

Supporting Information

Electrochemical Diselenylation of Indolizines via Intermolecular C-Se Formation of 2-Methylpyridines, α -Bromoketones with Diselenides

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A. General Methods

¹H and ¹³C NMR spectra were recorded using a Bruker DRX-400 spectrometer using CDCl₃ as solvent. The chemical shifts are referenced to signals at 7.26 and 77.0 ppm, respectively. The data of HRMS was carried out on a high-resolution mass spectrometer (LCMS-IT-TOF). IR spectra were obtained either as potassium bromide pellets or as liquid films between two potassium bromide pellets with a Bruker TENSOR 27 spectrometer. Melting points were determined with a Büchi Melting Point B-545 instrument. The instrument for electrolysis is dual display potentiostat (CHI 660E) (made in China). The anode electrode is vitreous carbon plate (10 mm×10 mm×1 mm) and cathodic electrode was platinum plate (10 mm×10 mm×0.1 mm). Unless otherwise noted, materials were obtained from commercial suppliers and used without further purification.

B. General Procedure for the Preparation of 4 and 5

2-Bromoacetophenone (0.8 mmol, 1.6 equiv), 2-methylpyridine (0.5 mmol, 1 equiv), diphenyl diselenide (0.5 mmol 1 equiv), KI (1 mmol), K_2CO_3 (0.5 mmol), DMF (6 mL) and H_2O (5 mL) were placed in a 20 mL undivided electrolytic cell with a vitreous carbon plate anode (10 mm×10 mm×1 mm) and a platinum plate cathode (10 mm×10 mm×0.1 mm). The electrolysis was carried out at 50 °C under a constant current of 10 mA for 8 hours (monitored by TLC). When the reaction was finished, the resulting reaction solution was quenched with 100 mL brine and extracted with 4×60 mL ethylacetate. The extract was dried with Na_2SO_4 . The solvent was removed with a rotary evaporator. The pure product was obtained by flash column chromatography on silica gel.

C. General procedure for cyclic voltammetry

Cyclic voltammetry was performed in a three-electrode cell connected to a schlenk line under nitrogen at room temperature. The working electrode was a steady vitreous carbon plate electrode, the counter electrode was a platinum plate. The reference was an Ag/AgCl electrode submerged in saturated aqueous KCl solution, and separated from reaction by a salt bridge. Then 10mL electrolyte solution containing 0.05 M *n*-Bu₄NPF₆ in the mixed solvent of DMF and H₂O was poured into electrochemical cell. The concentration of samples was 0.01 M. The scan rate was 0.1 V/s, ranging from 0 V to 2.0 V.

D. Cyclic voltammogram

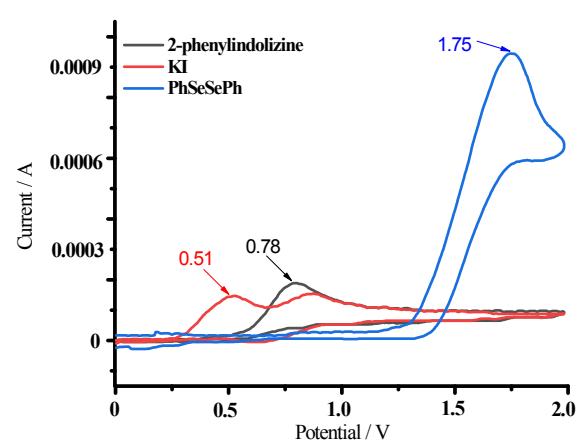
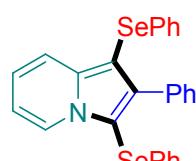


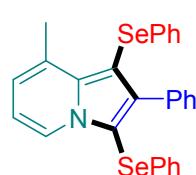
Figure S1 Cyclic voltammetry of 2-phenylindolizine **6aa**, KI and diphenyl diselenide **3a** in the mixed solvent of DMF and H₂O.

E. Analytical data

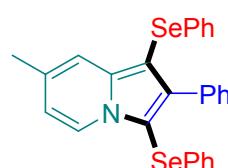
2-phenyl-1,3-bis(phenylselanyl)indolizine (4aa)

 Yellow solid (194.4 mg, 77%), mp 136.2-138.5 °C; IR (KBr): 3055, 2921, 1475, 1344, 747, 687 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.42 (d, *J* = 7.0 Hz, 1H), 7.74 (d, *J* = 8.9 Hz, 1H), 7.44-7.40 (m, 2H), 7.33-7.38 (m, 3H), 7.22-7.12 (m, 8H), 7.06-6.98 (m, 3H), 6.73 (t, *J* = 6.8 Hz, 1H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 142.4, 140.0, 134.9, 134.6, 132.2, 130.8, 129.4, 128.9, 128.1, 127.9, 127.5, 127.4, 126.2, 125.7, 125.4, 121.7, 118.8, 112.5, 105.5, 94.7. ⁷⁷Se NMR (76 MHz, Chloroform-d) δ ppm 227.8, 205.9. HRMS MALDI (m/z): calcd for C₂₆H₁₉NSe₂ [M + H]⁺: 505.9922, found: 505.9926.

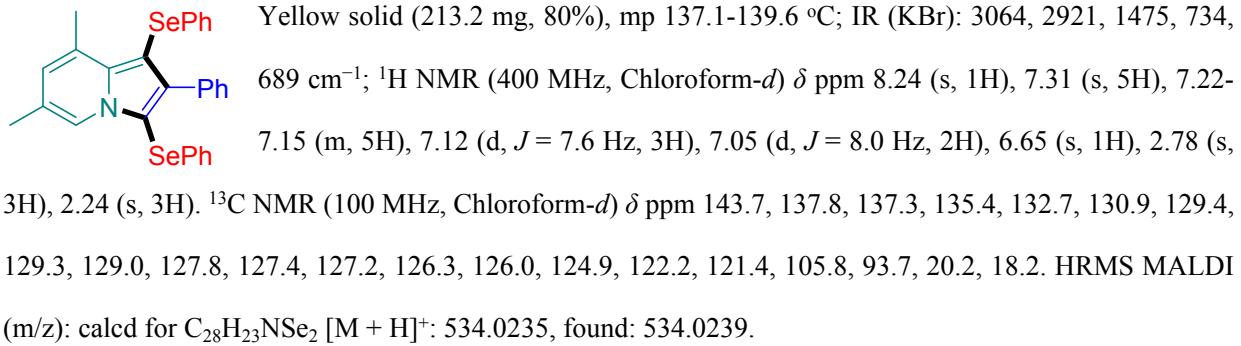
8-methyl-2-phenyl-1,3-bis(phenylselanyl)indolizine (4ab)

 Brown solid (199.8 mg, 77%), mp 92.3-95.8 °C; IR (KBr): 3067, 2917, 1578, 1476, 734, 691 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.34 (d, *J* = 6.9 Hz, 1H), 7.26 (t, *J* = 2.0 Hz, 5H), 7.14-7.03 (m, 8H), 6.98 (d, *J* = 8.1 Hz, 2H), 6.68 (d, *J* = 6.7 Hz, 1H), 6.54 (t, *J* = 6.9 Hz, 1H), 2.75 (s, 3H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 144.0, 138.4, 137.6, 135.3, 132.3, 130.9, 129.9, 129.4, 128.9, 128.0, 127.4, 127.3, 127.3, 126.1, 125.0, 124.5, 122.8, 111.9, 106.3, 94.2, 20.4. HRMS MALDI (m/z): calcd for C₂₇H₂₁NSe₂ [M + H]⁺: 520.0079, found: 520.0082.

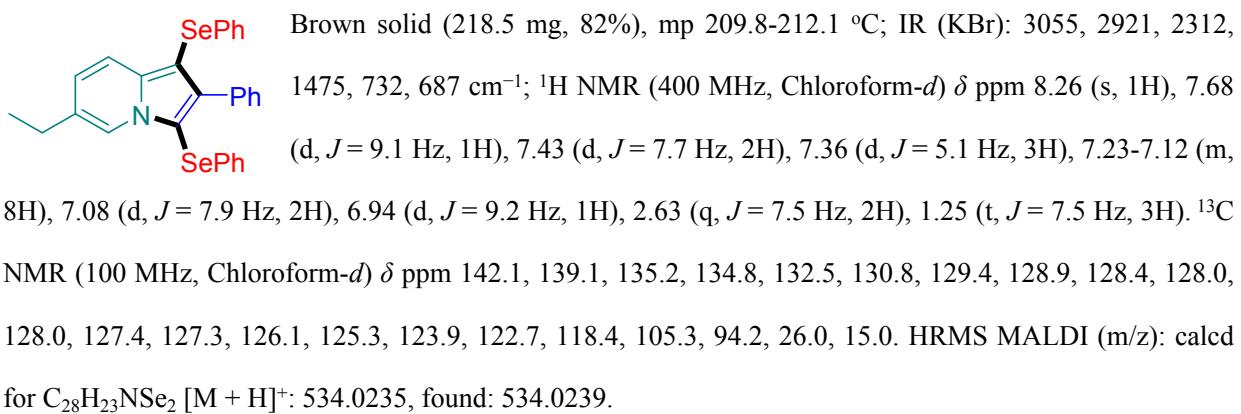
7-methyl-2-phenyl-1,3-bis(phenylselanyl)indolizine (4ac)

 Brown solid (210.2 mg, 81%), mp 185.3-187.4 °C; IR (KBr): 3054, 2922, 2382, 1508, 735, 686 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.28 (d, *J* = 7.5 Hz, 1H), 7.47 (s, 1H), 7.38-7.34 (m, 2H), 7.33-7.29 (m, 3H), 7.19-7.09 (m, 8H), 7.02 (d, *J* = 8.0 Hz, 2H), 6.55 (d, *J* = 7.2 Hz, 1H), 2.35 (s, 3H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 142.6, 140.5, 135.3, 134.8, 132.6, 132.6, 130.8, 129.4, 129.0, 127.9, 127.9, 127.5, 127.5, 127.4, 126.1, 125.3, 125.3, 117.1, 115.2, 104.6, 93.0, 21.1. HRMS MALDI (m/z): calcd for C₂₇H₂₁NSe₂ [M + H]⁺: 505.0079, found: 505.0082.

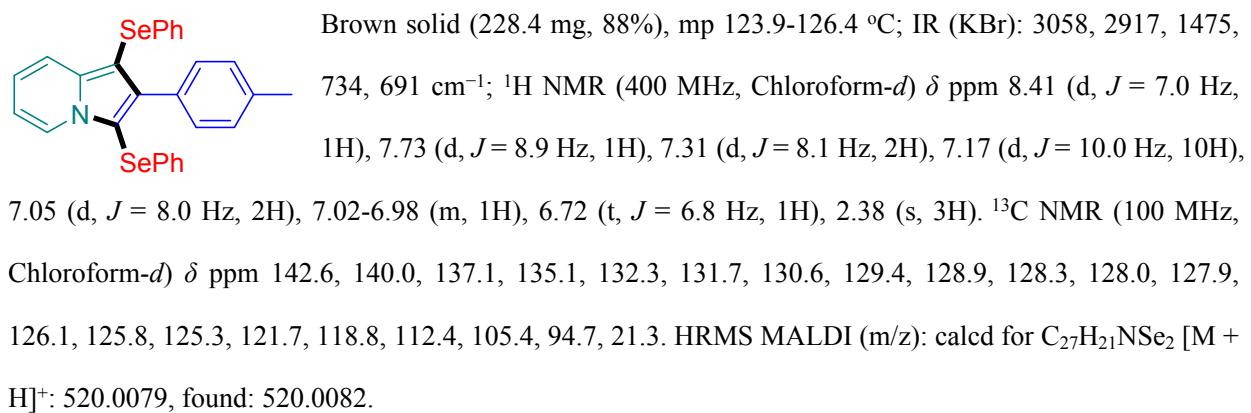
6,8-dimethyl-2-phenyl-1,3-bis(phenylselanyl)indolizine (4ad)



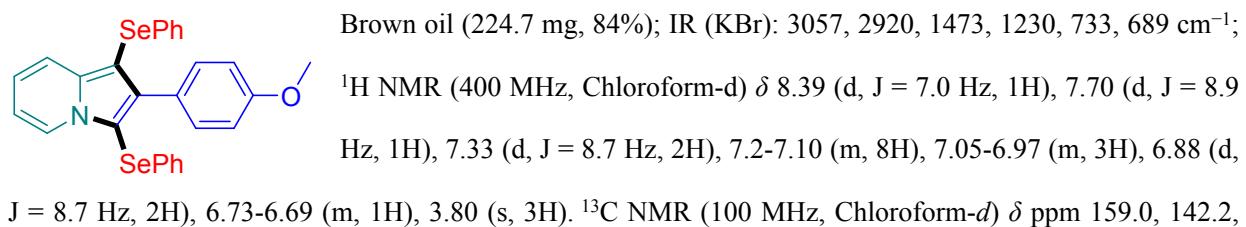
6-ethyl-2-phenyl-1,3-bis(phenylselanyl)indolizine (4ae)



1,3-bis(phenylselanyl)-2-(p-tolyl)indolizine (4ba)

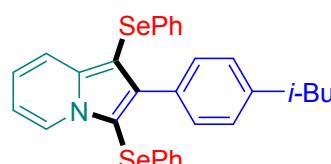


2-(4-methoxyphenyl)-1,3-bis(phenylselanyl)indolizine (4ca)



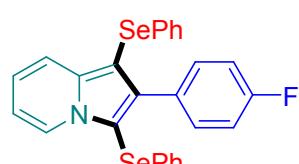
140.1, 135.1, 132.4, 131.9, 129.5, 129.0, 128.0, 127.9, 127.0, 126.2, 125.8, 125.4, 121.7, 118.8, 113.1, 112.4, 105.4, 94.7, 55.1. HRMS MALDI (m/z): calcd for $C_{27}H_{21}NOSe_2$ [M + H]⁺: 536.0028, found: 536.0032.

2-(4-isobutylphenyl)-1,3-bis(phenylselanyl)indolizine (4da)



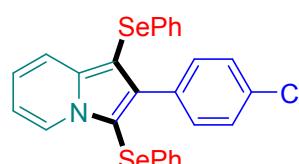
Brown oil (224.4 mg, 80%); IR (KBr): 3057, 2951, 2312, 1474, 731, 689 cm⁻¹; ¹H NMR (400 MHz, Chloroform-d) δ ppm 8.42 (d, *J* = 6.9 Hz, 1H), 7.74 (d, *J* = 8.9 Hz, 1H), 7.37 (d, *J* = 8.0 Hz, 2H), 7.23-7.12 (m, 10H), 7.07 (d, *J* = 7.9 Hz, 2H), 7.02-6.98 (m, 1H), 6.72 (t, *J* = 6.9 Hz, 1H), 2.52 (d, *J* = 7.1 Hz, 2H), 1.88-1.98 (m, 1H), 0.96 (d, *J* = 6.5 Hz, 6H). ¹³C NMR (100 MHz, Chloroform-d) δ ppm 142.3, 140.8, 140.0, 135.0, 132.3, 131.8, 130.5, 129.4, 128.9, 128.3, 128.2, 128.0, 126.1, 125.7, 125.4, 121.6, 118.8, 112.3, 105.4, 94.8, 45.2, 30.0, 22.5. HRMS MALDI (m/z): calcd for $C_{30}H_{27}NSe_2$ [M + H]⁺: 562.0548, found: 562.0553.

2-(4-fluorophenyl)-1,3-bis(phenylselanyl)indolizine (4ea)



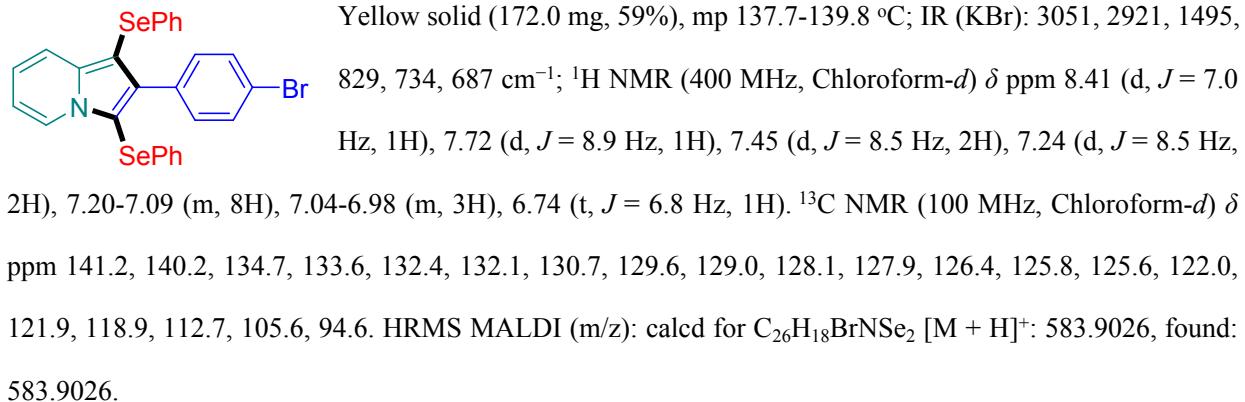
Brown solid (185.7 mg, 71%), mp 152.9-155.7 °C; IR (KBr): 3053, 2921, 1471, 837, 734, 688 cm⁻¹; ¹H NMR (400 MHz, Chloroform-d) δ ppm 8.40 (d, *J* = 7.0 Hz, 1H), 7.71 (d, *J* = 8.9 Hz, 1H), 7.36-7.30 (m, 2H), 7.21-7.09 (m, 8H), 7.04-6.97 (m, 5H), 6.73 (t, *J* = 6.8 Hz, 1H). ¹³C NMR (100 MHz, Chloroform-d) δ ppm 162.4 (*J* = 246.5 Hz), 141.5, 140.1, 134.8, 132.4 (*J* = 8.0 Hz), 132.2, 130.6 (*J* = 3.3 Hz), 129.5, 129.0, 128.1, 127.9, 126.3, 125.8, 125.5, 121.9, 118.9, 114.6 (*J* = 21.5 Hz), 112.6, 105.7, 94.8. HRMS MALDI (m/z): calcd for $C_{26}H_{18}FNSe_2$ [M + H]⁺: 523.9835, found: 523.9831.

2-(4-chlorophenyl)-1,3-bis(phenylselanyl)indolizine (4fa)

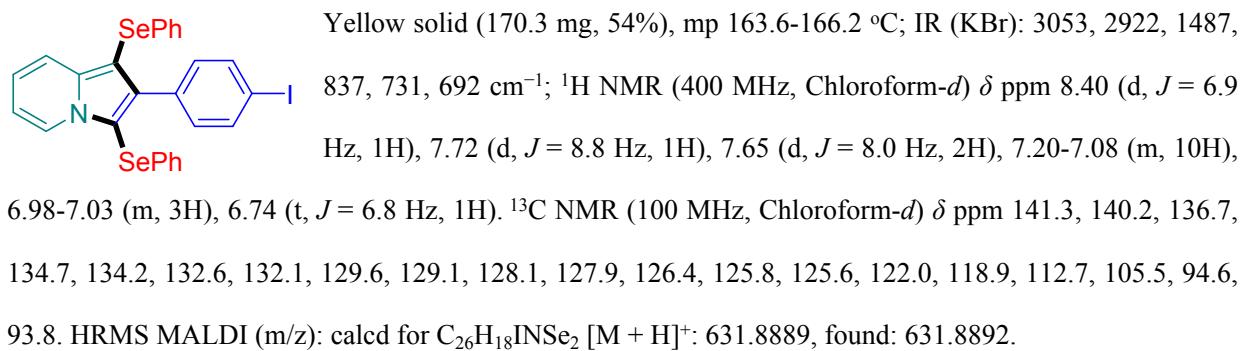


Yellow solid (164.4 mg, 61%), mp 139.7-142.1 °C; IR (KBr): 3067, 2925, 1476, 833, 734, 691 cm⁻¹; ¹H NMR (400 MHz, Chloroform-d) δ ppm 8.31 (d, *J* = 7.1 Hz, 1H), 7.62 (d, *J* = 8.9 Hz, 1H), 7.20 (s, 4H), 7.10-6.99 (m, 8H), 6.92 (t, *J* = 7.5 Hz, 3H), 6.64 (t, *J* = 7.2 Hz, 1H). ¹³C NMR (100 MHz, Chloroform-d) δ ppm 141.2, 140.2, 134.7, 133.5, 133.1, 132.1, 129.6, 129.0, 128.1, 127.9, 127.8, 126.3, 125.8, 125.6, 122.0, 118.9, 112.7, 105.6, 94.7. HRMS MALDI (m/z): calcd for $C_{26}H_{18}ClNSe_2$ [M + H]⁺: 539.9529, found: 539.9532.

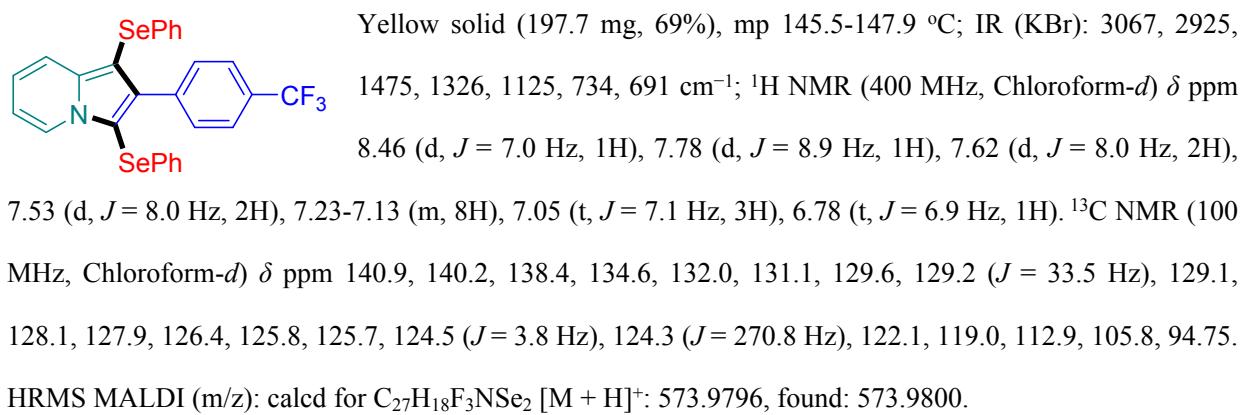
2-(4-bromophenyl)-1,3-bis(phenylselanyl)indolizine (4ga)



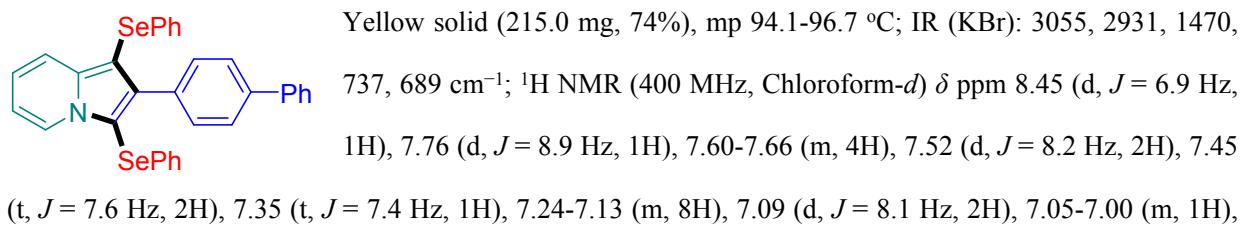
2-(4-iodophenyl)-1,3-bis(phenylselanyl)indolizine (4ha)



1,3-bis(phenylselanyl)-2-(4-(trifluoromethyl)phenyl)indolizine (4ia)

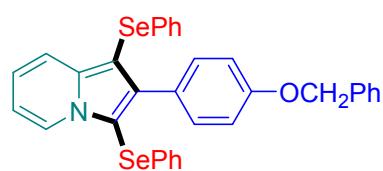


2-([1,1'-biphenyl]-4-yl)-1,3-bis(phenylselanyl)indolizine (4ja)



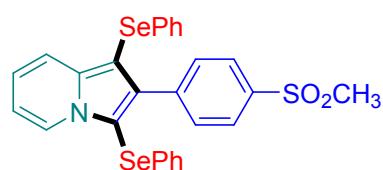
6.75 (t, $J = 6.8$ Hz, 1H). ^{13}C NMR (100 MHz, Chloroform-*d*) δ ppm 142.0, 140.8, 140.2, 140.0, 135.0, 133.6, 132.3, 131.2, 129.5, 129.0, 128.6, 128.1, 127.9, 127.1, 127.0, 126.2, 125.8, 125.4, 121.8, 118.8, 112.5, 105.5, 94.7. HRMS MALDI (m/z): calcd for $\text{C}_{32}\text{H}_{23}\text{NSe}_2$ [M + H] $^+$: 582.0235, found: 582.0240.

2-(4-(benzyloxy)phenyl)-1,3-bis(phenylselanyl)indolizine (4ka)



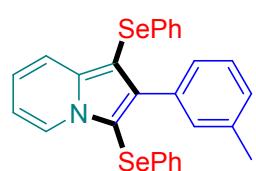
Brown oil (235.2 mg, 77%); IR (KBr): 3064, 2924, 1475, 1238, 733, 689 cm $^{-1}$; ^1H NMR (400 MHz, Chloroform-*d*) δ ppm 8.44 (d, $J = 7.0$ Hz, 1H), 7.76 (d, $J = 8.9$ Hz, 1H), 7.49 (d, $J = 7.2$ Hz, 2H), 7.36-7.46 (m, 5H), 7.25-7.14 (m, 8H), 7.09 (d, $J = 7.9$ Hz, 2H), 7.02 (d, $J = 8.7$ Hz, 3H), 6.74 (t, $J = 6.8$ Hz, 1H), 5.09 (s, 2H). ^{13}C NMR (100 MHz, Chloroform-*d*) δ ppm 158.3, 142.1, 140.0, 136.9, 135.0, 132.3, 131.9, 129.4, 128.9, 128.5, 128.0, 127.8, 127.5, 127.2, 126.1, 125.7, 125.4, 121.7, 118.7, 113.9, 112.4, 105.3, 94.6, 69.8. HRMS MALDI (m/z): calcd for $\text{C}_{33}\text{H}_{25}\text{NOSe}_2$ [M + H] $^+$: 612.0341, found: 612.0346.

2-(4-(methylsulfonyl)phenyl)-1,3-bis(phenylselanyl)indolizine (4la)



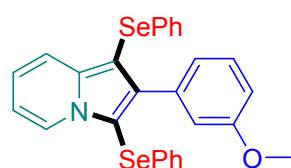
Brown oil (195.3 mg, 67%); IR (KBr): 3067, 2924, 1475, 1311, 1145, 734, 688, 544 cm $^{-1}$; ^1H NMR (400 MHz, Chloroform-*d*) δ ppm 8.44 (d, $J = 7.0$ Hz, 1H), 7.88 (d, $J = 8.3$ Hz, 2H), 7.75 (d, $J = 8.9$ Hz, 1H), 7.57 (d, $J = 8.4$ Hz, 2H), 7.21-7.08 (m, 8H), 7.04 (d, $J = 9.0$ Hz, 1H), 7.02-6.97 (m, 2H), 6.78 (t, $J = 6.8$ Hz, 1H), 3.06 (s, 3H). ^{13}C NMR (100 MHz, Chloroform-*d*) δ ppm 140.5, 140.3, 140.1, 138.9, 134.4, 131.8, 131.7, 129.6, 129.1, 128.0, 127.8, 126.5, 126.5, 125.8, 125.7, 122.3, 119.0, 113.1, 106.0, 94.7, 44.4. HRMS MALDI (m/z): calcd for $\text{C}_{27}\text{H}_{21}\text{NO}_2\text{SSe}_2$ [M + H] $^+$: 583.9697, found: 583.9700.

1,3-bis(phenylselanyl)-2-(m-tolyl)indolizine (4ma)



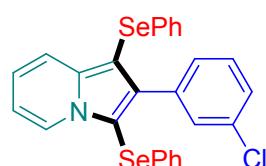
Brown oil (210.2 mg, 81%); IR (KBr): 3057, 2912, 1473, 736, 694 cm $^{-1}$; ^1H NMR (400 MHz, Chloroform-*d*) δ ppm 8.41 (d, $J = 7.0$ Hz, 1H), 7.74 (d, $J = 8.9$ Hz, 1H), 7.24-7.11 (m, 12H), 7.05 (d, $J = 8.0$ Hz, 2H), 7.00-6.95 (m, 1H), 6.69 (t, $J = 6.8$ Hz, 1H), 2.29 (s, 3H). ^{13}C NMR (100 MHz, Chloroform-*d*) δ ppm 142.6, 140.0, 136.8, 135.1, 134.5, 132.4, 131.7, 129.4, 128.9, 128.3, 128.2, 128.0, 127.8, 127.3, 126.2, 125.7, 125.4, 121.6, 118.8, 112.4, 105.6, 94.9, 21.4. HRMS MALDI (m/z): calcd for $\text{C}_{27}\text{H}_{21}\text{NSe}_2$ [M + H] $^+$: 520.0079, found: 520.0082.

2-(3-methoxyphenyl)-1,3-bis(phenylselanyl)indolizine (4na)



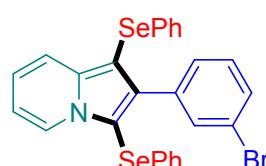
Brown oil (214.0 mg, 80%); IR (KBr): 3067, 2921, 1475, 736, 688 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.47 (d, *J* = 7.0 Hz, 1H), 7.79 (d, *J* = 8.9 Hz, 1H), 7.29 (t, *J* = 7.9 Hz, 1H), 7.25-7.18 (m, 7H), 7.17-7.13 (m, 1H), 7.11-7.01 (m, 4H), 6.96 (s, 1H), 6.93 (d, *J* = 8.1 Hz, 1H), 6.75 (t, *J* = 6.8 Hz, 1H), 3.54 (s, 3H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 158.6, 142.3, 140.2, 135.9, 135.2, 132.5, 129.4, 128.9, 128.4, 127.9, 127.8, 126.1, 125.7, 125.3, 123.1, 121.8, 118.8, 115.6, 113.9, 112.5, 105.6, 94.5, 54.7. HRMS MALDI (m/z): calcd for C₂₇H₂₁NOSe₂ [M + H]⁺: 536.0028, found: 536.0032.

2-(3-chlorophenyl)-1,3-bis(phenylselanyl)indolizine (4oa)



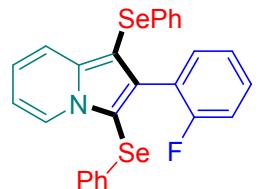
Brown oil (207.5 mg, 77%); IR (KBr): 3052, 2920, 1475, 731, 691 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.47 (d, *J* = 6.5 Hz, 1H), 7.80 (d, *J* = 9.3 Hz, 1H), 7.43 (d, *J* = 12.5 Hz, 1H), 7.36-7.27 (m, 3H), 7.25-7.13 (m, 8H), 7.04 (dd, *J* = 13.4, 5.3 Hz, 3H), 6.75 (t, *J* = 6.8 Hz, 1H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 140.8, 140.0, 136.5, 134.6, 133.1, 131.9, 130.9, 129.5, 129.0, 128.9, 128.7, 128.3, 128.1, 127.4, 126.4, 125.7, 125.6, 121.9, 118.9, 112.7, 105.8, 95.0. HRMS MALDI (m/z): calcd for C₂₆H₁₈ClNSe₂ [M + H]⁺: 539.9529, found: 539.9531.

2-(3-bromophenyl)-1,3-bis(phenylselanyl)indolizine (4pa)



Brown solid (215.7 mg, 74%), mp 96.9-99.2 °C; IR (KBr): 3055, 2926, 1473, 734, 691 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.44 (d, *J* = 7.1 Hz, 1H), 7.77 (d, *J* = 8.7 Hz, 1H), 7.53 (s, 1H), 7.46 (d, *J* = 8.1 Hz, 1H), 7.34 (d, *J* = 7.6 Hz, 1H), 7.23-7.14 (m, 9H), 7.06-7.00 (m, 3H), 6.75 (t, *J* = 6.8 Hz, 1H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 140.7, 140.0, 136.7, 134.6, 133.8, 132.0, 130.3, 129.5, 129.4, 129.0, 129.0, 128.4, 128.1, 126.4, 125.7, 125.6, 121.9, 121.4, 118.9, 112.7, 105.9, 95.0. HRMS MALDI (m/z): calcd for C₂₆H₁₈BrNSe₂ [M + H]⁺: 583.9021, found: 583.9026.

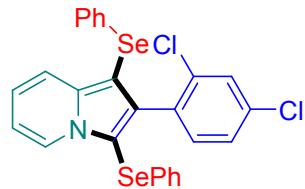
2-(2-fluorophenyl)-1,3-bis(phenylselanyl)indolizine (4qa)



Green solid (185.7 mg, 71%), mp 107.9-110.4 °C; IR (KBr): 3054, 2921, 1475, 741, 687 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.37 (d, *J* = 7.1 Hz, 1H), 7.72 (d, *J* = 8.9 Hz, 1H), 7.38-7.33 (m, 1H), 7.29 (t, *J* = 7.5 Hz, 1H), 7.20-7.11 (m, 10H),

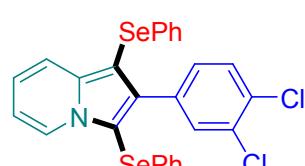
7.07 (d, $J = 7.8$ Hz, 2H), 7.02-6.97 (m, 1H), 6.72 (t, $J = 6.8$ Hz, 1H). ^{13}C NMR (100 MHz, Chloroform-*d*) δ ppm 161.3 ($J = 246.5$ Hz), 139.6, 137.0, 134.3, 133.0 ($J = 2.9$ Hz), 131.5, 129.6 ($J = 8.1$ Hz), 129.3, 128.8, 128.4, 128.3, 126.3, 125.6, 125.5, 123.3 ($J = 3.6$ Hz), 122.8 ($J = 15.4$ Hz), 121.5, 118.8, 115.4 ($J = 22.3$ Hz), 112.5, 106.7, 95.9. HRMS MALDI (m/z): calcd for $\text{C}_{26}\text{H}_{18}\text{FNSe}_2$ [M + H] $^+$: 523.9828, found: 523.9831.

2-(2,4-dichlorophenyl)-1,3-bis(phenylselanyl)indolizine (4ra)



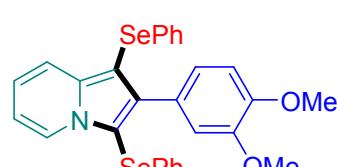
Brown oil (209.1 mg, 73%); IR (KBr): 3069, 2923, 1461, 732, 691 cm^{-1} ; ^1H NMR (400 MHz, Chloroform-*d*) δ ppm 8.39 (d, $J = 7.1$ Hz, 1H), 7.74 (d, $J = 9.0$ Hz, 1H), 7.51 (d, $J = 1.9$ Hz, 1H), 7.24-7.13 (m, 10H), 7.10-7.05 (m, 2H), 7.04-6.99 (m, 1H), 6.74 (t, $J = 6.6$ Hz, 1H). ^{13}C NMR (100 MHz, Chloroform-*d*) δ ppm 139.3, 139.2, 135.4, 134.2, 133.9, 133.5, 132.9, 131.2, 129.3, 129.0, 128.8, 128.6, 128.5, 126.4, 126.3, 125.6, 125.6, 121.6, 118.8, 112.6, 106.6, 95.9. HRMS MALDI (m/z): calcd for $\text{C}_{26}\text{H}_{17}\text{Cl}_2\text{NSe}_2$ [M + H] $^+$: 573.9134, found: 573.9138.

2-(3,4-dichlorophenyl)-1,3-bis(phenylselanyl)indolizine (4sa)



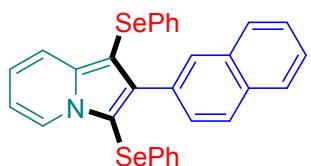
Brown oil (223.4 mg, 78%), IR (KBr): 3052, 2921, 1469, 731, 696 cm^{-1} ; ^1H NMR (400 MHz, Chloroform-*d*) δ ppm 8.44 (d, $J = 7.0$ Hz, 1H), 7.76 (d, $J = 8.9$ Hz, 1H), 7.45 (d, $J = 2.0$ Hz, 1H), 7.39 (d, $J = 8.3$ Hz, 1H), 7.22-7.12 (m, 9H), 7.06-7.00 (m, 3H), 6.76 (t, $J = 6.8$ Hz, 1H). ^{13}C NMR (100 MHz, Chloroform-*d*) δ ppm 140.2, 139.8, 134.7, 134.4, 132.7, 131.9, 131.6, 131.5, 130.1, 129.6, 129.5, 129.1, 128.4, 128.1, 126.5, 125.8, 125.8, 122.1, 119.0, 112.9, 105.9, 94.9. HRMS MALDI (m/z): calcd for $\text{C}_{26}\text{H}_{17}\text{Cl}_2\text{NSe}_2$ [M + H] $^+$: 573.9134, found: 573.9138.

2-(3,4-dimethoxyphenyl)-1,3-bis(phenylselanyl)indolizine (4ta)



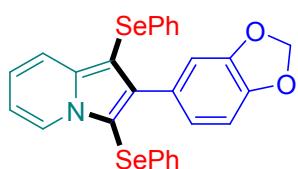
Brown solid (231.7 mg, 82%), mp 155.9-159.0 $^\circ\text{C}$; IR (KBr): 3057, 2924, 1476, 1233, 731, 692 cm^{-1} ; ^1H NMR (400 MHz, Chloroform-*d*) δ ppm 8.46 (d, $J = 7.0$ Hz, 1H), 7.75 (d, $J = 8.9$ Hz, 1H), 7.22-7.11 (m, 8H), 7.05 (d, $J = 6.5$ Hz, 2H), 7.01 (d, $J = 9.0$ Hz, 1H), 6.97 (d, $J = 8.2$ Hz, 1H), 6.84 (d, $J = 8.3$ Hz, 2H), 6.74 (t, $J = 6.2$ Hz, 1H), 3.87 (s, 3H), 3.40 (s, 3H). ^{13}C NMR (100 MHz, Chloroform-*d*) δ ppm 148.3, 147.8, 142.2, 140.5, 135.5, 132.8, 129.5, 129.0, 127.7, 127.6, 127.2, 126.1, 125.8, 125.3, 122.7, 121.9, 118.7, 114.0, 112.5, 110.2, 105.4, 94.3, 55.6, 55.0. HRMS MALDI (m/z): calcd for $\text{C}_{28}\text{H}_{23}\text{NO}_2\text{Se}_2$ [M + H] $^+$: 566.0133, found: 566.0138.

2-(naphthalen-2-yl)-1,3-bis(phenylselanyl)indolizine (**4ua**)



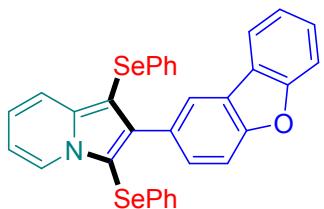
Brown solid (185.9 mg, 67%), mp 113.9-115.4 °C; IR (KBr): 3067, 2912, 1475, 733, 689 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.43 (d, *J* = 7.0 Hz, 1H), 7.87 (d, *J* = 8.2 Hz, 2H), 7.76 (d, *J* = 8.9 Hz, 1H), 7.55 (d, *J* = 8.5 Hz, 1H), 7.46 (d, *J* = 7.1 Hz, 1H), 7.41 (d, *J* = 7.2 Hz, 1H), 7.32 (d, *J* = 6.5 Hz, 1H), 7.25-7.21 (m, 1H), 7.15-7.11 (m, 3H), 7.07 (s, 5H), 7.02 (t, *J* = 9.8 Hz, 3H), 6.76 (t, *J* = 6.6 Hz, 1H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 141.3, 139.5, 134.2, 133.2, 133.1, 133.0, 131.7, 129.2, 128.9, 128.8, 128.7, 128.6, 128.1, 128.0, 126.4, 126.3, 125.7, 125.6, 125.5, 125.4, 124.8, 121.5, 118.7, 112.4, 107.3, 97.1. HRMS MALDI (m/z): calcd for C₃₀H₂₁NSe₂ [M + H]⁺: 556.0079, found: 556.0083.

2-(benzo[d][1,3]dioxol-5-yl)-1,3-bis(phenylselanyl)indolizine (**4va**)



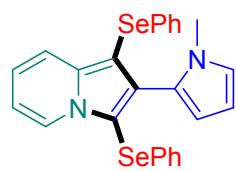
Brown oil (208.6 mg, 76%); IR (KBr): 3053, 2931, 1453, 737, 682 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.41 (d, *J* = 7.0 Hz, 1H), 7.73 (d, *J* = 8.9 Hz, 1H), 7.22-7.11 (m, 8H), 7.06-6.98 (m, 3H), 6.93-6.87 (m, 2H), 6.81 (d, *J* = 7.9 Hz, 1H), 6.72 (t, *J* = 6.8 Hz, 1H), 5.95 (s, 2H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 147.0, 146.8, 142.1, 139.9, 134.8, 132.2, 129.5, 129.0, 128.3, 128.0, 127.9, 126.2, 125.7, 125.4, 124.5, 121.8, 118.8, 112.5, 111.2, 107.6, 105.5, 100.9, 94.7. HRMS MALDI (m/z): calcd for C₂₇H₁₉NO₂Se₂ [M + H]⁺: 549.9820, found: 549.9824.

2-(dibenzo[b,d]furan-2-yl)-1,3-bis(phenylselanyl)indolizine (**4wa**)



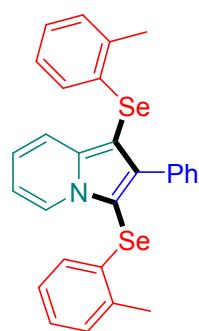
Brown oil (229.1 mg, 77%); IR (KBr): 3061, 2924, 1462, 735, 683 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.49 (d, *J* = 7.0 Hz, 1H), 7.90-7.85 (m, 1H), 7.80 (d, *J* = 8.9 Hz, 1H), 7.67 (d, *J* = 7.5 Hz, 1H), 7.58-7.53 (m, 2H), 7.50 (d, *J* = 8.5 Hz, 1H), 7.46-7.41 (m, 1H), 7.29 (t, *J* = 7.1 Hz, 1H), 7.21 (d, *J* = 3.1 Hz, 8H), 7.09-7.03 (m, 3H), 6.77 (t, *J* = 6.2 Hz, 1H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 156.4, 155.7, 142.3, 140.2, 135.2, 132.5, 131.4, 130.0, 129.5, 129.0, 128.4, 128.1, 126.9, 126.3, 125.8, 125.6, 124.3, 123.6, 123.2, 122.5, 121.9, 120.6, 118.9, 112.6, 111.5, 110.7, 106.0, 95.3. HRMS MALDI (m/z): calcd for C₃₂H₂₁NOSe₂ [M + H]⁺: 596.0028, found: 596.0033.

2-(1-methyl-1H-pyrrol-2-yl)-1,3-bis(phenylselanyl)indolizine (4xa)



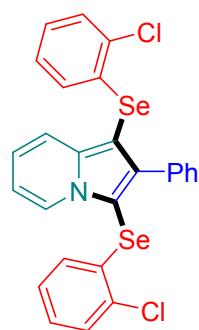
Brown oil (147.3 mg, 58%); IR (KBr): 3068, 2920, 1544, 1020, 735, 685 cm^{-1} ; ^1H NMR (400 MHz, Chloroform-*d*) δ ppm 8.35 (d, $J = 7.0$ Hz, 1H), 7.68 (d, $J = 8.9$ Hz, 1H), 7.20-7.03 (m, 10H), 7.01-6.95 (m, 3H), 6.77-6.73 (m, 1H), 6.10 (d, $J = 1.2$ Hz, 1H), 3.16 (s, 3H). ^{13}C NMR (100 MHz, Chloroform-*d*) δ ppm 139.4, 133.4, 133.0, 131.3, 129.5, 129.4, 128.9, 128.8, 128.1, 126.7, 126.0, 125.7, 121.8, 119.0, 113.6, 112.9, 108.6, 104.5, 97.8, 97.4, 34.6. HRMS MALDI (m/z): calcd for $\text{C}_{25}\text{H}_{20}\text{N}_2\text{Se}_2$ [M + H]⁺: 509.0031, found: 509.0034.

