

Supporting information

Metal-Free, Iodine/DMSO-Promoted Oxidative C(sp²)-H Difunctionalization: Synthesis of Aminochalcogenylated indoles

Satyajit Pal^a, Subhankar Sarkar^a, Tanmay Pramanik^a, Anindita Mukherjee^b, Sougata Santra^b,
Grigory V. Zyryanov^{b,c} and Adinath Majee*^a

^aDepartment of Chemistry, Visva-Bharati (A Central University), Santiniketan-731235,
India; E-mail: adinath.majee@visva-bharati.ac.in

^bDepartment of Organic & Biomolecular Chemistry, Chemical Engineering Institute, Ural
Federal University, 19 Mira Str., Yekaterinburg, 620002, Russian Federation

^cI. Ya. Postovskiy Institute of Organic Synthesis, Ural Division of the Russian Academy of
Sciences 22 S. Kovalevskoy Street, Yekaterinburg 620219, Russian Federation

Contents

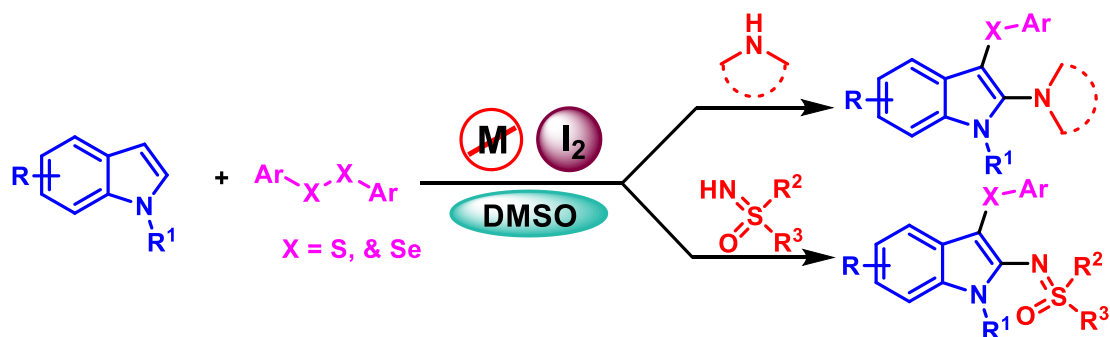
Sl. No.	Topics	Page No.
1	General information.	S2
2	General experimental procedure for the synthesis of 4aa-4aw .	S2
3	General procedure for the application	S3
4	Characterization data of the synthesized compounds	S3-S16
5	References.	S16
6	NMR spectra [¹ H, and ¹³ C{ ¹ H}] of synthesized products.	S18

1. Experimental Section

General Information

^1H NMR spectra were determined on a Bruker 400 (400 MHz) spectrometer as solutions in CDCl_3 . Chemical shifts are expressed in parts per million (δ) and the signals were reported as s (singlet), d (doublet), t (triplet), m (multiplet), and coupling constants J were given in Hz. $^{13}\text{C}\{^1\text{H}\}$ NMR spectra were recorded at 100 MHz in CDCl_3 solution. Chemical shifts are expressed in parts per million (δ) and are referenced to CDCl_3 ($\delta = 77.16$) as an internal standard. TLC was done on a silica gel-coated glass slide (Merck, Silica gel G for TLC). Silica gel (60-120 mesh, SRL, India) was used for column chromatography. Petroleum ether refers to the fraction boiling in the range of 60-80 °C unless otherwise mentioned. All solvents were dried and distilled before use. Commercially available substrates were freshly distilled before the reaction. Solvents, reagents, and chemicals were purchased from Aldrich, Merck, and Spectrochem Chemicals. All reactions involving moisture-sensitive reactants were executed using oven-dried glassware. All the *NH*-sulfoximines¹ were prepared according to the reported method. A borosilicate glass tube was used as a reaction tube.

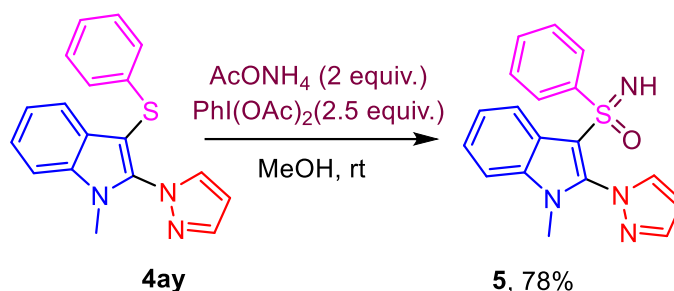
2. Scheme 1. General experimental procedure for the synthesis of 4aa-4ay.



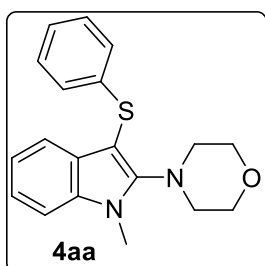
A mixture of indole (**1**, 1 mmol), amines (**2**, 1 mmol) and diaryl disulfide (**3**, 0.75 mmol) was stirred in the presence of 2 mL of DMSO as solvent and I_2 catalyst (20 mol%) at 80 °C for 8 h. After the completion of the reaction, confirmed by TLC, the mixture was cooled to room temperature and then diluted with saturated saline water (3×15 mL), saturated sodium thiosulfate solution (2×15 mL) and extracted with ethyl acetate. The combined organic layer was collected and dried over anhydrous Na_2SO_4 . The residue was purified by column chromatography on silica gel to afford the desired products (eluent: ethyl acetate/petroleum ether).

3. General procedure for the application (5)¹:

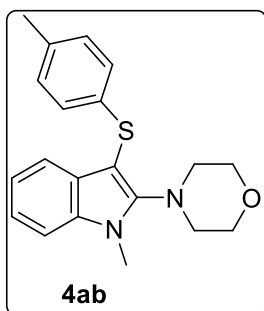
A mixture of **4ay**, PhI(OAc)₂ (2.5 equiv) and AcONH₄ (2 equiv) was stirred in the presence of MeOH as solvent at room temperature. After the completion of the reaction, confirmed by TLC, the solvent is removed by rotary evaporator. Then, the mixture was cooled to room temperature and then diluted with saturated saline water (3×15 mL), saturated sodium thiosulfate solution (2×15 mL) and extracted with ethyl acetate. The combined organic layer was collected and dried over anhydrous Na₂SO₄. The residue was purified by column chromatography on silica gel to afford the desired products (eluent: DCM/MeOH).



