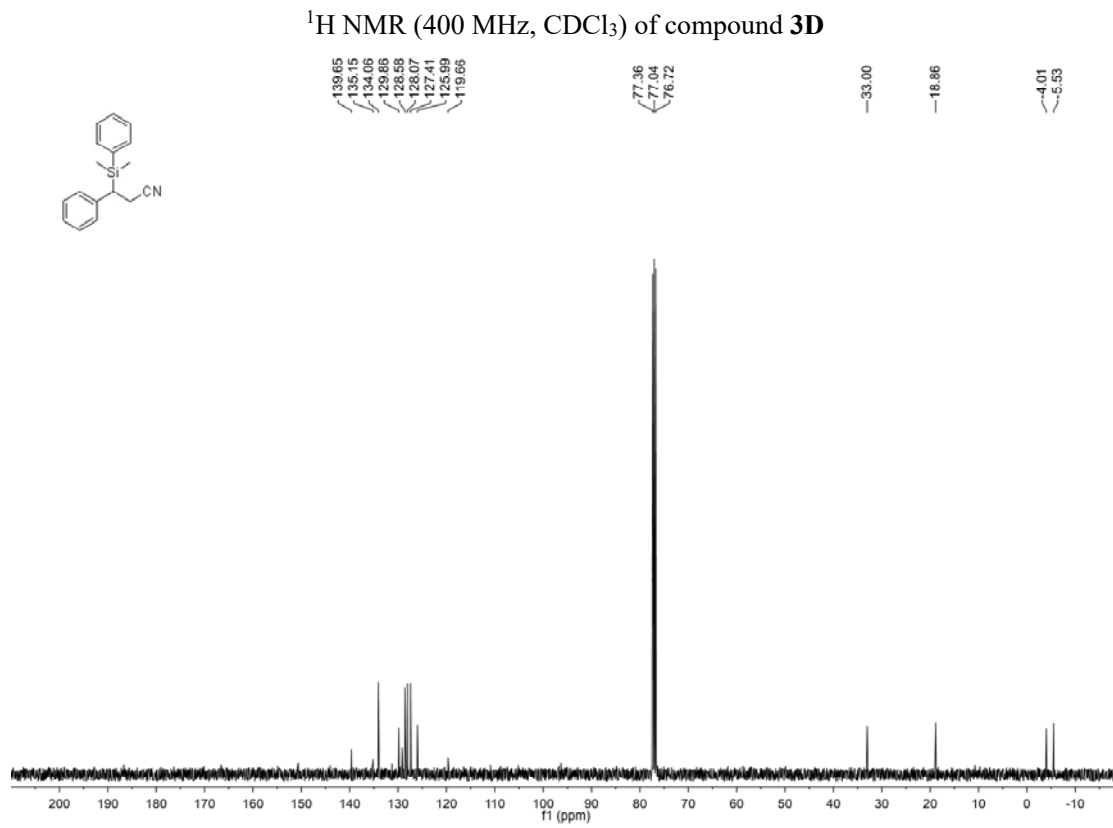
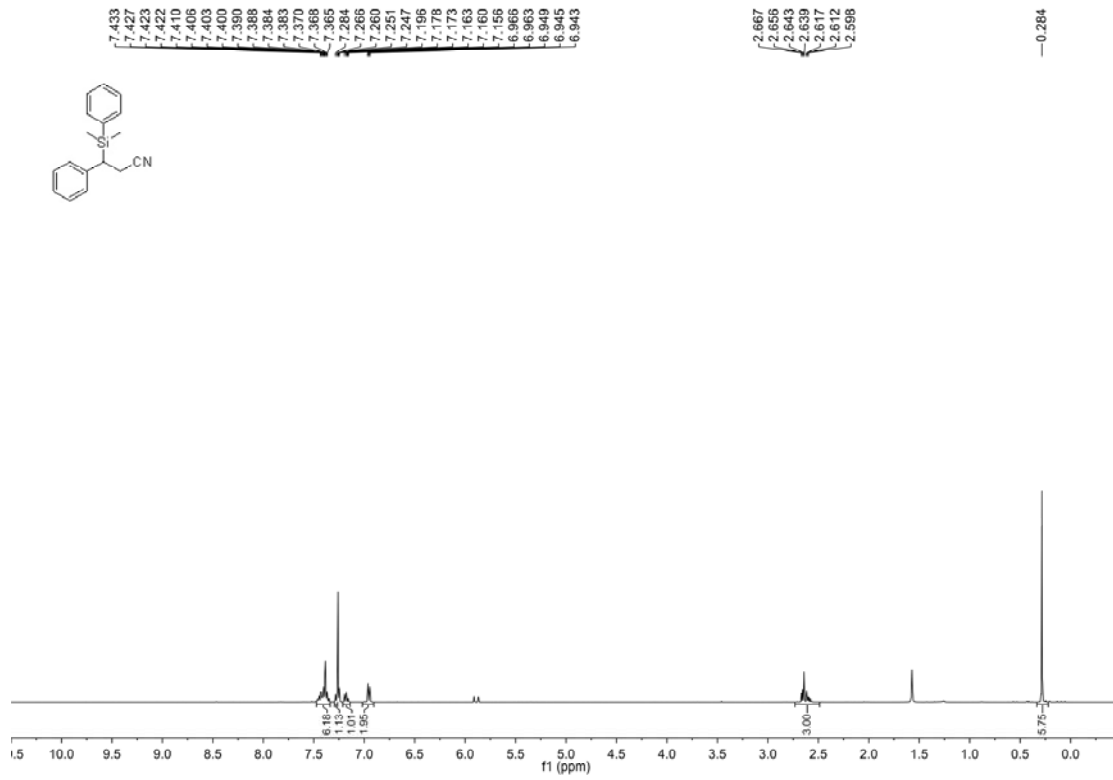
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **3y**

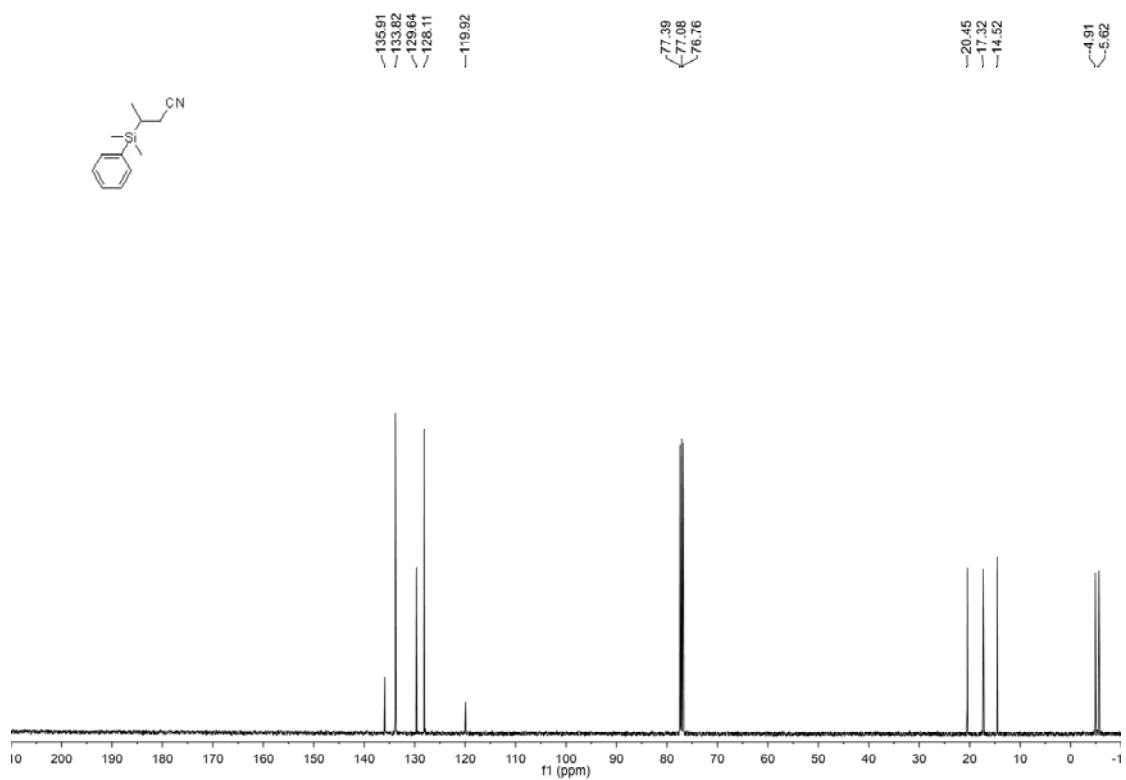
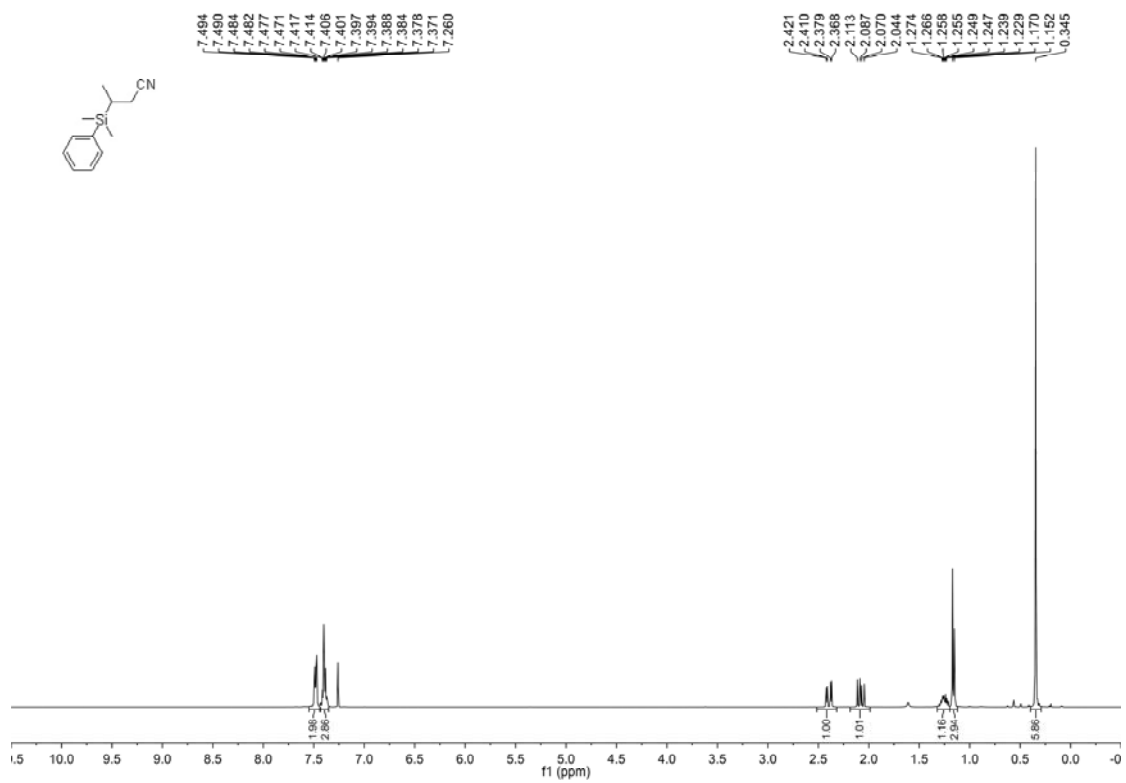


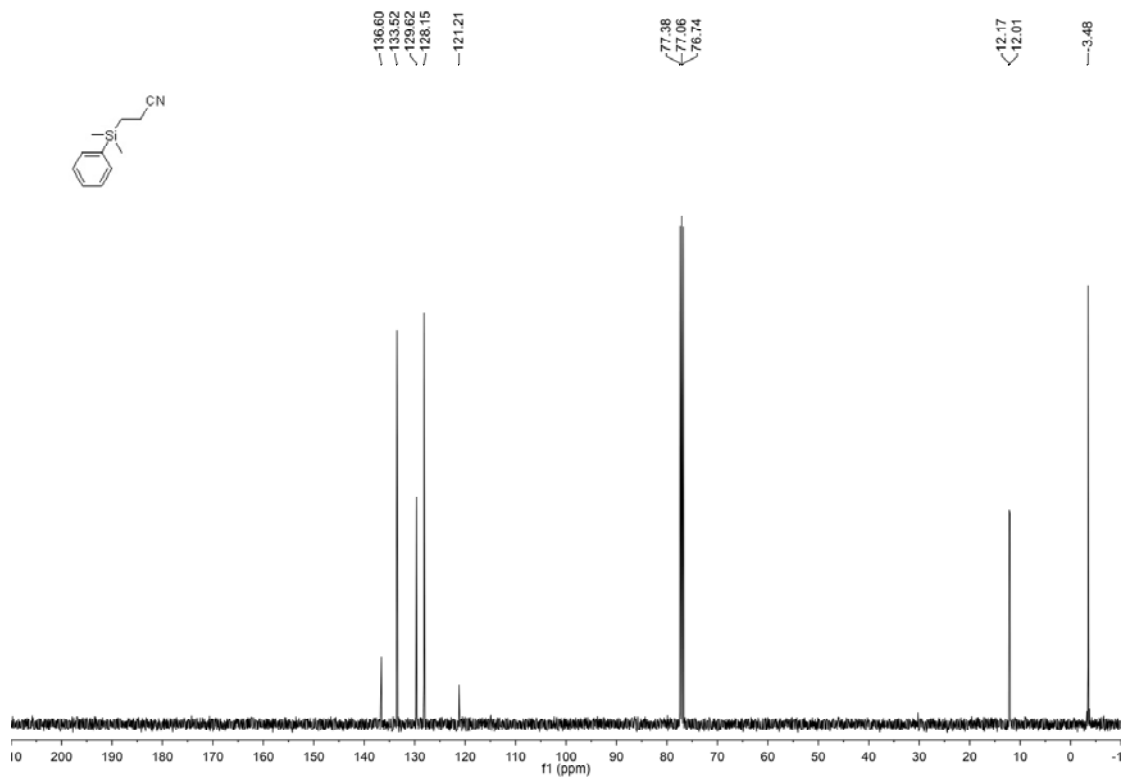
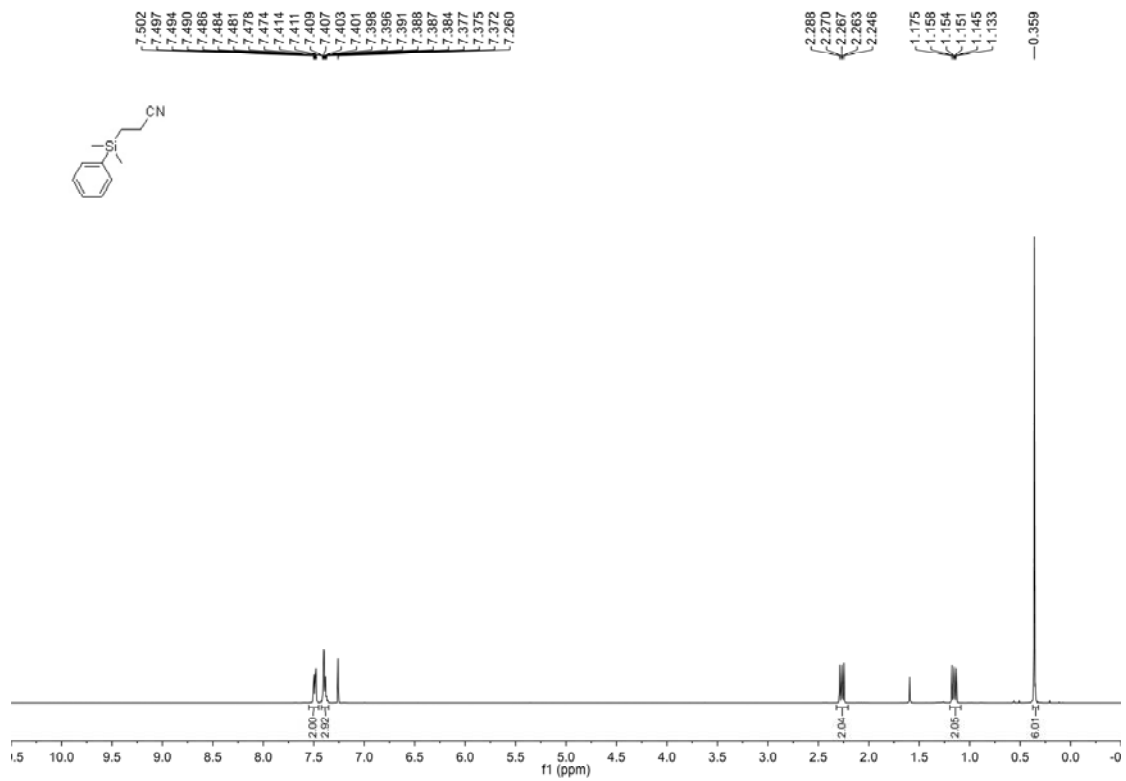


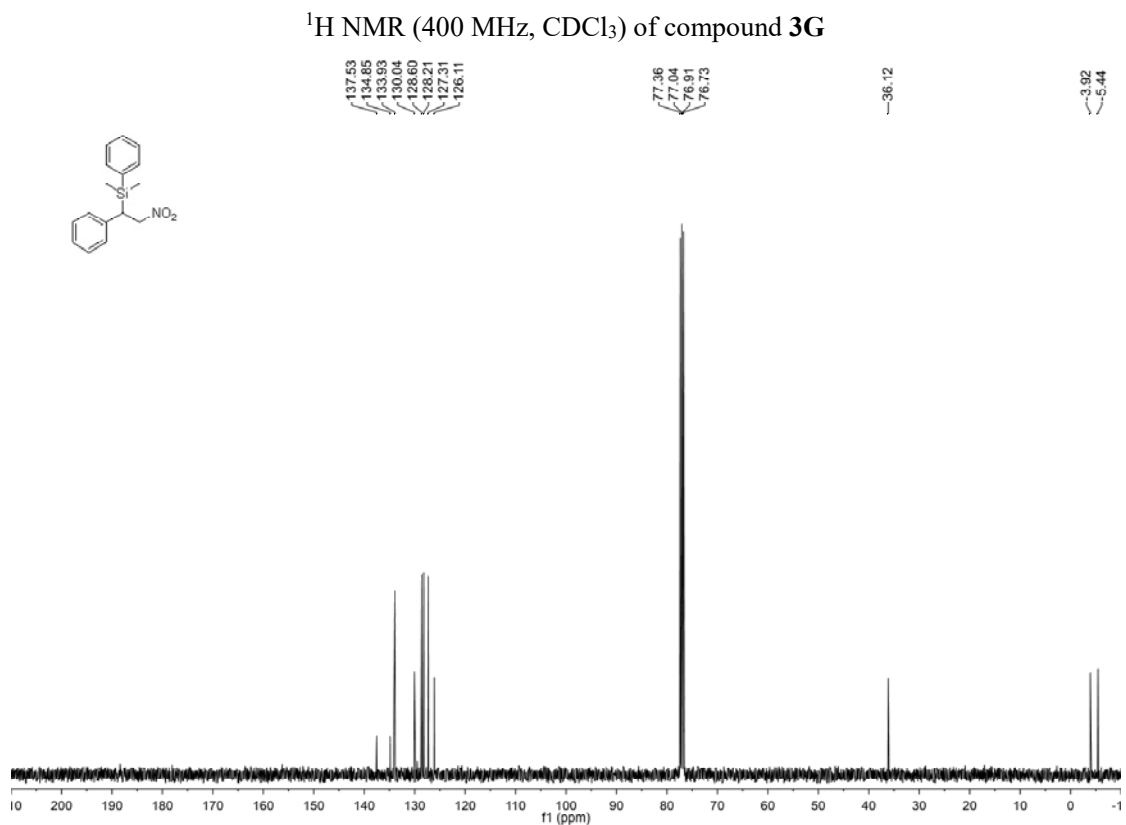
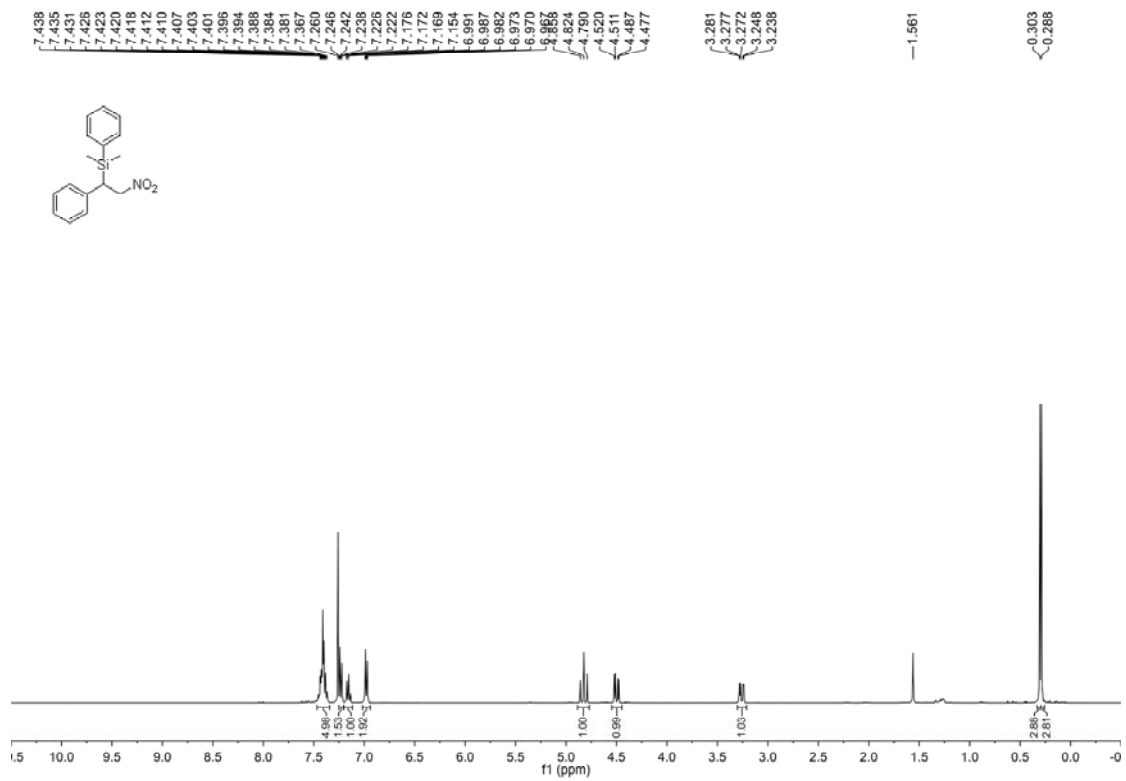




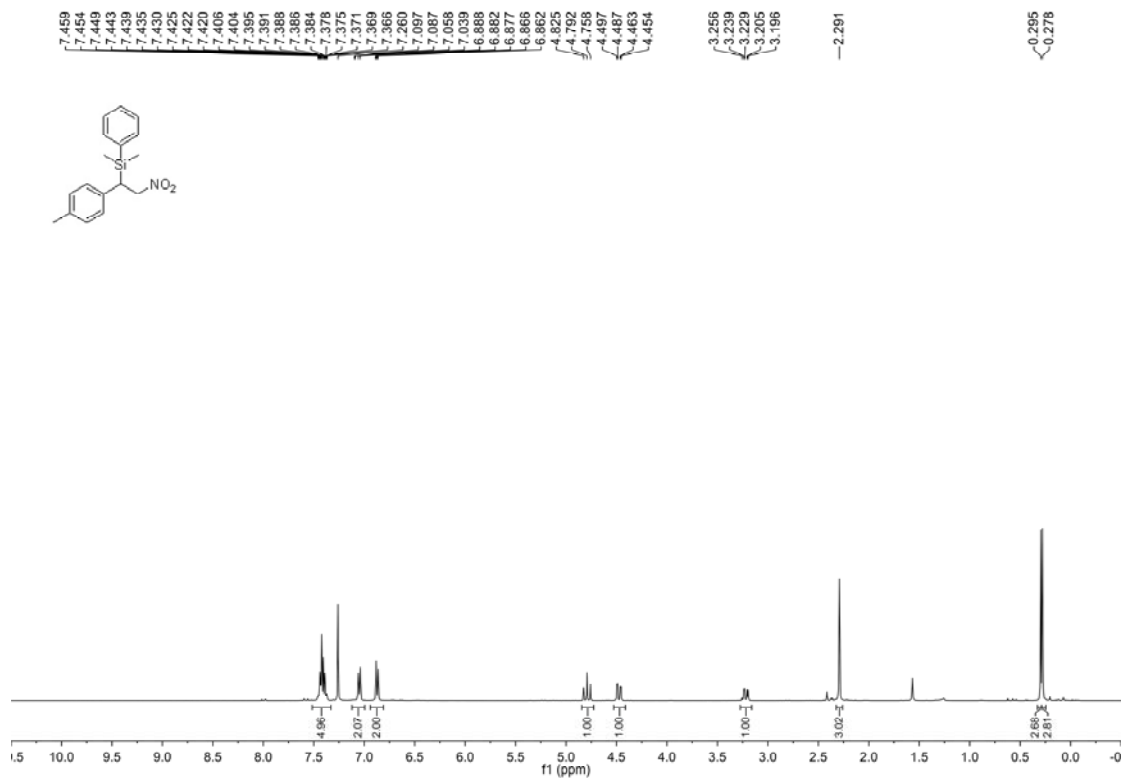




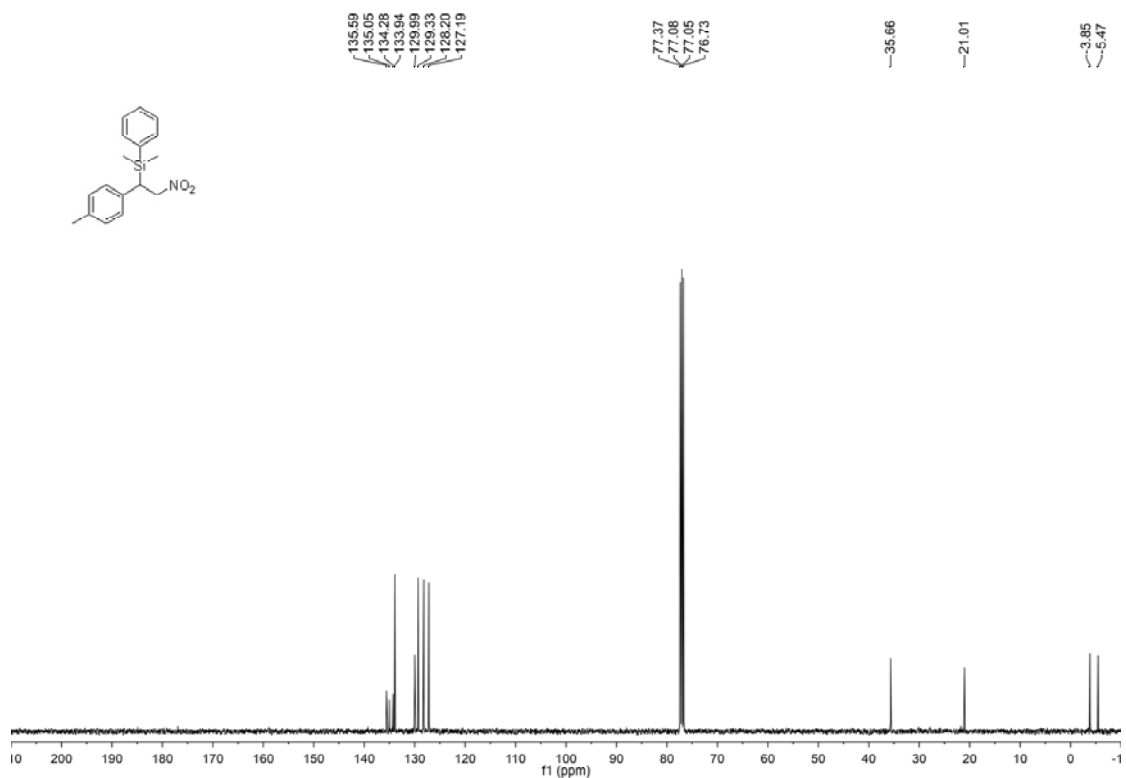




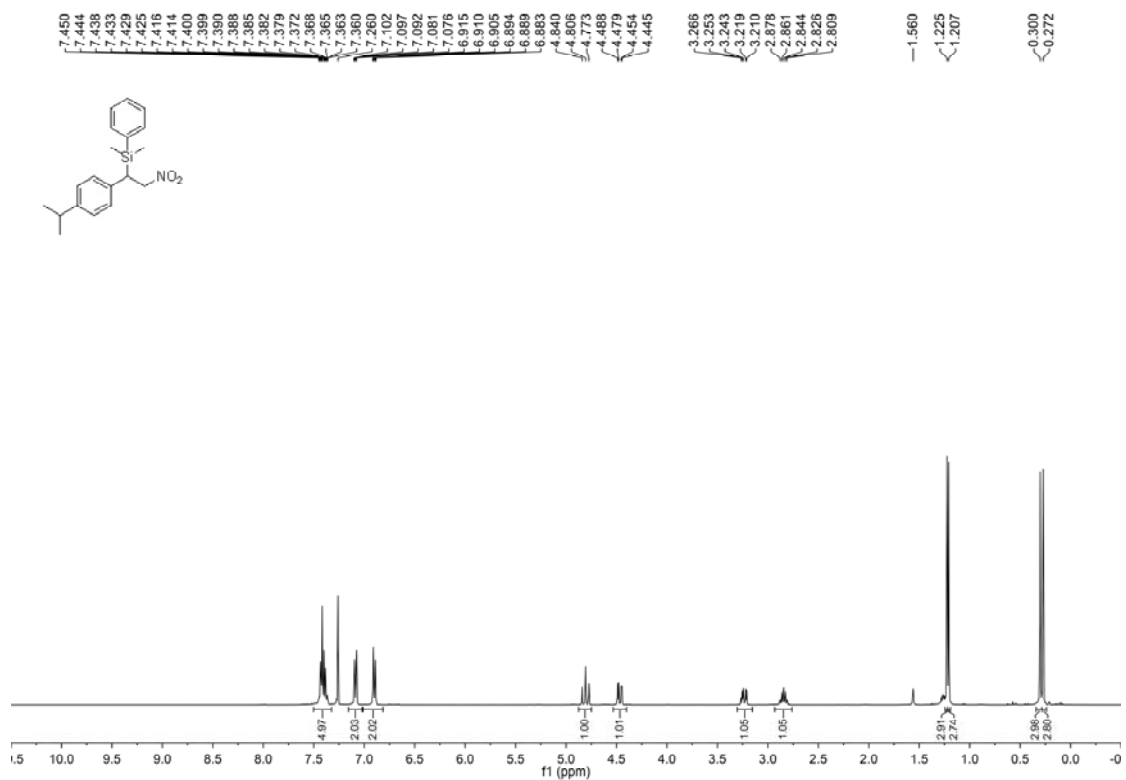
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound 3G



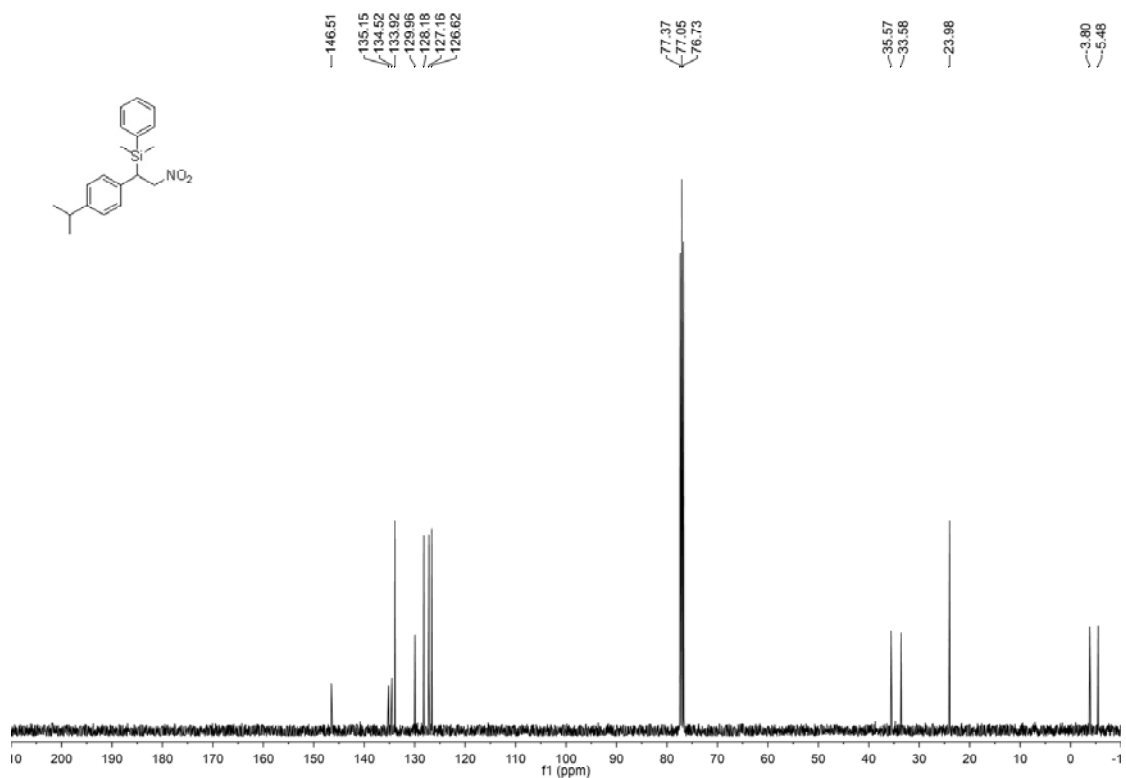
¹H NMR (400 MHz, CDCl₃) of compound 3H



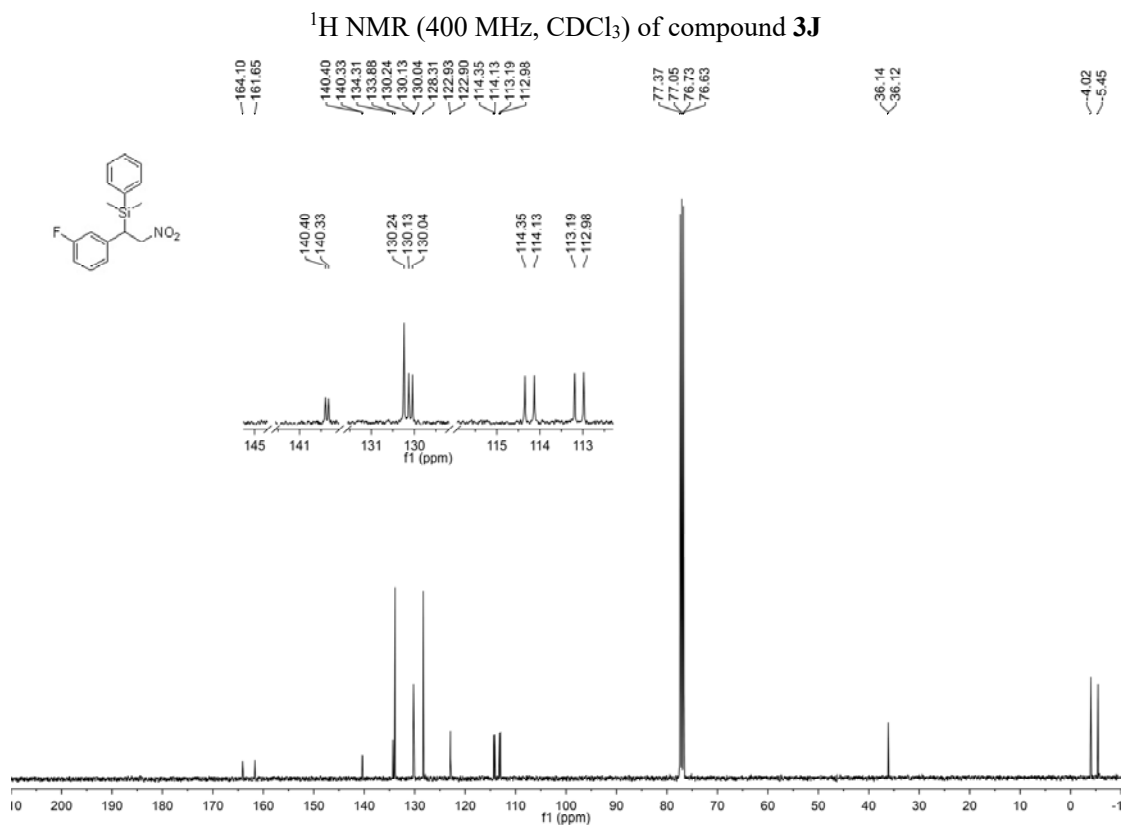
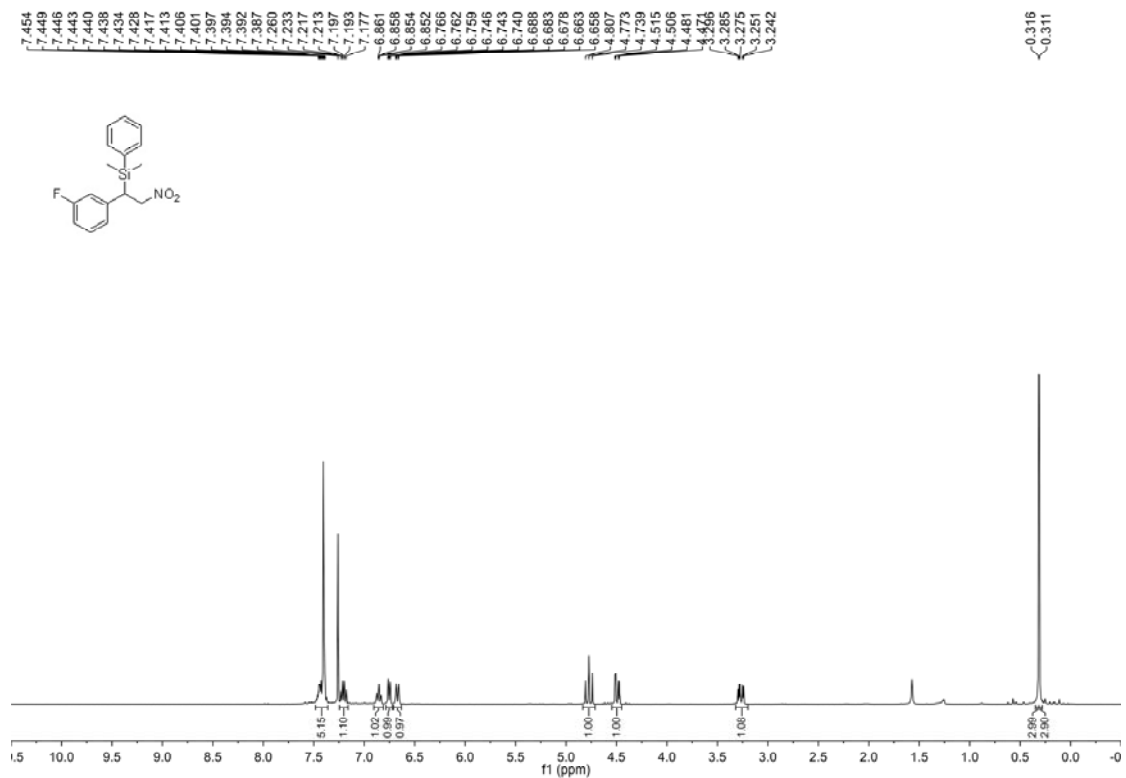
¹³C {¹H} NMR (100 MHz, CDCl₃) of compound 3H

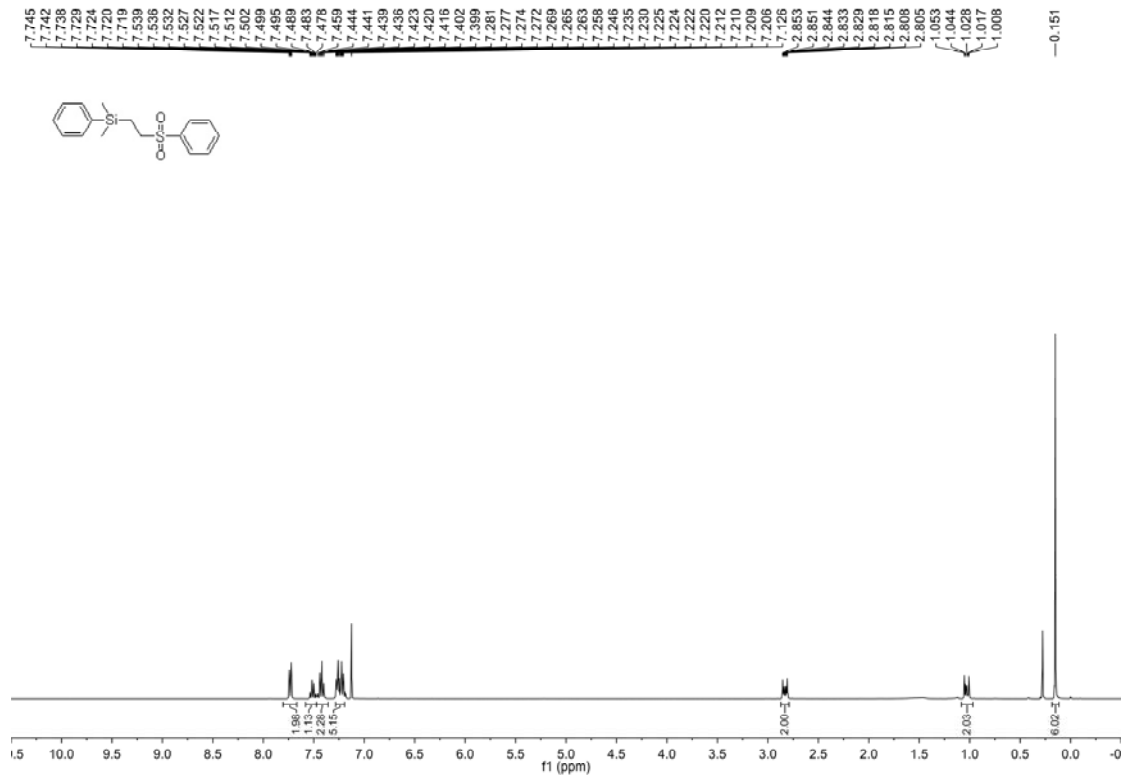


^1H NMR (400 MHz, CDCl_3) of compound **3I**



^{13}C NMR (100 MHz, CDCl_3) of compound **3I**

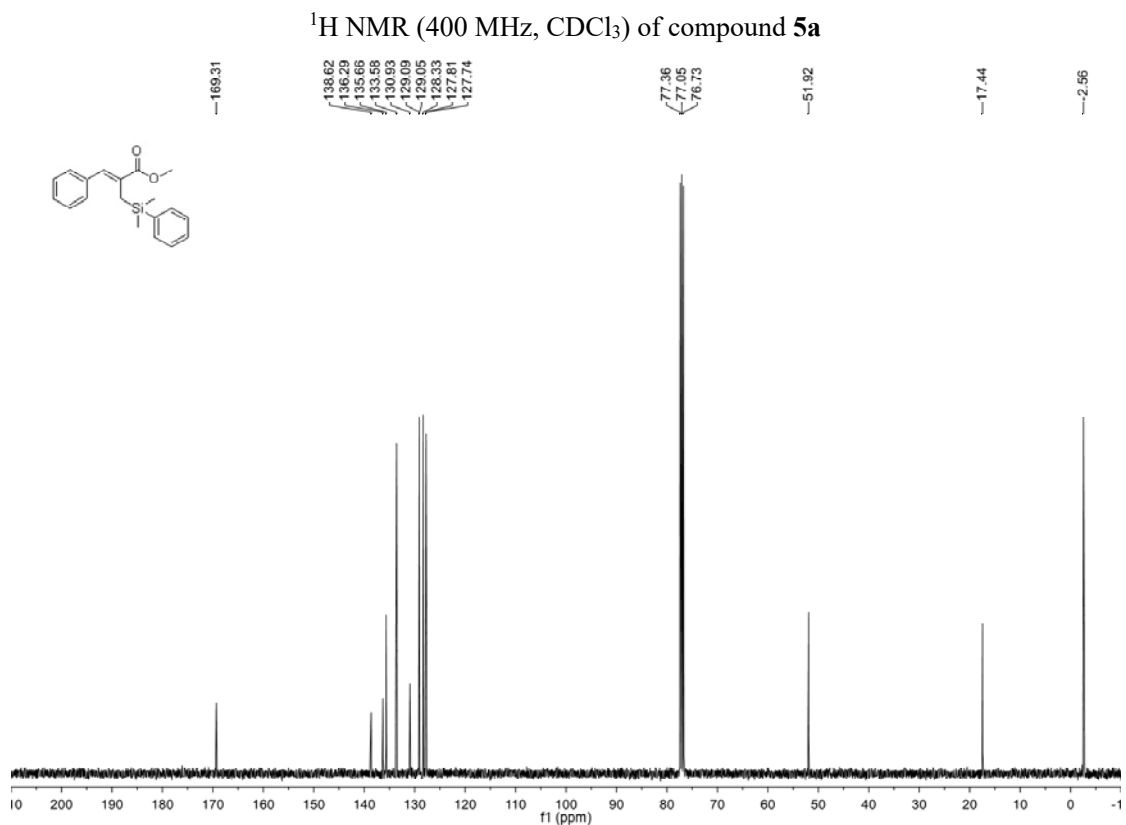
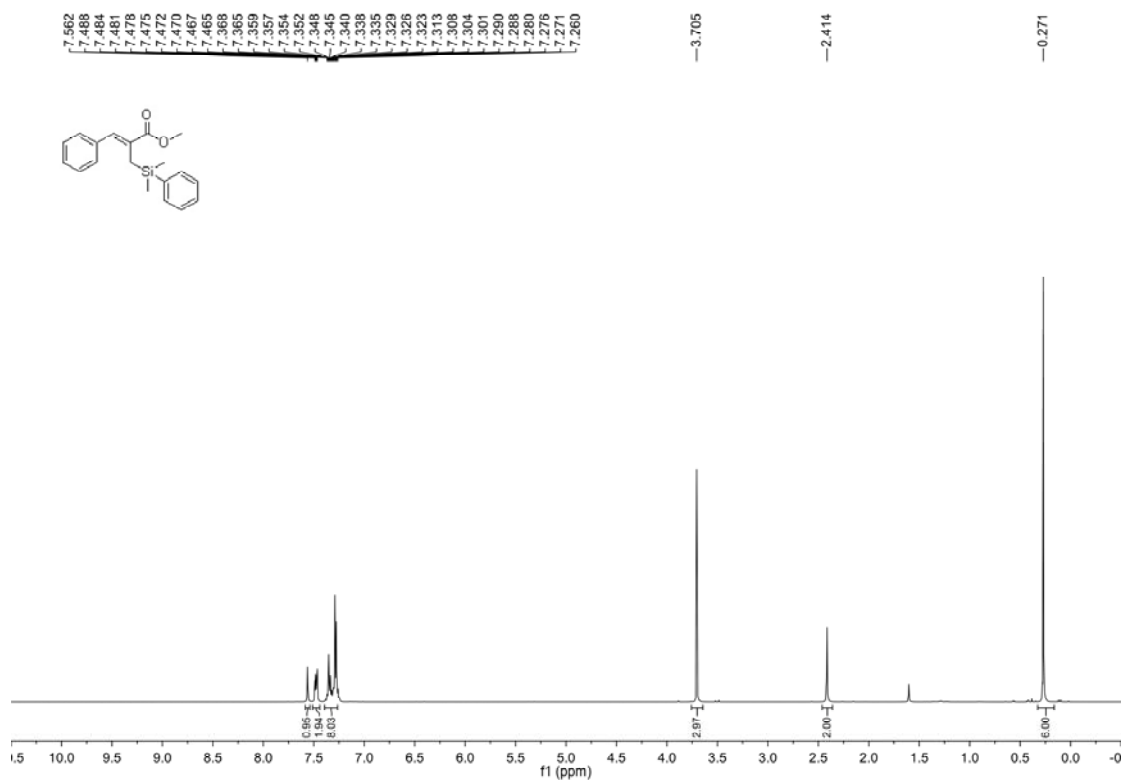


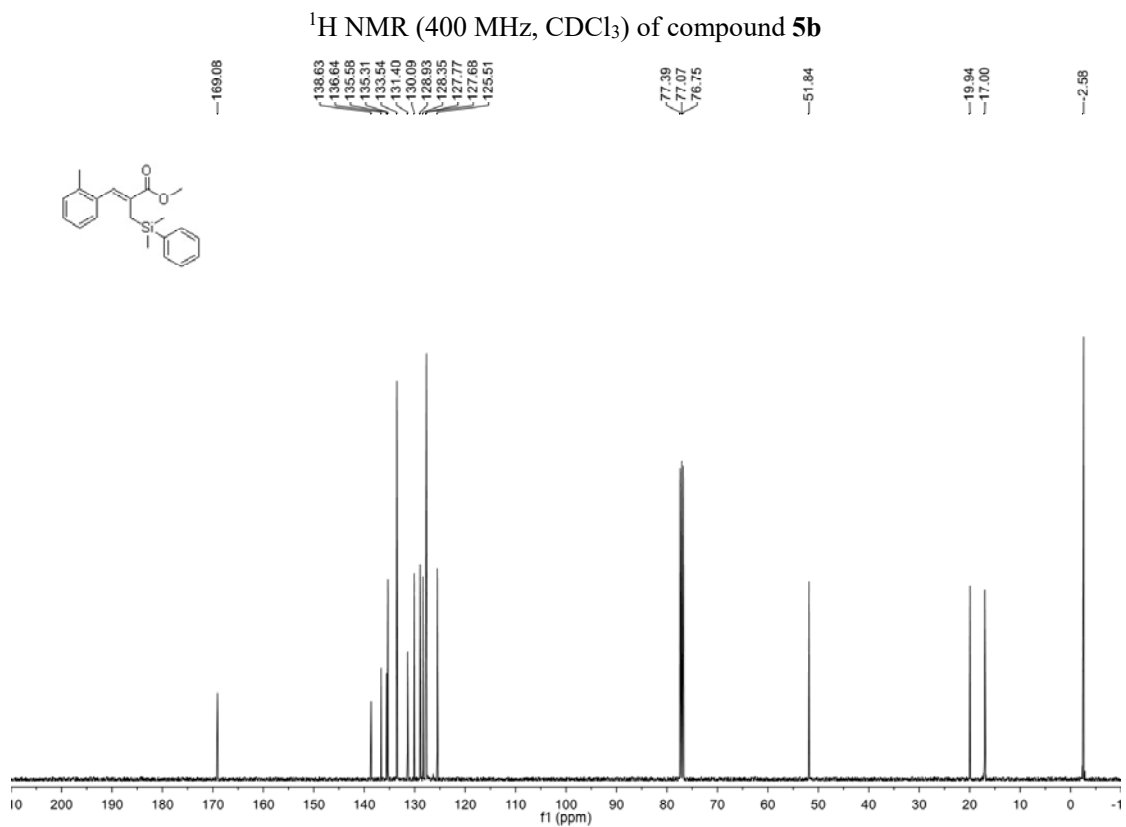
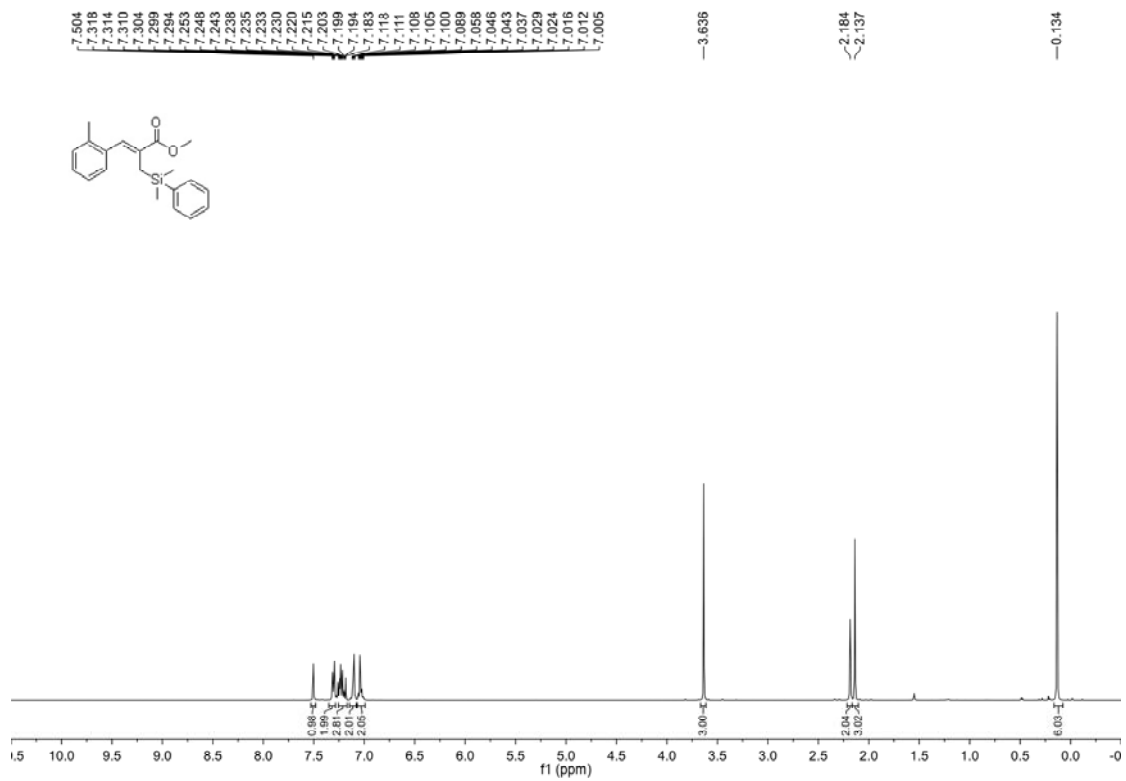


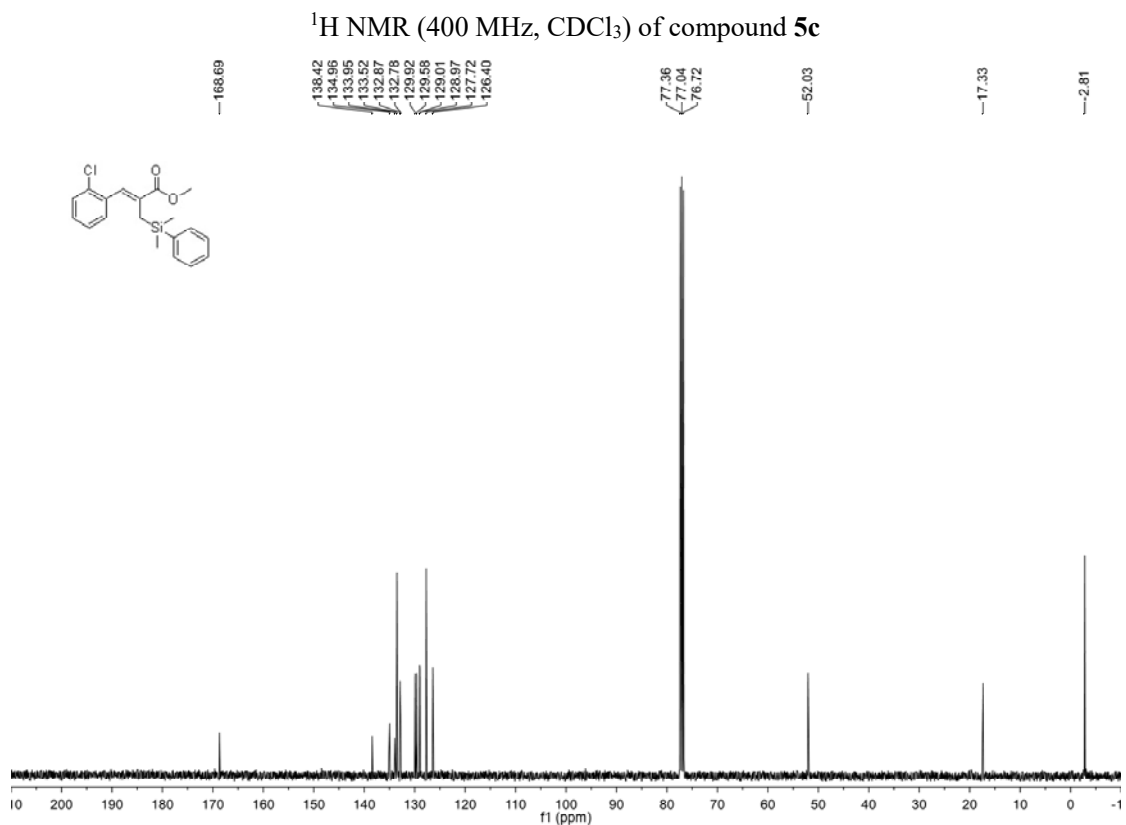
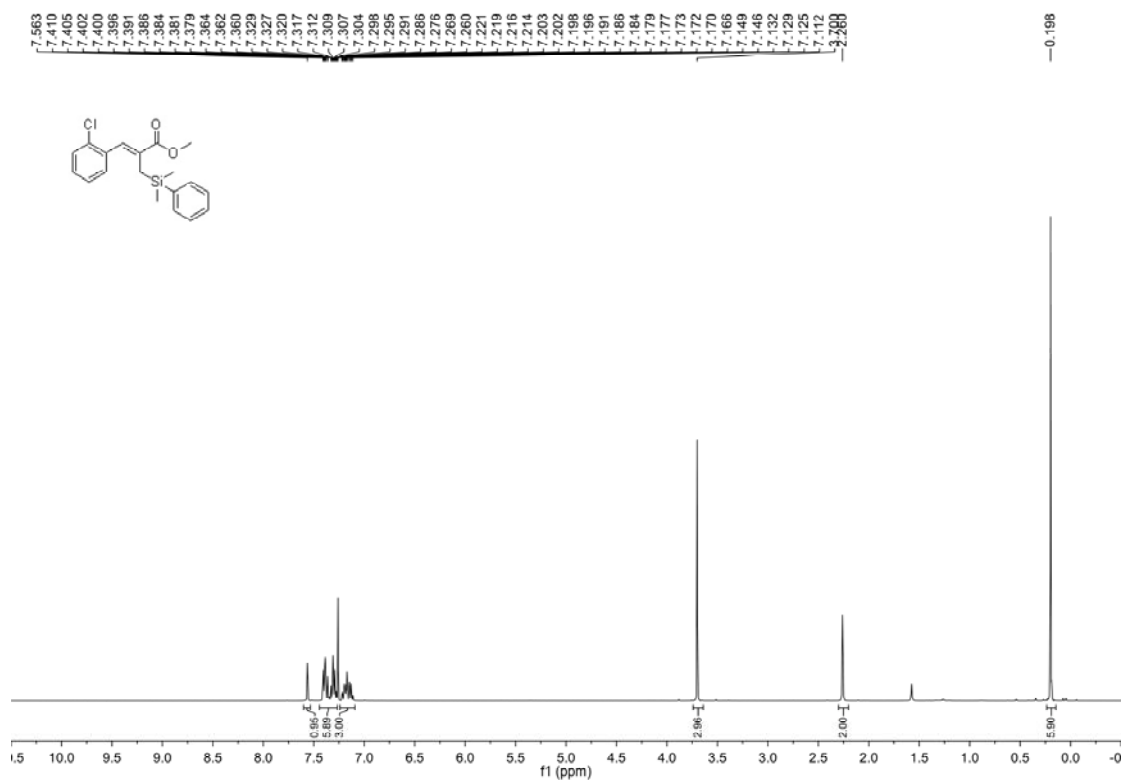
^1H NMR (400 MHz, CDCl_3) of compound **3K**