2-phenyl-1,3-bis(o-tolylselanyl)indolizine (5ab)



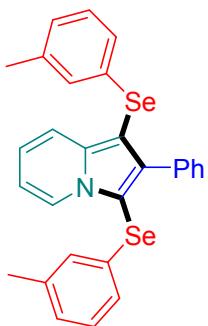
Brown solid (207.9 mg, 78%), mp 115.8-118.3 °C; IR (KBr): 3051, 2923, 1475, 731, 688 cm^{-1} ; ^1H NMR (400 MHz, Chloroform-*d*) δ ppm 8.31 (d, $J = 7.0$ Hz, 1H), 7.66 (d, $J = 8.9$ Hz, 1H), 7.37-7.33 (m, 2H), 7.28-7.32 (m, 3H), 7.17-7.07 (m, 3H), 7.06-7.02 (m, 1H), 7.01-6.96 (m, 1H), 6.92 (t, $J = 7.6$ Hz, 2H), 6.76-6.69 (m, 2H), 6.48 (d, $J = 7.7$ Hz, 1H), 2.39 (d, $J = 2.4$ Hz, 6H). ^{13}C NMR (100 MHz, Chloroform-*d*) δ ppm 142.9, 140.3, 136.3, 136.0, 135.5, 134.6, 132.8, 130.6, 130.4, 129.8, 127.5, 127.4, 127.2, 126.9, 126.6, 126.5, 126.0, 125.9, 125.1, 121.7, 118.9, 112.5, 104.6, 93.9, 21.2, 21.0. HRMS MALDI (m/z): calcd for $\text{C}_{28}\text{H}_{23}\text{NSe}_2$ [M + H]⁺: 534.0235, found: 534.0239.

1,3-bis((2-chlorophenyl)selanyl)-2-phenylindolizine (5ac)



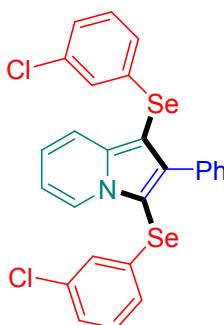
Yellow solid (212.0 mg, 74%), mp 152.9-154.7 °C; IR (KBr): 3062, 2921, 1476, 734, 687 cm^{-1} ; ^1H NMR (400 MHz, Chloroform-*d*) δ ppm 8.35 (d, $J = 6.9$ Hz, 1H), 7.68 (d, $J = 8.9$ Hz, 1H), 7.36-7.28 (m, 7H), 7.13-7.09 (m, 1H), 7.07-7.02 (m, 2H), 6.99 (d, $J = 8.5$ Hz, 1H), 6.95 (d, $J = 7.2$ Hz, 1H), 6.77 (t, $J = 6.8$ Hz, 1H), 6.64 (d, $J = 7.9$ Hz, 1H), 6.39 (d, $J = 7.9$ Hz, 1H). ^{13}C NMR (100 MHz, Chloroform-*d*) δ ppm 143.6, 140.7, 135.1, 134.0, 132.6, 132.5, 132.1, 130.5, 129.8, 129.3, 128.1, 127.7, 127.7, 127.2, 127.2, 126.3, 126.0, 122.4, 118.9, 113.0, 104.7, 93.7. HRMS MALDI (m/z): calcd for $\text{C}_{26}\text{H}_{17}\text{Cl}_2\text{NSe}_2$ [M + H]⁺: 509.0031, found: 509.0034.

2-phenyl-1,3-bis(m-tolylselanyl)indolizine (5ad)



Brown oil (218.5 mg, 82%); IR (KBr): 3057, 2927, 1475, 727, 692 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.41 (d, *J* = 7.0 Hz, 1H), 7.72 (d, *J* = 8.9 Hz, 1H), 7.40-7.43 (m, 2H), 7.37-7.32 (m, 3H), 7.08-6.97 (m, 5H), 6.93-6.88 (m, 3H), 6.78 (d, *J* = 7.7 Hz, 1H), 6.72 (t, *J* = 6.8 Hz, 1H), 2.25 (d, *J* = 3.6 Hz, 6H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 142.4, 140.0, 139.3, 138.6, 134.9, 134.7, 132.1, 130.9, 129.2, 128.8, 128.7, 128.6, 127.5, 127.4, 127.1, 126.3, 125.8, 125.1, 124.9, 121.6, 118.9, 112.4, 105.6, 94.7, 21.4, 21.4. HRMS MALDI (m/z): calcd for C₂₈H₂₃NSe₂ [M + H]⁺: 534.0235, found: 534.0239.

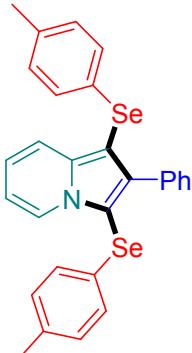
1,3-bis((3-chlorophenyl)selanyl)-2-phenylindolizine (5ae)



Yellow solid (226.3 mg, 79%), mp 109.1-111.7 °C; IR (KBr): 3062, 2921, 1476, 734, 687 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.39 (d, *J* = 6.9 Hz, 1H), 7.70 (d, *J* = 8.8 Hz, 1H), 7.38-7.33 (m, 5H), 7.09-7.15 (m, 2H), 7.08-7.02 (m, 5H), 6.97 (d, *J* = 4.9 Hz, 1H), 6.83 (d, *J* = 7.8 Hz, 1H), 6.77 (t, *J* = 6.8 Hz, 1H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 142.9, 140.2, 136.8, 135.3, 134.9, 134.3, 134.1, 130.7, 130.5, 129.9, 127.7, 126.5, 126.0, 125.9, 125.7, 125.6, 122.4, 118.8, 112.9, 105.2, 94.3.

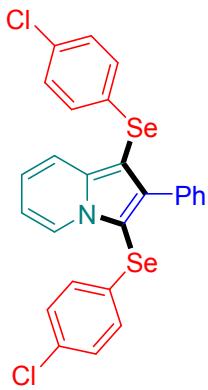
HRMS MALDI (m/z): calcd for C₂₆H₁₇Cl₂NSe₂ [M + H]⁺: 509.0031, found: 509.0034.

2-phenyl-1,3-bis(p-tolylselanyl)indolizine (5af)



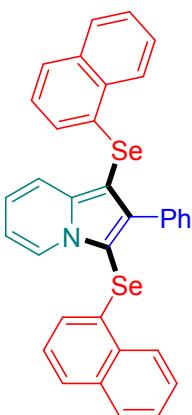
Green solid (229.2 mg, 86%), mp 116.9-119.4 °C; IR (KBr): 3052, 2929, 1470, 729, 682 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.44 (d, *J* = 7.0 Hz, 1H), 7.76 (d, *J* = 8.9 Hz, 1H), 7.44-7.48 (m, 2H), 7.41-7.36 (m, 3H), 7.07 (t, *J* = 9.0 Hz, 3H), 7.03-6.97 (m, 6H), 6.72 (t, *J* = 6.2 Hz, 1H), 2.30 (d, *J* = 2.5 Hz, 6H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 142.2, 139.8, 136.1, 135.1, 134.8, 131.0, 130.9, 130.2, 129.7, 128.3, 128.2, 127.4, 127.3, 125.7, 121.5, 118.8, 112.3, 105.8, 95.0, 20.9, 20.9. HRMS MALDI (m/z): calcd for C₂₈H₂₃NSe₂ [M + H]⁺: 534.0235, found: 534.0239.

1,3-bis((4-chlorophenyl)selanyl)-2-phenylindolizine (5ag)



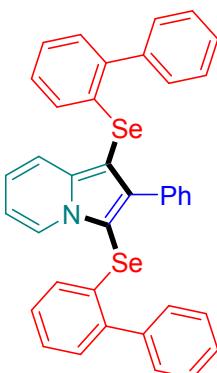
Yellow solid (226.3 mg, 79%), mp 130.2-132.9 °C; IR (KBr): 3068, 2929, 1475, 729, 693 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.37 (d, *J* = 6.9 Hz, 1H), 7.68 (d, *J* = 8.9 Hz, 1H), 7.36-7.31 (m, 5H), 7.14 (d, *J* = 8.5 Hz, 2H), 7.10 (d, *J* = 8.6 Hz, 2H), 7.04-6.99 (m, 3H), 6.92 (d, *J* = 8.5 Hz, 2H), 6.76 (t, *J* = 6.8 Hz, 1H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 142.6, 140.0, 134.4, 133.0, 132.4, 131.5, 130.7, 130.3, 129.6, 129.5, 129.4, 129.1, 127.7, 127.6, 125.7, 122.2, 118.8, 112.8, 105.4, 94.7. HRMS MALDI (m/z): calcd for C₂₆H₁₇Cl₂NSe₂ [M + H]⁺: 509.0031, found: 509.0034.

1,3-bis(naphthalen-1-ylselanyl)-2-phenylindolizine (5ah)



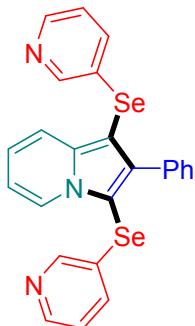
Yellow solid (245.0 mg, 81%), mp 135.2-137.5 °C; IR (KBr): 3041, 2941, 1470, 733, 692 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.26 (d, *J* = 6.9 Hz, 1H), 8.08 (d, *J* = 8.8 Hz, 1H), 8.01-7.95 (m, 1H), 7.78-7.71 (m, 2H), 7.62-7.53 (m, 3H), 7.46-7.37 (m, 6H), 7.22-7.17 (m, 3H), 7.10 (t, *J* = 7.8 Hz, 2H), 6.96 (d, *J* = 7.3 Hz, 1H), 6.88-6.82 (m, 1H), 6.76 (d, *J* = 7.3 Hz, 1H), 6.56 (t, *J* = 6.8 Hz, 1H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 143.0, 140.4, 134.5, 134.1, 133.8, 133.7, 132.2, 132.1, 130.7, 130.5, 128.6, 128.4, 127.6, 127.5, 126.8, 126.4, 126.3, 126.2, 126.0, 126.0, 125.9, 125.7, 125.5, 125.4, 121.8, 118.9, 112.5, 104.4, 93.9. HRMS MALDI (m/z): calcd for C₃₄H₂₃NSe₂ [M + H]⁺: 606.0235, found: 606.0241.

1,3-bis([1,1'-biphenyl]-2-ylselanyl)-2-phenylindolizine (5ai)



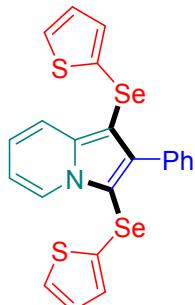
Brown solid (243.1 mg, 74%), mp 162.3-164.4 °C; IR (KBr): 3053, 2911, 1483, 740, 689 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.22 (d, *J* = 7.1 Hz, 1H), 7.60 (d, *J* = 8.9 Hz, 1H), 7.52-7.35 (m, 13H), 7.33-7.29 (m, 3H), 7.24-7.19 (m, 2H), 7.16 (t, *J* = 6.7 Hz, 1H), 7.07-7.02 (m, 2H), 6.98-6.94 (m, 1H), 6.82 (d, *J* = 8.0 Hz, 1H), 6.68 (t, *J* = 6.8 Hz, 1H), 6.55 (d, *J* = 7.7 Hz, 1H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 142.8, 141.8, 141.4, 141.3, 140.8, 140.3, 134.8, 134.5, 132.1, 130.7, 130.4, 129.9, 129.0, 128.8, 128.5, 128.4, 128.3, 127.9, 127.9, 127.7, 127.6, 127.5, 127.3, 127.1, 126.1, 125.9, 125.2, 121.7, 118.9, 112.5, 105.9, 95.1. HRMS MALDI (m/z): calcd for C₃₈H₂₇NSe₂ [M + H]⁺: 658.0548, found: 658.0555.

2-phenyl-1,3-bis(pyridin-3-ylselanyl)indolizine (5aj)



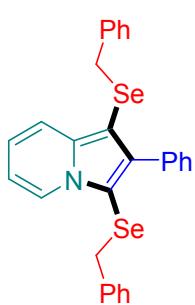
Yellow solid (212.9 mg, 84%), mp 120.8-123.5 °C; IR (KBr): 3026, 2922, 1540, 1406, 1014, 744, 702 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.39 (d, *J* = 7.0 Hz, 1H), 8.36 (d, *J* = 4.7 Hz, 1H), 8.30 (d, *J* = 6.8 Hz, 3H), 7.69 (d, *J* = 8.9 Hz, 1H), 7.38-7.29 (m, 6H), 7.24 (d, *J* = 8.4 Hz, 1H), 7.09-7.00 (m, 3H), 6.76 (t, *J* = 6.9 Hz, 1H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 149.0, 147.6, 146.5, 142.8, 140.0, 136.4, 136.1, 134.1, 131.7, 130.7, 127.8, 127.7, 125.6, 124.4, 124.0, 122.4, 118.7, 113.0, 104.5, 93.8. HRMS MALDI (m/z): calcd for C₂₄H₁₇N₃Se₂ [M + H]⁺: 507.9827, found: 507.9830.

2-phenyl-1,3-bis(thiophen-2-ylselanyl)indolizine (5ak)



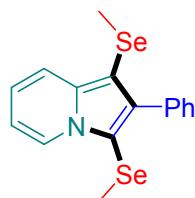
Brown solid (183.5 mg, 71%), mp 95.3-97.8 °C; IR (KBr): 3020, 2922, 1629, 1492, 741, 694 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.57 (d, *J* = 6.4 Hz, 1H), 7.85 (d, *J* = 8.5 Hz, 1H), 7.54-7.44 (m, 5H), 7.23 (d, *J* = 5.3 Hz, 1H), 7.18 (d, *J* = 5.1 Hz, 1H), 7.03 (d, *J* = 7.0 Hz, 1H), 7.00 (d, *J* = 3.3 Hz, 1H), 6.93 (d, *J* = 2.8 Hz, 1H), 6.89-6.85 (m, 1H), 6.84-6.80 (m, 1H), 6.76 (t, *J* = 6.8 Hz, 1H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 141.0, 138.8, 134.8, 132.4, 131.8, 131.5, 129.3, 128.7, 128.6, 127.8, 127.5, 127.5, 125.5, 125.4, 121.6, 118.8, 112.3, 107.5, 97.8. HRMS MALDI (m/z): calcd for C₂₂H₁₅NS₂Se₂ [M + H]⁺: 517.9048, found: 517.9051.

1,3-bis(benzylselanyl)-2-phenylindolizine (5al)



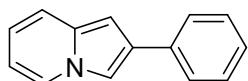
Brown oil (213.2 mg, 80%); IR (KBr): 3024, 2926, 1492, 1342, 743, 692 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.23 (d, *J* = 7.0 Hz, 1H), 7.47-7.43 (m, 3H), 7.38 (d, *J* = 8.9 Hz, 1H), 7.28 (d, *J* = 6.6 Hz, 2H), 7.18-7.11 (m, 6H), 6.94 (d, *J* = 7.8 Hz, 2H), 6.89 (d, *J* = 7.4 Hz, 2H), 6.82-6.78 (m, 1H), 6.54 (t, *J* = 6.8 Hz, 1H), 3.76 (s, 2H), 3.67 (s, 2H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 142.0, 139.3, 139.2, 138.5, 135.4, 131.0, 128.6, 128.4, 128.3, 128.0, 127.3, 127.0, 126.7, 126.2, 125.2, 120.2, 118.5, 111.3, 106.5, 95.8, 33.0, 32.8. ⁷⁷Se NMR (76 MHz, Chloroform-d) δ ppm 202.2, 184.4. HRMS MALDI (m/z): calcd for C₂₈H₂₃NSe₂ [M + H]⁺: 534.0235, found: 534.0239.

1,3-bis(methylselanyl)-2-phenylindolizine (5am)



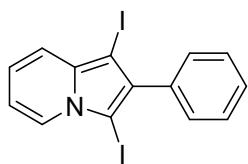
Brown oil (117.8 mg, 62%); IR (KBr): 3057, 2920, 1492, 1340, 748, 698 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.51 (d, *J* = 7.1 Hz, 1H), 7.72 (d, *J* = 9.0 Hz, 1H), 7.45-7.55 (m, 4H), 7.42 (t, *J* = 7.0 Hz, 1H), 7.00-6.89 (m, 1H), 6.73 (t, *J* = 6.3 Hz, 1H), 2.00 (s, 3H), 1.94 (s, 3H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 140.1, 138.0, 135.6, 130.9, 127.5, 127.2, 125.5, 120.0, 118.9, 111.6, 107.1, 96.6, 9.8, 8.7. ⁷⁷Se NMR (76 MHz, Chloroform-d) δ ppm 21.2, 1.6. HRMS MALDI (m/z): calcd for C₁₆H₁₅NSe₂ [M + H]⁺: 381.9609., found: 381.9607.

2-phenylindolizine (6aa)



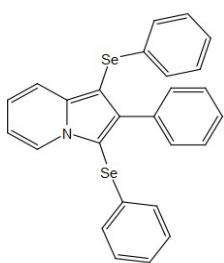
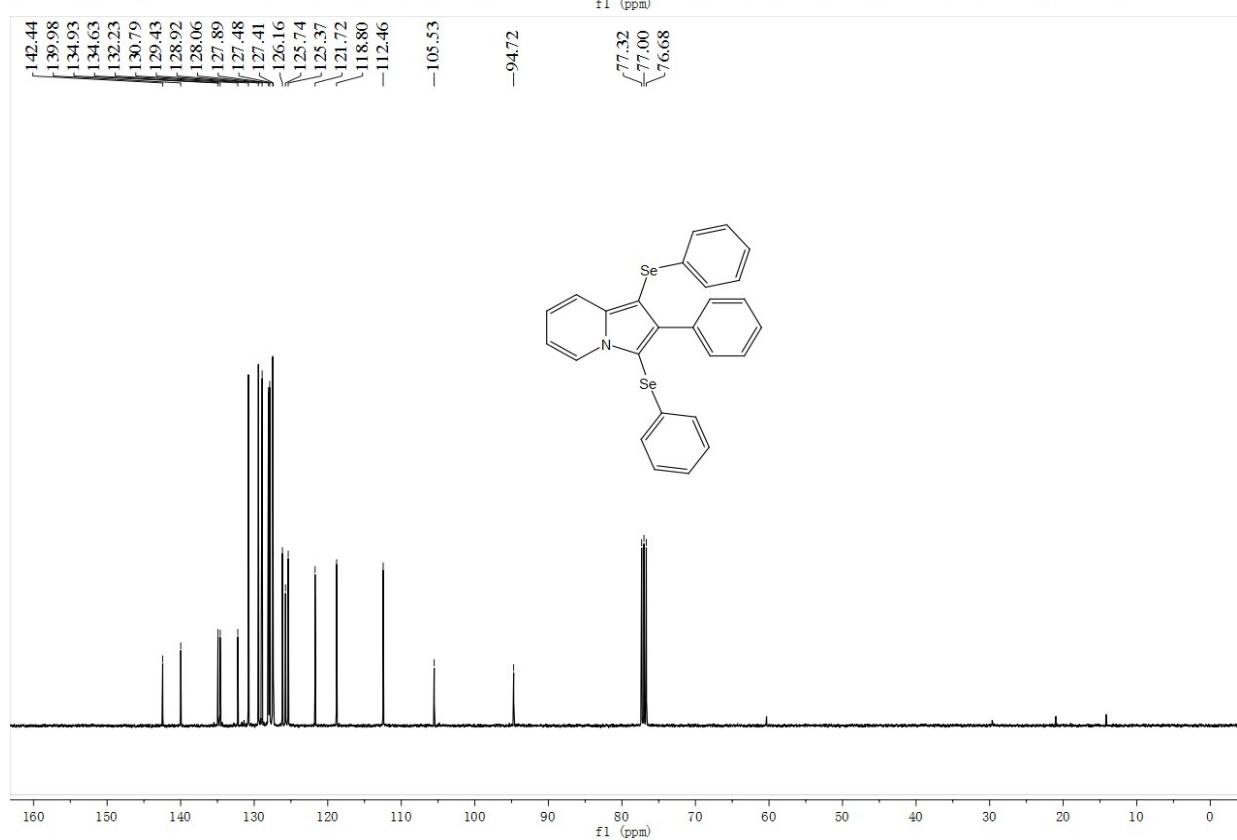
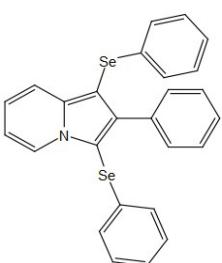
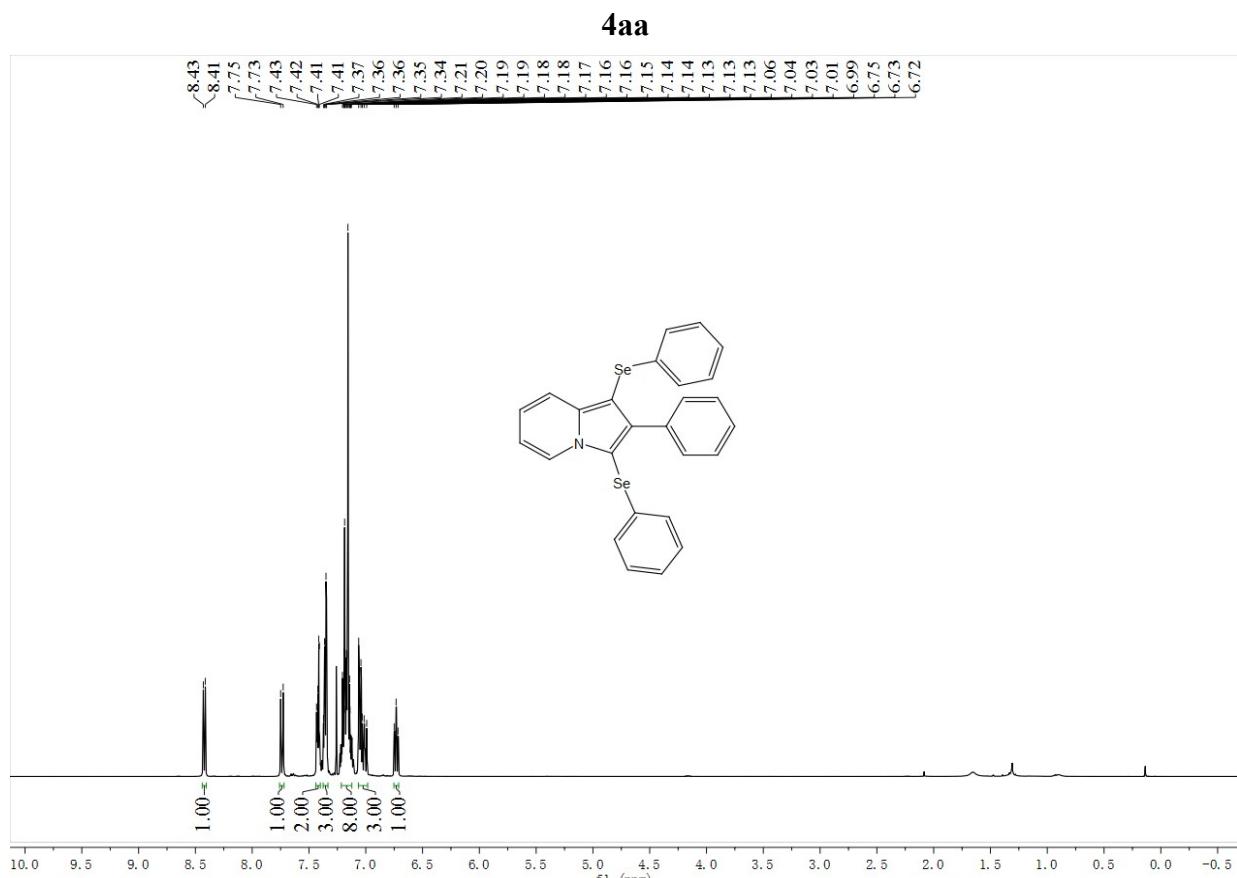
Brown solid (57.9 mg, 60%); mp 213.6-215.2 °C; IR (KBr): 3101, 1631, 1454, 727, 688 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 7.89 (d, *J* = 7.0 Hz, 1H), 7.67 (d, *J* = 7.2 Hz, 2H), 7.58 (s, 1H), 7.40 (t, *J* = 7.7 Hz, 2H), 7.35 (d, *J* = 9.0 Hz, 1H), 7.29-7.23 (m, 1H), 6.70 (s, 1H), 6.68-6.60 (m, 1H), 6.45 (t, *J* = 6.5 Hz, 1H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 135.3, 133.6, 129.4, 128.7, 126.5, 126.2, 125.0, 119.0, 117.3, 110.5, 109.2, 96.6. MS (EI) calcd for C₁₄H₁₁N: 193.1, found: 193.1.

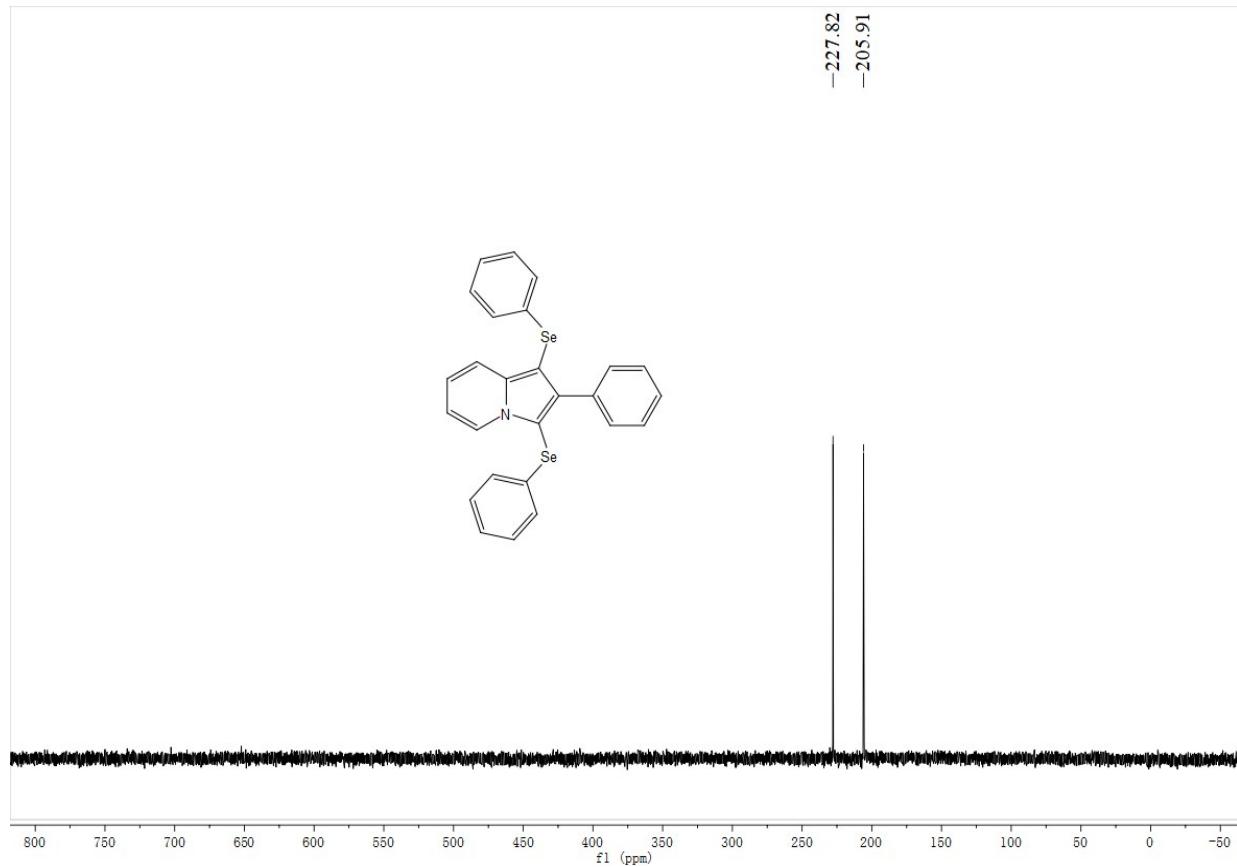
1,3-diiodo-2-phenylindolizine (7aa)



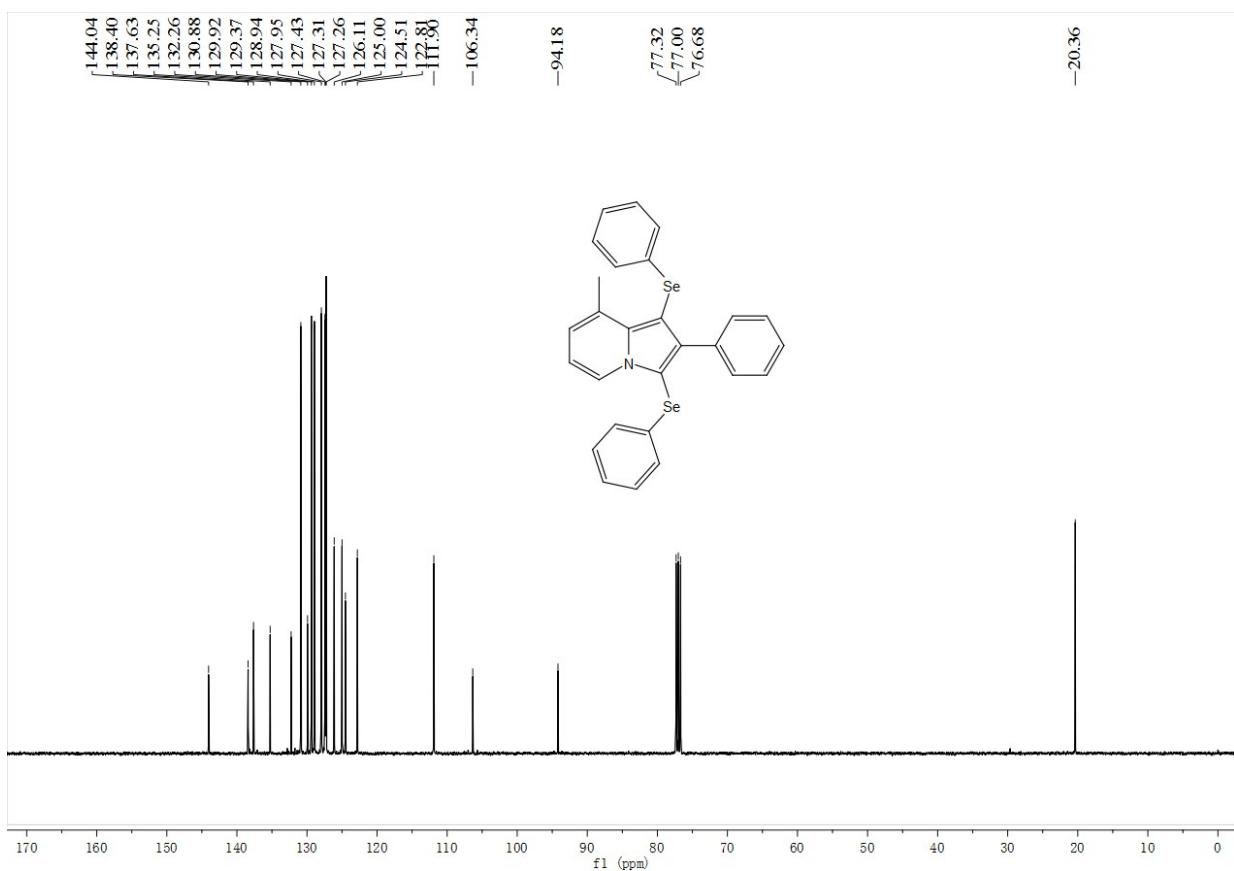
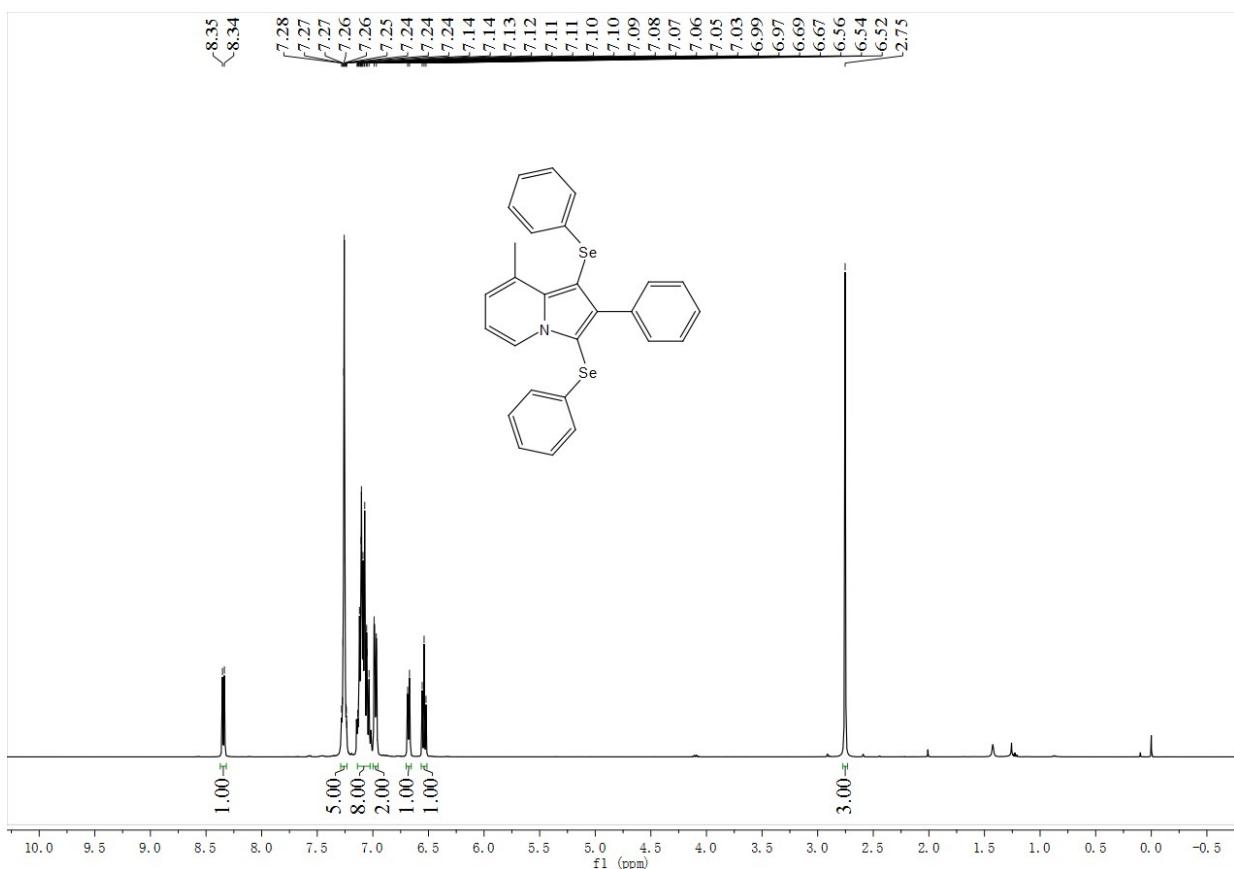
Unstable white solid (193.5 mg, 87%); IR (KBr): 2920, 2848, 1631, 1348, 725, 702 cm⁻¹; ¹H NMR (400 MHz, Chloroform-*d*) δ ppm 8.06 (d, *J* = 7.0 Hz, 1H), 7.55-7.40 (m, 5H), 7.35 (d, *J* = 9.0 Hz, 1H), 6.91-6.82 (m, 1H), 6.71 (t, *J* = 6.8 Hz, 1H). ¹³C NMR (100 MHz, Chloroform-*d*) δ ppm 137.9, 136.4, 135.4, 130.7, 128.0, 127.9, 127.4, 119.8, 119.7, 112.4, 62.2, 55.4. MS (EI) calcd for C₁₄H₉I₂N: 444.9, found: 444.9.

F. NMR Spectra

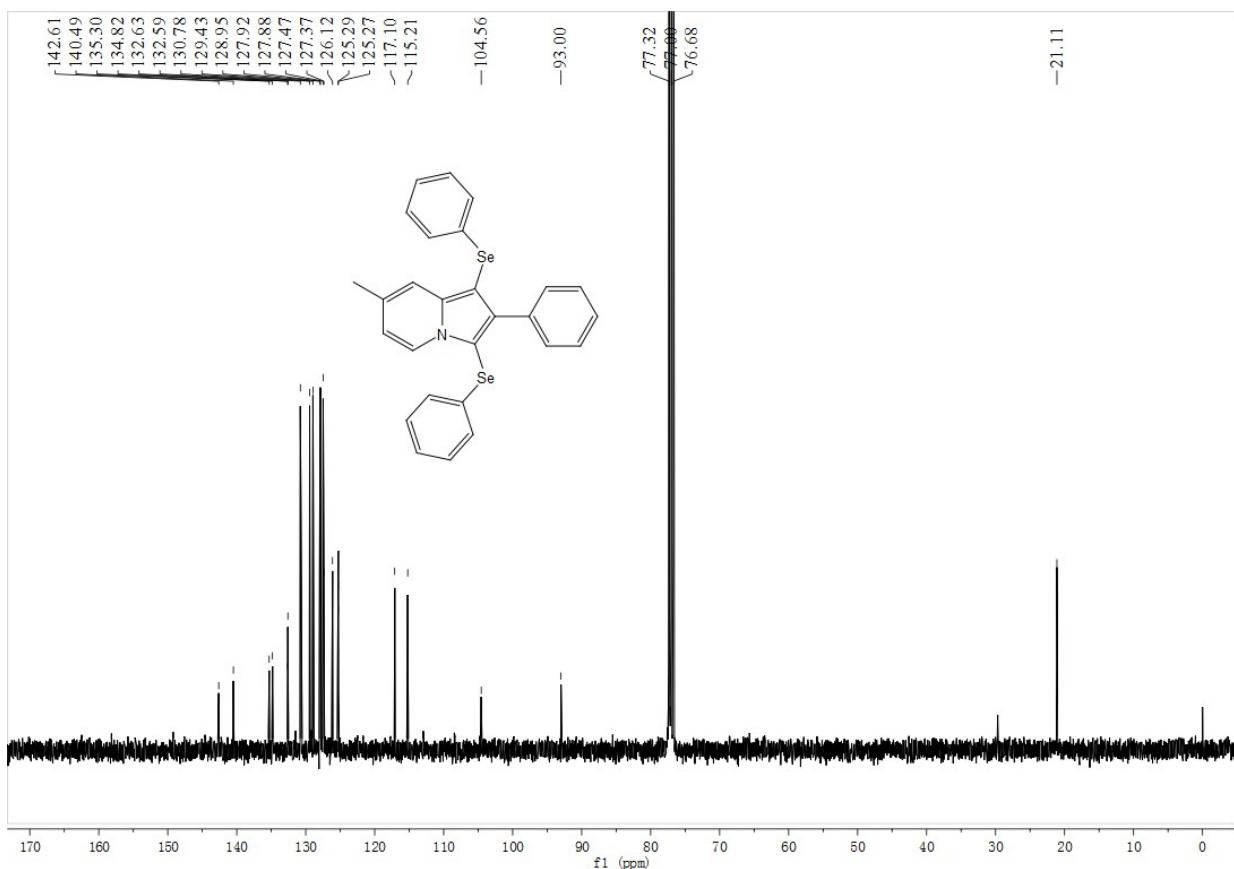
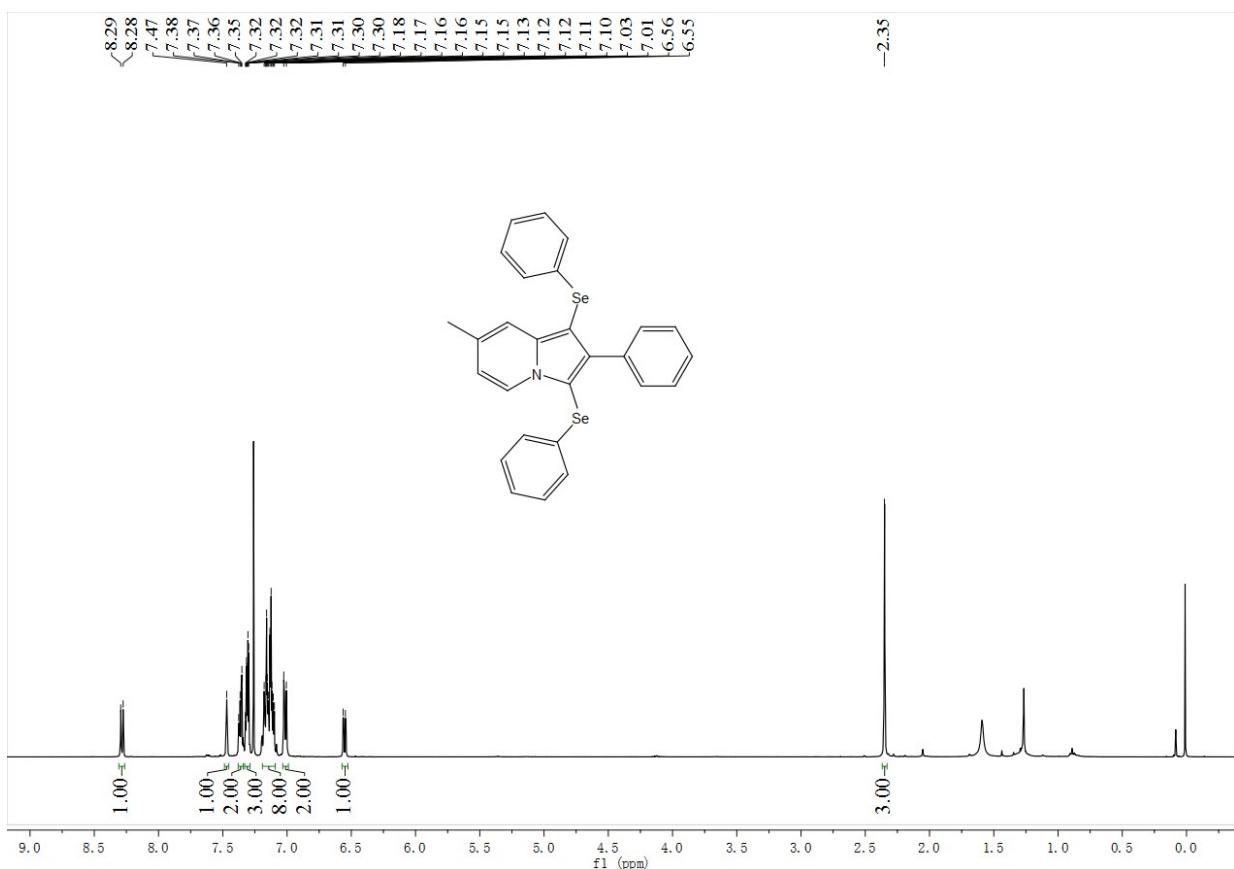