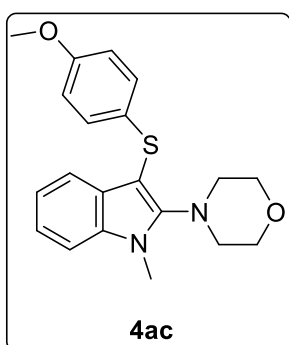
4. Characterization data of the synthesized compounds:



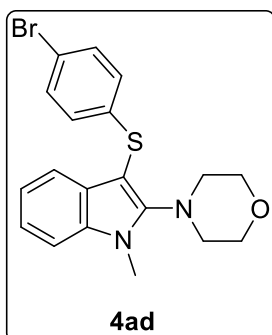
4-(1-methyl-3-(phenylthio)-1H-indol-2-yl)morpholine (4aa)²: Yield: 83%; 268 mg; yellowish white solid; m.p.: 121-122 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/14); ¹H NMR (CDCl₃, 400 MHz): δ 7.62 (d, *J*=7.6 Hz, 1H), 7.34-7.32 (m, 2H), 7.26-7.20 (m, 3H), 7.15-7.11 (m, 3H), 3.90- 3.88 (m, 4H), 3.80 (s, 3H), 3.42- 3.40 (m, 4H); ¹³C {¹H} NMR (CDCl₃, 100 MHz): δ 150.7, 140.4, 134.8, 129.6, 128.8, 125.2, 124.5, 122.2, 120.8, 118.7, 109.5, 90.9, 67.7, 51.6, 29.1.



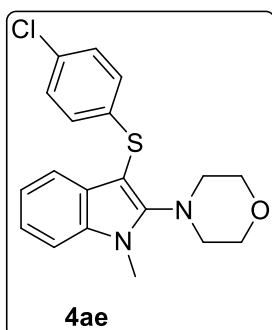
4-(1-methyl-3-(p-tolylthio)-1H-indol-2-yl)morpholine (4ab): Yield: 80%, 270 mg; white solid; m.p.: 127-129 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/10); ^1H NMR (CDCl_3 , 400 MHz): δ 7.54 (d, $J=7.6$ Hz, 1H), 7.30-7.22 (m, 2H), 7.15-7.11 (m, 1H), 6.99-6.95 (m, 4H), 3.82-3.80 (m, 4H), 3.71 (s, 3H), 3.33- 3.31 (m, 4H), 2.26 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (CDCl_3 , 100 MHz): δ 150.6, 136.8, 134.8, 134.2, 129.7, 129.6, 125.3, 122.1, 120.7, 118.8, 109.5, 91.4, 67.7, 51.6, 29.1, 20.9. HRMS (ESI-TOF) m/z: $[\text{M}+\text{H}]^+$ Calculated for $[\text{C}_{20}\text{H}_{23}\text{N}_2\text{OS}]^+$: 339.1526; Found : 339.1532.



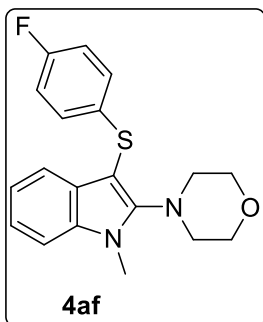
4-(3-((4-methoxyphenyl)thio)-1-methyl-1H-indol-2-yl)morpholine (4ac): Yield: 80%, 270 mg; yellowish white solid; m.p.: 158-160 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/8); ^1H NMR (CDCl_3 , 400 MHz): δ 7.60 (d, $J=7.6$ Hz, 1H), 7.31-7.25 (m, 2H), 7.18-7.14 (m, 1H), 7.06-7.04 (m, 2H), 6.77-6.75 (m, 2H), 3.84-3.82 (m, 4H), 3.74 (s, 3H), 3.71 (s, 3H), 3.35- 3.33 (m, 4H); $^{13}\text{C}\{^1\text{H}\}$ NMR (CDCl_3 , 100 MHz): δ 156.4, 149.4, 133.7, 130.1, 128.7, 126.1, 121.0, 119.6, 117.7, 113.6, 108.4, 91.2, 66.6, 54.4, 50.6, 28.0. HRMS (ESI-TOF) m/z: $[\text{M}+\text{H}]^+$ Calculated for $[\text{C}_{20}\text{H}_{23}\text{N}_2\text{O}_2\text{S}]^+$: 355.1475; Found : 355.1482.



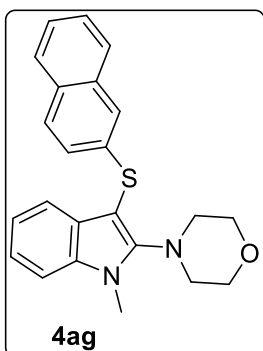
4-(3-((4-bromophenyl)thio)-1-methyl-1H-indol-2-yl)morpholine (4ad): Yield: 79%, 317 mg; white solid; m.p.: 166-168 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/8); ^1H NMR (CDCl_3 , 400 MHz): δ 7.46 (d, $J=7.6$ Hz, 1H), 7.28-7.21 (m, 4H), 7.13-7.09 (m, 1H), 6.89-6.86 (m, 2H), 3.78-3.76 (m, 4H), 3.67 (s, 3H), 3.27- 3.25 (m, 4H); ^{13}C $\{^1\text{H}\}$ NMR (CDCl_3 , 100 MHz): δ 150.8, 139.8, 134.9, 131.8, 129.4, 126.7, 122.3, 121.0, 118.5, 118.0, 109.6, 90.2, 67.6, 51.6, 29.2. HRMS (ESI-TOF) m/z : $[\text{M}+\text{H}]^+$ Calculated for $[\text{C}_{19}\text{H}_{20}\text{BrN}_2\text{OS}]^+$: 403.0474; Found : 403.0488.



4-(3-((4-chlorophenyl)thio)-1-methyl-1H-indol-2-yl)morpholine (4ae): Yield: 80%, 286 mg; yellowish white solid; m.p.: 138-140 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/8); ^1H NMR (CDCl_3 , 400 MHz): δ 7.50 (d, $J=7.6$ Hz, 1H), 7.31-7.24 (m, 2H), 7.16-7.10 (m, 3H), 6.98-6.96 (m, 2H), 3.82-3.80 (m, 4H), 3.71 (s, 3H), 3.31- 3.19 (m, 4H); ^{13}C $\{^1\text{H}\}$ NMR (CDCl_3 , 100 MHz): δ 150.8, 139.1, 134.8, 130.2, 129.4, 128.9, 126.4, 122.3, 120.9, 118.6, 109.6, 90.4, 67.6, 51.6, 29.2. HRMS (ESI-TOF) m/z : $[\text{M}+\text{H}]^+$ Calculated for $[\text{C}_{19}\text{H}_{20}\text{ClN}_2\text{OS}]^+$: 359.0979; Found : 359.0981.