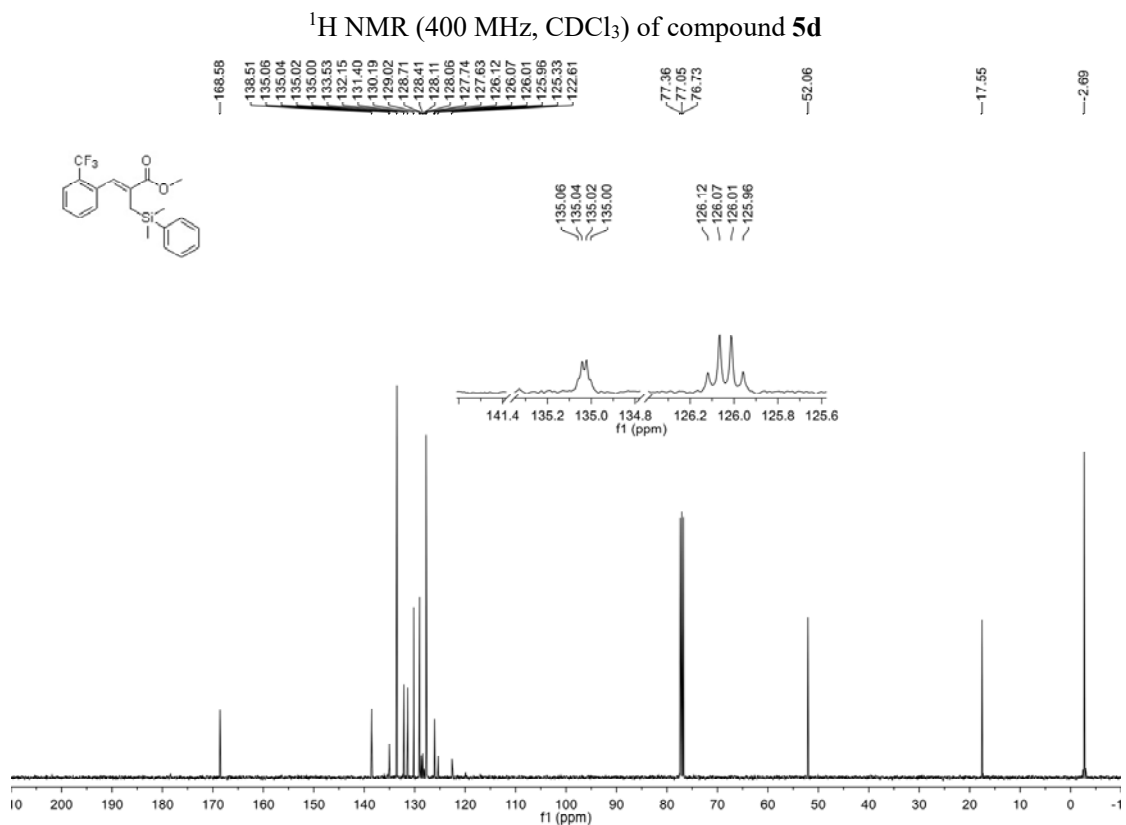
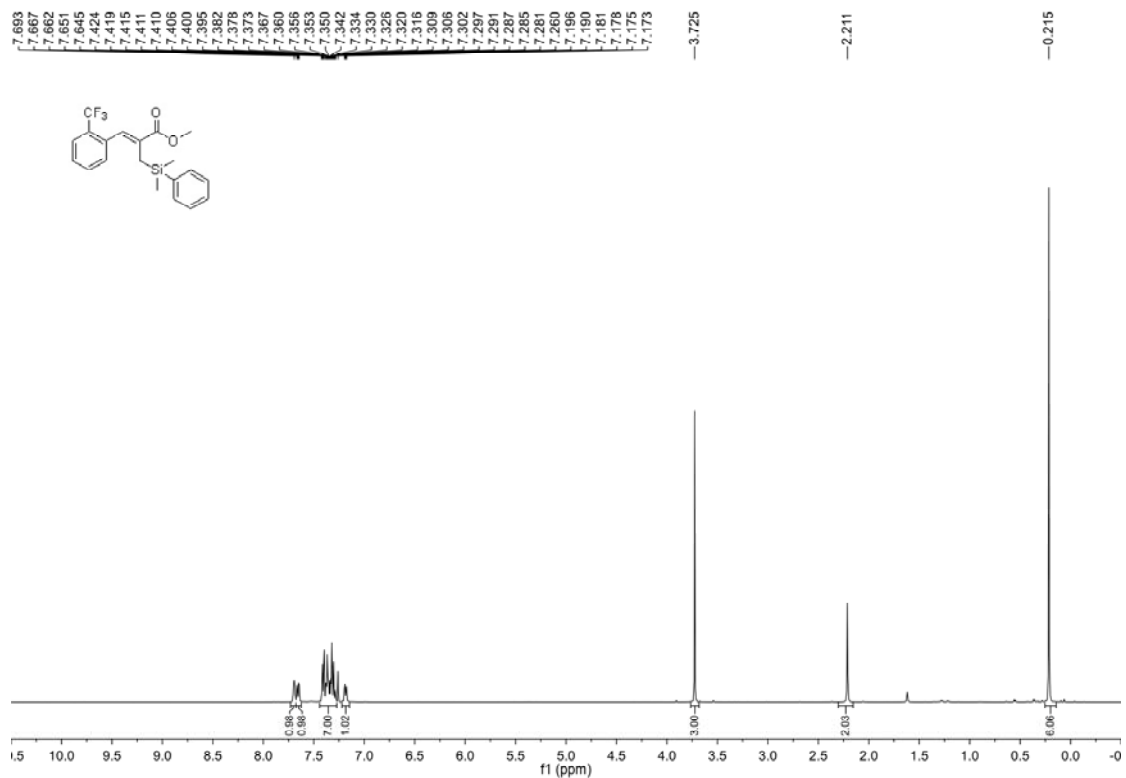


$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **3K**

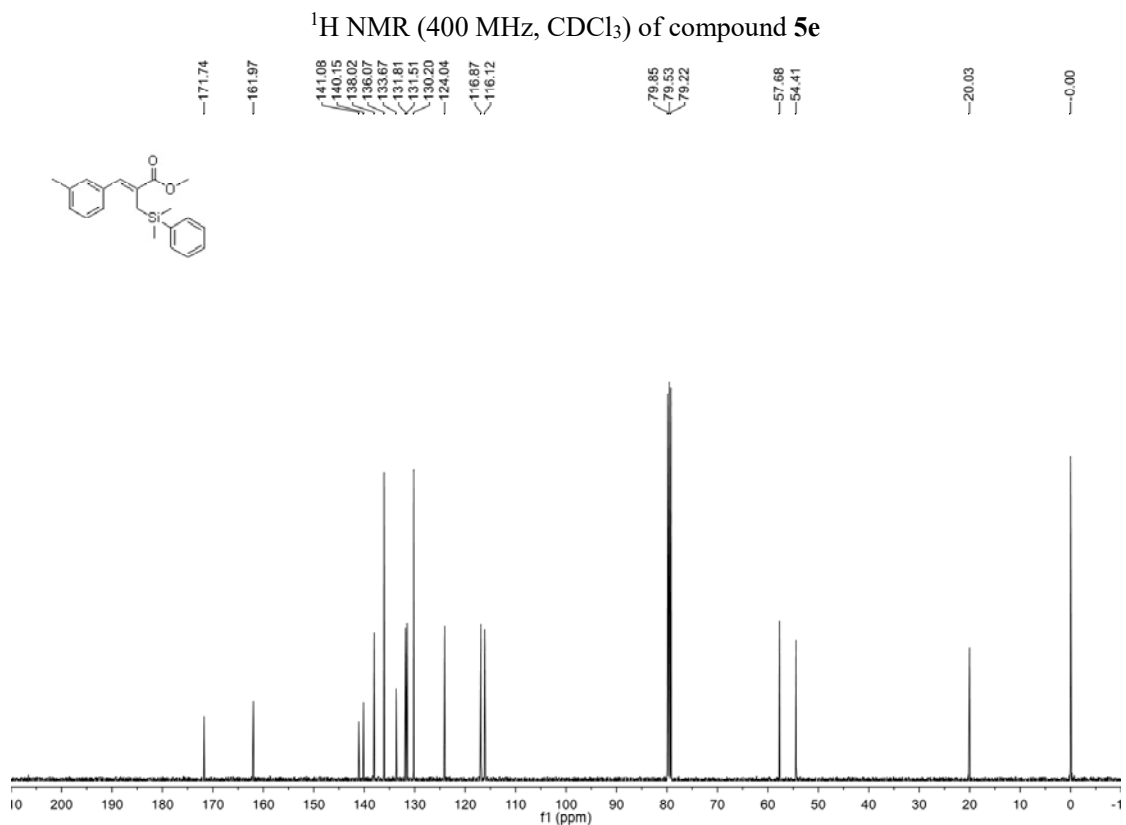
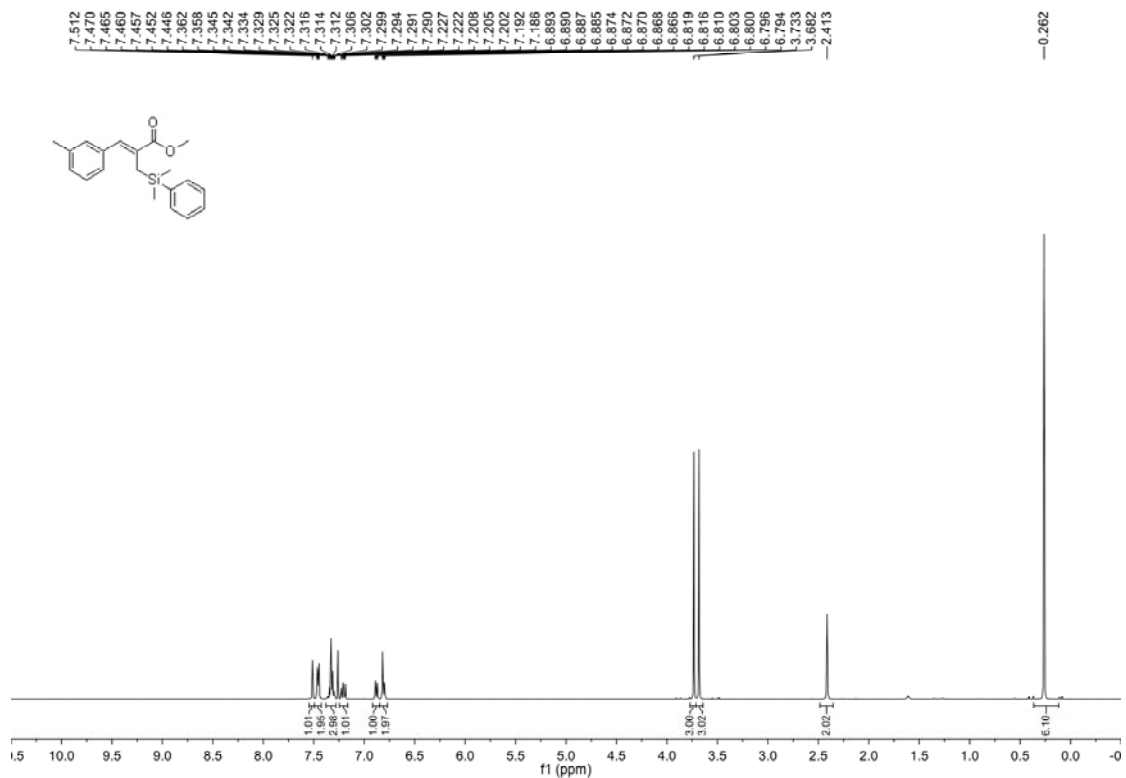


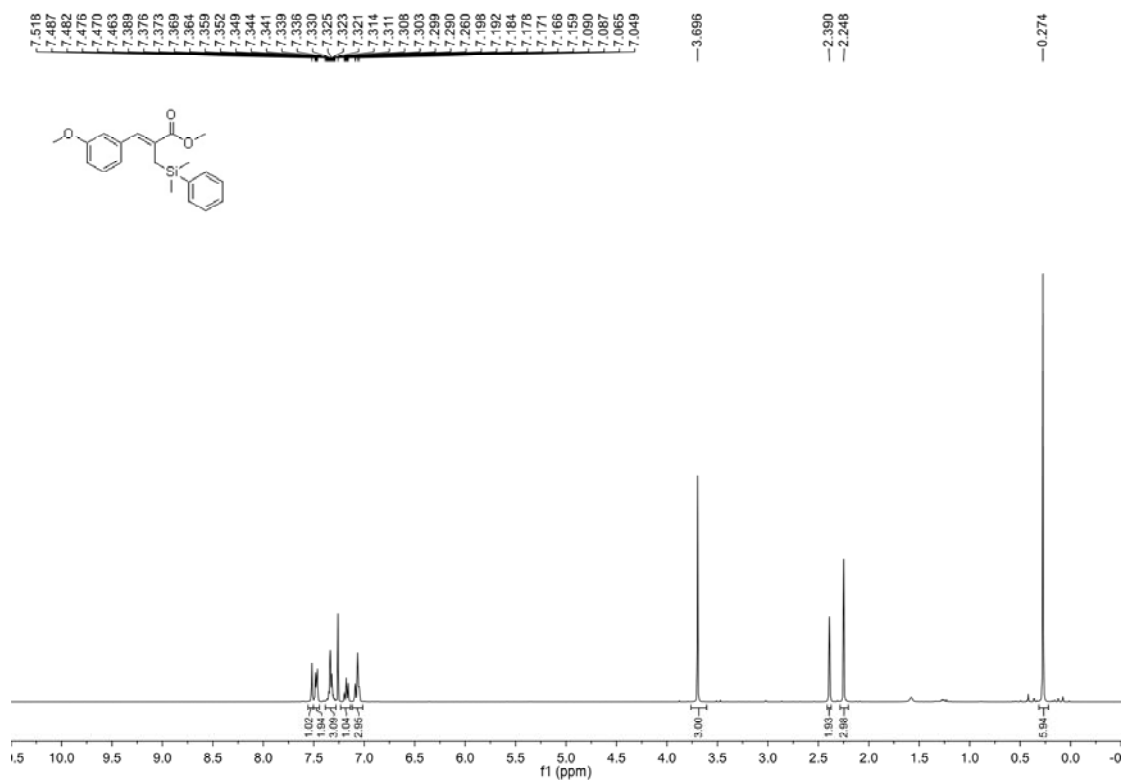




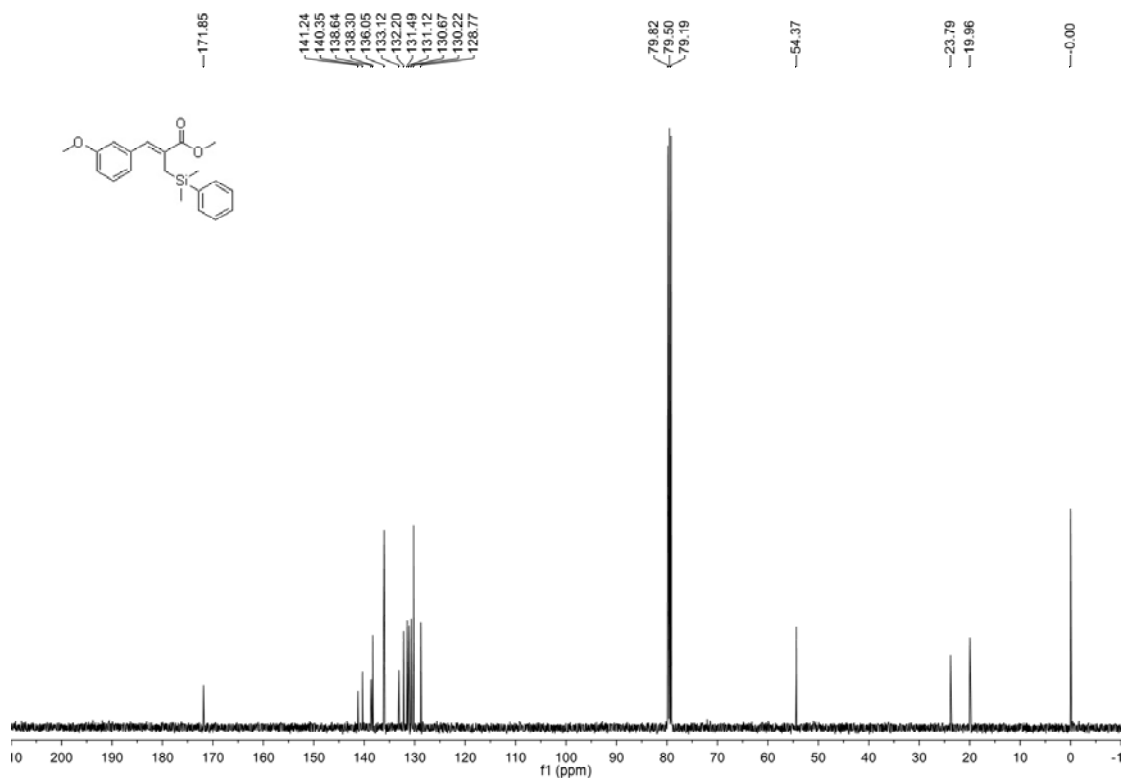


$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **5d**

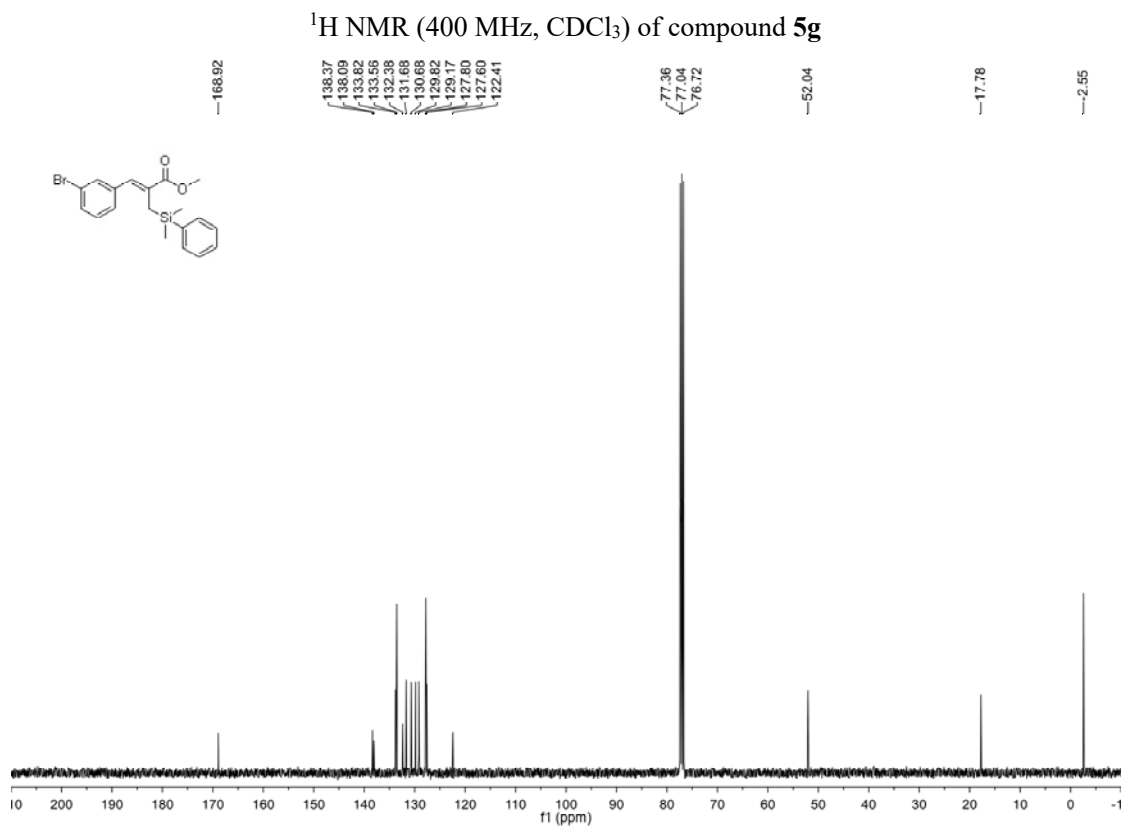
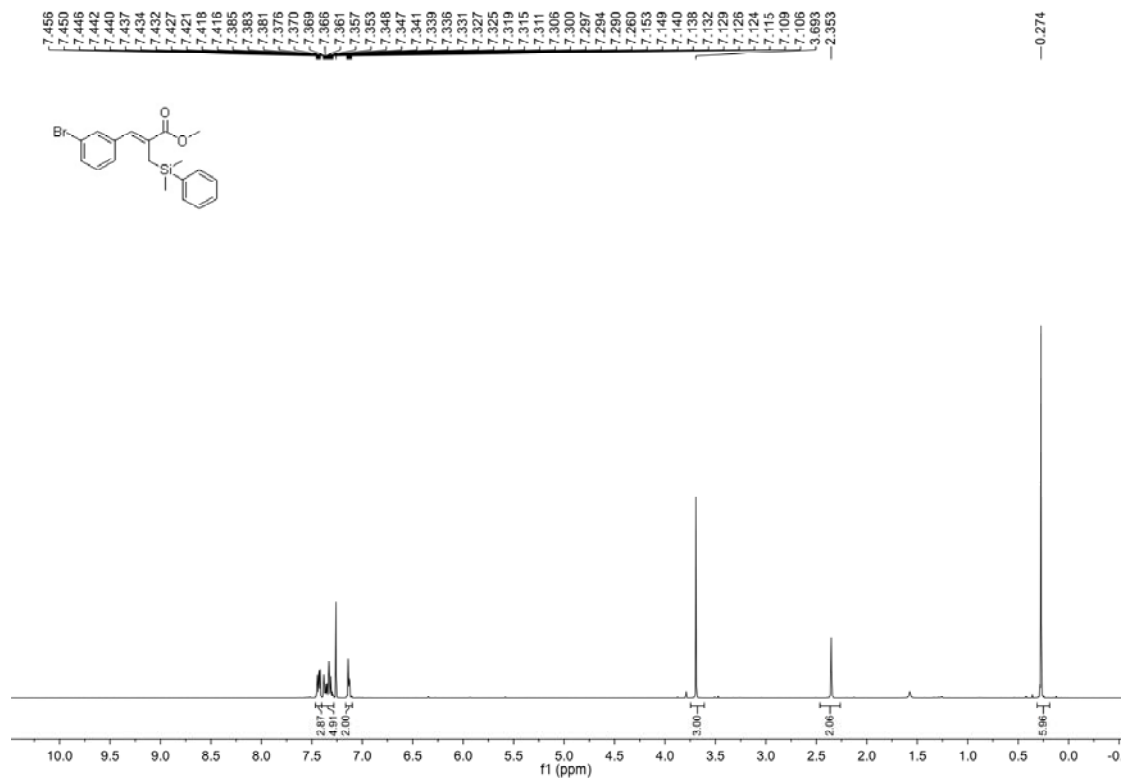


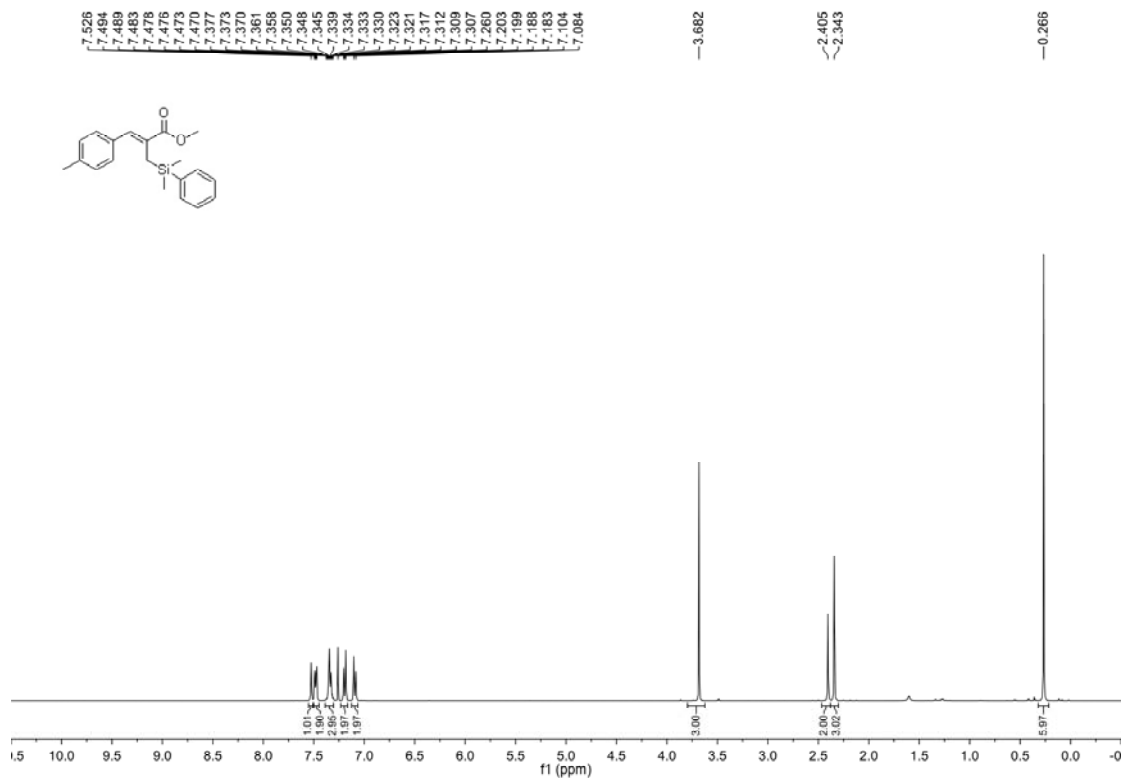


^1H NMR (400 MHz, CDCl_3) of compound **5f**

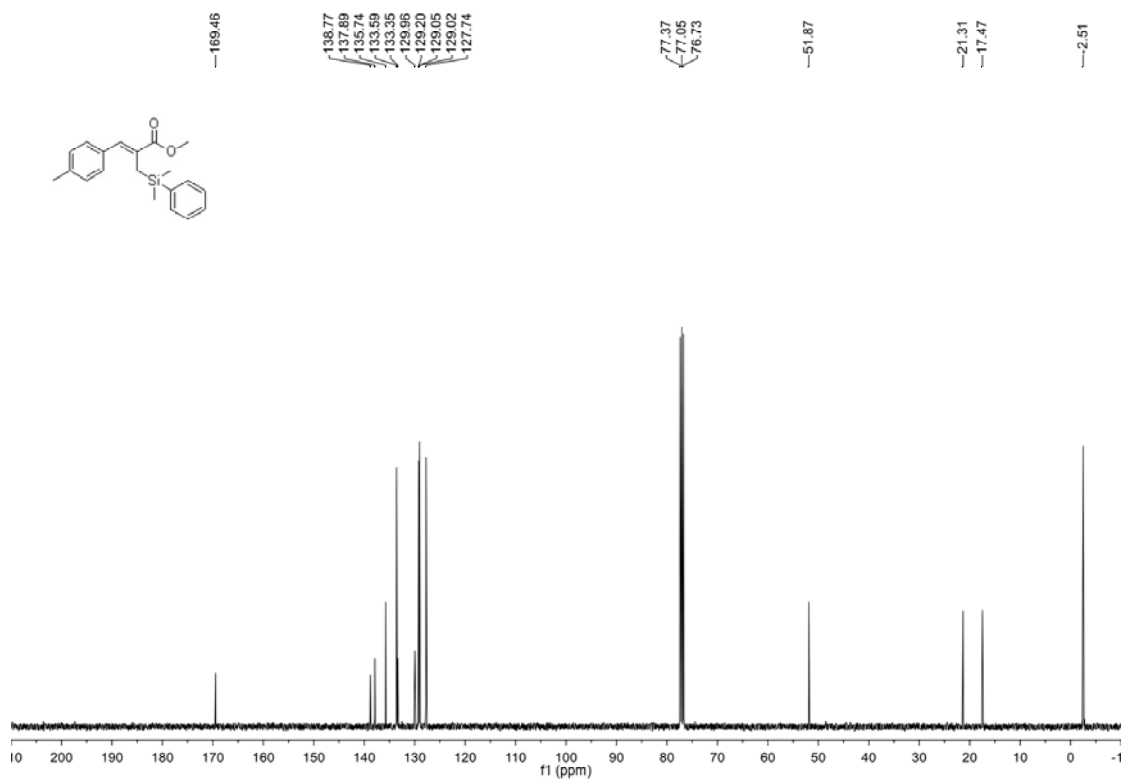


$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **5f**

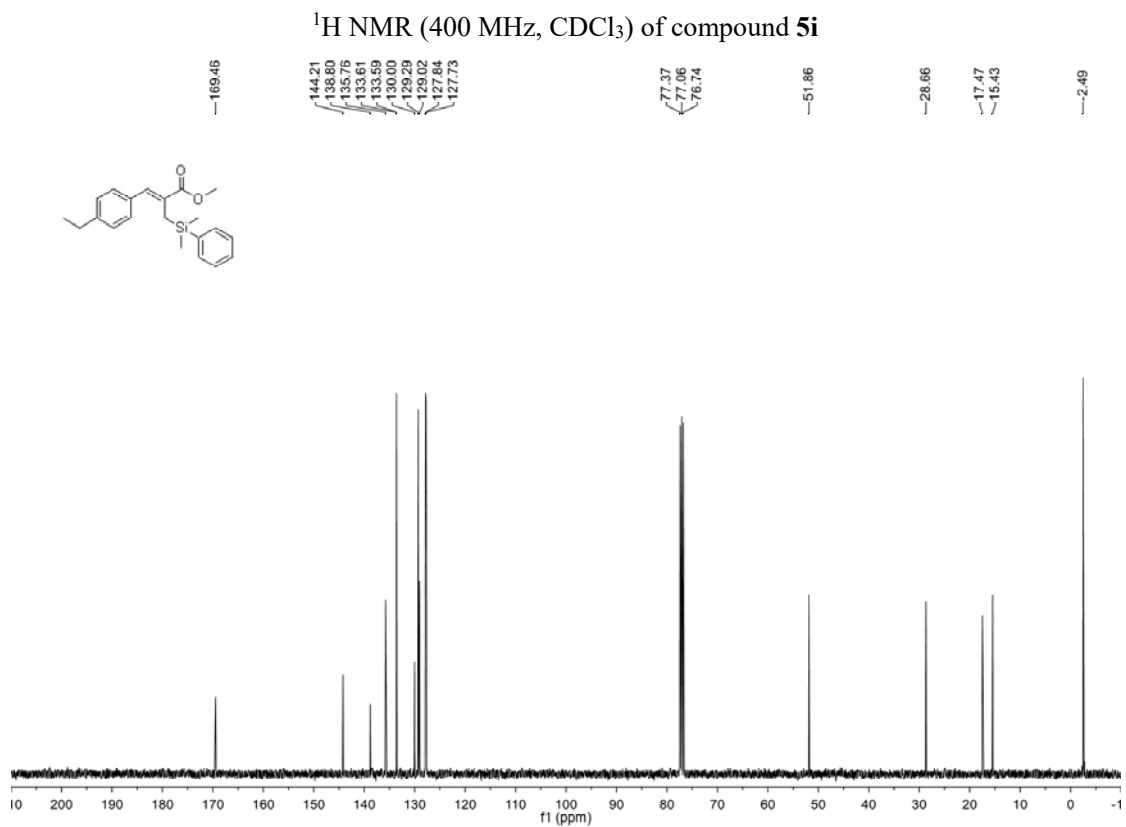
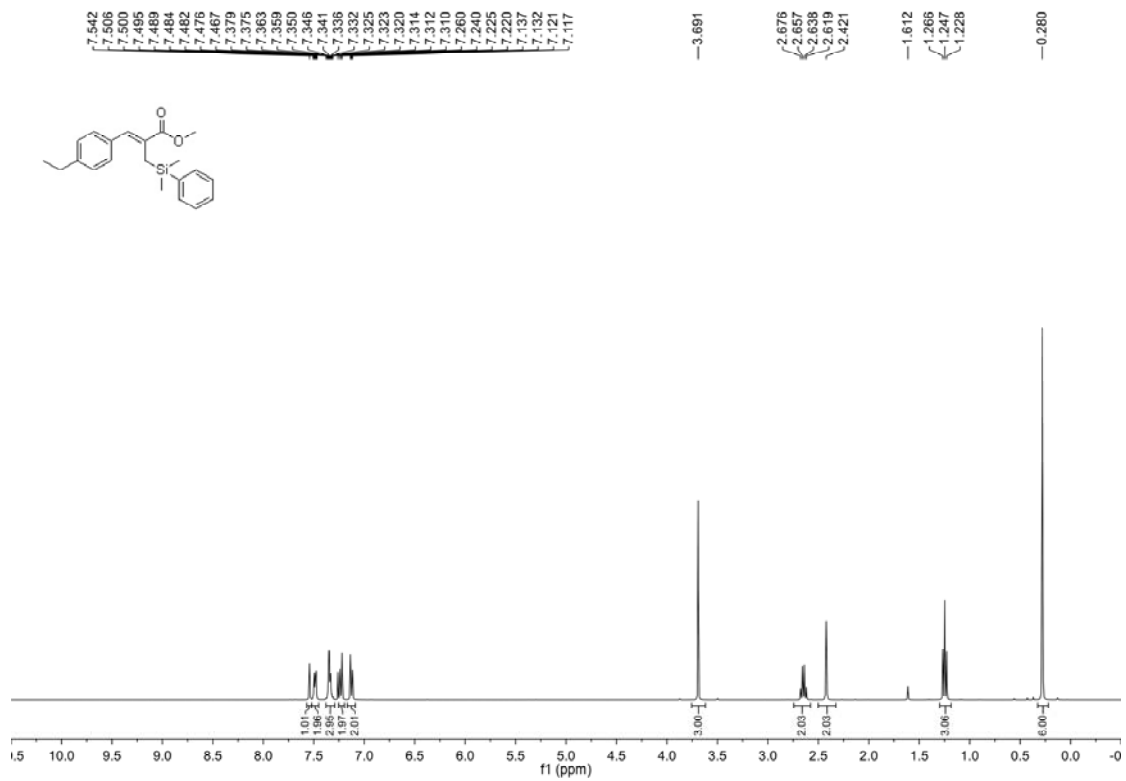


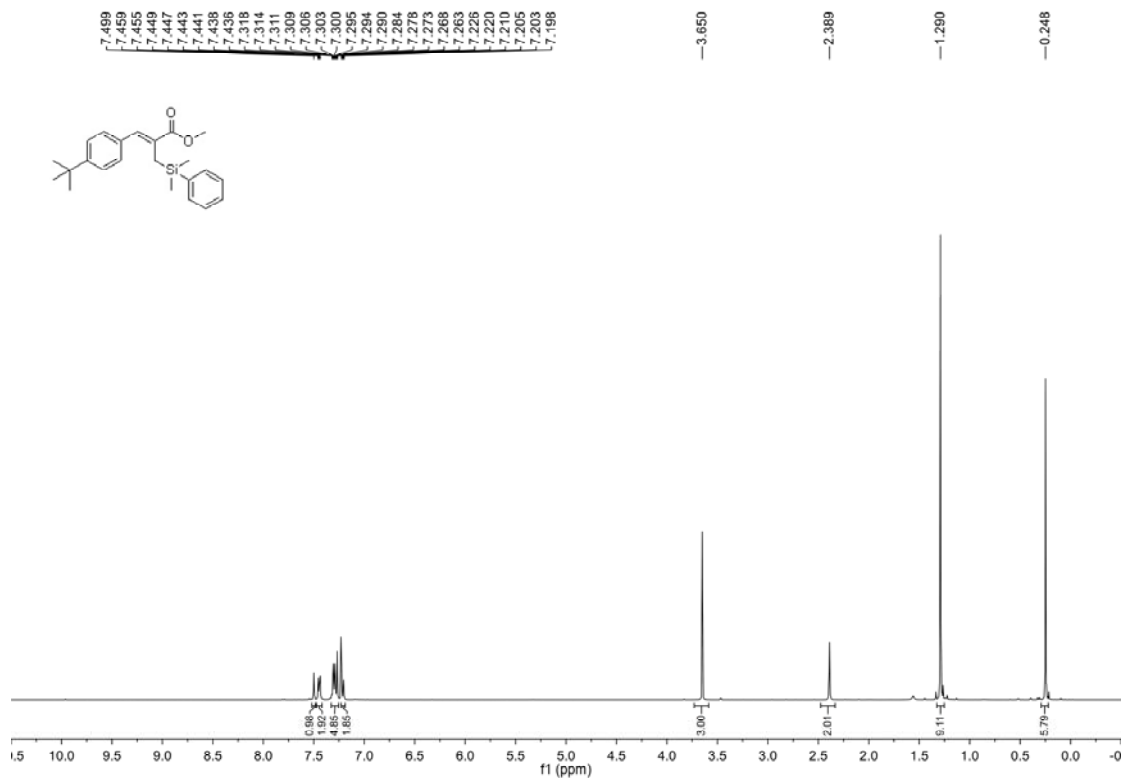


¹H NMR (400 MHz, CDCl₃) of compound **5h**

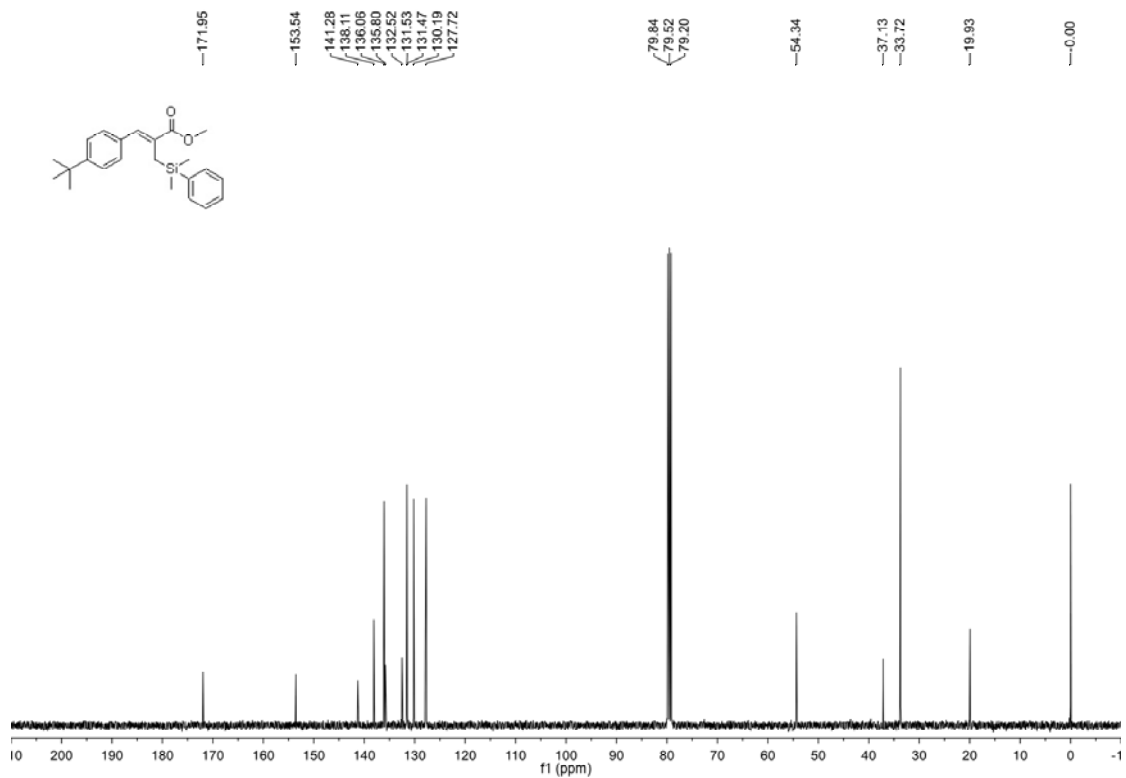


¹³C{¹H} NMR (100 MHz, CDCl₃) of compound **5h**

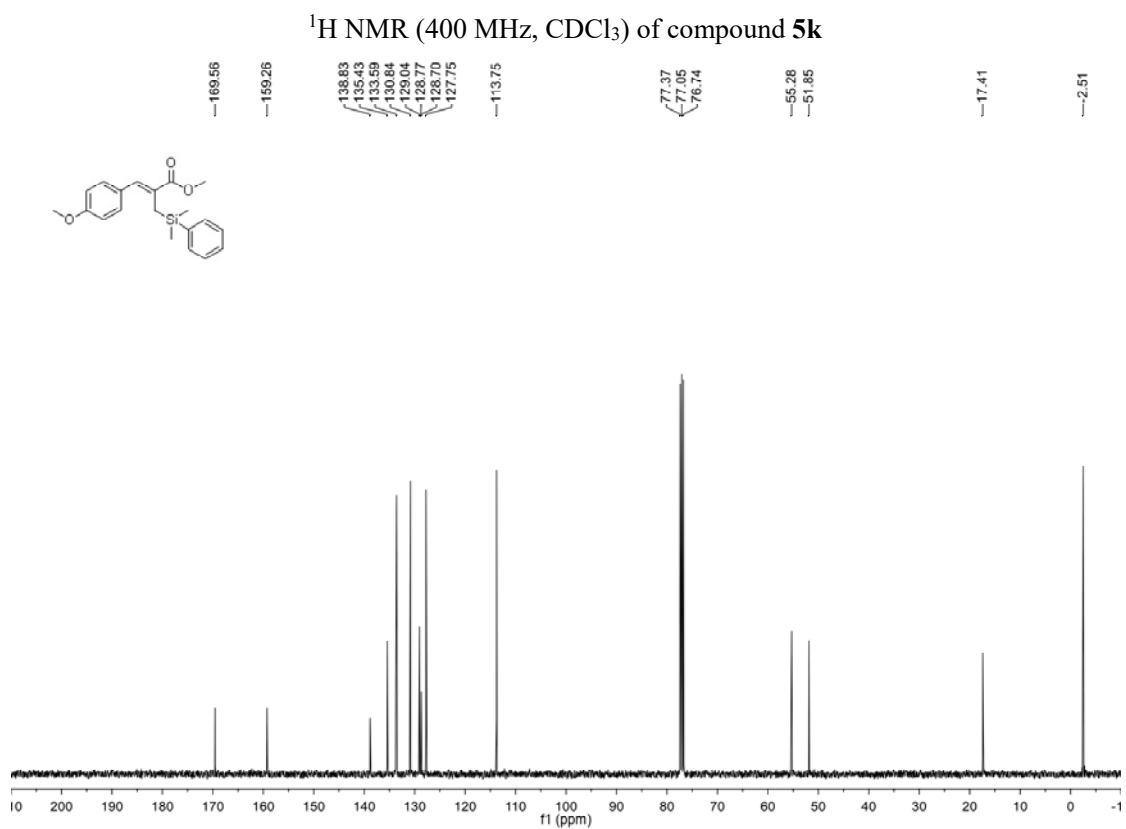
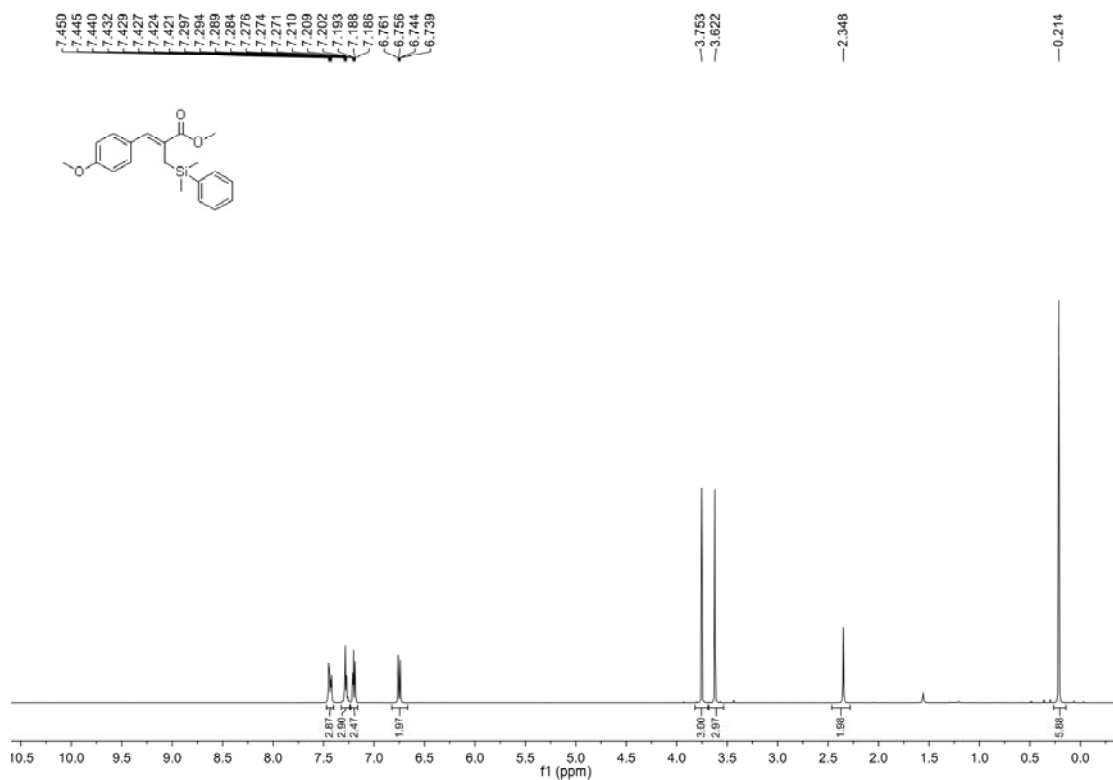


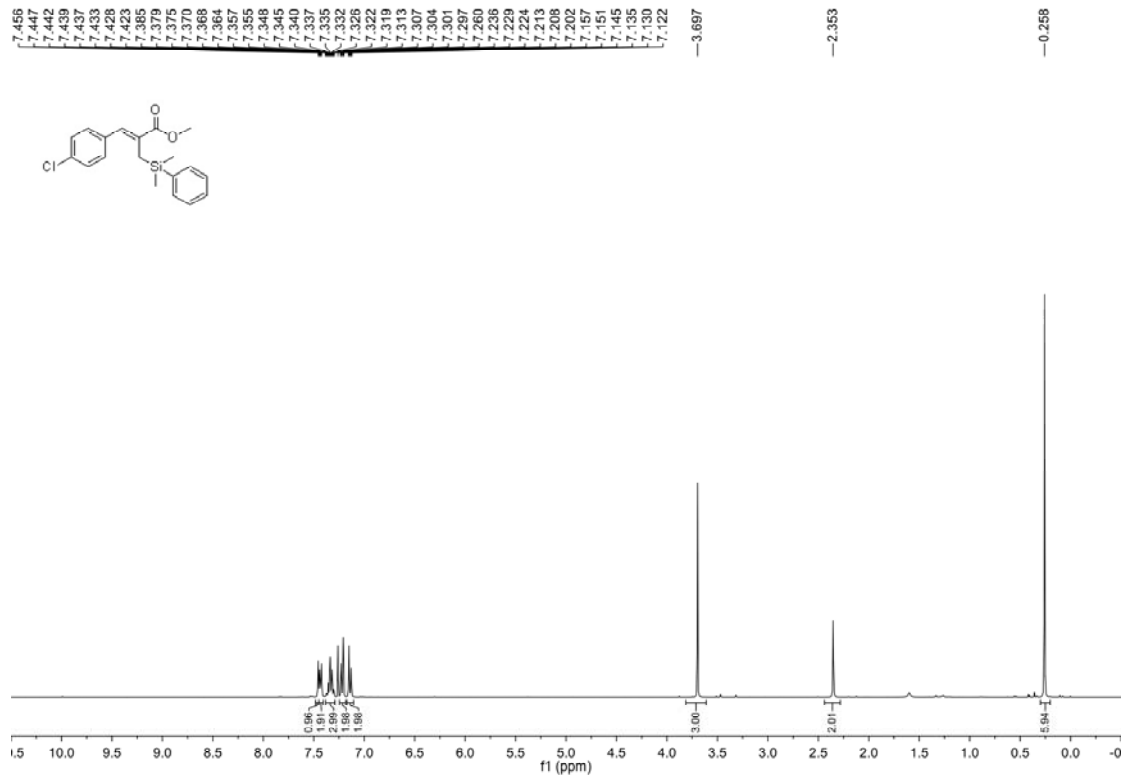


^1H NMR (400 MHz, CDCl_3) of compound **5j**

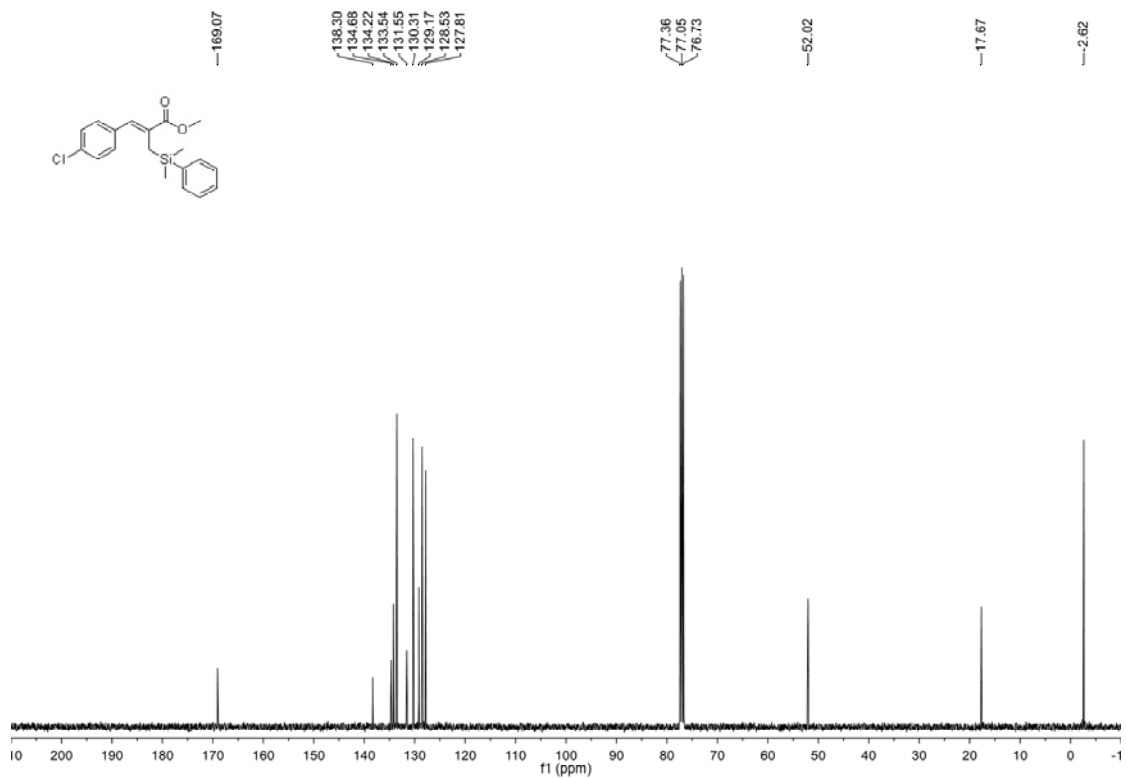


$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **5j**

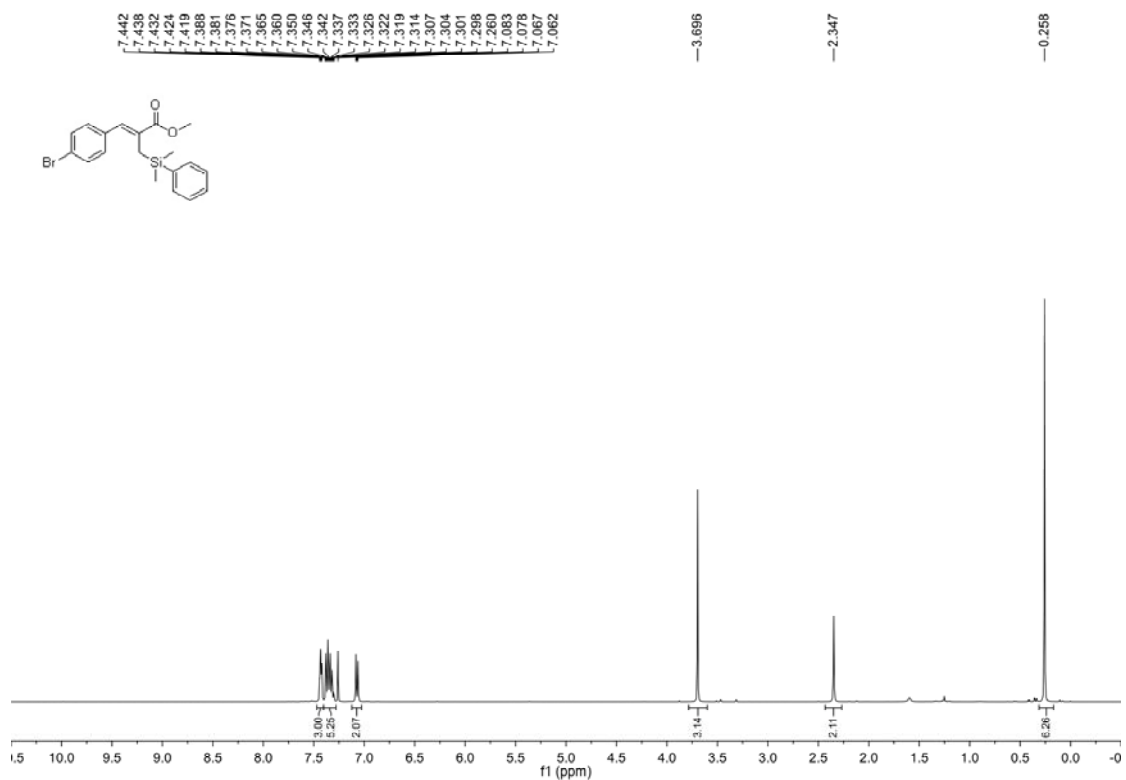




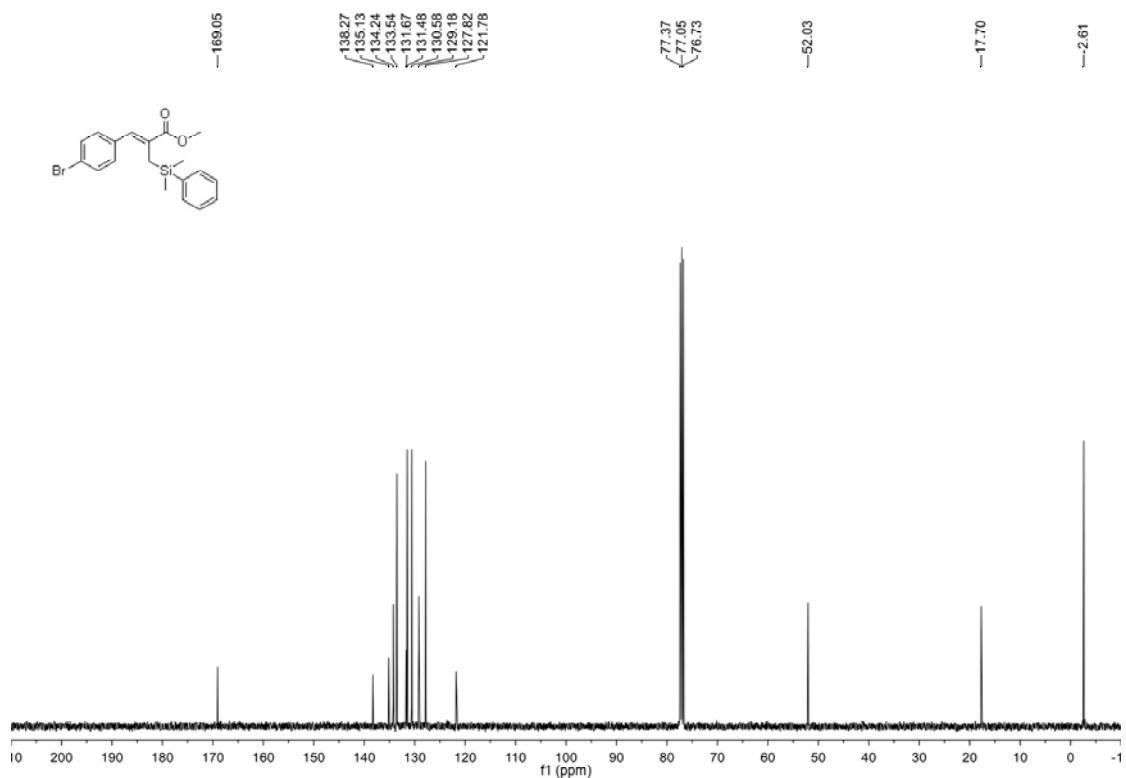
^1H NMR (400 MHz, CDCl_3) of compound **51**



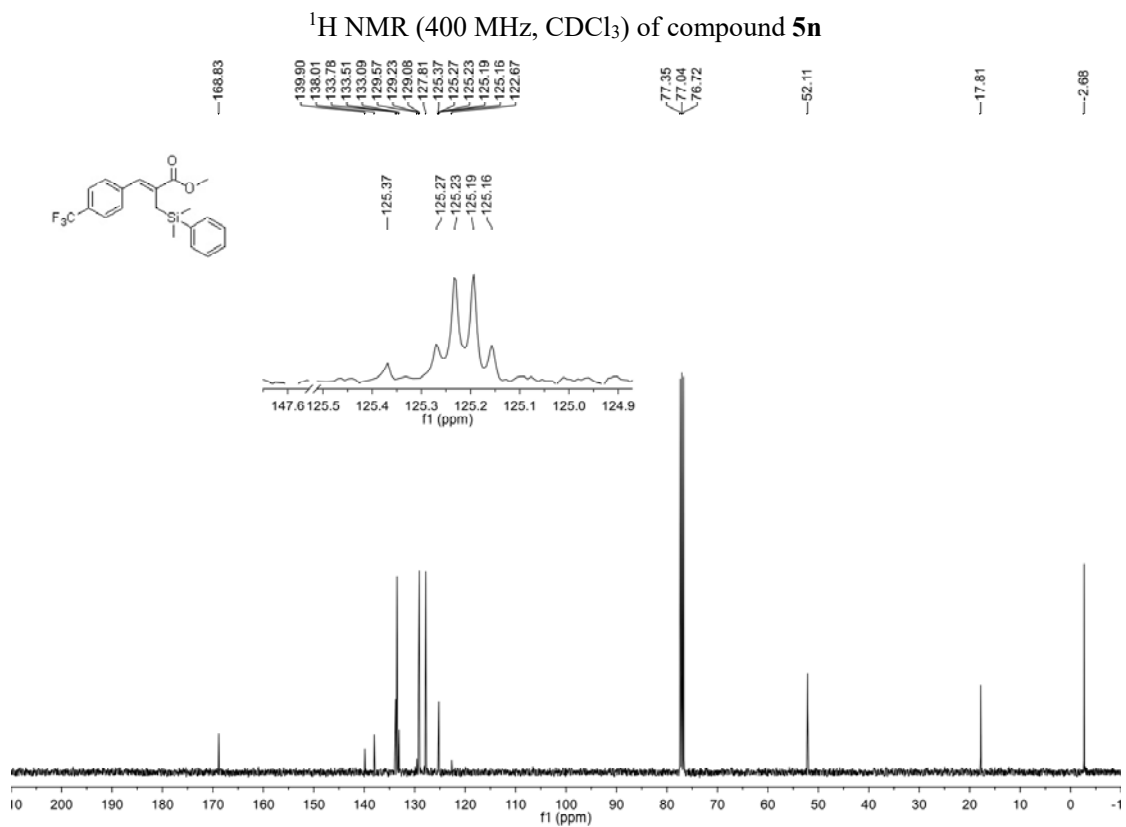
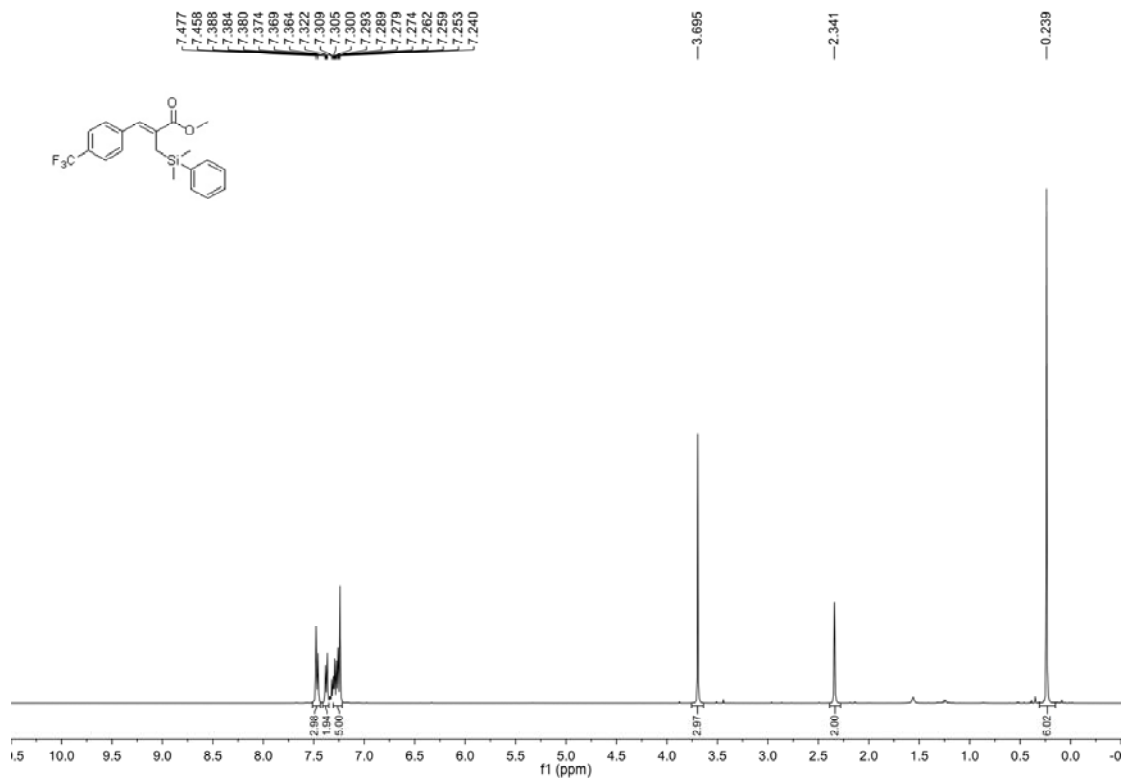
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **51**