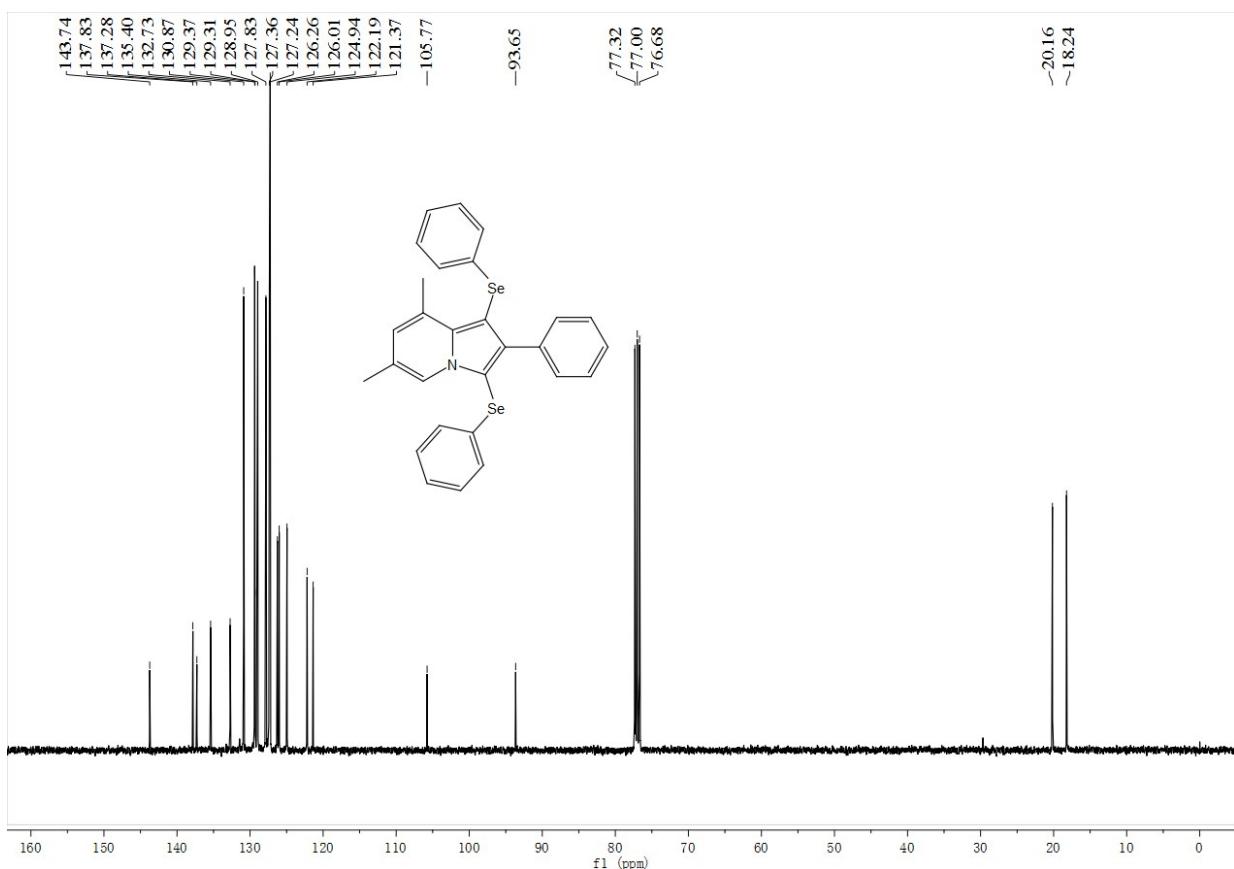
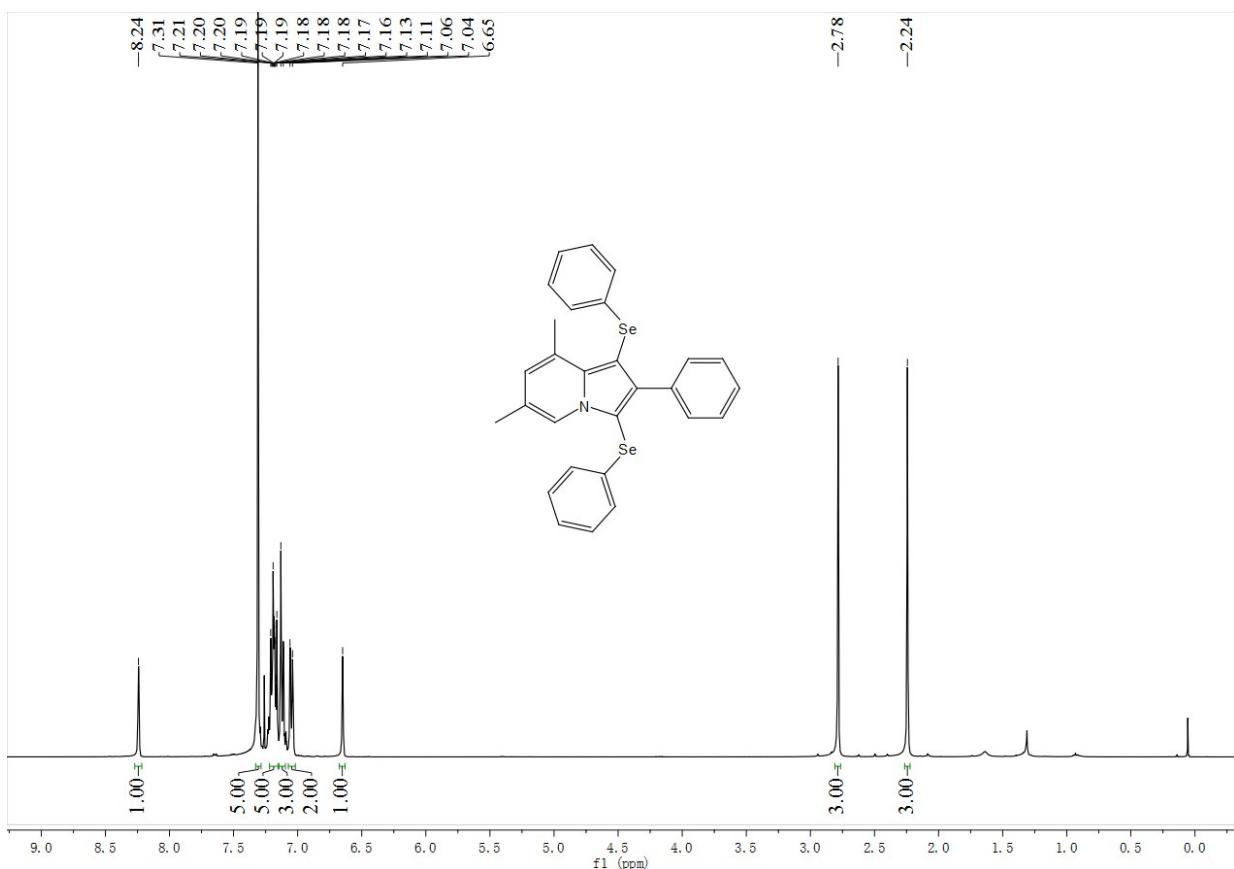
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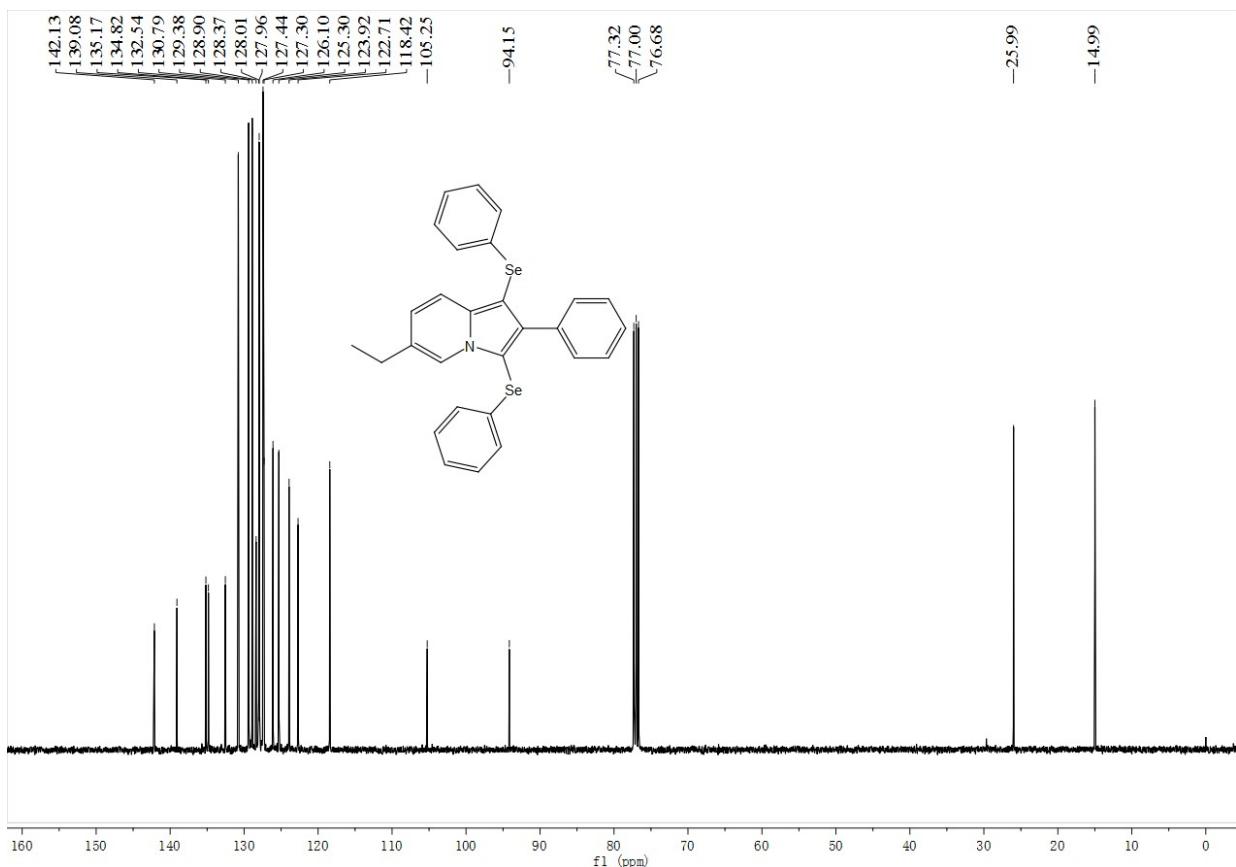
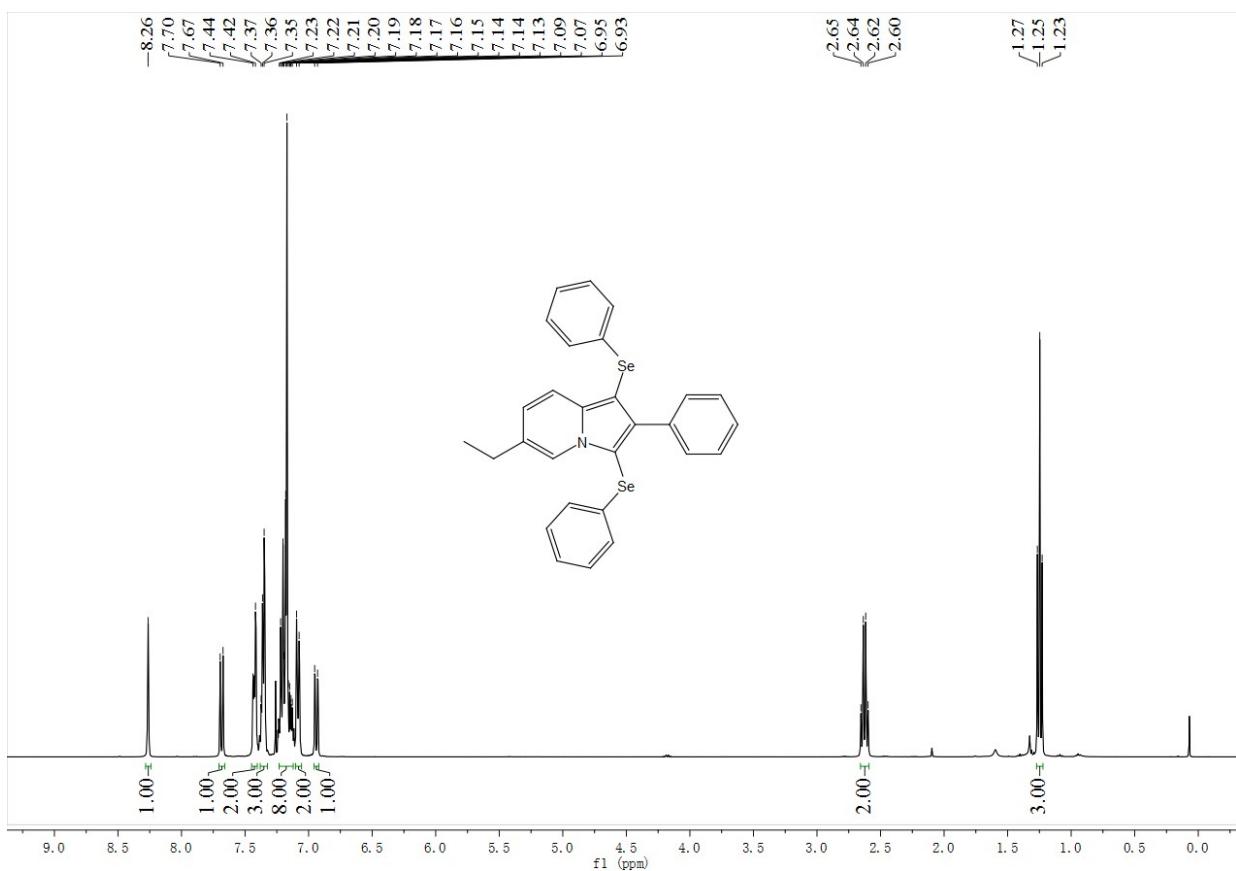
4ac



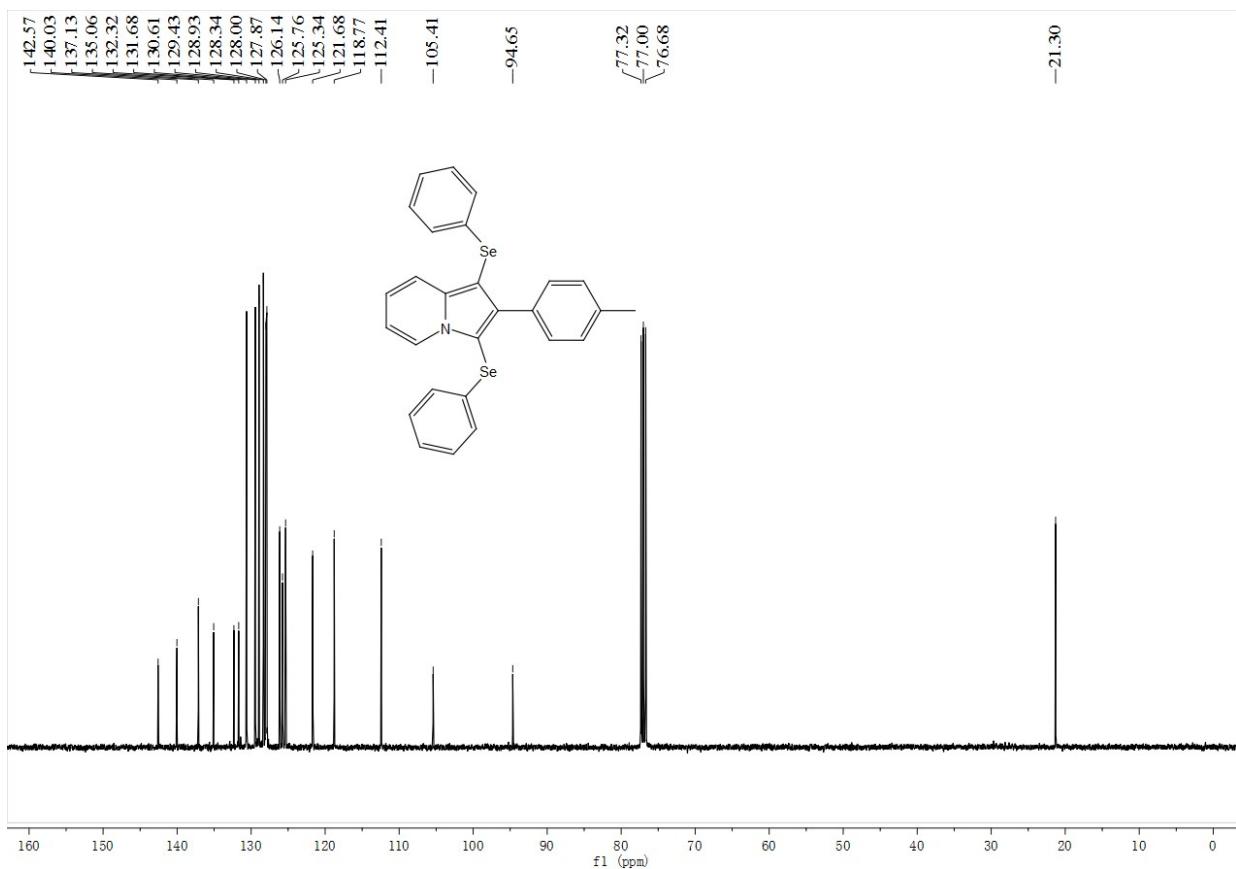
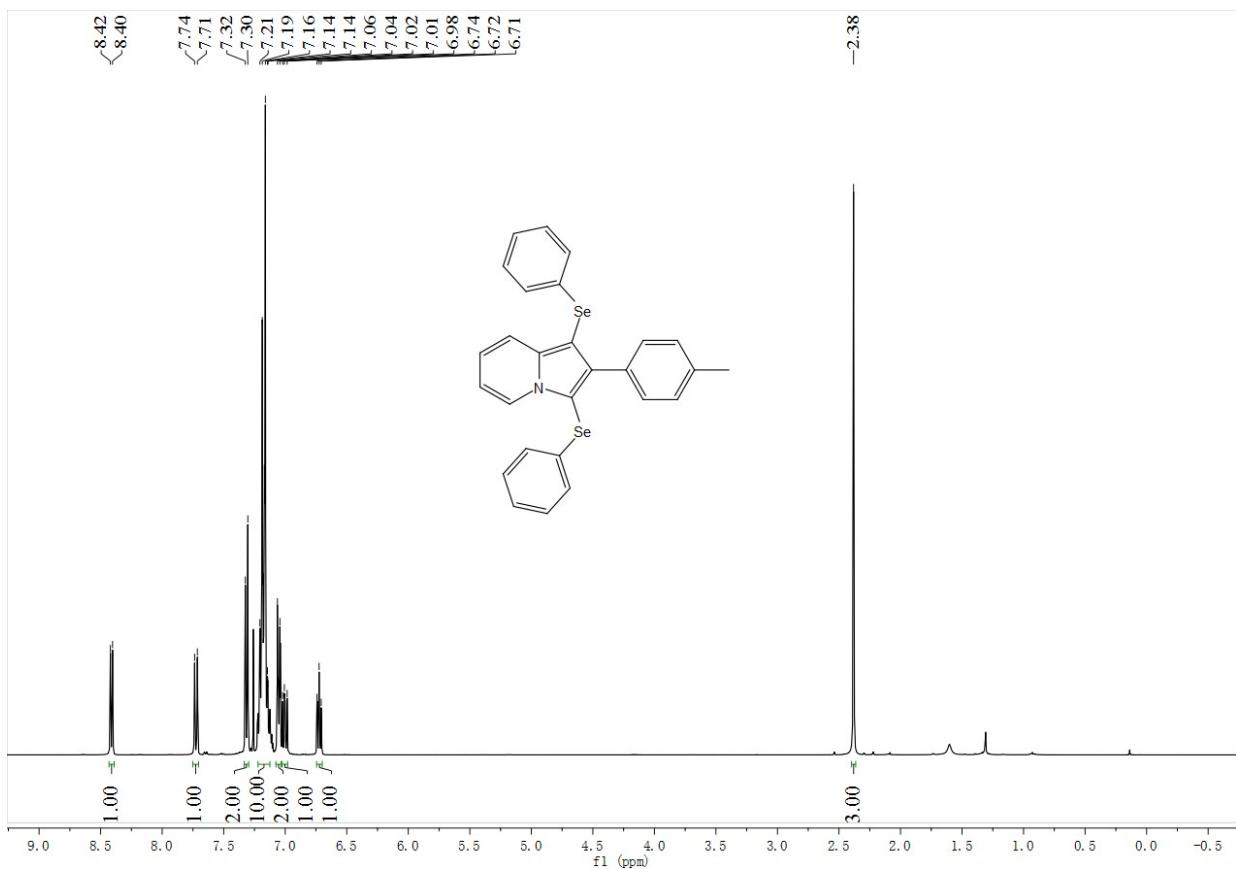
4ad



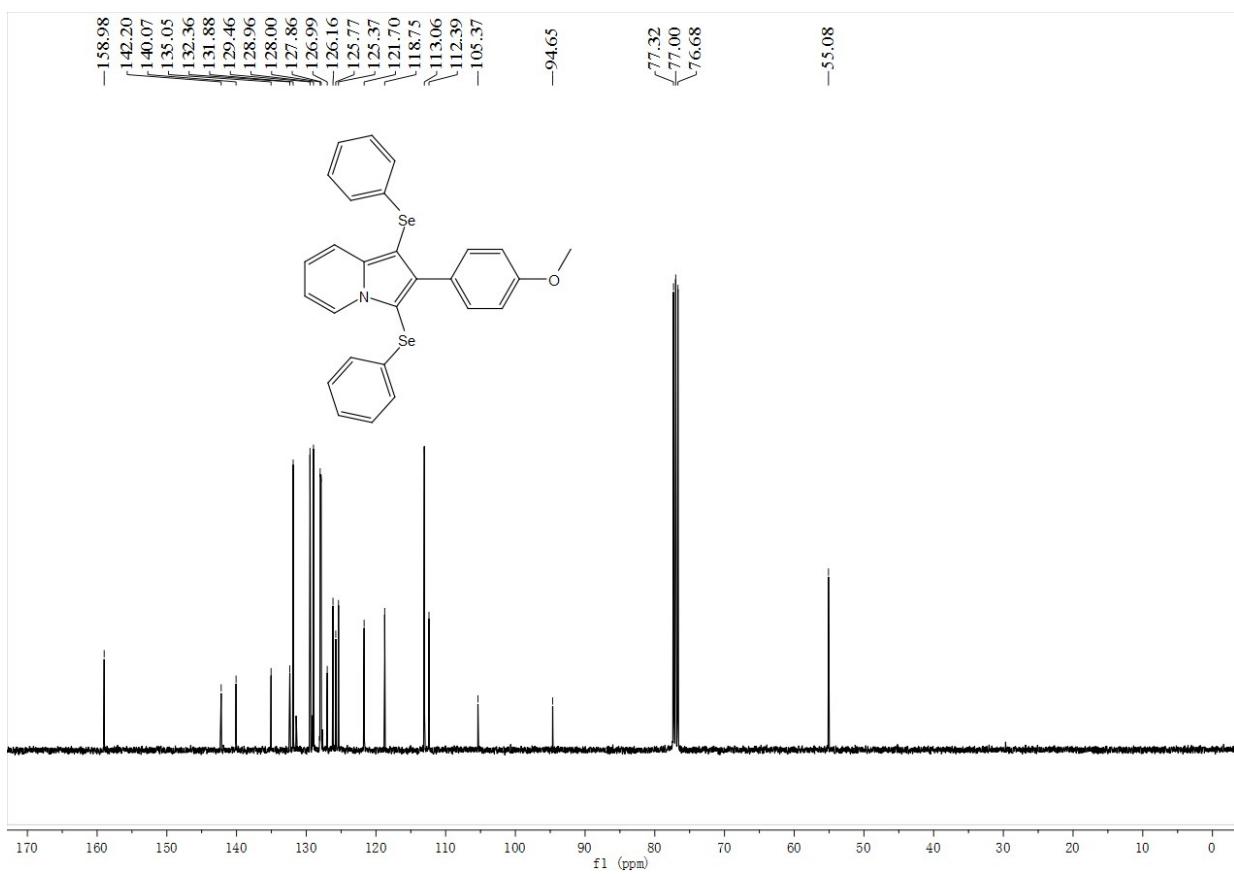
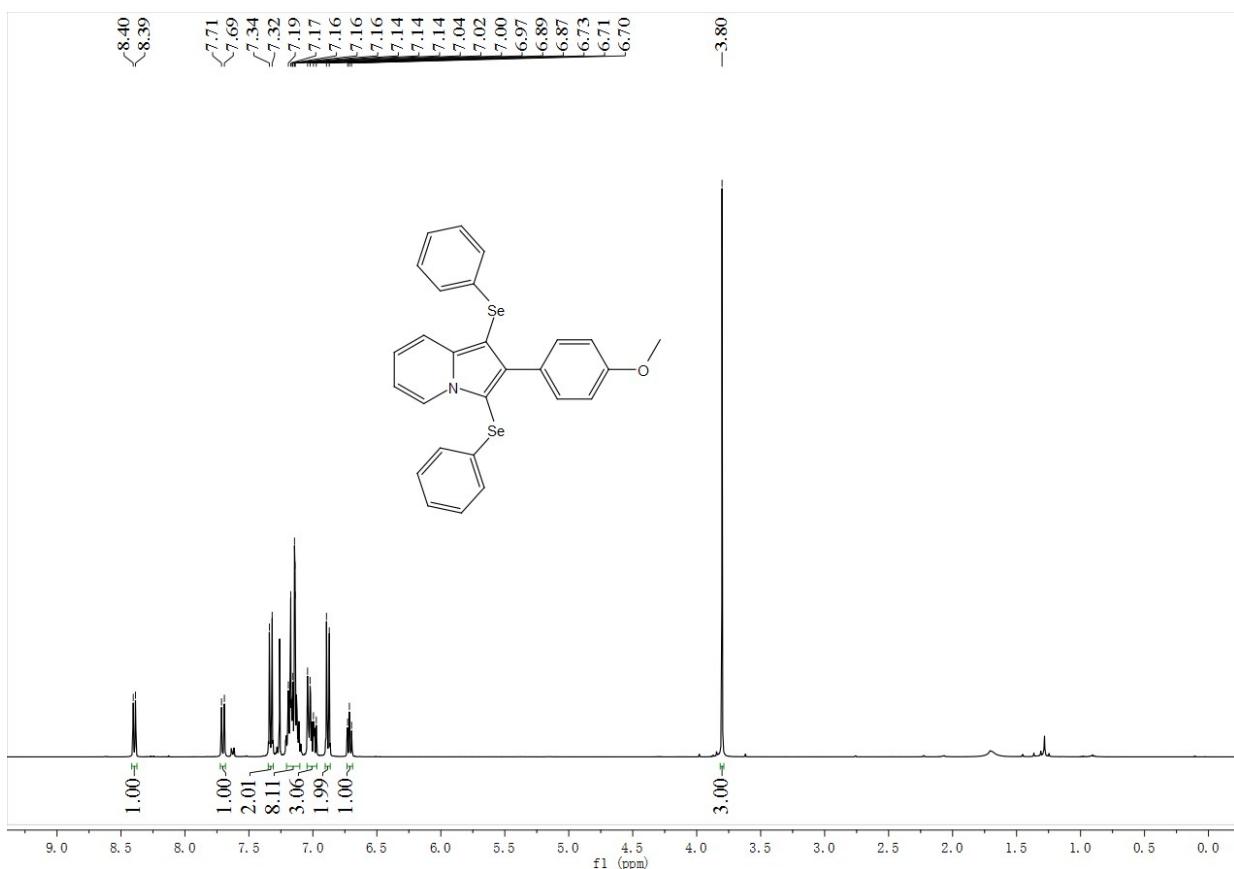
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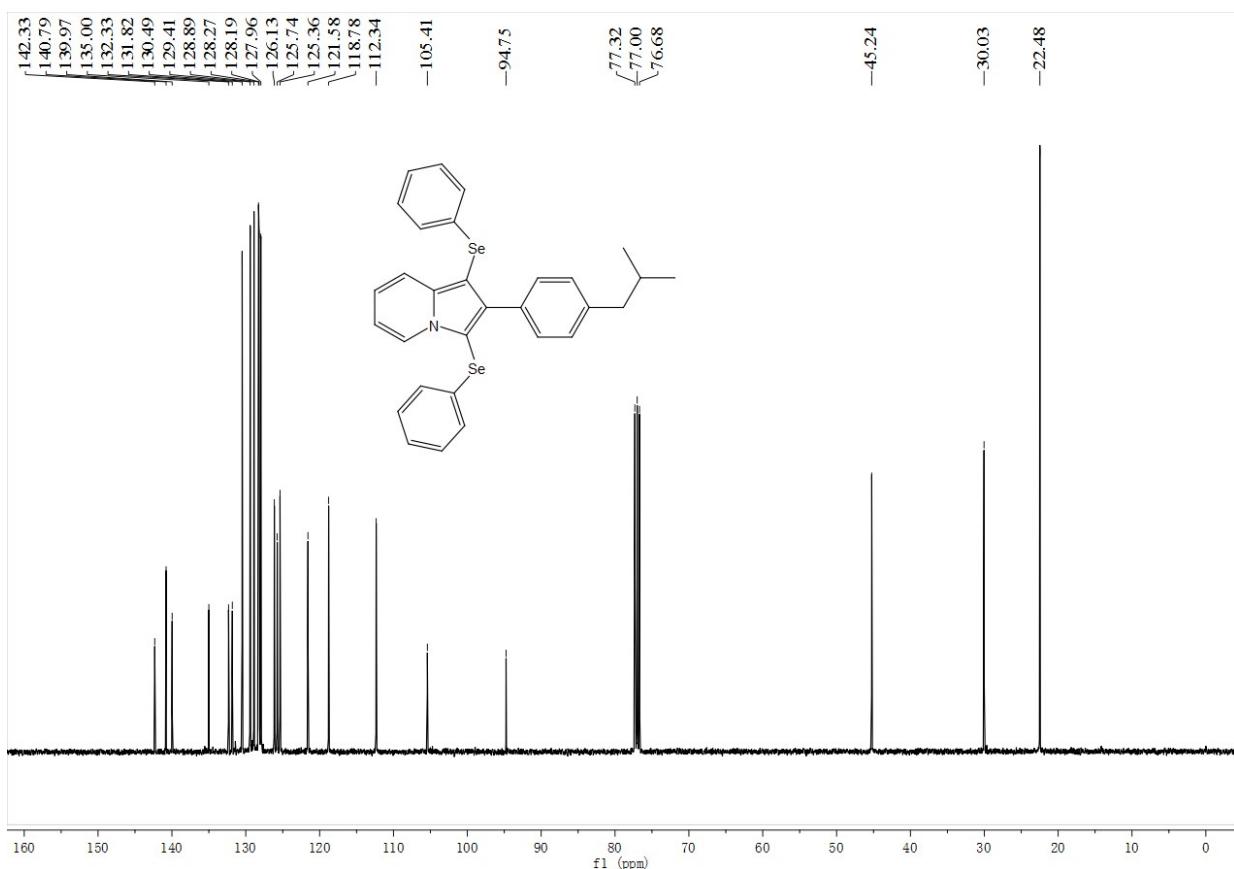
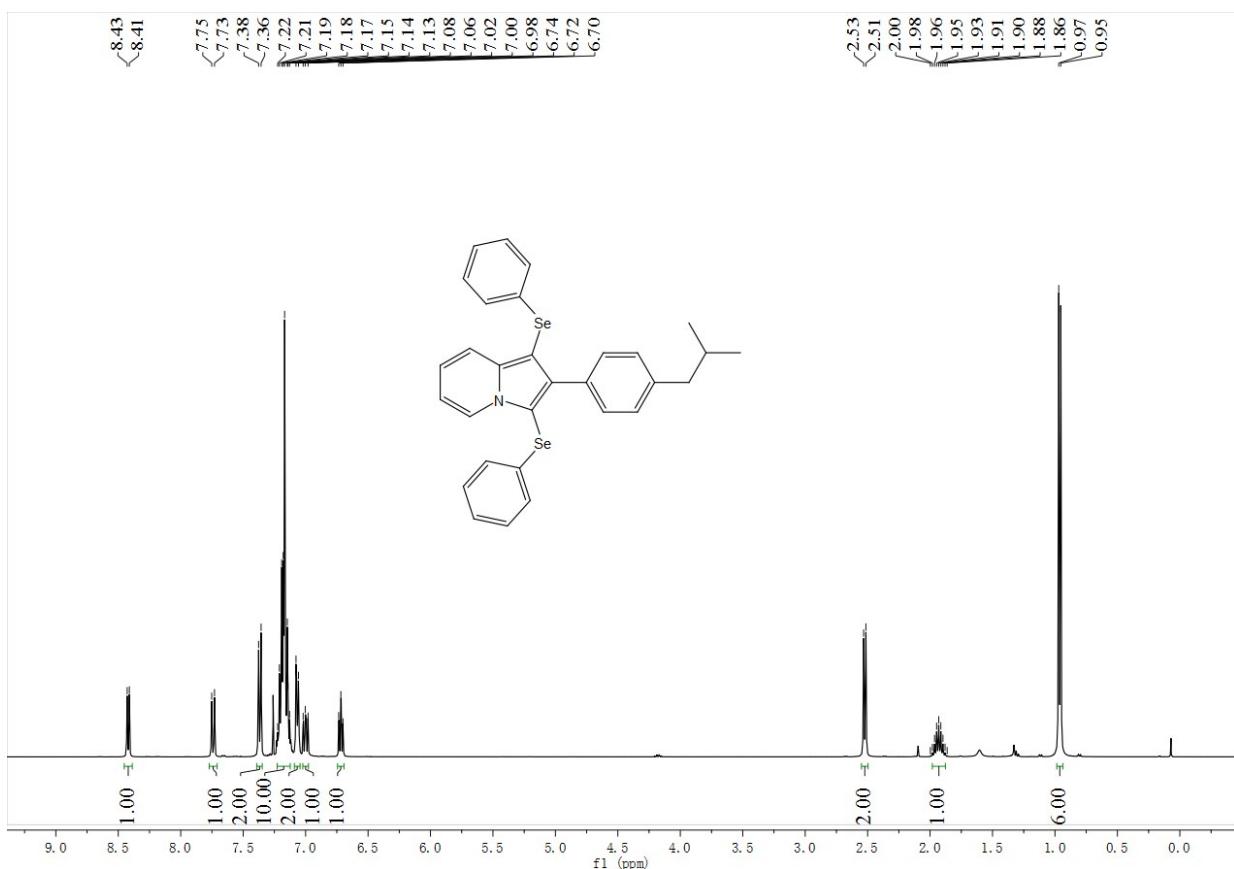
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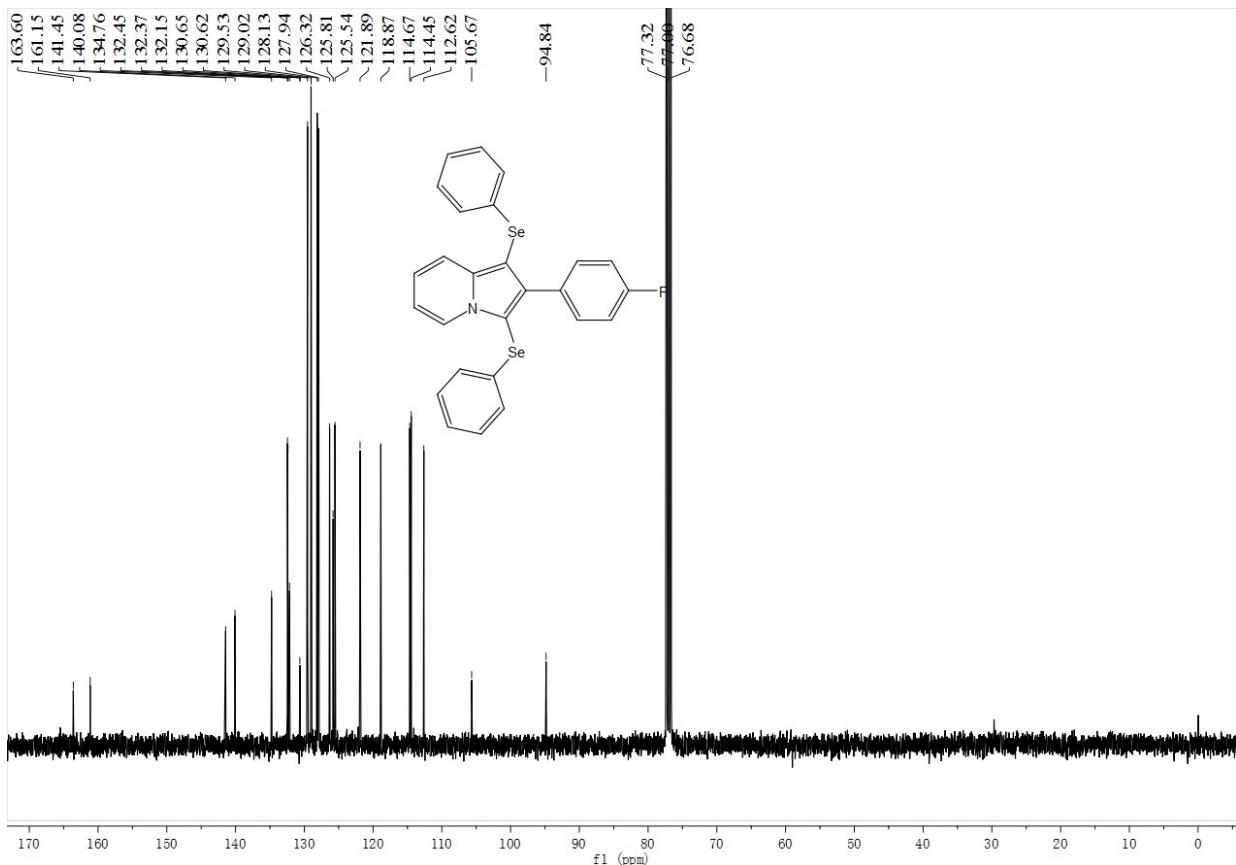
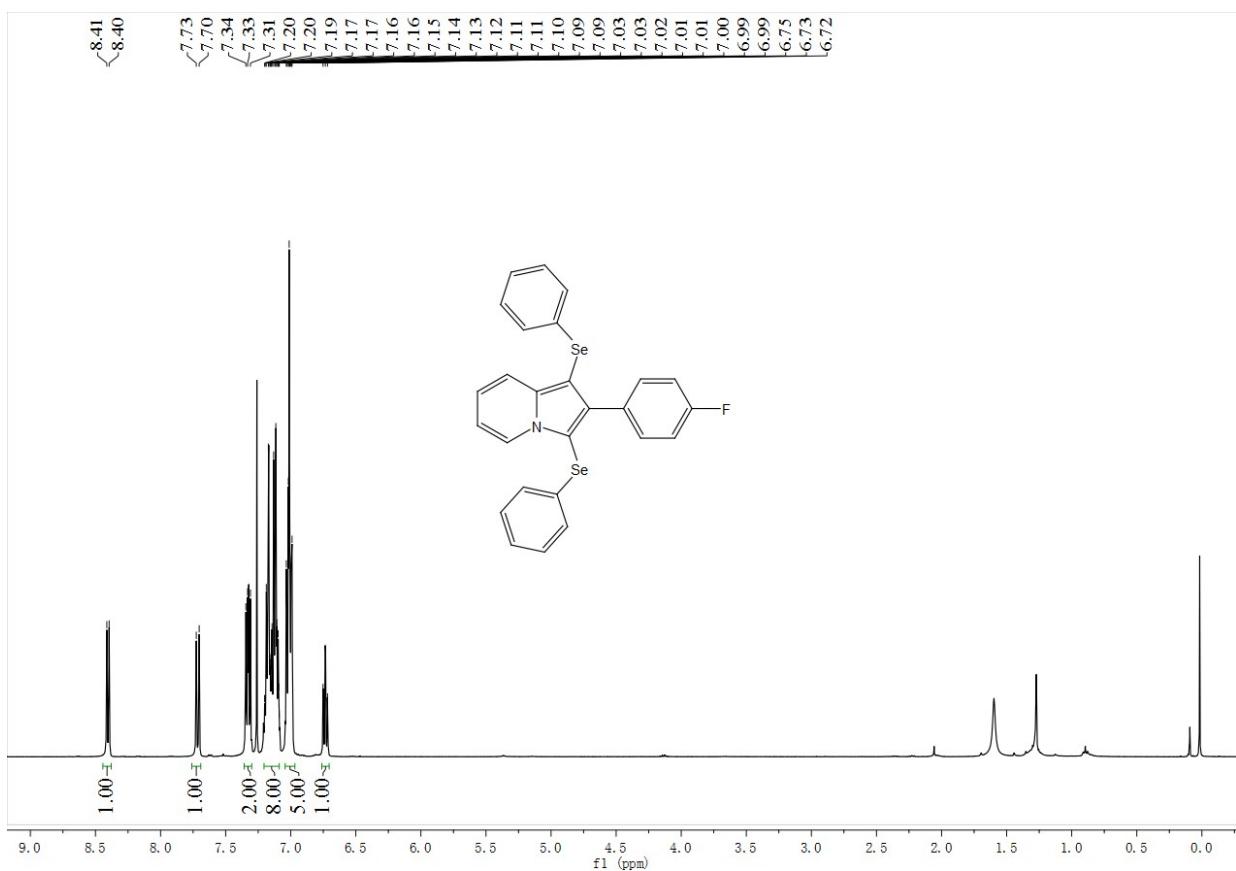
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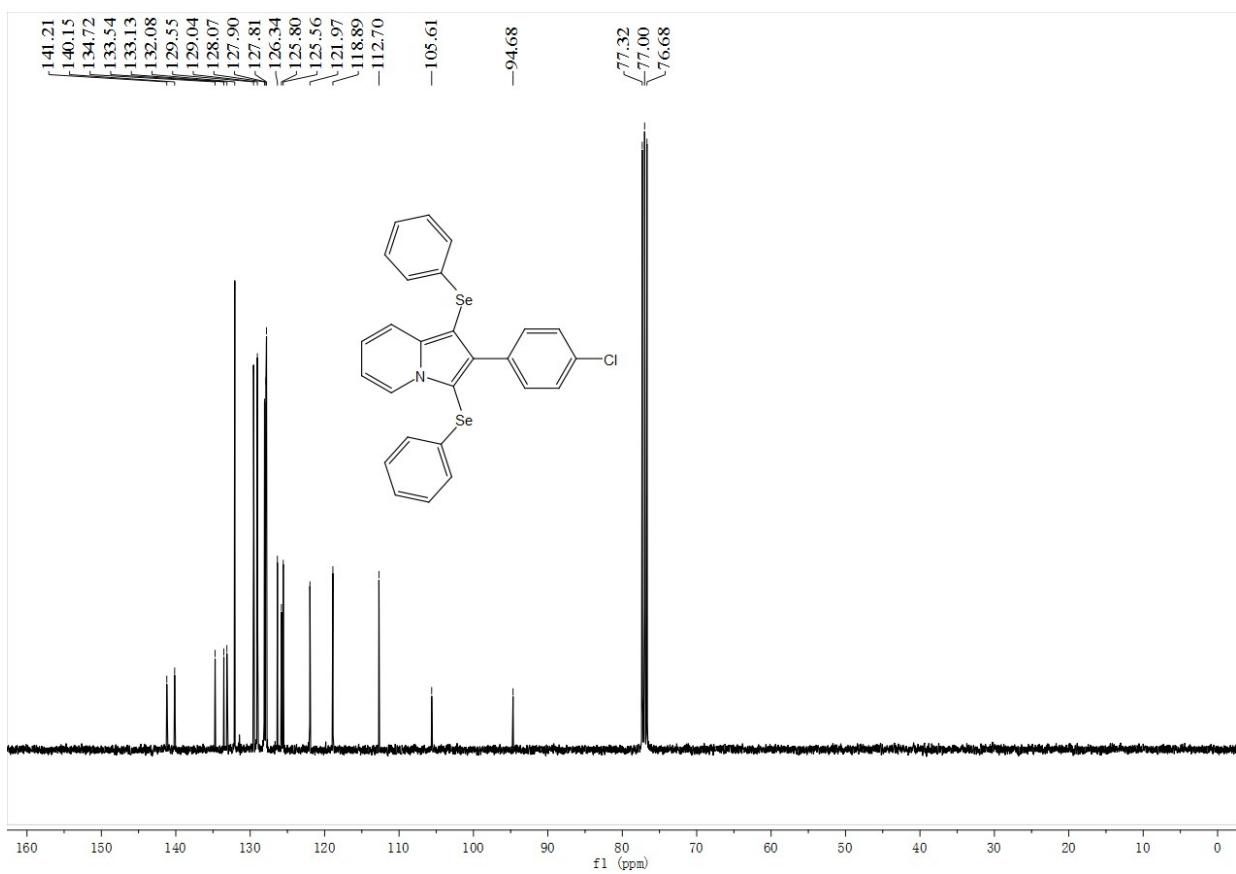
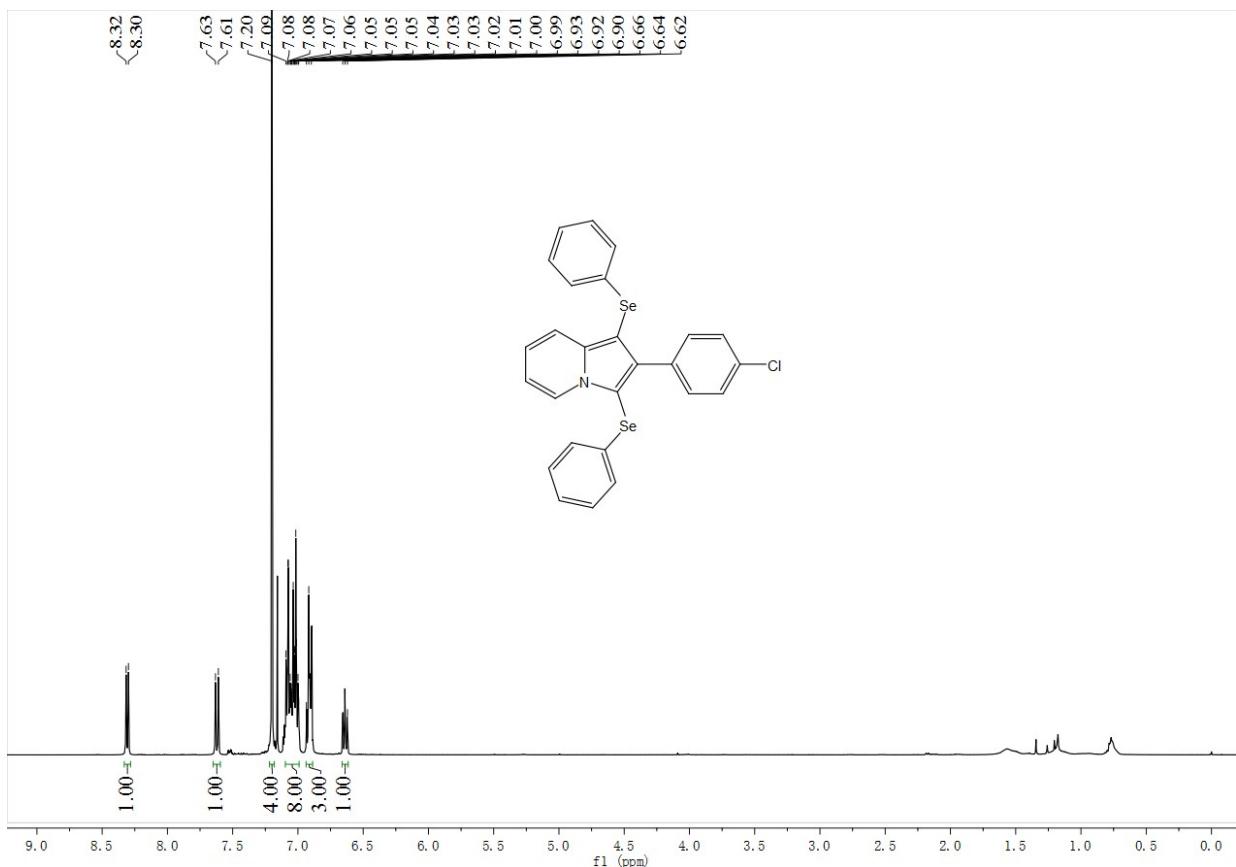
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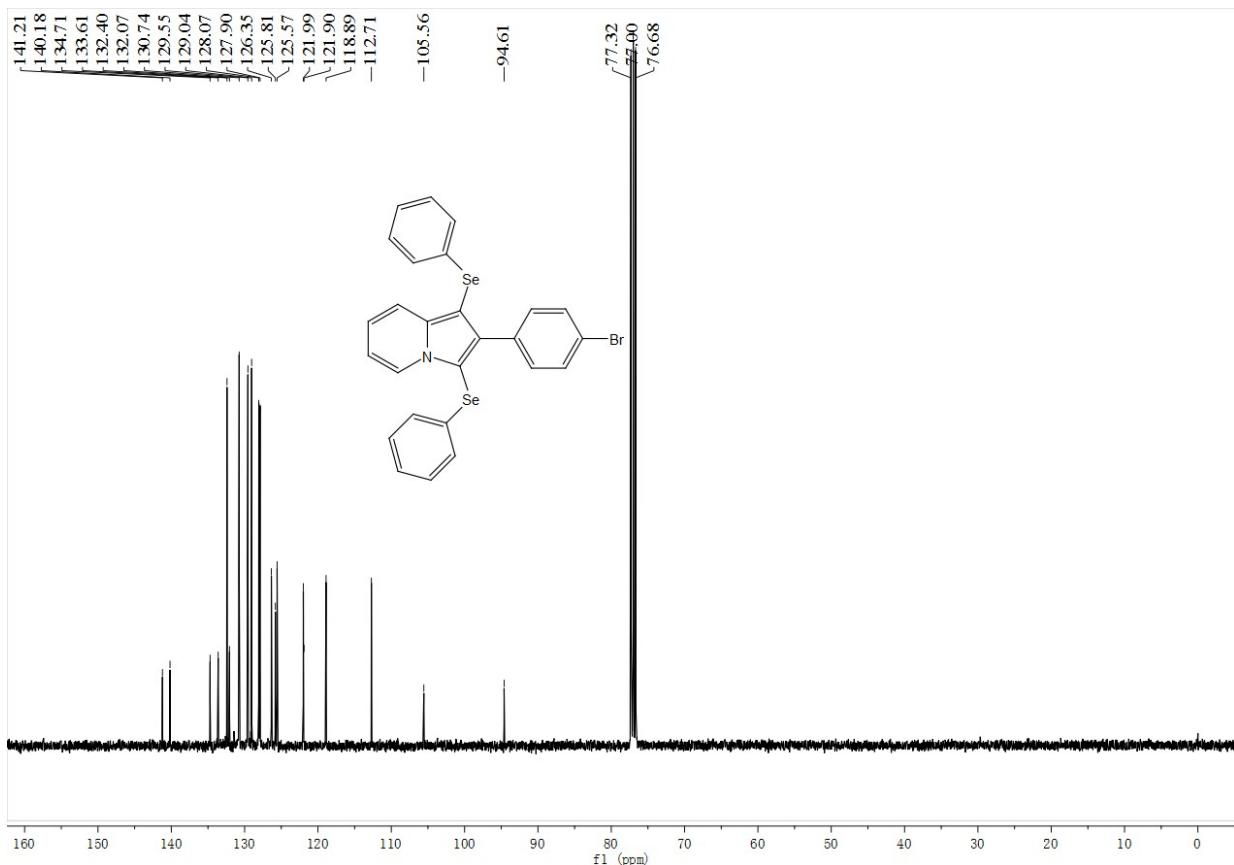
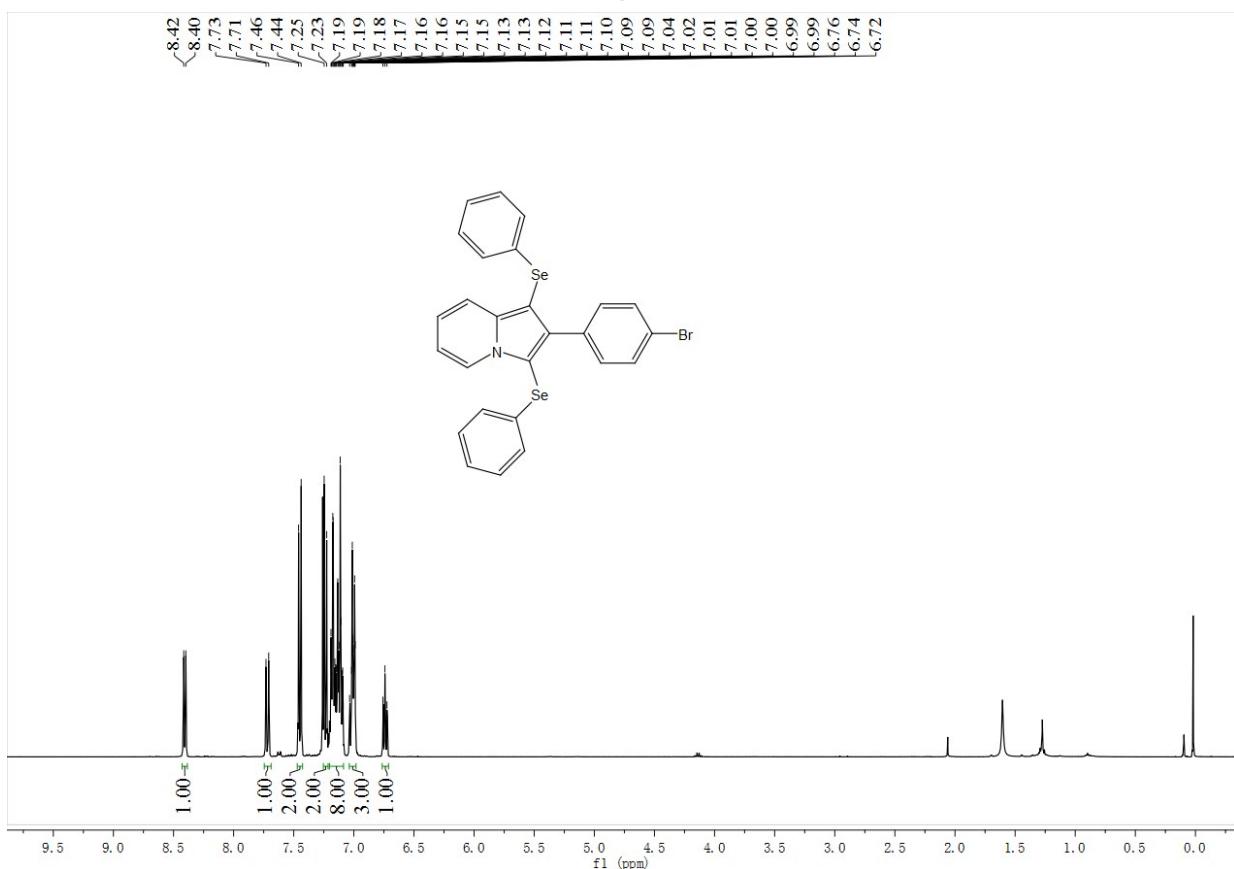
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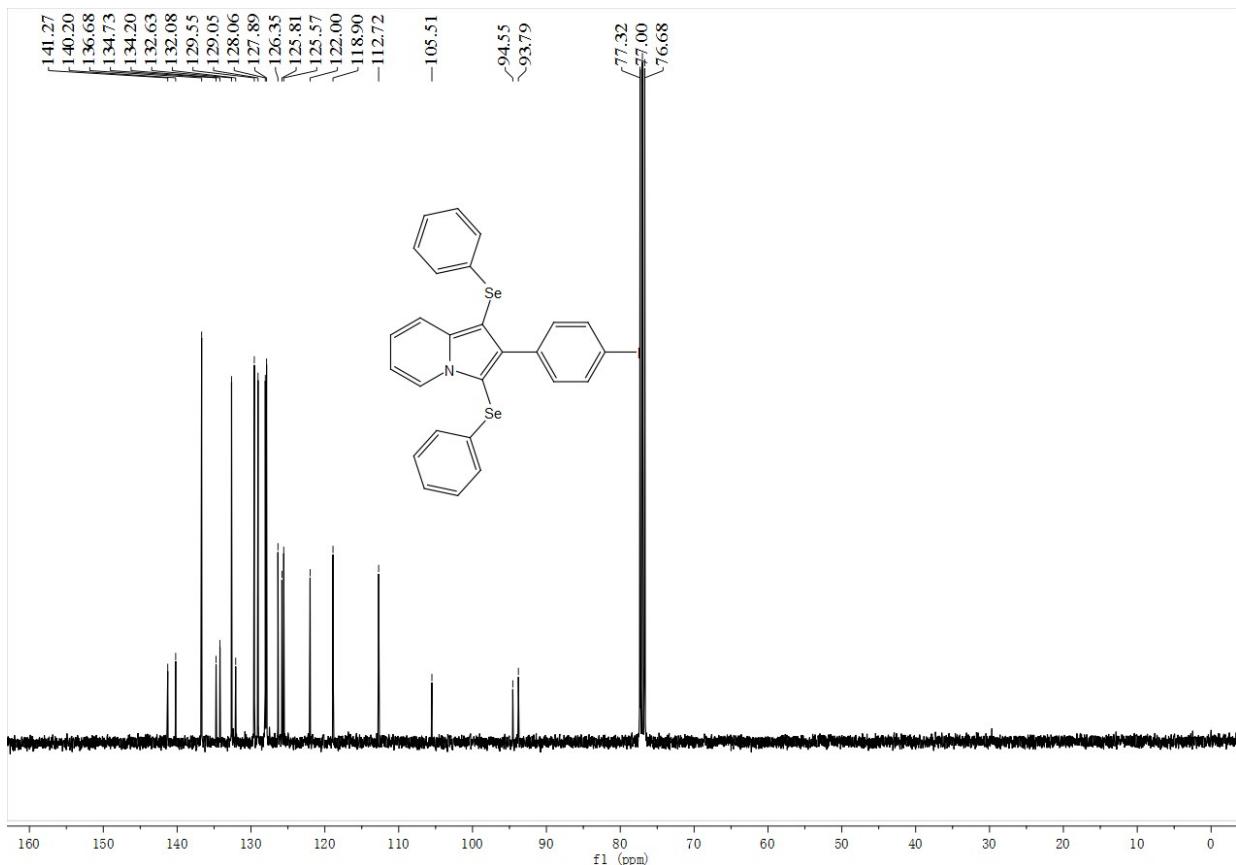
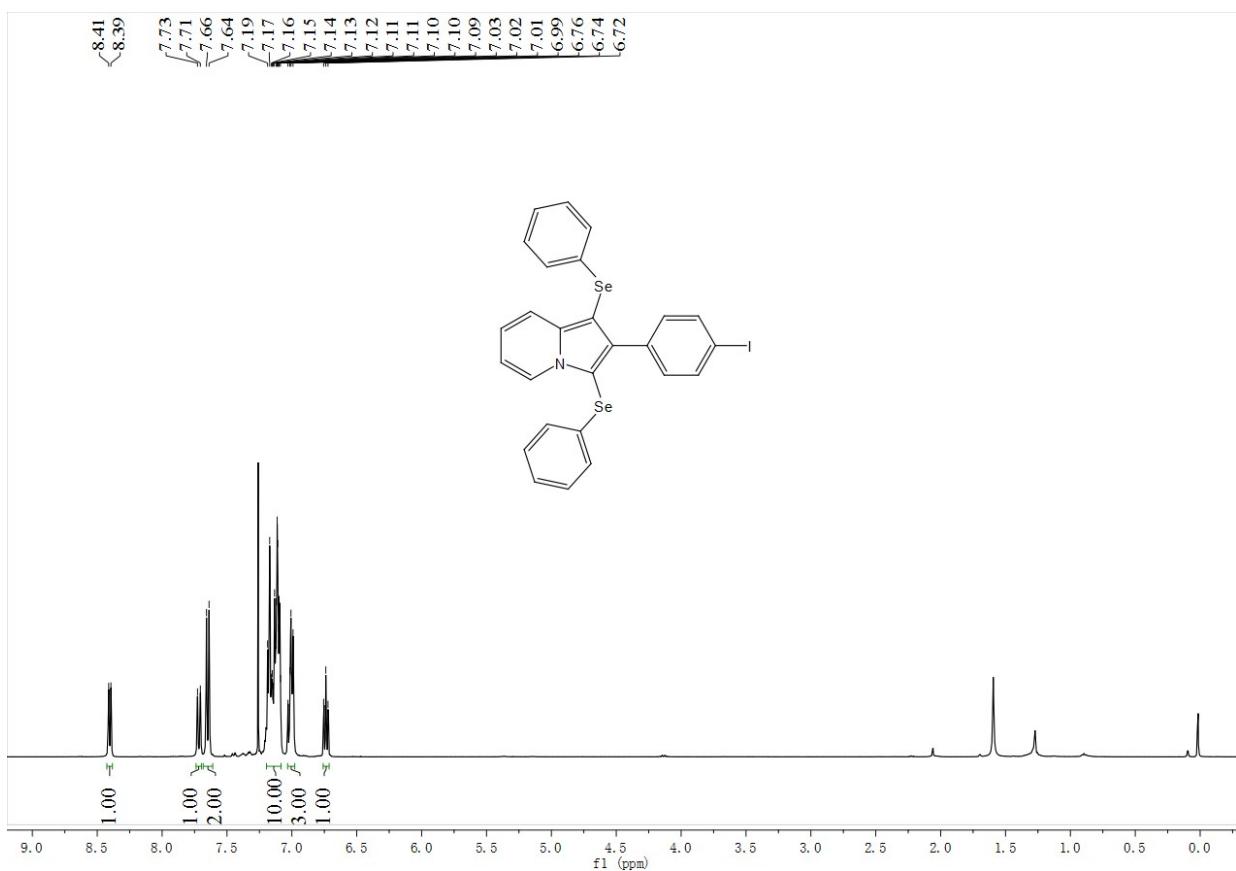
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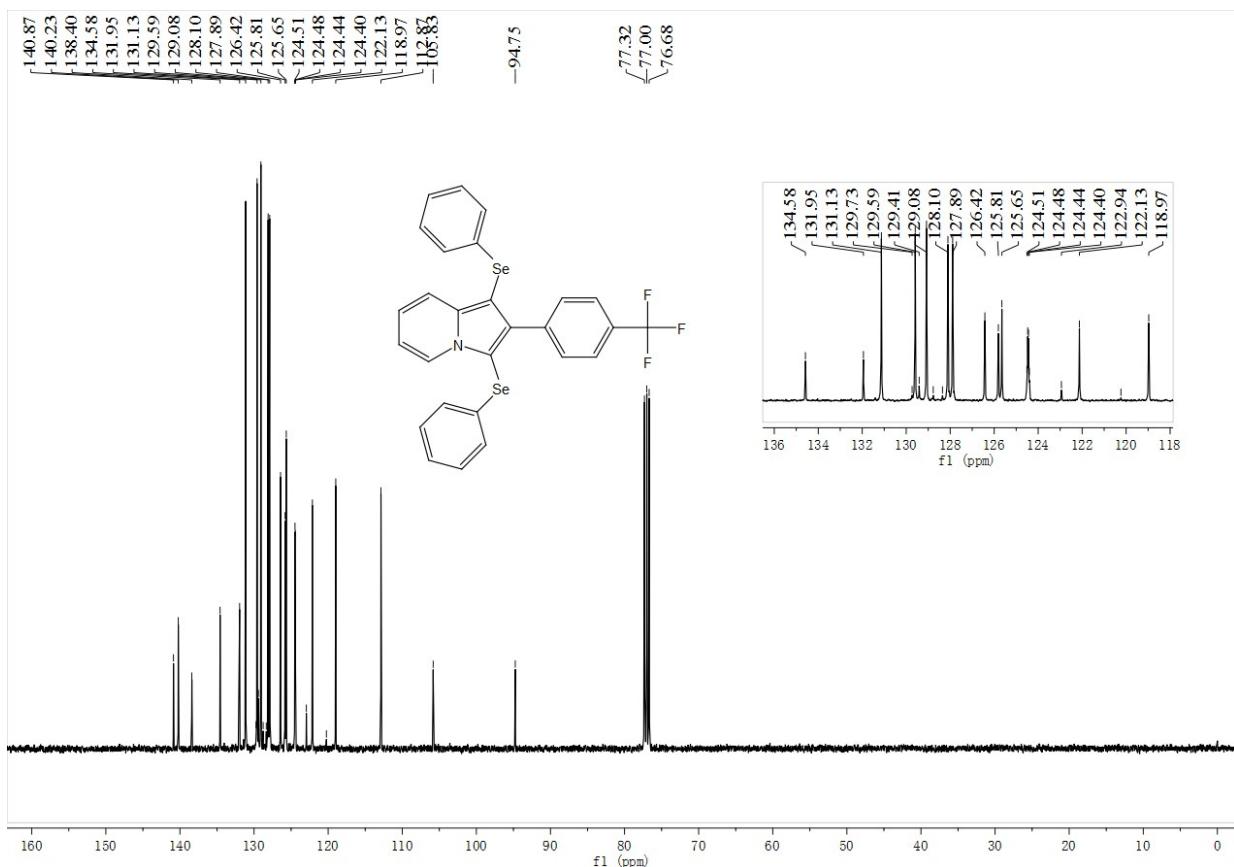
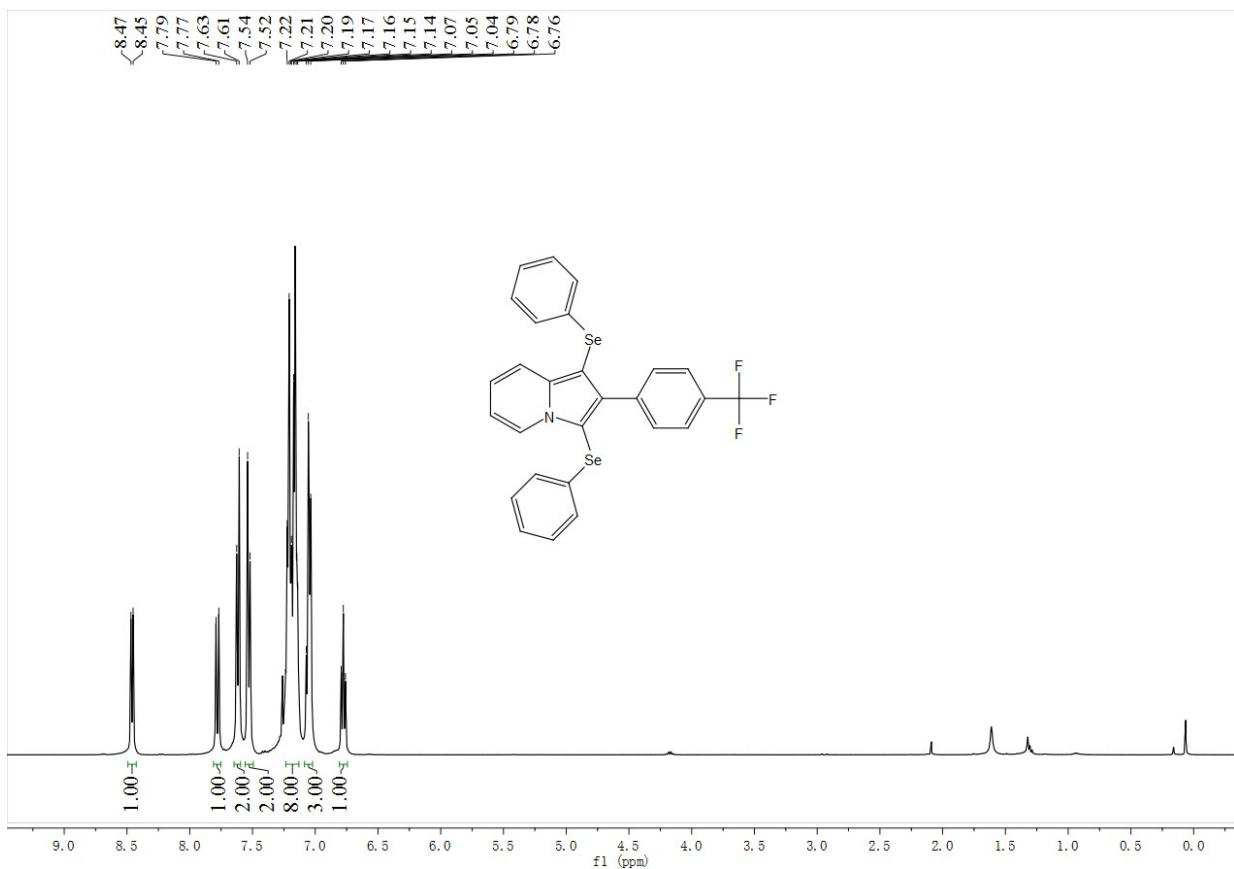
4ga