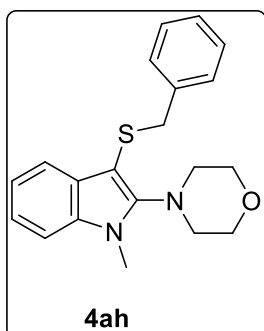


4-(3-((4-fluorophenyl)thio)-1-methyl-1H-indol-2-yl)morpholine (4af): Yield: 82%, 280 mg; brown solid; m.p.: 117-119 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/12); ¹H NMR (CDCl₃, 400 MHz): δ 7.45 (d, *J*=7.6 Hz, 1H), 7.22-7.15 (m, 2H), 7.05-7.05 (m, 1H), 6.95-6.92 (m, 2H), 6.80-6.76 (m, 2H), 3.74-3.72 (m, 4H), 3.62 (s, 3H), 3.22 (m, 4H); ¹³C {¹H} NMR (CDCl₃, 100 MHz): δ 160.7 (d, *J*=241Hz), , 159.5, 150.6, 135.4, 134.8, 129.5, 126.8 (d, *J*=7 Hz), , 122.2, 120.8, 118.5, 115.8 (d, *J*=22 Hz), , 109.6, 91.2, 67.6, 51.5, 29.1. HRMS (ESI-TOF) m/z: [M+H]⁺ Calculated for [C₁₉H₂₀FN₂OS]⁺ : 343.1275; Found : 343.1283.

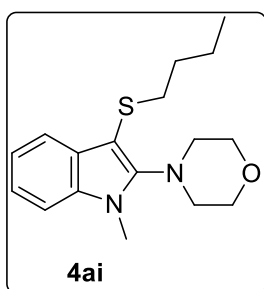


4-(1-methyl-3-(naphthalen-2-ylthio)-1H-indol-2-yl)morpholine (4ag): Yield: 76%, 284 mg; brown solid; m.p.: 96-98 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/12); ¹H NMR (CDCl₃, 400 MHz): δ 7.63 (d, *J*=7.6 Hz, 1H), 7.55 (d, *J*=8.4 Hz, 1H), 7.47 (d, *J*=7.6 Hz, 2H), 7.32 (s, 1H), 7.29-7.22 (m, 3H), 7.19-7.14 (m, 2H), 7.06-7.02 (m, 1H), 3.70-3.68 (m, 4H), 3.64 (s, 3H), 3.26-3.24 (m, 4H); ¹³C {¹H} NMR (CDCl₃, 100 MHz): δ 150.9, 138.1, 134.9, 133.9, 131.3, 129.8, 129.4, 128.4, 127.8, 126.9, 126.4, 125.0, 124.3, 122.5, 122.2,

120.9, 118.8, 109.6, 67.6, 51.6, 29.2. HRMS (ESI-TOF) m/z : $[M+H]^+$ Calculated for $[C_{23}H_{23}N_2OS]^+$: 375.1526; Found : 375.1526.

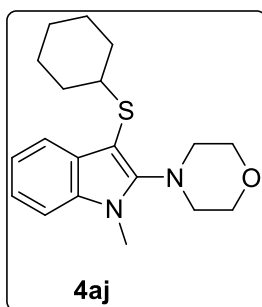


4-(3-(benzylthio)-1-methyl-1H-indol-2-yl)morpholine (4ah): Yield: 79%, 267 mg; brown solid; m.p.: 128-130 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/12); 1H NMR ($CDCl_3$, 400 MHz): δ 7.65 (d, $J=7.6$ Hz, 1H), 7.25-7.21 (m, 3H), 7.20-7.18 (m, 2H), 7.17-7.15 (m, 1H), 7.03-7.01 (m, 2H), 3.85 (s, 2H), 3.71-3.69 (m, 4H), 3.58 (m, 3H), 2.97 (s, 4H); $^{13}C\{^1H\}$ NMR ($CDCl_3$, 100 MHz): δ 150.6, 138.9, 134.8, 132.6, 129.2, 128.3, 126.7, 121.8, 120.3, 118.5, 109.5, 93.5, 67.7, 51.9, 41.1, 28.9. HRMS (ESI-TOF) m/z : $[M+H]^+$ Calculated for $[C_{20}H_{23}N_2OS]^+$: 339.1526; Found : 339.1537.

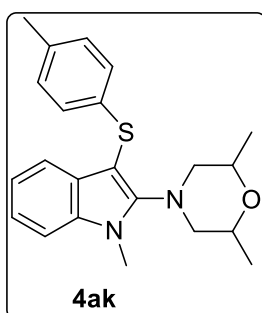


4-(3-(butylthio)-1-methyl-1H-indol-2-yl)morpholine (4ai): Yield: 77%, 234 mg; colourless oil; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/15); 1H NMR ($CDCl_3$, 400 MHz): δ 7.71 (d, $J=7.6$ Hz, 1H), 7.25-7.23 (m, 2H), 7.21-7.19 (m, 1H), 3.89-3.87 (m, 4H), 3.65 (s, 3H), 3.42-4.40 (m, 4H), 2.7-2.69 (m, 2H), 1.64-1.57 (m, 2H), 1.49-1.40 (m, 2H), 0.92 (t, $J=7.6$, 3H); $^{13}C\{^1H\}$ NMR ($CDCl_3$, 100 MHz): δ 149.6, 134.5, 129.9, 121.7, 120.1, 118.7,

109.3, 95.4, 67.7, 52.1, 37.4, 32.1, 28.8, 22.1, 13.8. HRMS (ESI-TOF) m/z : $[M+H]^+$ Calculated for $[C_{17}H_{25}N_2OS]^+$: 305.1682; Found : 305.1693.