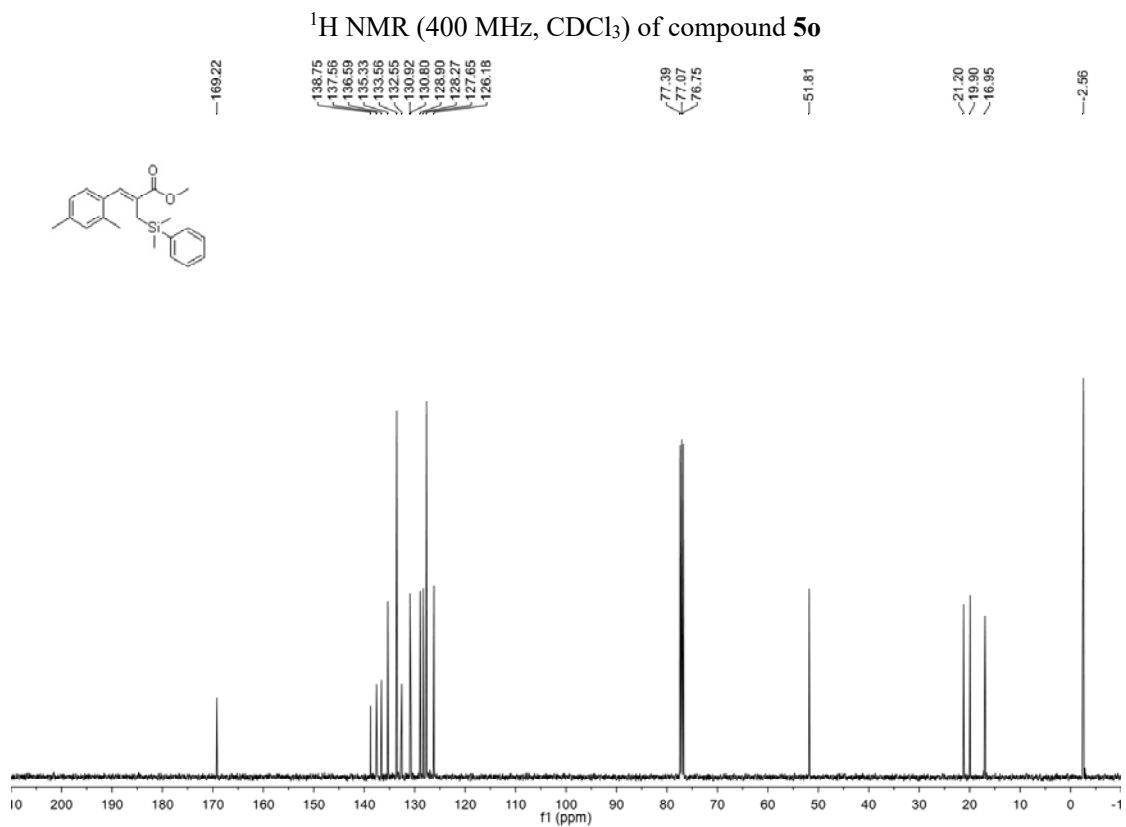
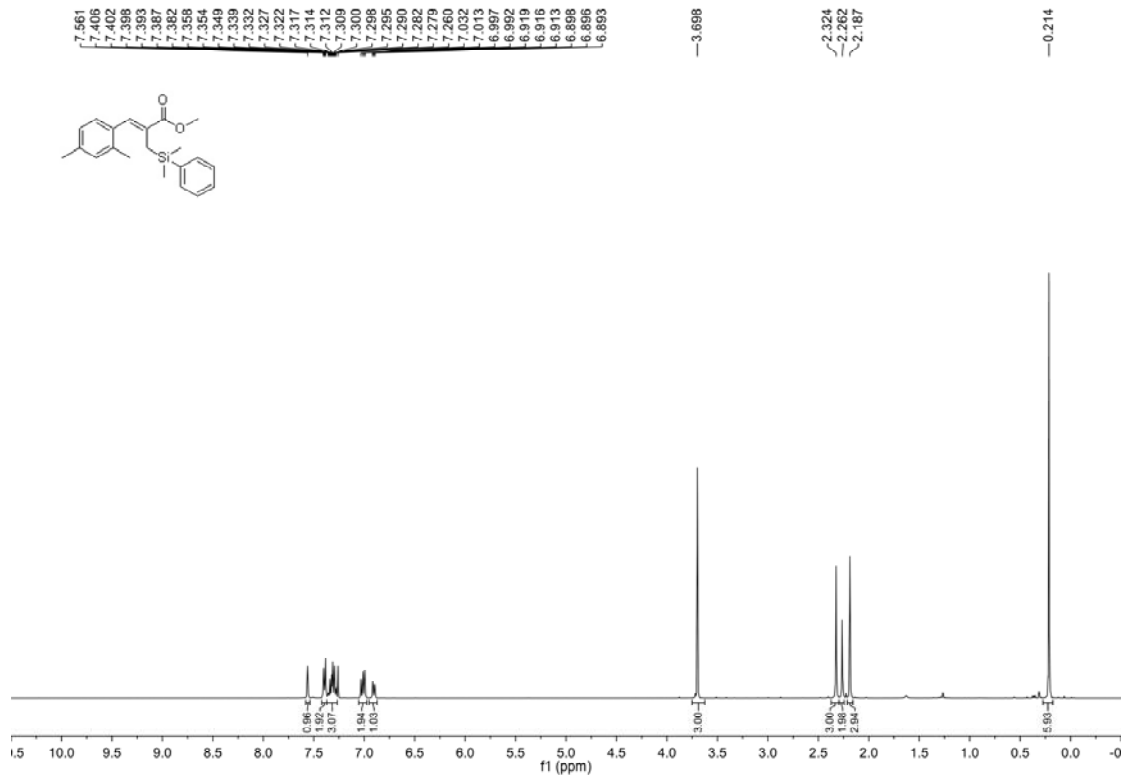


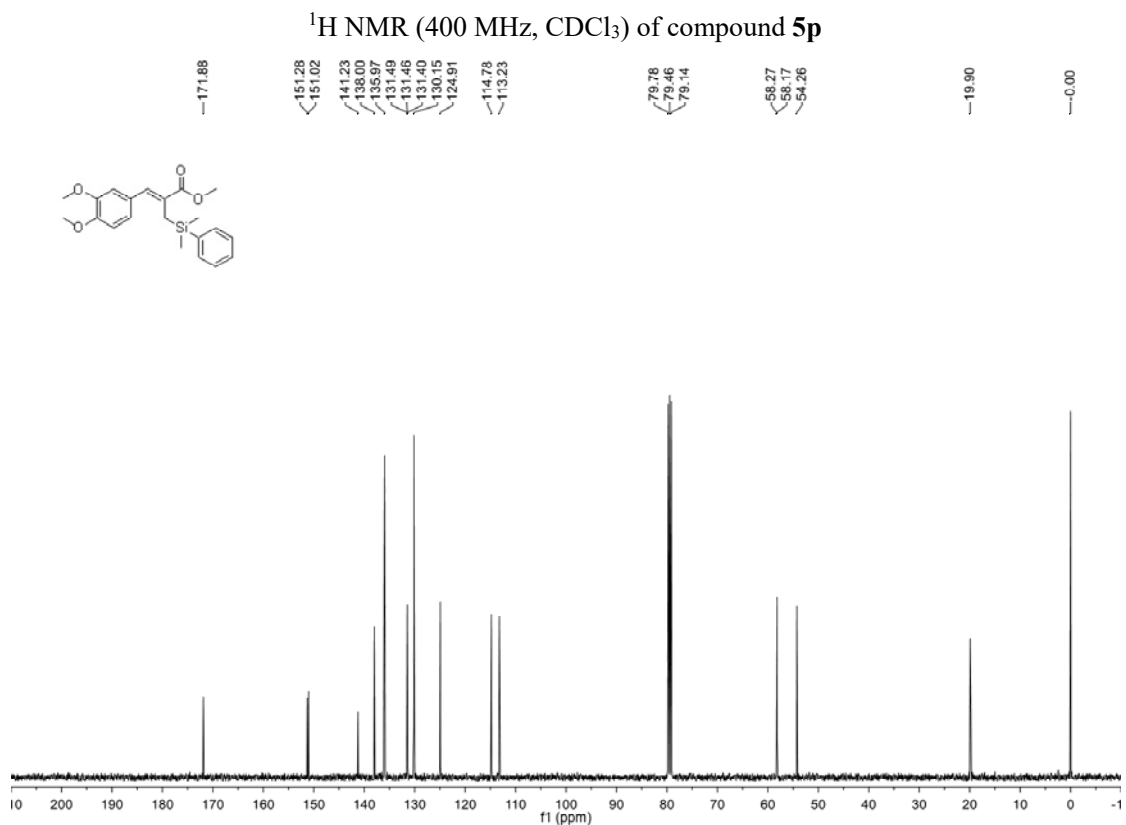
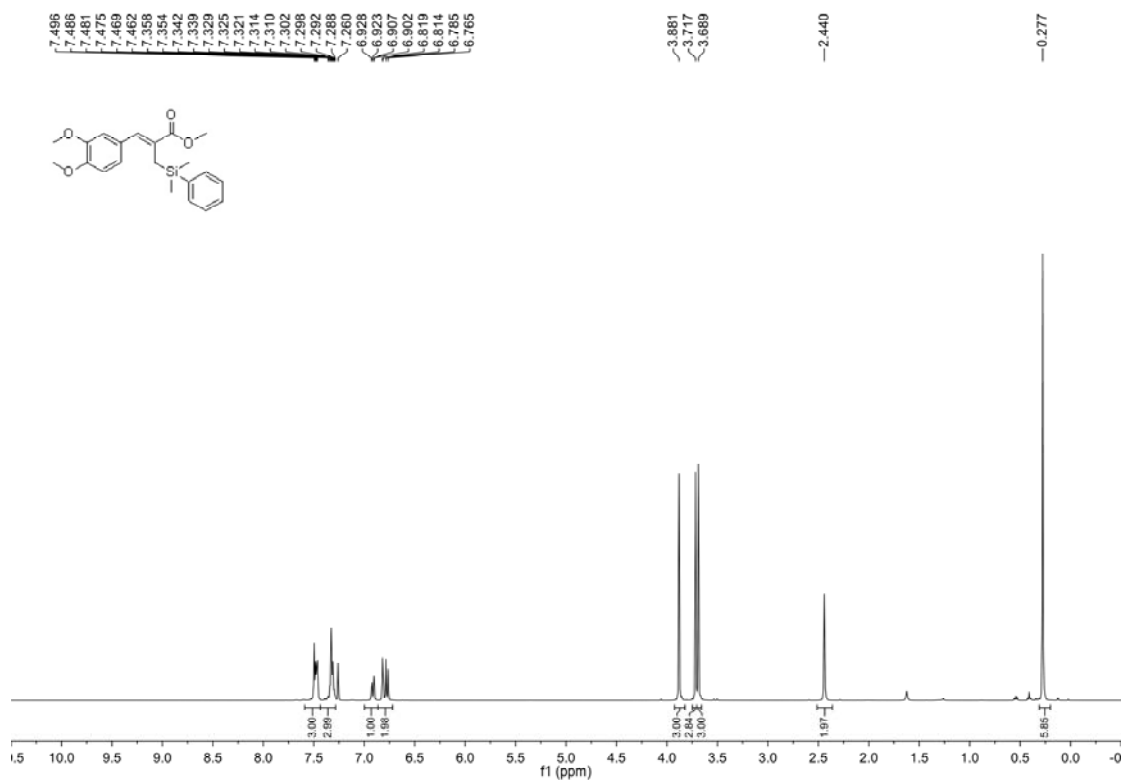
^1H NMR (400 MHz, CDCl_3) of compound **5m**

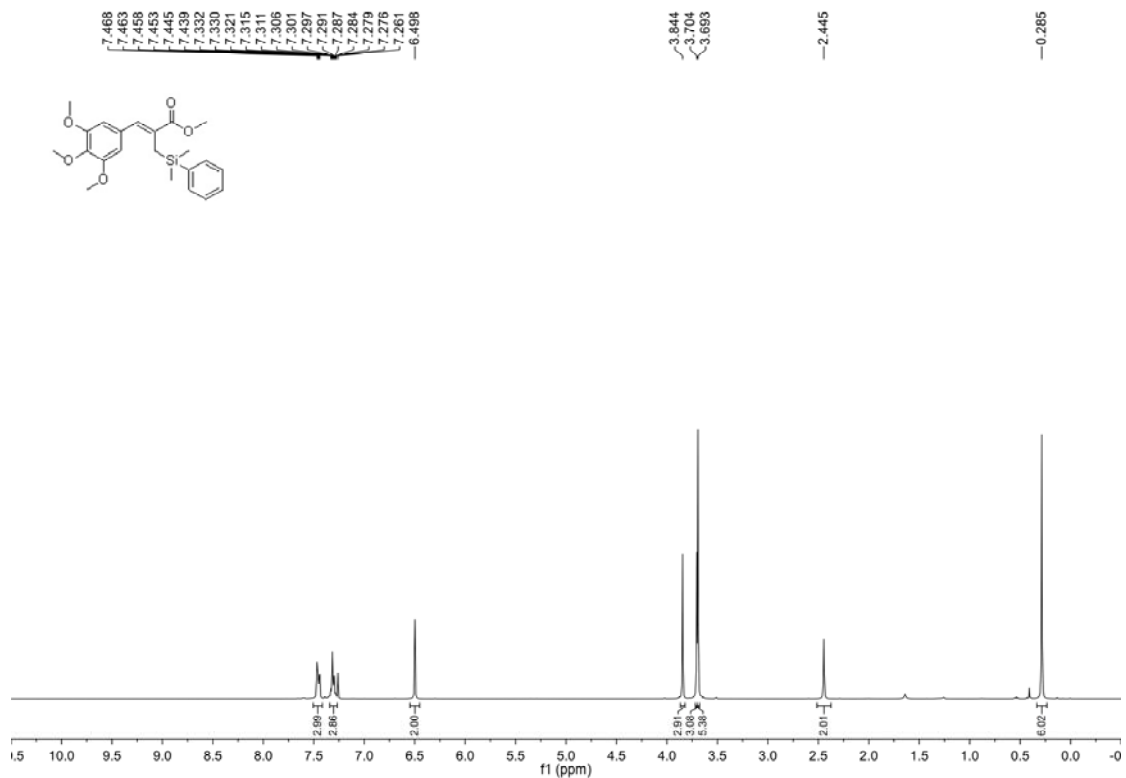


$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **5m**

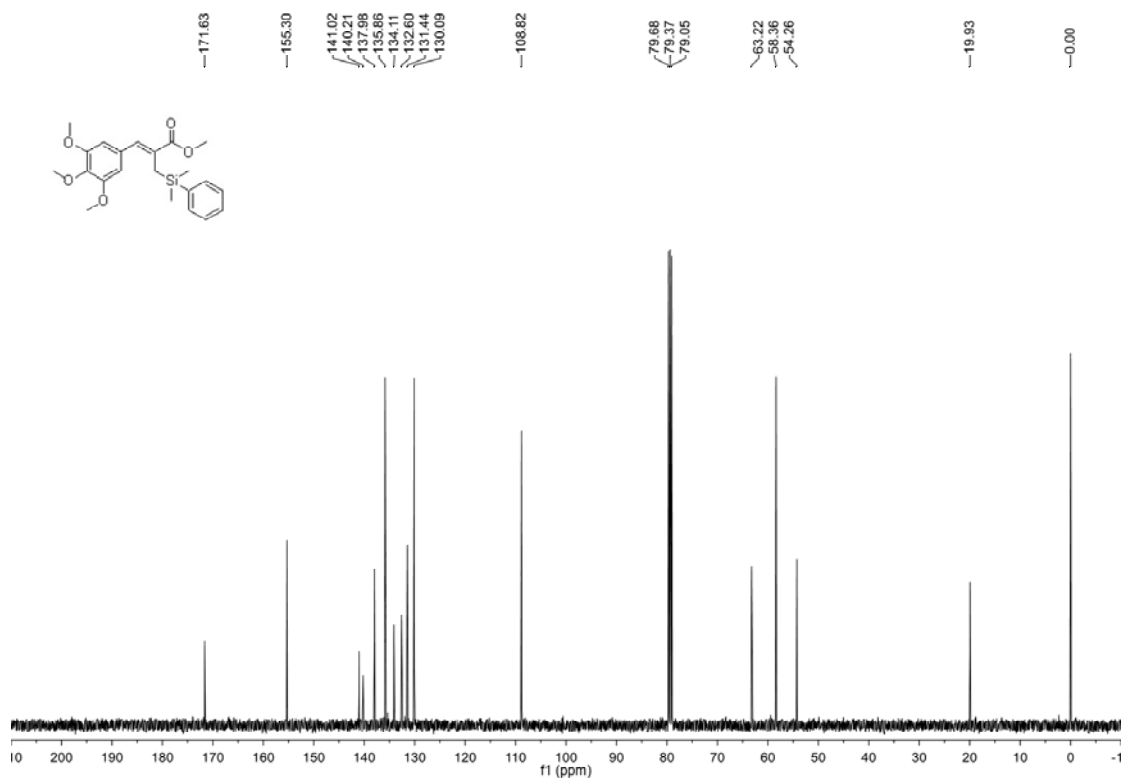




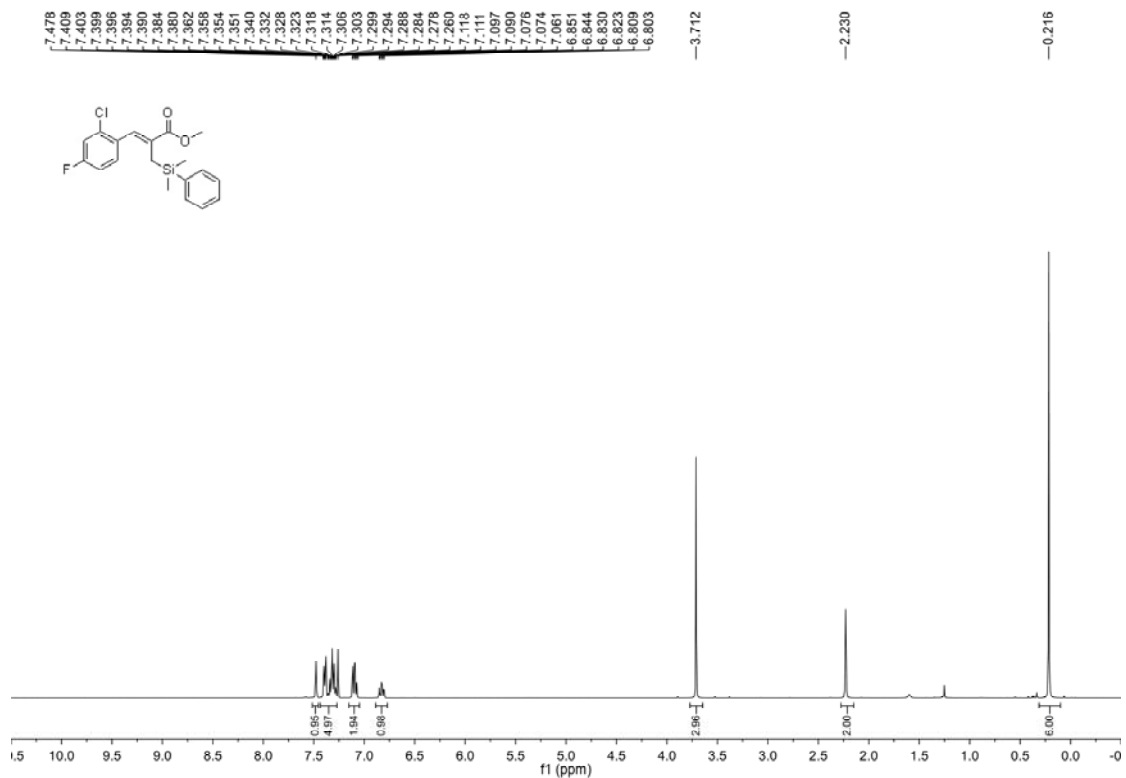




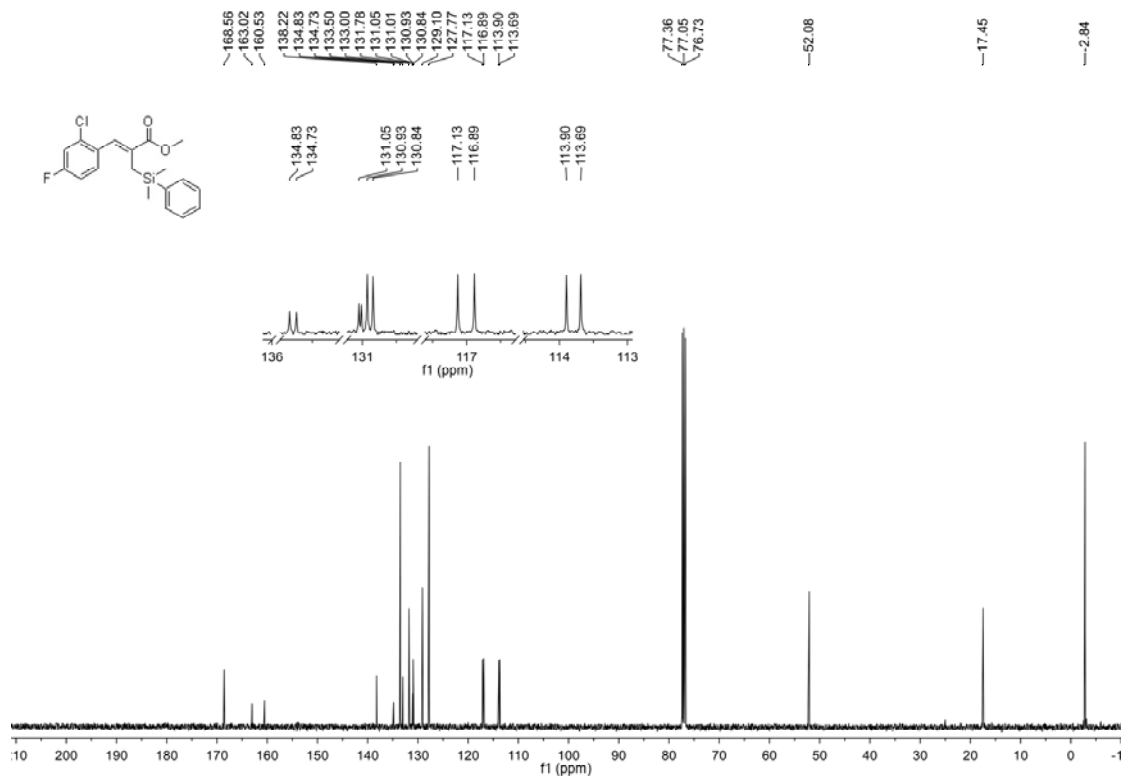
^1H NMR (400 MHz, CDCl_3) of compound **5q**



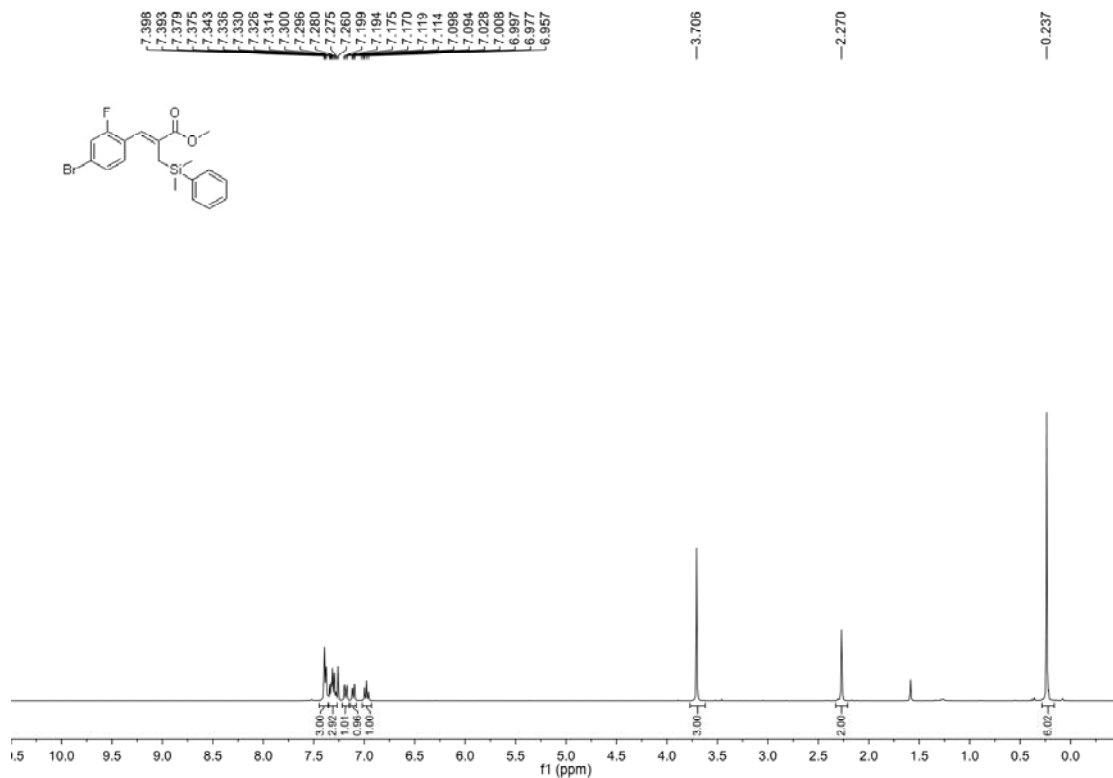
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **5q**



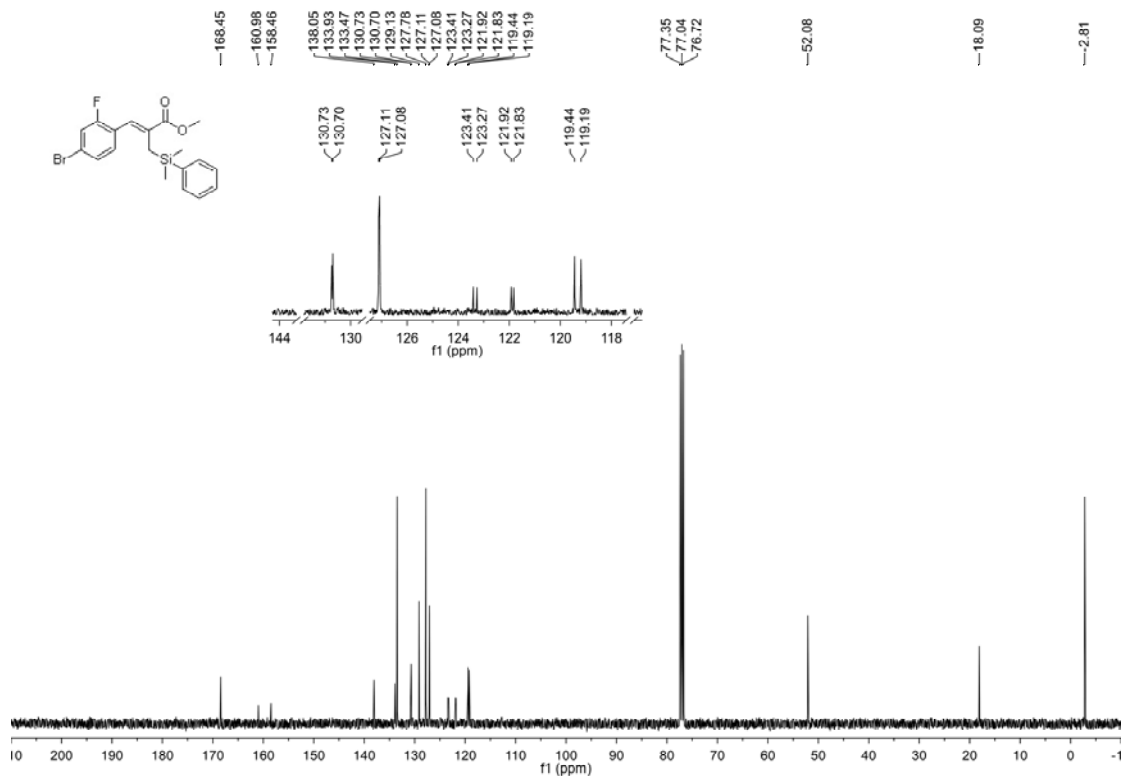
$^1\text{H NMR}$ (400 MHz, CDCl_3) of compound **5r**



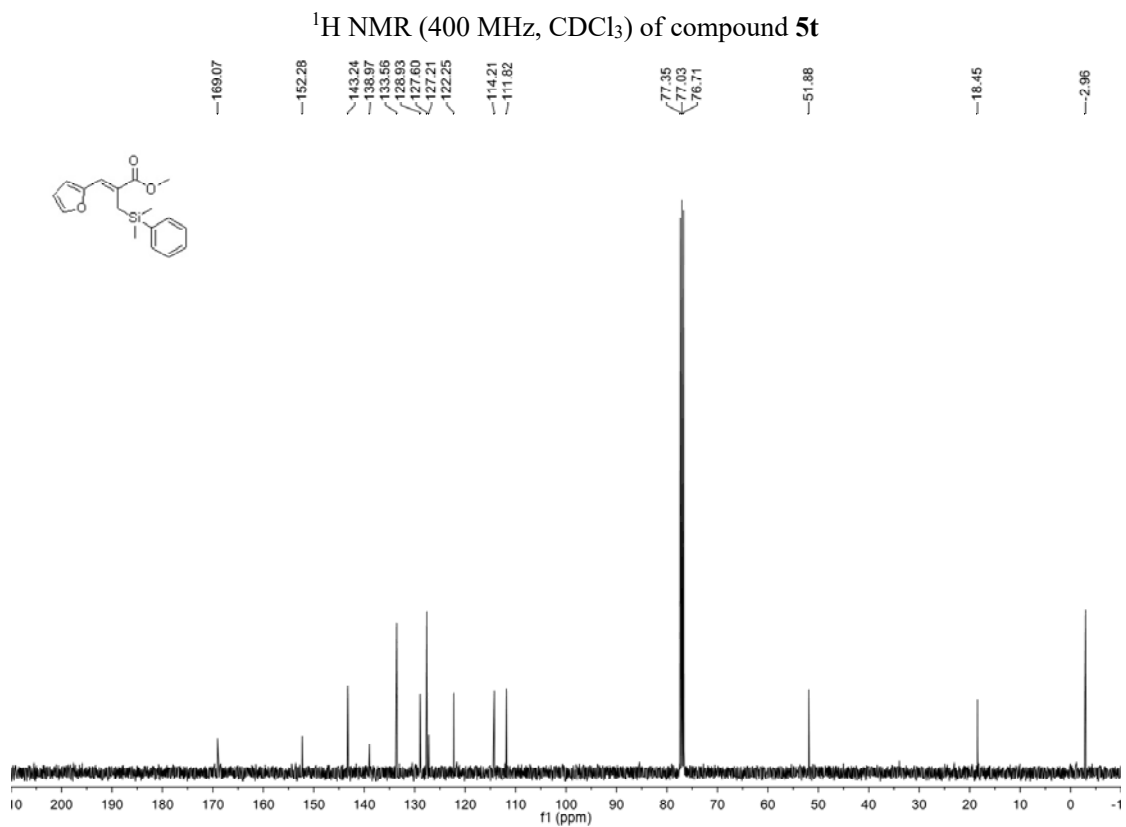
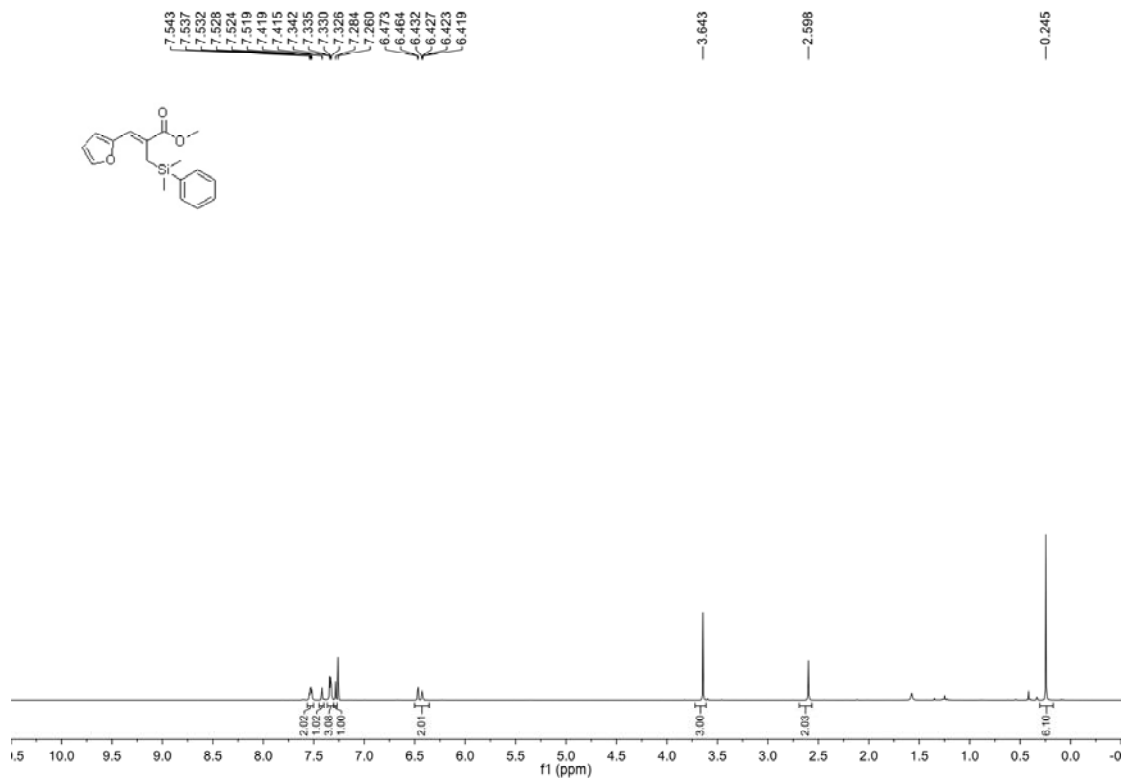
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **5r**



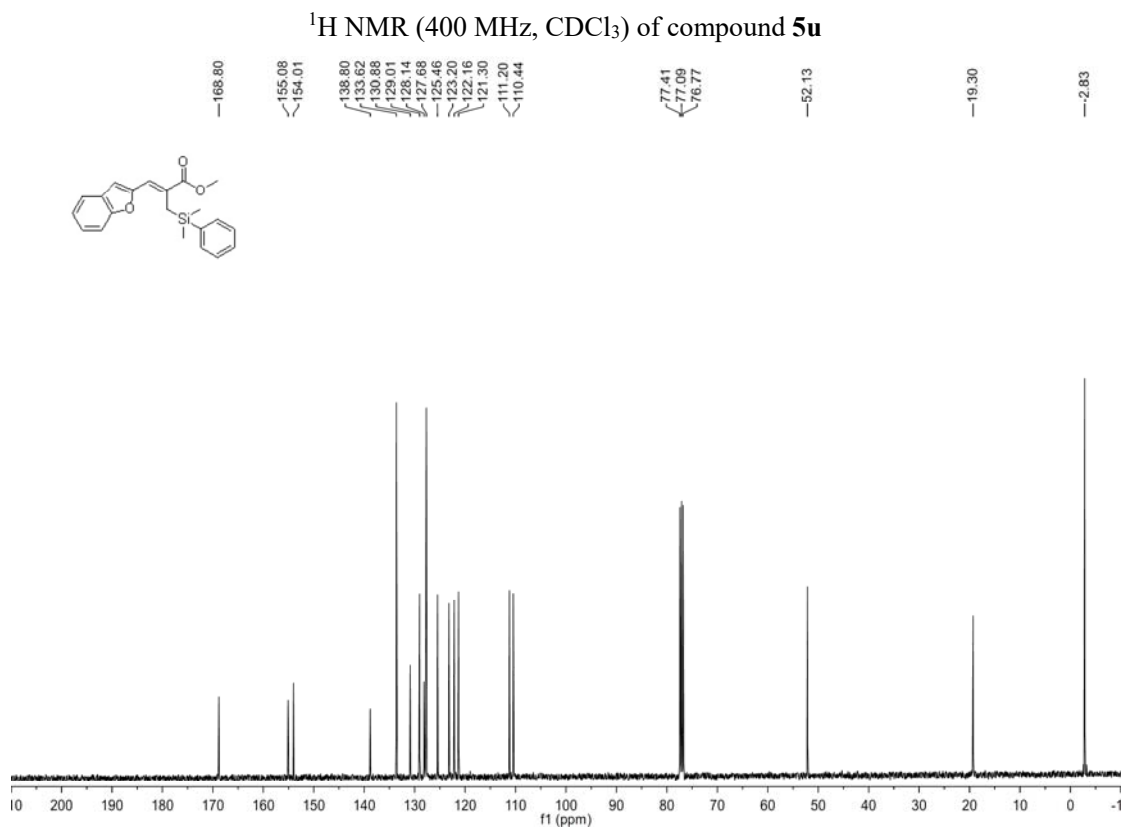
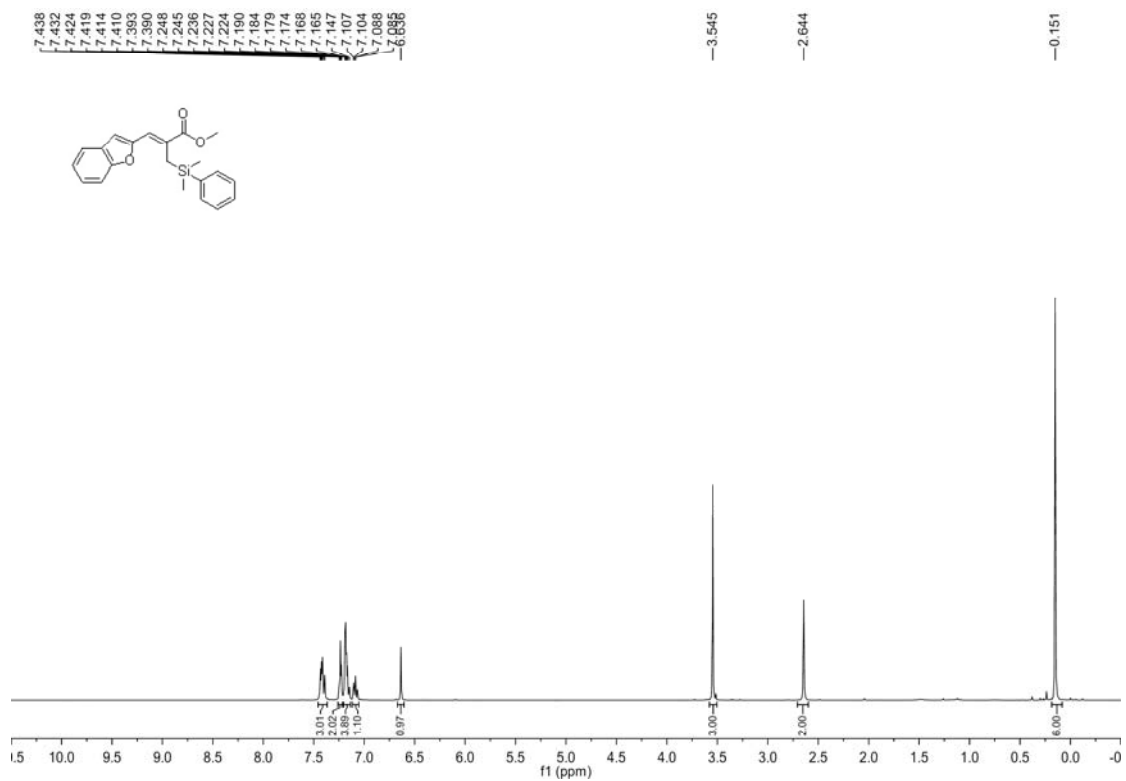
^1H NMR (400 MHz, CDCl_3) of compound **5s**

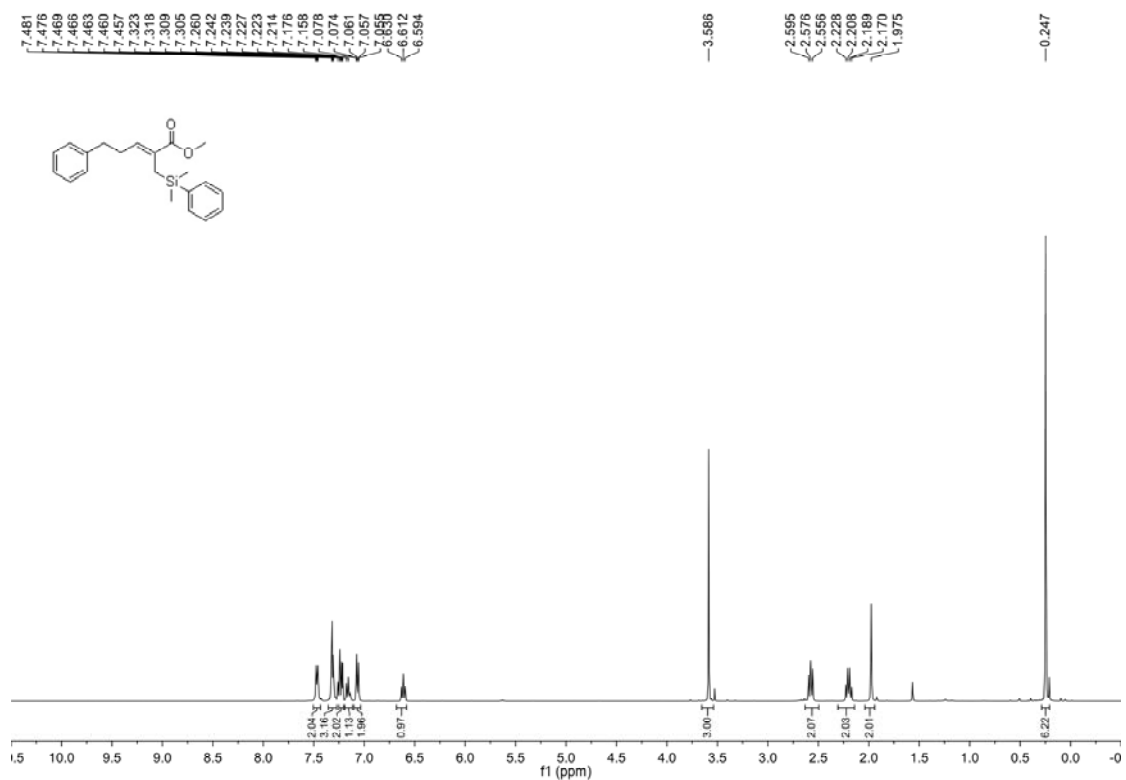


$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **5s**

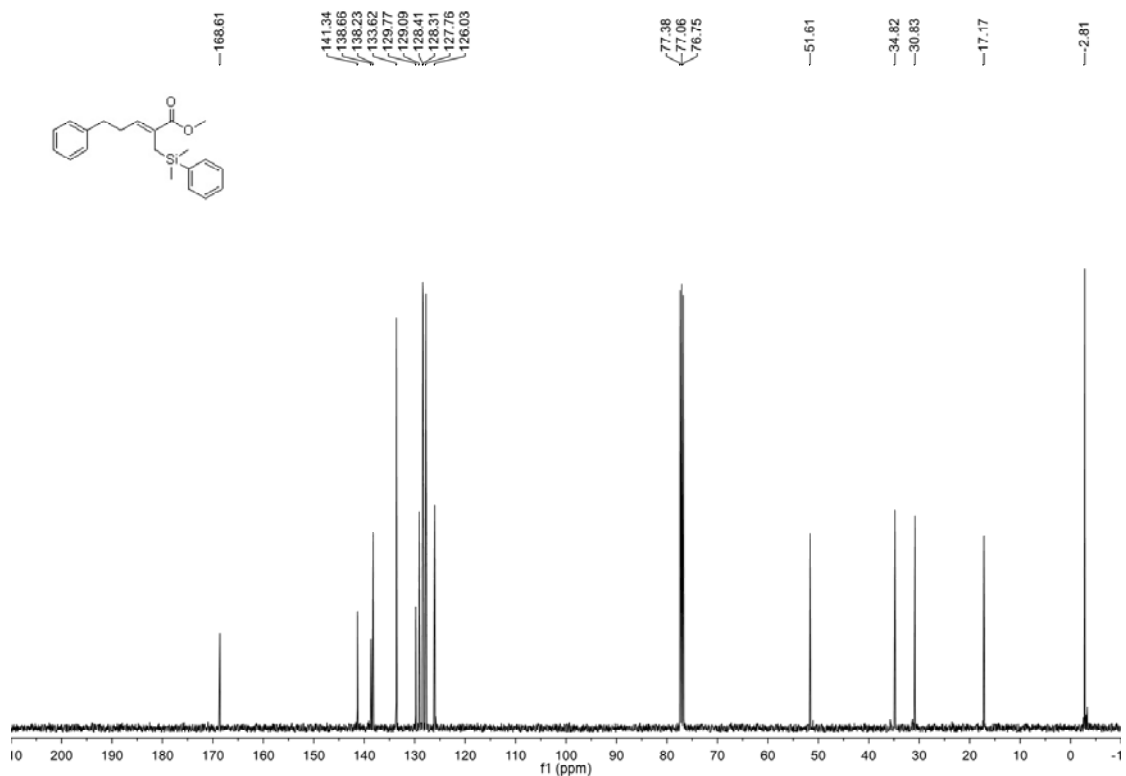


$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **5t**

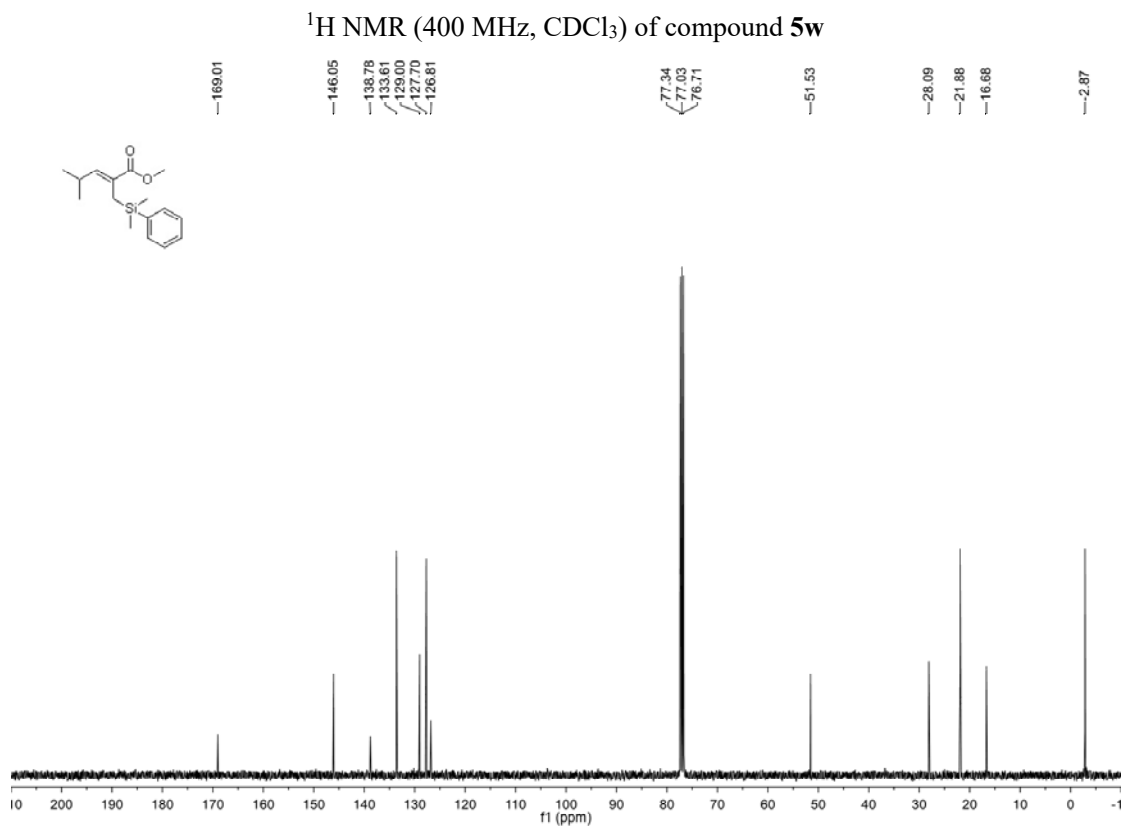
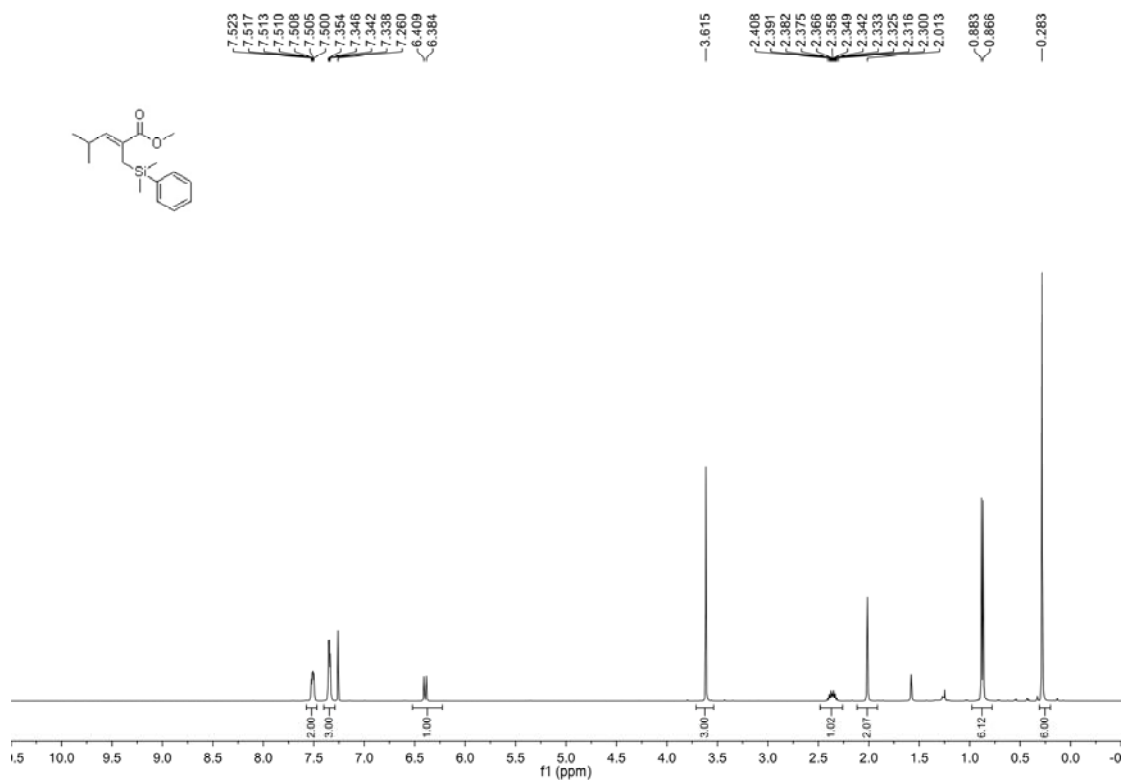




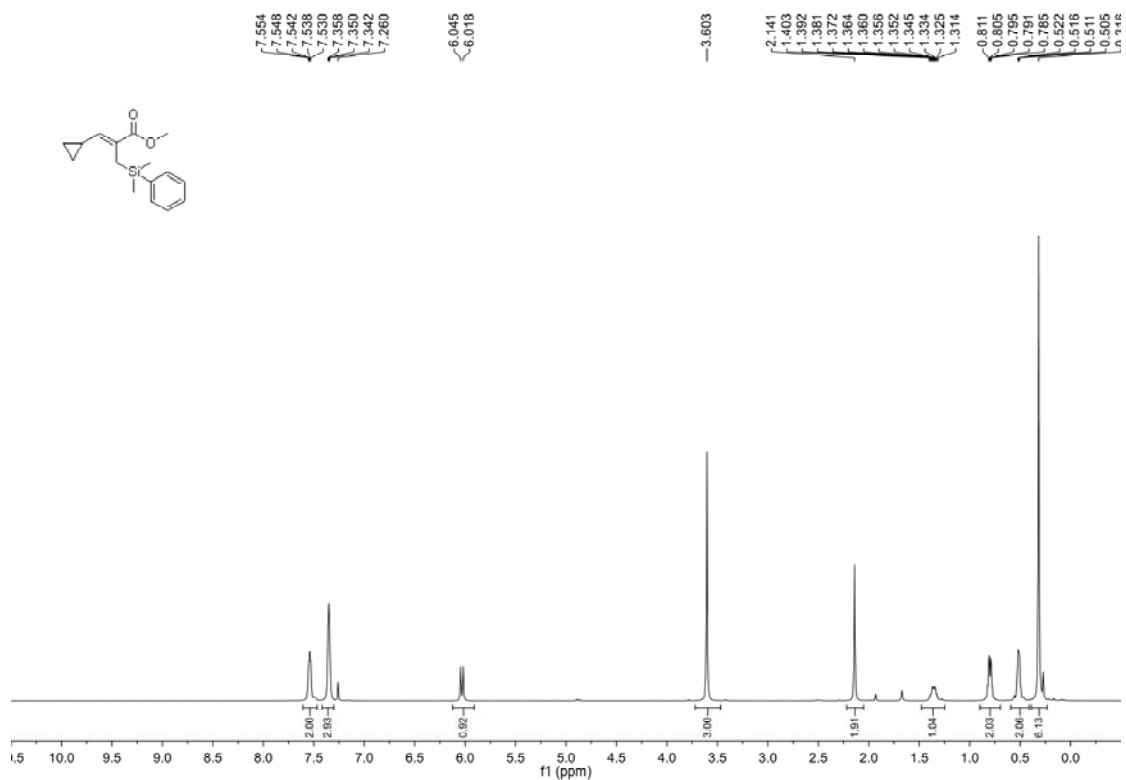
^1H NMR (400 MHz, CDCl_3) of compound **5v**



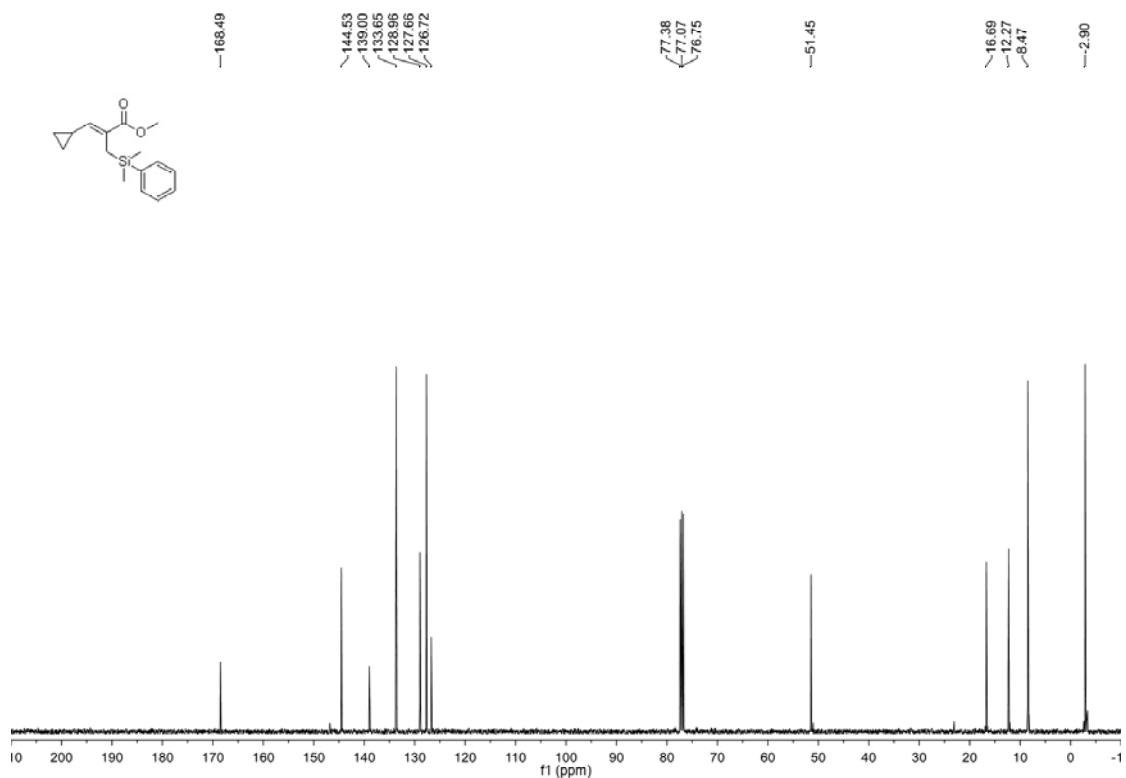
^{13}C NMR (100 MHz, CDCl_3) of compound **5v**



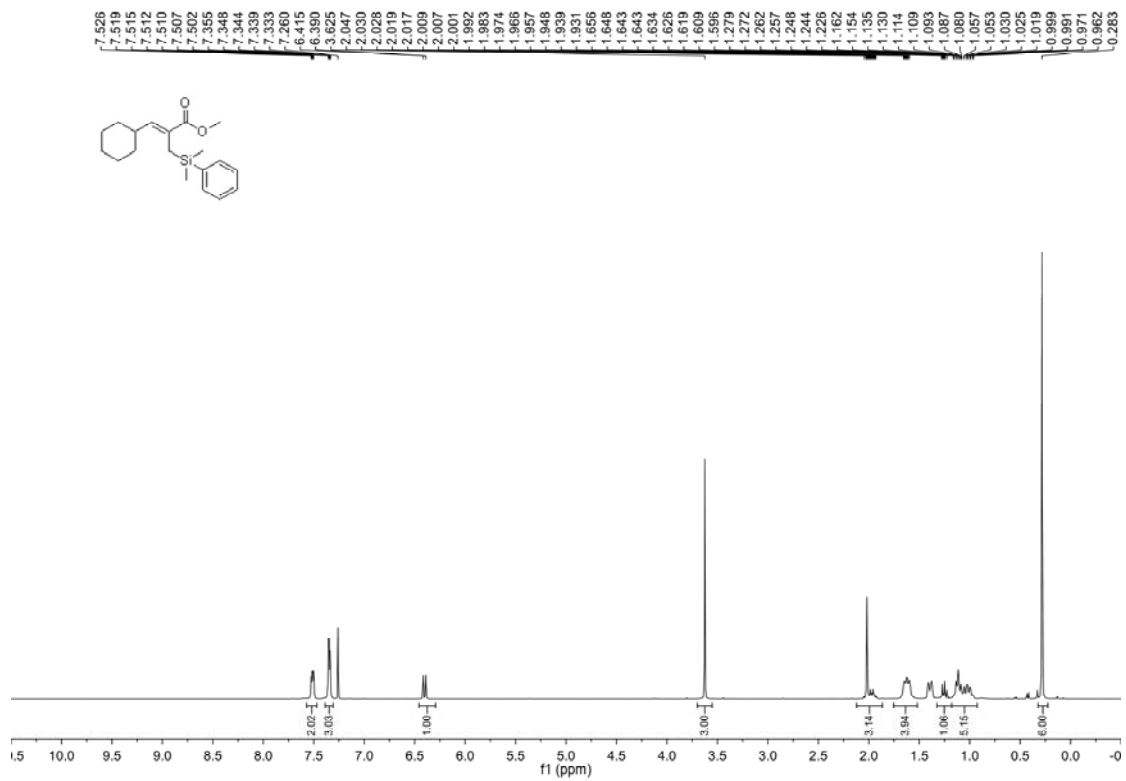
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **5w**