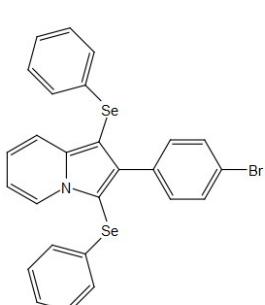
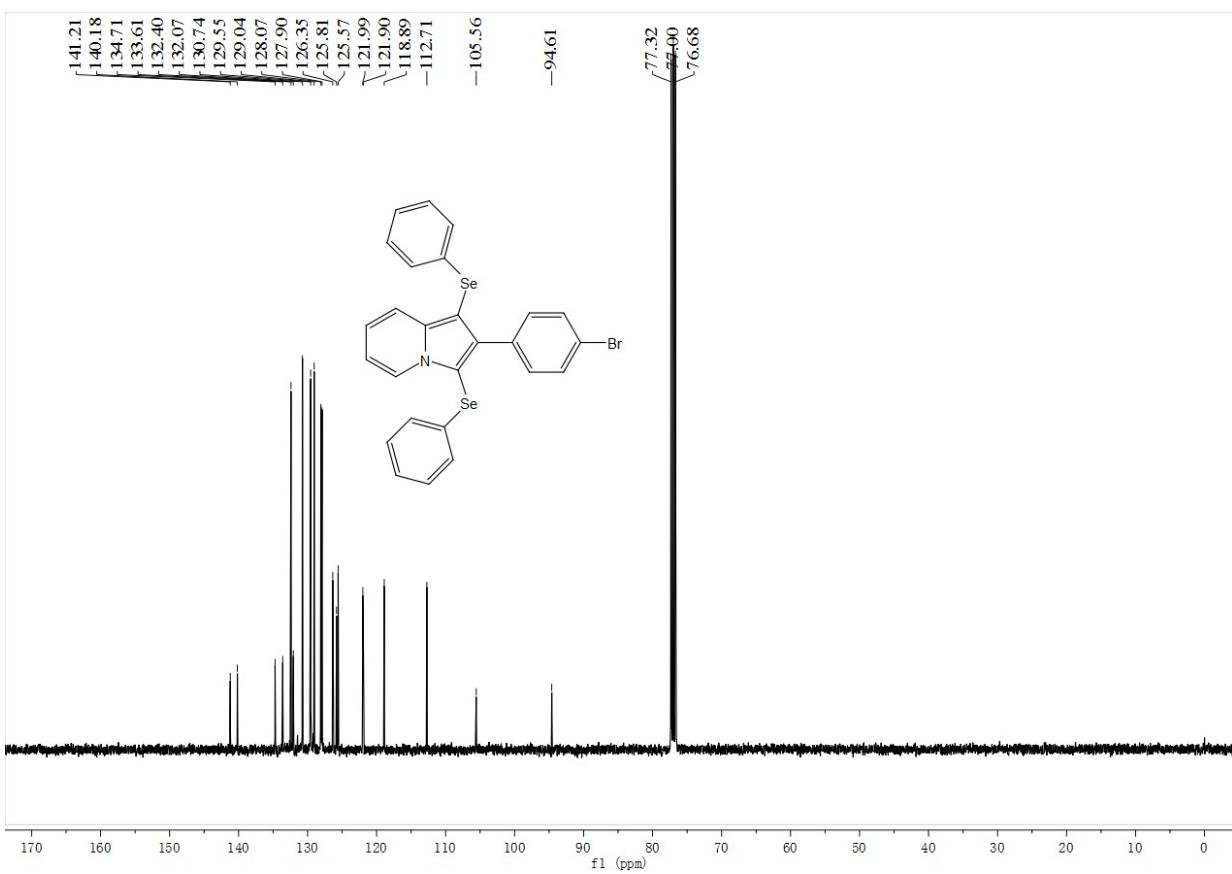
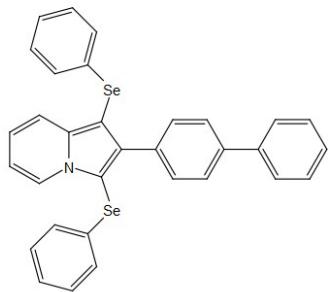
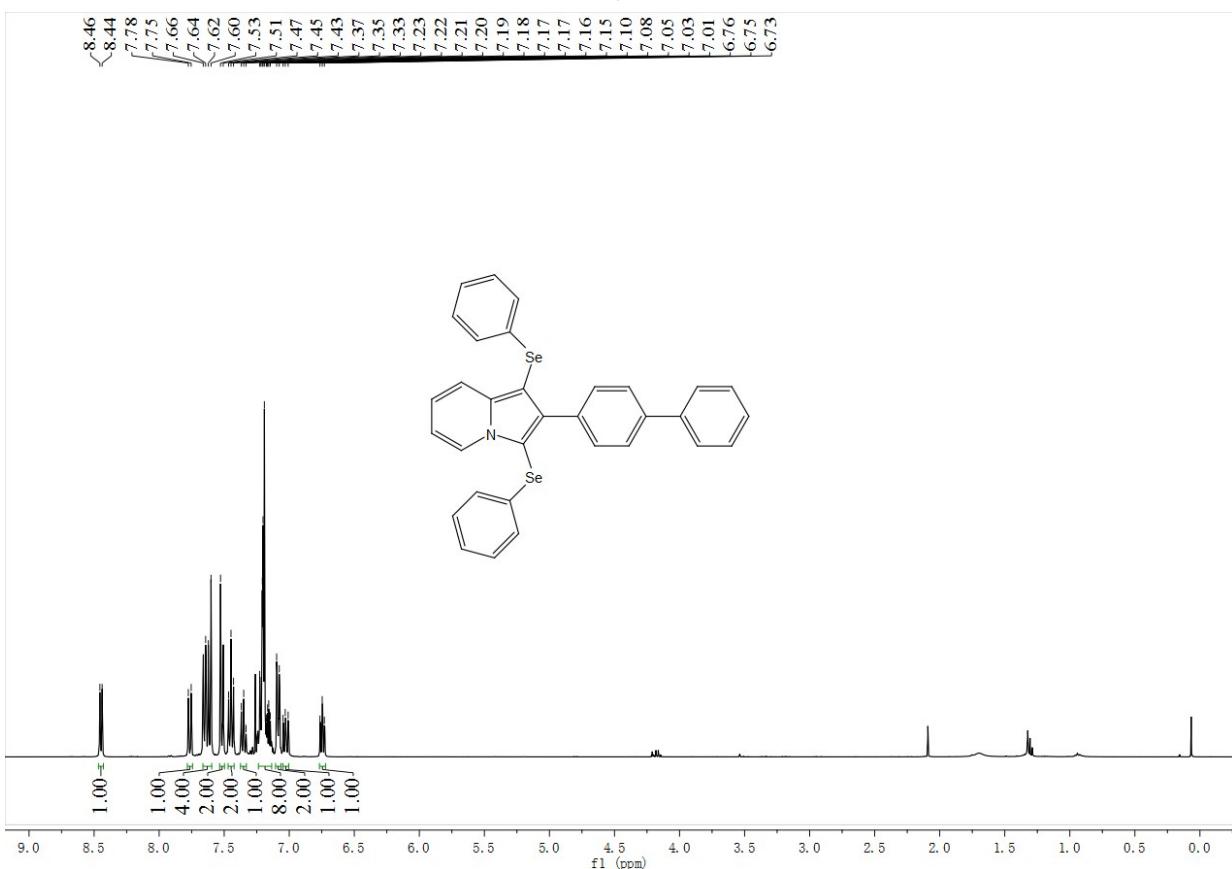
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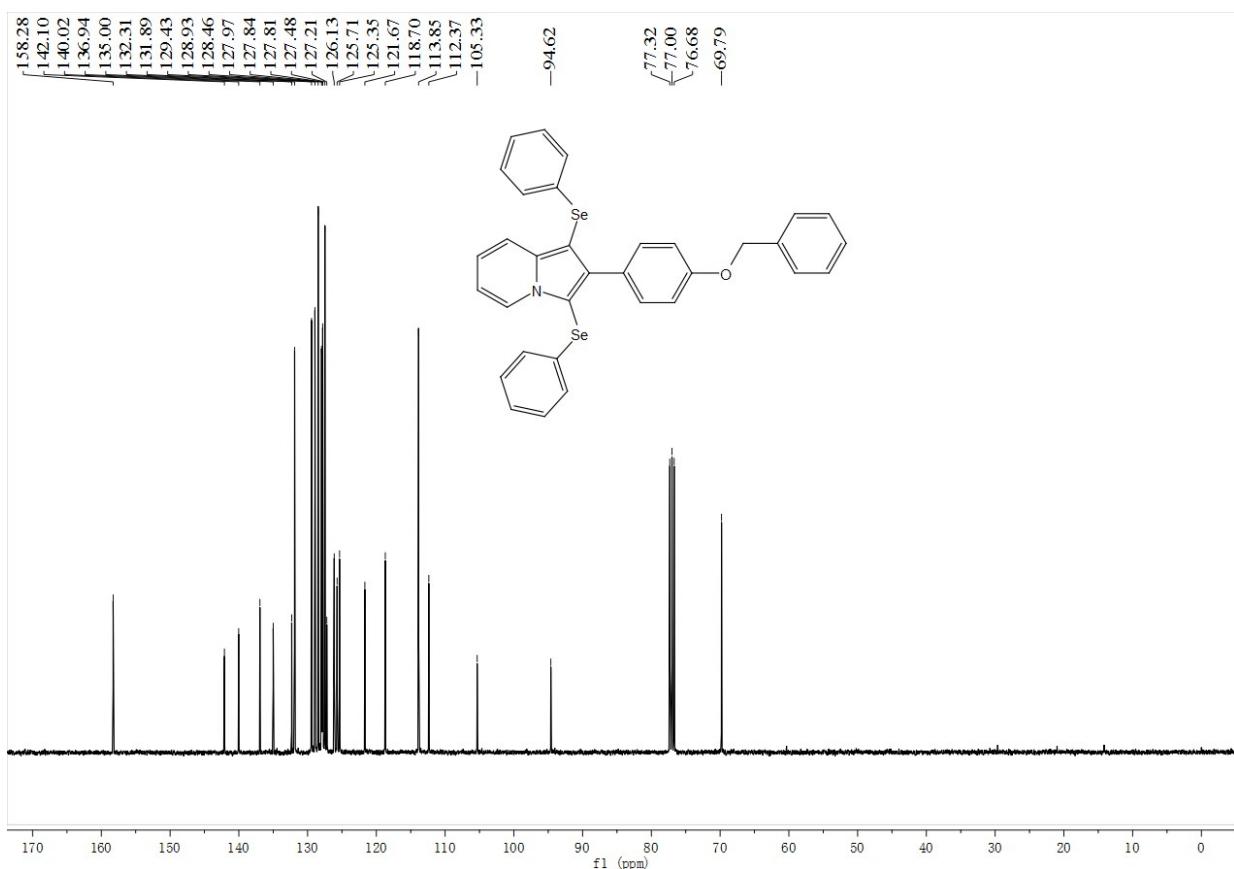
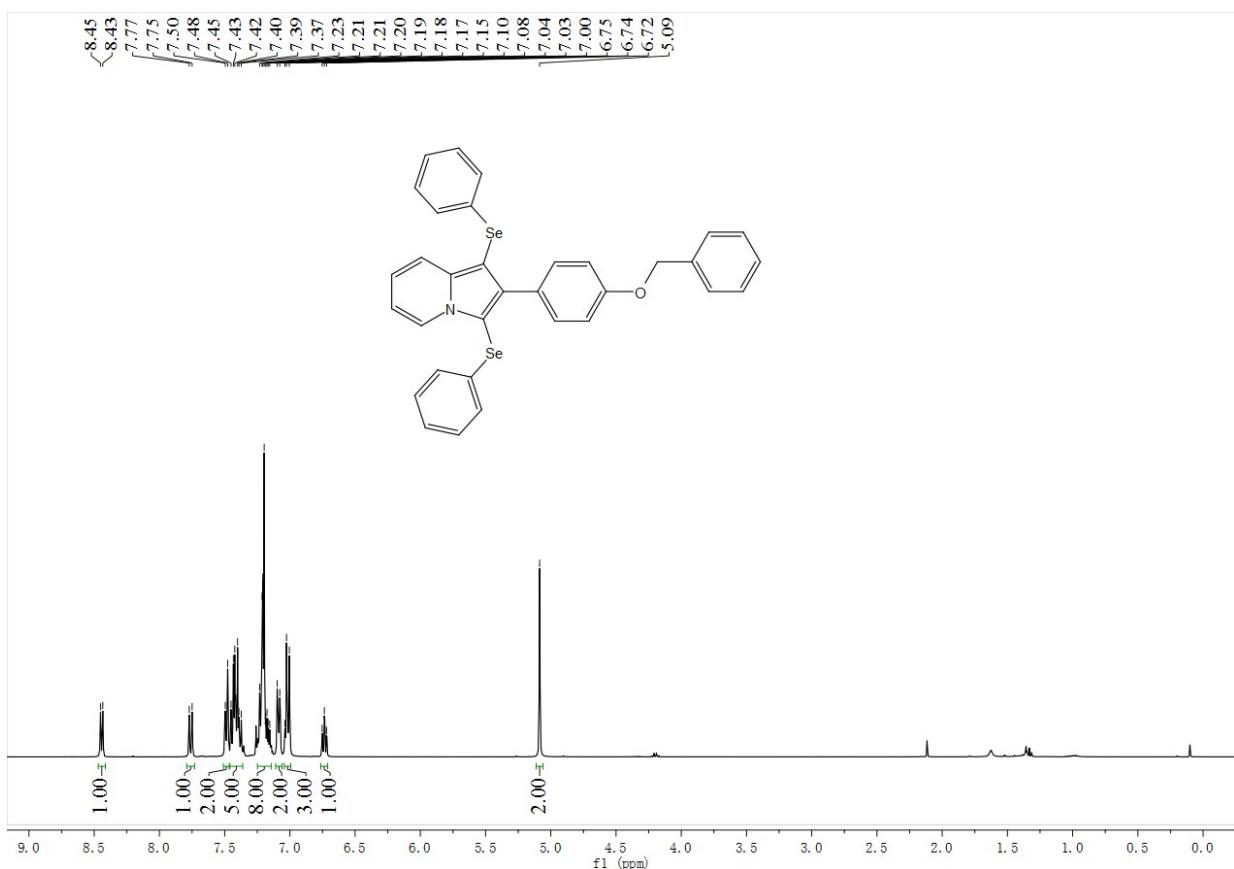
4ia



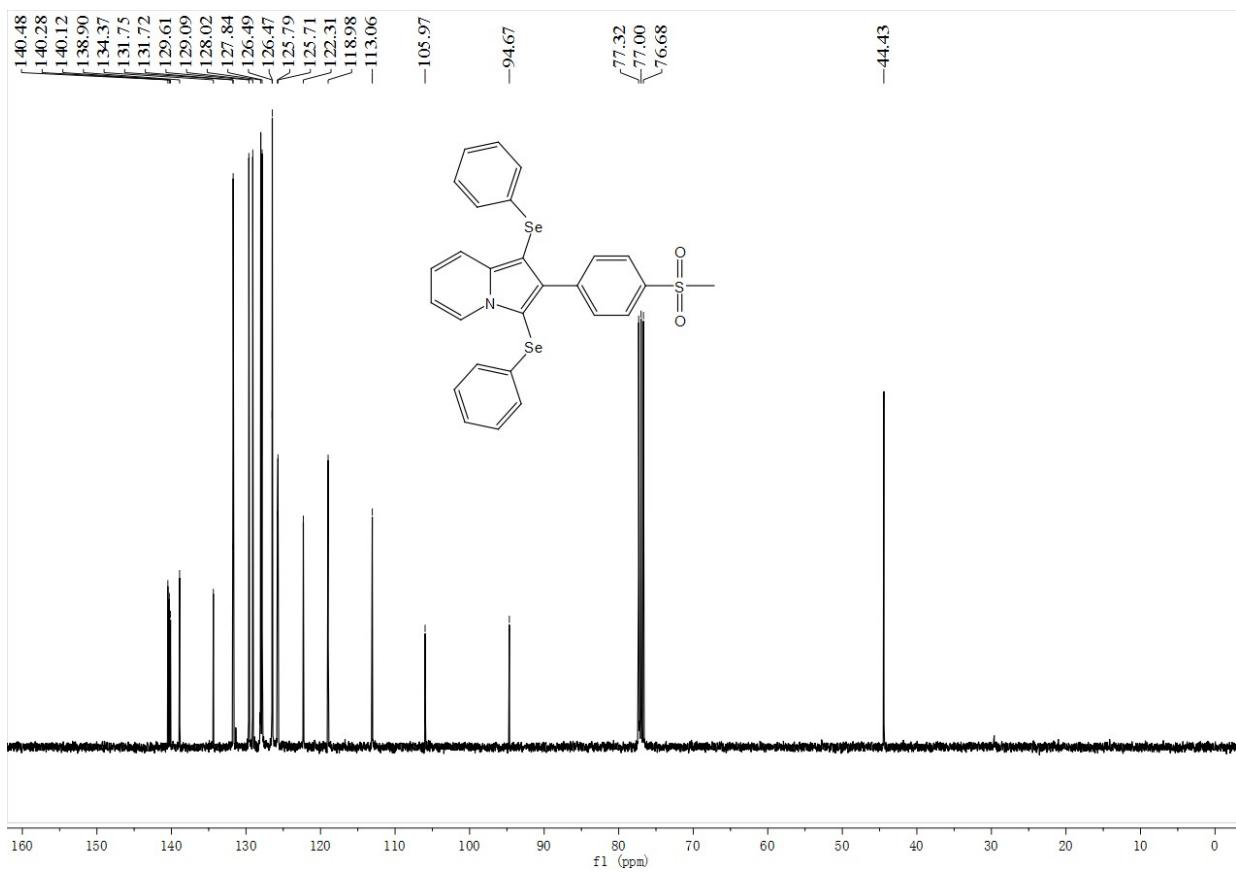
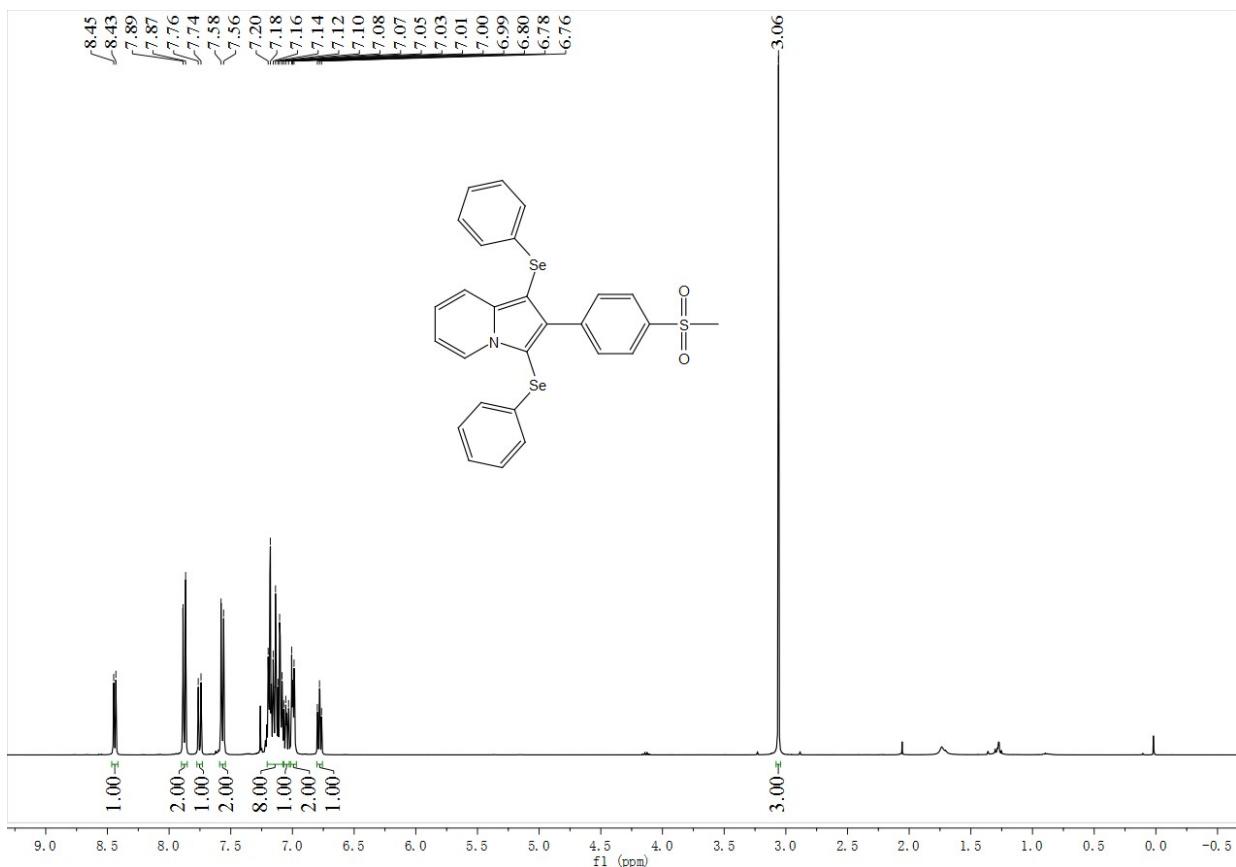
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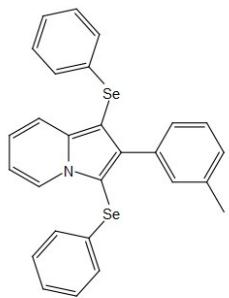
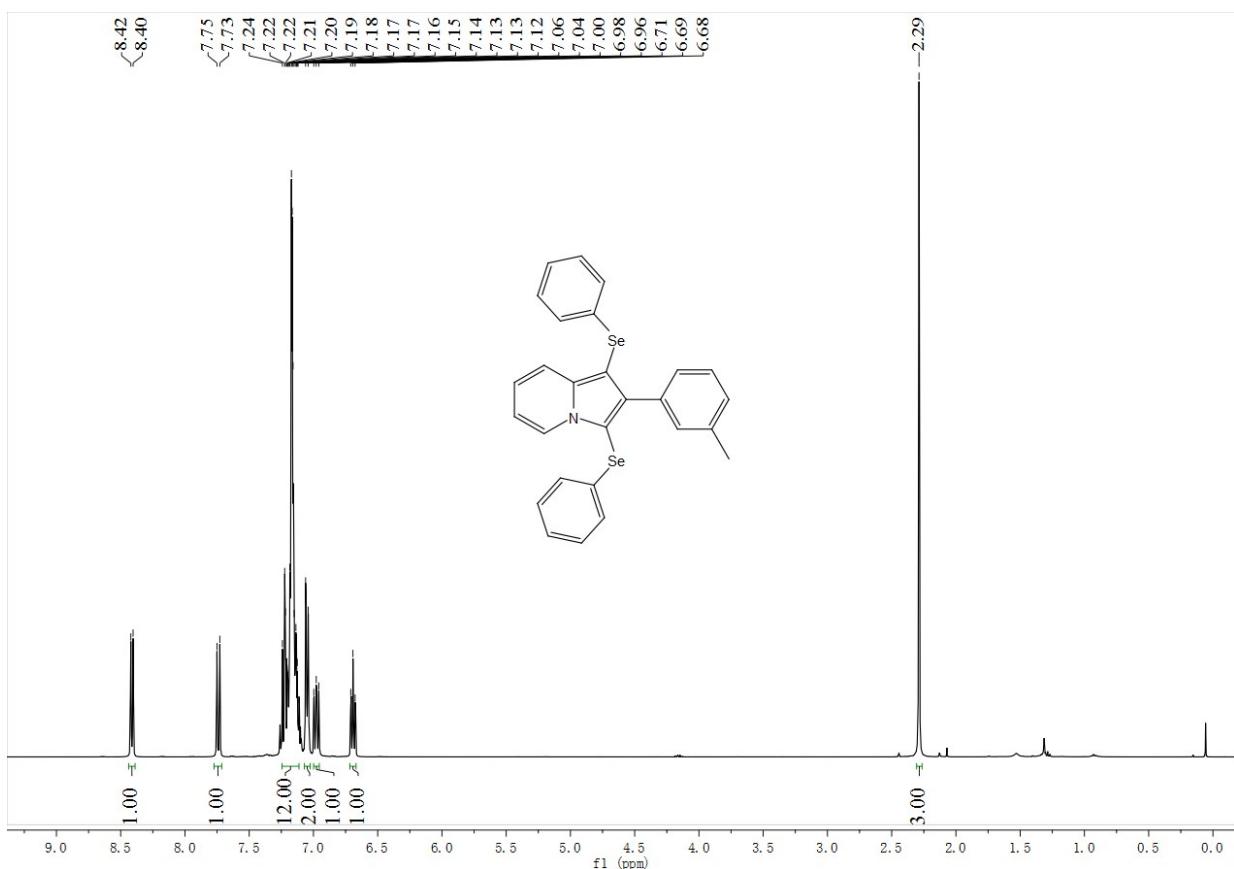
4ka