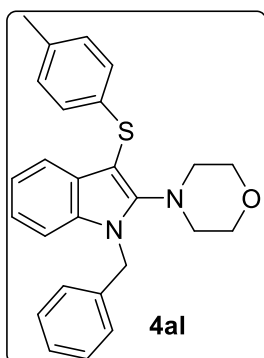


4-(3-(cyclohexylthio)-1-methyl-1H-indol-2-yl)morpholine (4aj): Yield: 76%; 250 mg; colourless oil; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/15); 1H NMR ($CDCl_3$, 400 MHz): δ 7.70 (d, $J=7.6$ Hz, 1H), 7.23-7.17 (m, 3H), 3.88-3.86 (m, 4H), 3.65 (s, 3H), 3.41-4.39 (m, 4H), 2.87-2.79 (m, 1H), 1.98-1.93 (m, 2H), 1.79-1.74 (m, 2H), 1.63-1.58 (m, 1H), 1.45-1.35 (m, 2H), 1.28-1.19 (m, 3H); $^{13}C\{^1H\}$ NMR ($CDCl_3$, 100 MHz): δ 149.9, 134.5, 130.6, 121.7, 120.1, 119.1, 109.2, 94.0, 67.8, 52.3, 48.8, 33.9, 29.0, 26.4, 25.9. HRMS (ESI-TOF) m/z : $[M+H]^+$ Calculated for $[C_{19}H_{27}N_2OS]^+$: 331.1839; Found : 331.1856.

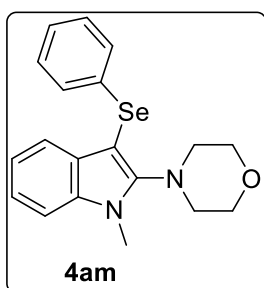


2,6-dimethyl-4-(1-methyl-3-(p-tolylthio)-1H-indol-2-yl)morpholine (4ak): Yield: 78%, 285 mg; yellow gummy mass; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/8); 1H NMR ($CDCl_3$, 400 MHz): δ 7.54 (d, $J=8$ Hz, 1H), 7.30-7.23 (m, 2H), 7.16-7.12 (m, 1H), 6.99 (s, 6H), 3.85-3.82 (m, 4H), 3.69 (s, 3H), 3.27-3.21 (m, 2H), 3.00-2.97 (m, 2H), 2.27 (s, 3H), 1.20 (s, 3H), 1.18 (s, 3H); $^{13}C\{^1H\}$ NMR ($CDCl_3$, 100 MHz): δ 150.4, 136.7, 134.7,

134.1, 129.5, 121.9, 120.6, 118.7, 109.4, 91.2, 72.6, 56.9, 29.2, 20.9, 18.9. HRMS (ESI-TOF) m/z: $[M+H]^+$ Calculated for $[C_{22}H_{27}N_2OS]^+$: 367.1839; Found : 367.1851.

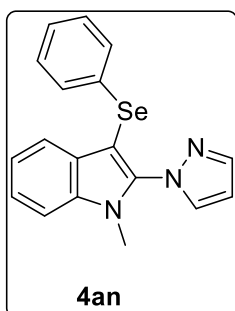


4-(1-benzyl-3-(p-tolylthio)-1H-indol-2-yl)morpholine (4al): Yield: 78 %; 298 mg; yellowish white solid; m.p.: 132-134 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/10); 1H NMR ($CDCl_3$, 400 MHz): δ 7.32-7.29 (m, 2H), 7.25-7.22 (m, 1H), 7.20-7.16 (m, 2H), 7.14-7.08 (m, 4H), 6.98-6.95 (m, 3H), 5.40 (s, 2H), 3.67-3.65 (m, 4H), 3.20 (s, 4H), 2.25 (s, 3H); $^{13}C\{^1H\}$ NMR ($CDCl_3$, 100 MHz): δ 150.7, 138.1, 136.4, 134.6, 134.3, 129.7, 129.7, 128.9, 127.5, 126.3, 125.4, 122.6, 120.9, 119.0, 110.4, 93.5, 67.7, 51.8, 46.2, 21.0. HRMS (ESI-TOF) m/z: $[M+H]^+$ Calculated for $[C_{26}H_{27}N_2OS]^+$: 415.1839; Found : 415.1851.

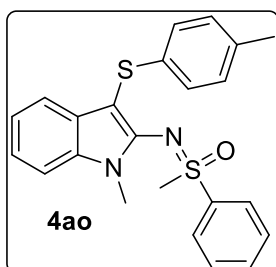


4-(1-methyl-3-(phenylselanyl)-1H-indol-2-yl)morpholine (4am)²: Yield: 78%; 290 mg; red liquid; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/15); 1H NMR ($CDCl_3$, 400 MHz): δ 7.59 (d, $J=7.6$ Hz, 1H), 7.32-7.26 (m, 2H), 7.22-7.18 (m, 2H), 7.16-7.10 (m, 4H), 3.84-3.82 (m, 4H), 3.74 (s, 3H), 3.34 (s, 4H); $^{13}C\{^1H\}$ NMR ($CDCl_3$, 100 MHz): δ

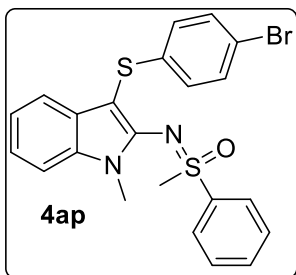
150.5, 135.2, 135.0, 130.4, 129.0, 128.1, 125.4, 122.1, 120.6, 119.6, 109.4, 87.2, 67.7, 51.8, 29.0.



1-methyl-3-(phenylselanyl)-2-(1H-pyrazol-1-yl)-1H-indole (4an)²: Yield: 81 %; 285 mg; white solid; m.p.: 97-98 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/8); ¹H NMR (CDCl₃, 400 MHz): δ 7.88 (d, *J*=1.6 Hz, 1H), 7.76 (d, *J*=8 Hz, 1H), 7.71 (d, *J*=2.4 Hz, 1H), 7.47-7.43 (m, 2H), 7.32-7.25 (m, 4H), 7.21-7.19 (m, 3H), 3.78 (s, 3H); ¹³C{¹H} NMR (CDCl₃, 100 MHz): δ 142.3, 139.2, 135.7, 133.7, 133.4, 129.3, 128.9, 125.9, 124.2, 123.8, 121.5, 121.4, 121.2, 110.1, 106.7, 92.5, 30.7.

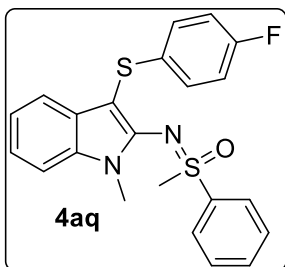


methyl((1-methyl-3-(p-tolylthio)-1H-indol-2-yl)imino)(phenyl)-λ⁶-sulfanone (4ao): Yield: 81%; 328 mg; yellowish solid; m.p.: 129-131 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/5); ¹H NMR (CDCl₃, 400 MHz): δ 8.43 (d, *J*=7.6 Hz, 1H), 7.93-7.89 (m, 1H), 7.83-7.79 (m, 4H), 7.64-7.61 (m, 2H), 7.56-7.52 (m, 2H), 7.47-7.43 (m, 2H), 4.16 (s, 3H), 3.56 (s, 3H), 2.58 (s, 3H); ¹³C{¹H} NMR (CDCl₃, 100 MHz): δ 139.7, 136.2, 135.2, 133.9, 129.7, 129.4, 129.4, 128.3, 125.5, 121.1, 120.4, 118.5, 109.1, 45.0, 29.9, 21.0. HRMS (ESI-TOF) m/z: [M+H]⁺ Calculated for [C₂₃H₂₃N₂OS₂]⁺ : 407.1246; Found : 407.1259.