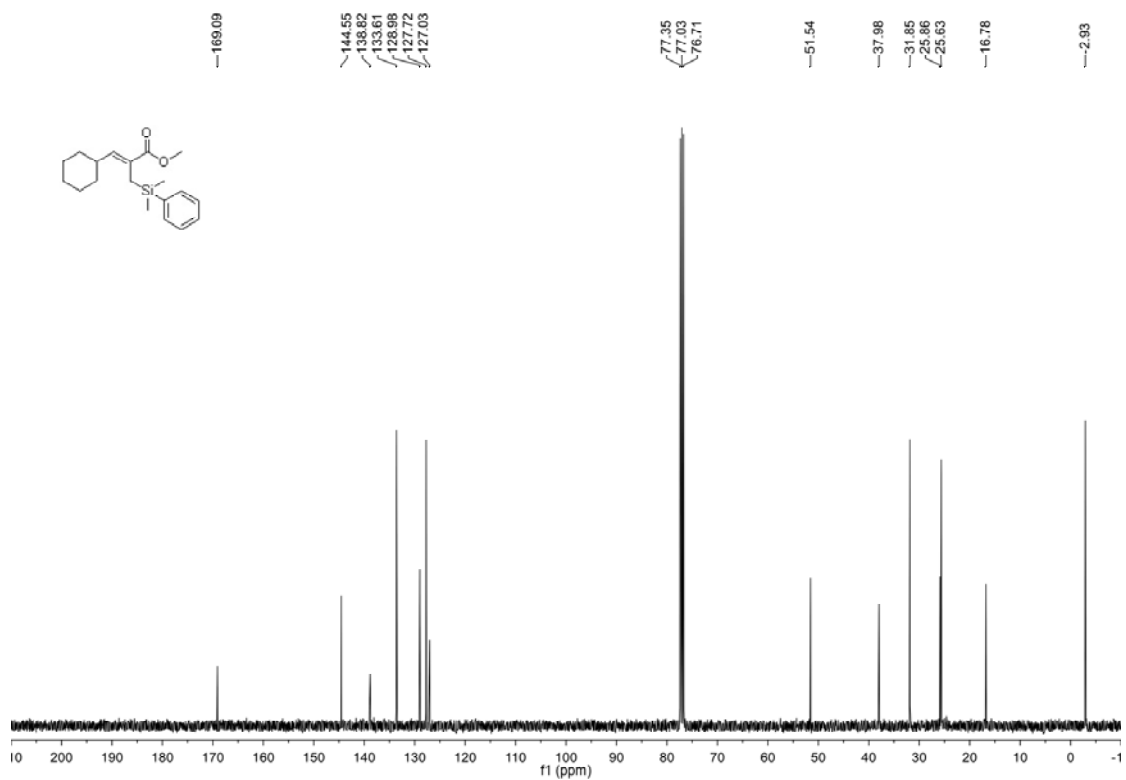
¹H NMR (400 MHz, CDCl₃) of compound **5x**



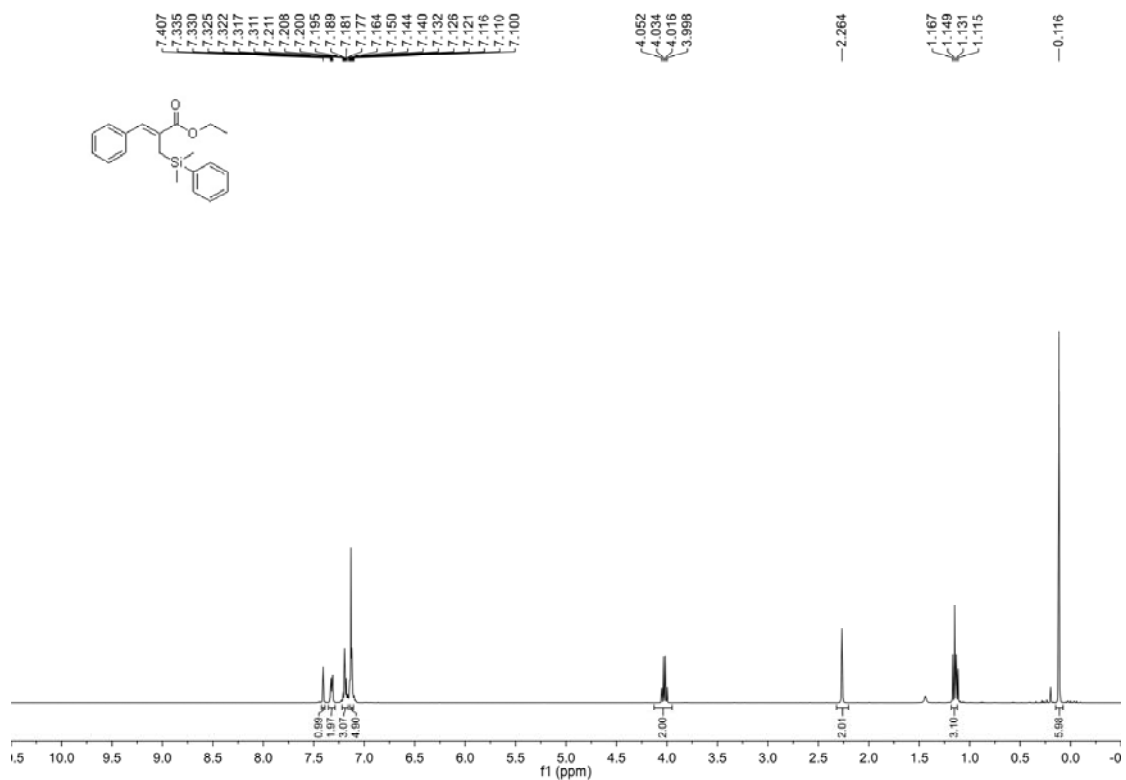
¹³C NMR (100 MHz, CDCl₃) of compound **5x**



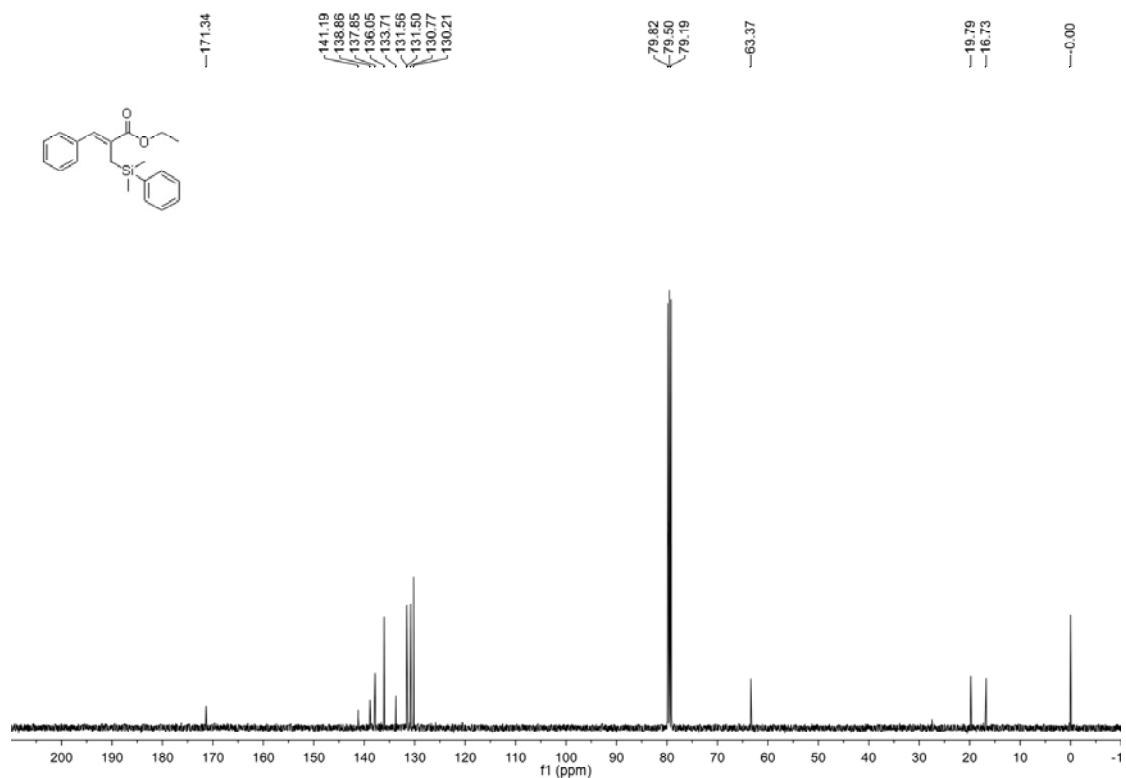
^1H NMR (400 MHz, CDCl_3) of compound **5y**



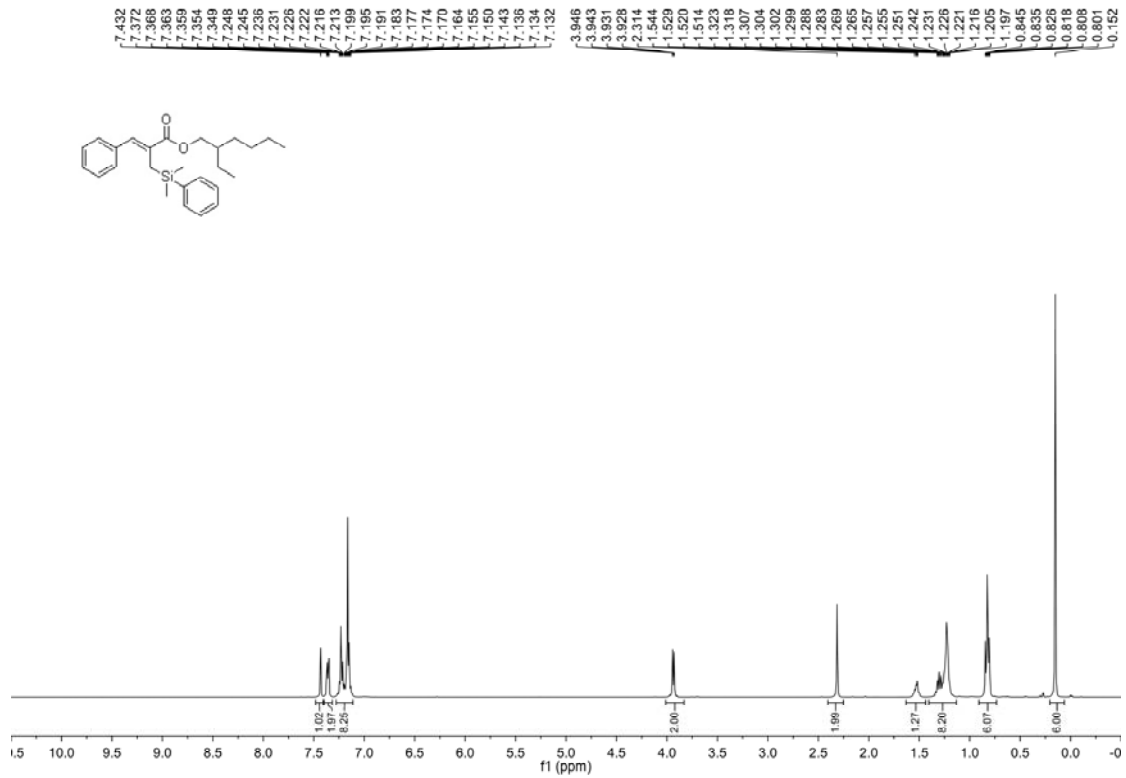
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **5y**



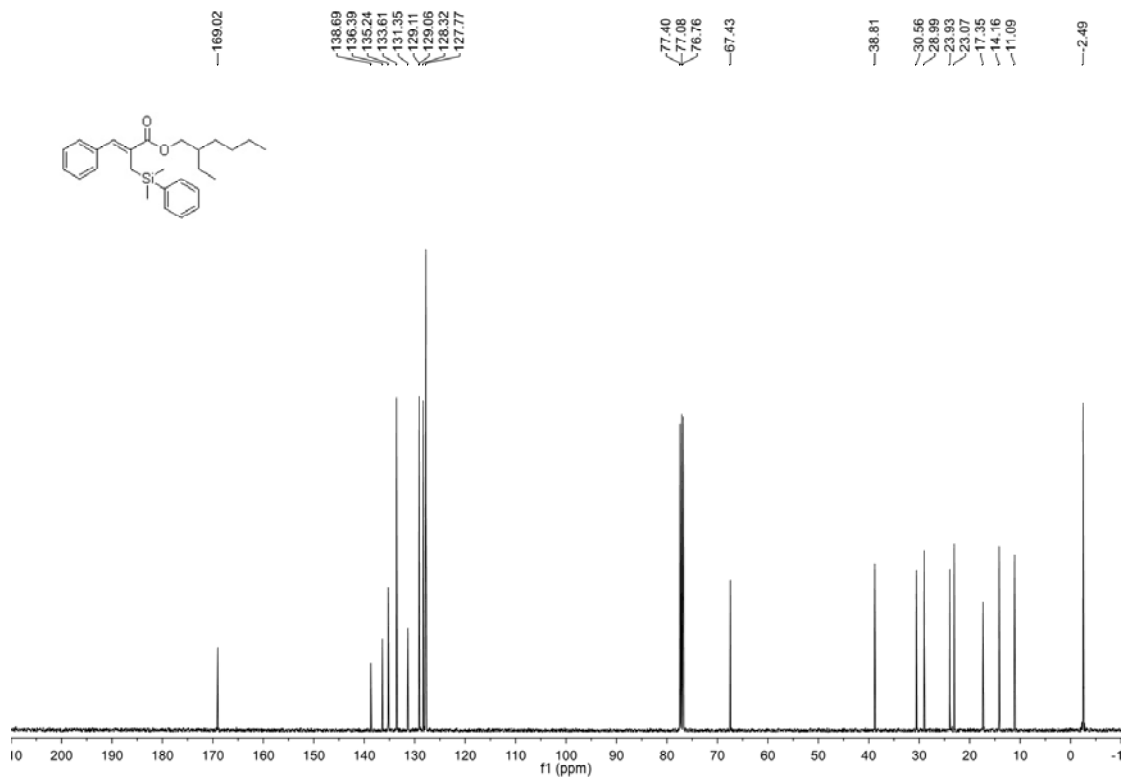
^1H NMR (400 MHz, CDCl_3) of compound **5z**



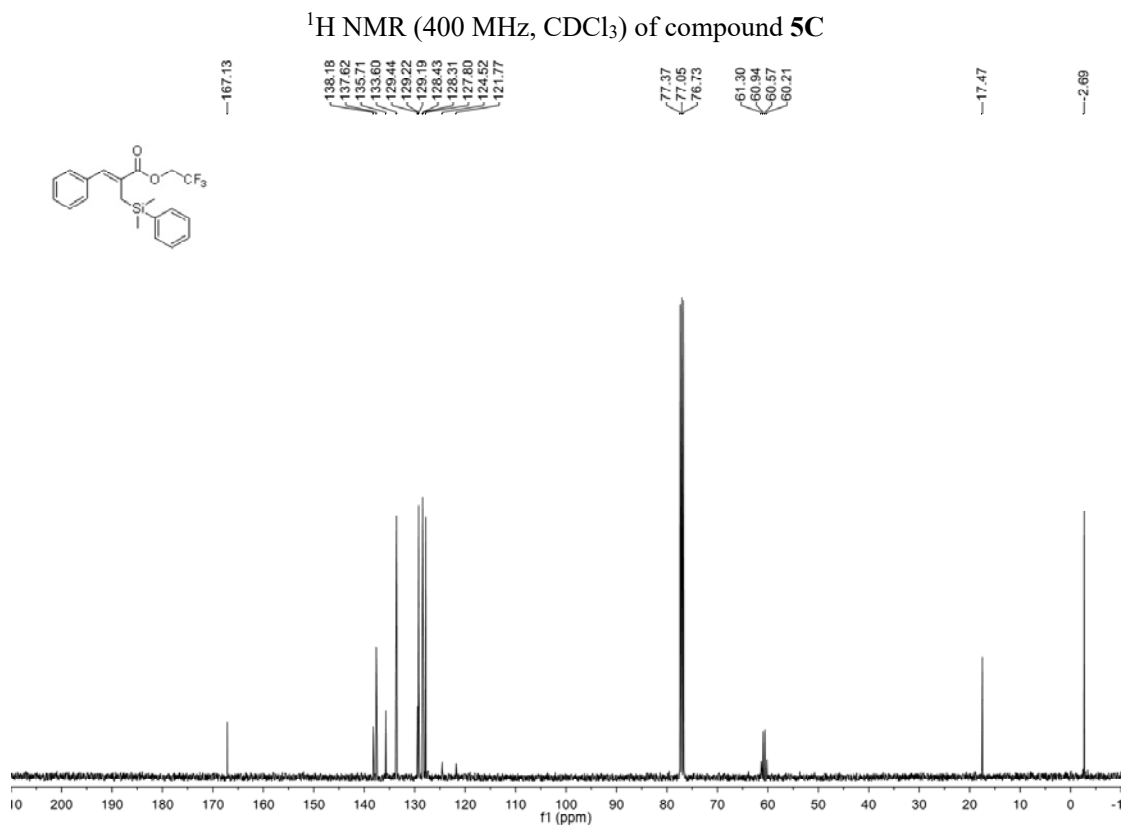
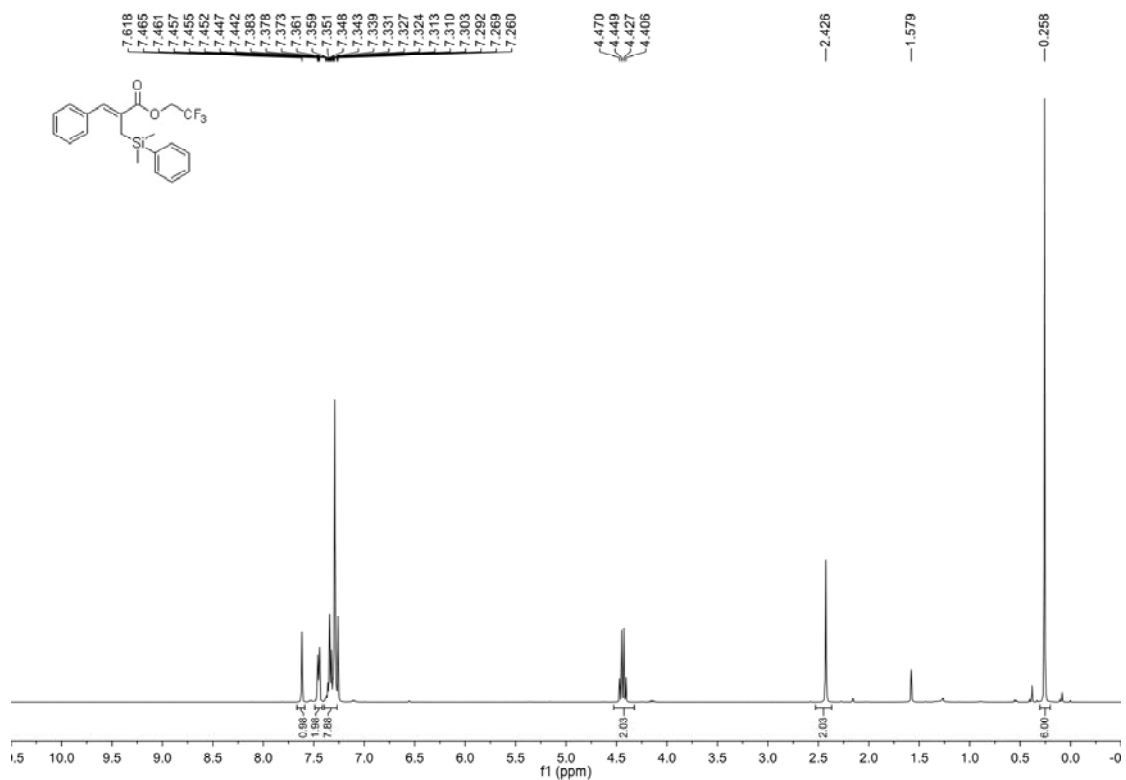
^{13}C NMR (100 MHz, CDCl_3) of compound **5z**

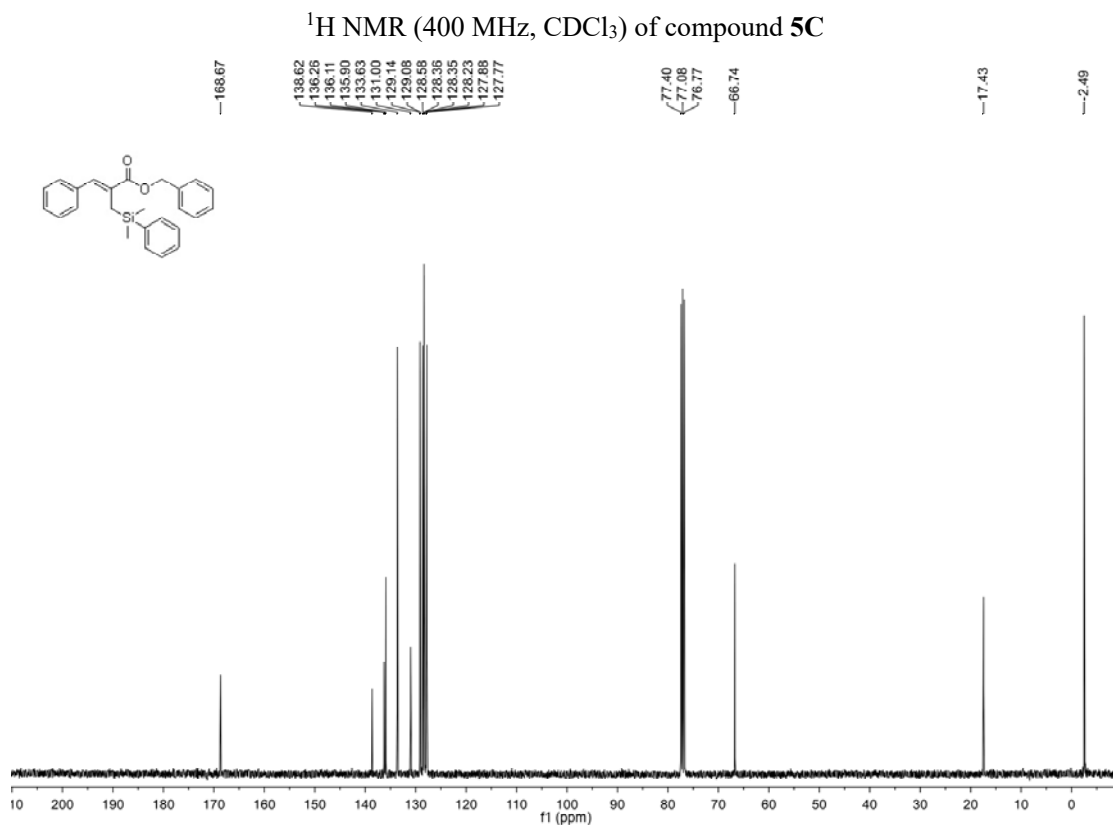
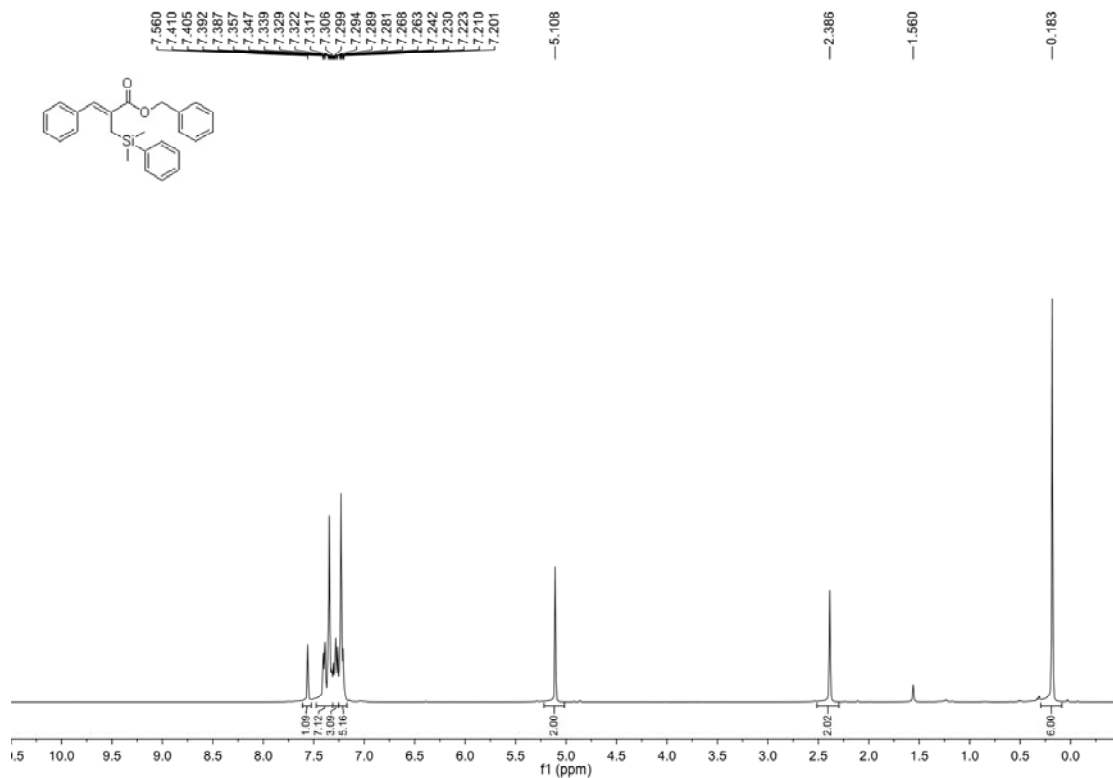


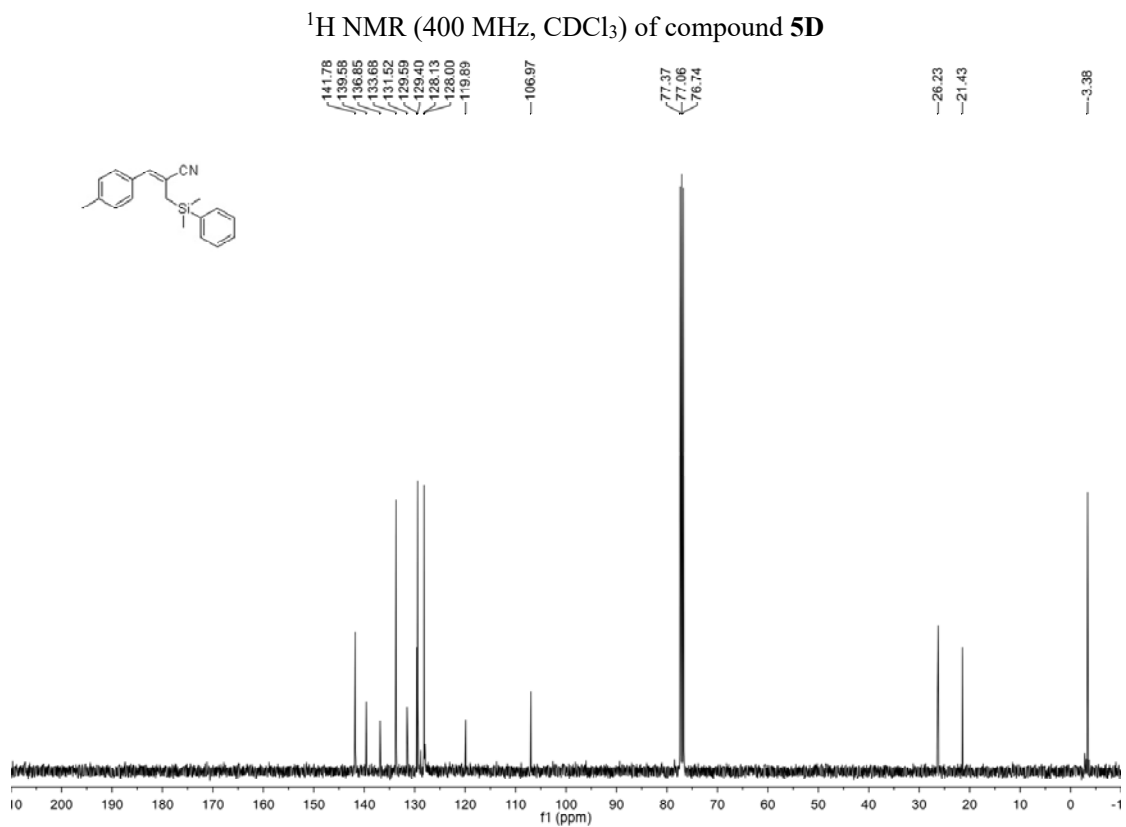
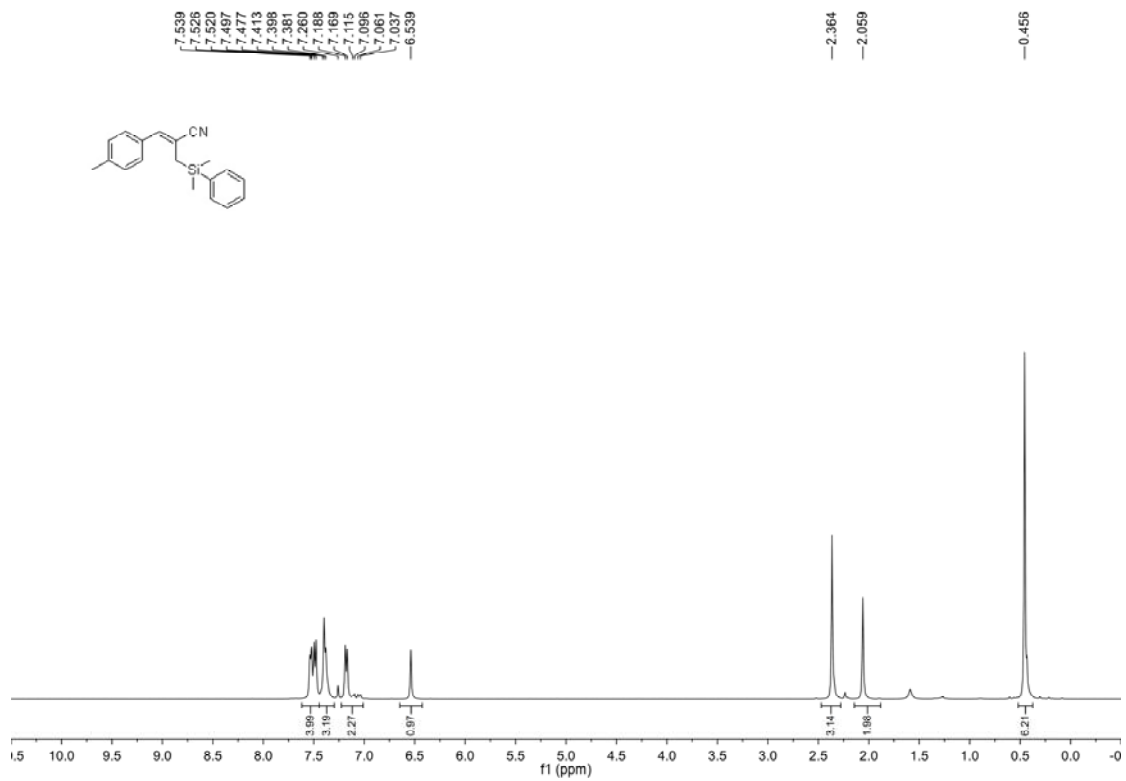
^1H NMR (400 MHz, CDCl_3) of compound 5A

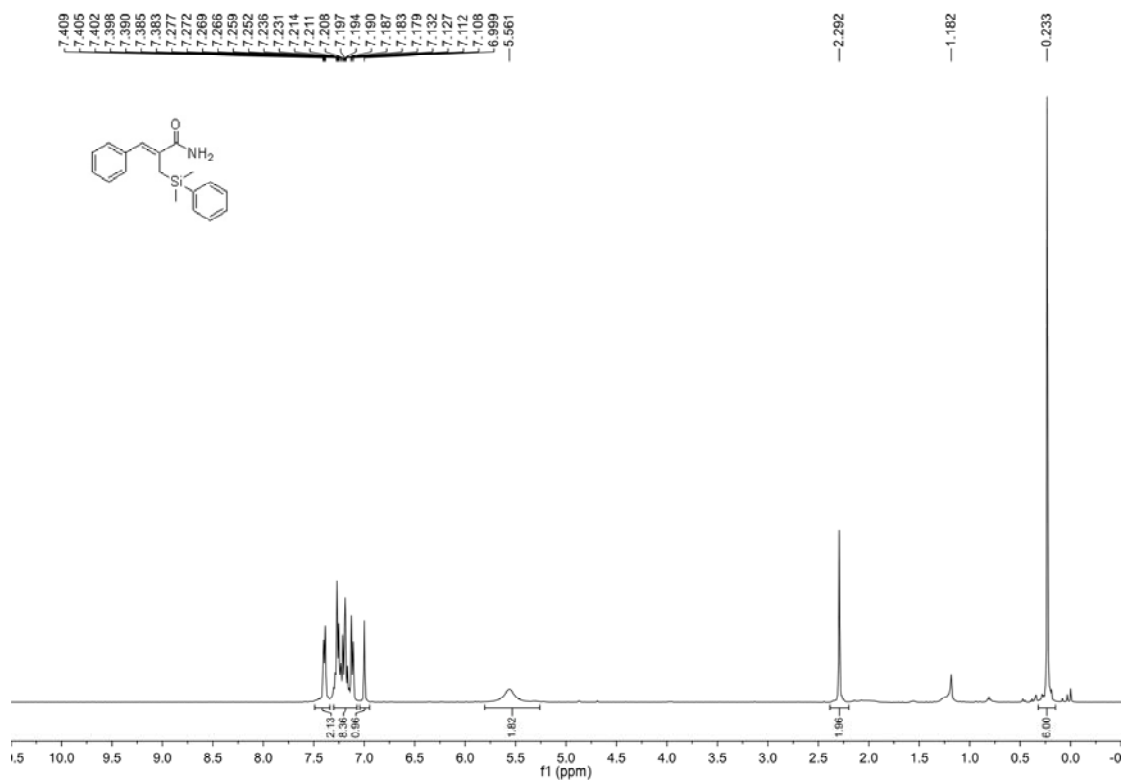


$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound 5A

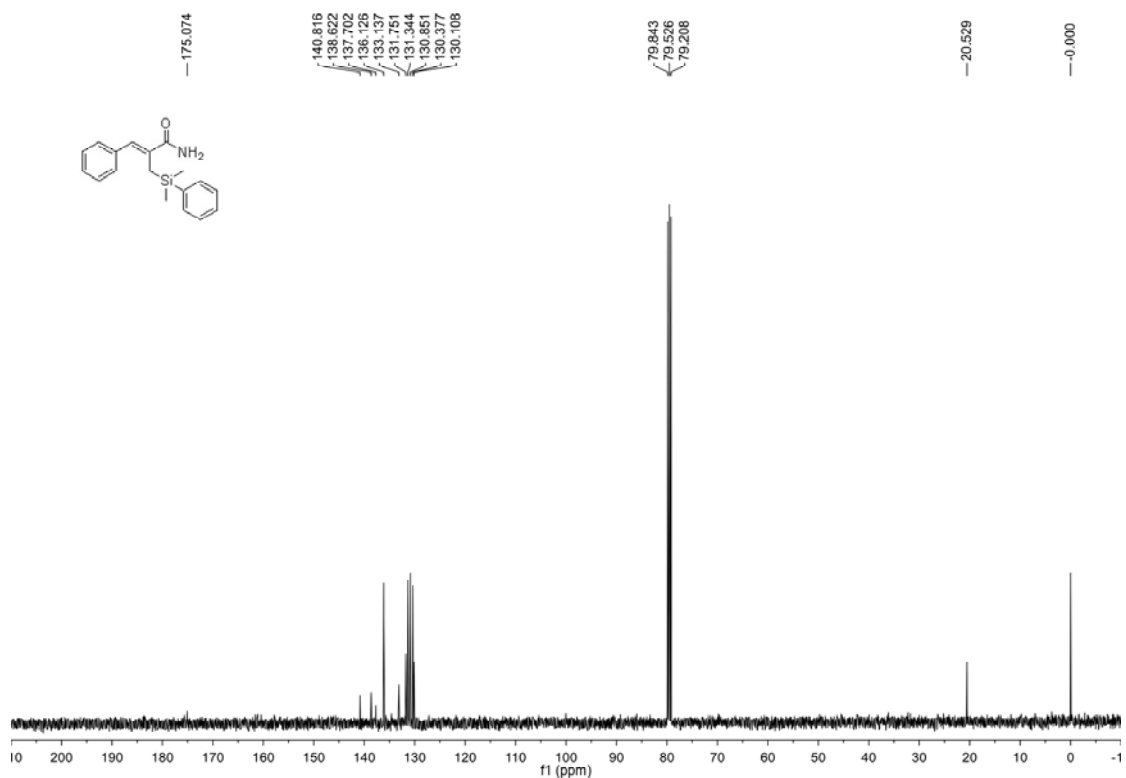




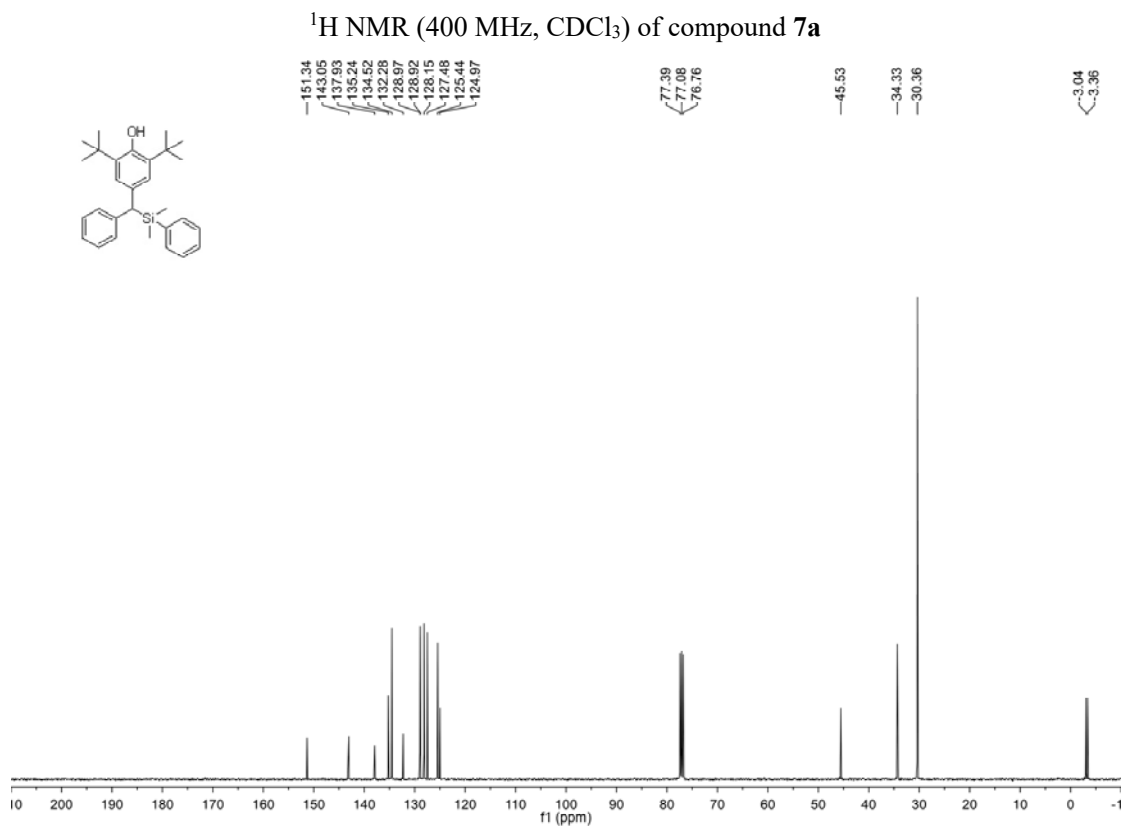
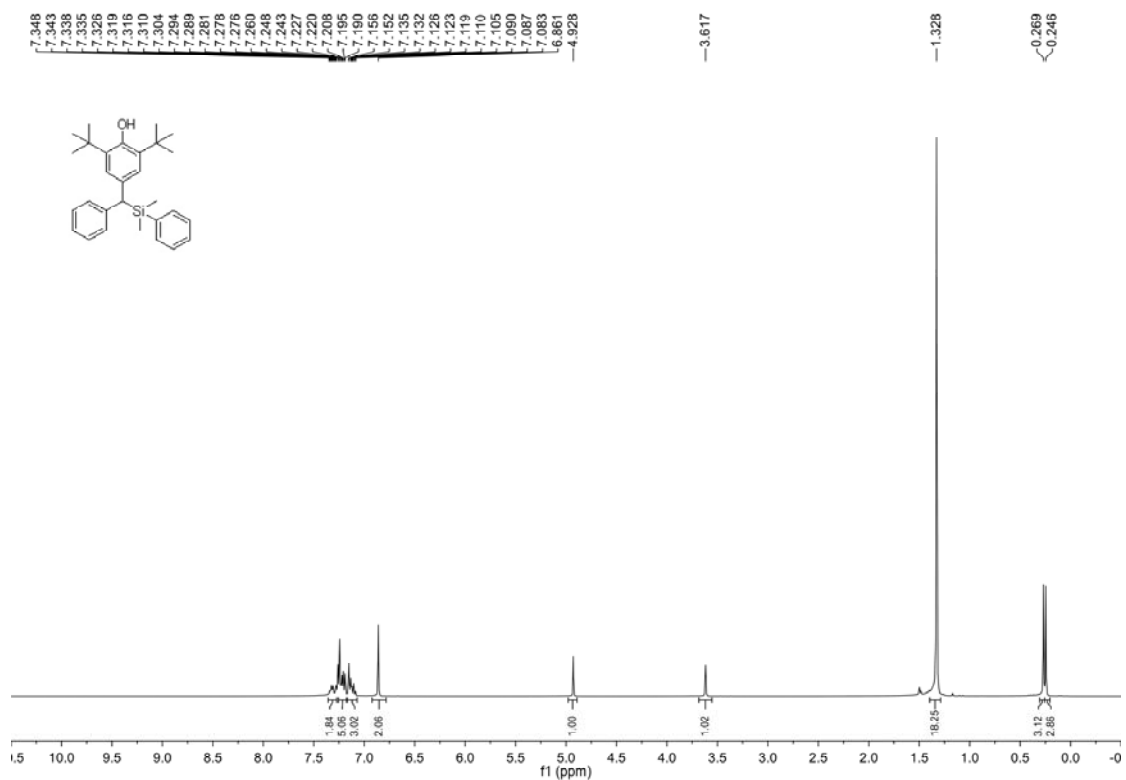


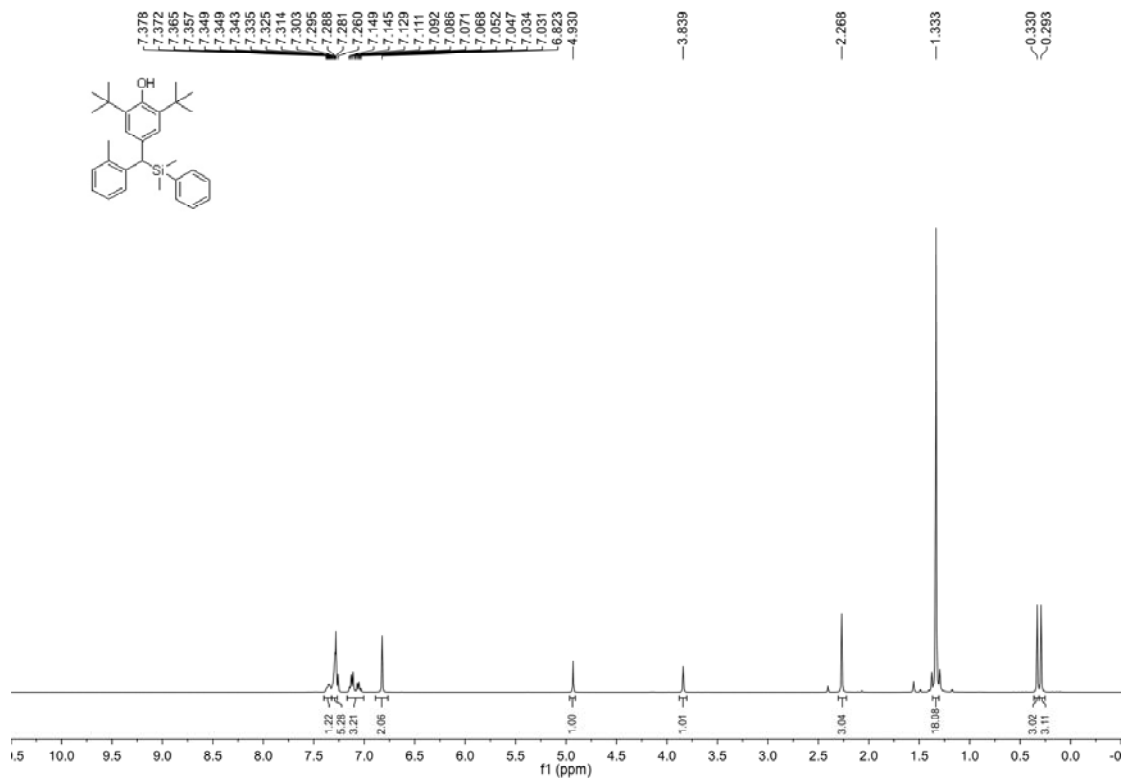


^1H NMR (400 MHz, CDCl_3) of compound **5E**

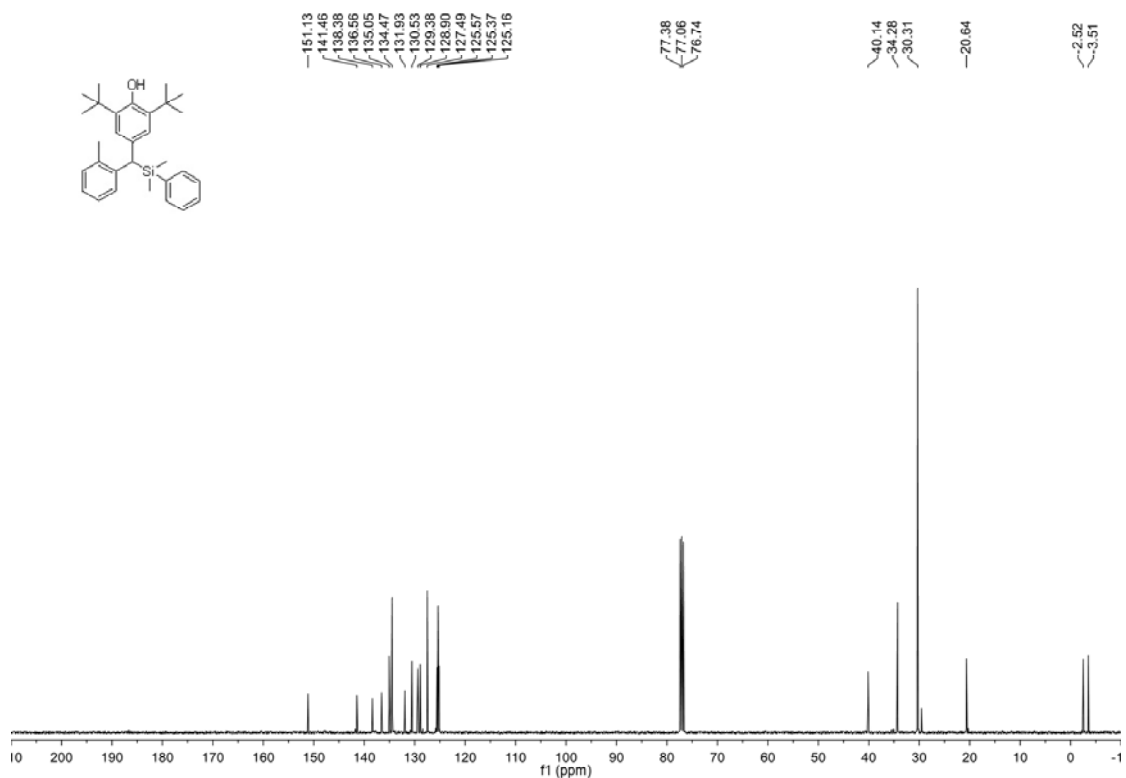


^{13}C NMR (100 MHz, CDCl_3) of compound **5E**

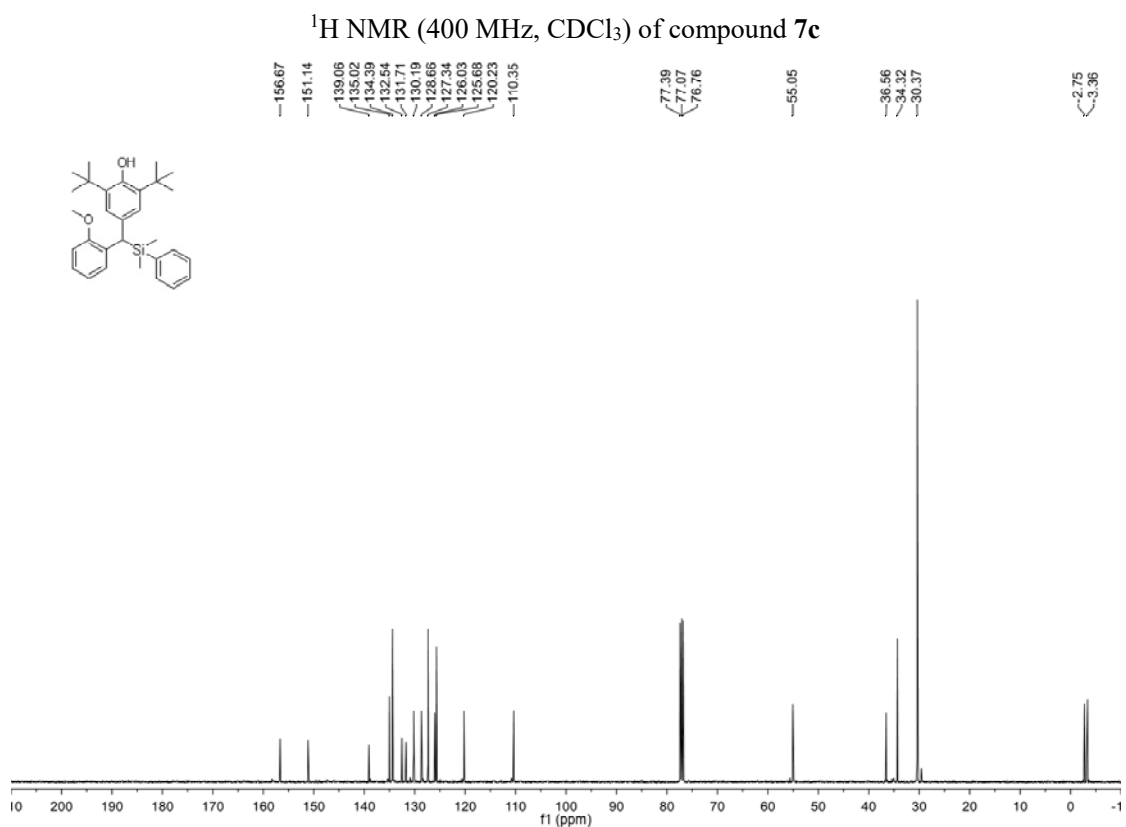
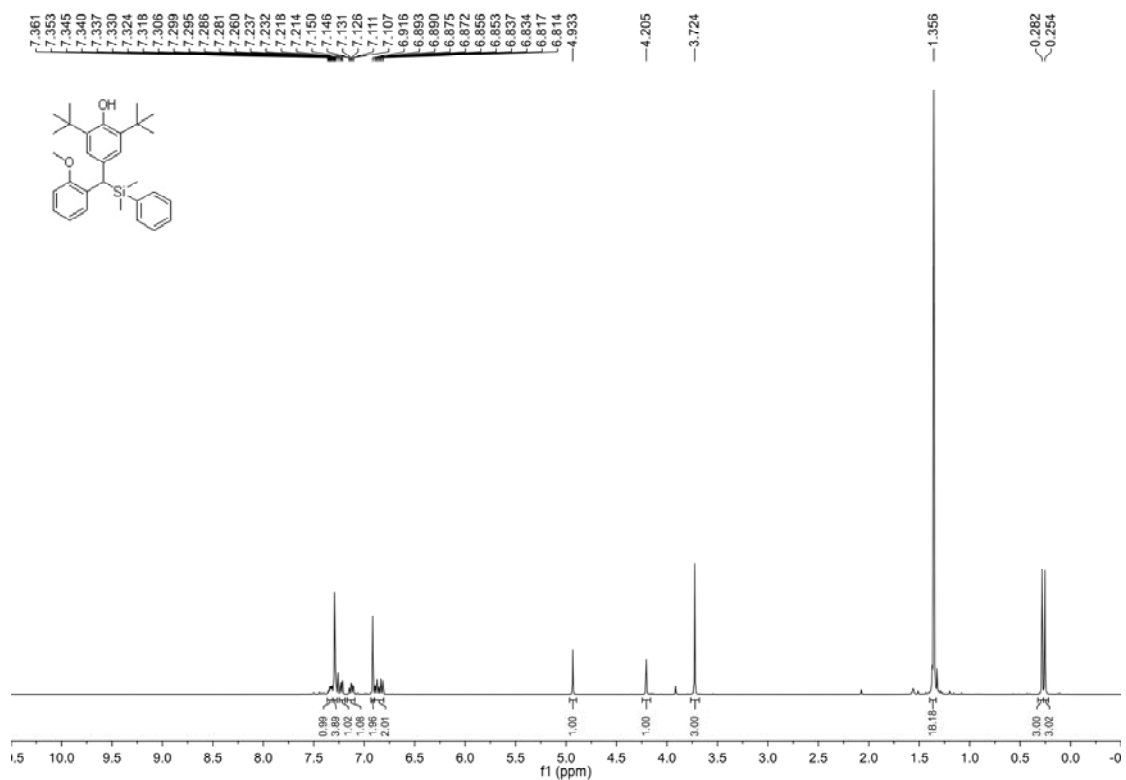


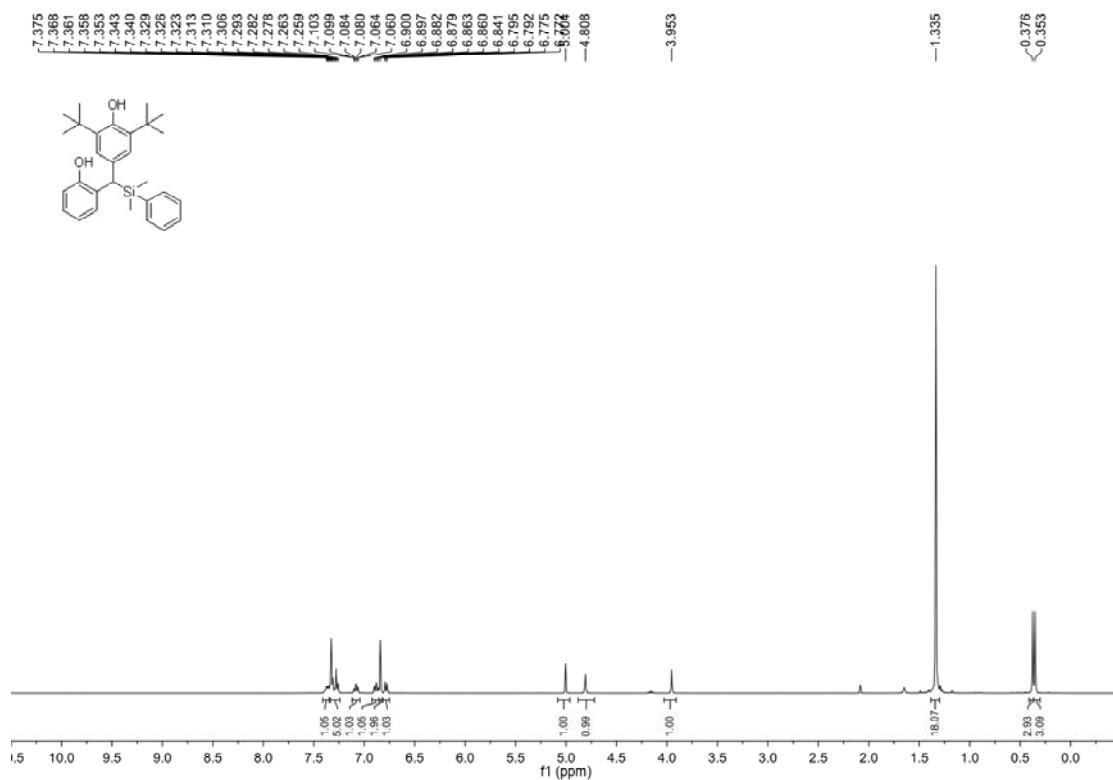


^1H NMR (400 MHz, CDCl_3) of compound **7b**

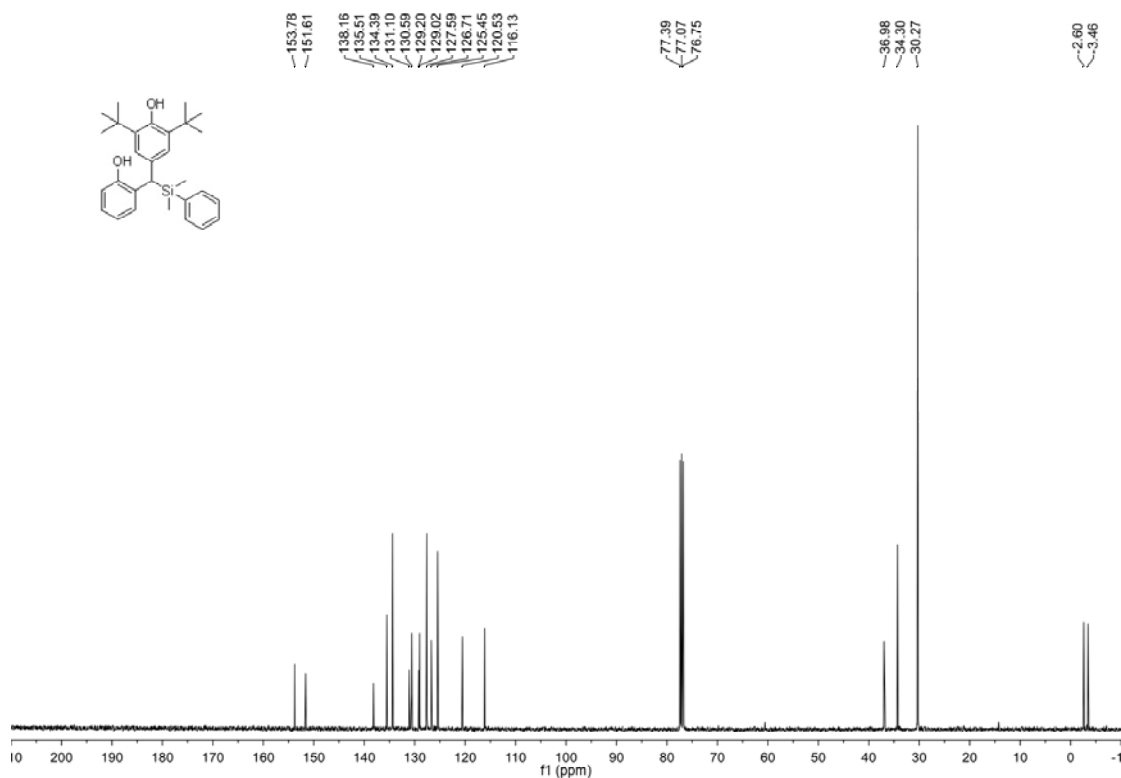


$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **7b**

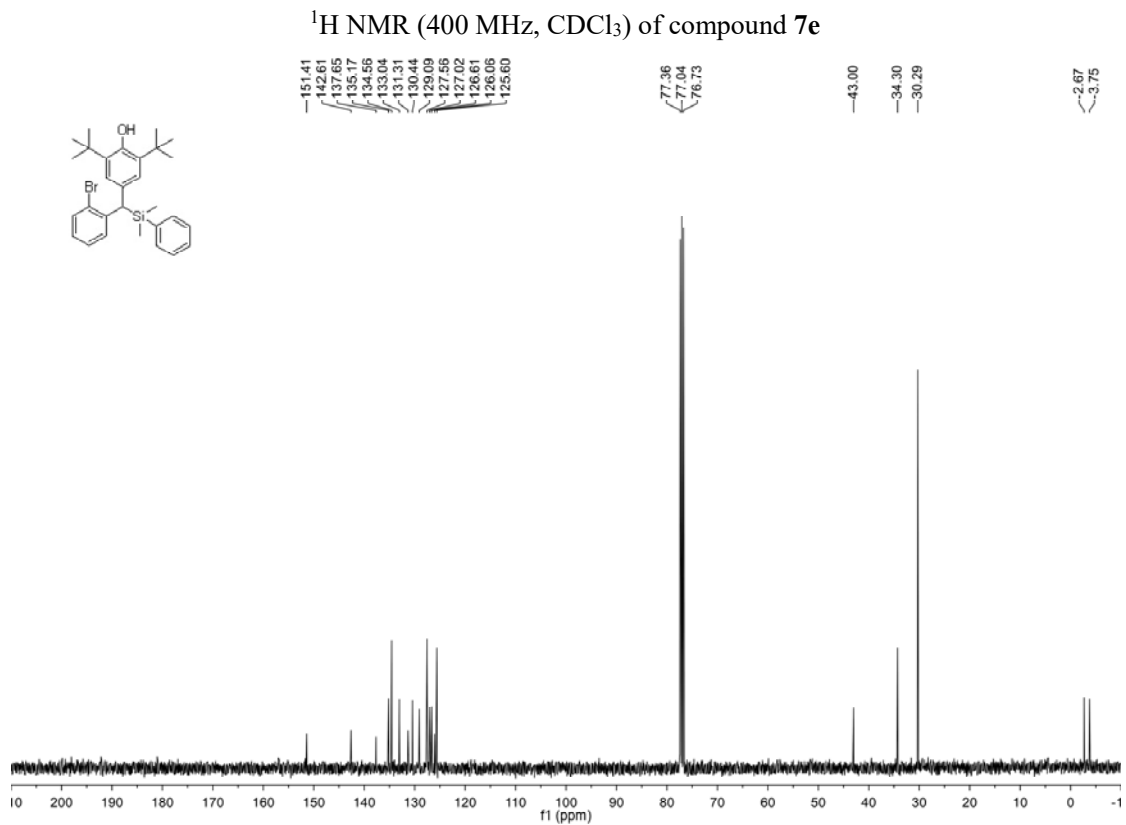
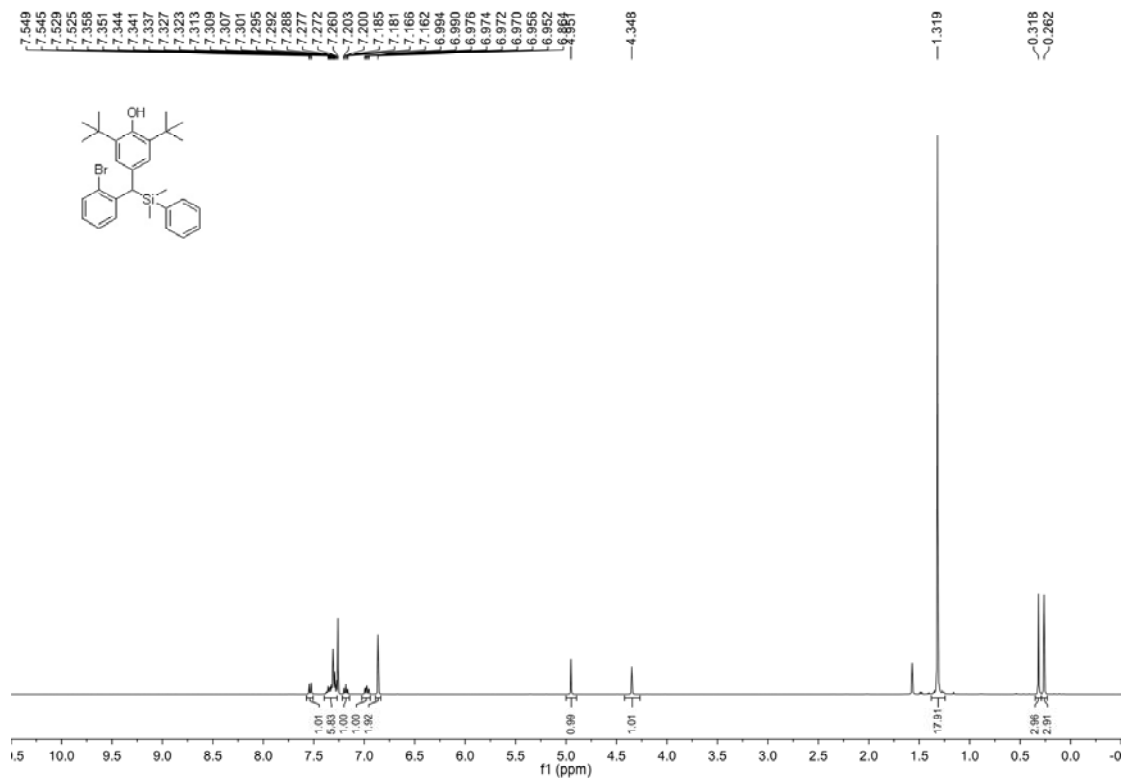