4la

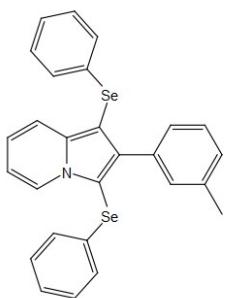


4ma

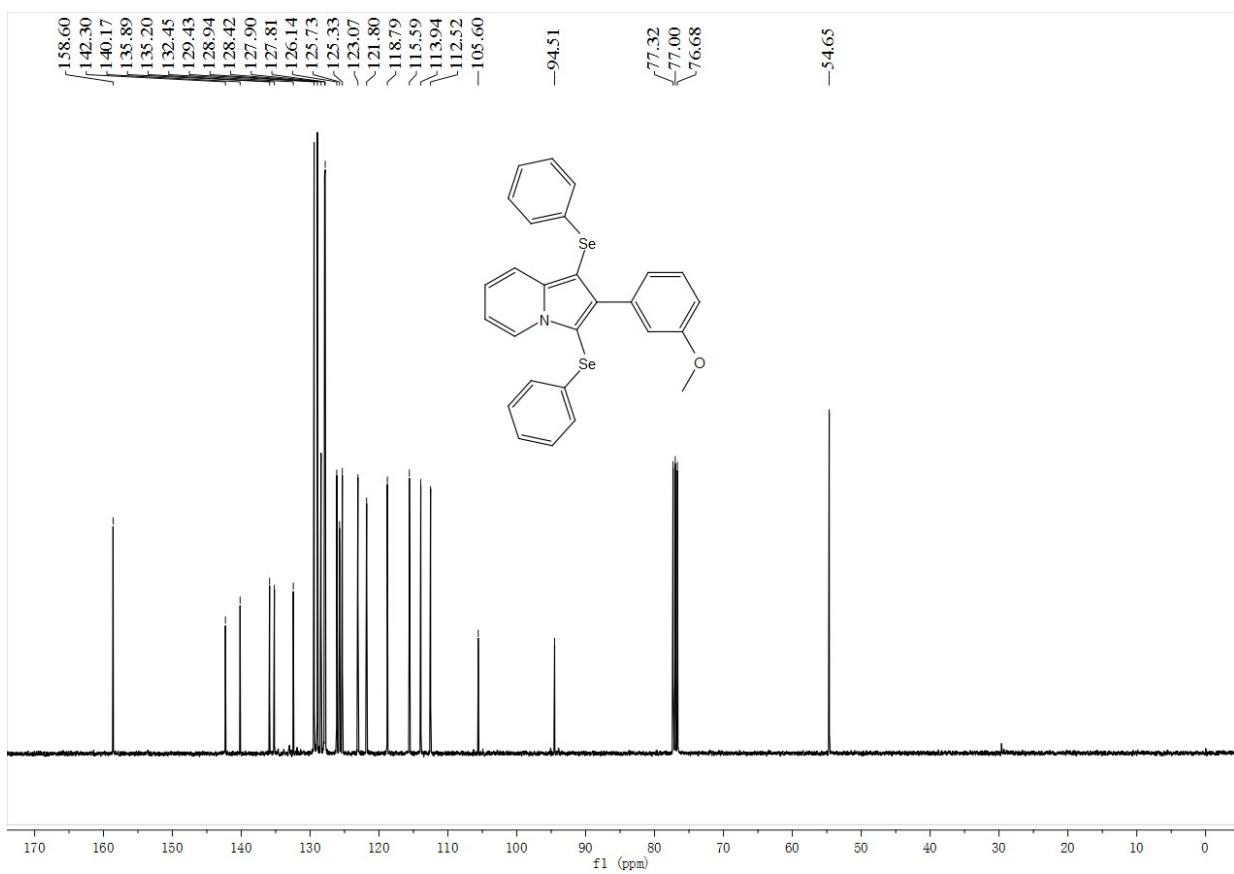
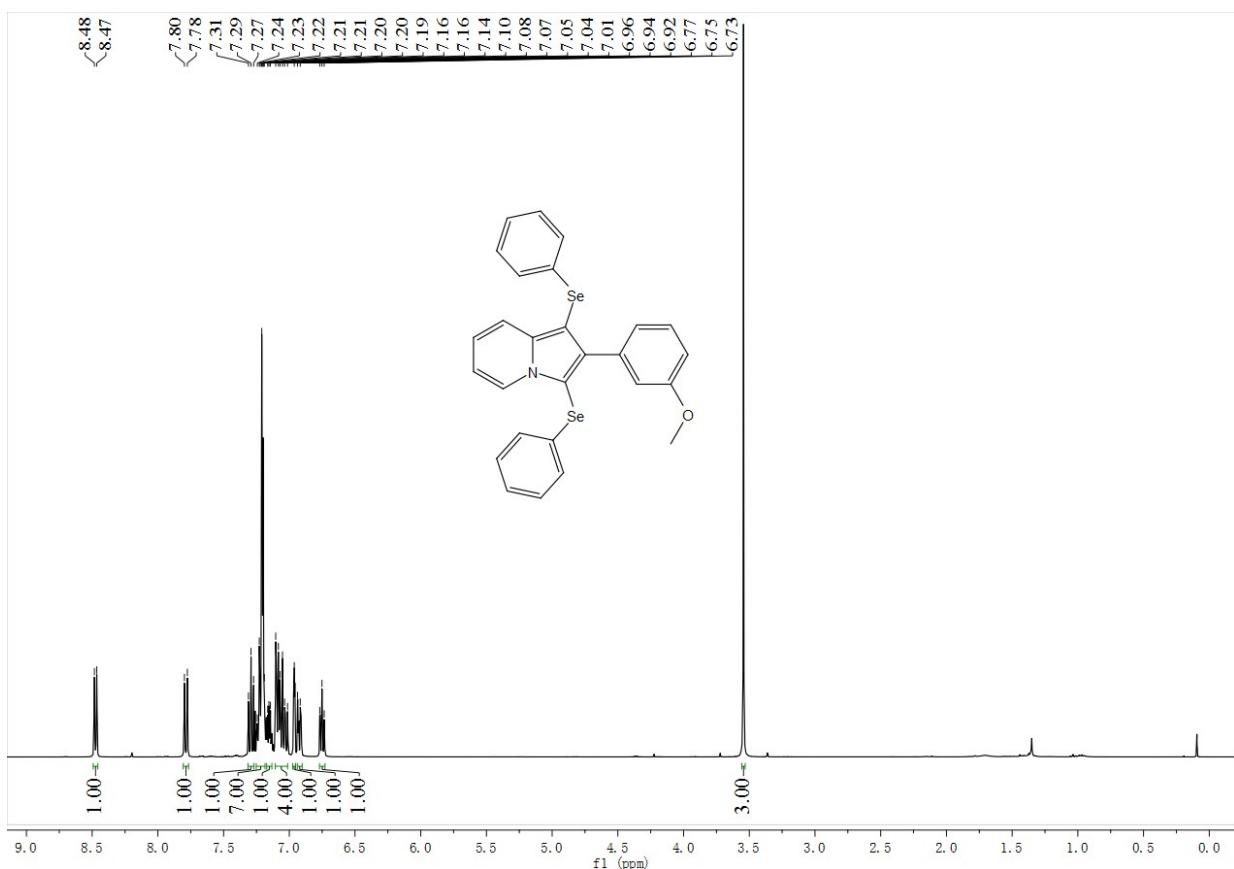


The figure displays the ^1H NMR spectrum of compound 1. The chemical structure of the compound is shown in the center, featuring a tricyclic core with two selenocyclotriphosphazene rings fused to the system. The spectrum shows several distinct sets of peaks corresponding to different chemical environments. Key peak assignments include:

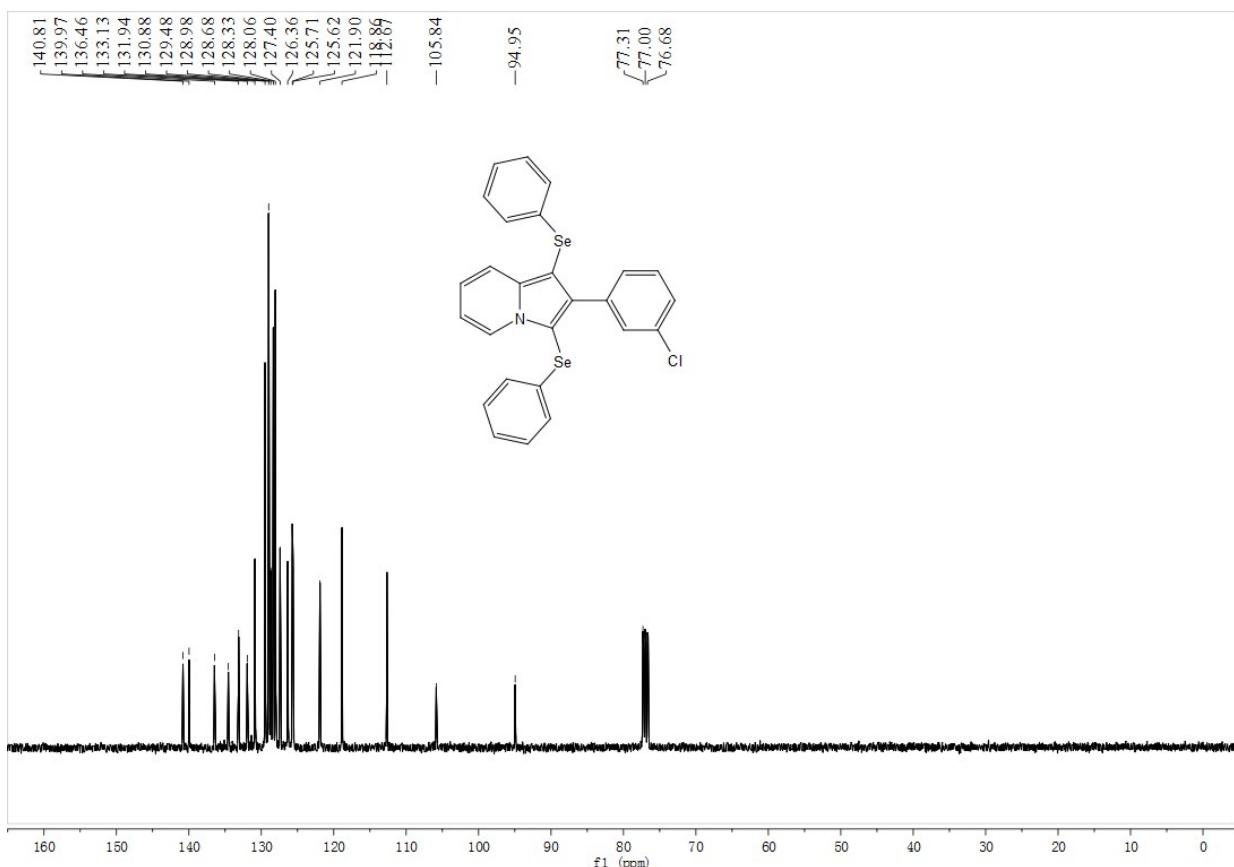
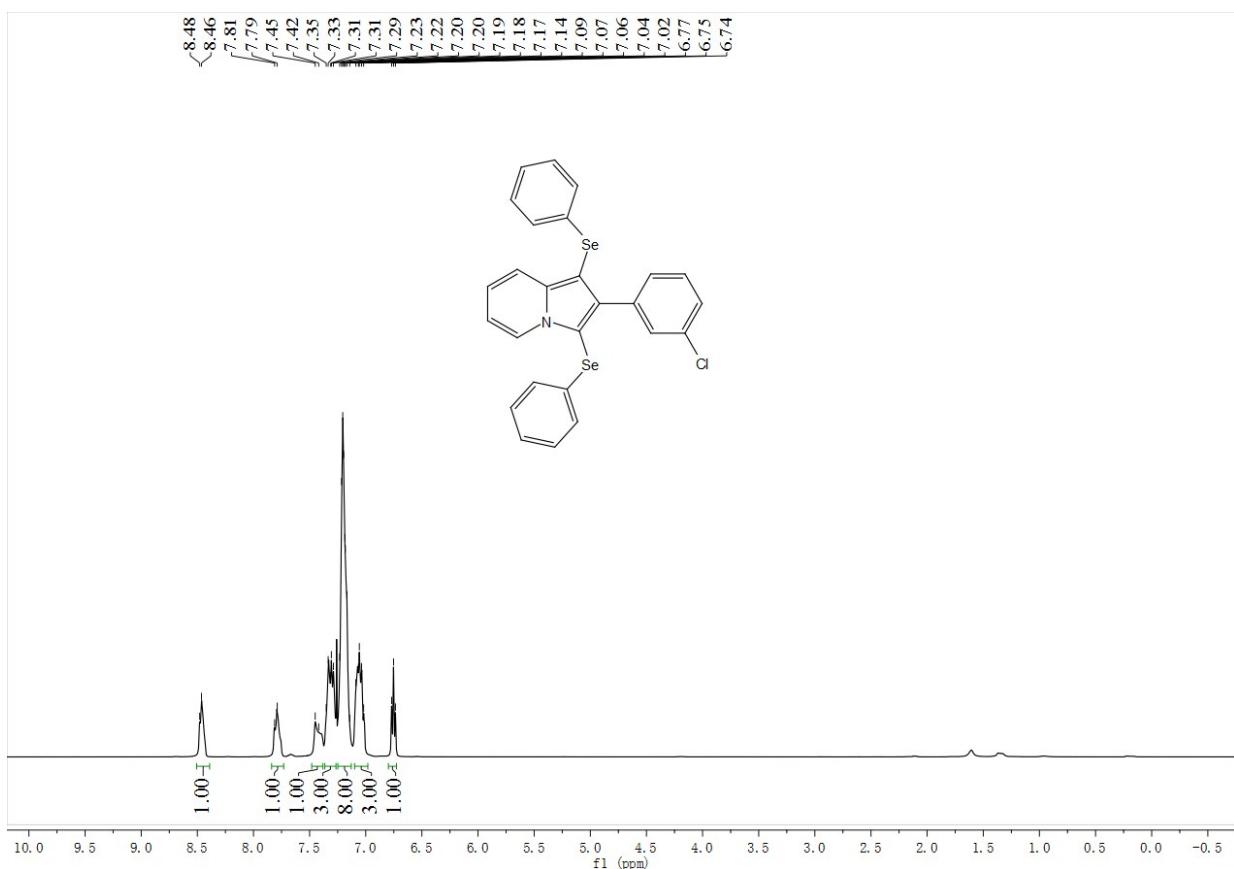
- Upfield region (0-40 ppm): A sharp peak at 21.42 ppm is assigned to the methyl protons (CH_3) of the 4,4'-dimethylbiphenyl group.
- Middle region (60-100 ppm): A complex multiplet between 76.68 and 77.32 ppm is assigned to the aromatic protons of the phenyl groups. A singlet at 94.94 ppm is assigned to the selenium atom (SeH).
- Downfield region (110-170 ppm): A large cluster of peaks between 112.38 and 142.62 ppm is assigned to the aliphatic protons of the tricyclic core and the 4,4'-dimethylbiphenyl group.



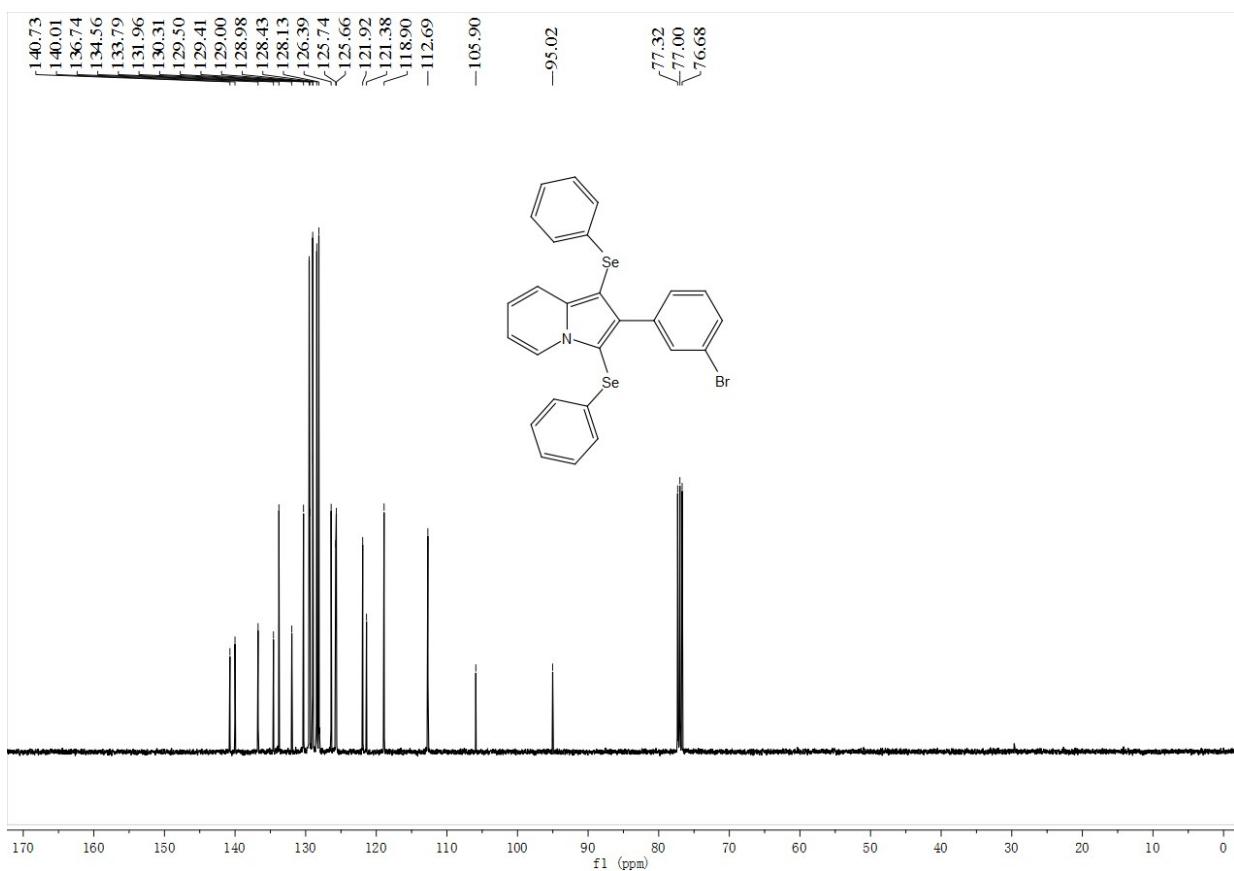
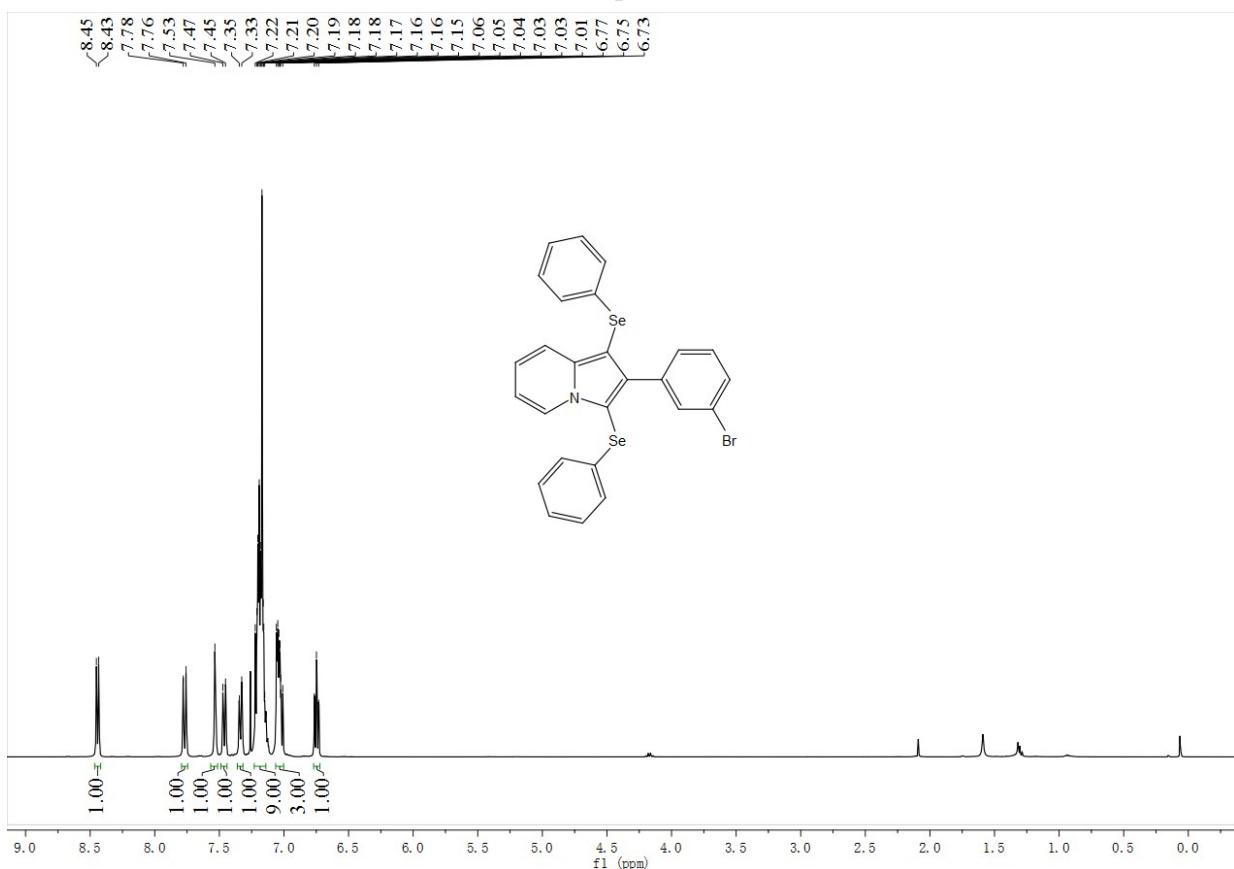
4na



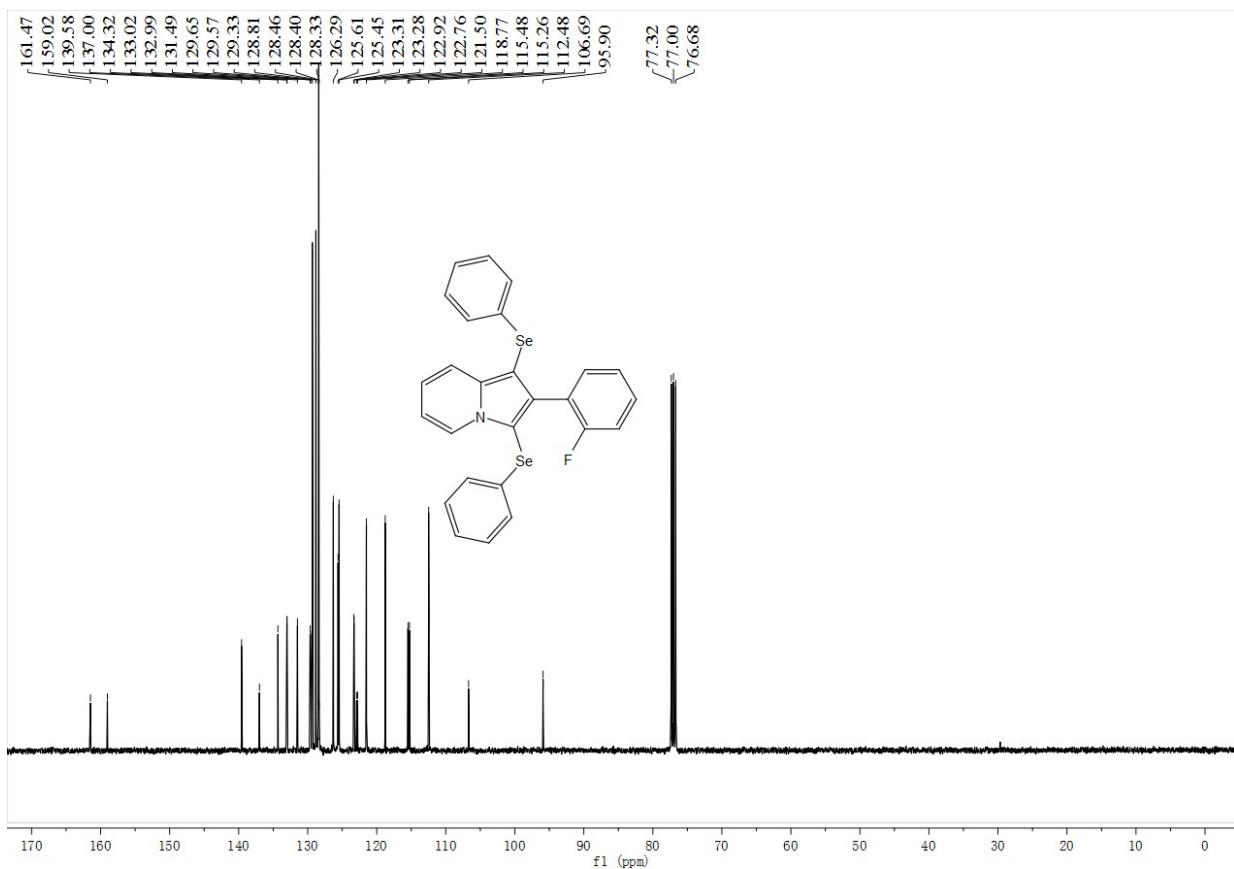
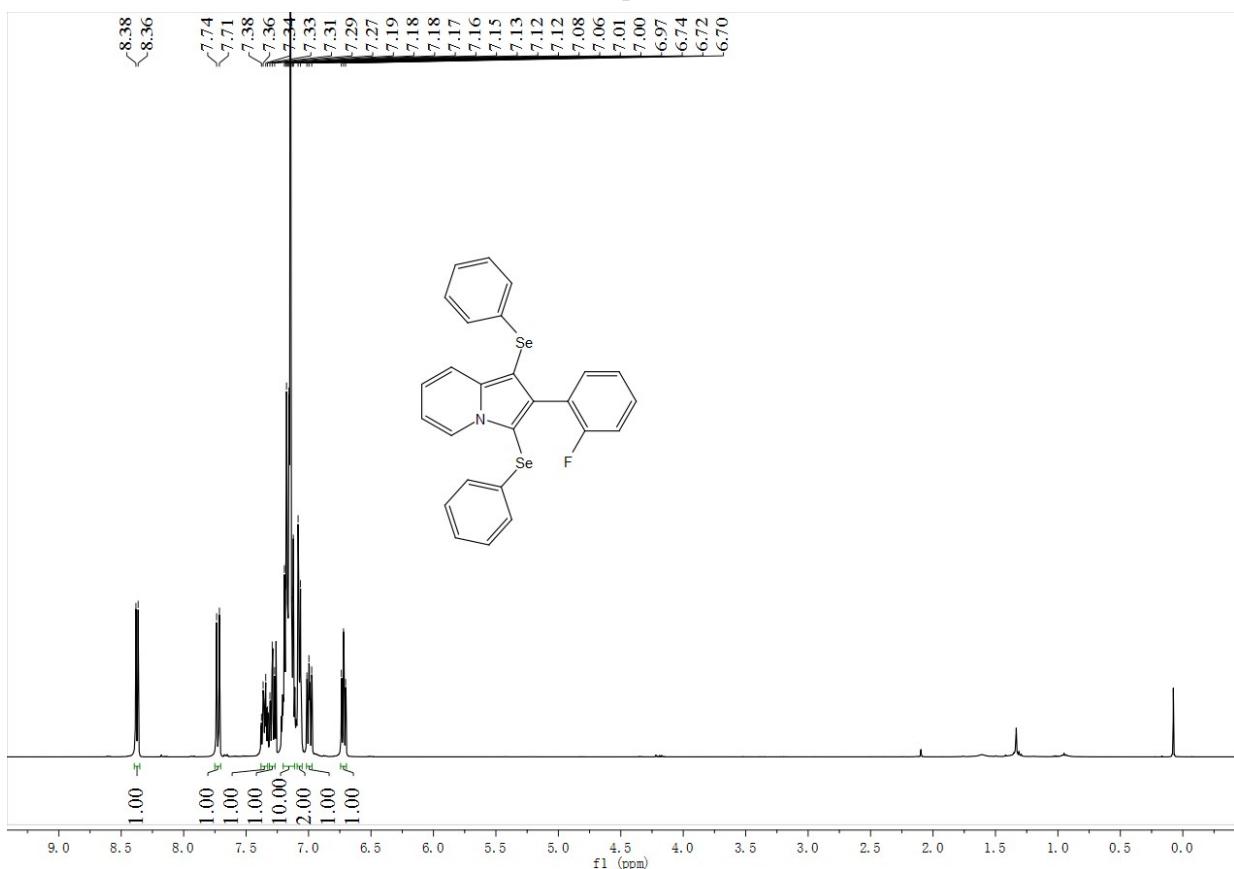
4oa



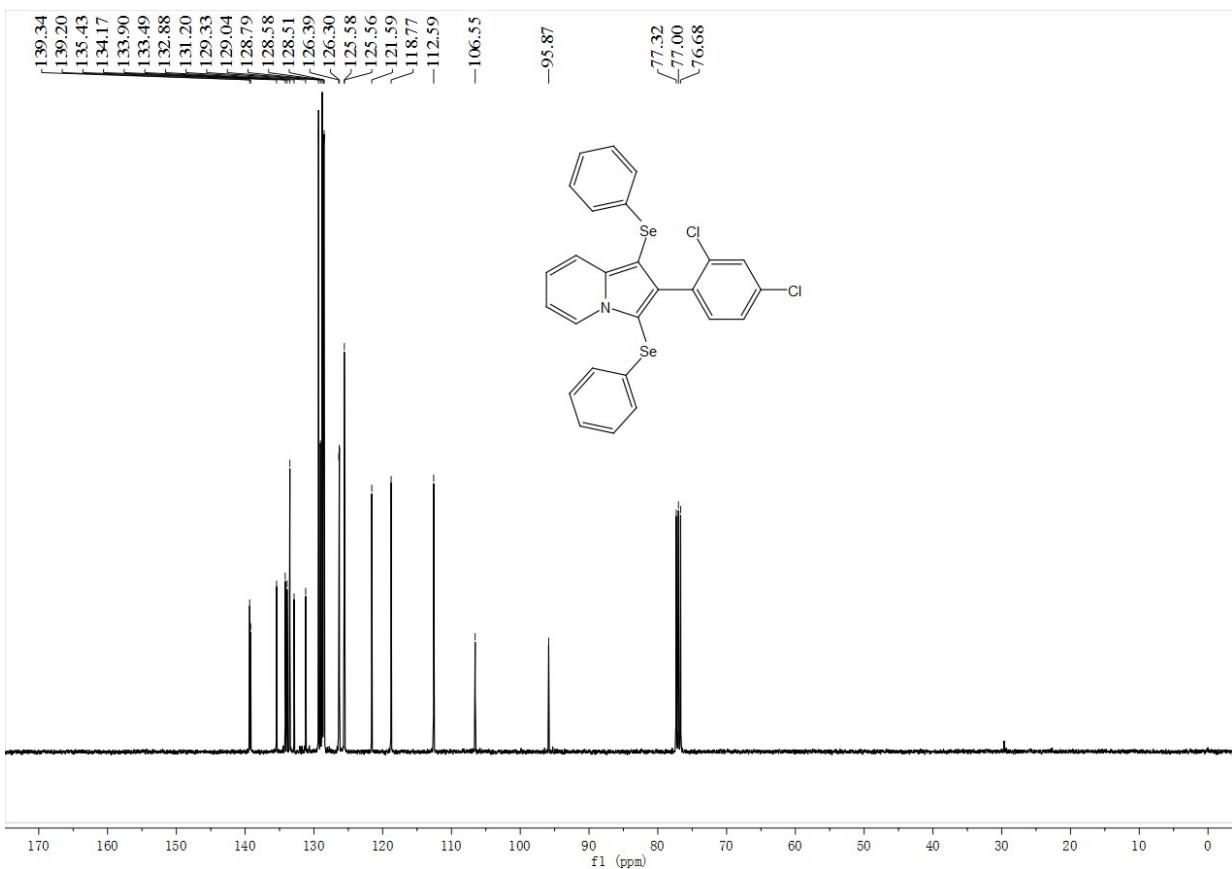
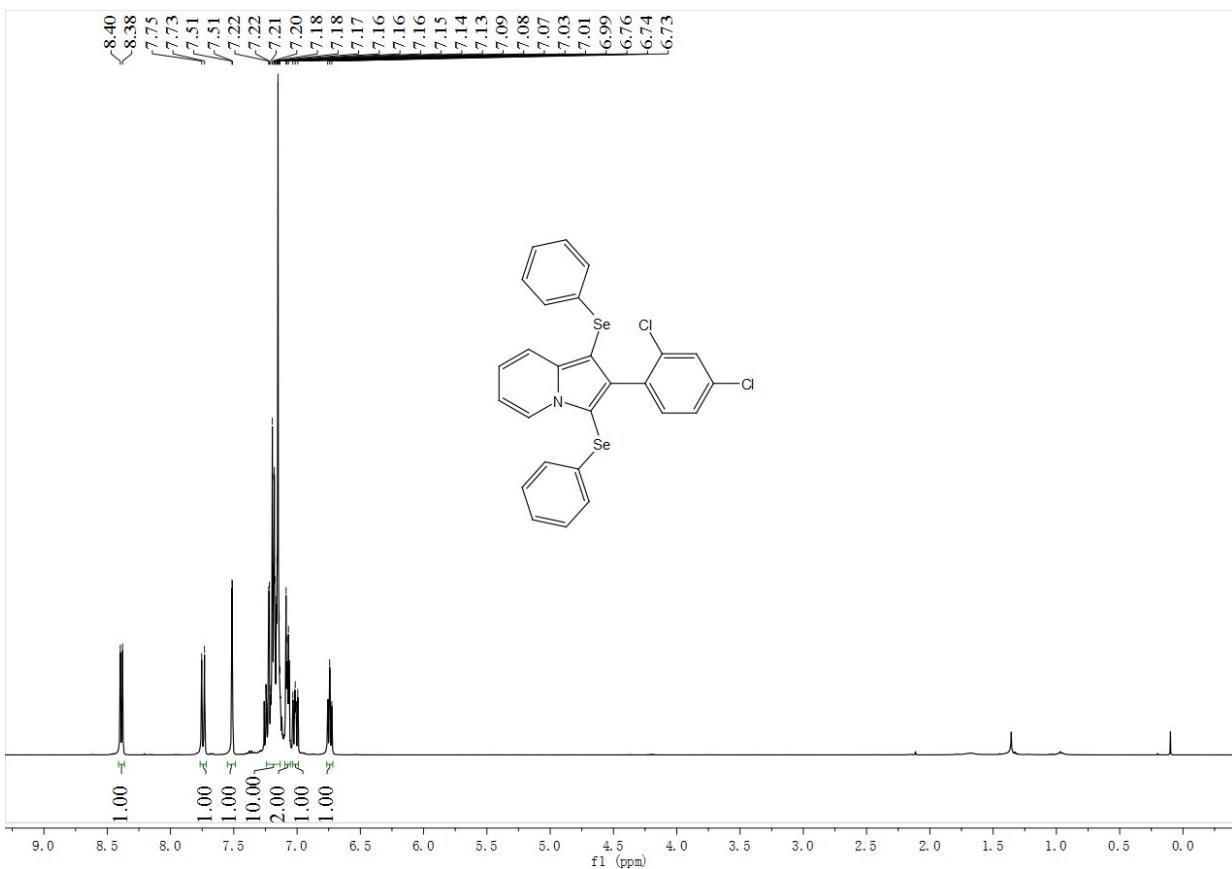
4pa



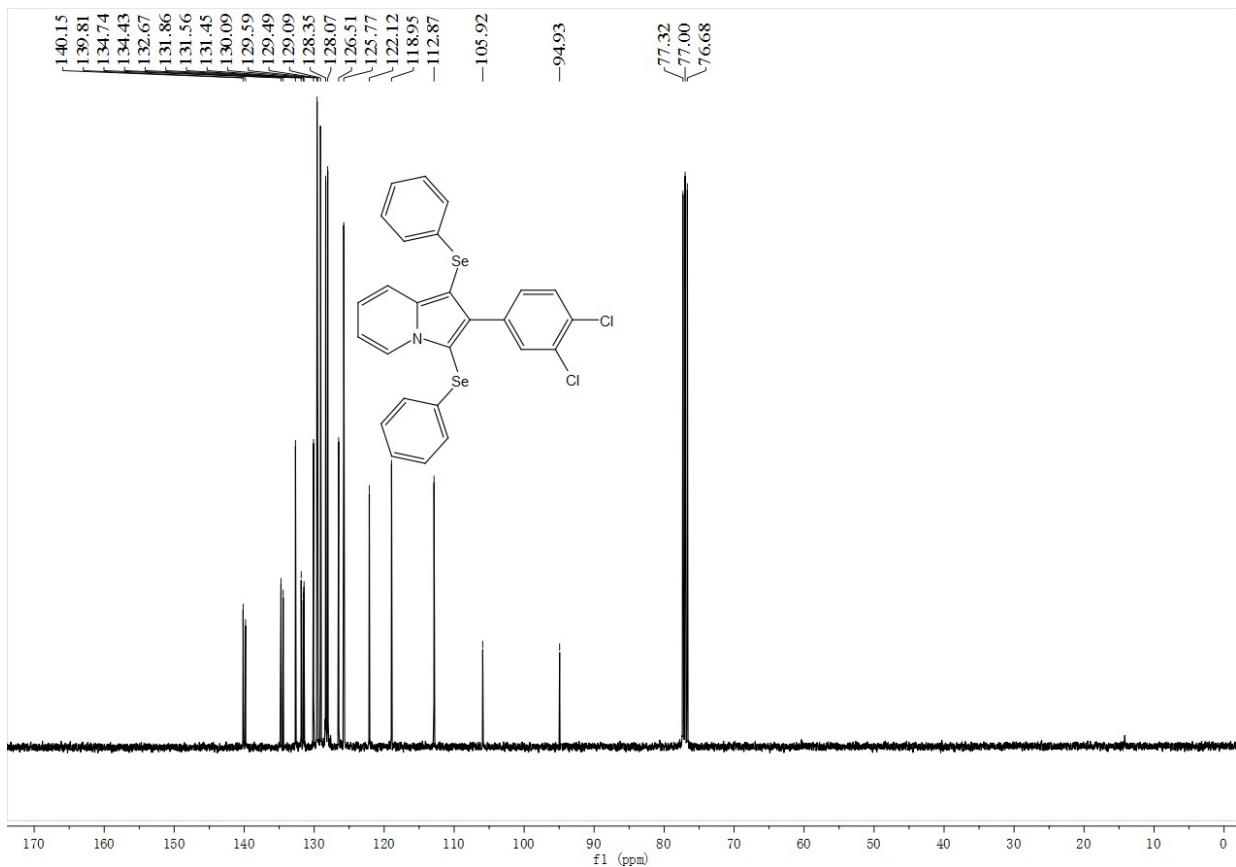
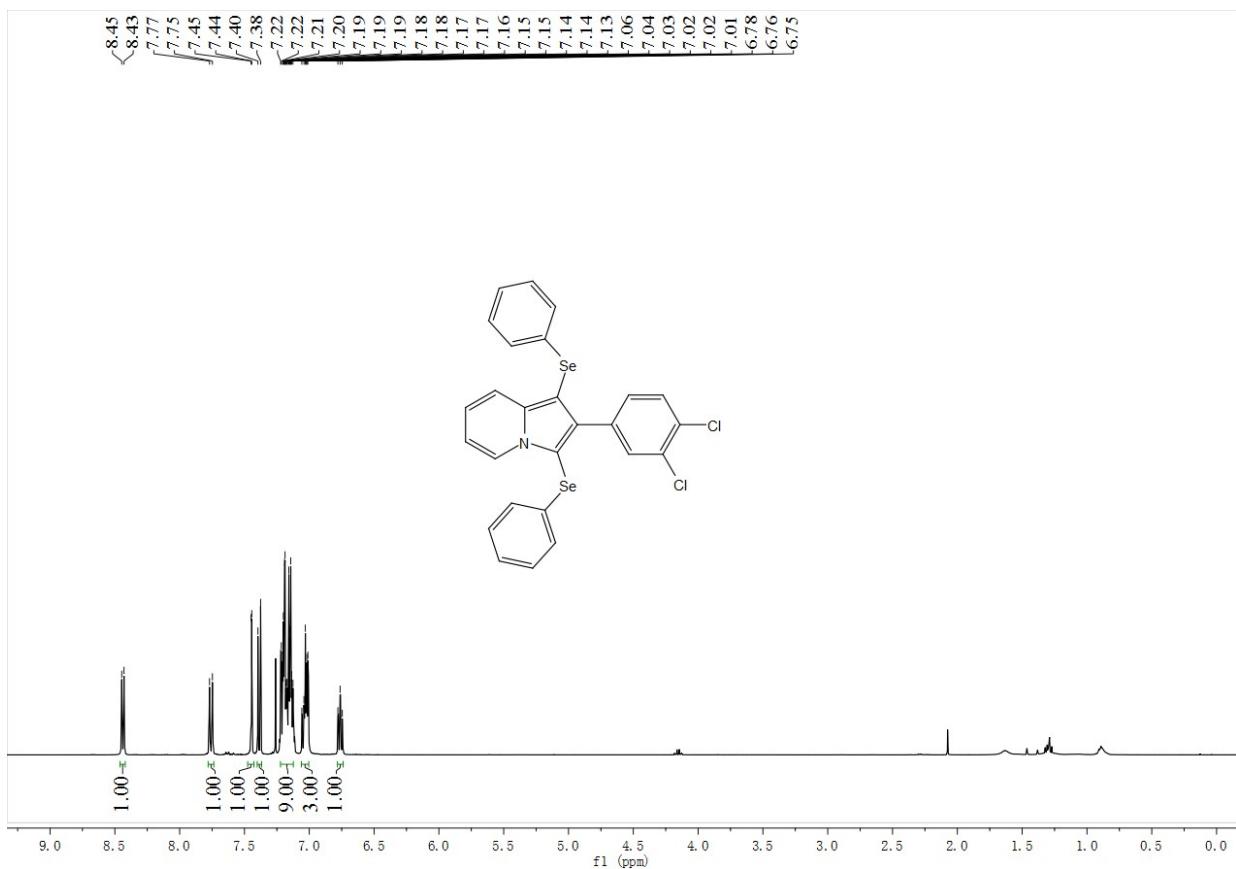
4qa