((3-((4-bromophenyl)thio)-1-methyl-1H-indol-2-yl)imino)(methyl)(phenyl)-λ⁶-sulfanone

(4ap): Yield: 76%; 357 mg; brown solid; m.p.: 150-152 °C column chromatography done (eluent: ethyl acetate/petroleum ether = 1/5); ¹H NMR (CDCl₃, 400 MHz): δ 7.98-7.96 (m, 2H), 7.54-7.50 (m, 1H), 7.42-7.36 (m, 3H), 7.29-7.26 (m, 1H), 7.21-7.19 (m, 1H), 7.15-7.13 (m, 2H), 7.11-7.07 (m, 1H), 6.76-6.74 (m, 2H), 3.80 (s, 3H), 3.24 (s, 3H); ¹³C{¹H} NMR (CDCl₃, 100 MHz): δ 139.5, 139.4, 135.1, 133.5, 131.5, 129.5, 129.3, 128.0, 126.7, 126.6, 121.2, 120.7, 118.1, 117.6, 109.1, 45.8, 29.8. HRMS (ESI-TOF) m/z: [M+H]⁺ Calculated for [C₂₂H₂₀BrN₂OS₂]⁺: 471.0195; Found : 471.0201.

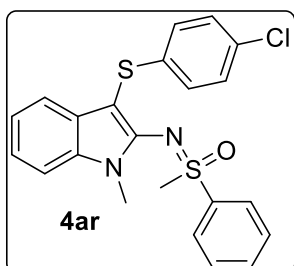


((3-((4-fluorophenyl)thio)-1-methyl-1H-indol-2-yl)imino)(methyl)(phenyl)-λ⁶-sulfanone

(4aq): Yield: 80%; 328 mg; brown solid; m.p.: 119-121 °C column chromatography done (eluent: ethyl acetate/petroleum ether = 1/5); ¹H NMR (CDCl₃, 400 MHz): δ 8.02 (d, *J*=7.6 Hz, 2H), 7.56-7.52 (m, 1H), 7.45-7.41 (m, 3H), 7.28-7.26 (m, 1H), 7.21-7.17 (m, 1H), 7.12-7.08 (m, 1H), 6.90-6.87 (m, 2H), 6.80-6.75 (m, 2H), 3.80 (s, 3H), 3.25 (s, 3H); ¹³C{¹H} NMR (CDCl₃, 100 MHz): δ 144.3, 139.6, 137.6, 135.2, 133.6, 130.4, 129.5, 129.4, 128.1, 127.6, 126.8

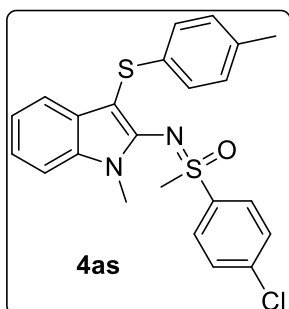
(d, $J=5$ Hz), 121.3 (d, $J=13$ Hz), 120.6, 118.2, 115.9, 115.6 (d, $J=22$ Hz), 109.1, 45.6, 29.9.

HRMS (ESI-TOF) m/z : $[M+H]^+$ Calculated for $[C_{22}H_{20}FN_2OS_2]^+$: 411.0996; Found : 411.1004.



((3-((4-chlorophenyl)thio)-1-methyl-1H-indol-2-yl)imino)(methyl)(phenyl)- λ^6 -sulfanone

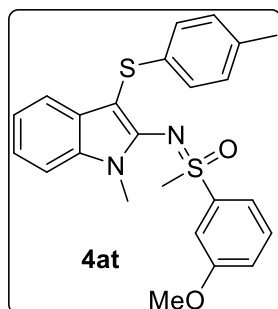
(4ar): Yield: 81%; 345 mg; brown solid; m.p.: 132-134 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/5); 1H NMR ($CDCl_3$, 400 MHz): δ 7.98 (d, $J=8$ Hz, 2H), 7.54-7.51 (m, 1H), 7.43-7.37 (m, 3H), 7.29-7.26 (m, 1H), 7.21-7.17 (m, 1H), 7.11-7.08 (m, 1H), 7.02-6.99 (m, 2H), 6.83-6.81 (m, 2H), 3.80 (s, 3H), 3.25 (s, 3H); $^{13}C\{^1H\}$ NMR ($CDCl_3$, 100 MHz): δ 139.5, 135.2, 133.5, 129.9, 129.5, 129.3, 128.6, 128.0, 126.4, 122.5, 121.3, 121.2, 120.6, 118.2, 109.1, 108.2, 45.8, 29.8. HRMS (ESI-TOF) m/z : $[M+H]^+$ Calculated for $[C_{22}H_{20}ClN_2OS_2]^+$: 427.0700; Found : 427.0709.



(4-chlorophenyl)(methyl)((1-methyl-3-(p-tolylthio)-1H-indol-2-yl)imino)- λ^6 -sulfanone (4as)

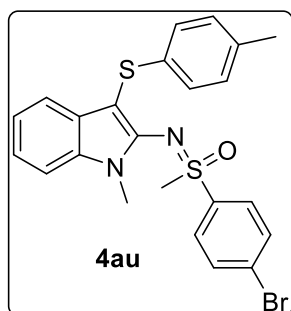
Yield: 77%; 338 mg; yellow solid; m.p.: 128-130 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/5); 1H NMR ($CDCl_3$, 400 MHz): δ 7.91 (d, $J=8.4$ Hz, 2H), 7.41 (d, $J=7.6$ Hz, 2H), 7.29-7.26 (m, 3H), 7.20-7.16 (m, 1H), 7.09-7.06 (m, 1H), 6.86 (d, $J=7.6$, 2H), 6.74 (d, $J=8.4$, 2H), 3.79 (s, 3H), 3.23 (s, 3H), 2.24 (s, 3H); $^{13}C\{^1H\}$ NMR ($CDCl_3$,

100 MHz): δ 140.2, 138.1, 136.2, 135.1, 133.9, 129.7, 129.6, 129.4, 125.0, 121.1, 120.5, 118.4, 109.0, 45.9, 29.8, 21.0. HRMS (ESI-TOF) m/z : $[M+H]^+$ Calculated for $[C_{23}H_{20}ClN_2OS_2]^+$: 441.0857; Found : 441.0861.