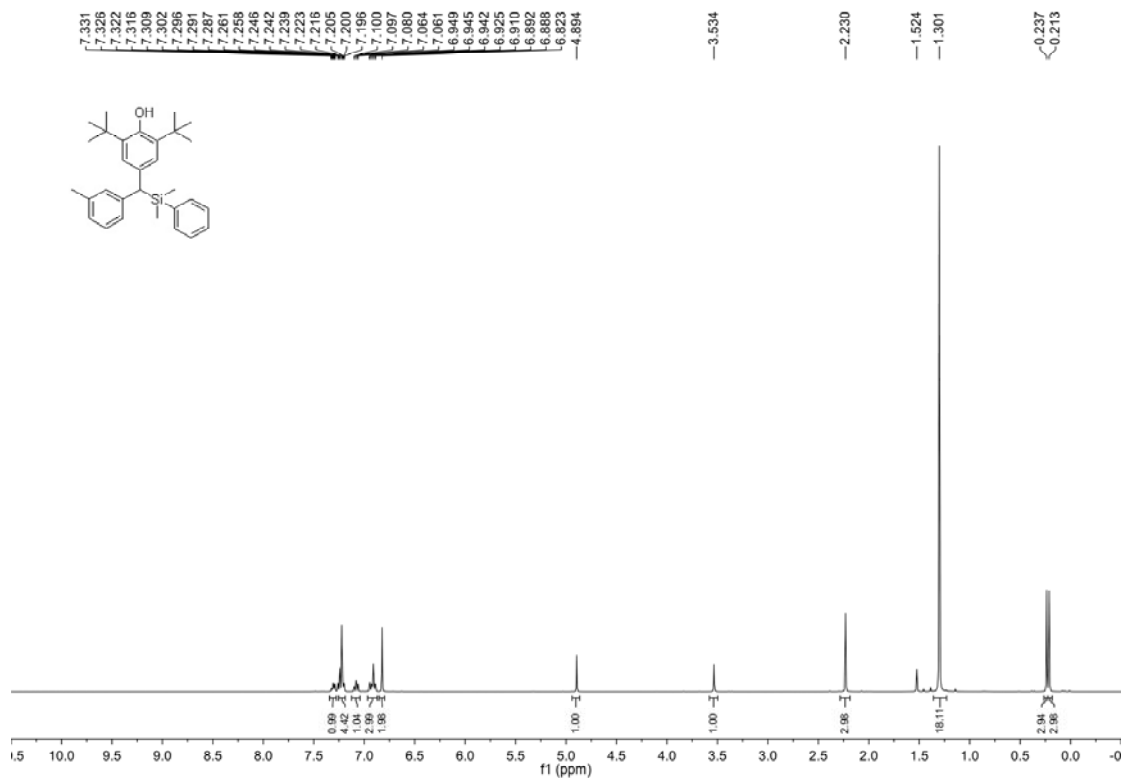


^1H NMR (400 MHz, CDCl_3) of compound **7d**

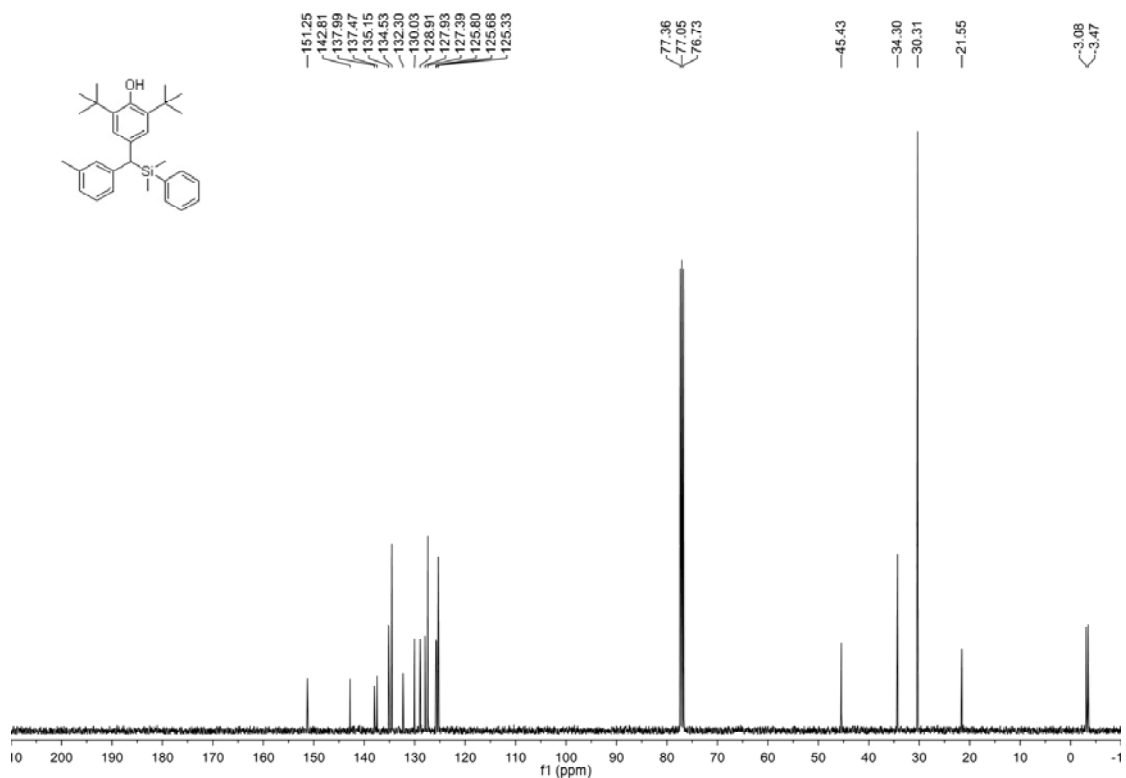


$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **7d**

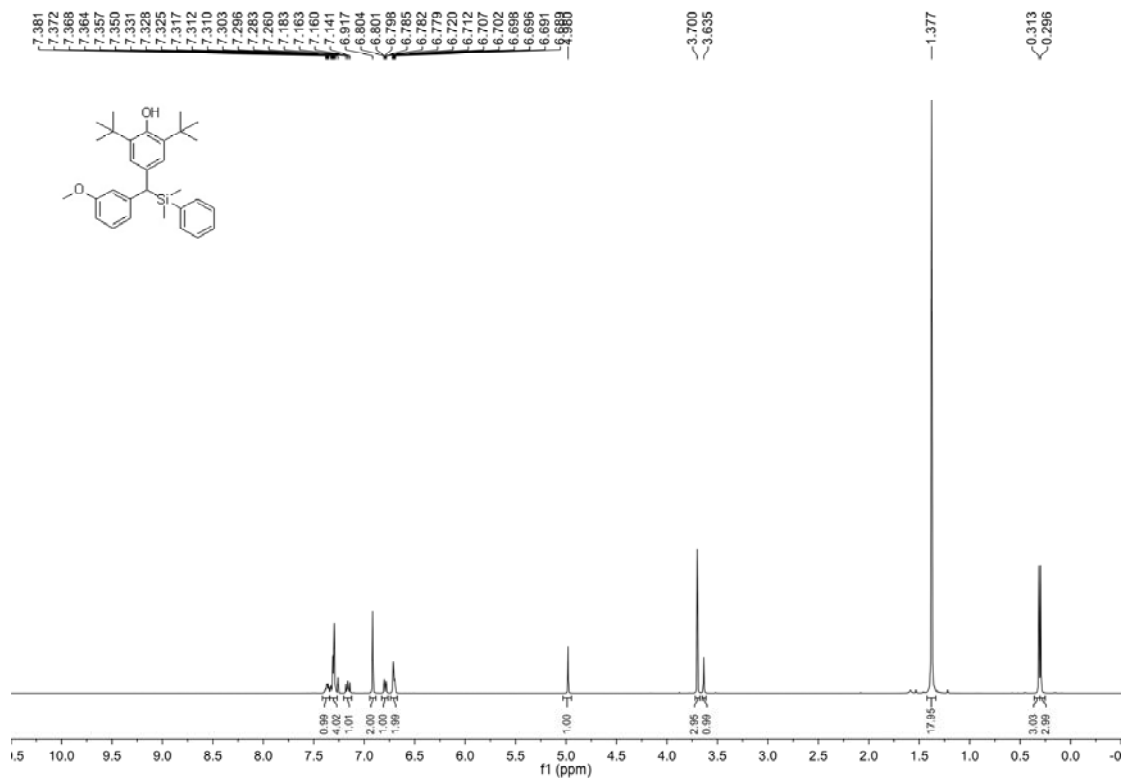




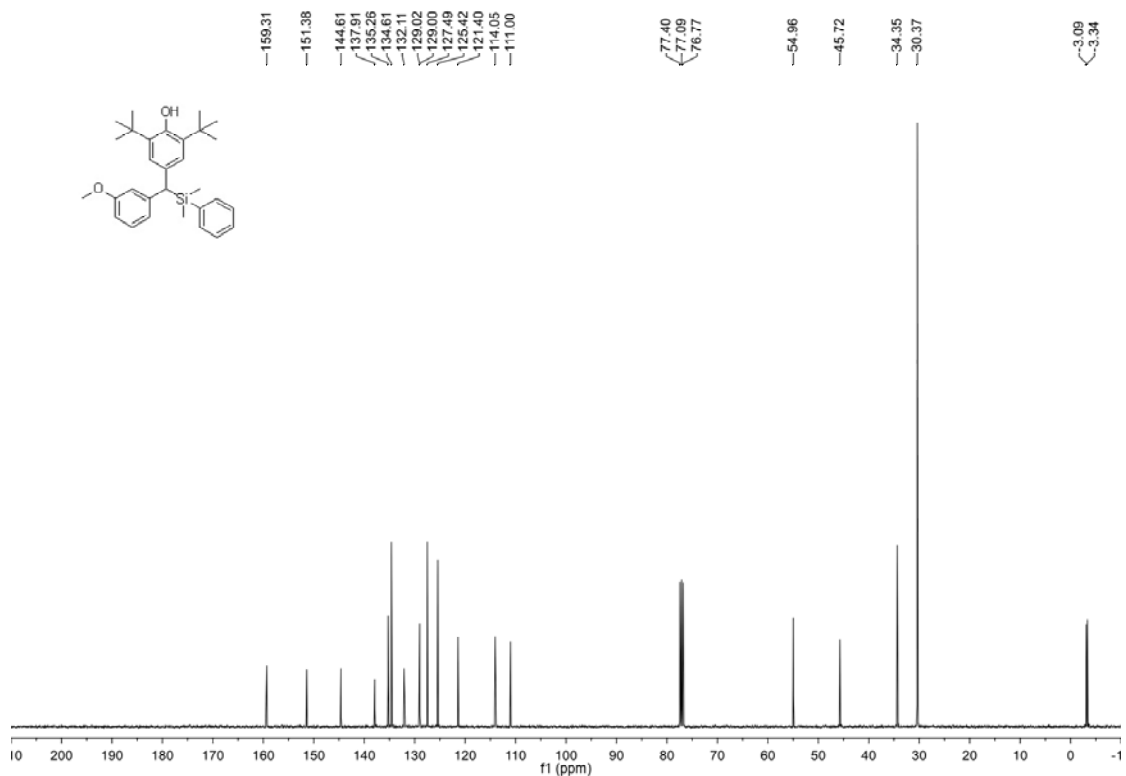
^1H NMR (400 MHz, CDCl_3) of compound **7f**



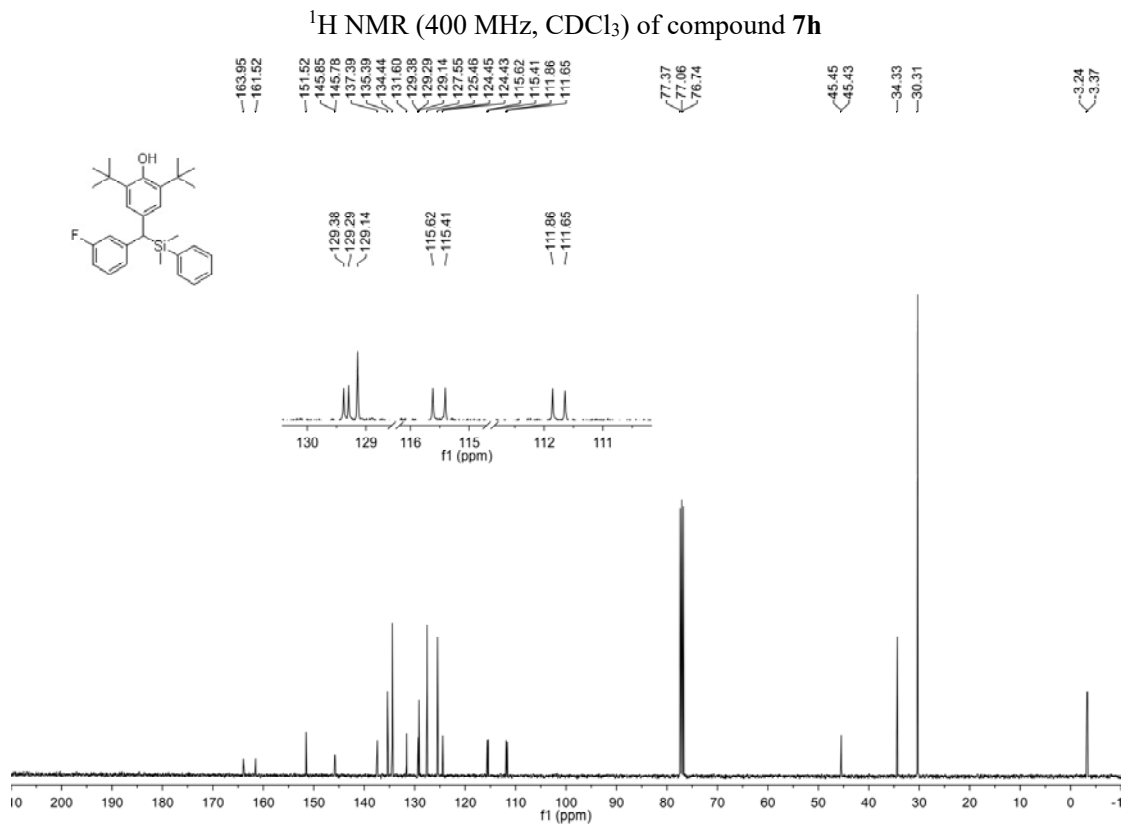
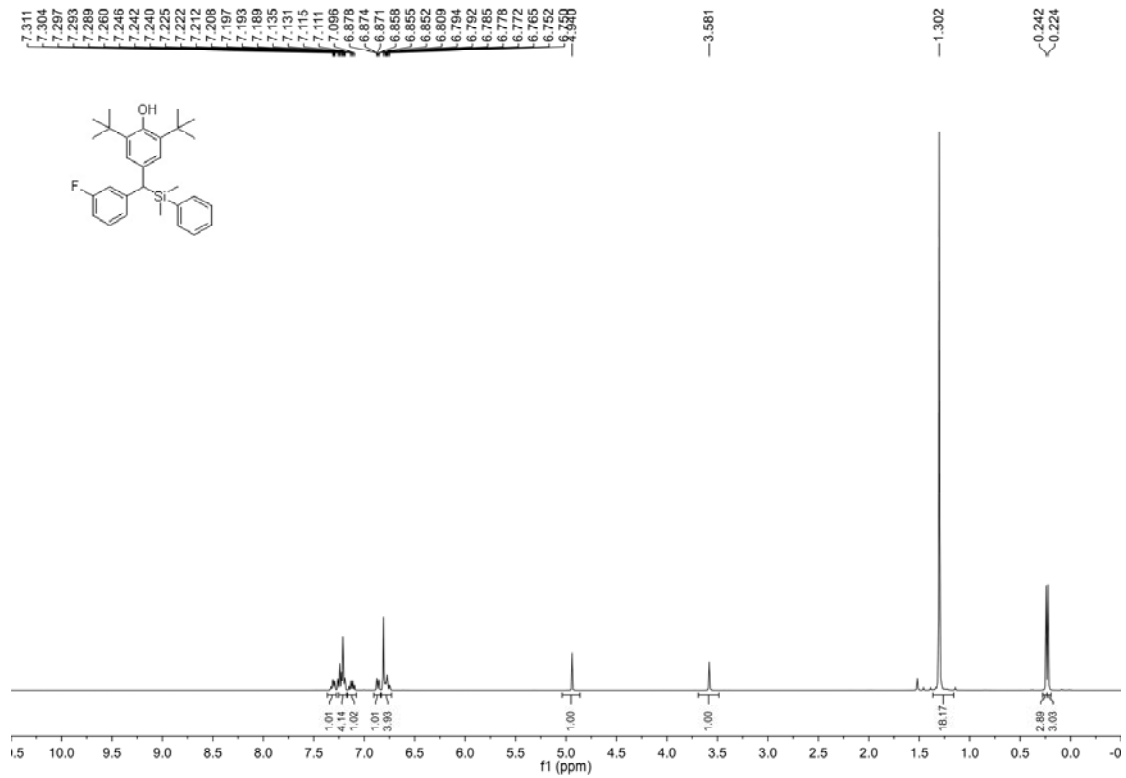
^{13}C NMR (100 MHz, CDCl_3) of compound **7f**

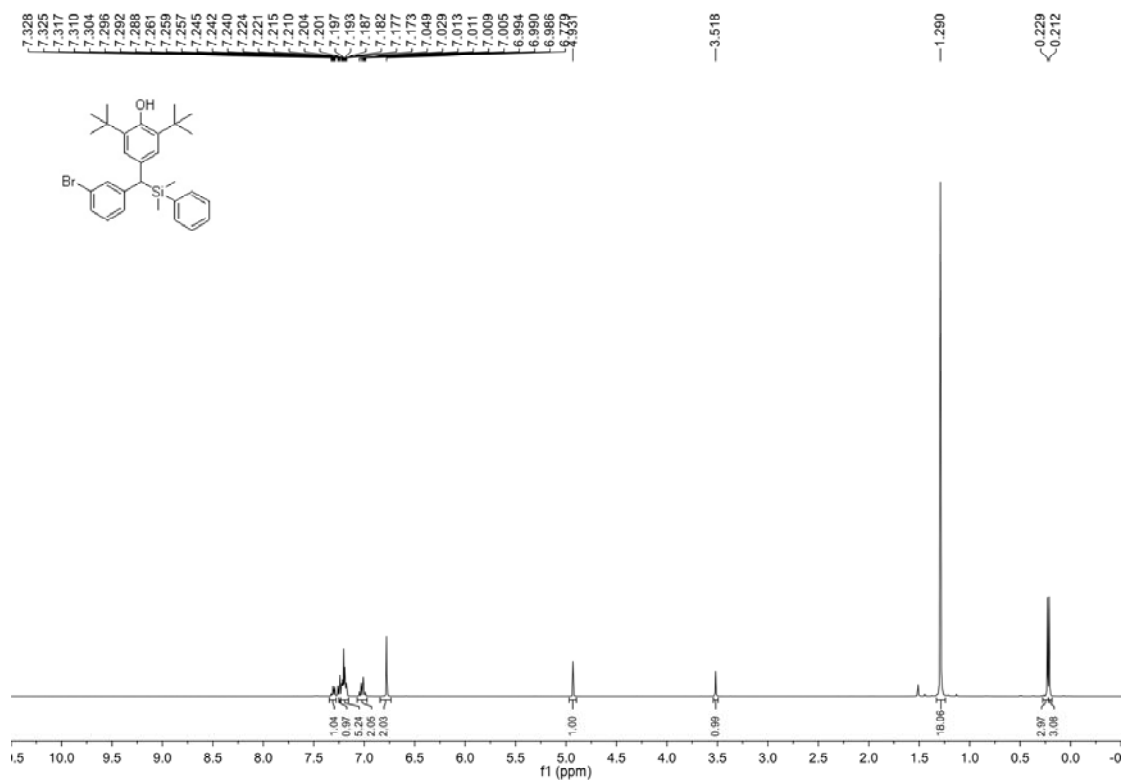


¹H NMR (400 MHz, CDCl₃) of compound 7g

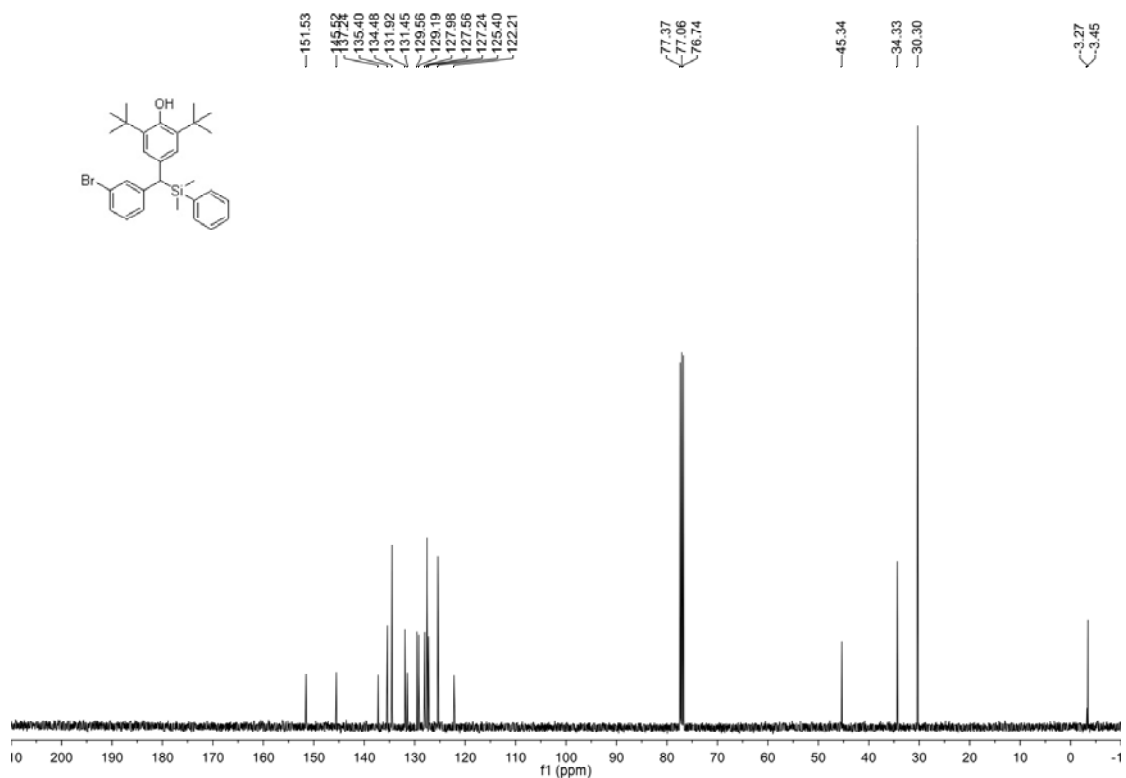


¹³C {¹H} NMR (100 MHz, CDCl₃) of compound 7g

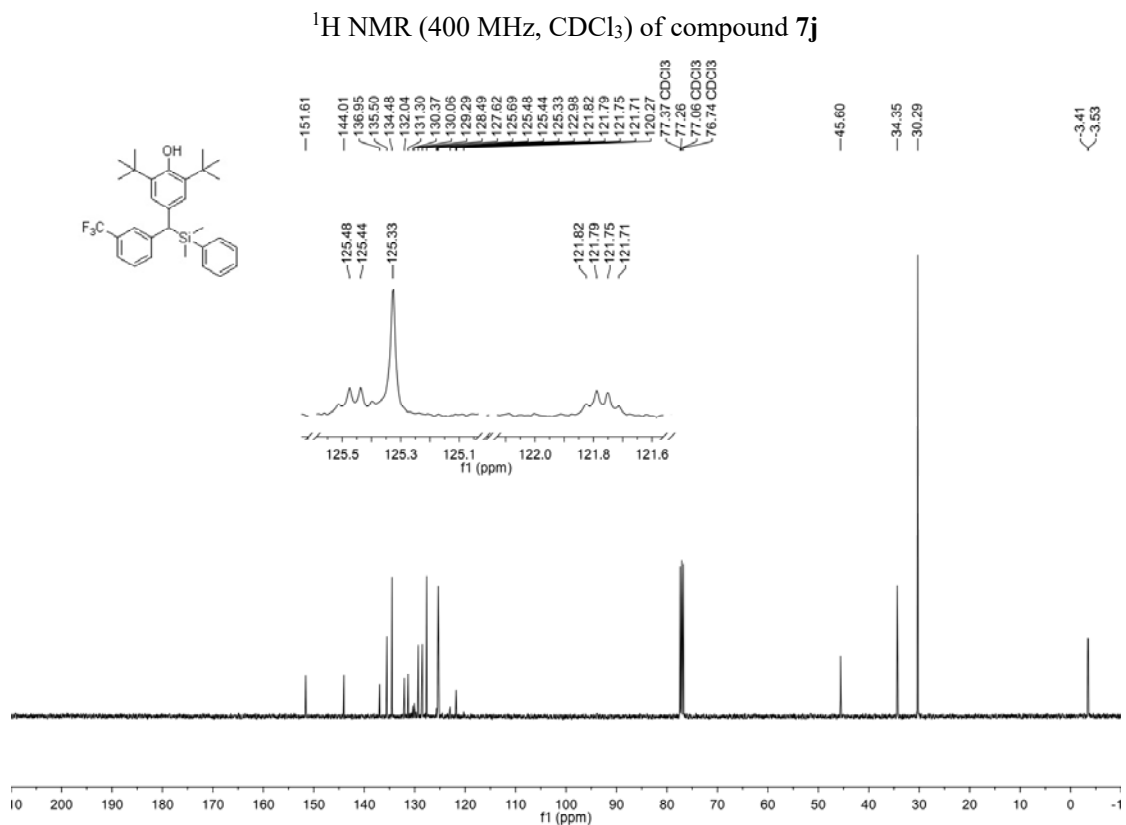
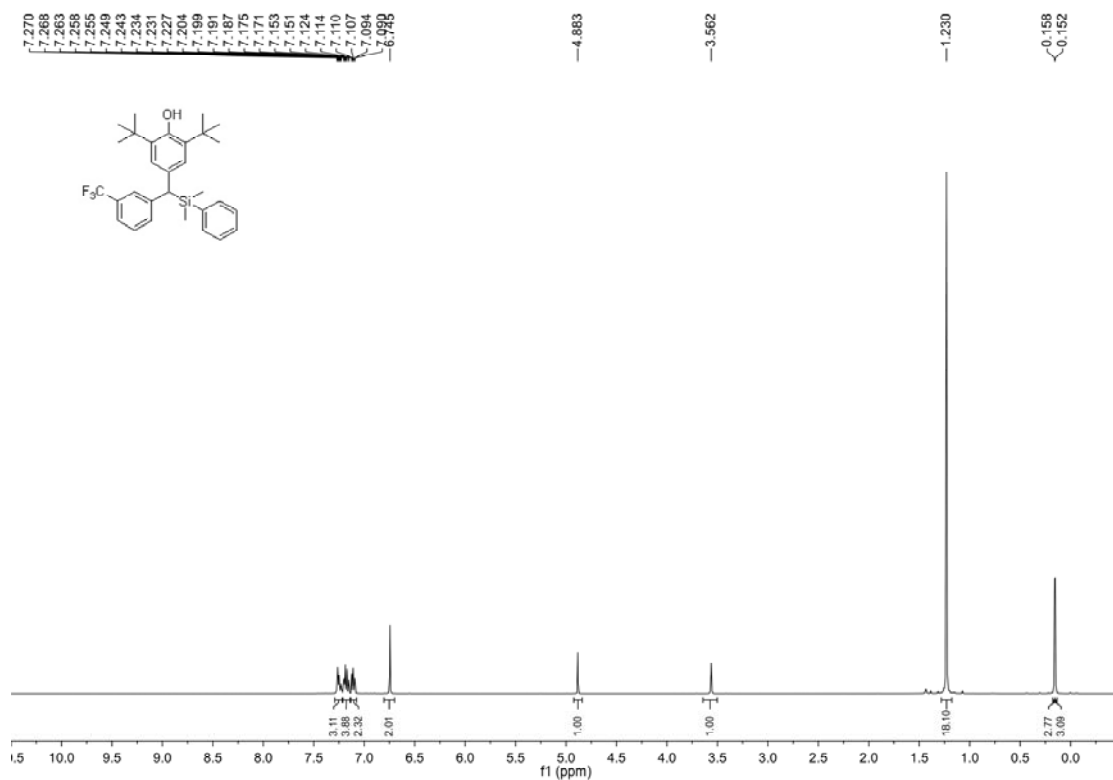




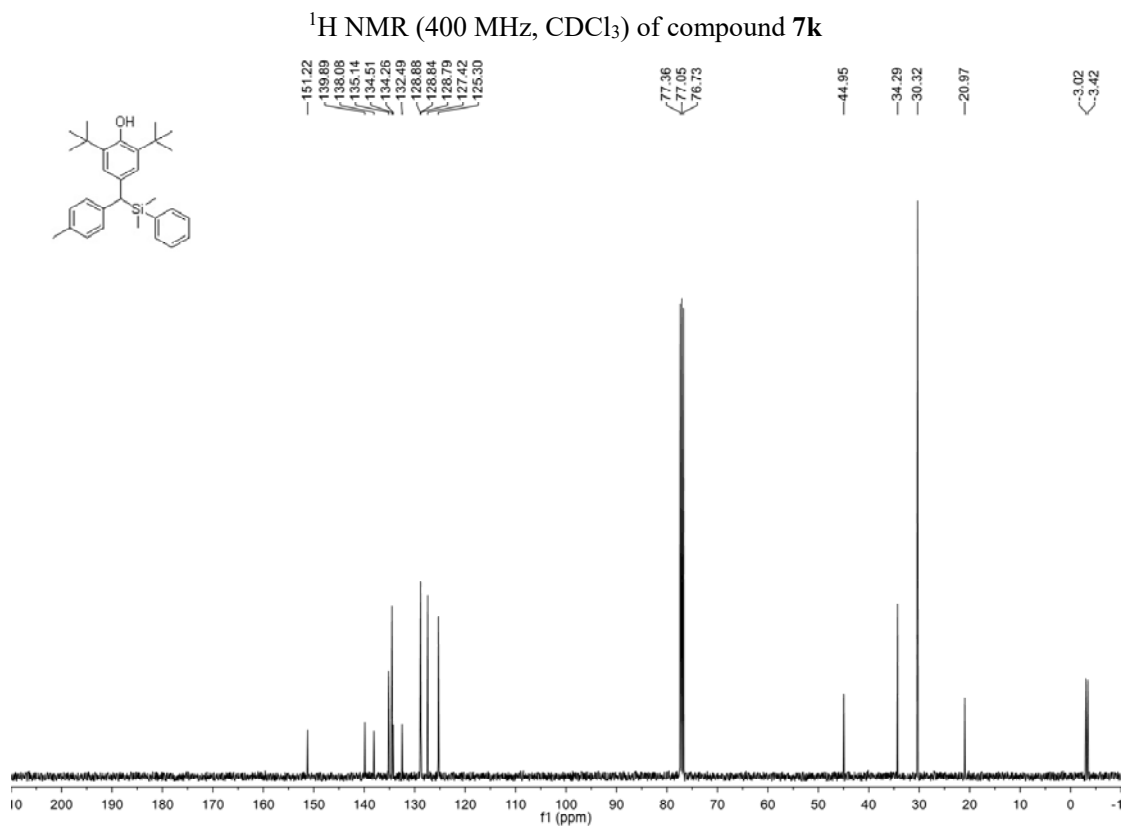
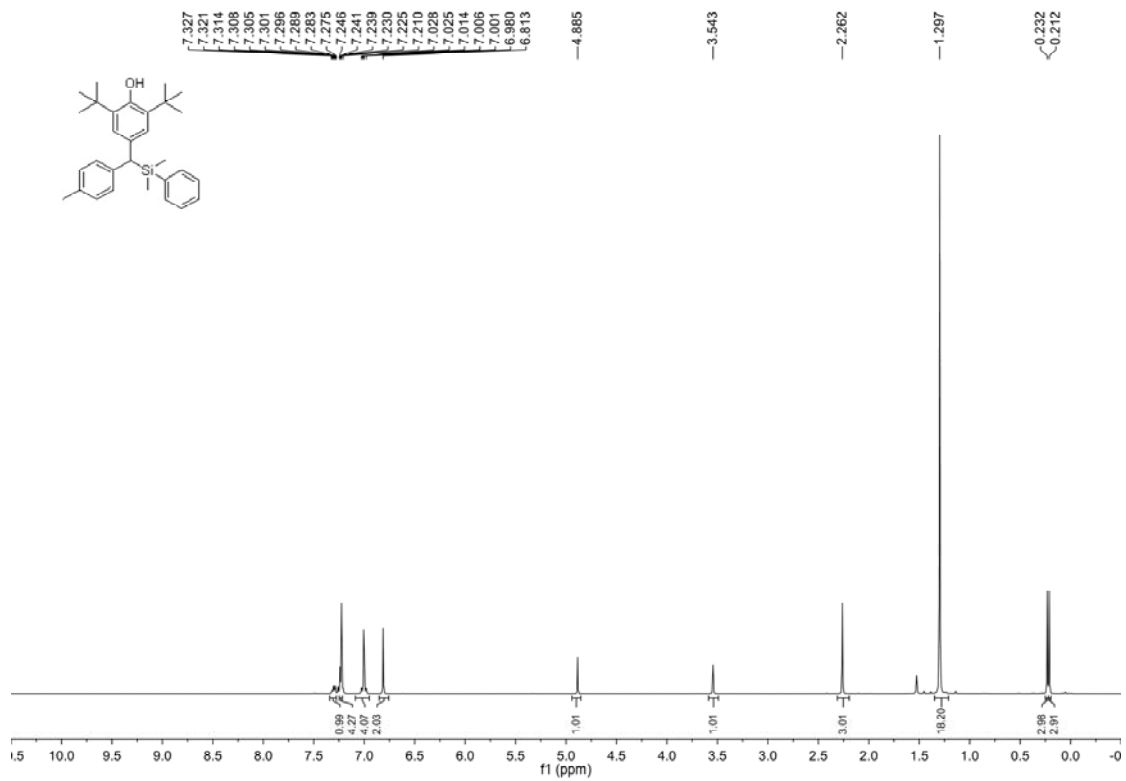
¹H NMR (400 MHz, CDCl₃) of compound **7i**

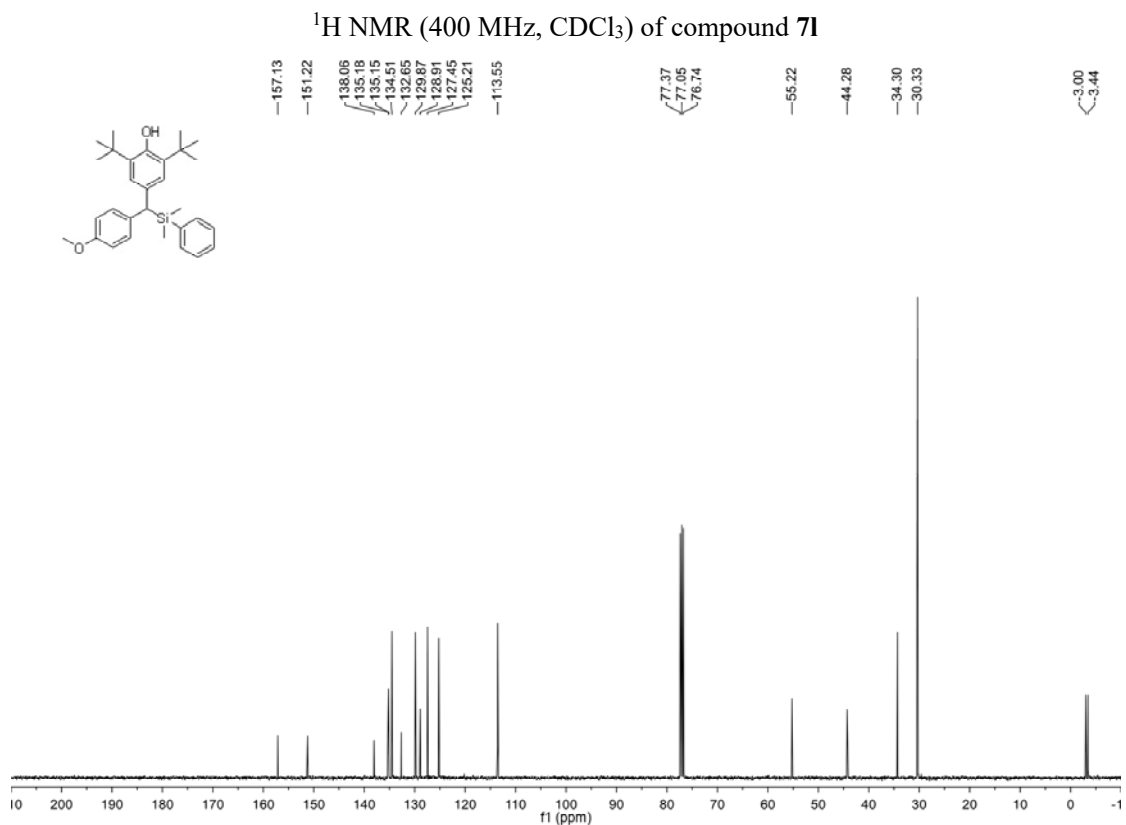
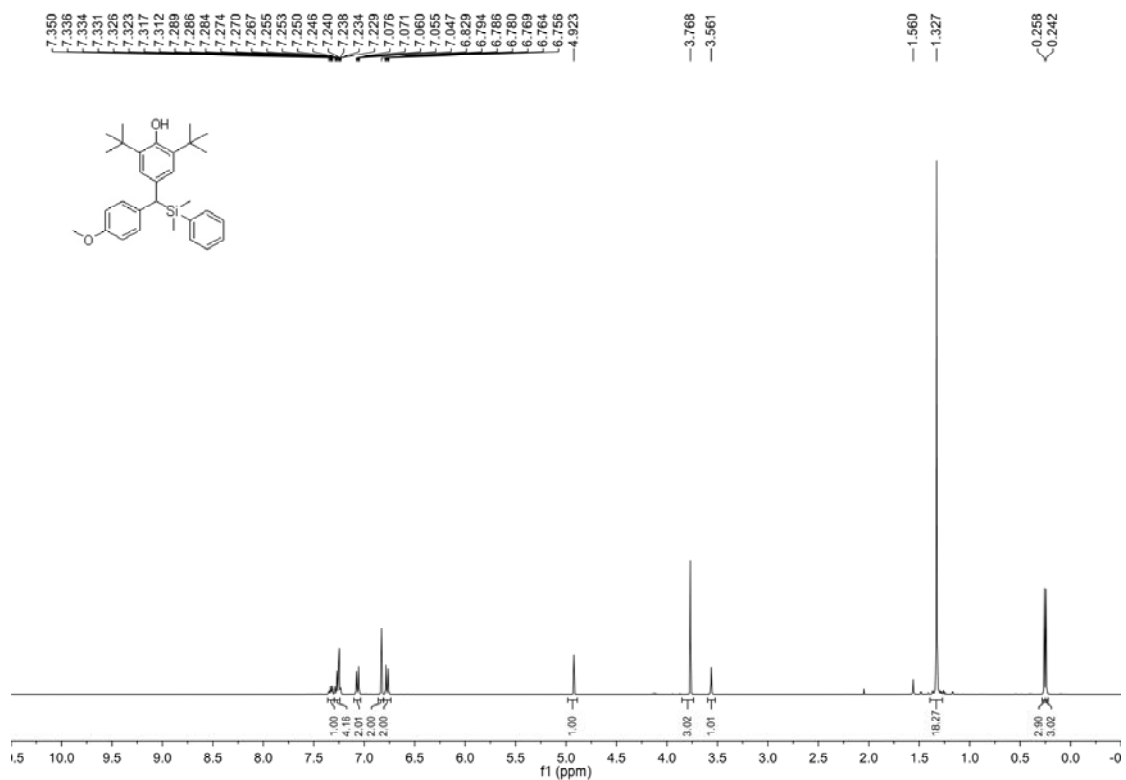


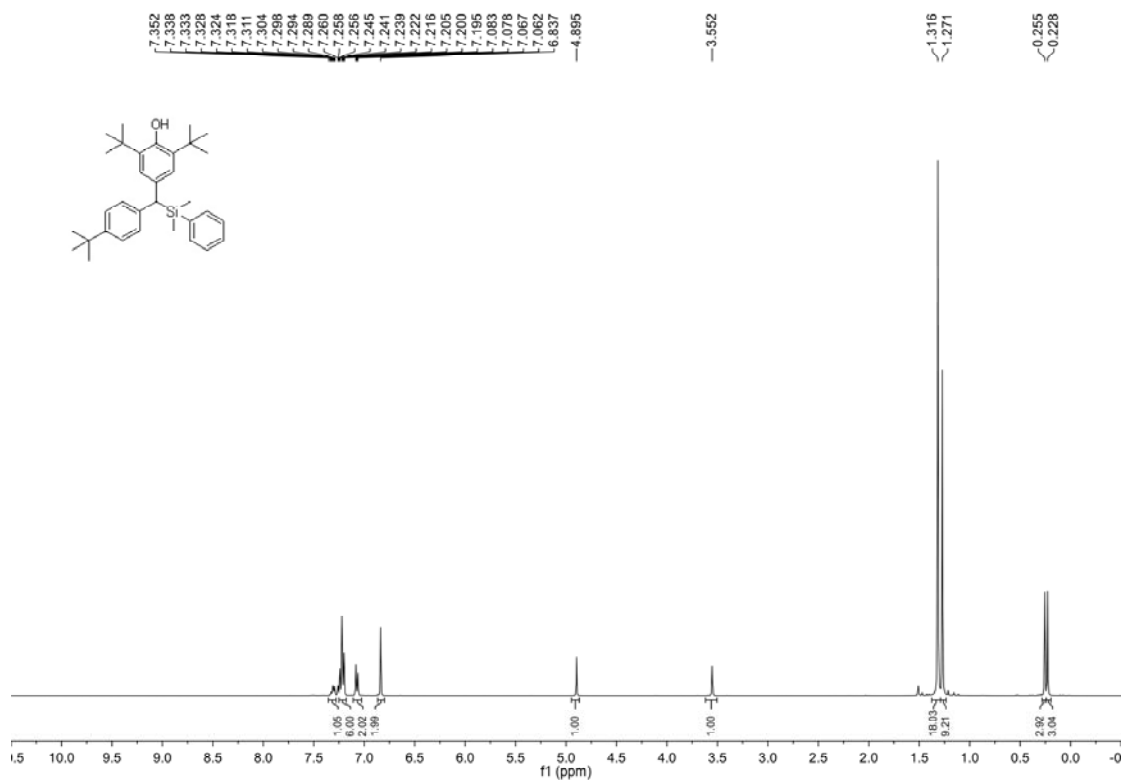
¹³C{¹H} NMR (100 MHz, CDCl₃) of compound **7i**



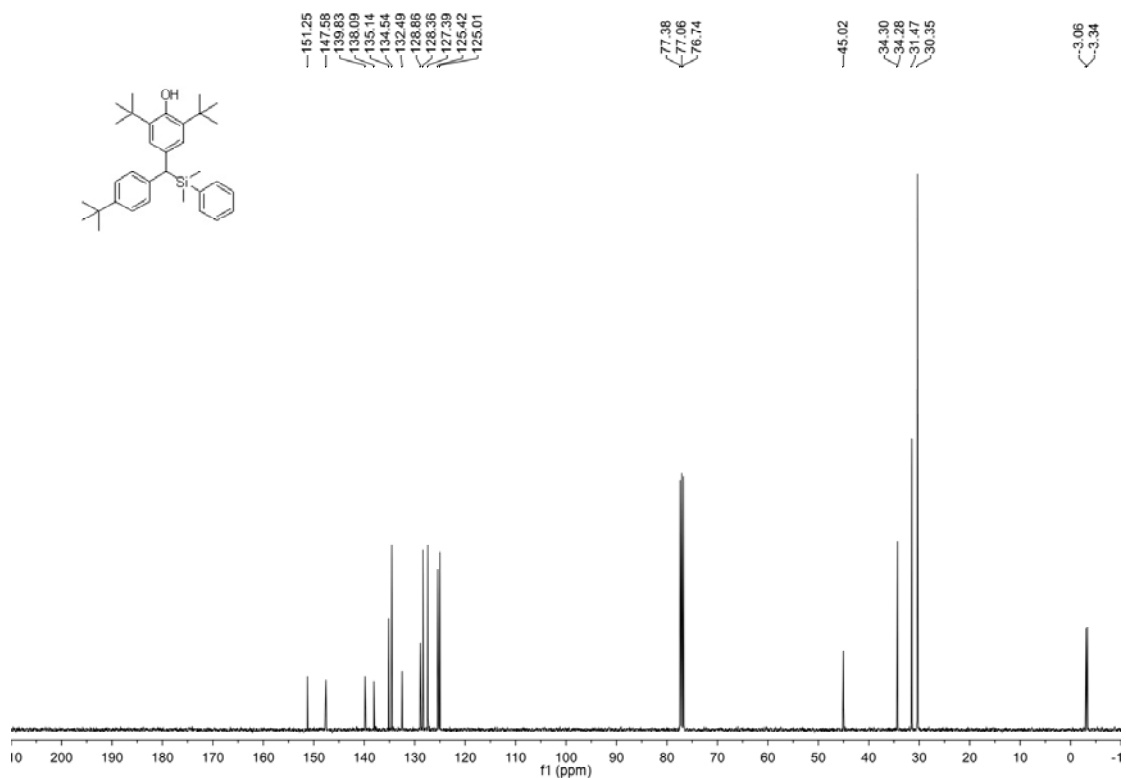
¹³C {¹H} NMR (100 MHz, CDCl₃) of compound **7j**



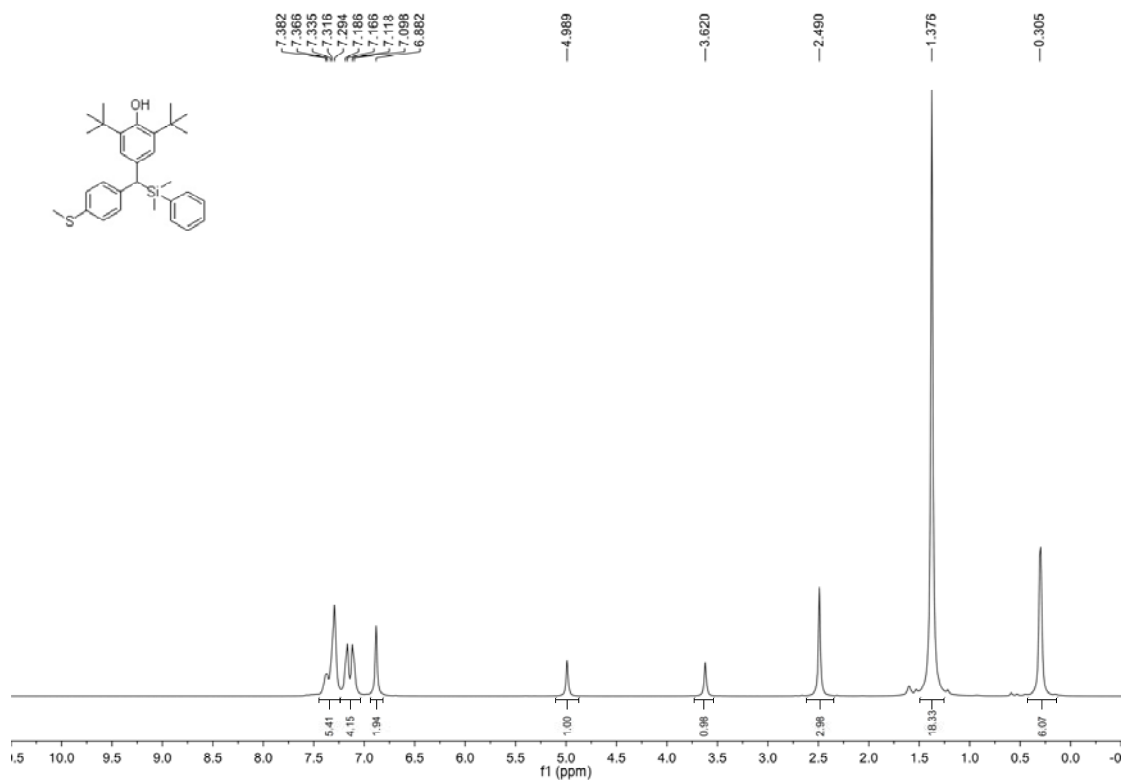




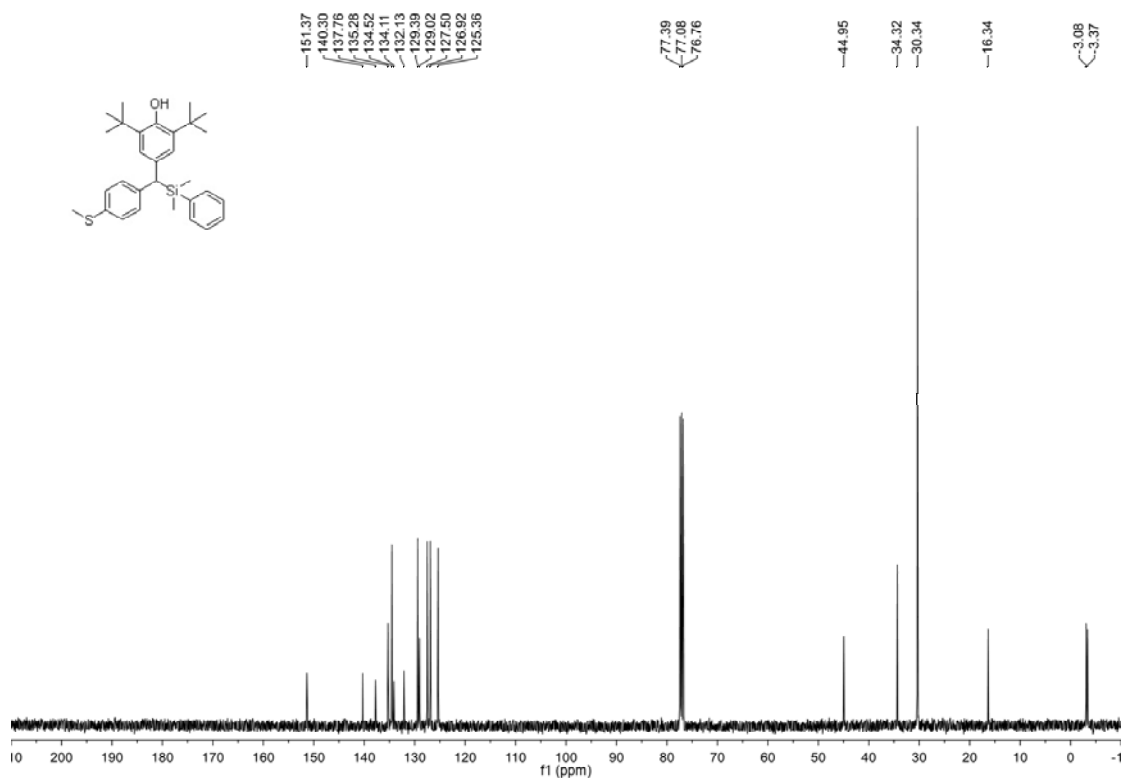
¹H NMR (400 MHz, CDCl₃) of compound **7m**



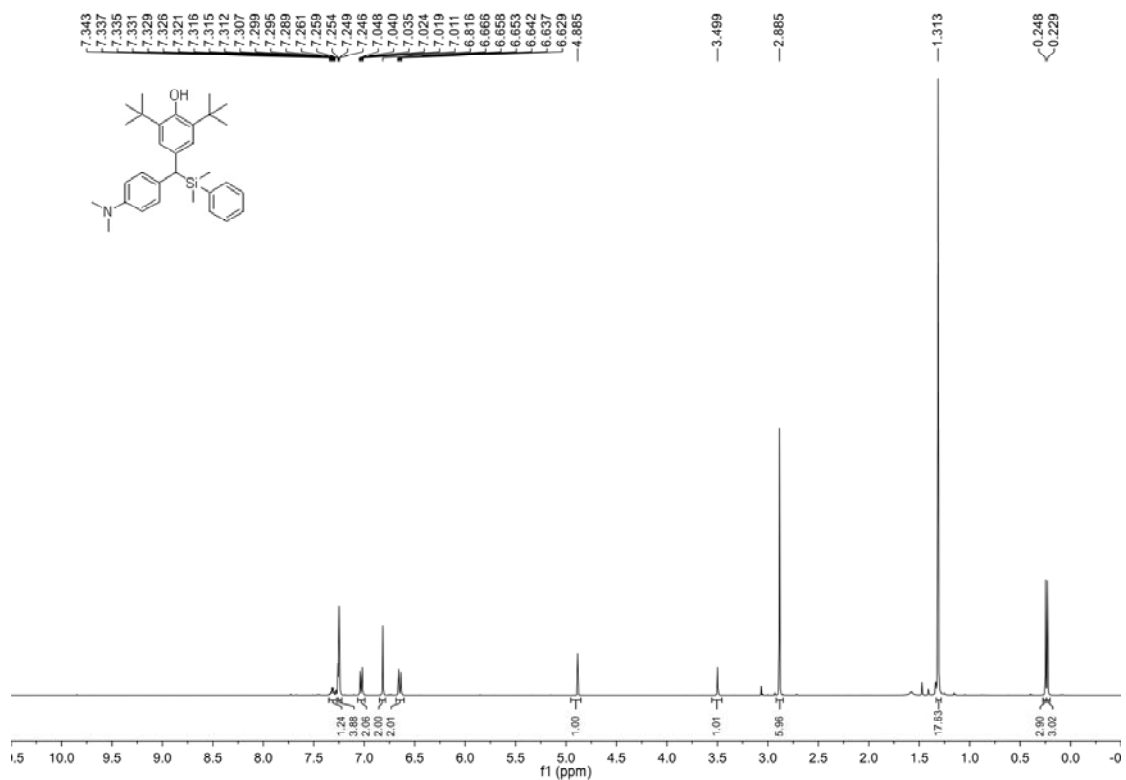
¹³C NMR (100 MHz, CDCl₃) of compound **7m**



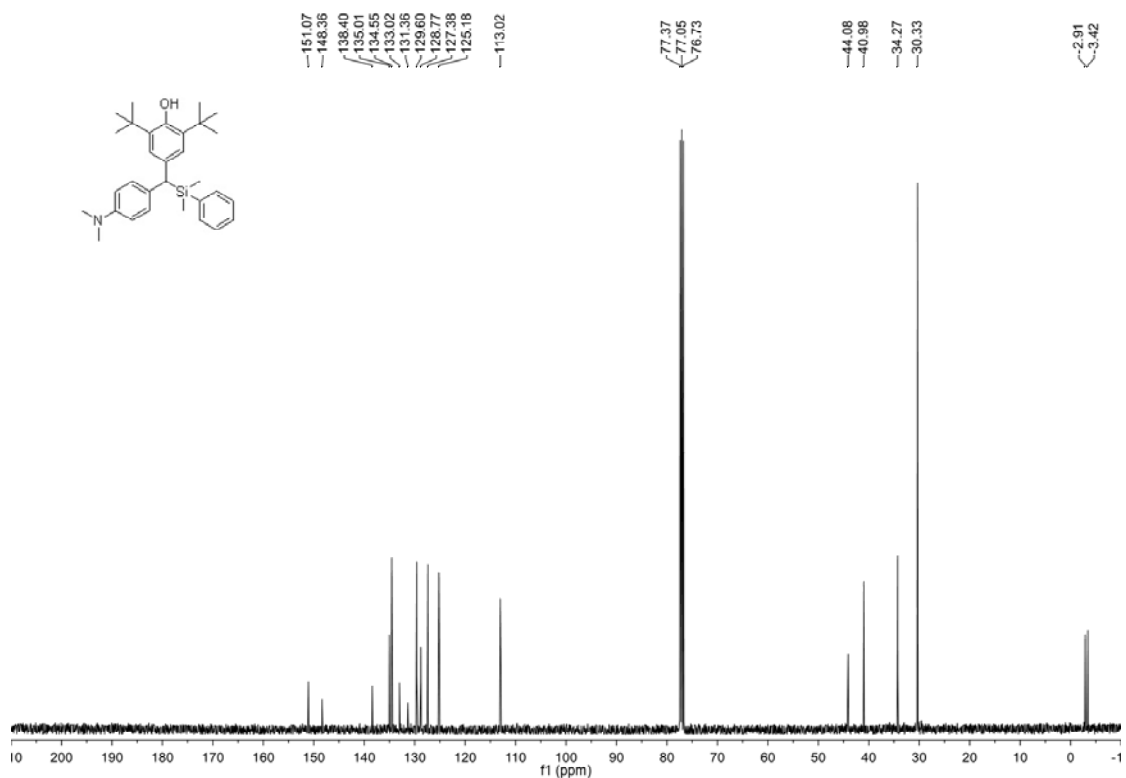
^1H NMR (400 MHz, CDCl_3) of compound **7n**