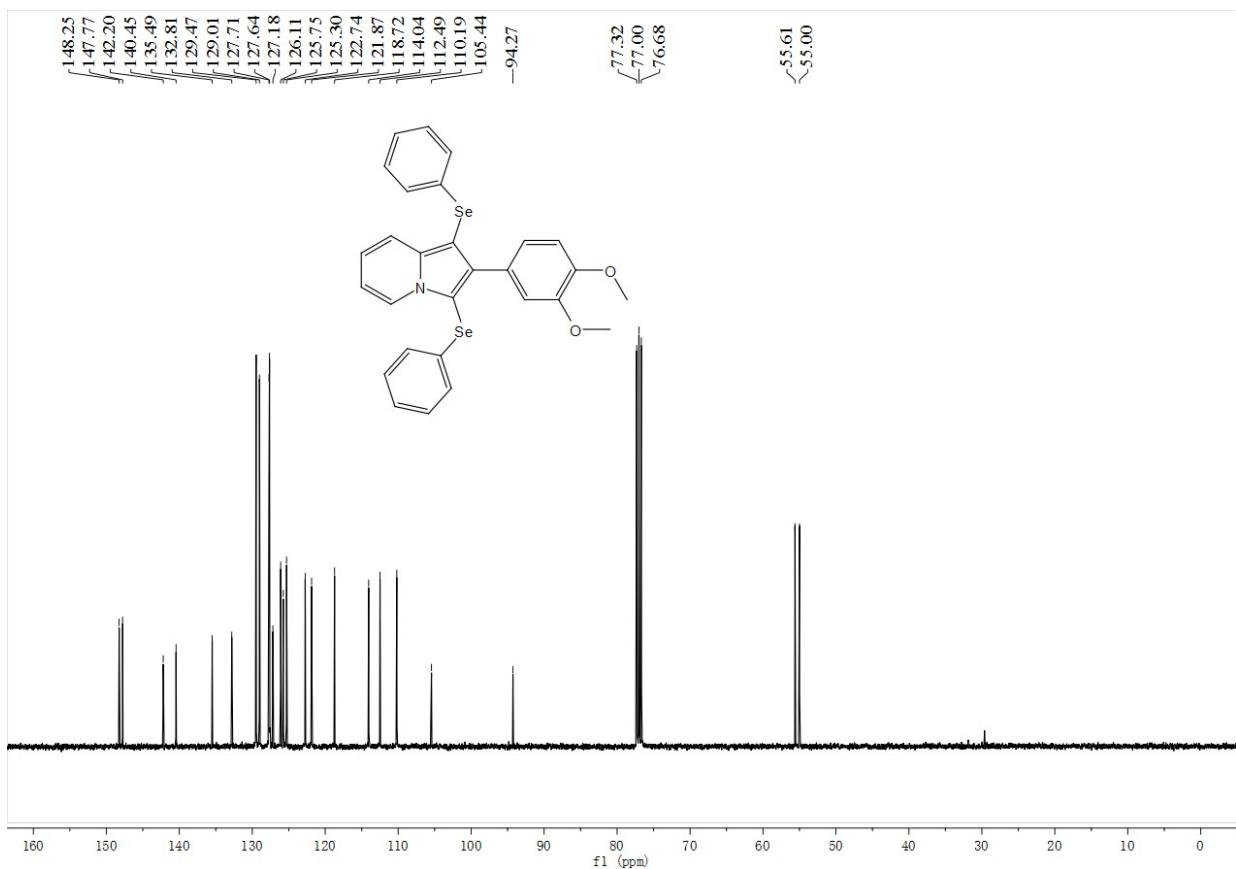
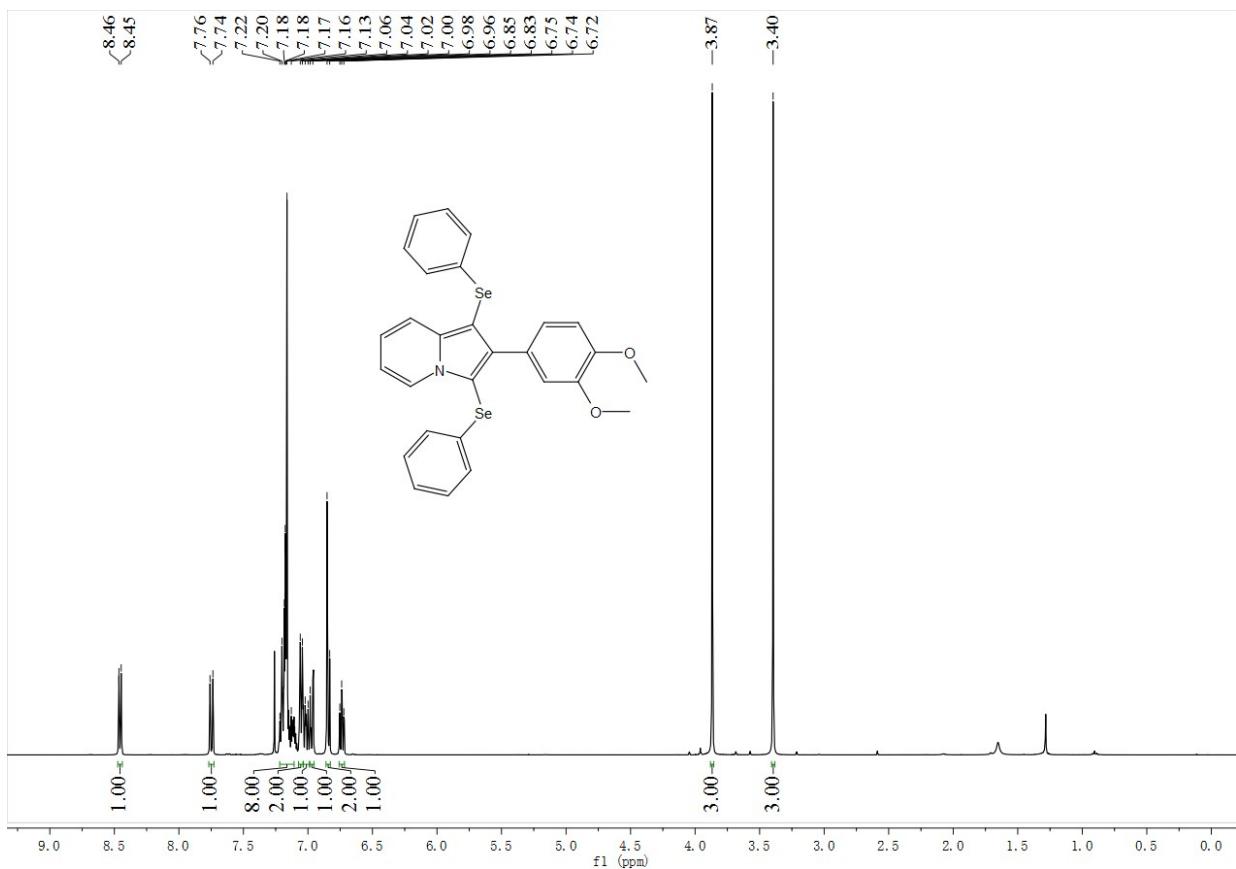
4ra



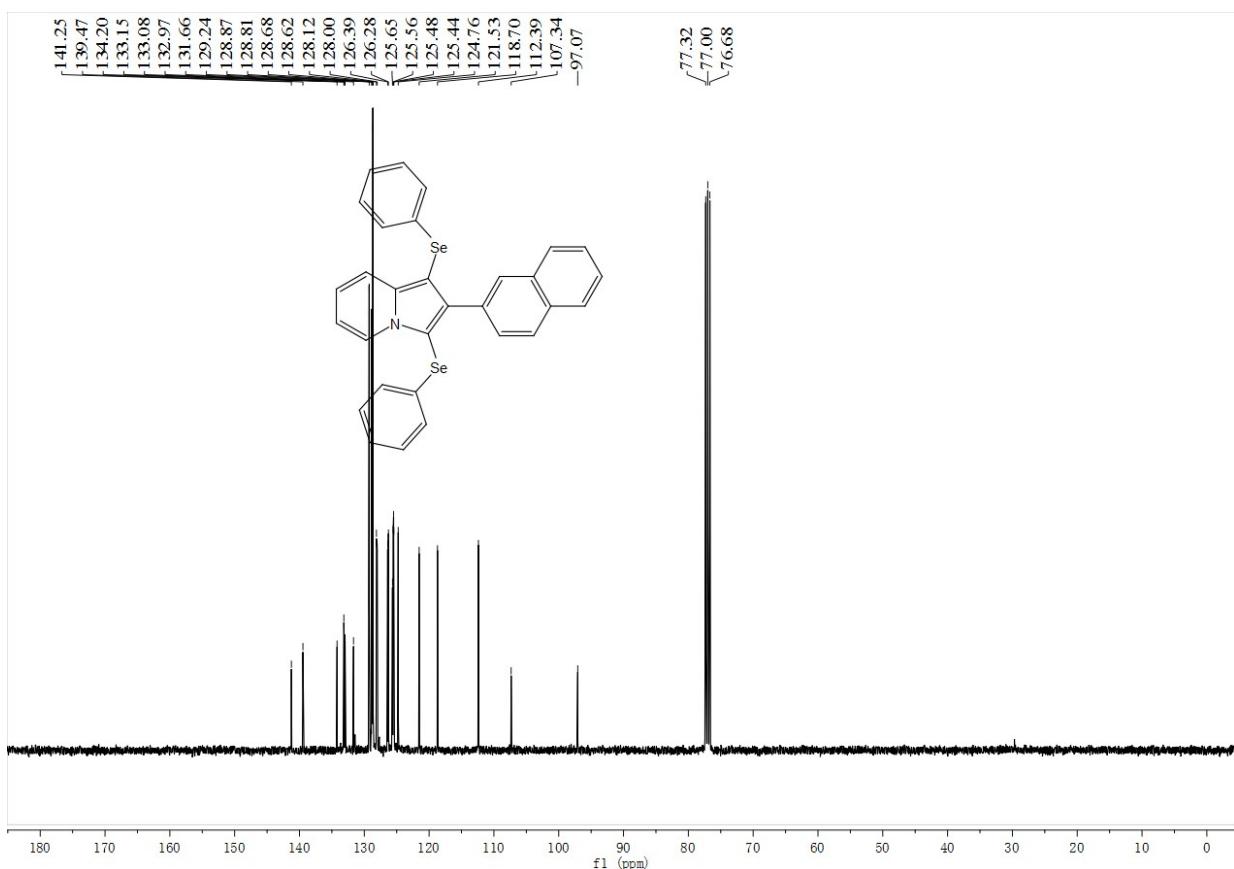
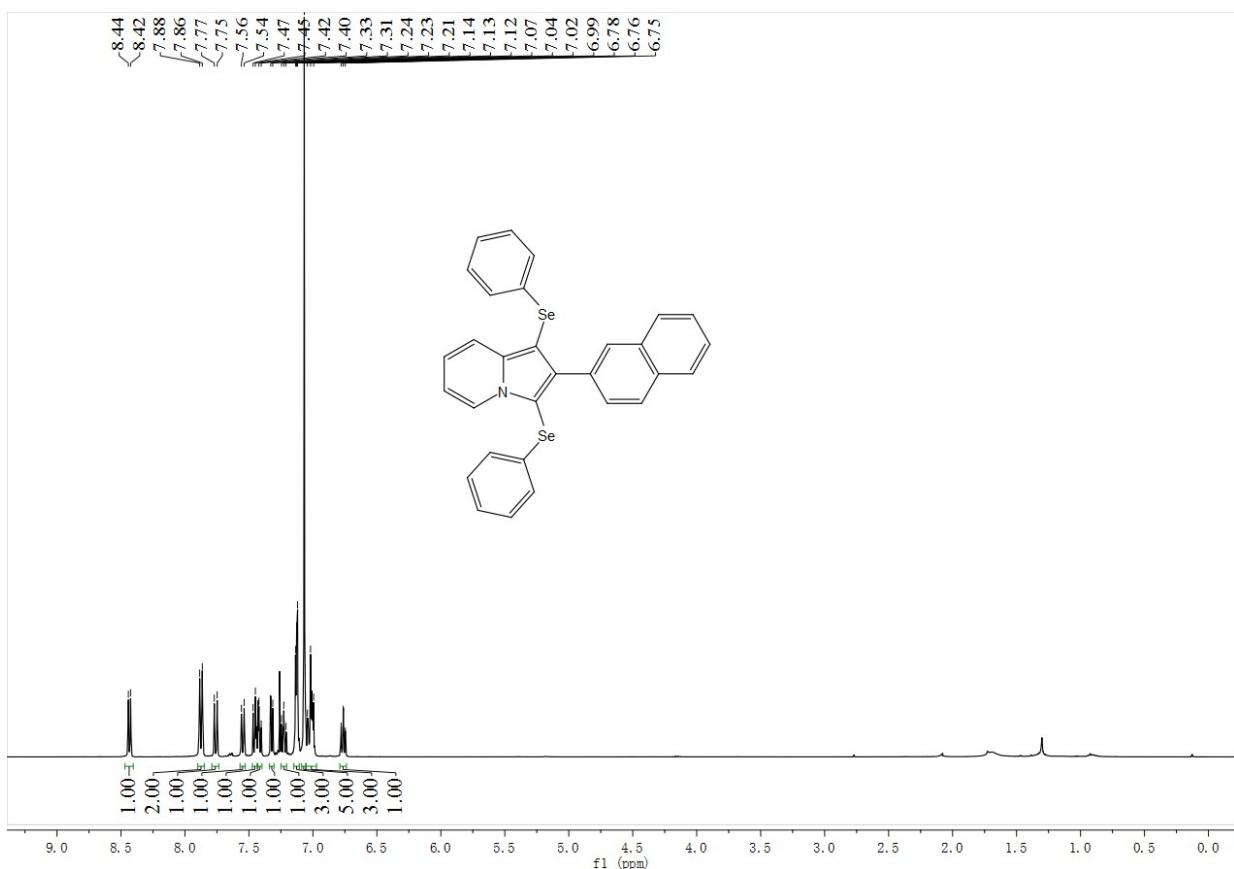
4sa



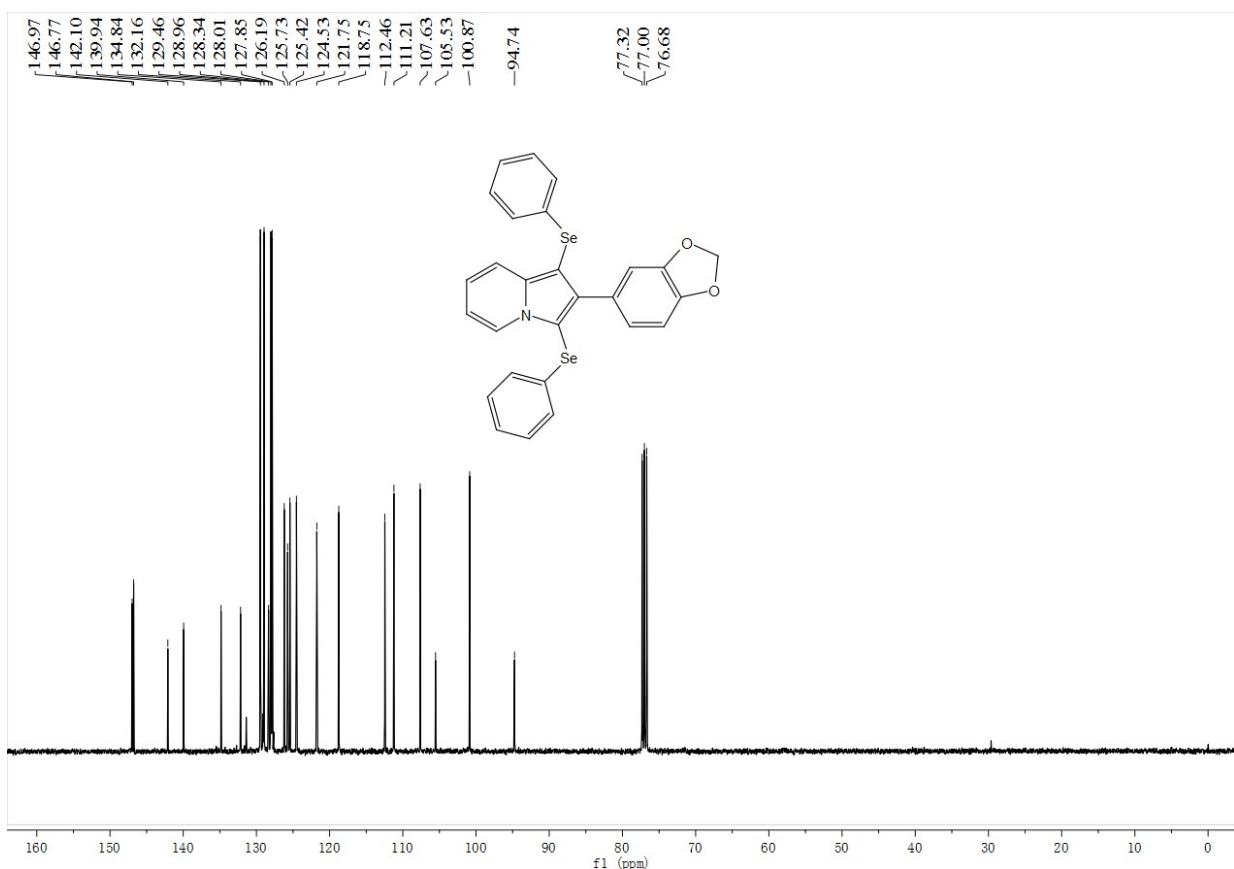
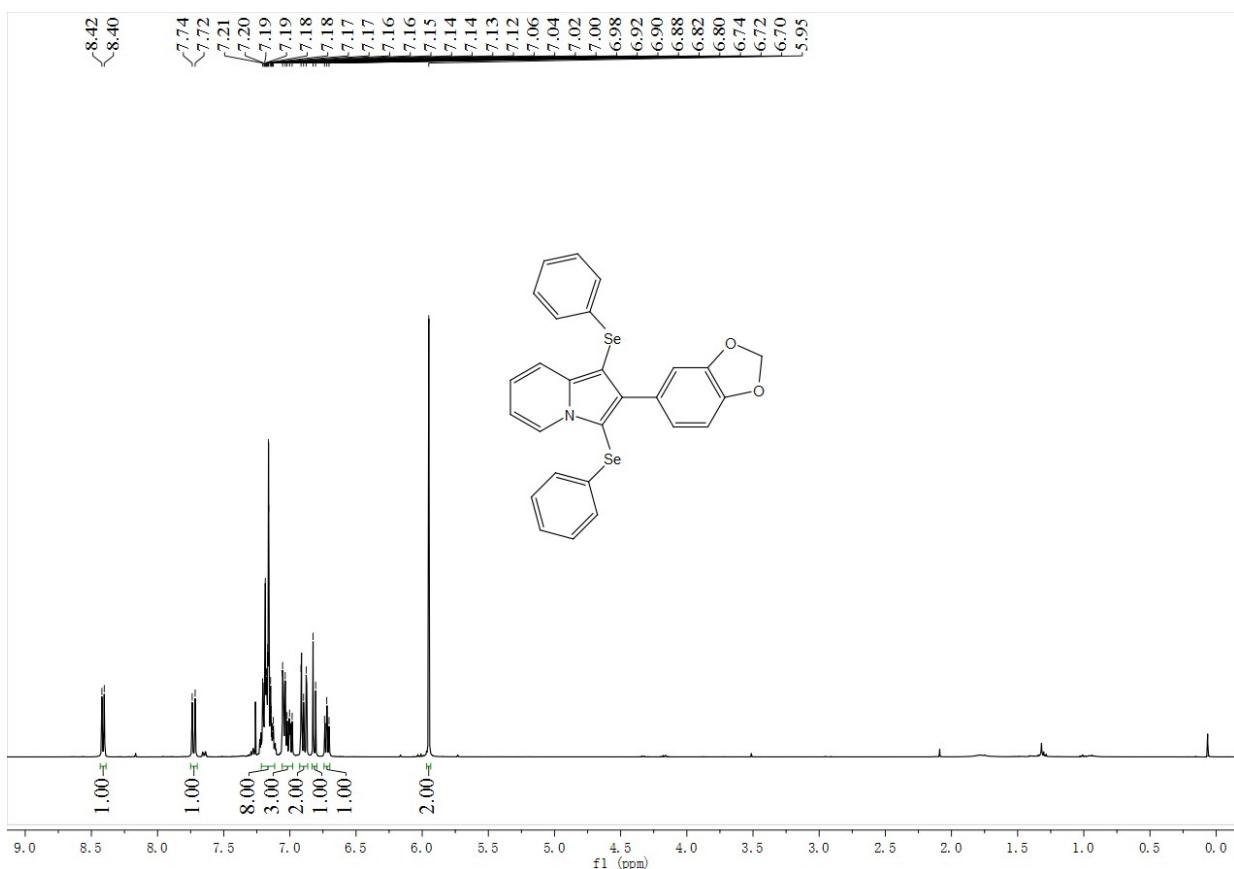
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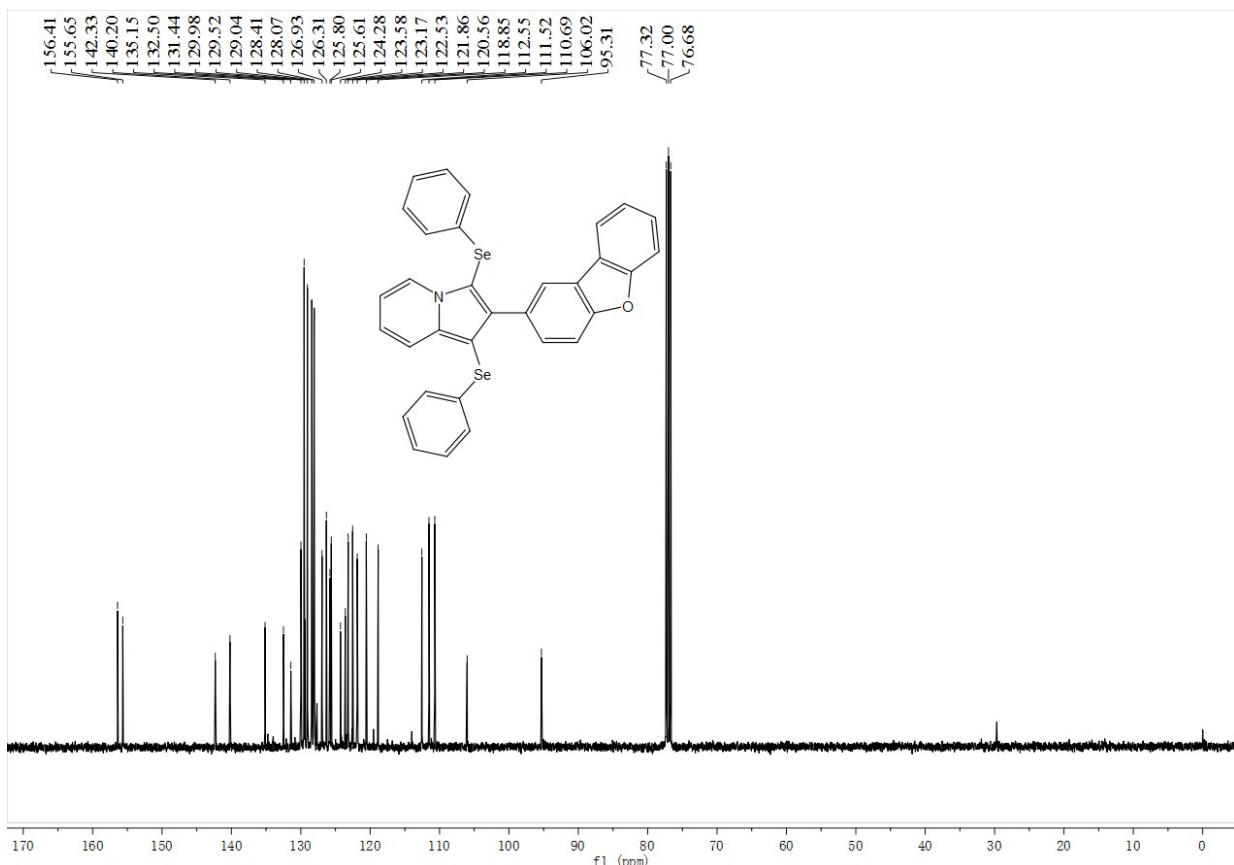
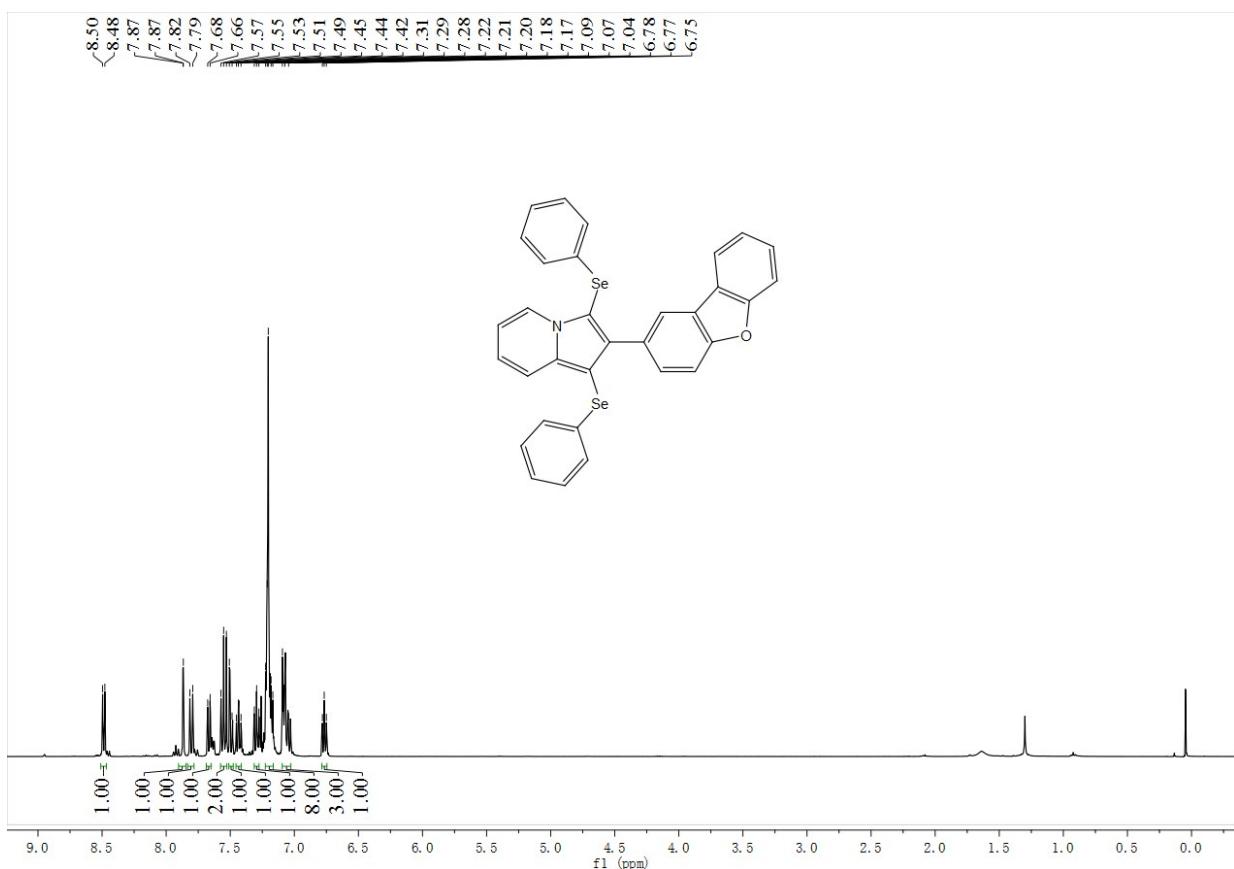
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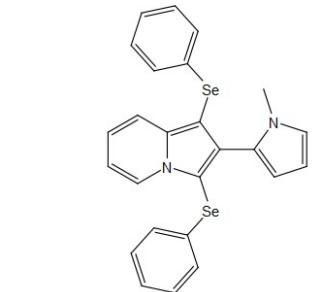
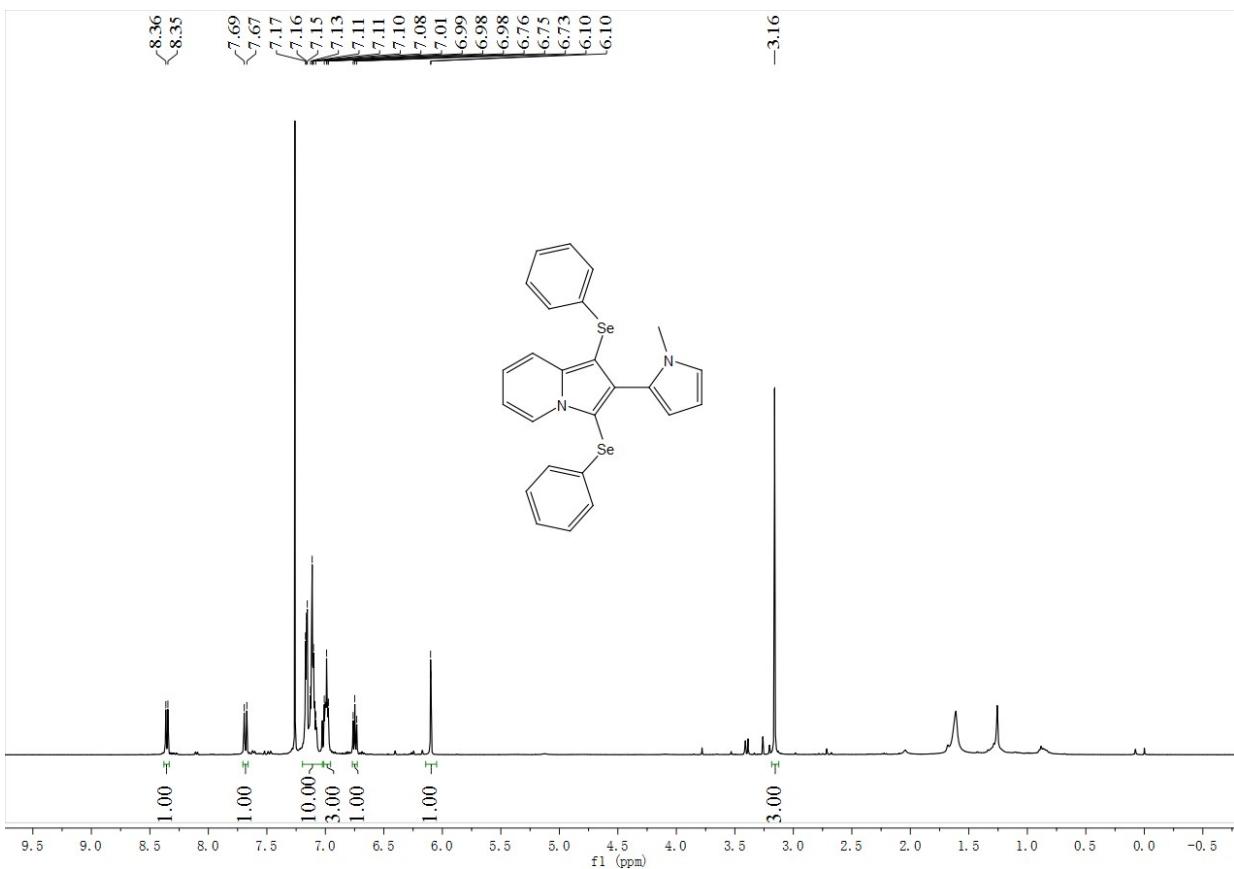
4va



4wa

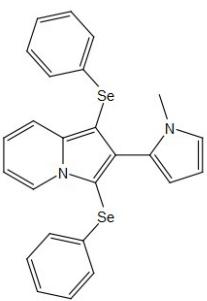
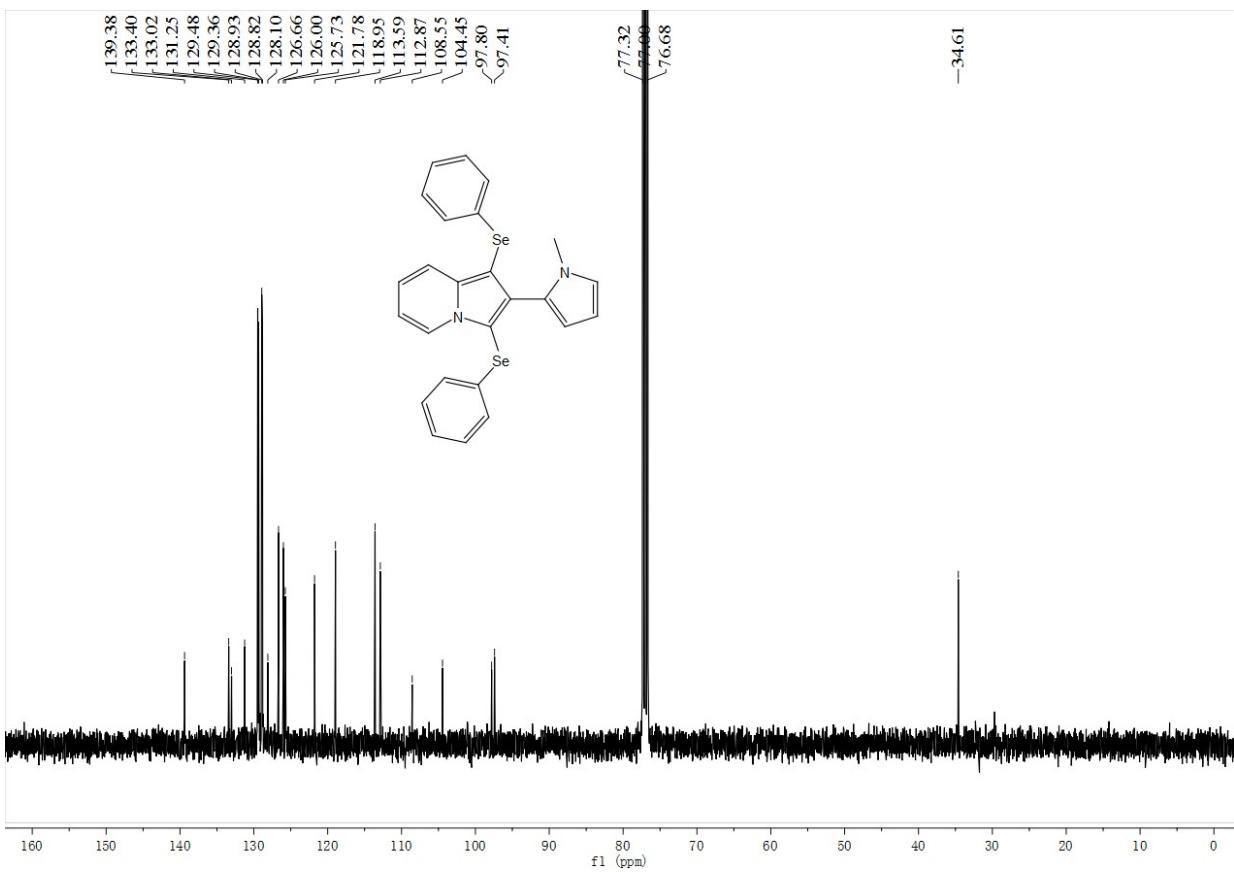


4xa

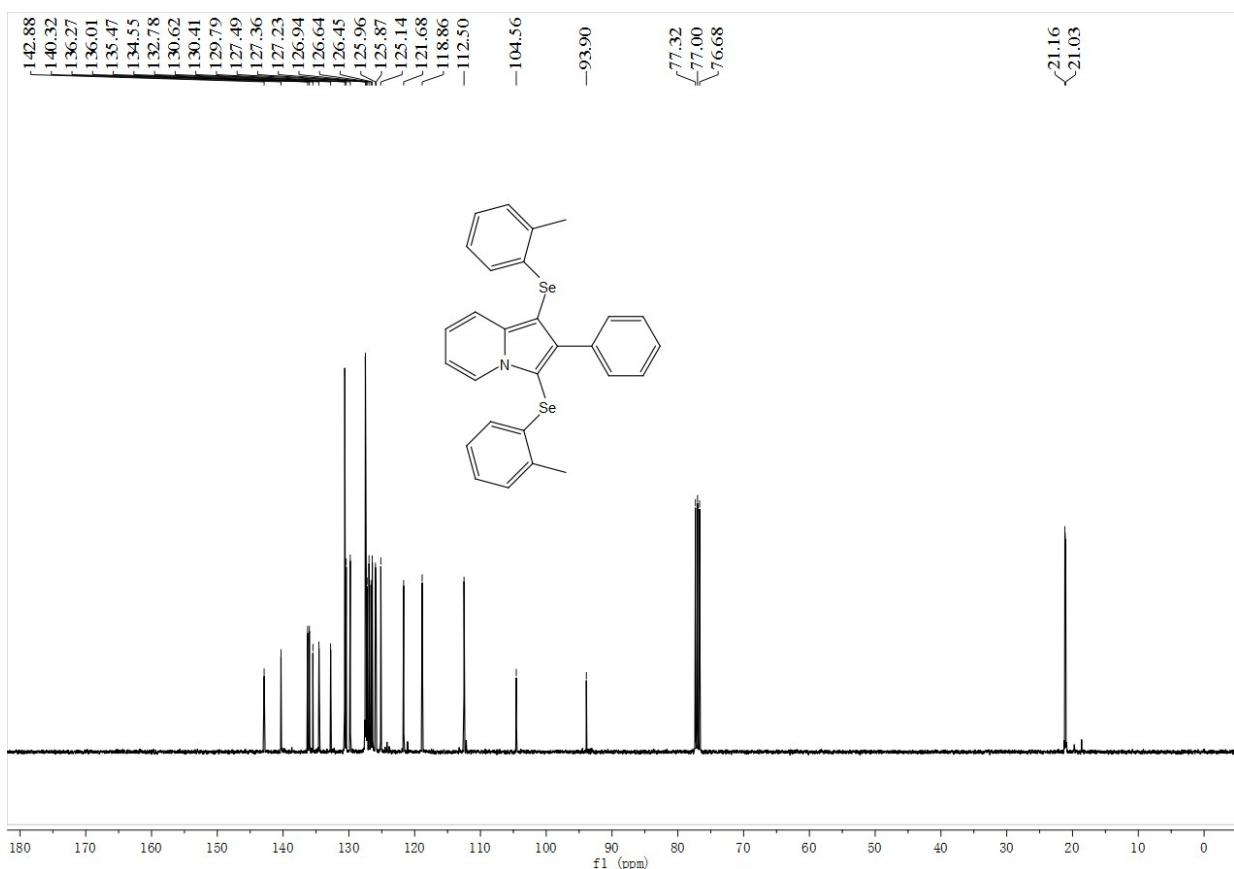
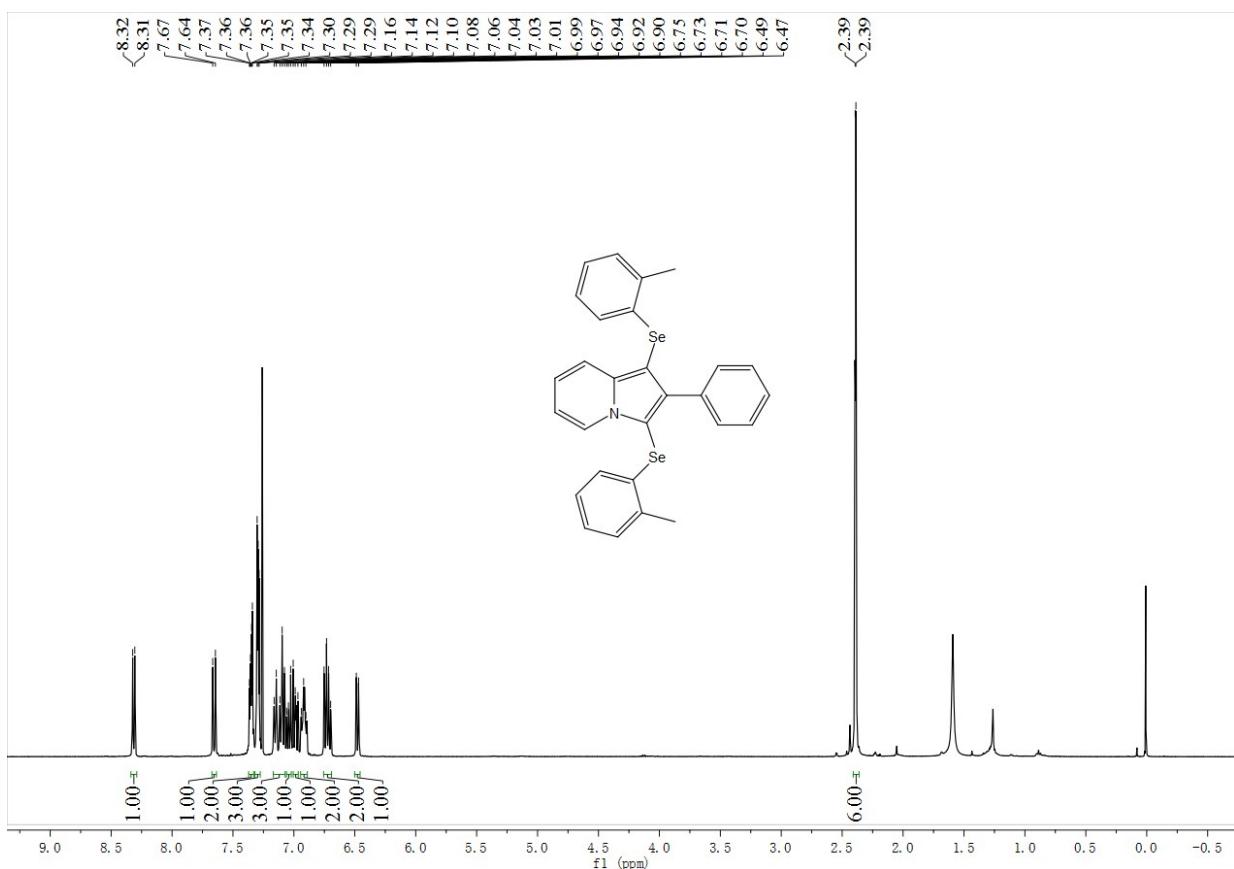