(3-methoxyphenyl)(methyl)((1-methyl-3-(p-tolylthio)-1H-indol-2-yl)imino)- λ^6 -sulfanone

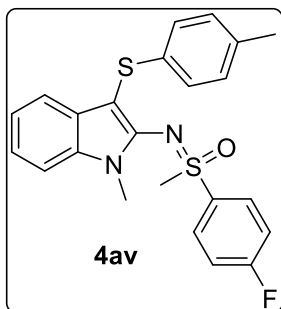
(4at): Yield: 79%; 344 mg; red gummy mass; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/5); 1H NMR ($CDCl_3$, 400 MHz): δ 7.64-7.62 (m, 2H), 7.47 (d, $J=8$ Hz, 2H), 7.37-7.33 (m, 1H), 7.30-7.26 (m, 1H), 7.21-7.18 (m, 1H), 7.12-7.09 (m, 1H), 7.06-7.04 (m, 1H), 6.91-6.87 (m, 4H), 3.82 (s, 3H), 3.69 (s, 3H), 3.21 (s, 3H), 2.23 (s, 3H); $^{13}C\{^1H\}$ NMR ($CDCl_3$, 100 MHz): δ 160.1, 144.3, 136.2, 135.1, 133.8, 130.4, 129.6, 129.4, 125.2, 121.0, 120.5, 120.4, 120.3, 118.4, 112.4, 109.0, 88.2, 55.5, 44.9, 29.8, 20.9. HRMS (ESI-TOF) m/z : $[M+H]^+$ Calculated for $[C_{24}H_{25}N_2O_2S_2]^+$: 437.1352; Found : 437.1363.



(4-bromophenyl)(methyl)((1-methyl-3-(p-tolylthio)-1H-indol-2-yl)imino)- λ^6 -sulfanone

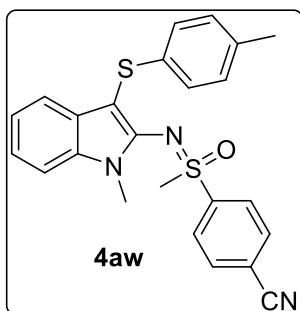
(4au): Yield: 77%; 344 mg; brown solid; m.p.: 162-164 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/5); 1H NMR ($CDCl_3$, 400 MHz): δ 7.66 (d, $J=7.6$ Hz,

2H), 7.29-7.26 (m, 3H), 7.28-7.25 (m, 3H), 7.11 (d, $J=9.2$ Hz, 1H), 7.05-7.01 (m, 1H), 6.95-6.91 (m, 1H), 6.70 (d, $J=8$ Hz, 2H), 6.56 (d, $J=8$ Hz, 2H), 3.64 (s, 3H), 3.08 (s, 3H), 2.10 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (CDCl_3 , 100 MHz): δ 143.8, 138.5, 136.2, 135.1, 133.9, 132.5, 129.6, 129.4, 128.7, 125.2, 124.9, 121.1, 120.5, 118.4, 108.9, 87.0, 46.0, 29.8, 21.0. HRMS (ESI-TOF) m/z : $[\text{M}+\text{H}]^+$ Calculated for $[\text{C}_{23}\text{H}_{22}\text{BrN}_2\text{OS}_2]^+$: 485.0351; Found : 485.0364.



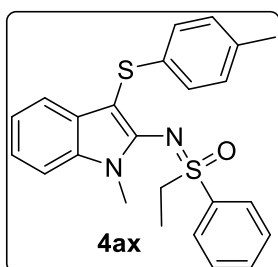
(4-fluorophenyl)(methyl)((1-methyl-3-(p-tolylthio)-1H-indol-2-yl)imino)- λ^6 -sulfanone

(4av): Yield: 75%; 331 mg; yellow solid; m.p.: 132-134 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/5); ^1H NMR (CDCl_3 , 400 MHz): δ 8.05-8.02 (m, 2H), 7.43 (d, $J=8.4$ Hz, 1H), 7.29-7.26 (m, 1H), 7.21-7.17 (m, 1H), 7.11-7.02 (m, 3H), 6.89 (d, $J=8$ Hz, 2H), 6.8 (d, $J=8$ Hz, 2H), 3.79 (s, 3H), 3.22 (s, 3H), 2.23 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (CDCl_3 , 100 MHz): δ 164.5, 143.8, 135.6 (d, $J=106$ Hz), 134.0, 131.2, 131.2, 129.6 (d, $J=10$ Hz), 128.5, 125.2, 120.9 (d, $J=67$ Hz), 118.5, 116.6 (d, $J=23$ Hz), 109.1, 88.1, 45.5, 29.9, 20.9. HRMS (ESI-TOF) m/z : $[\text{M}+\text{H}]^+$ Calculated for $[\text{C}_{23}\text{H}_{22}\text{FN}_2\text{OS}_2]^+$: 424.1079; Found : 424.1081.

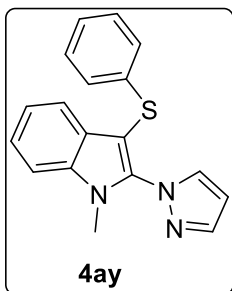


4-(S-methyl-N-(1-methyl-3-(p-tolylthio)-1H-indol-2-yl)sulfonimidoyl)benzonitrile (4aw):

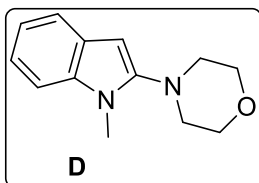
Yield: 47%; 202 mg; brown solid; m.p.: 152-153 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/4); ^1H NMR (CDCl_3 , 400 MHz): δ 7.99 (d, $J=7.6$ Hz, 2H), 7.46 (d, $J=8$ Hz, 2H), 7.34 (d, $J=8$ Hz, 1H), 7.27-7.26 (m, 1H), 7.20-7.16 (m, 1H), 7.08-7.04 (m, 1H), 6.81 (d, $J=8.4$ Hz, 2H), 6.54(d, $J=7.6$ Hz, 2H), 3.79 (s, 3H), 3.29 (s, 3H), 2.26 (s, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (CDCl_3 , 100 MHz): δ 143.9, 132.9, 129.5, 129.3, 128.5, 124.6, 121.2, 120.6, 118.4, 117.3, 116.5, 108.9, 46.4, 29.7, 20.9. HRMS (ESI-TOF) m/z : $[\text{M}+\text{H}]^+$ Calculated for $[\text{C}_{23}\text{H}_{22}\text{BrN}_2\text{OS}_2]^+$: 431.1126; Found : 431.1133.