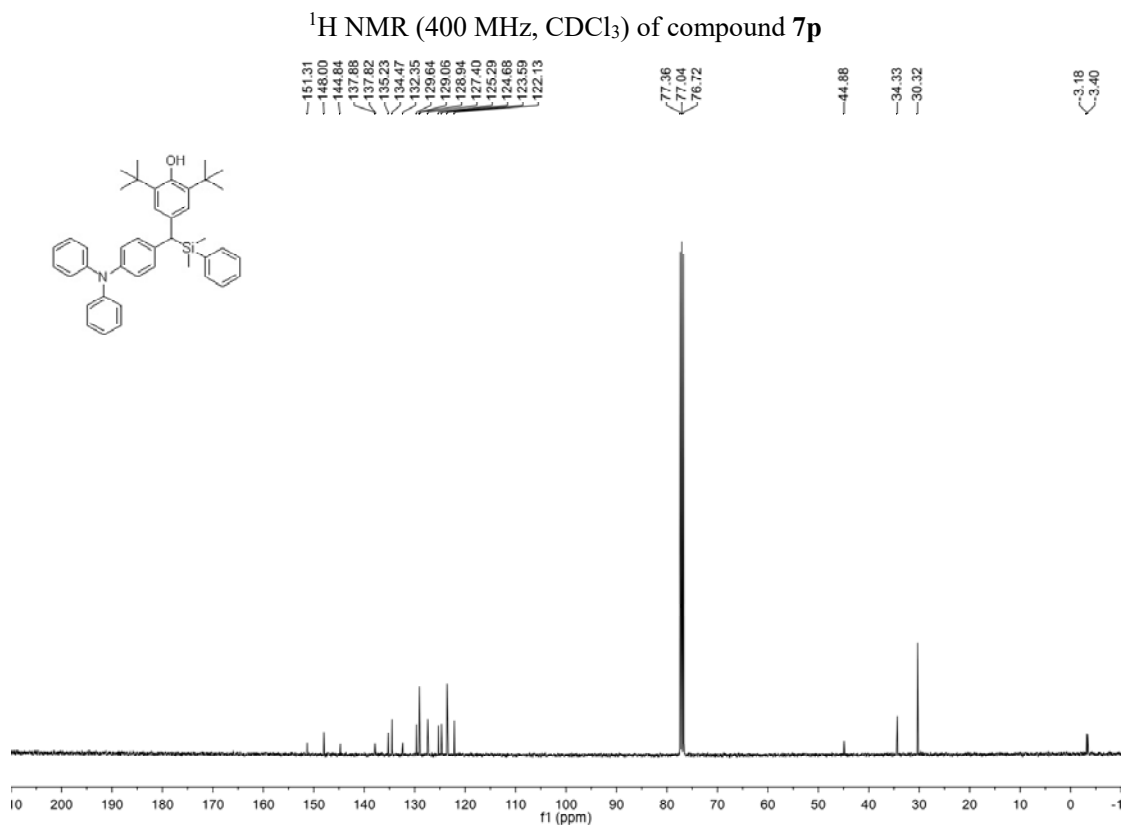
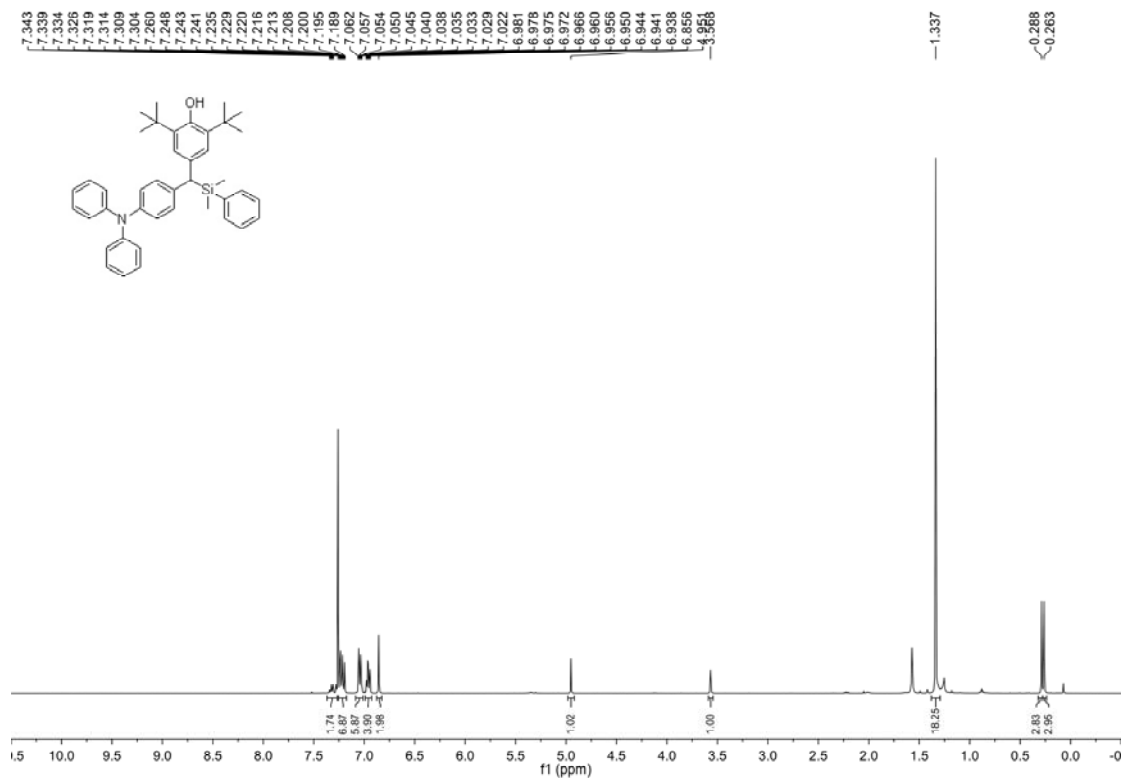
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **7n**

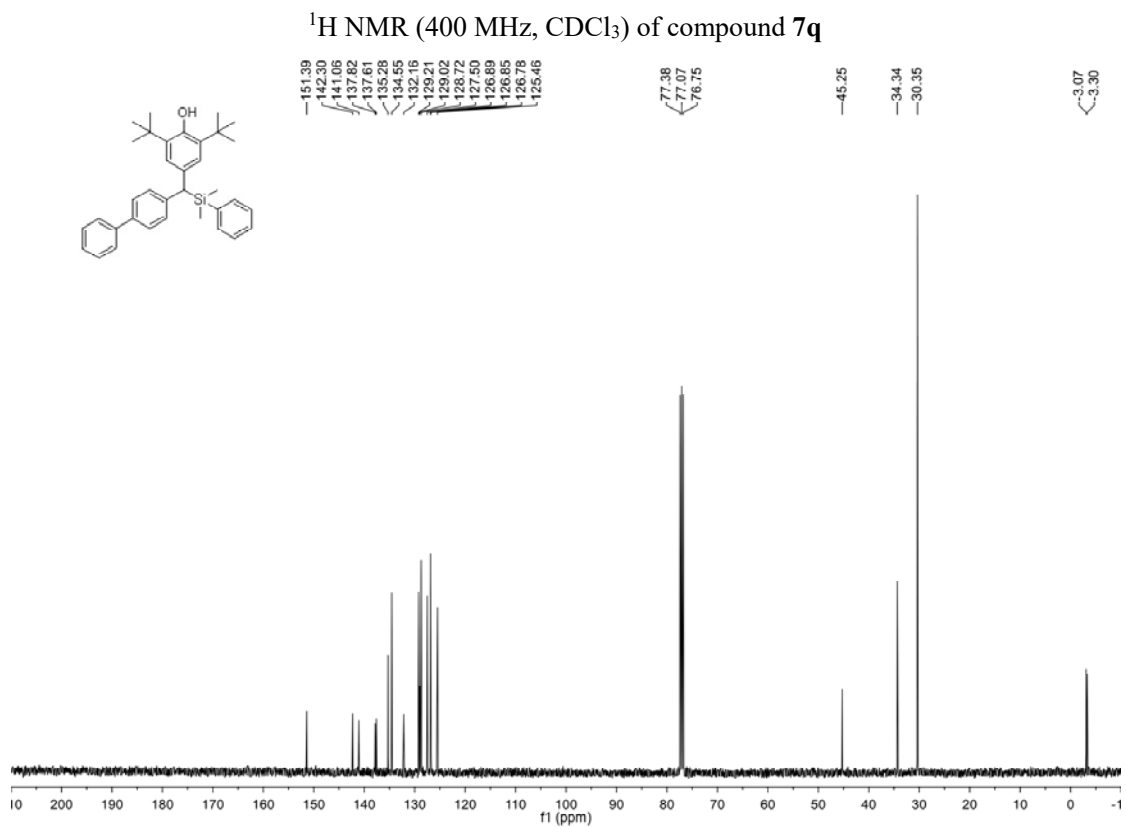
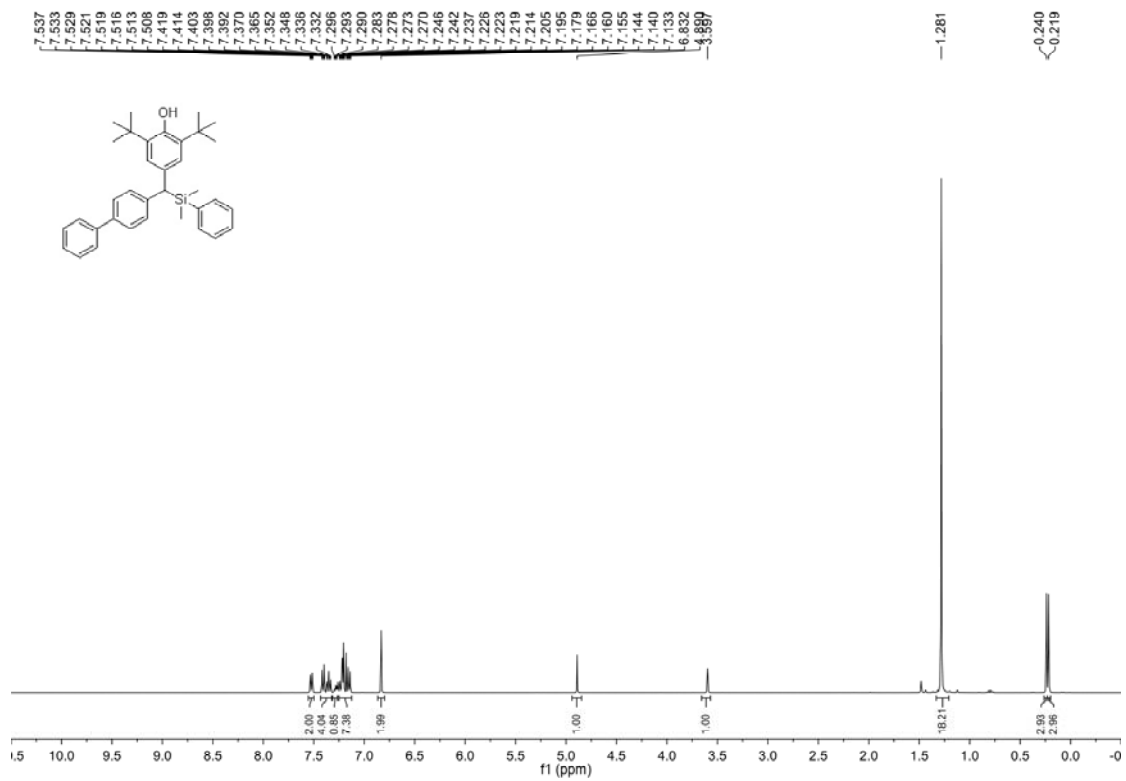


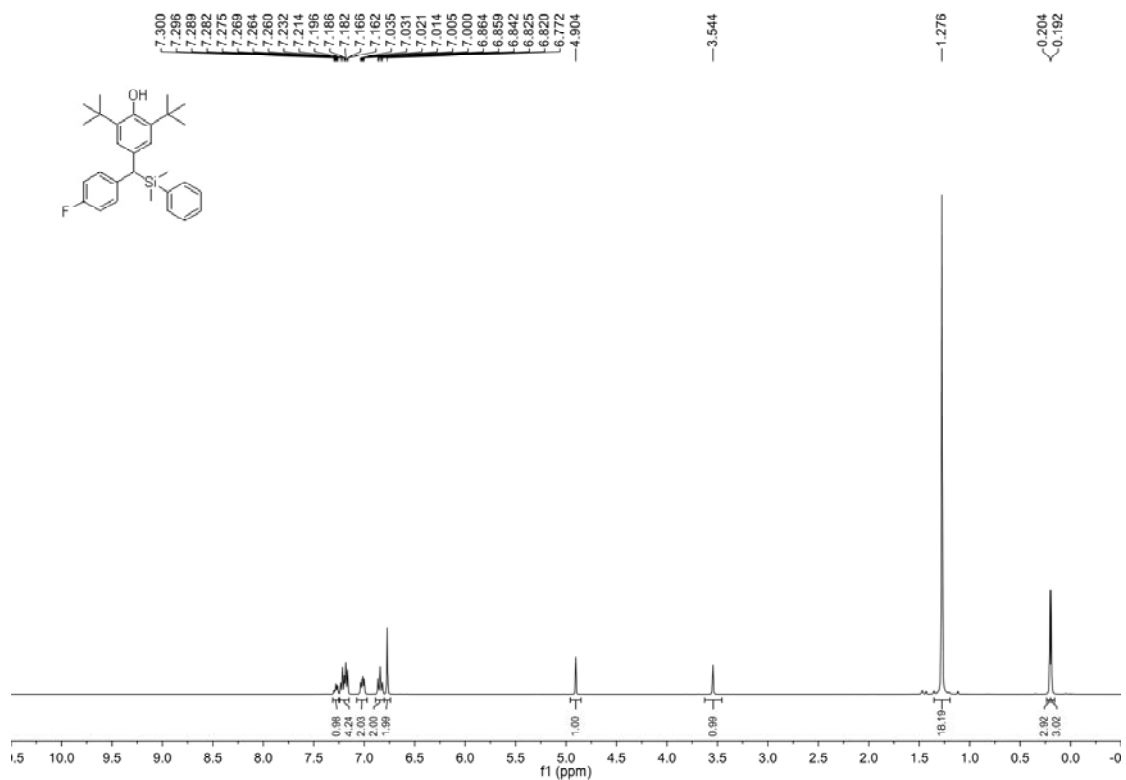
¹H NMR (400 MHz, CDCl₃) of compound **7o**



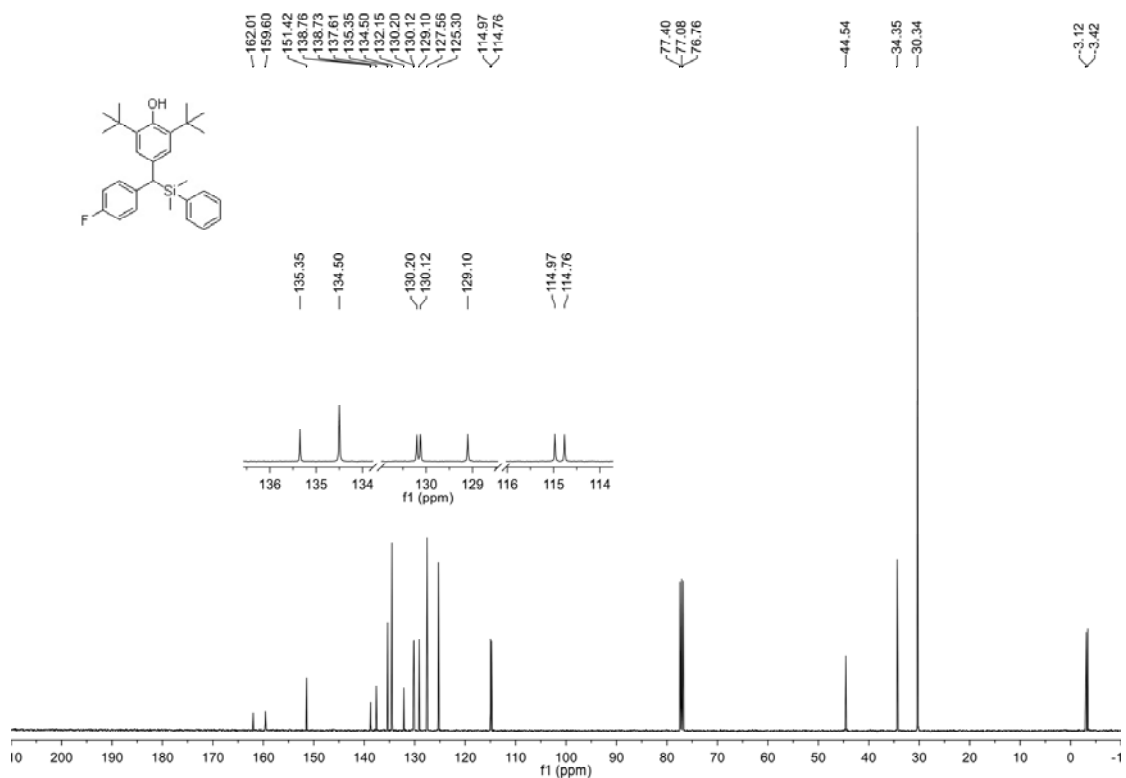
¹³C {¹H} NMR (100 MHz, CDCl₃) of compound **7o**



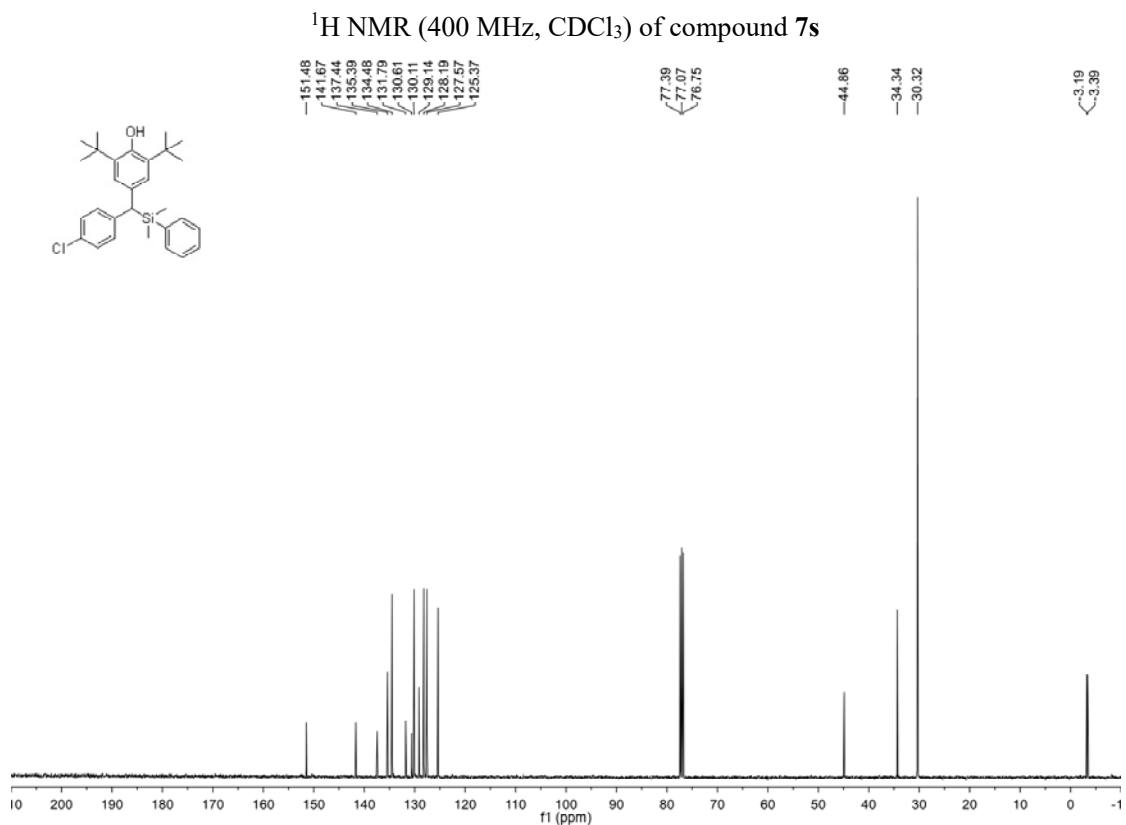
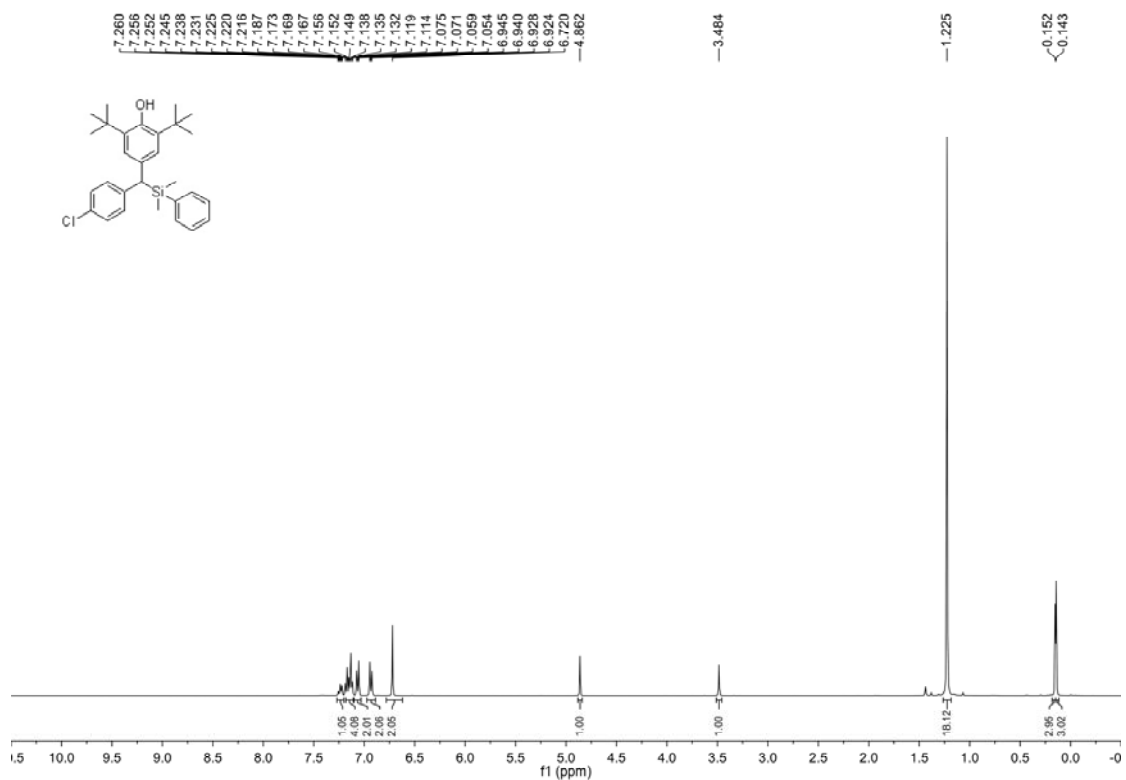




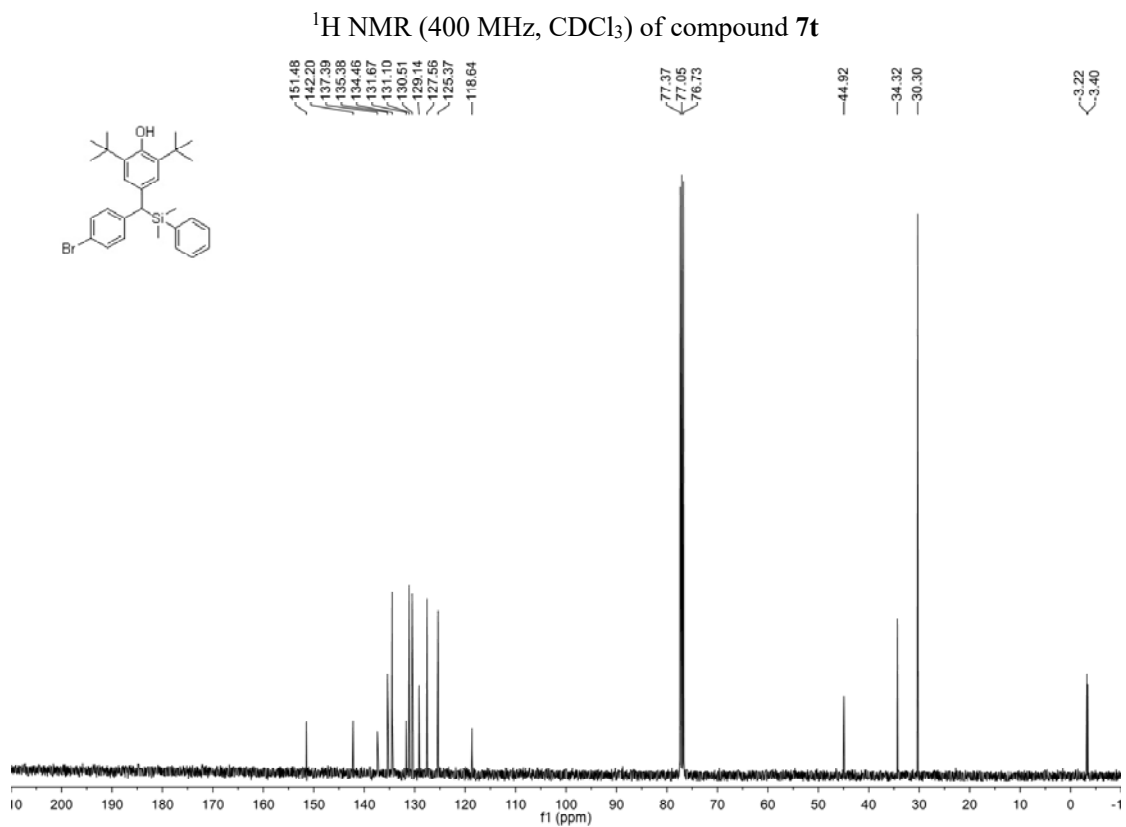
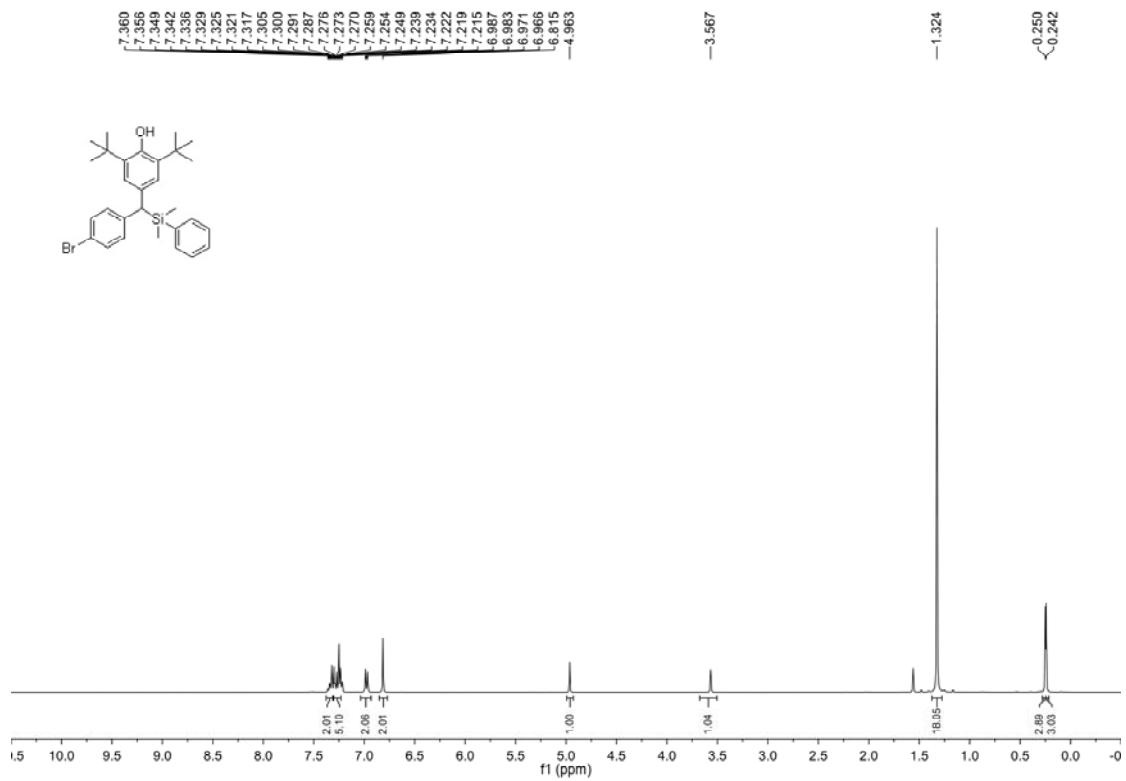
¹H NMR (400 MHz, CDCl₃) of compound **7r**

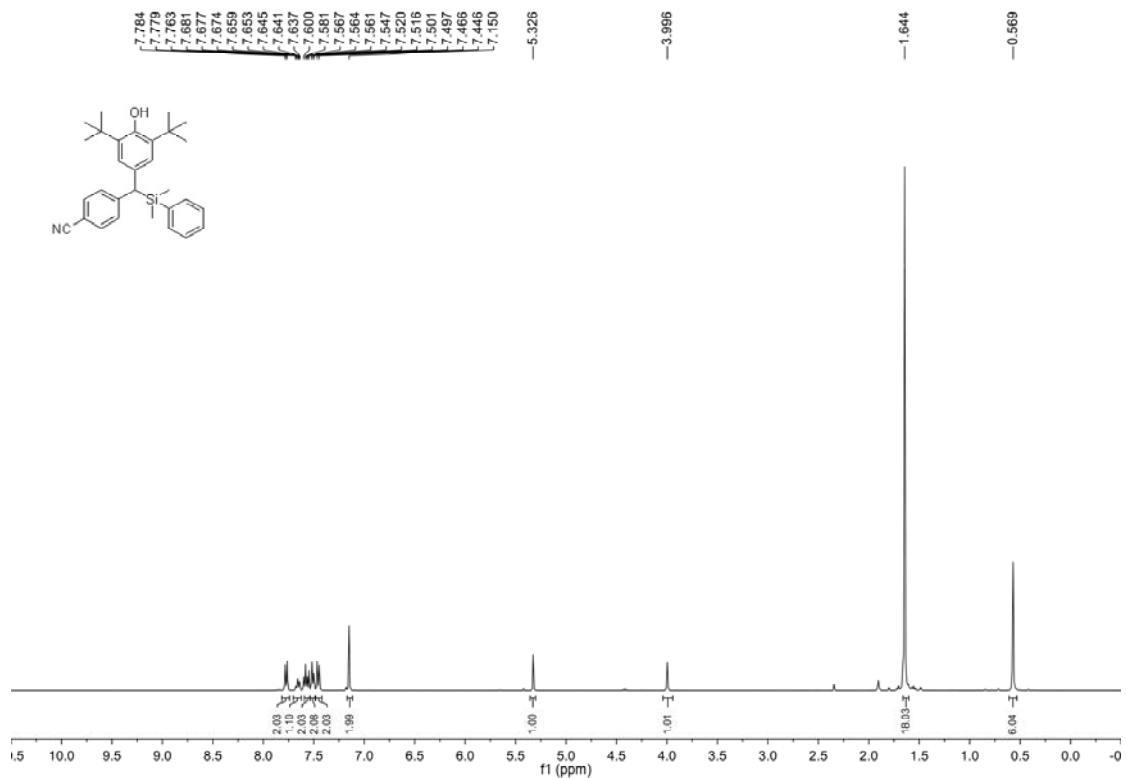


¹³C {¹H} NMR (100 MHz, CDCl₃) of compound **7r**

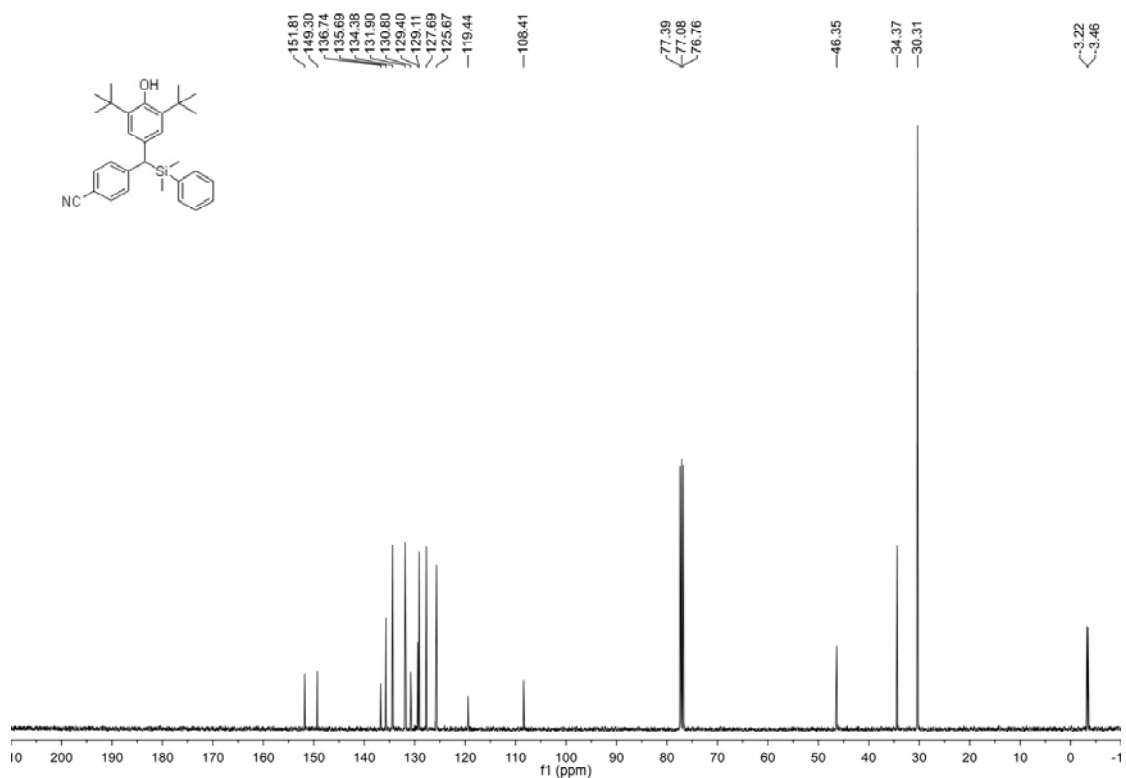


$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **7s**

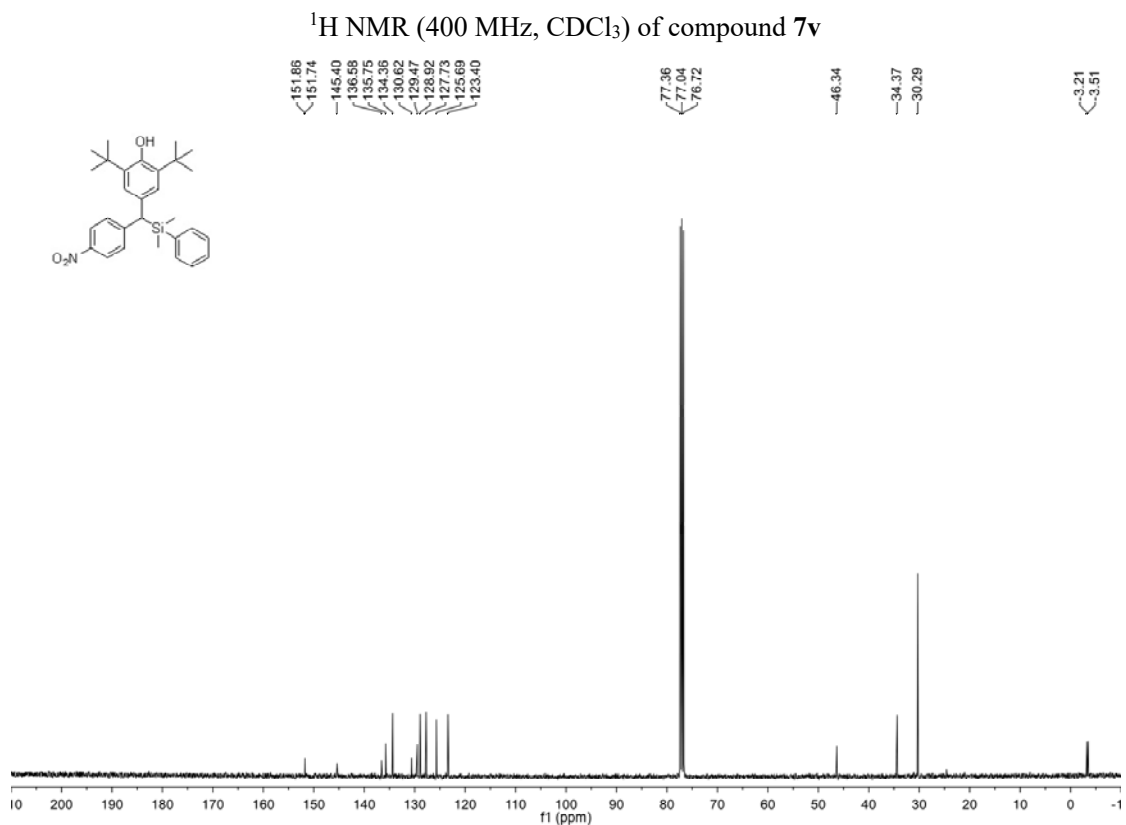
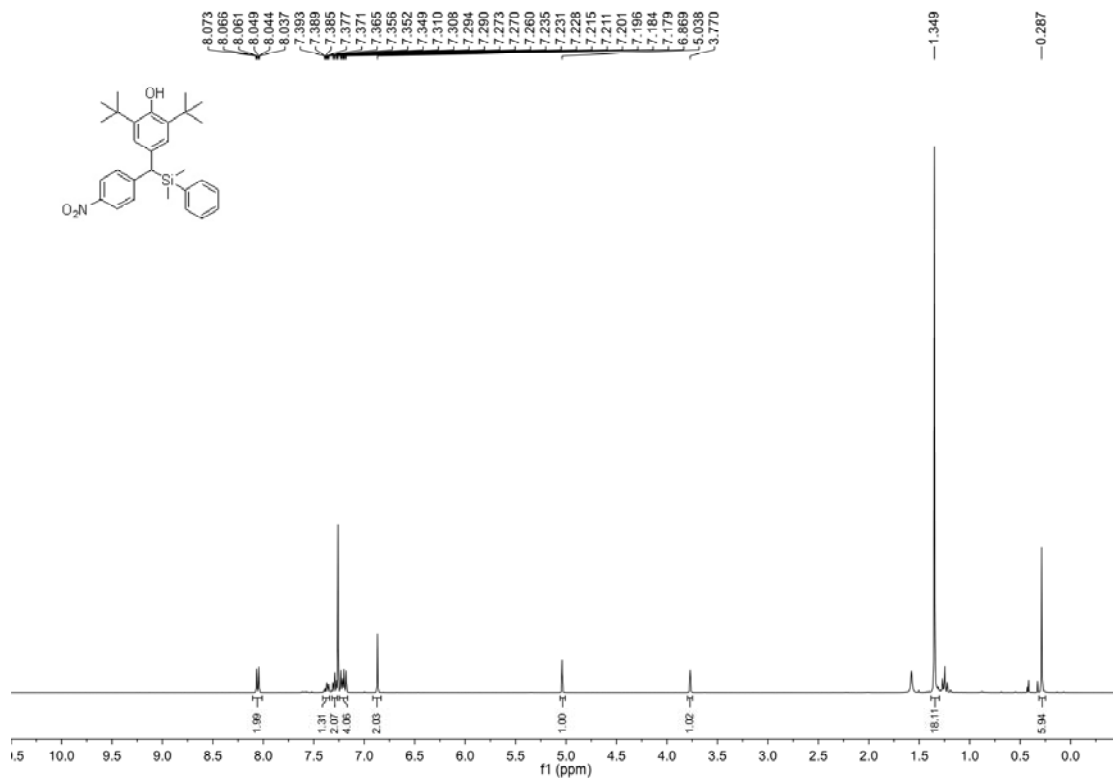


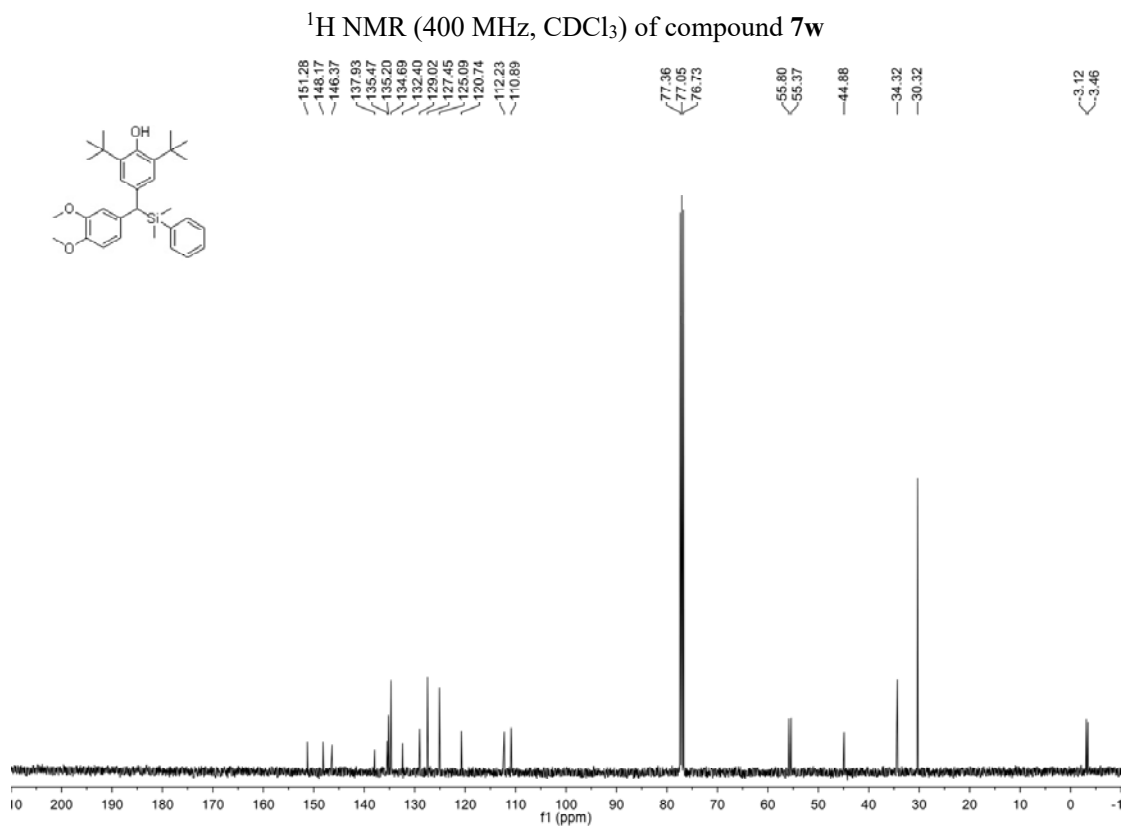
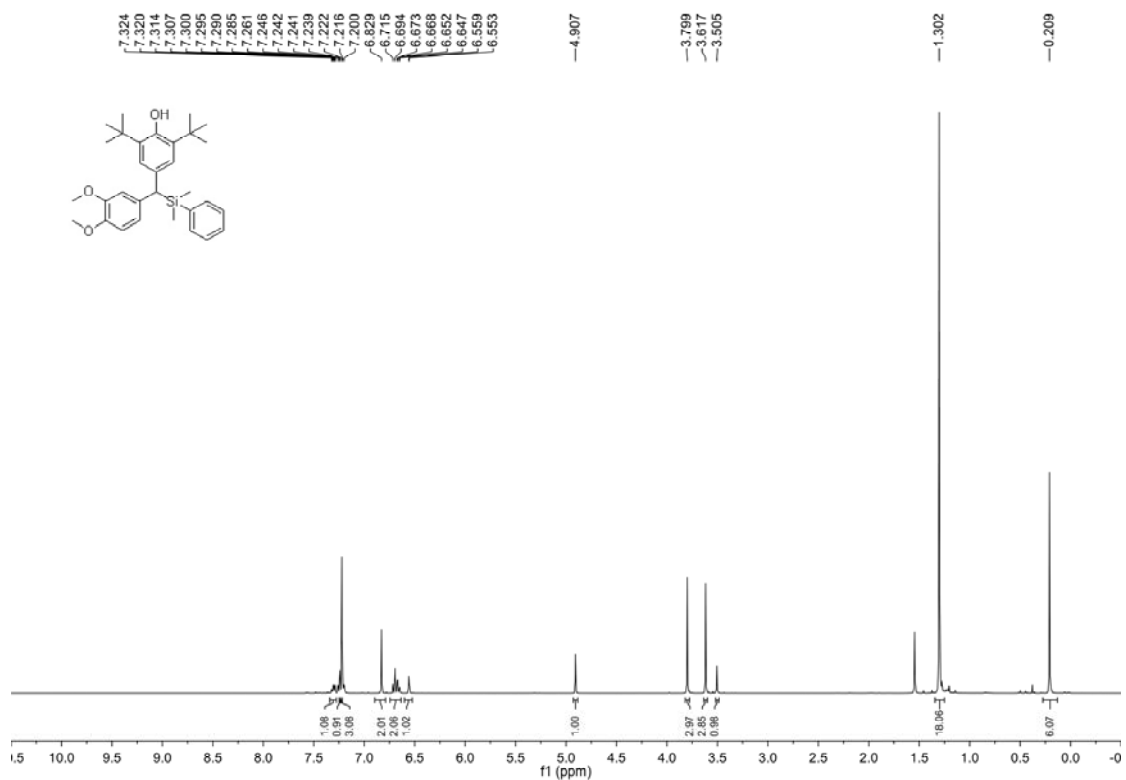


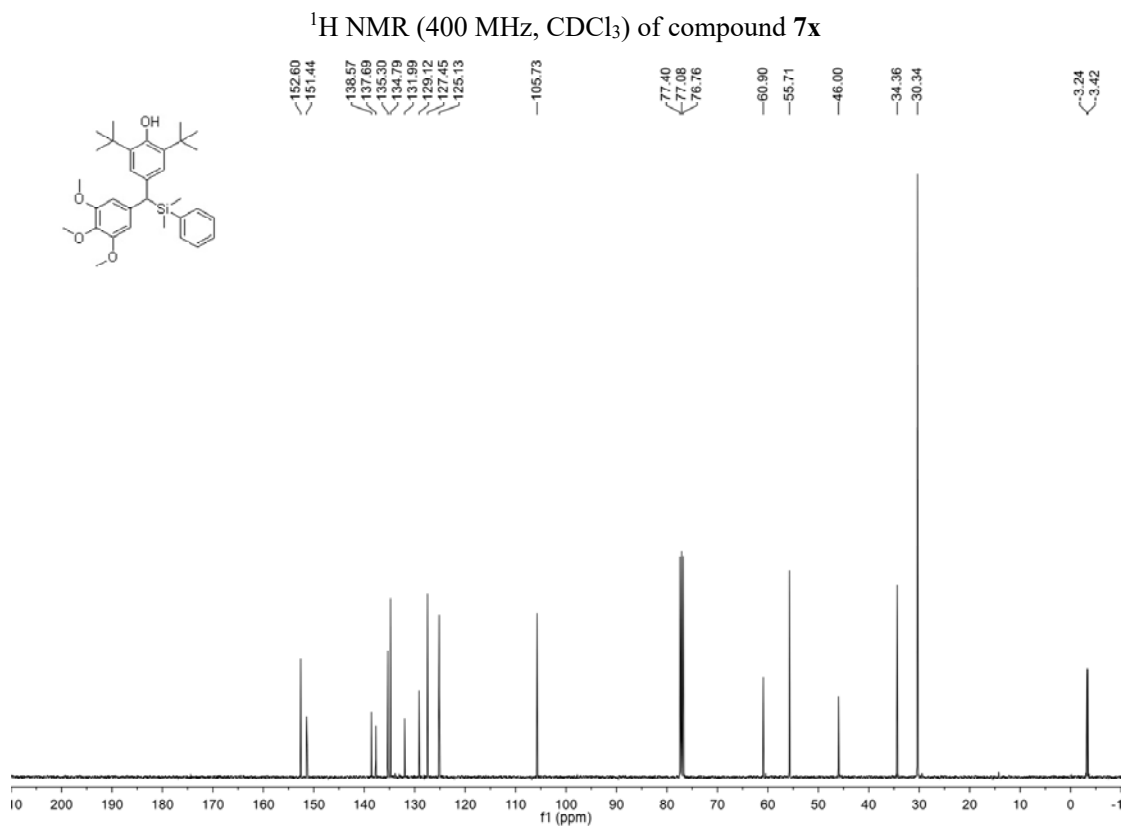
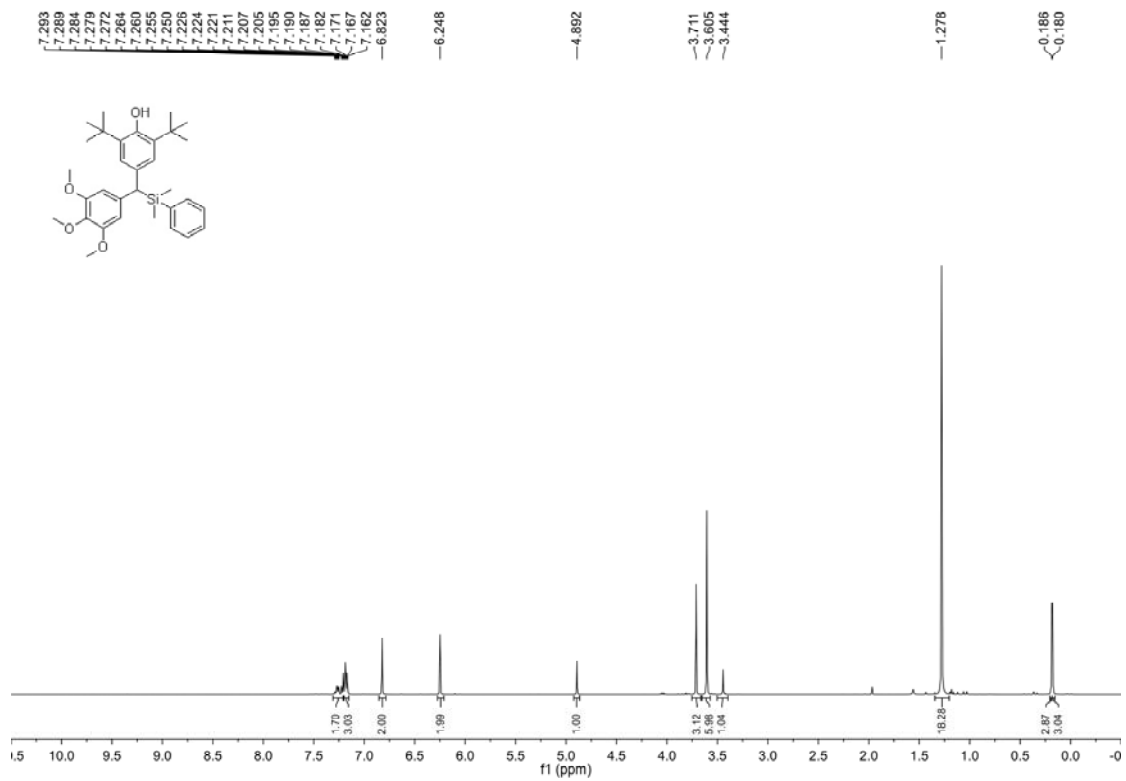
¹H NMR (400 MHz, CDCl₃) of compound **7u**

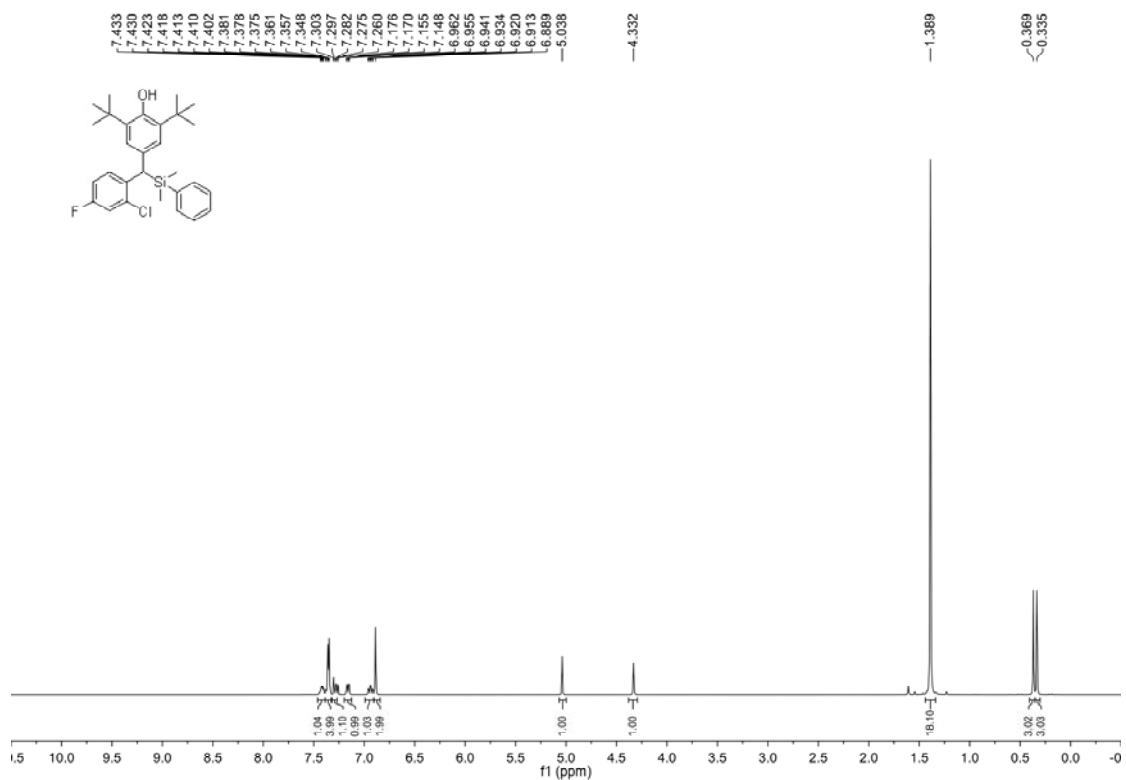


¹³C {¹H} NMR (100 MHz, CDCl₃) of compound **7u**

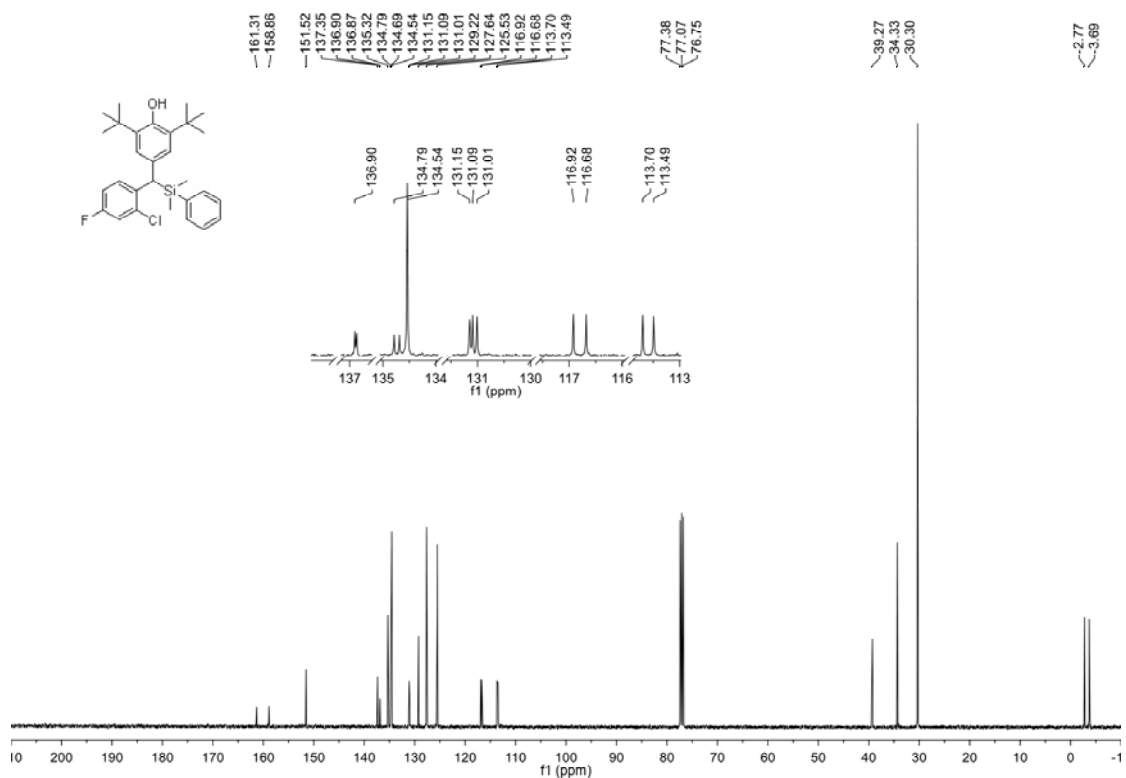




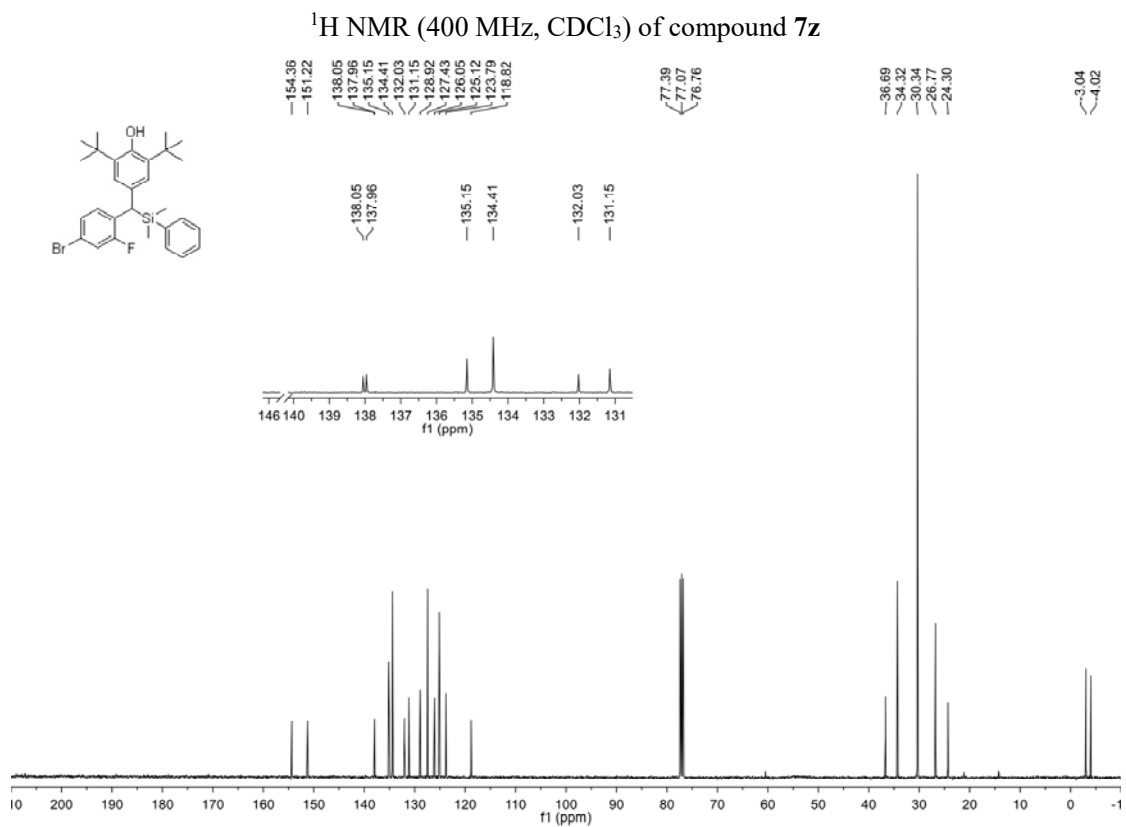
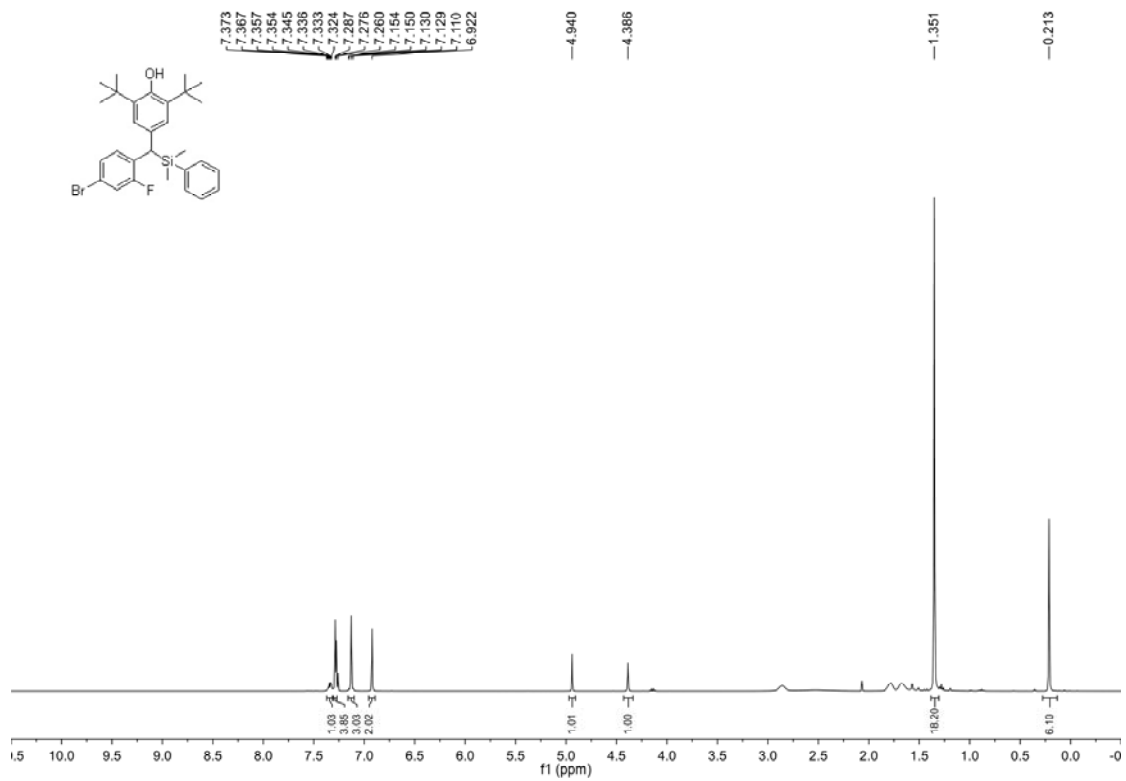