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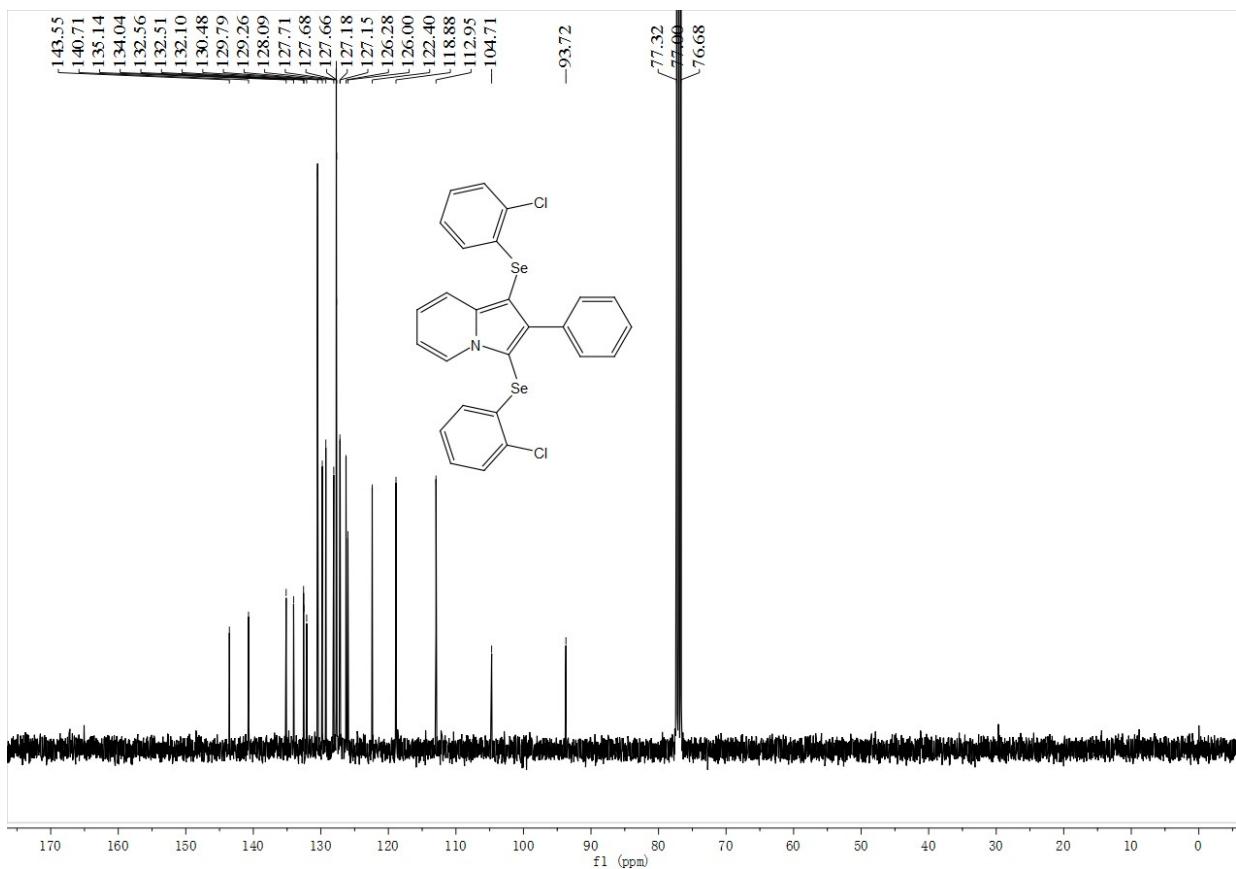
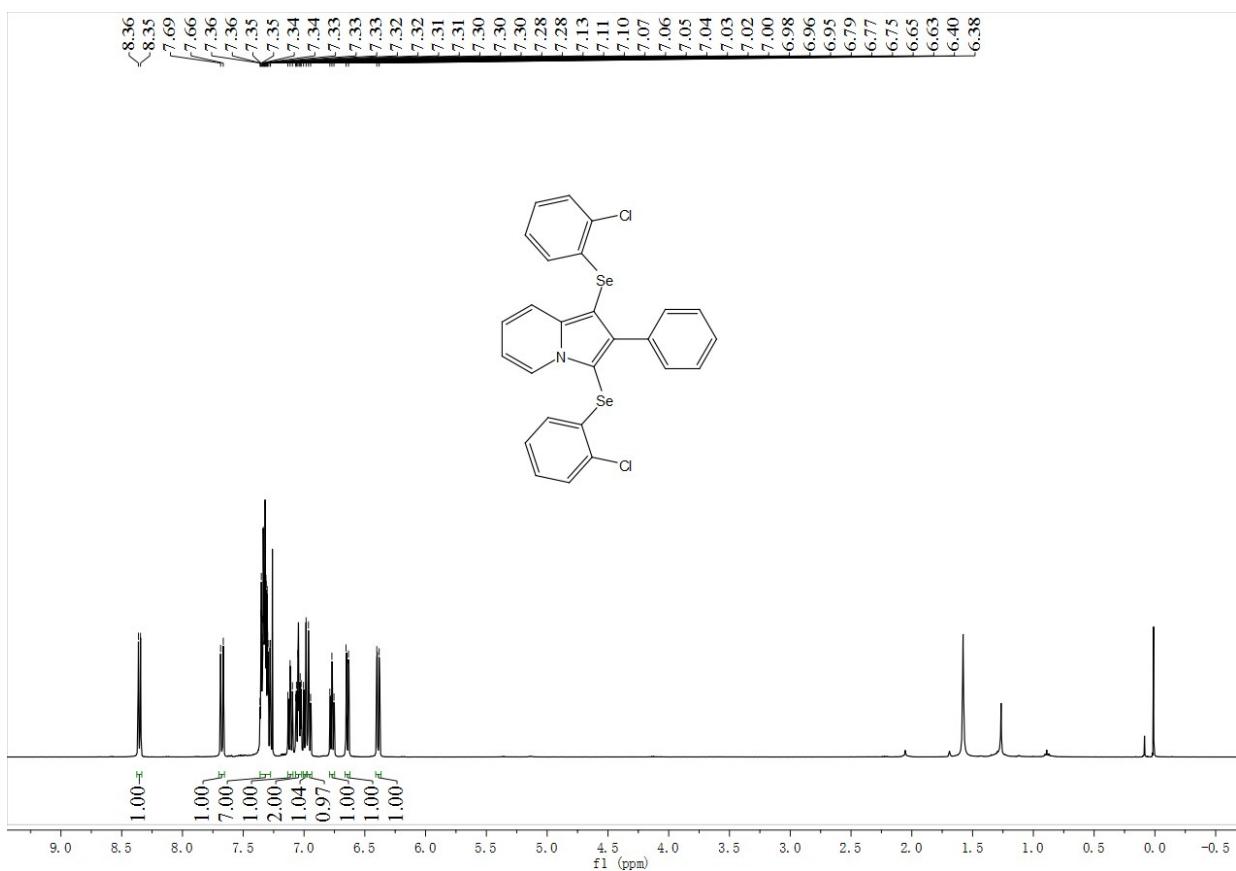
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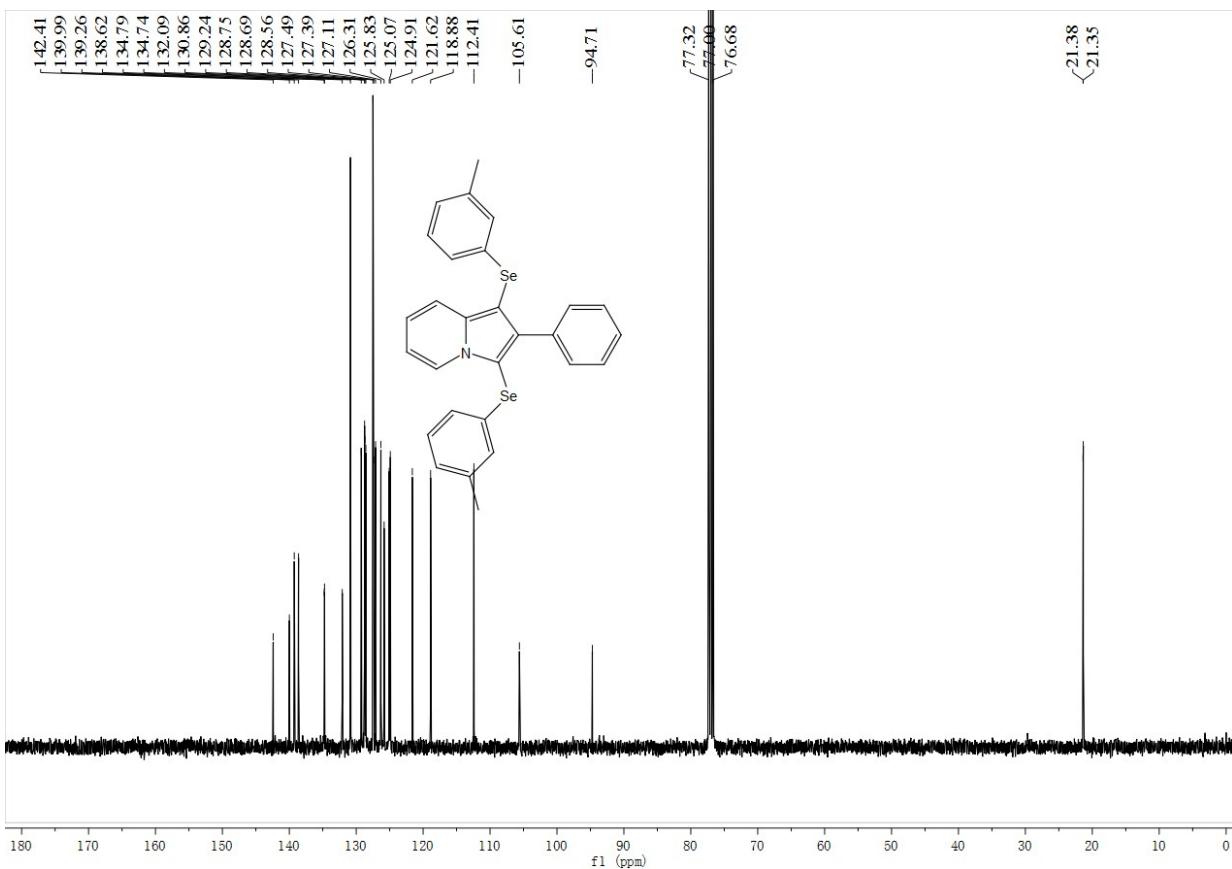
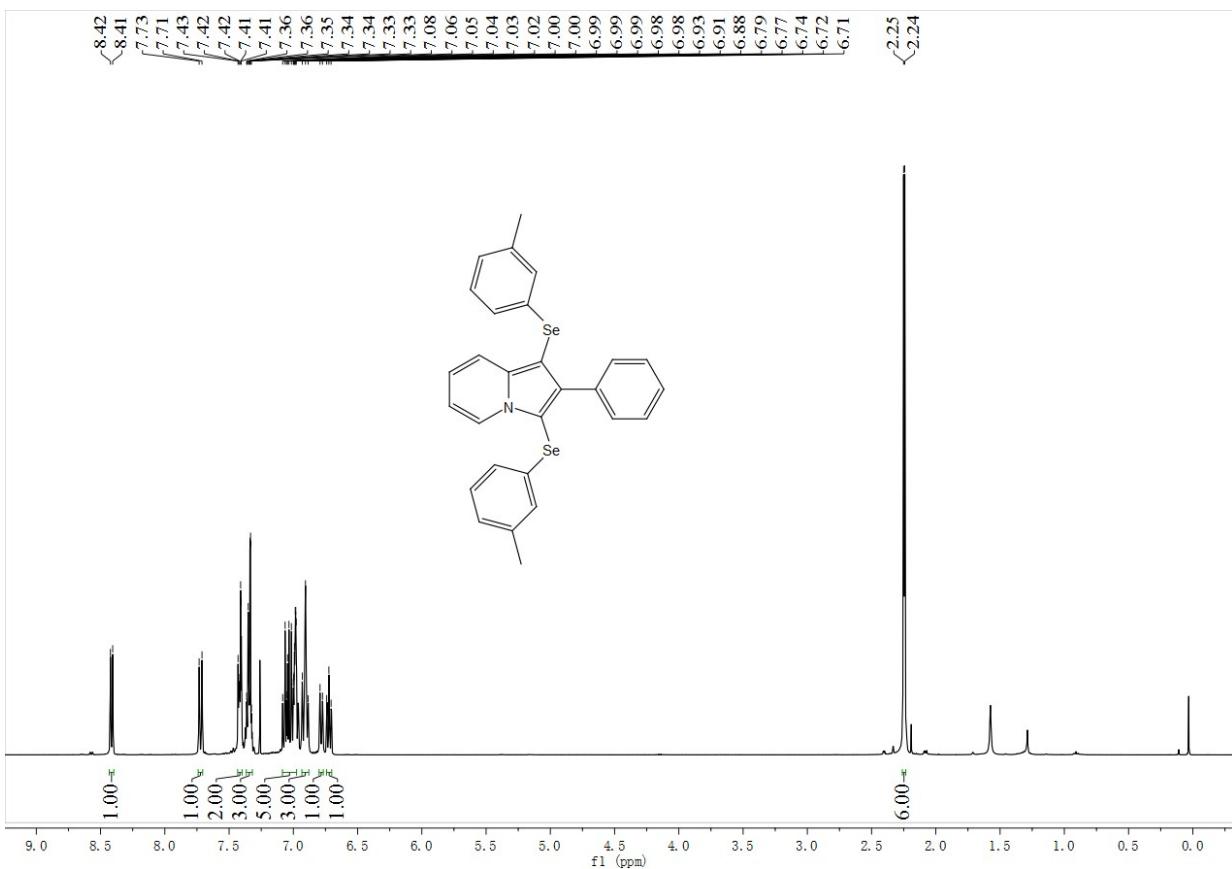
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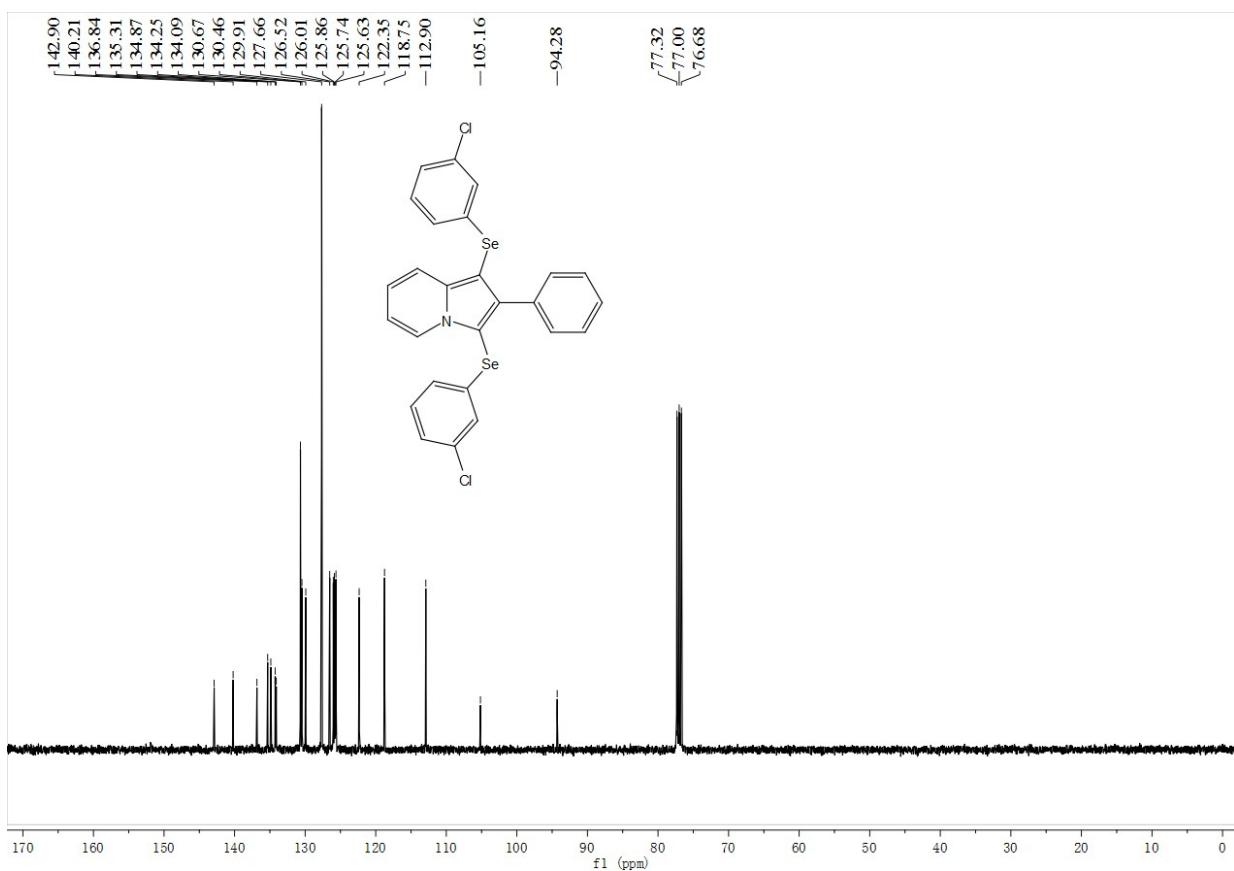
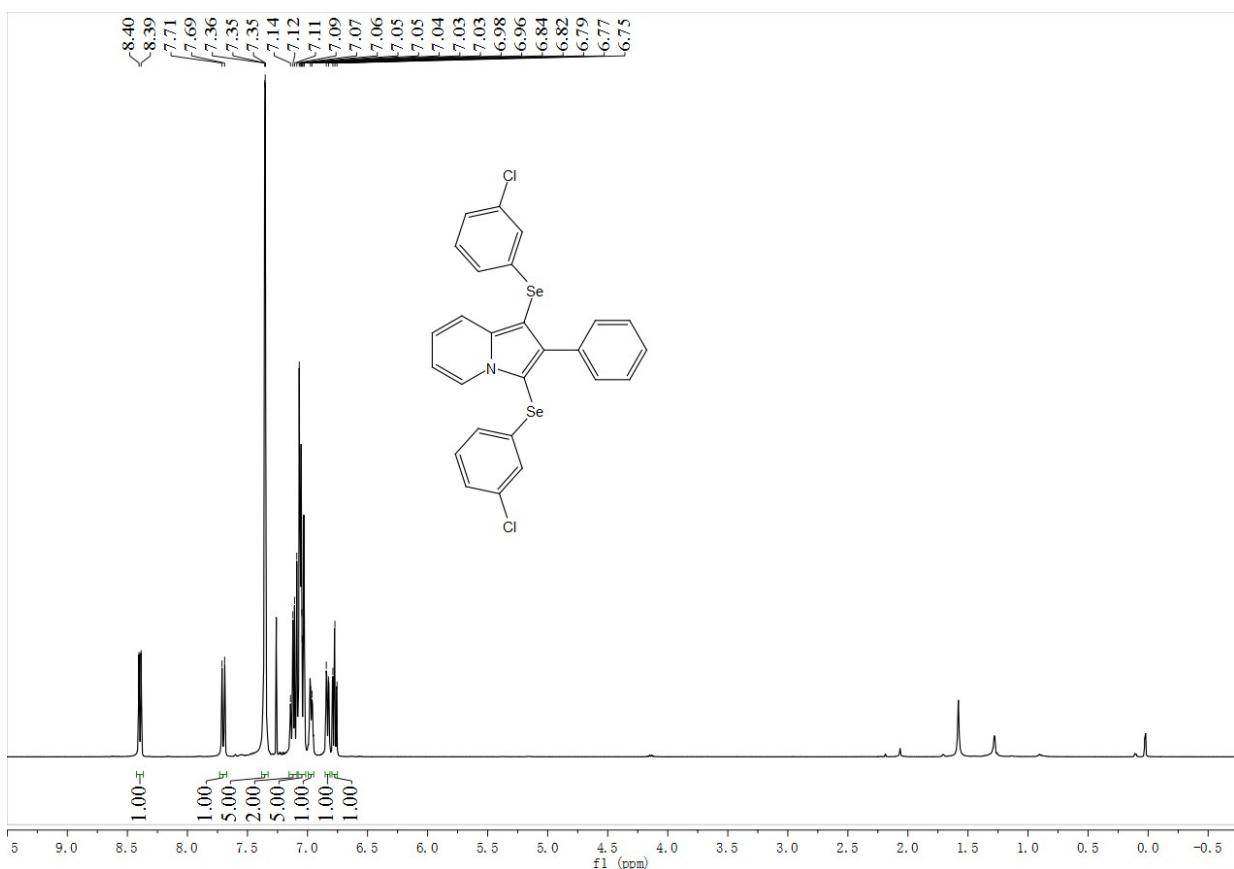
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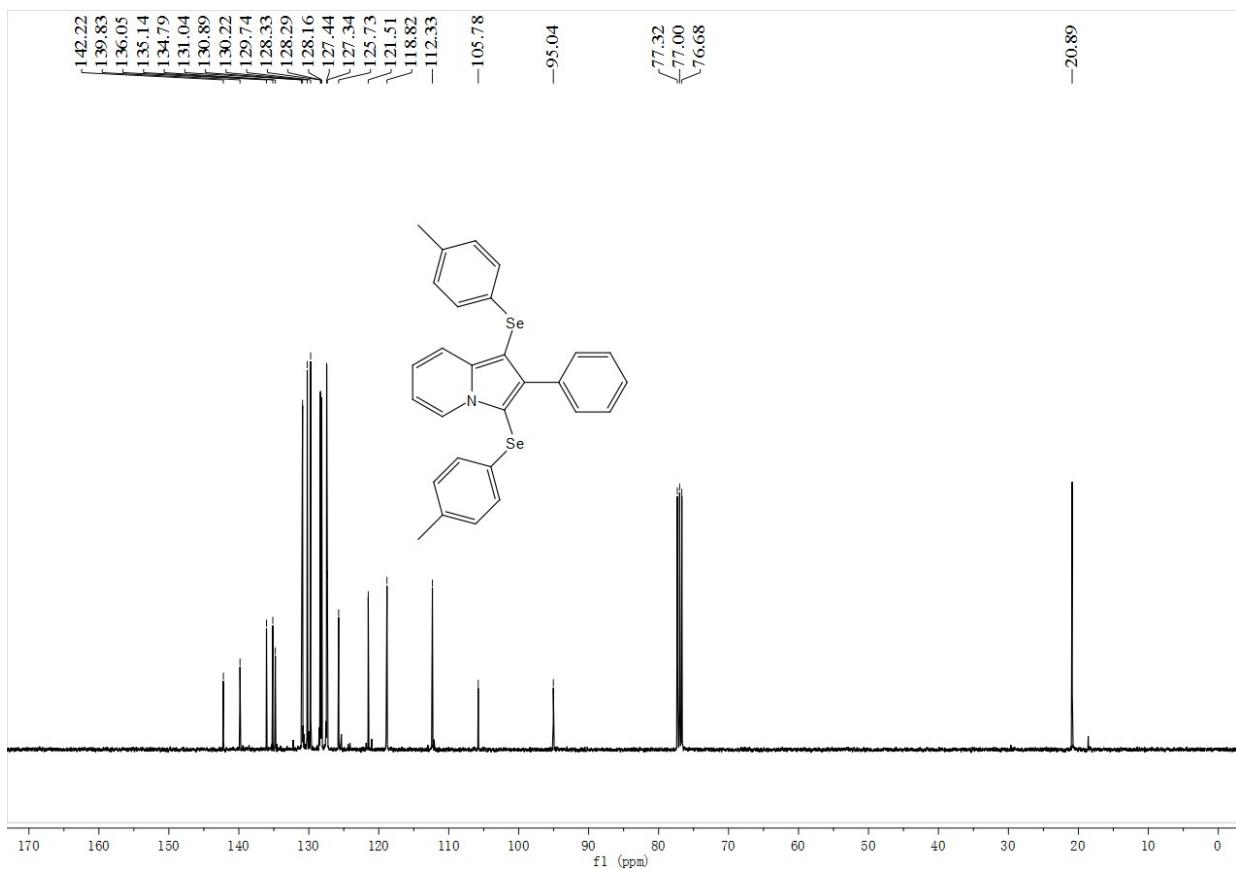
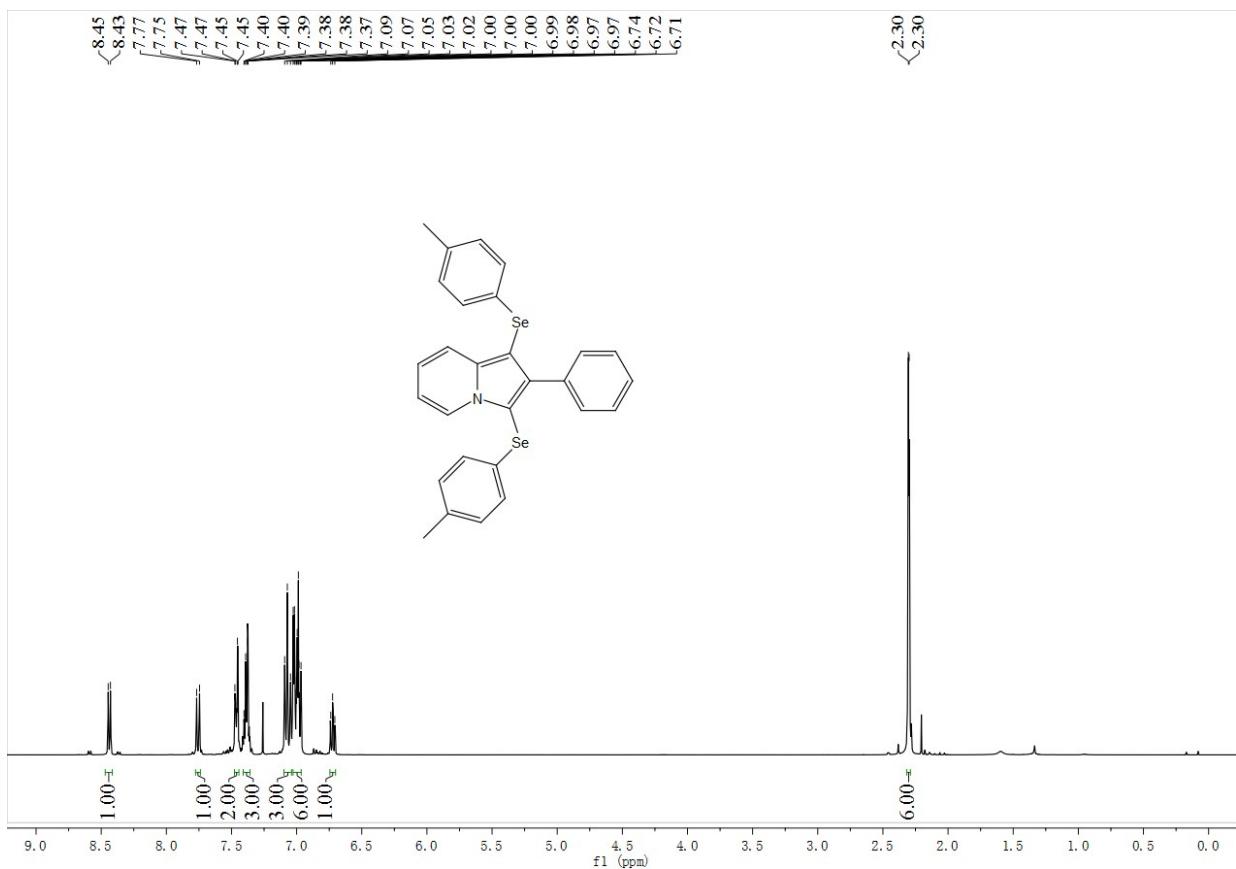
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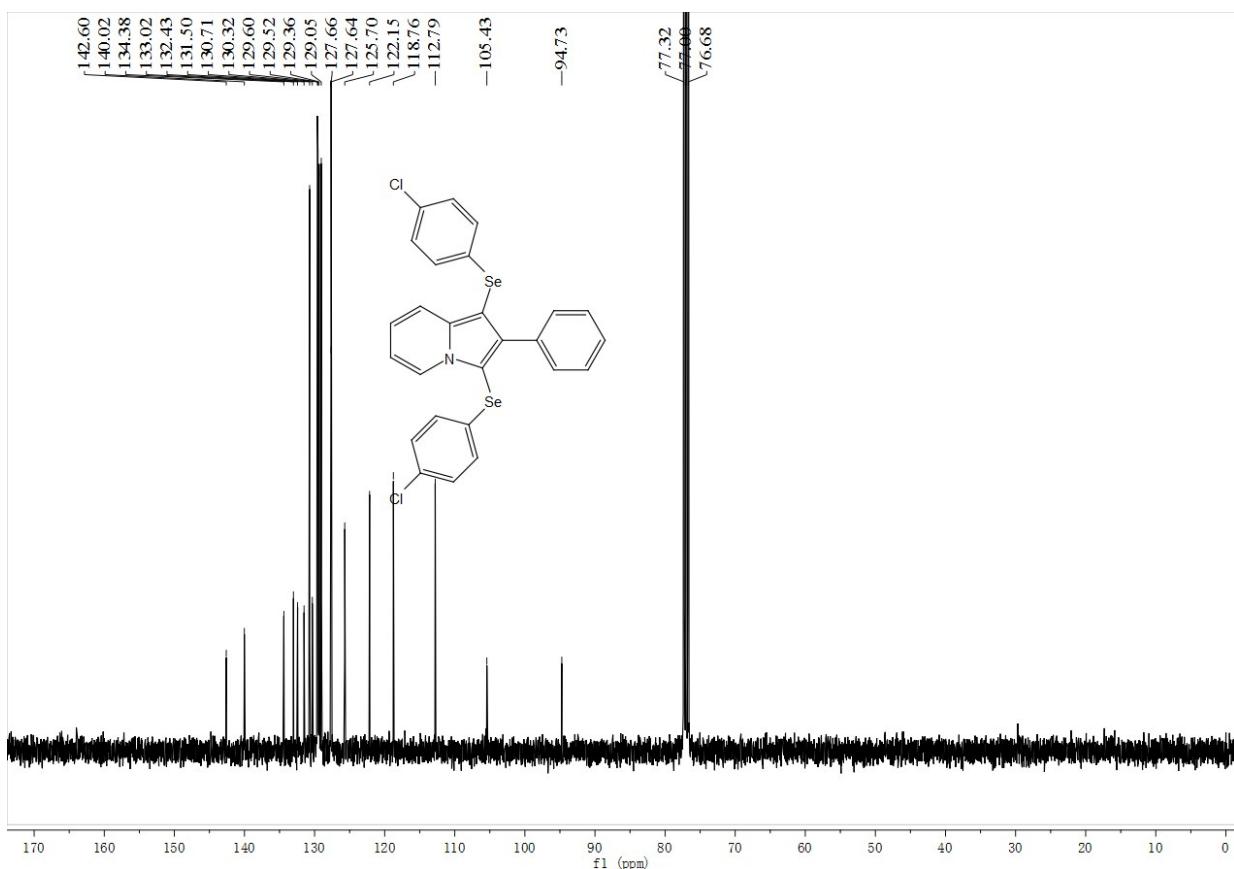
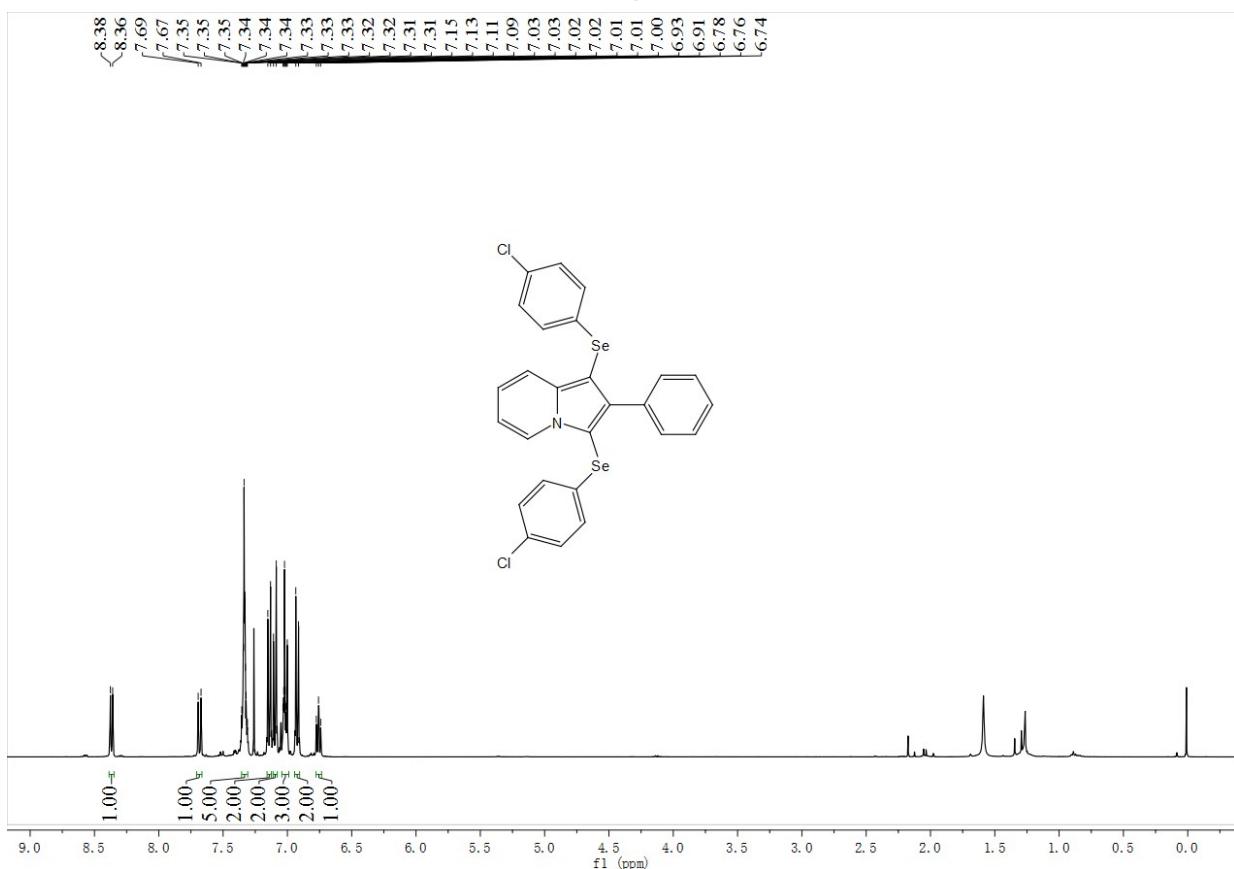
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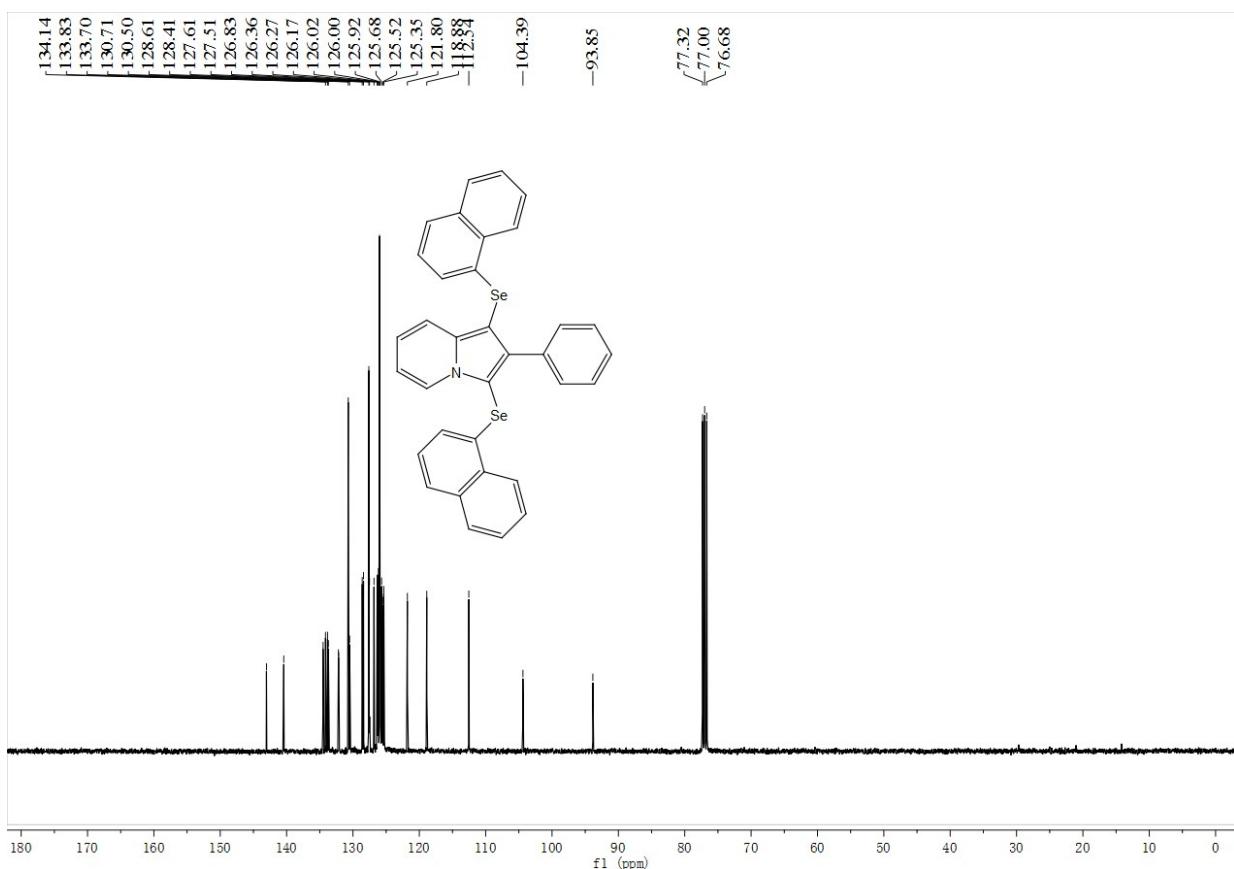
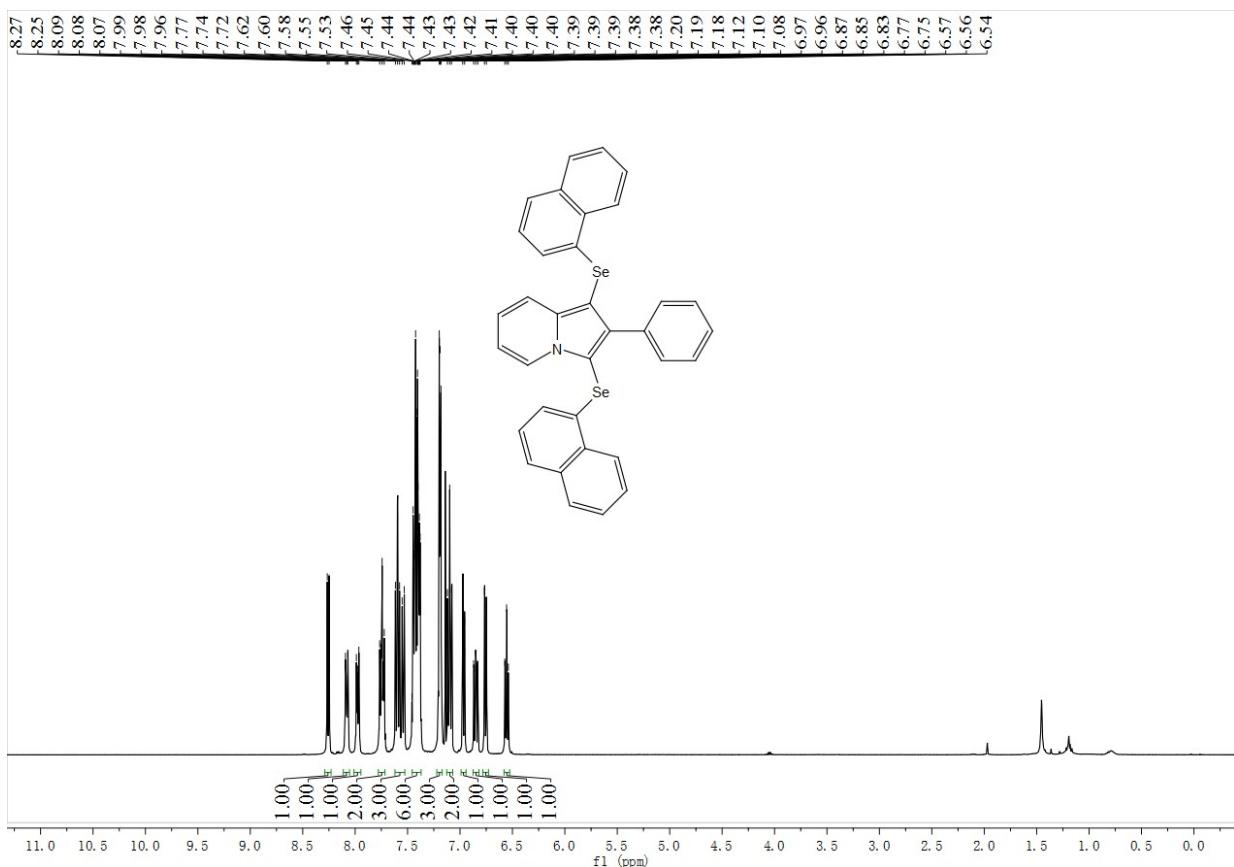
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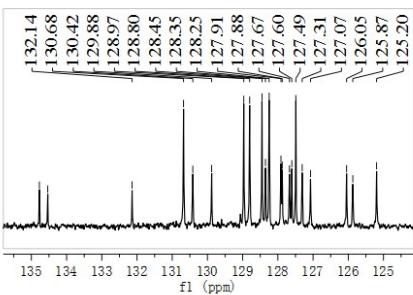
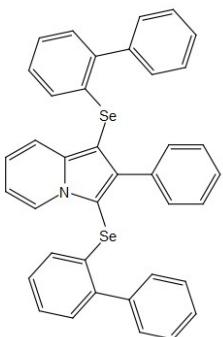
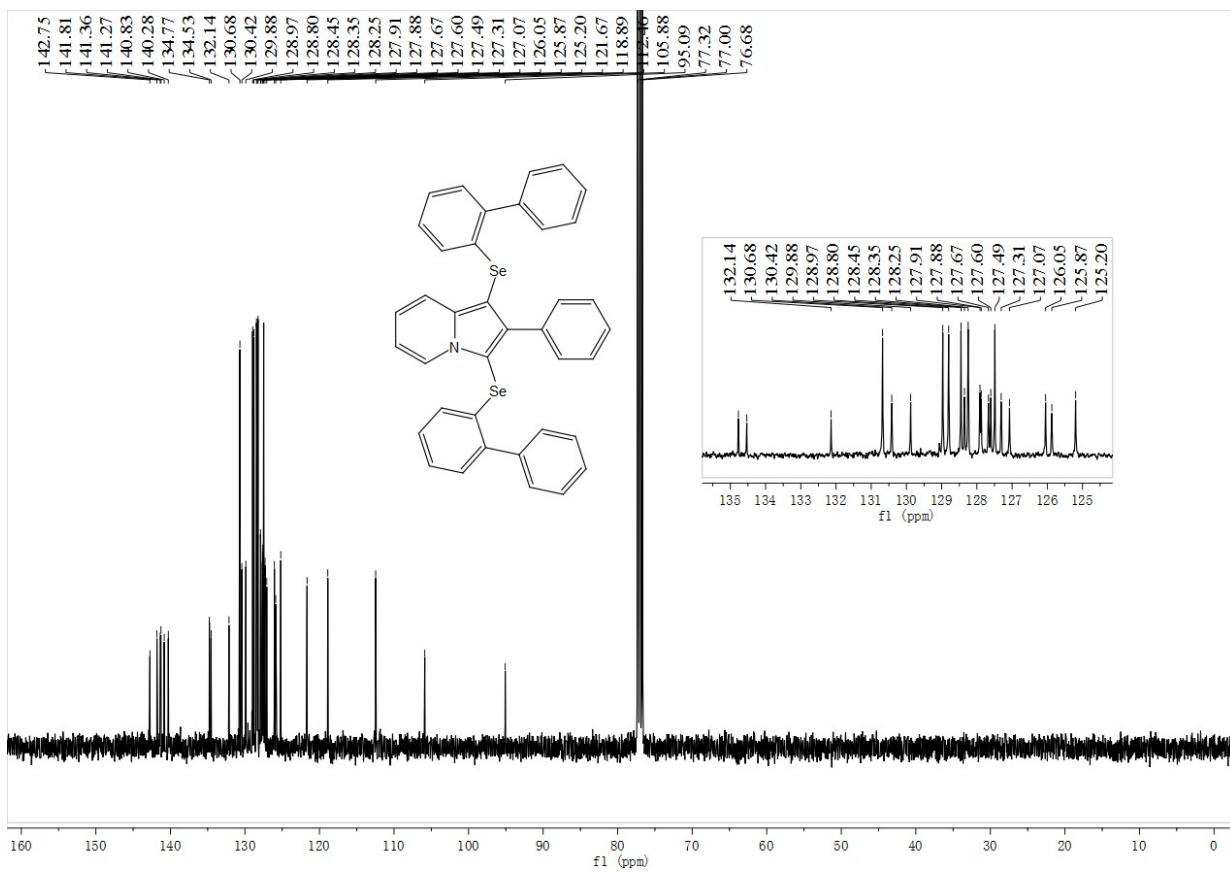
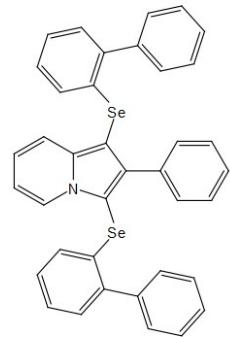
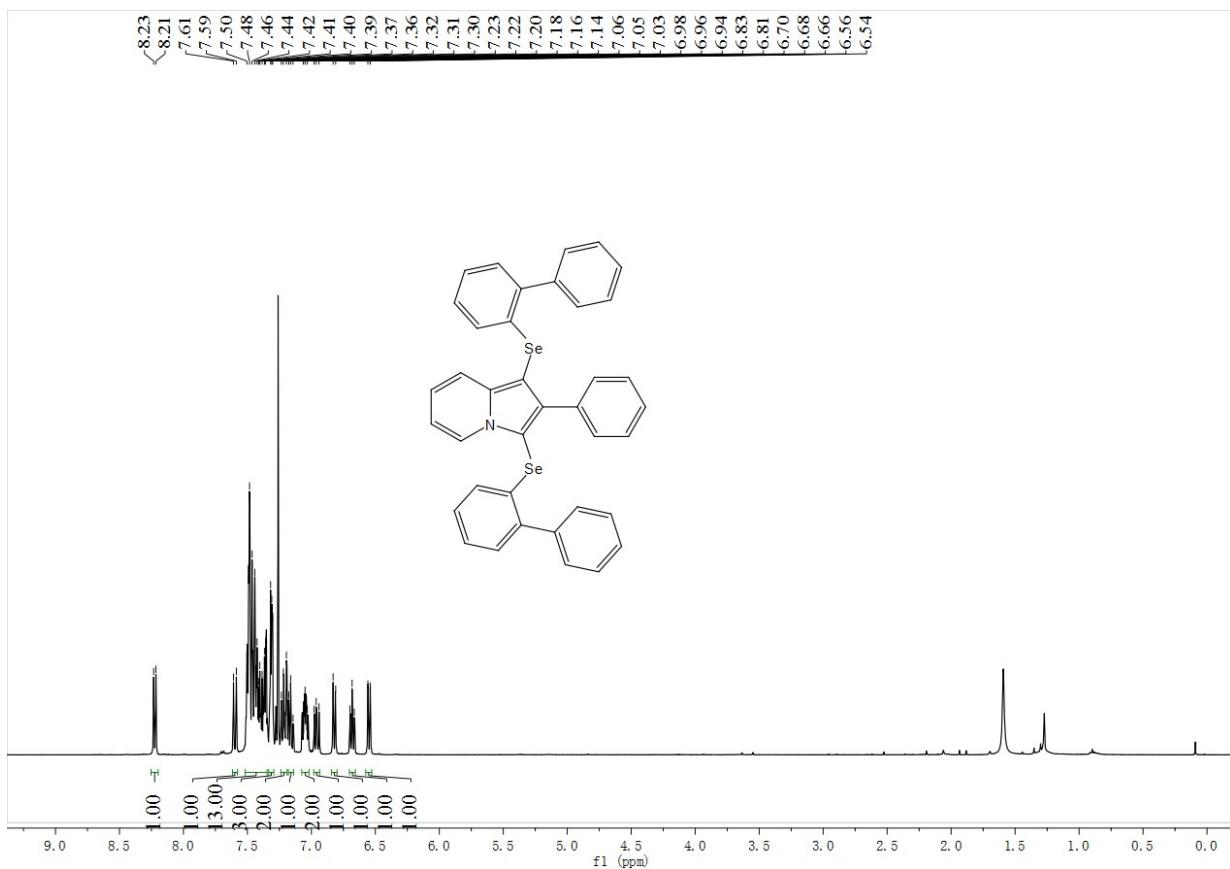
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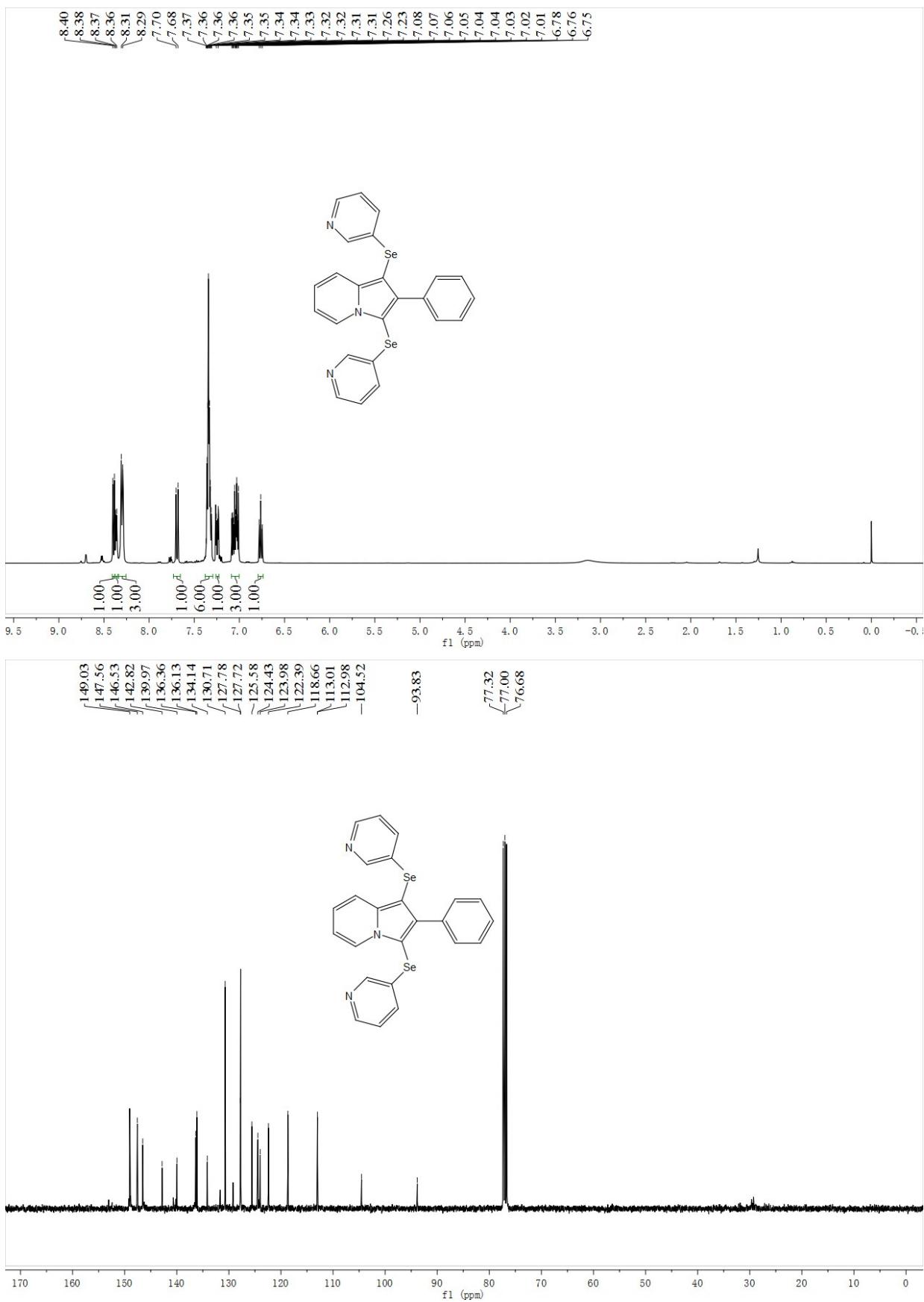
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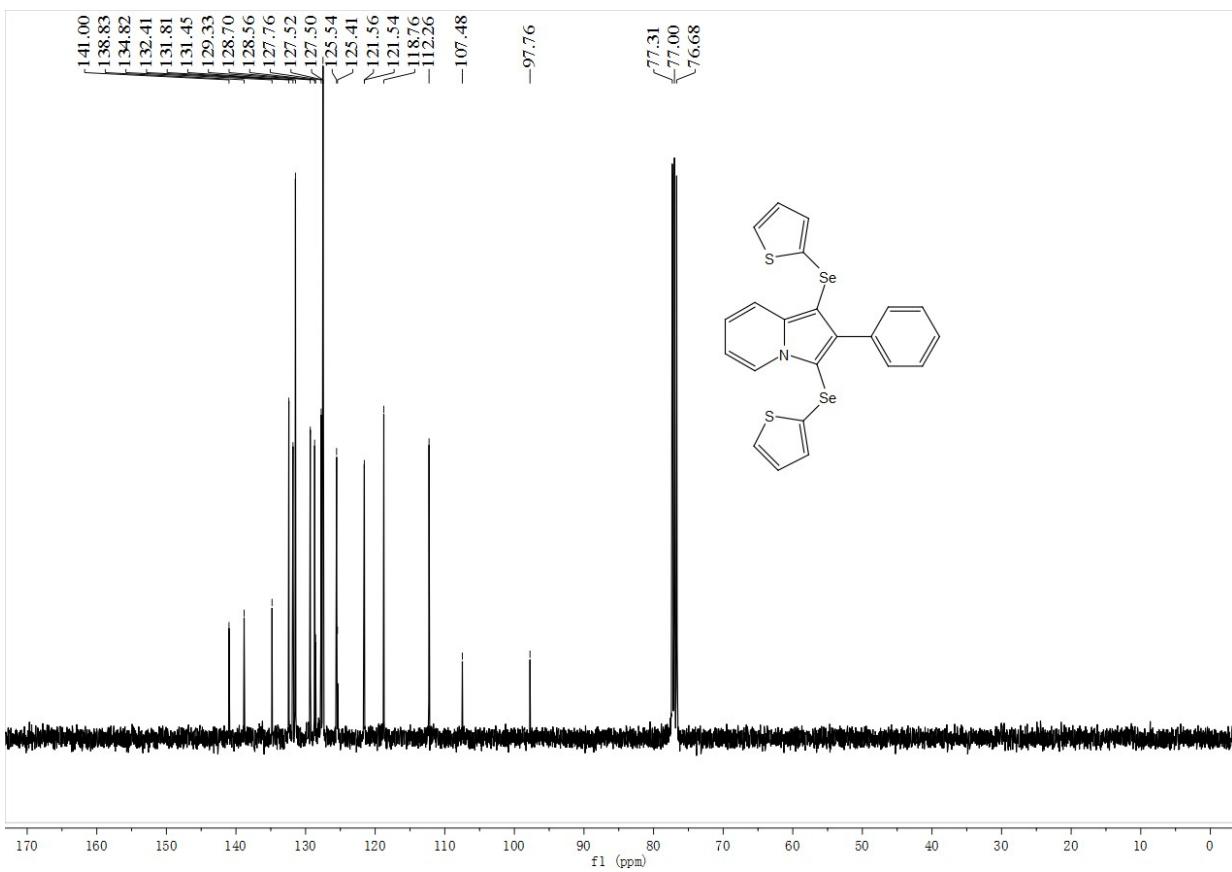
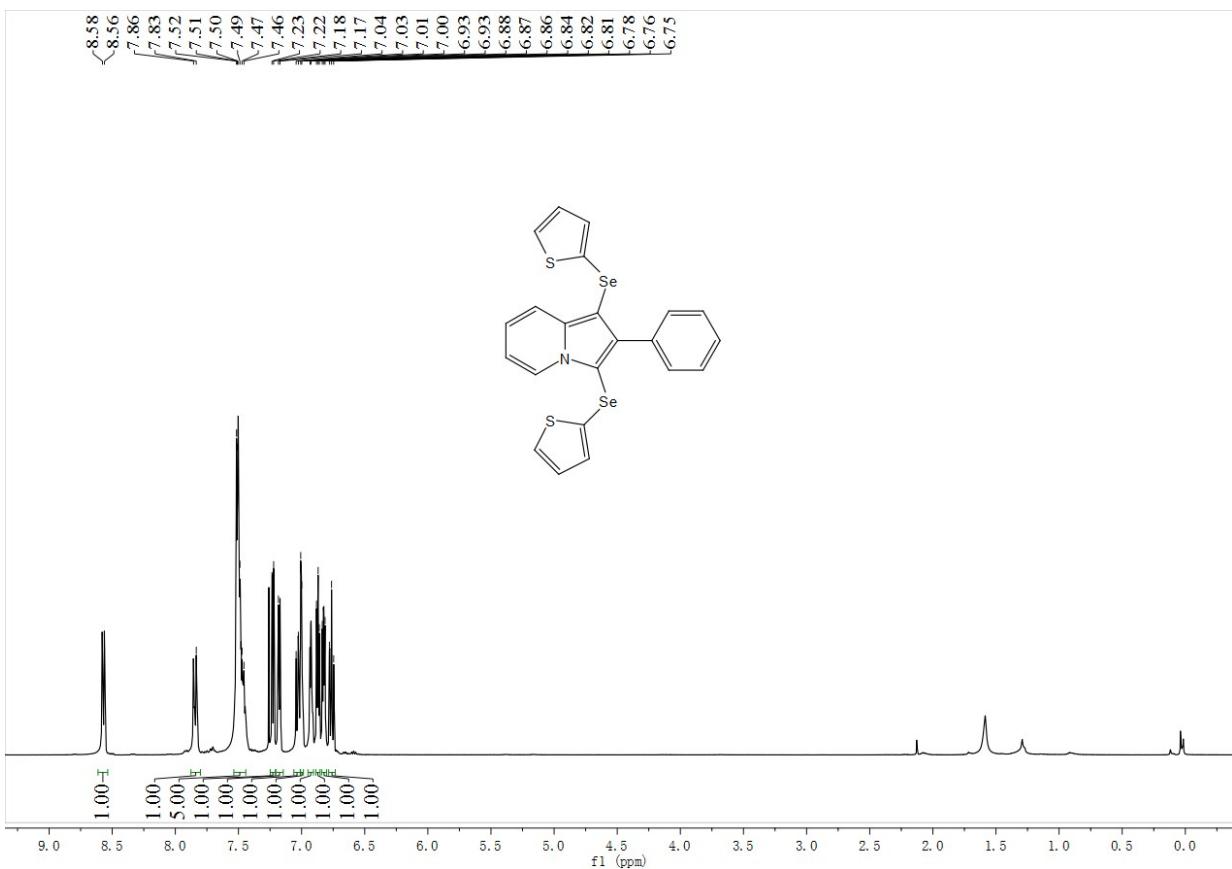
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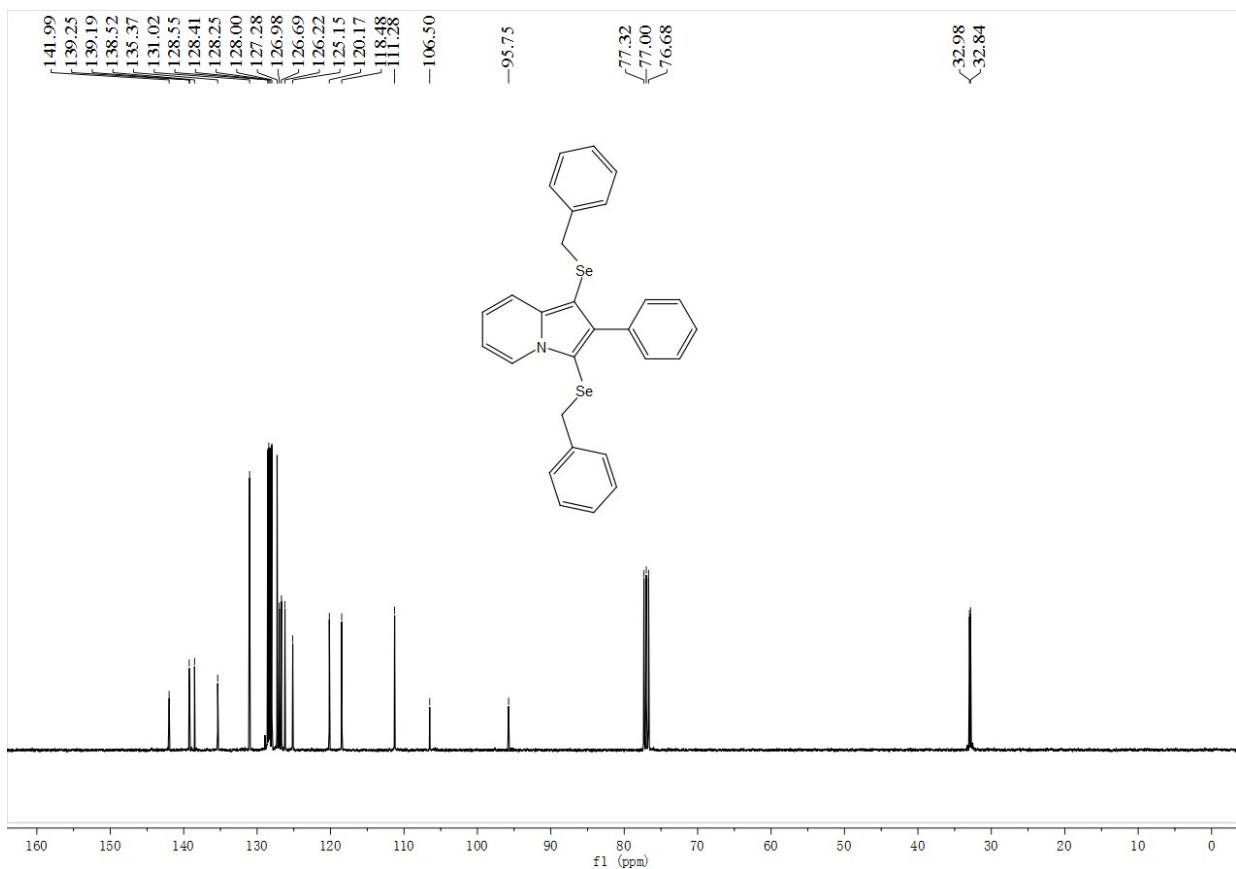
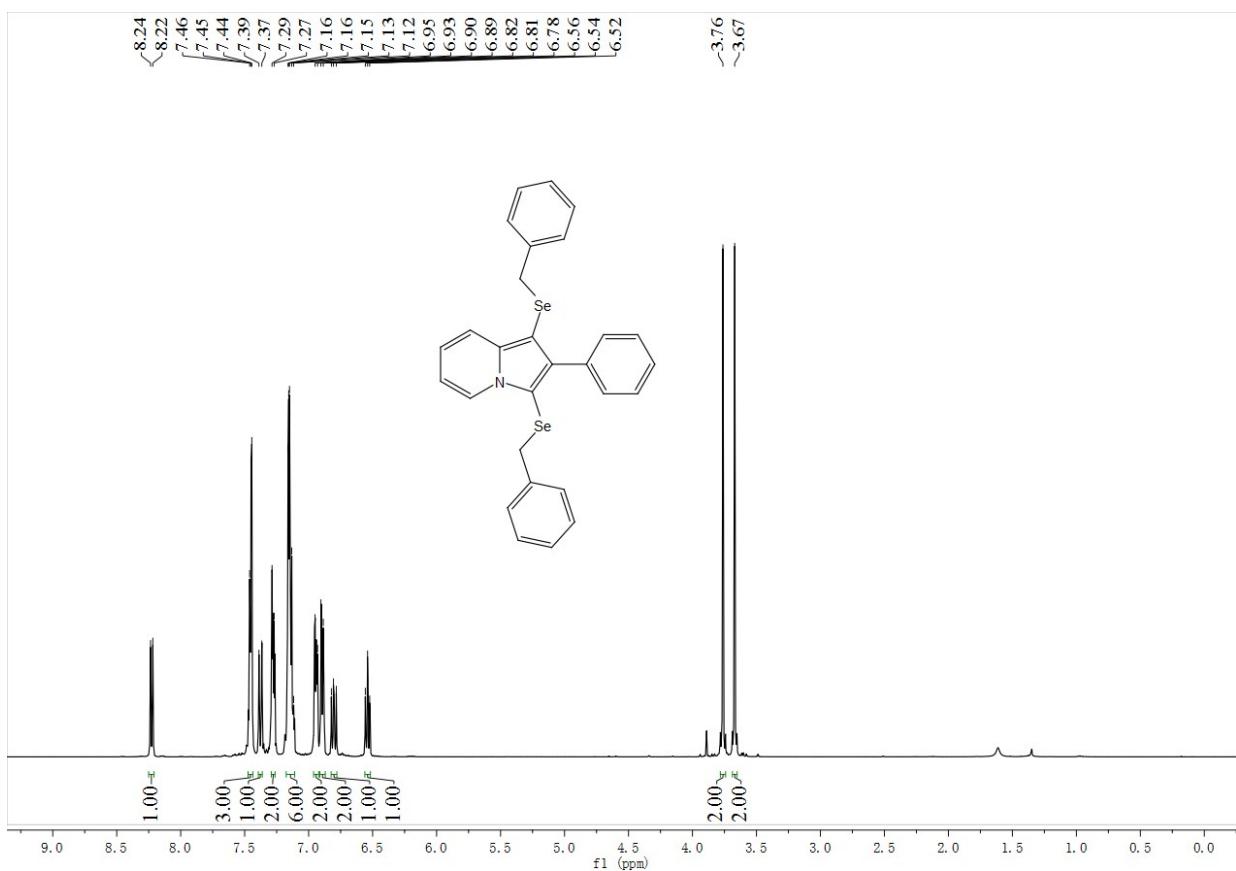
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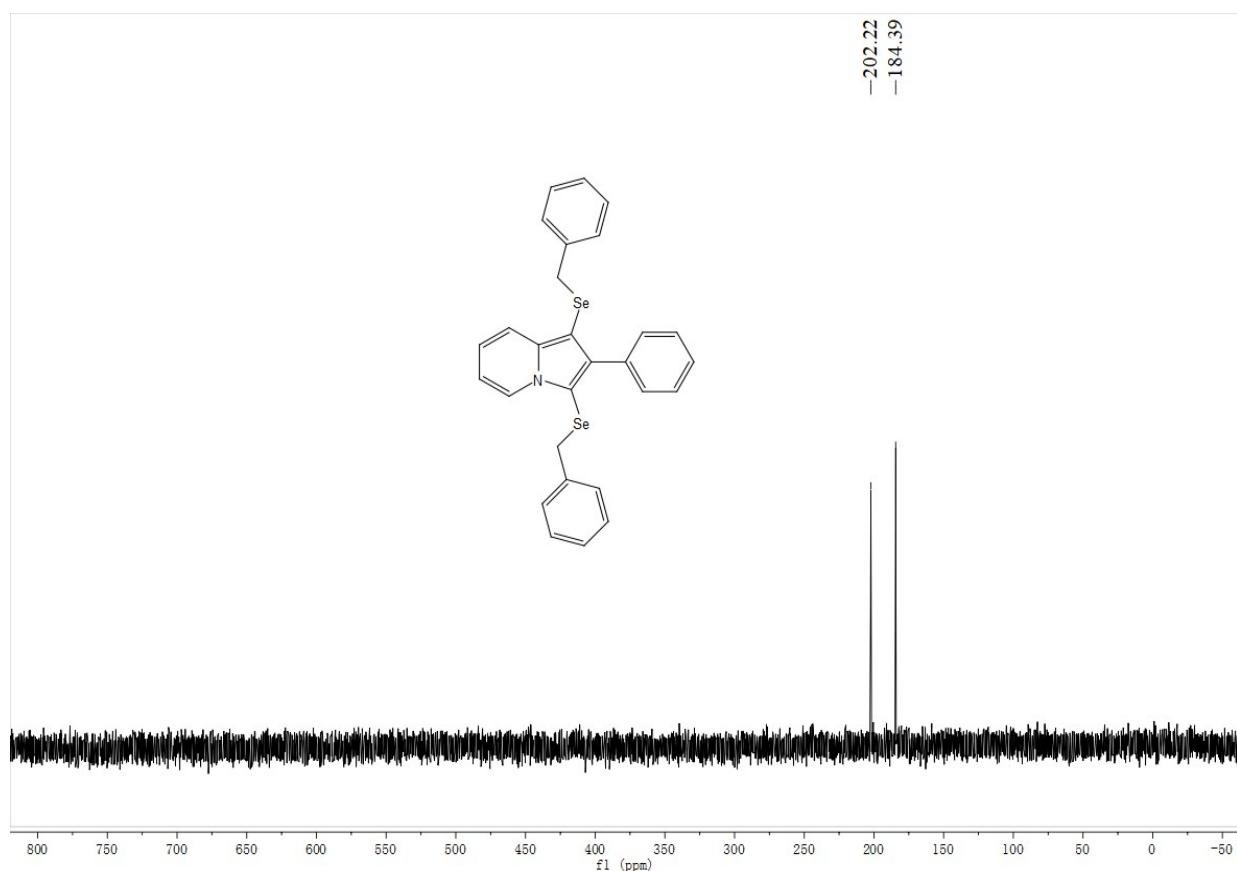