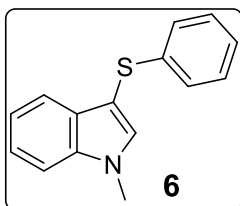
ethyl((1-methyl-3-(p-tolylthio)-1H-indol-2-yl)imino)(phenyl)- λ^6 -sulfanone (4ax): Yield: 76%; 319 mg; yellow solid; m.p.: 102-104 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/5); ^1H NMR (CDCl_3 , 400 MHz): δ 7.95 (d, $J=7.6$ Hz, 2H), 7.50-7.47 (m, 1H), 7.41-7.36 (m, 3H), 7.26-7.24 (m, 1H), 7.17-7.14 (m, 1H), 7.08-7.04 (m, 1H), 6.87-6.80 (m, 4H), 3.80 (s, 3H), 3.41-3.36 (m, 2H), 2.22 (s, 3H), 1.21 (t, $J=7.6$ Hz, 3H); $^{13}\text{C}\{^1\text{H}\}$ NMR (CDCl_3 , 100 MHz): δ 137.6, 136.5, 135.1, 133.7, 133.3, 129.8, 129.3, 129.3, 128.9, 125.4, 120.8, 120.3, 118.3, 108.8, 87.6, 52.0, 29.8, 21.0, 8.2. HRMS (ESI-TOF) m/z : $[\text{M}+\text{H}]^+$ Calculated for $[\text{C}_{24}\text{H}_{25}\text{N}_2\text{OS}_2]^+$: 421.1403; Found : 421.1415.



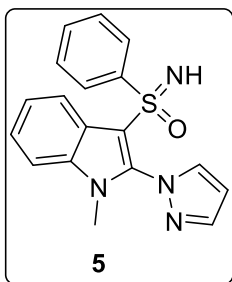
1-methyl-3-(phenylthio)-2-(1H-pyrazol-1-yl)-1H-indole (4ay)²: Yield: 78%; 337 mg; yellow solid; m.p.: 112-114 °C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/8); ¹H NMR (CDCl₃, 400 MHz): δ 7.86 (d, *J*=1.6 Hz, 1H), 7.71-7.69 (m, 1H), 7.47-7.40 (m, 3H), 7.27-7.24 (m, 1H), 7.20-7.16 (m, 2H), 7.09-7.08 (m, 3H), 6.47-6.46 (m, 1H), 3.78 (s, 3H); ¹³C{¹H} NMR (CDCl₃, 100 MHz): δ 142.4, 139.1, 138.6, 135.4, 133.6, 129.0, 128.2, 125.2, 123.9, 121.6, 120.3, 110.2, 106.9, 95.6, 30.8.



4-(1-methyl-1H-indol-2-yl)morpholine (D)³: Yield: 89%; 192 mg; white solid; mp: 101-102°C; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/10); ¹H NMR (CDCl₃, 400 MHz): δ 7.52 (d, *J*=7.6 Hz, 1H), 7.25 (d, *J*=8 Hz, 1H), 7.19-7.15 (m, 1H), 7.12-7.08 (m, 1H), 5.97 (s, 1H), 3.92-3.90 (m, 4H), 3.64 (s, 3H) 3.06-3.03 (m, 4H); ¹³C{¹H} NMR (CDCl₃, 100 MHz): δ 150.1, 135.4, 127.6, 120.5, 119.7, 119.5, 108.8, 87.1, 67.0, 52.9, 29.2.



1-methyl-3-(phenylthio)-1H-indole (6)⁴: Yield: 21%; 50 mg; colourless oil; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/25); ¹H NMR (CDCl₃, 400 MHz): δ 7.62 (d, *J* = 8 Hz, 1H), 7.39 (d, *J* = 8.4 Hz, 1H), 7.34-7.29 (m, 2H), 7.19-7.14 (m, 3H), 7.11-7.09 (m, 2H), 7.07-7.04 (m, 1H), 3.86 (s, 3H); ¹³C {¹H} NMR (CDCl₃, 100 MHz): δ 139.8, 137.6, 135.2, 129.9, 128.8, 125.8, 124.8, 122.7, 120.6, 119.9, 109.9, 100.5, 33.3.