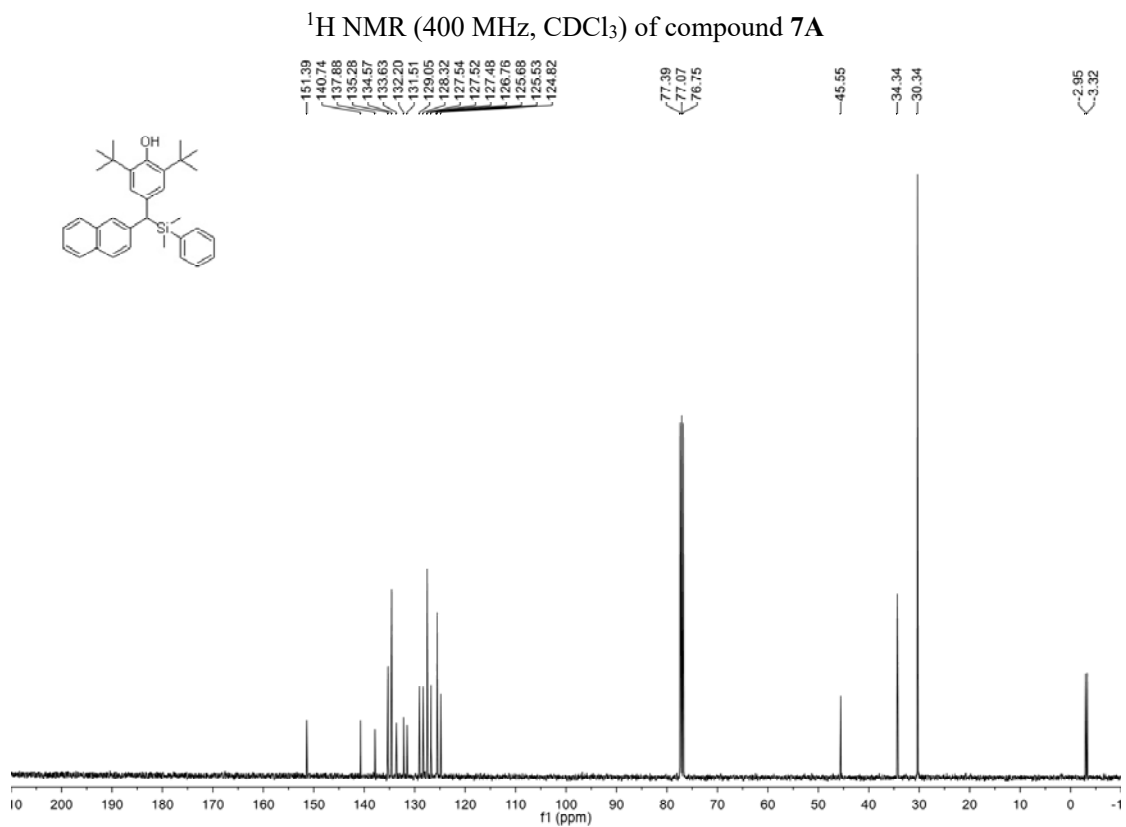
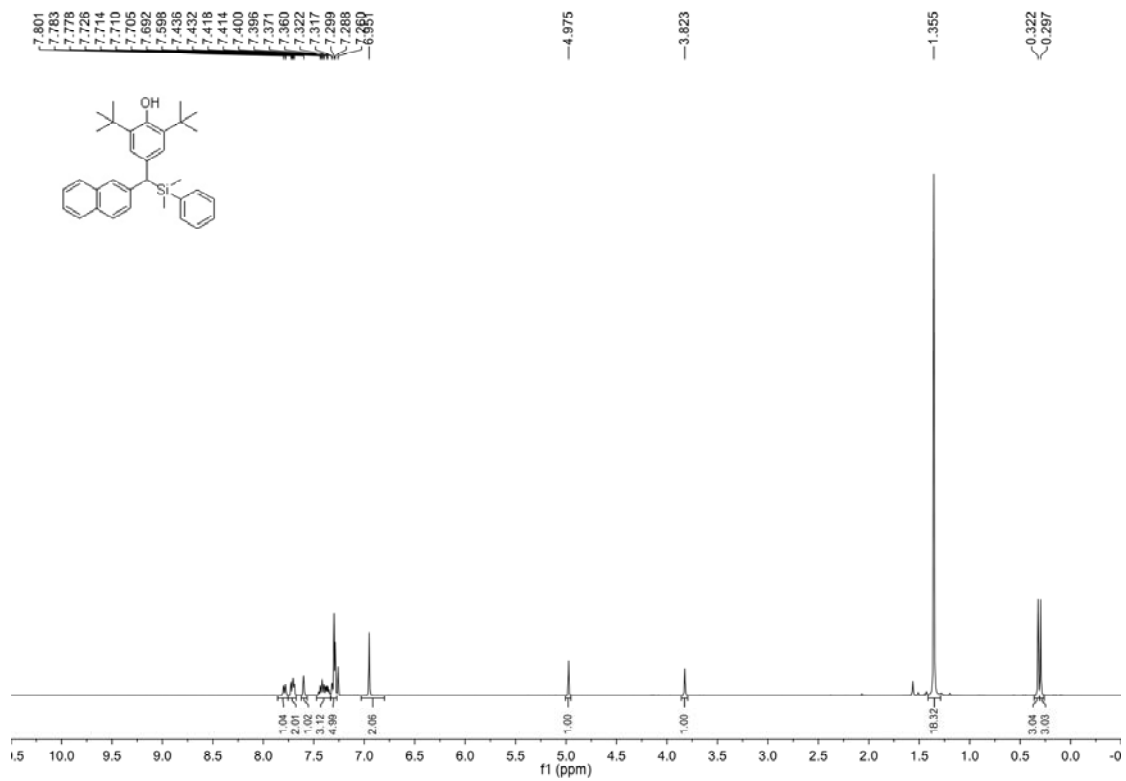


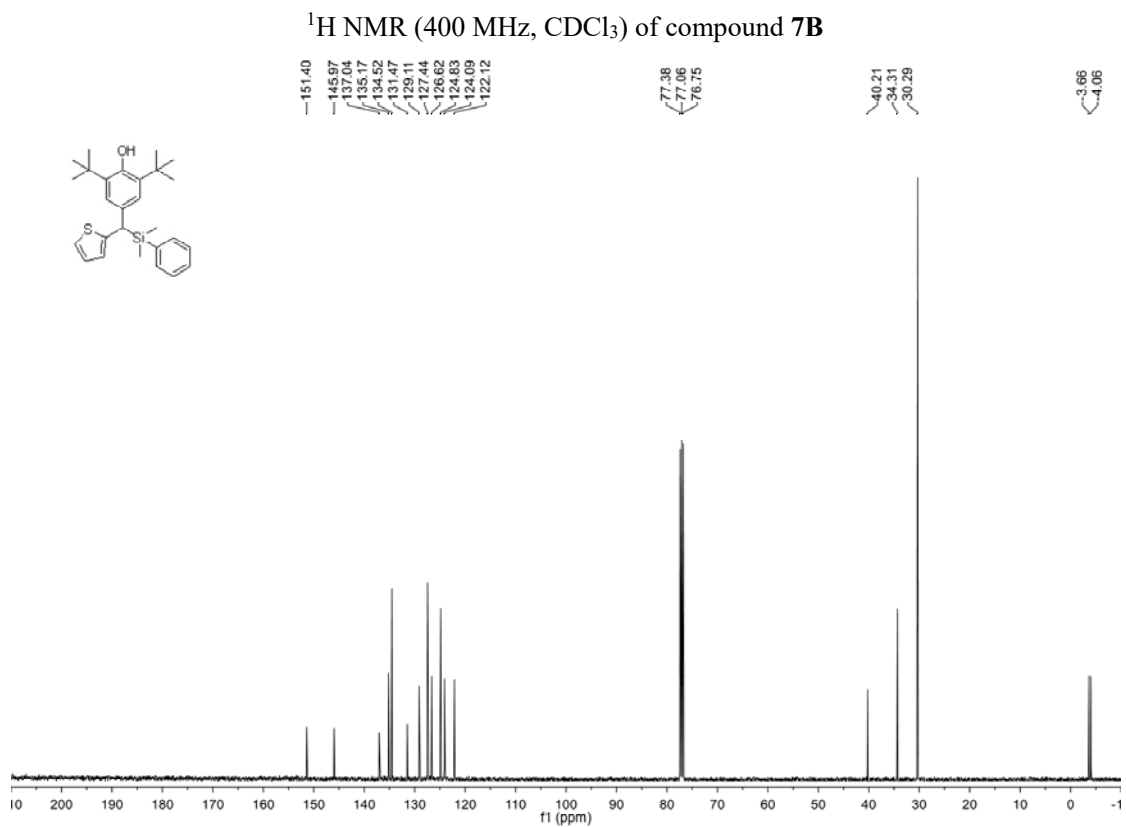
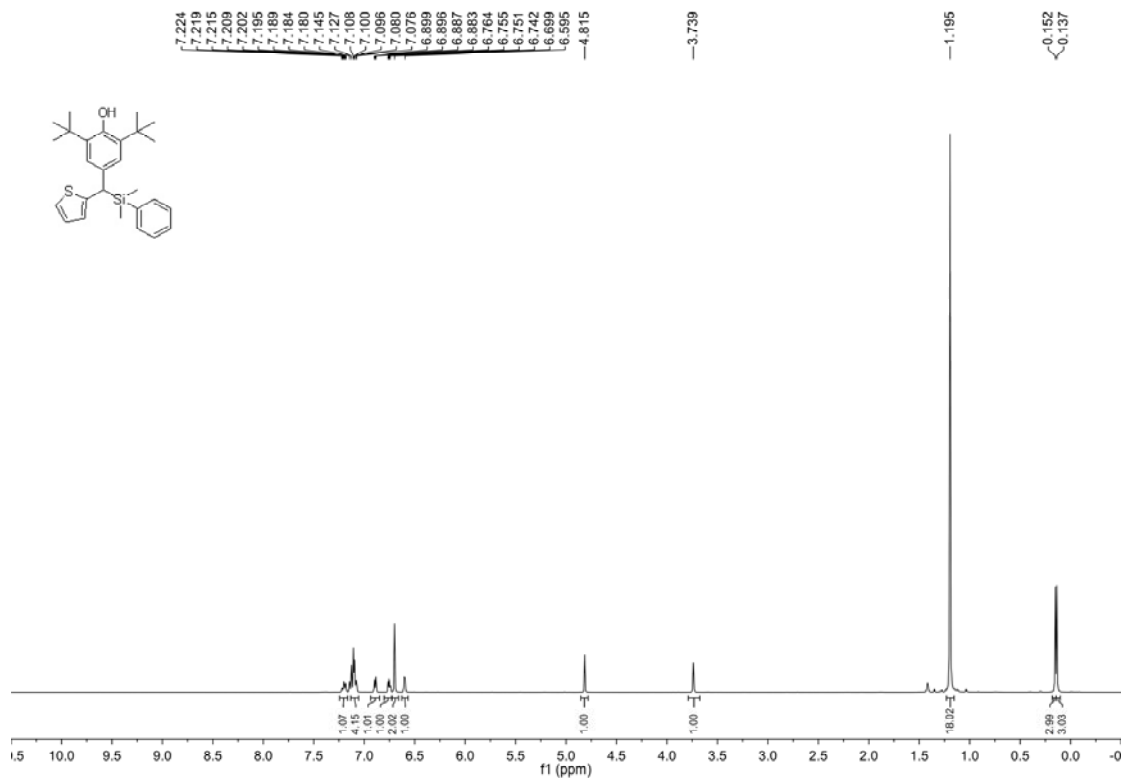
¹H NMR (400 MHz, CDCl₃) of compound **7y**

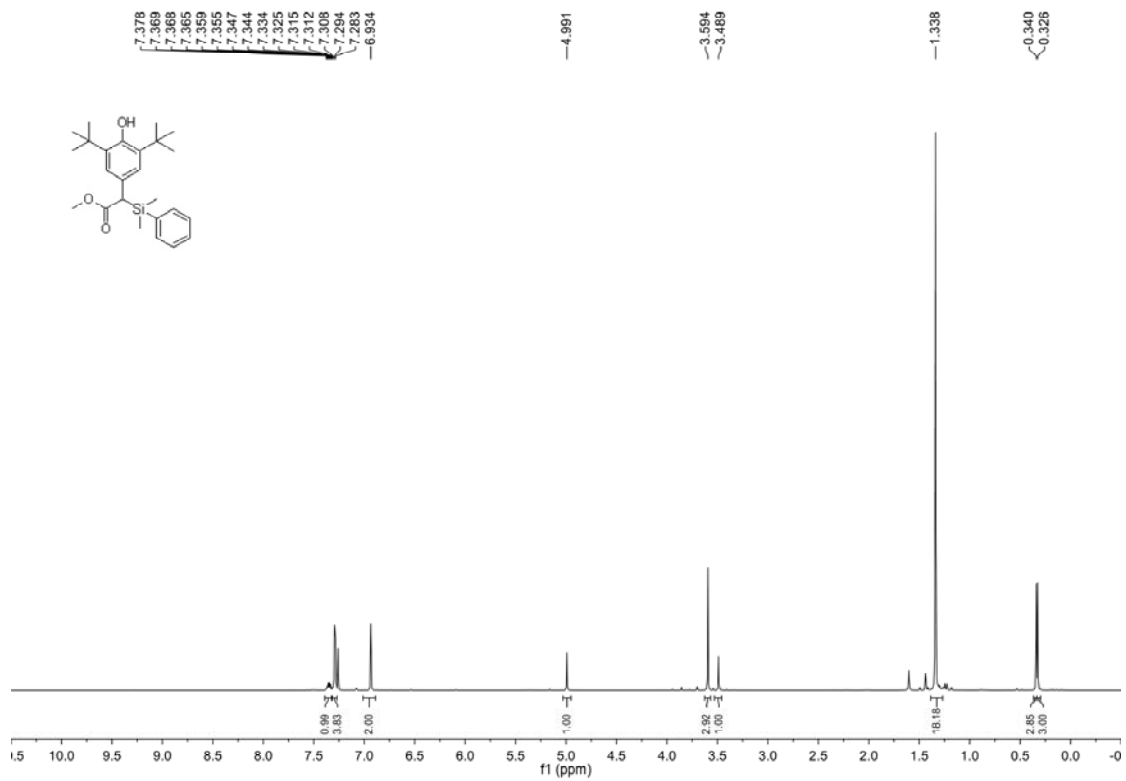


¹³C {¹H} NMR (100 MHz, CDCl₃) of compound **7y**

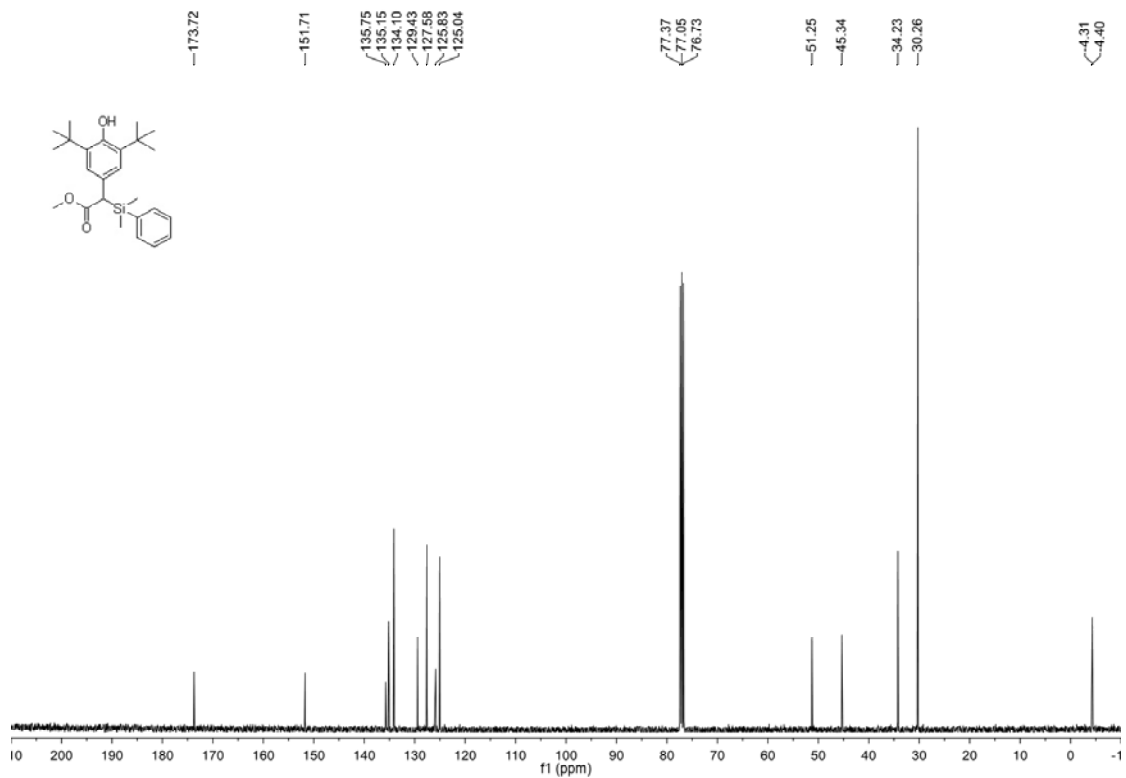




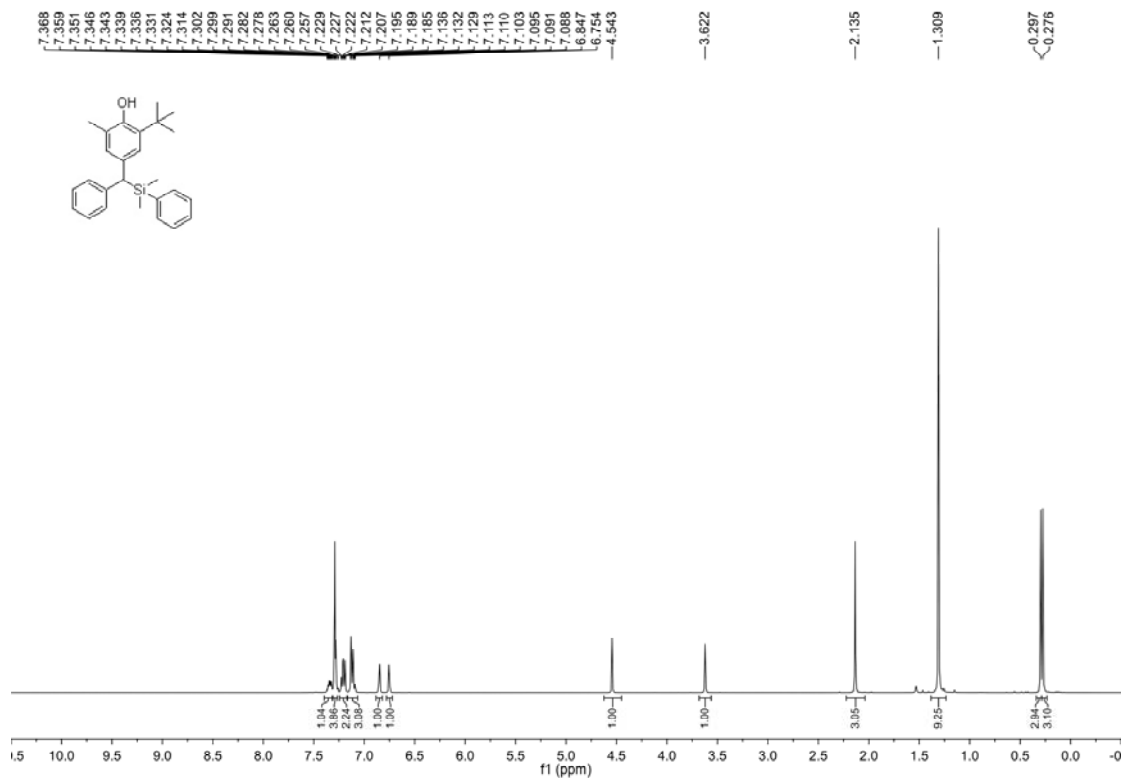




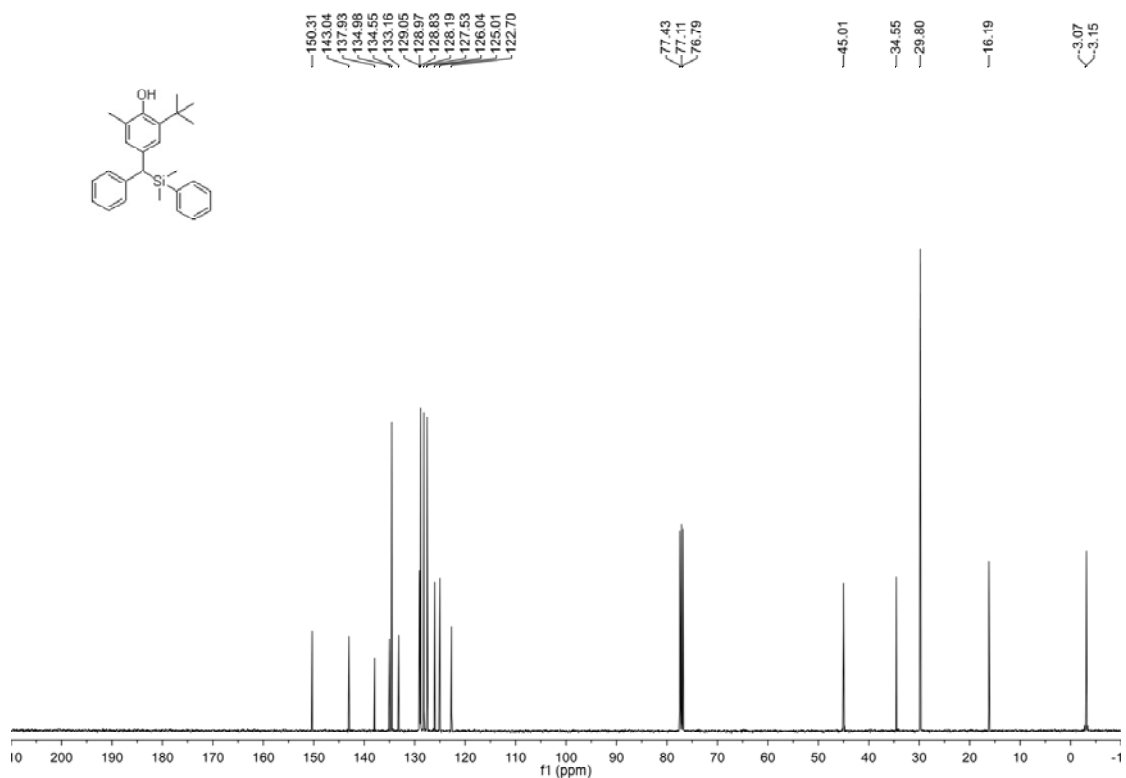
^1H NMR (400 MHz, CDCl_3) of compound **7C**



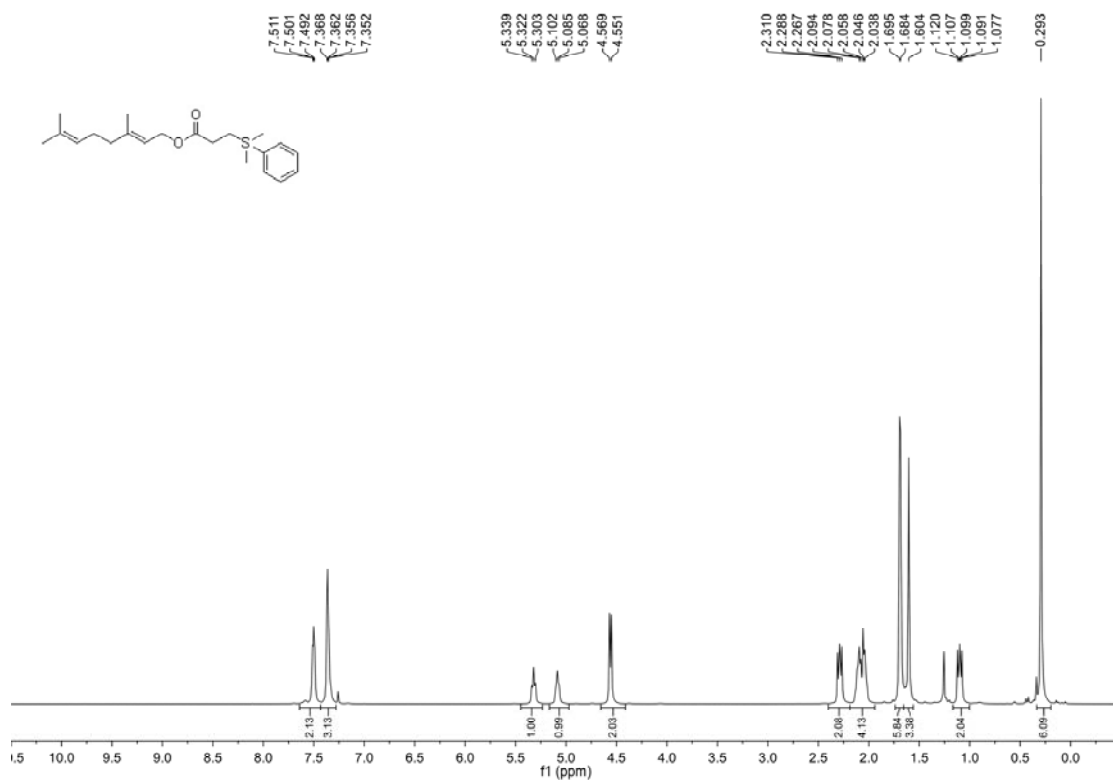
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **7C**



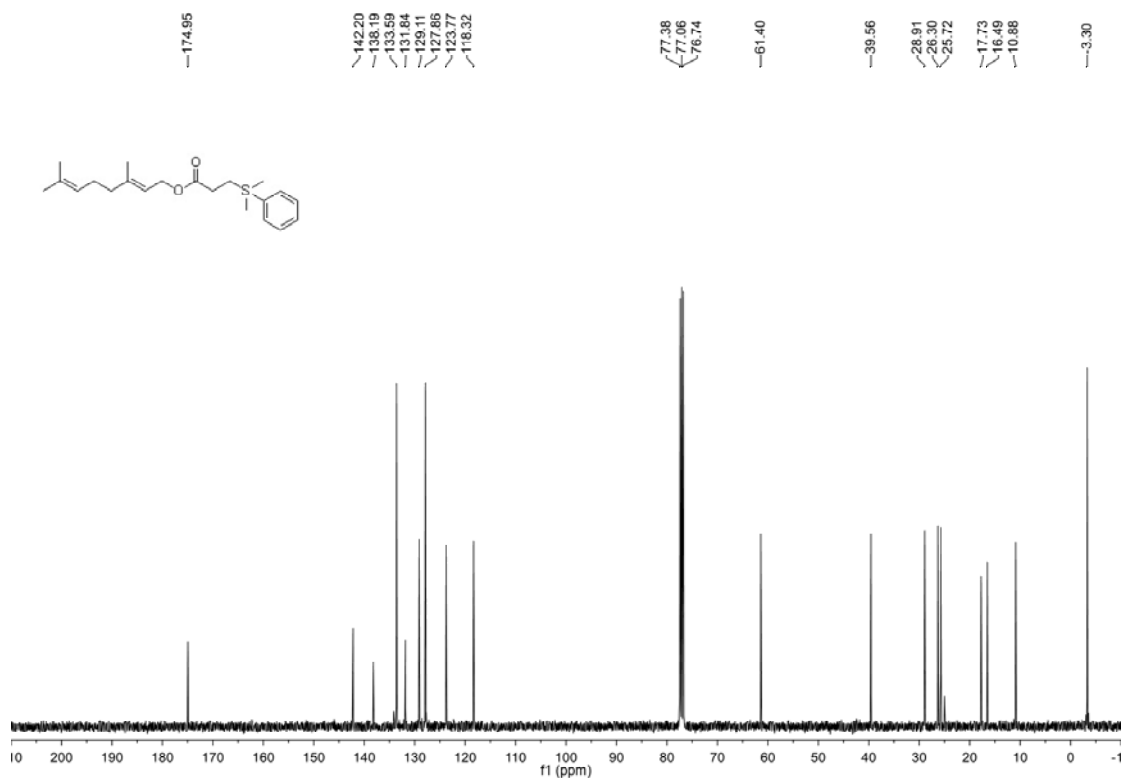
$^1\text{H NMR}$ (400 MHz, CDCl_3) of compound 7D



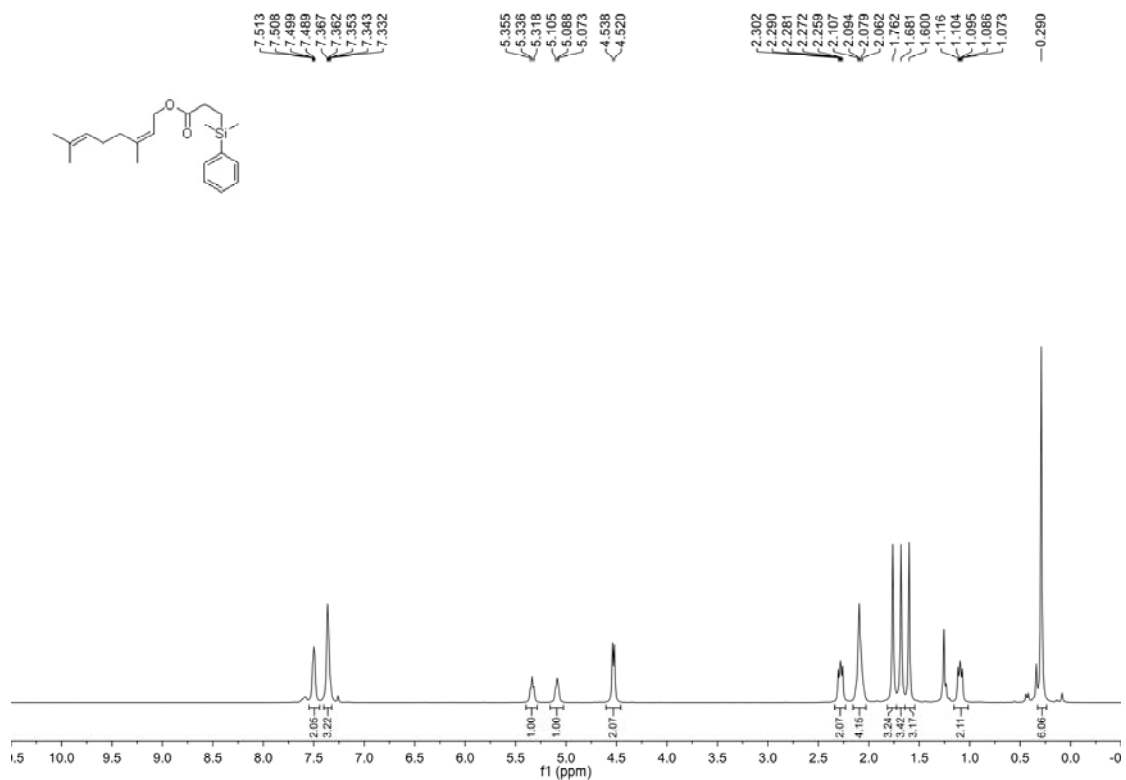
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound 7D



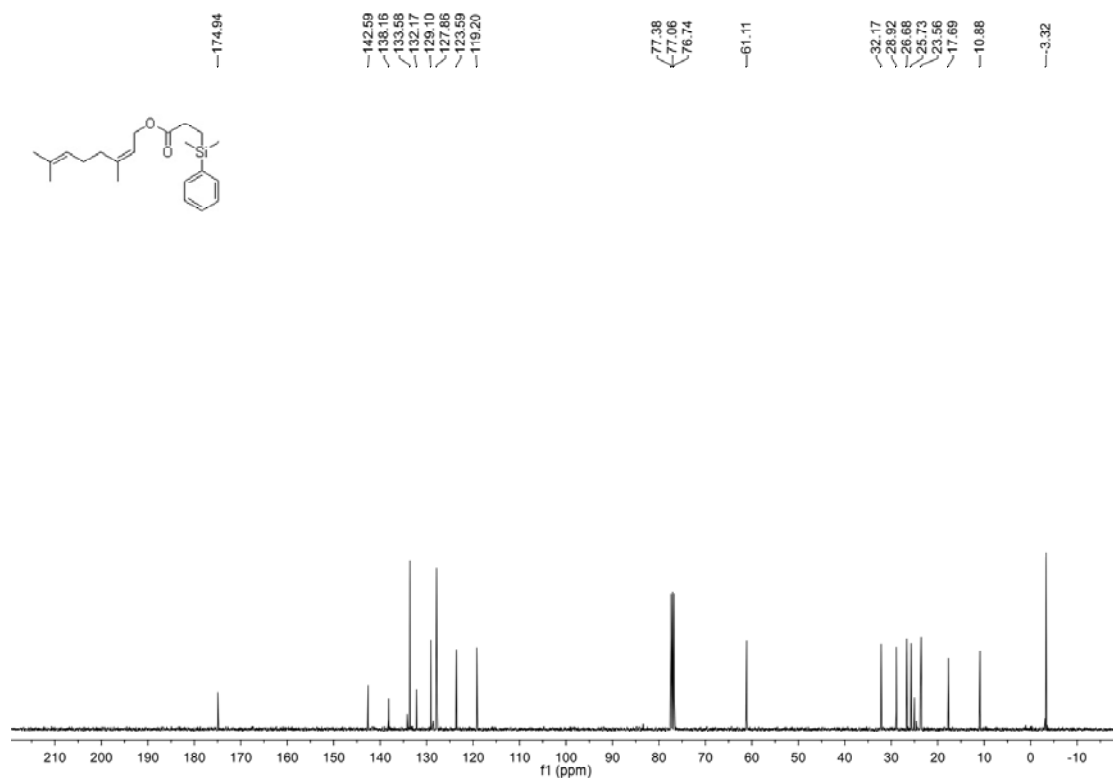
$^1\text{H NMR}$ (400 MHz, CDCl_3) of compound **9a**



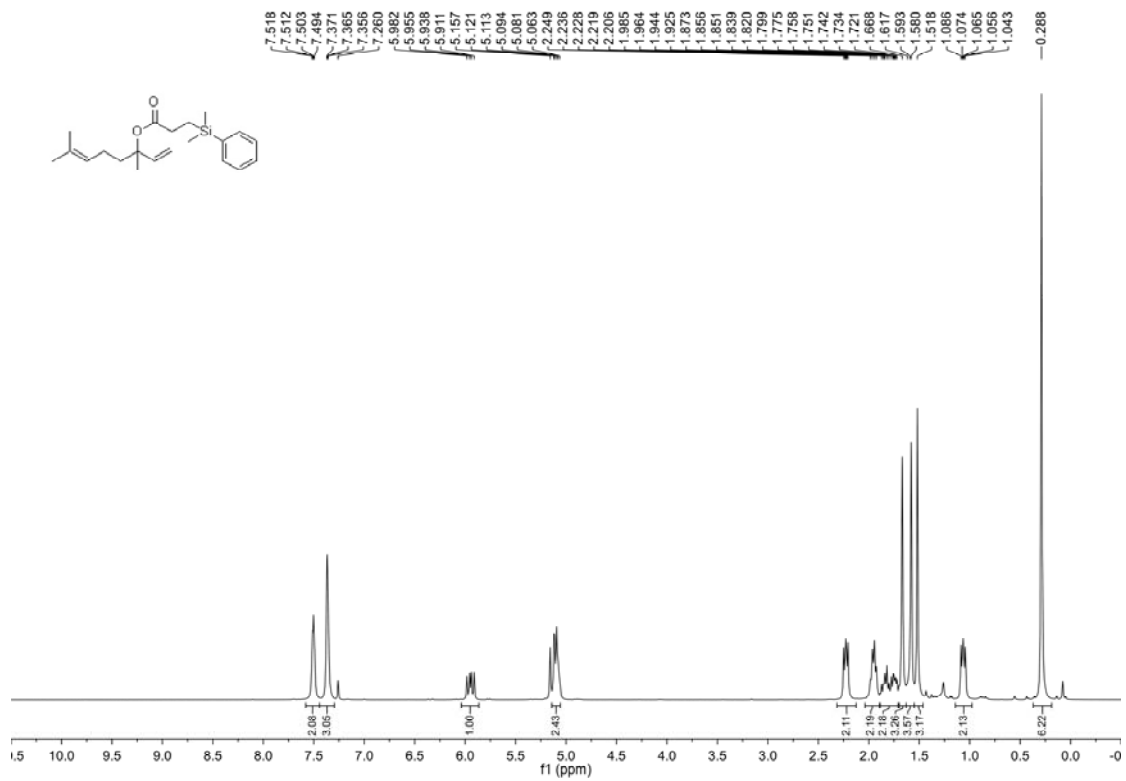
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **9a**



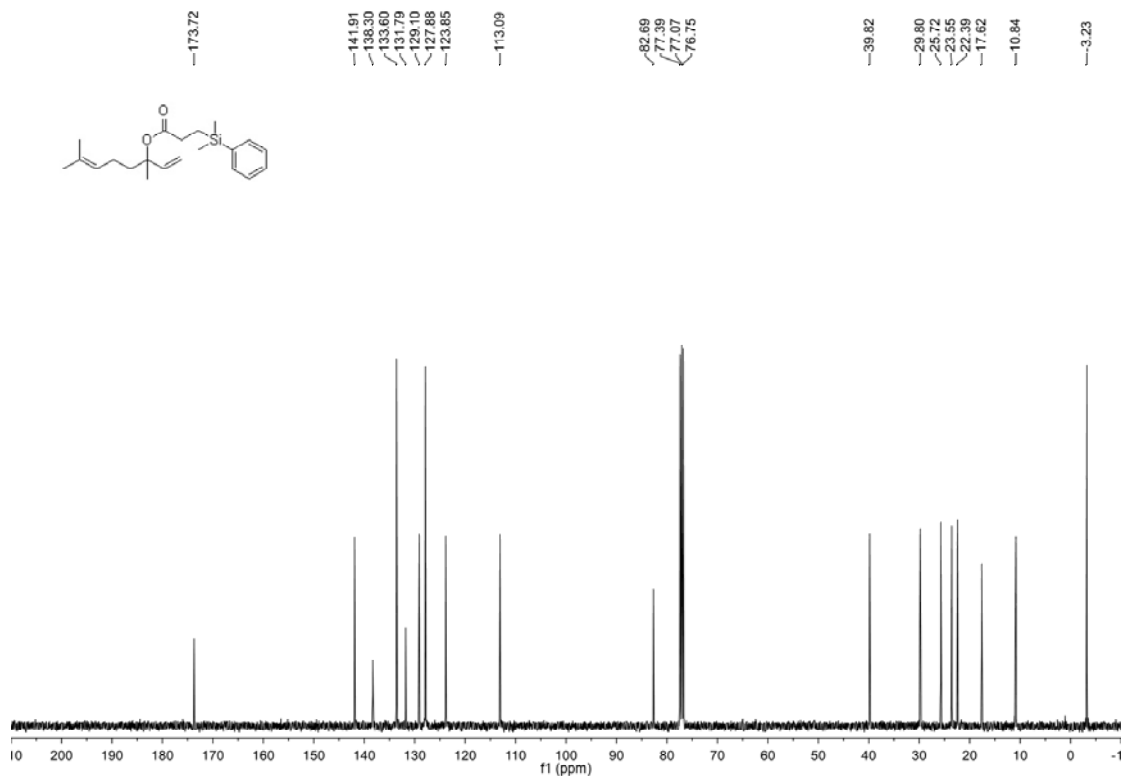
$^1\text{H NMR}$ (400 MHz, CDCl_3) of compound **9b**



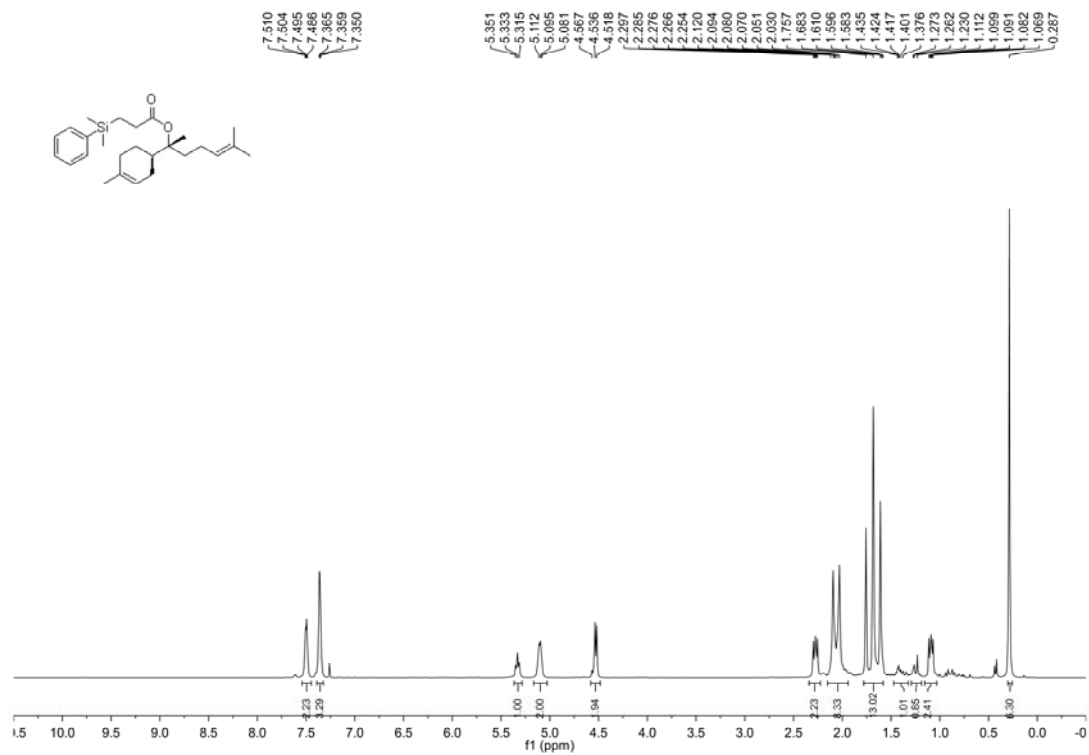
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **9b**



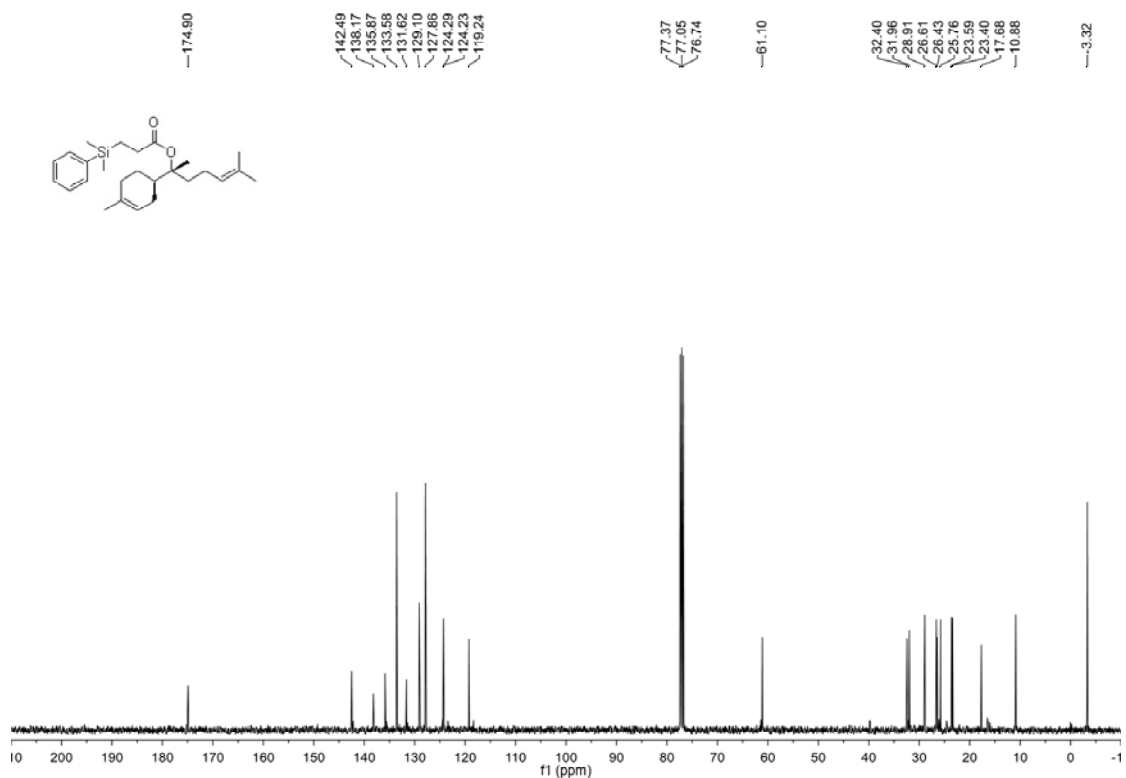
$^1\text{H NMR}$ (400 MHz, CDCl_3) of compound **9c**



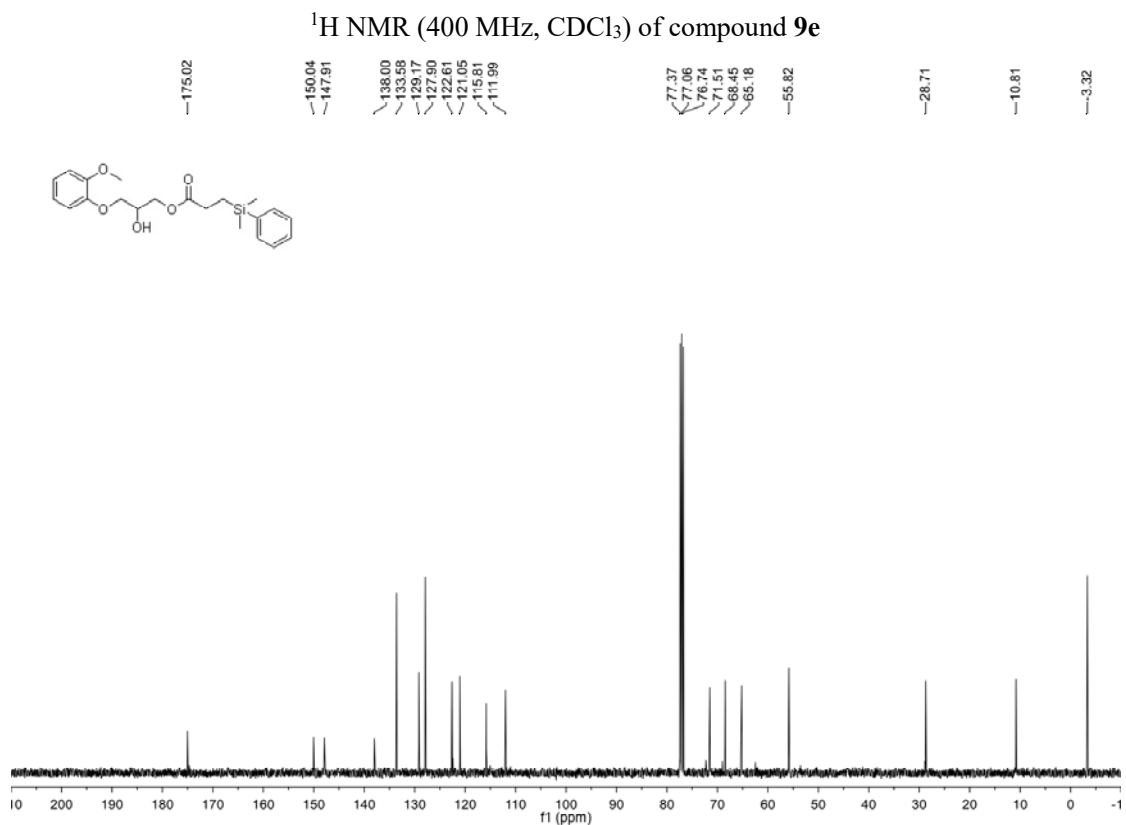
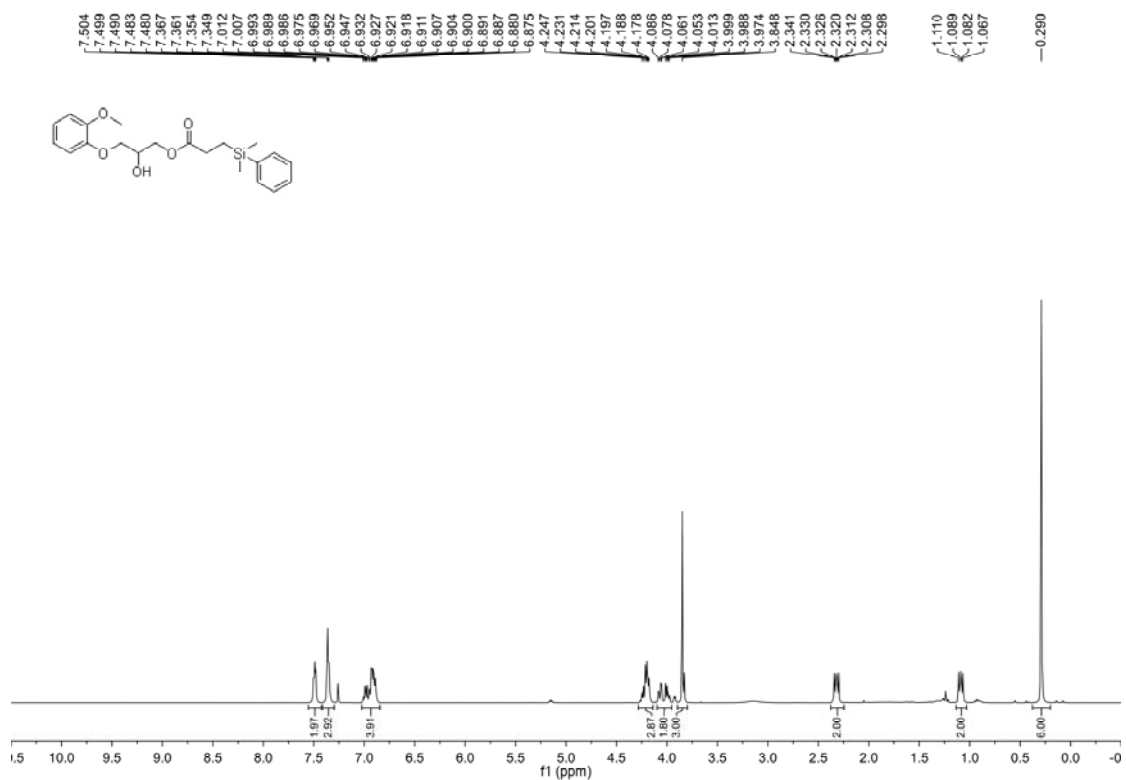
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **9c**

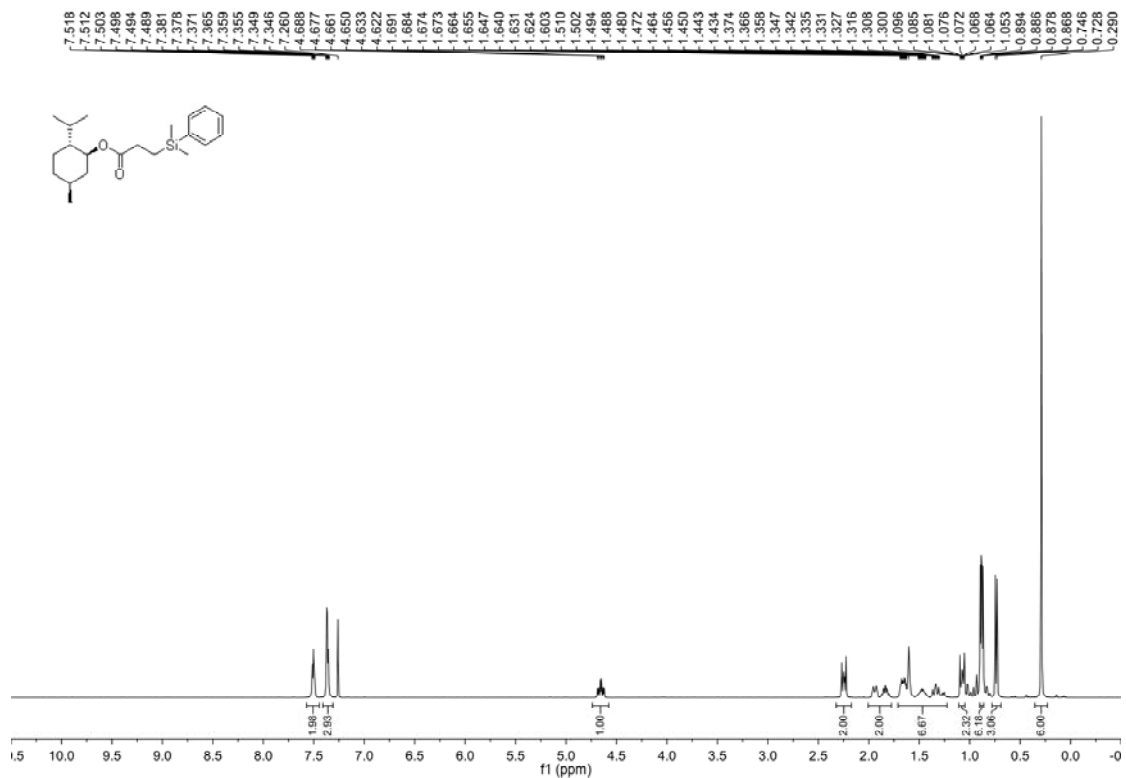


¹H NMR (400 MHz, CDCl₃) of compound **9d**

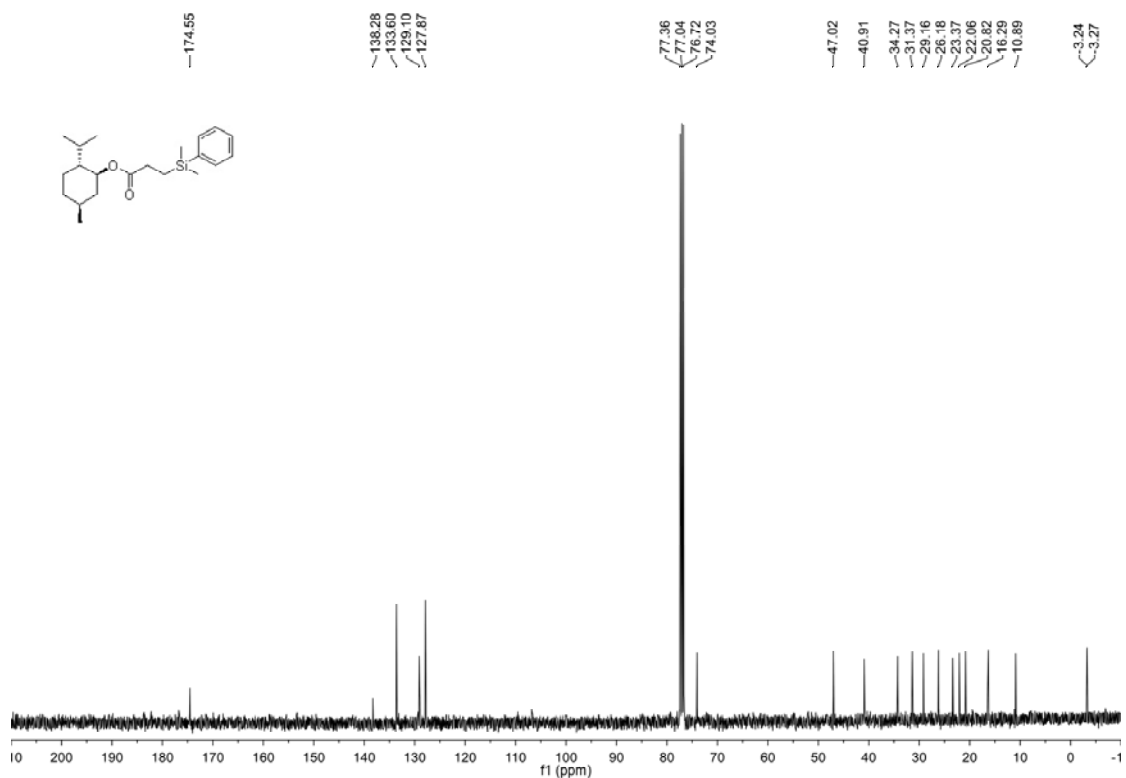


¹³C{¹H} NMR (100 MHz, CDCl₃) of compound **9d**

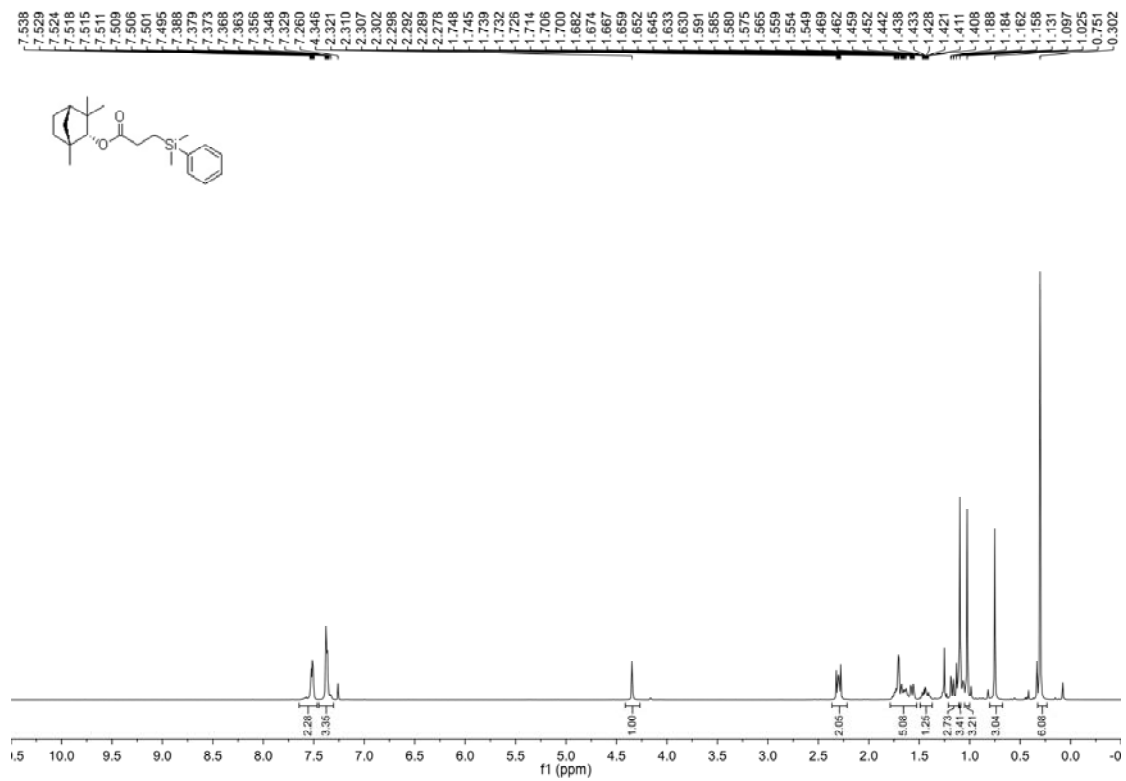




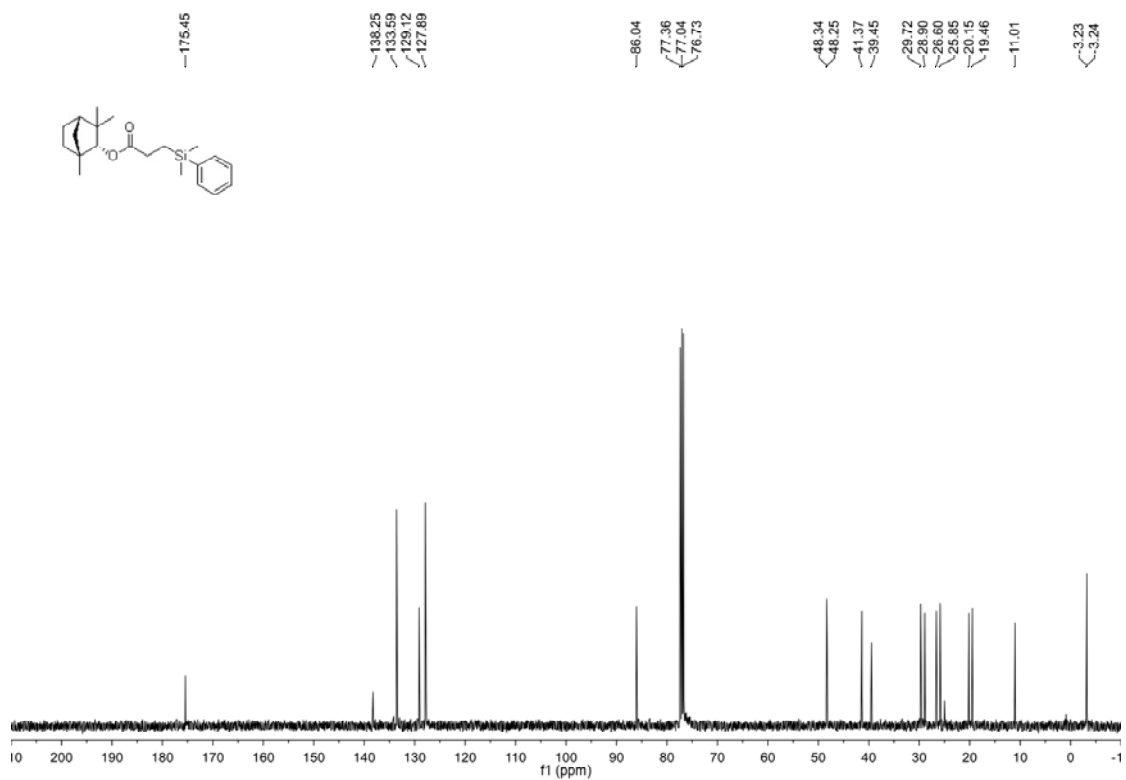
¹H NMR (400 MHz, CDCl₃) of compound **9f**