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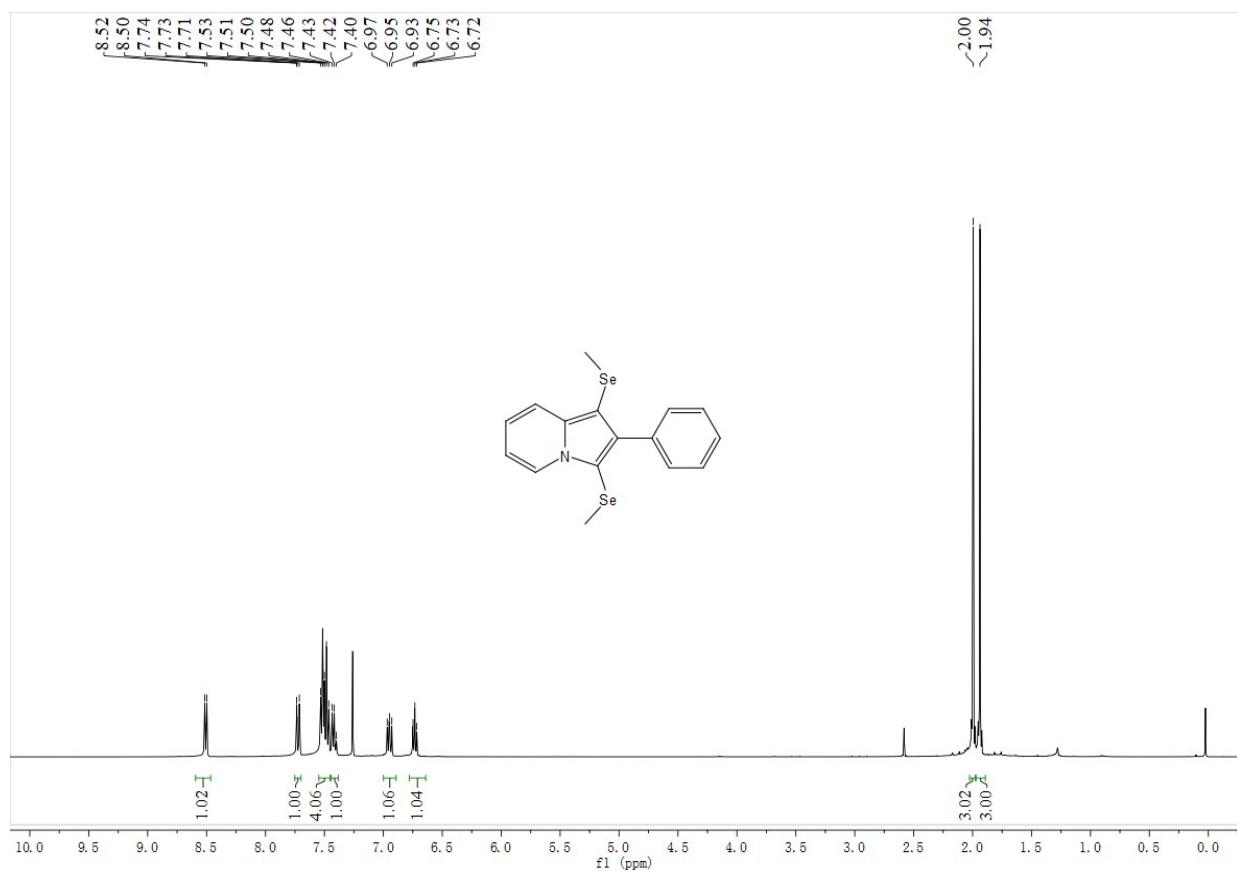


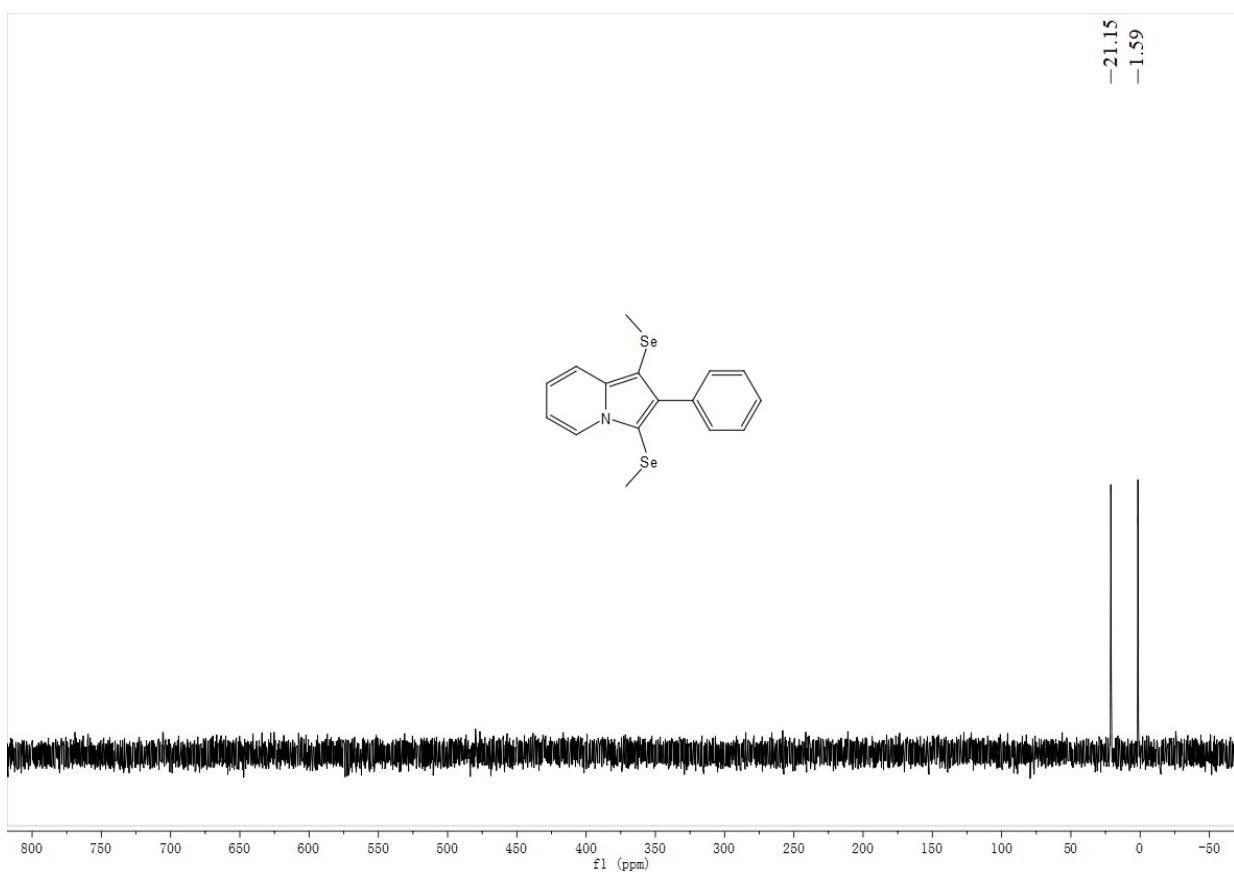
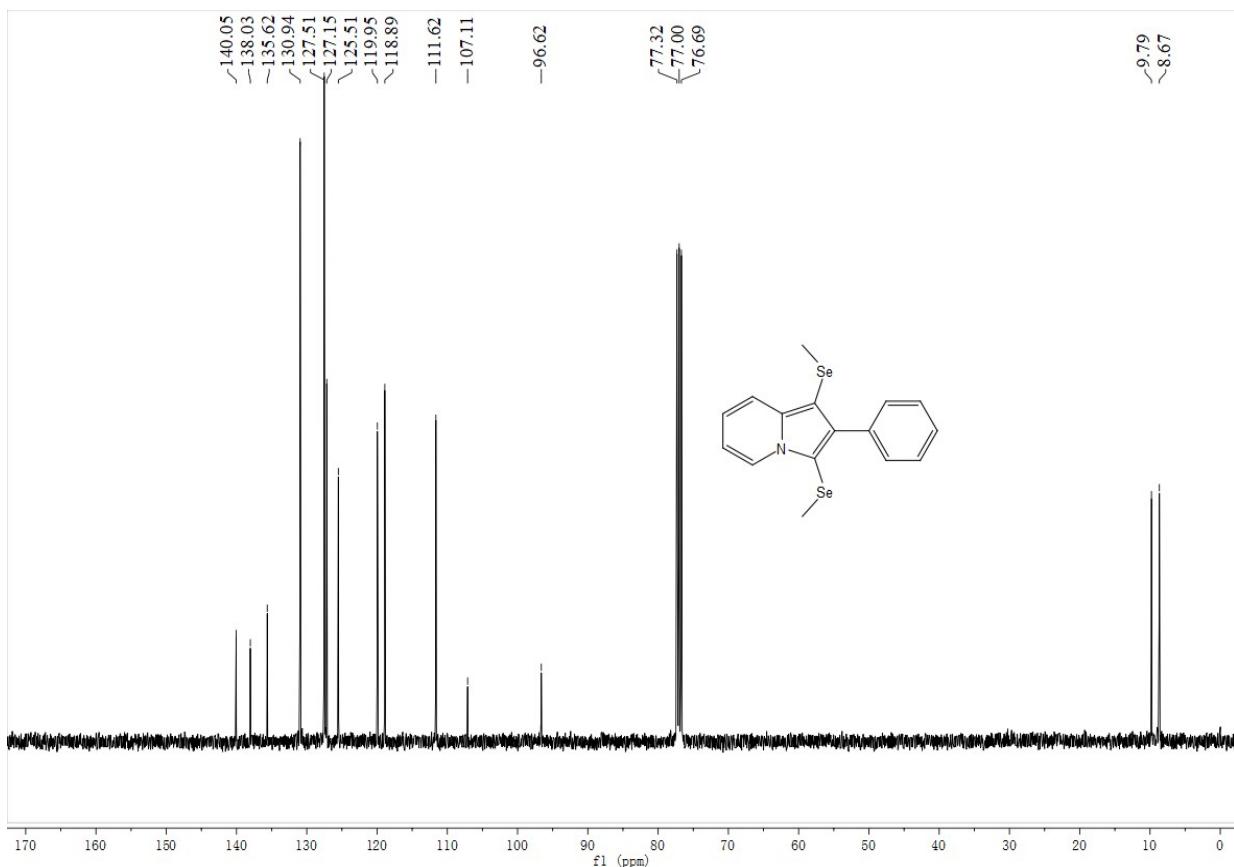
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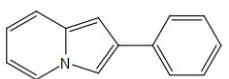
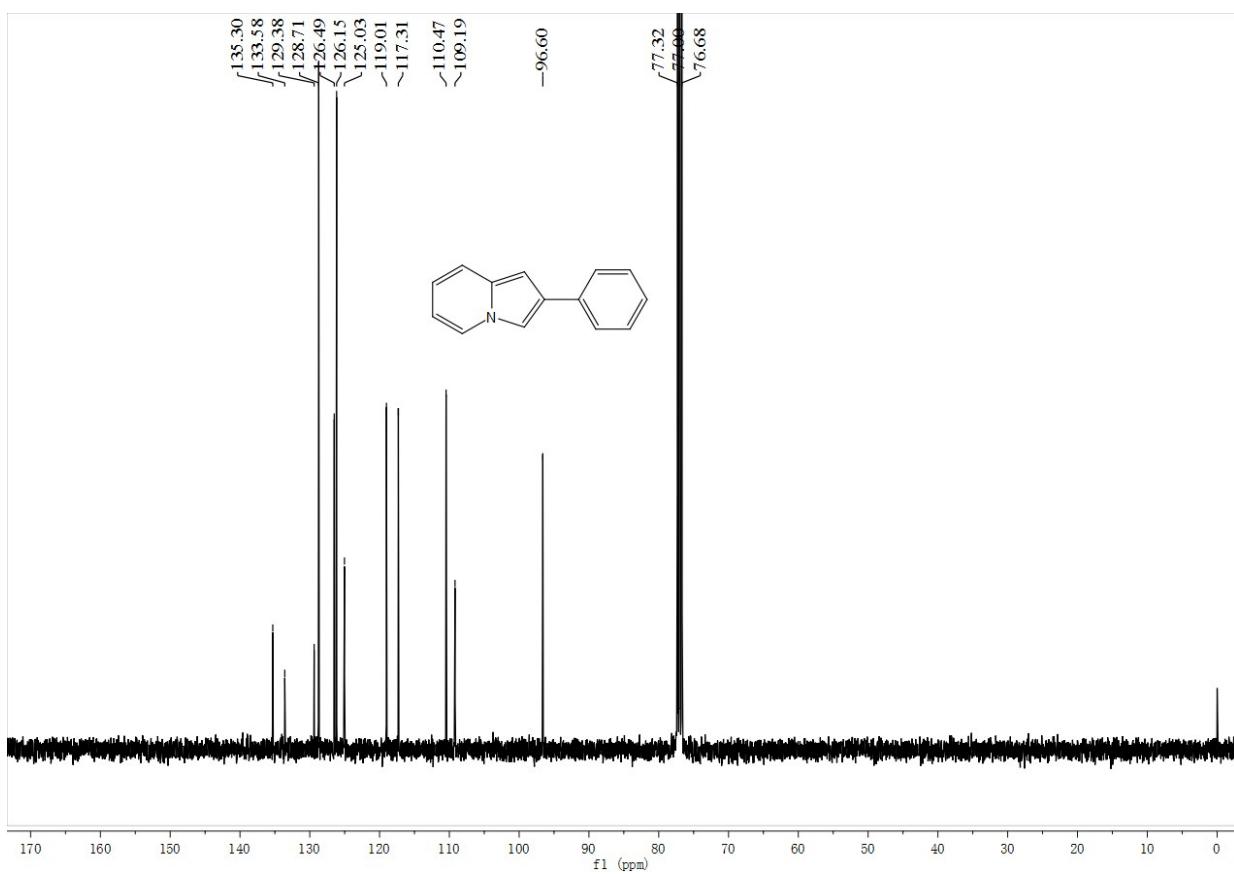
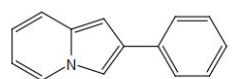
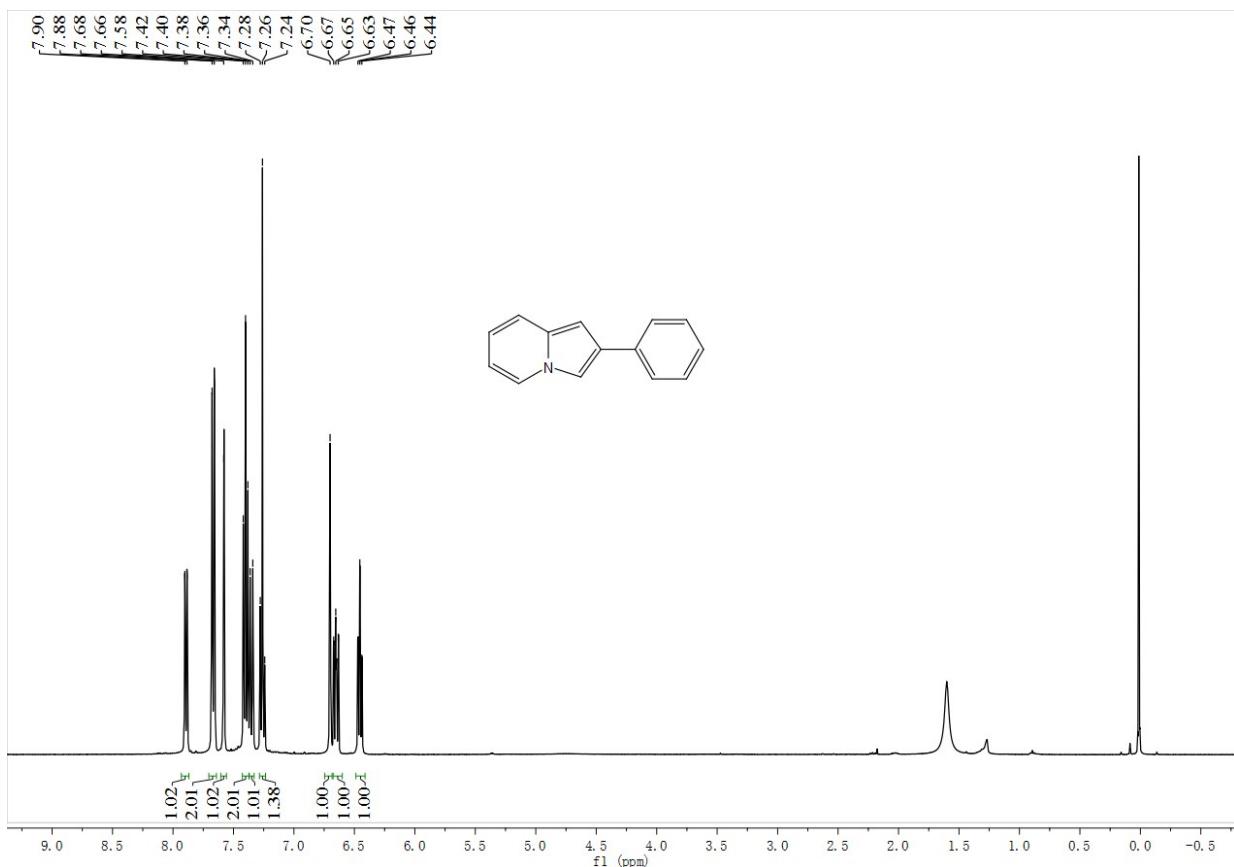


5am





6aa



7aa