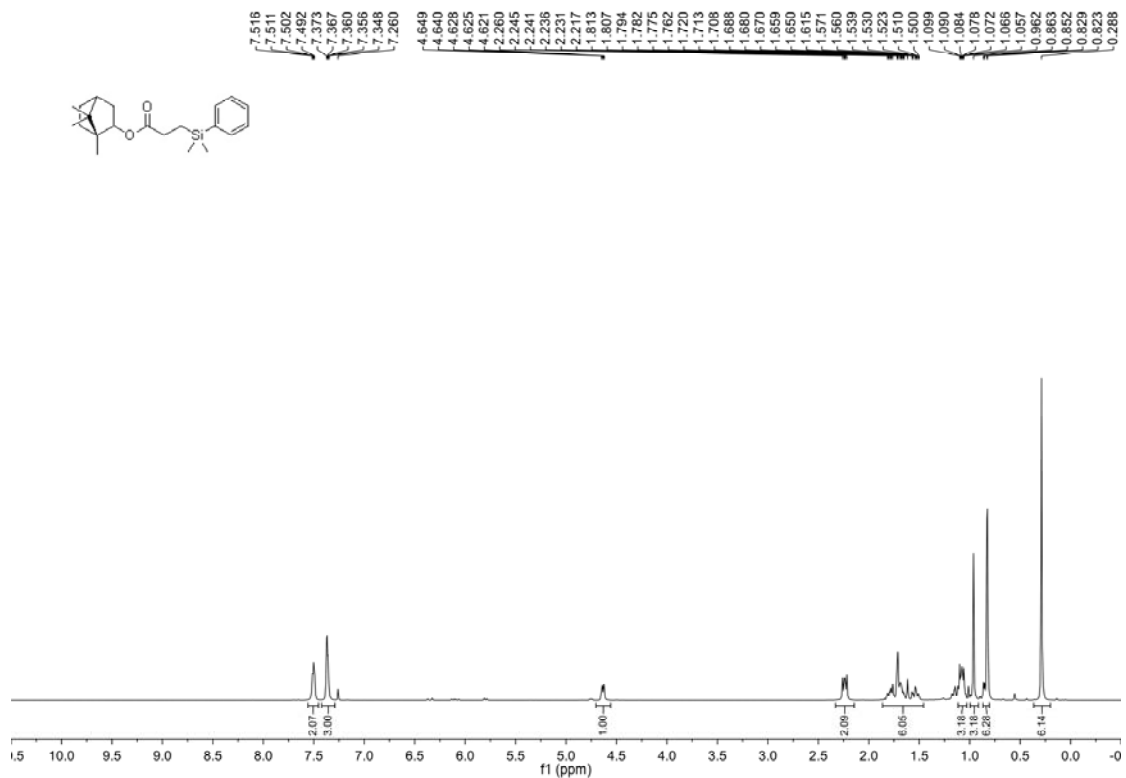
¹³C {¹H} NMR (100 MHz, CDCl₃) of compound **9f**



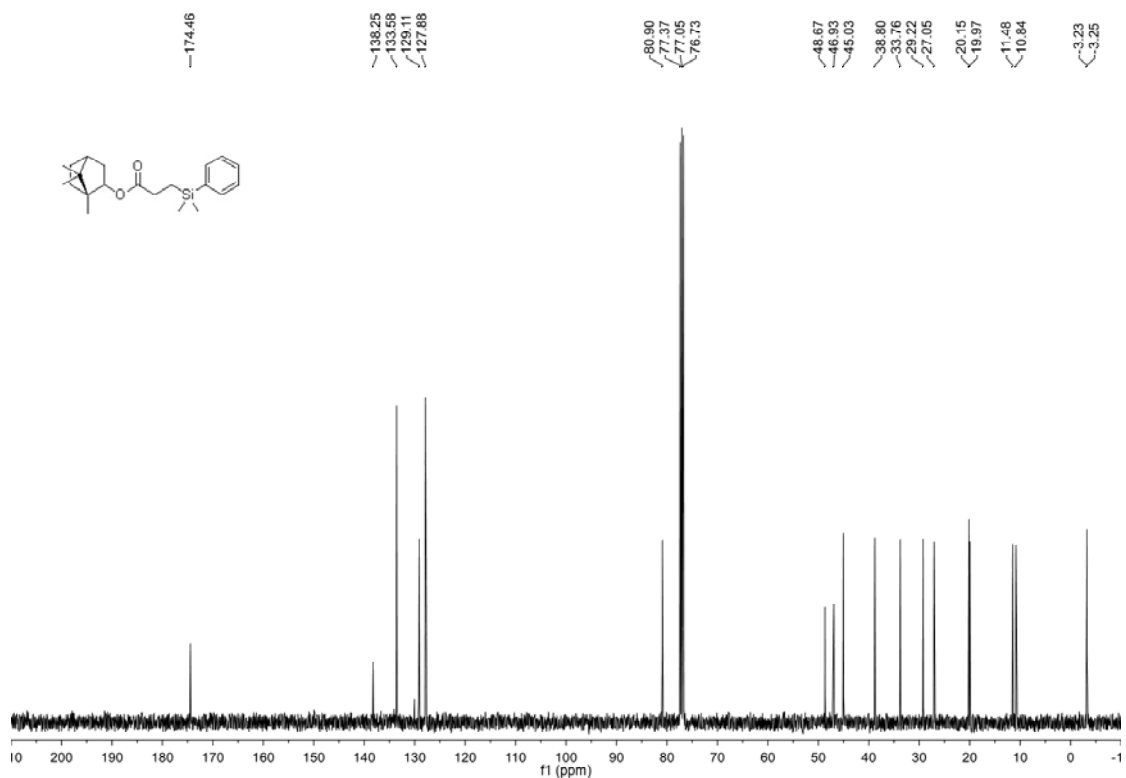
^1H NMR (400 MHz, CDCl_3) of compound **9g**



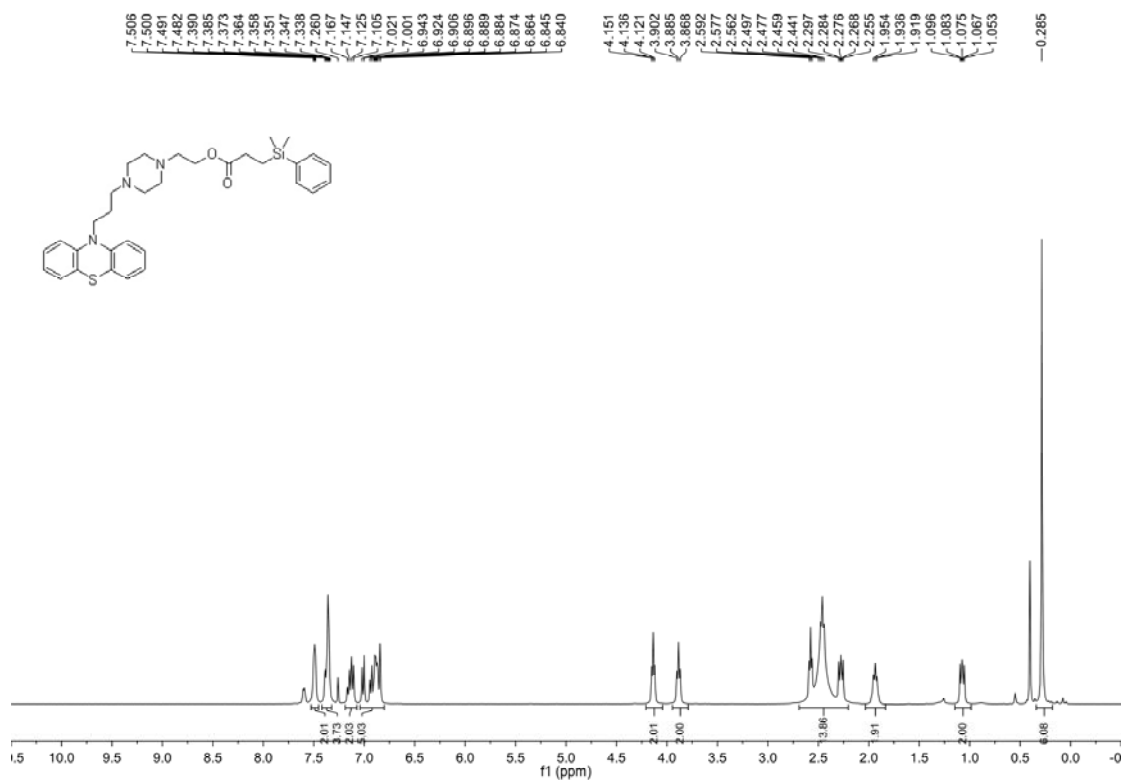
$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **9g**



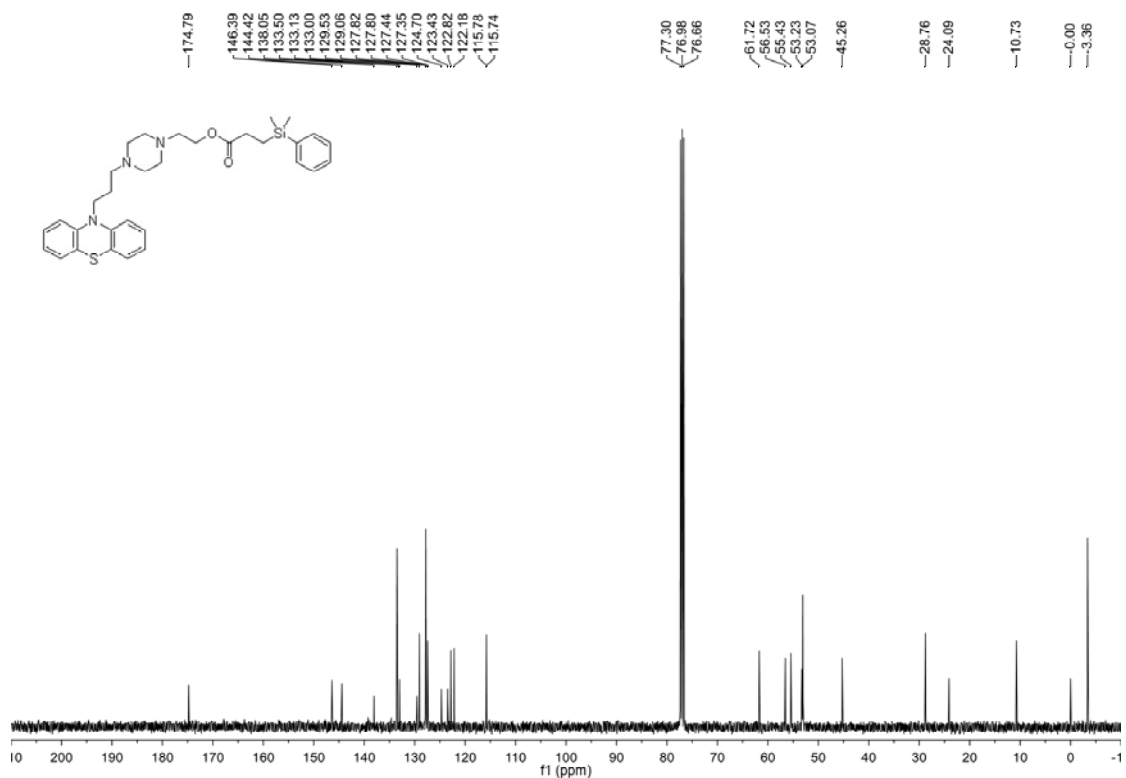
¹H NMR (400 MHz, CDCl₃) of compound **9h**



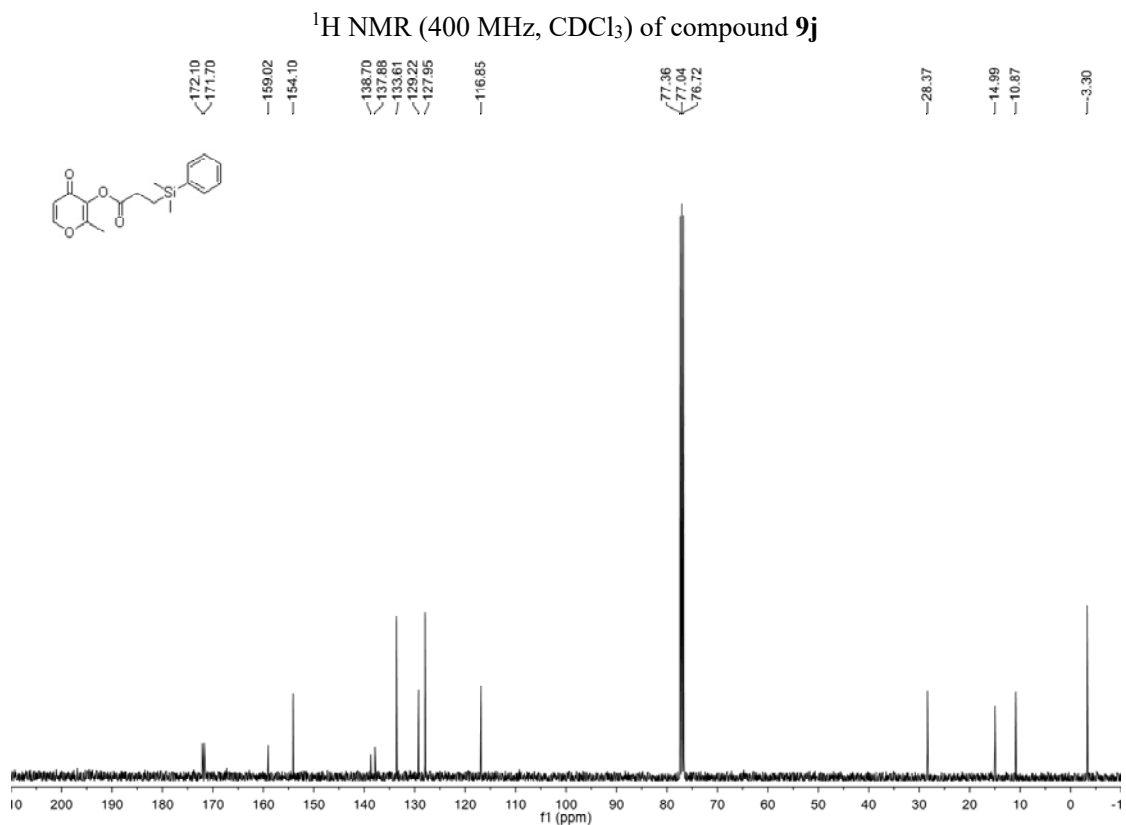
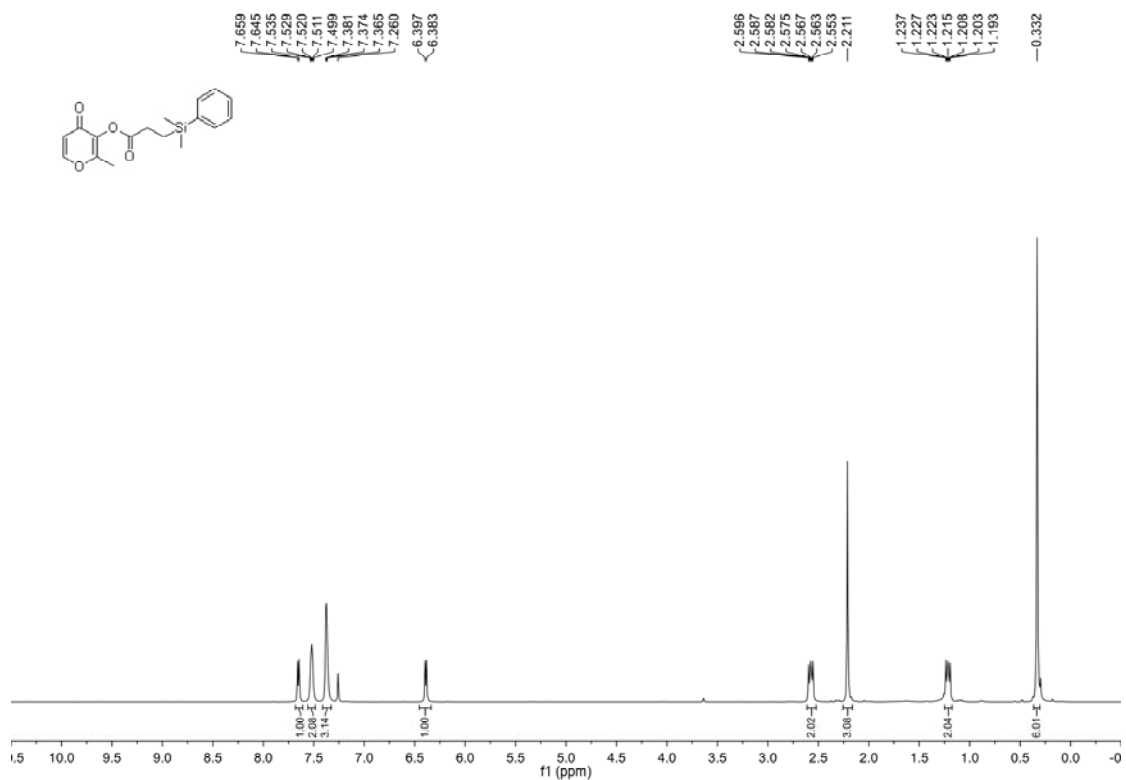
¹³C {¹H} NMR (100 MHz, CDCl₃) of compound **9h**

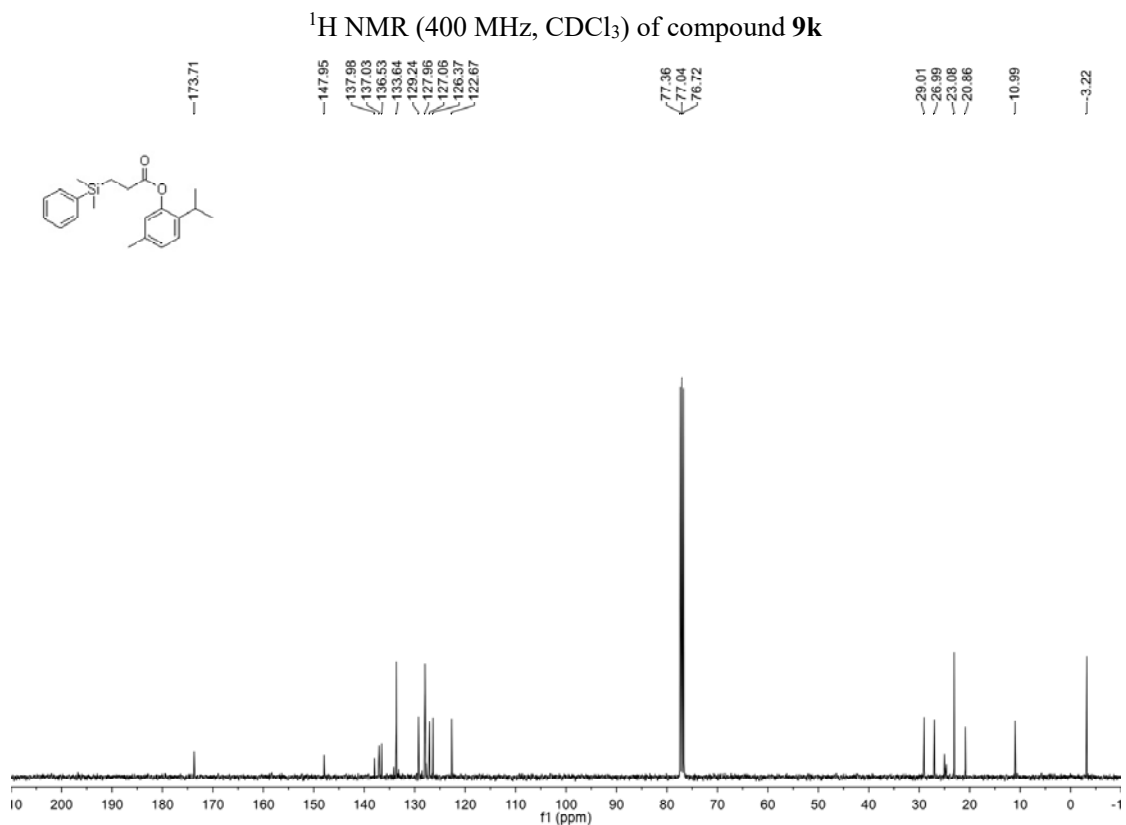
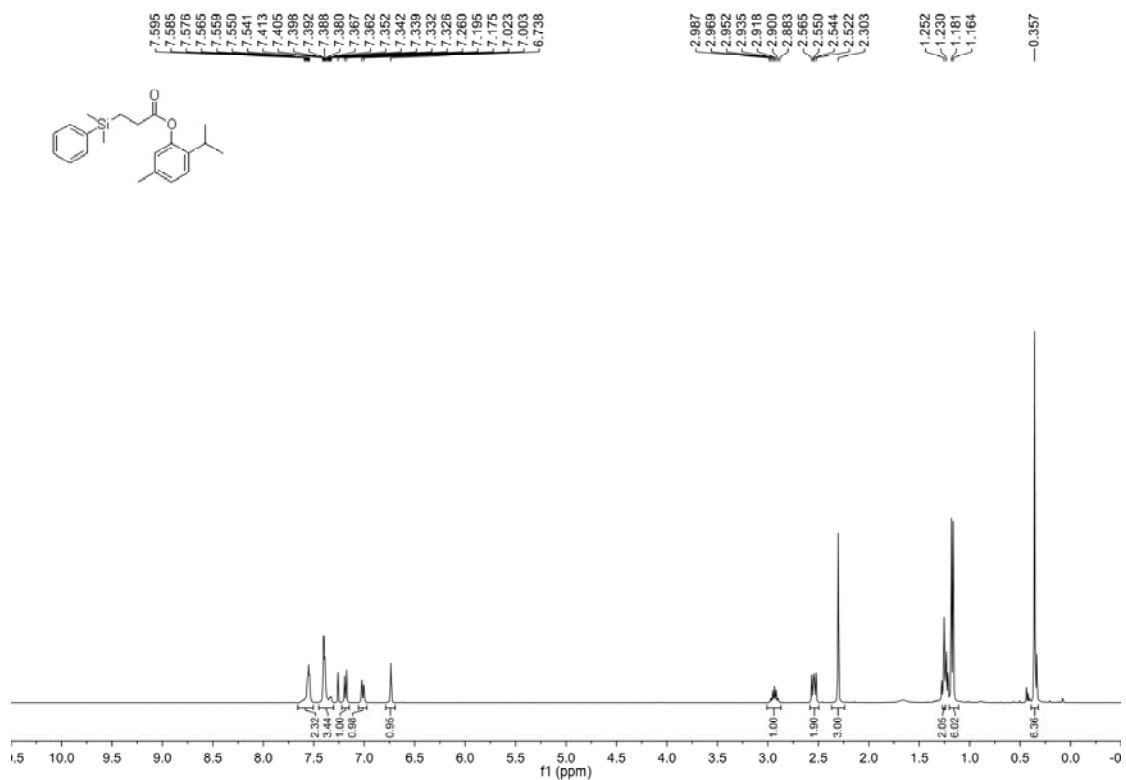


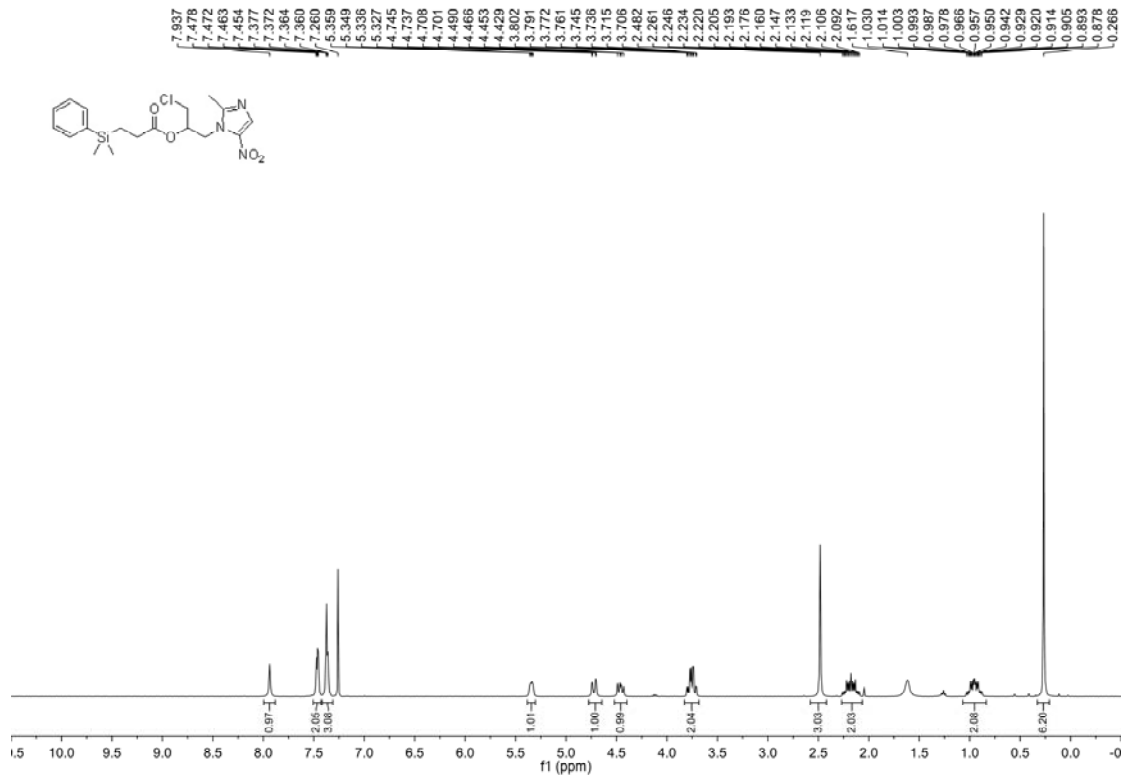
¹H NMR (400 MHz, CDCl₃) of compound **9i**



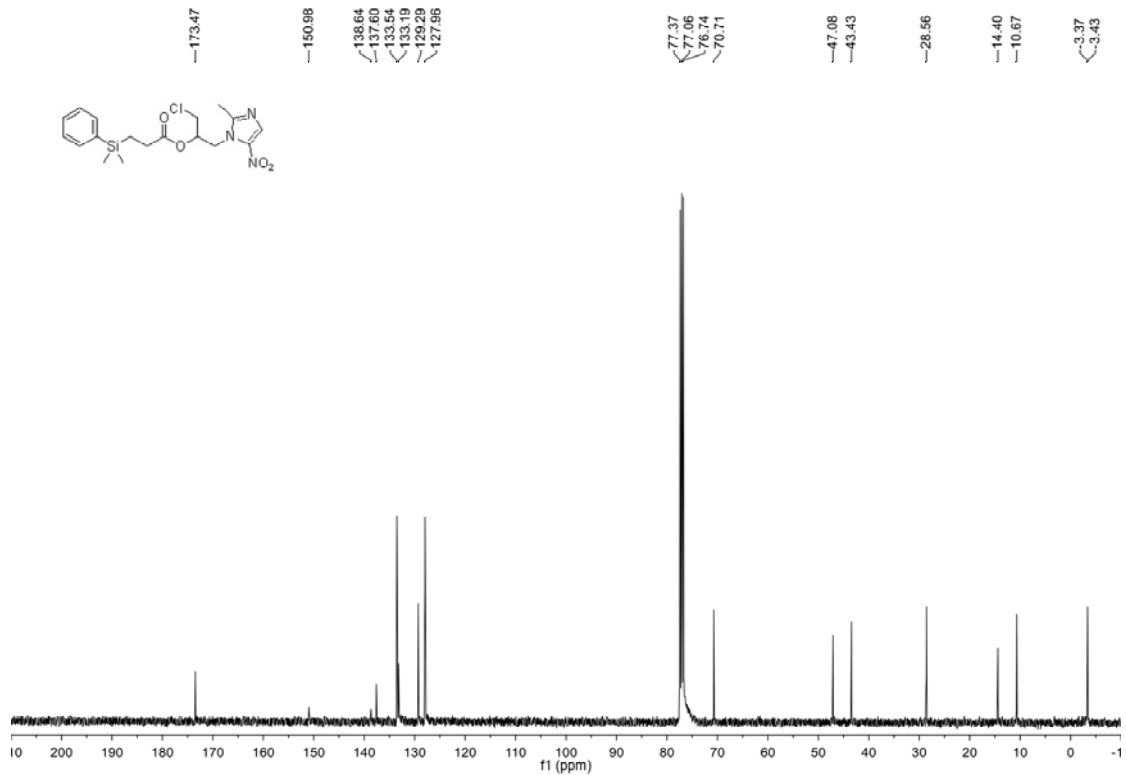
¹³C NMR (100 MHz, CDCl₃) of compound **9i**



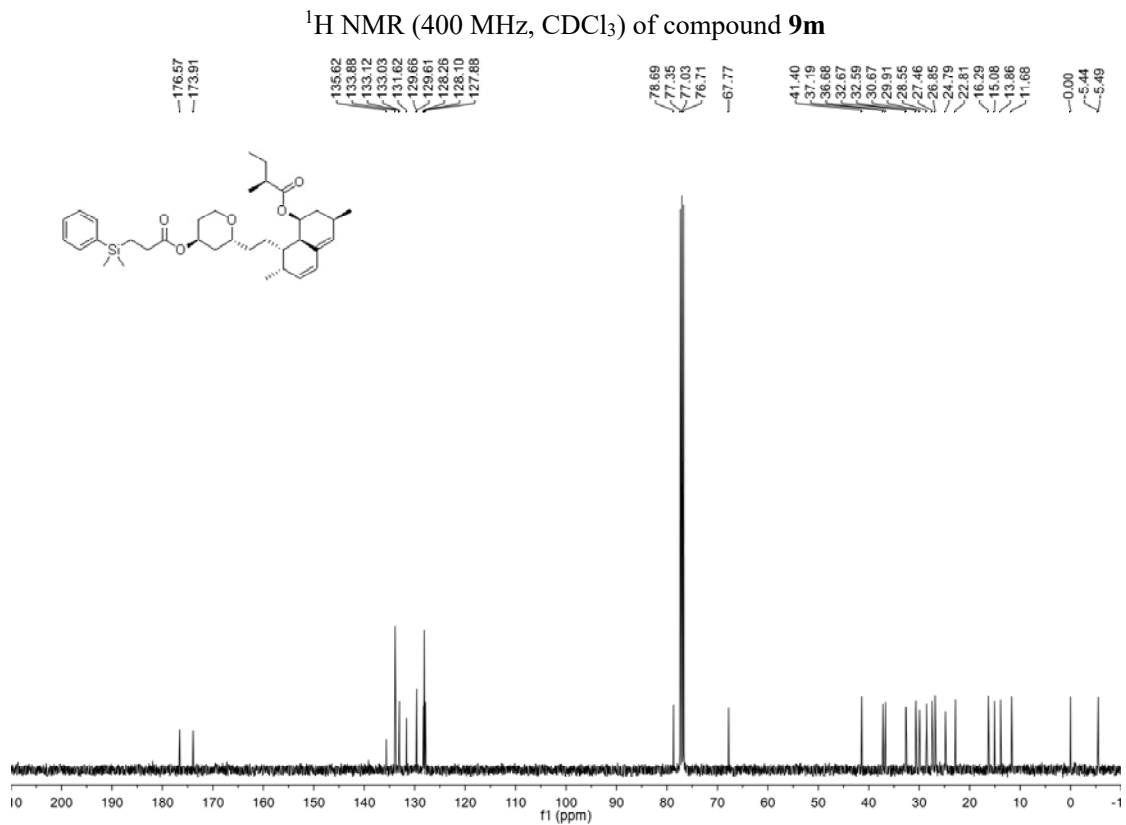
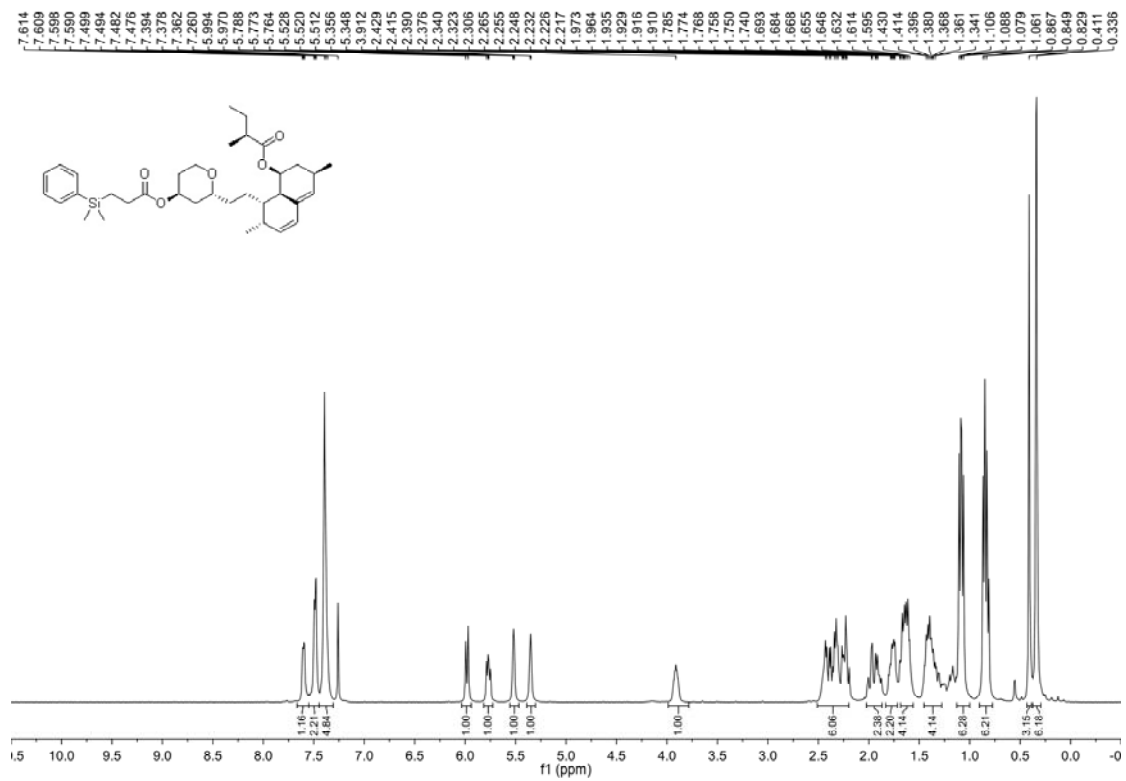


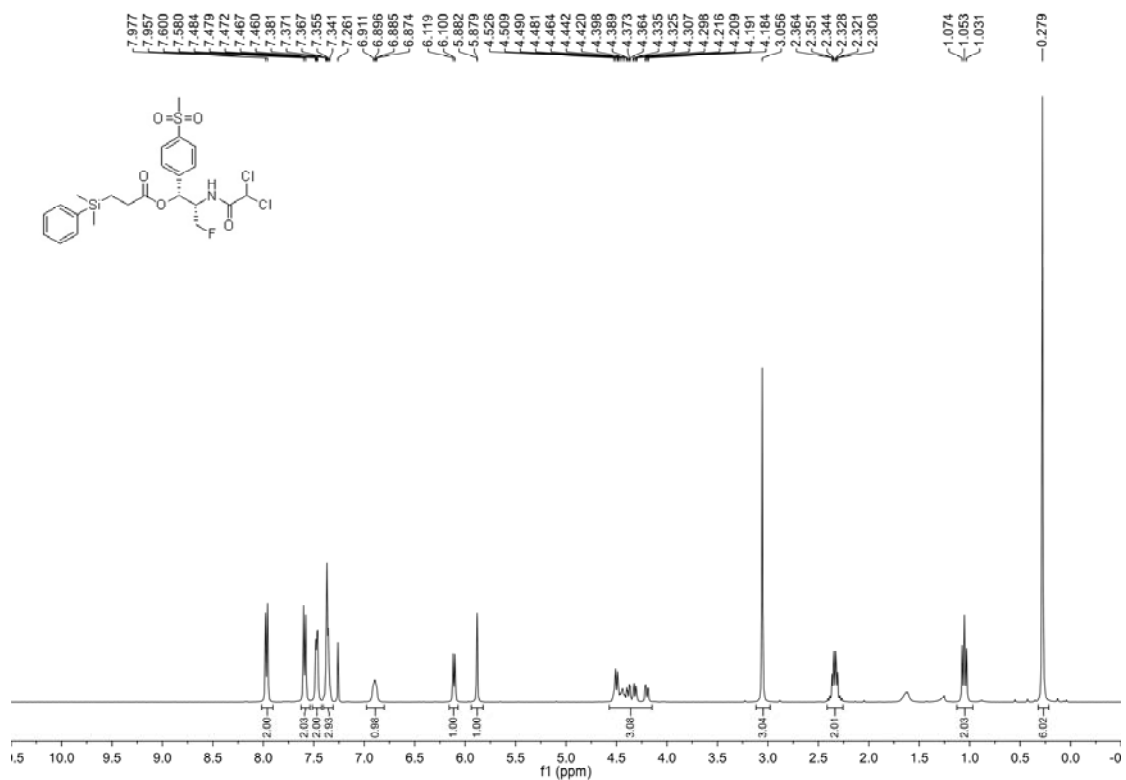


^1H NMR (400 MHz, CDCl_3) of compound **9I**

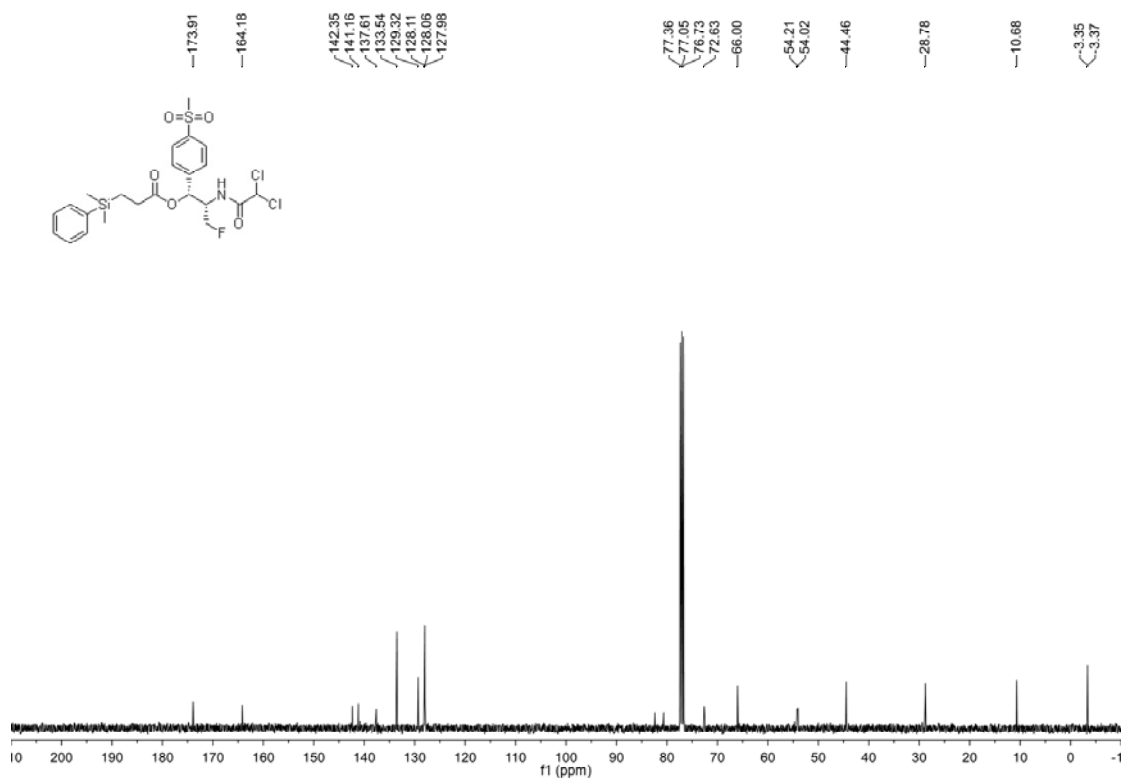


$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **9I**

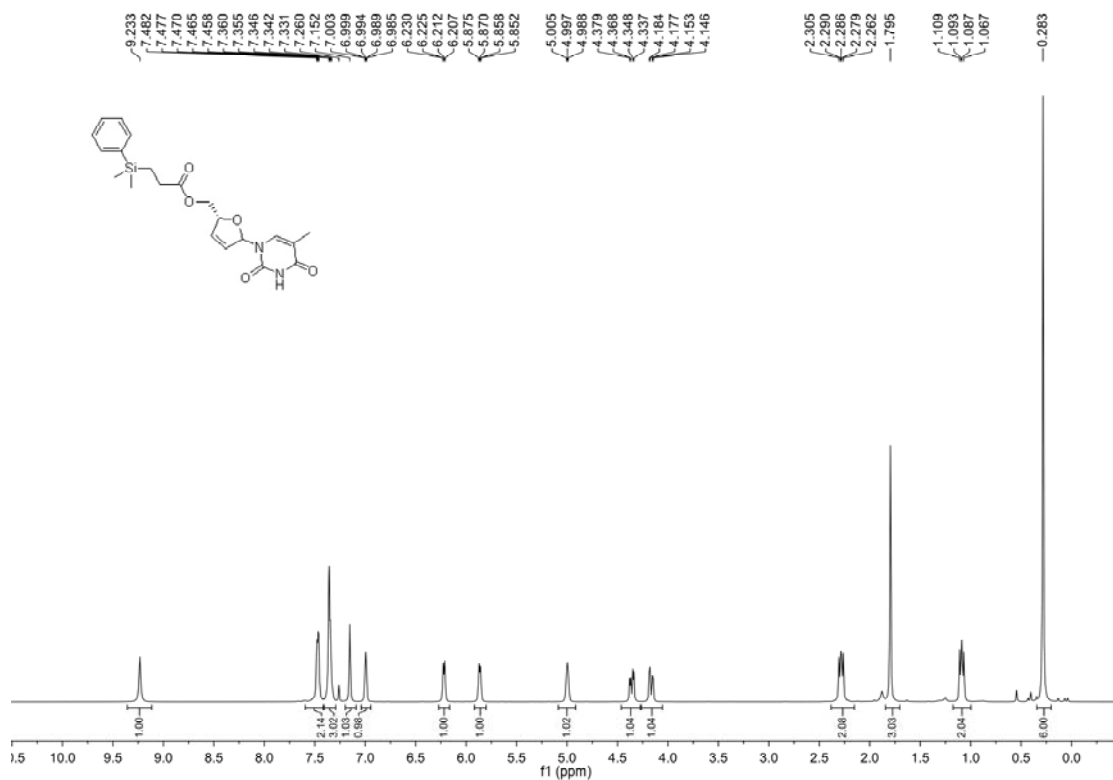




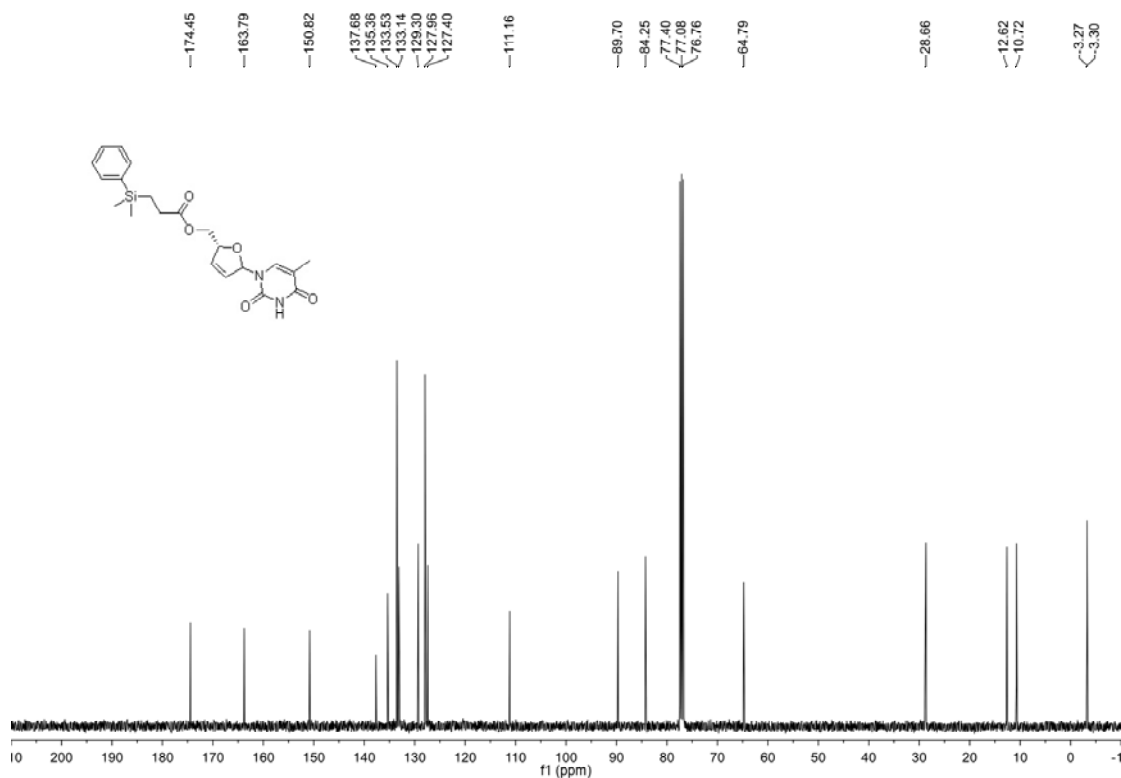
¹H NMR (400 MHz, CDCl₃) of compound **9n**



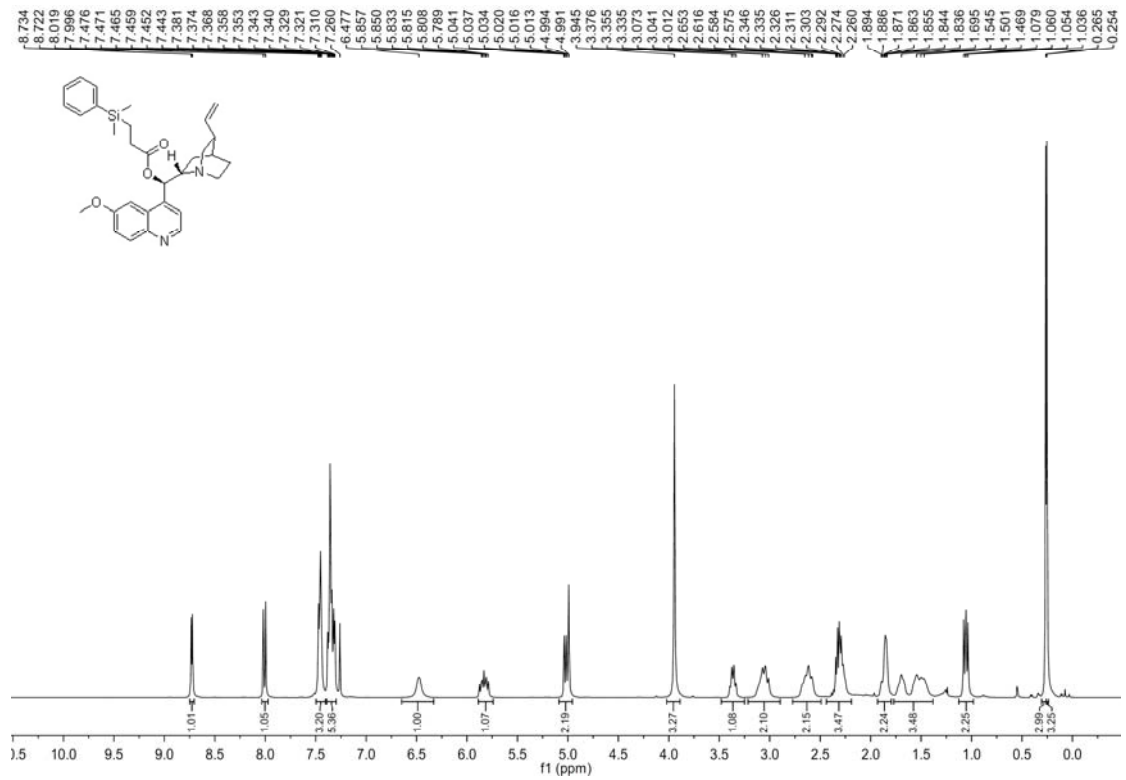
¹³C {¹H} NMR (100 MHz, CDCl₃) of compound **9n**



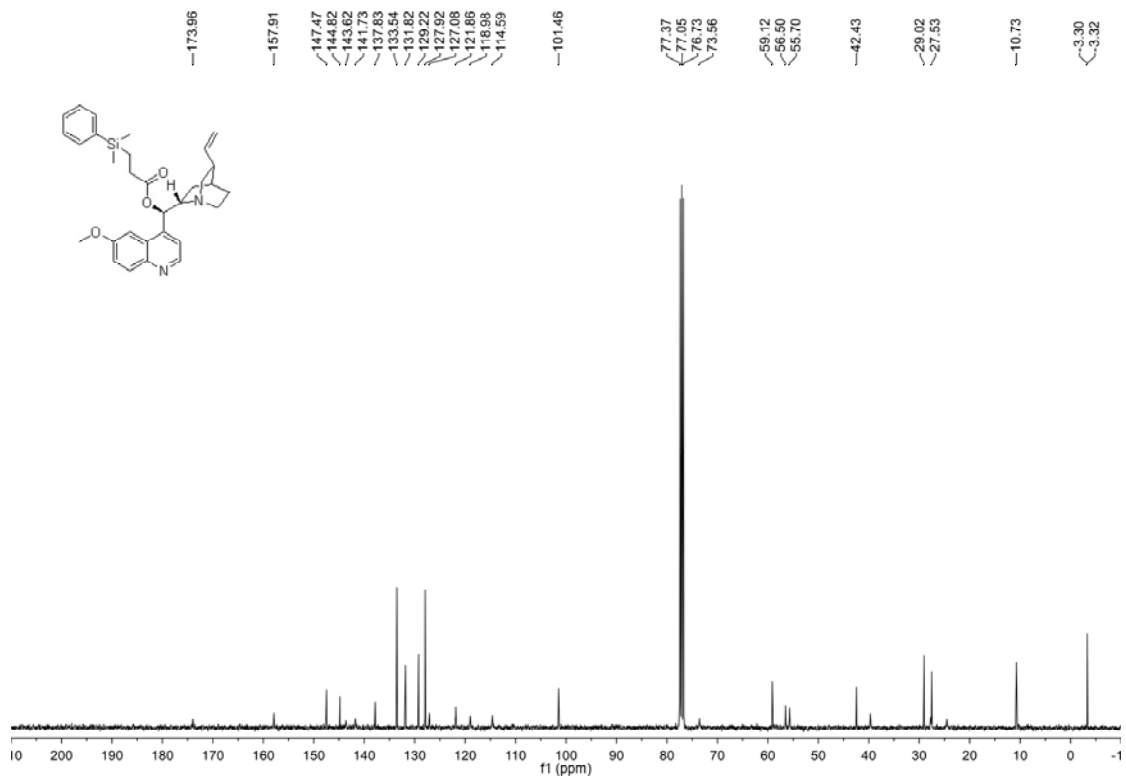
¹H NMR (400 MHz, CDCl₃) of compound **9o**



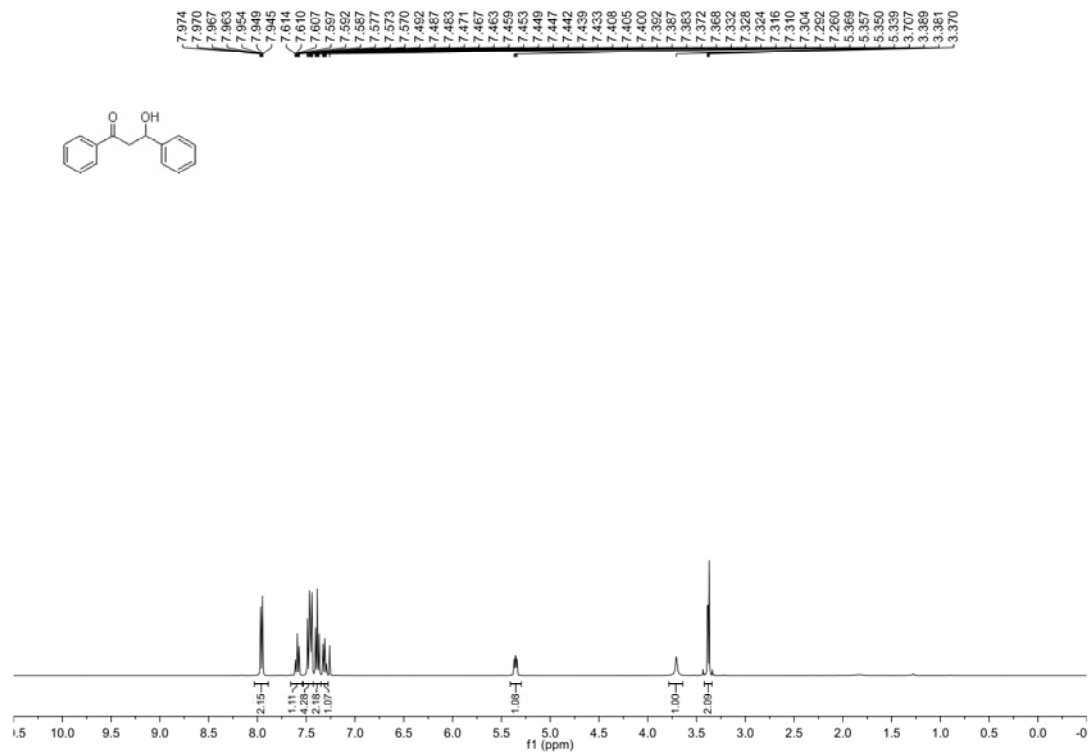
¹³C {¹H} NMR (100 MHz, CDCl₃) of compound **9o**



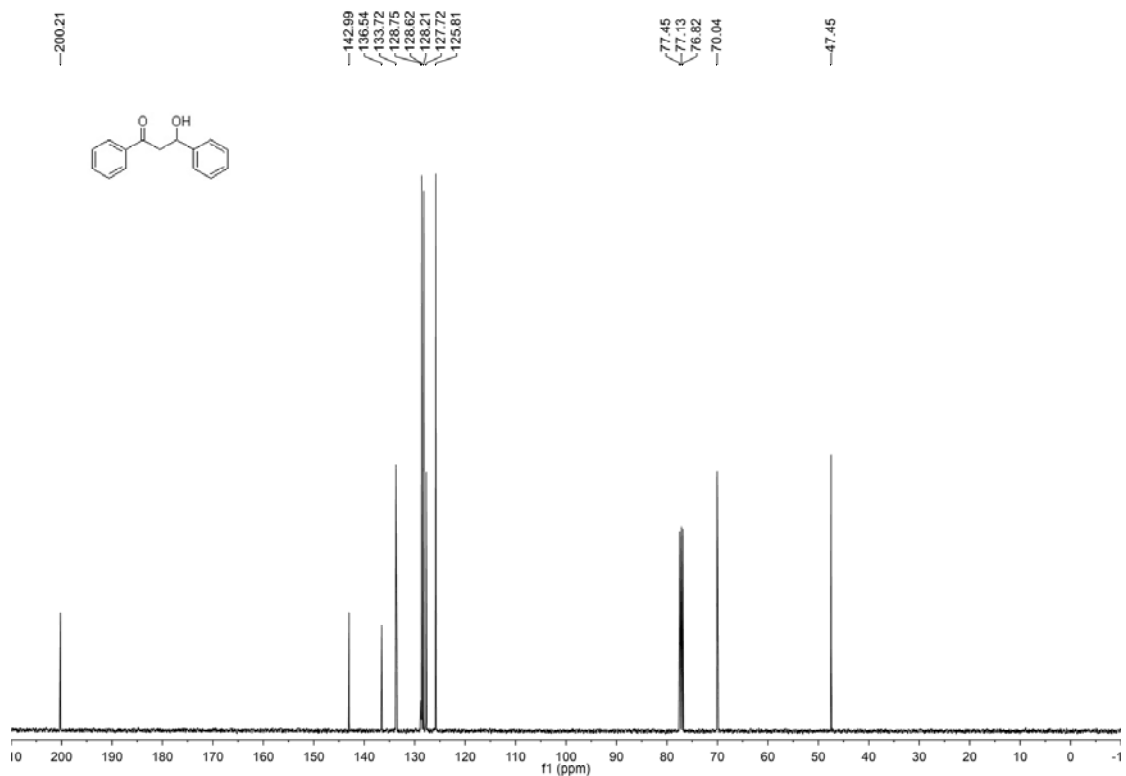
^1H NMR (400 MHz, CDCl_3) of compound **9p**



$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **9p**



^1H NMR (400 MHz, CDCl_3) of compound **10**



$^{13}\text{C}\{^1\text{H}\}$ NMR (100 MHz, CDCl_3) of compound **10**