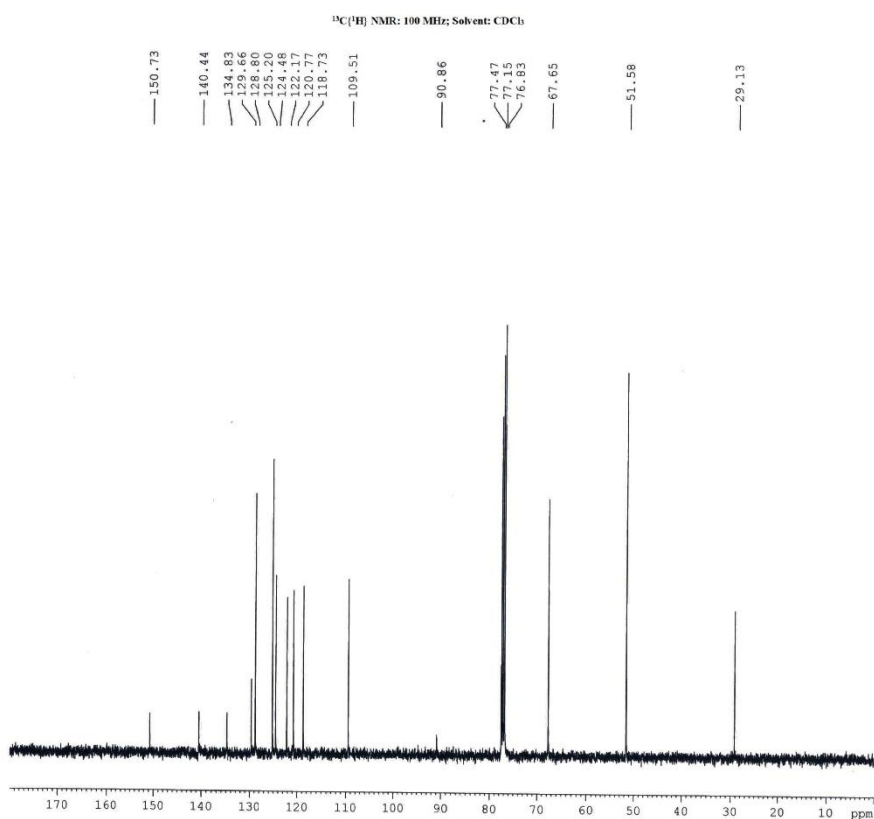
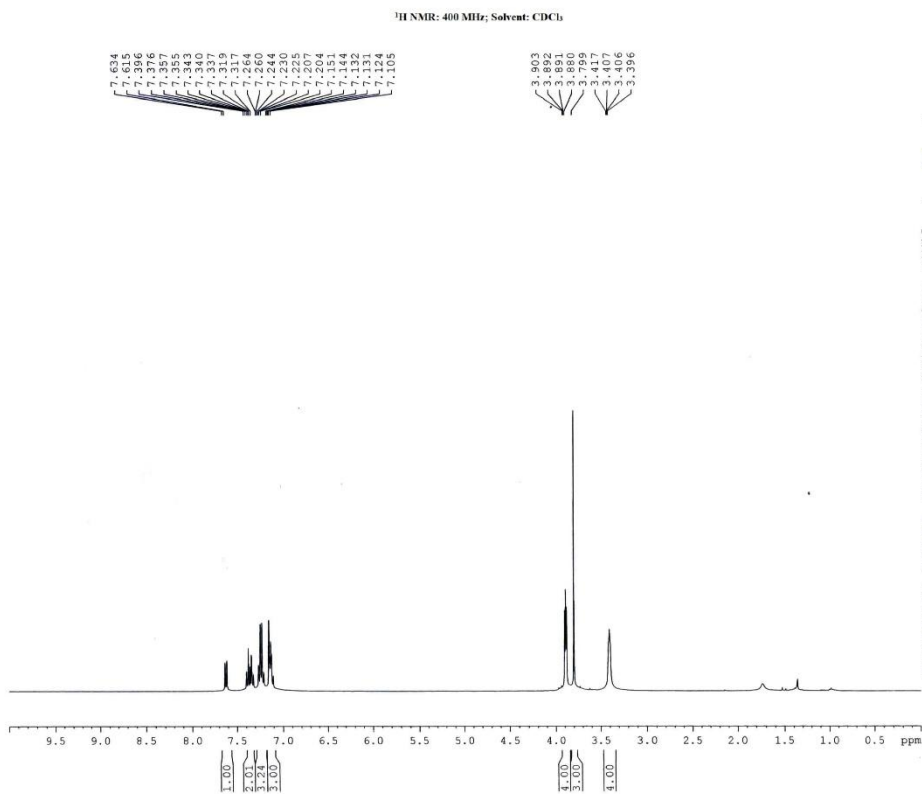


imino(1-methyl-2-(1H-pyrazol-1-yl)-1H-indol-3-yl)(phenyl)-16-sulfanone (5): Yield: 78%; 262 mg; yellow gummy mass; column chromatography done (eluent: ethyl acetate/petroleum ether = 1/8); ¹H NMR (CDCl₃, 400 MHz): δ 8.40-8.48 (m, 1H), 7.94-7.92 (m, 2H), 7.82 (d, *J* = 2.4 Hz, 1H), 7.43-7.33 (m, 6H), 6.57-6.56 (m, 1H), 3.43 (s, 3H), 2.62 (s, 1H); ¹³C {¹H} NMR (CDCl₃, 100 MHz): δ 133.9, 142.6, 135.9, 135.2, 133.9, 132.4, 128.9, 127.2, 124.8, 123.6, 123.1, 121.5, 111.9, 110.3, 107.2, 29.9. Anal. Calcd. For C₁₈H₁₆N₄OS: C, 64.27; H, 4.79; N, 16.65%; Found: C, 64.34; H, 4.89; N, 16.58%.

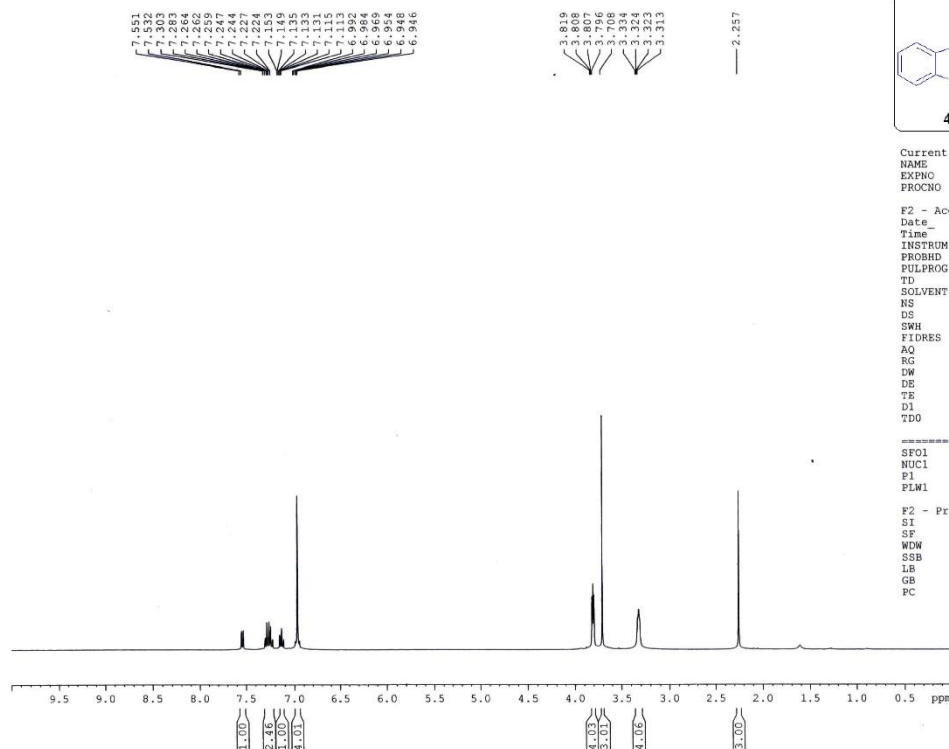
5. References:

1. A. Tota, M. Zenzola, S. J. Chawner, S. S. John-Campbell, C. Carlucci, G. Romanazzi, L. Degennaro, J. A. Bull and R. Luisi, *Chem. Commun. (Camb.)*, 2016, **53**, 348–351.
2. X. Zhang, C. Liu, Y. Zhang, F. Shen, W. Wei, Z. Zhang and T. Liang, *Org. Chem. Front.*, 2023, **10**, 5886–5894.
3. Y.-X. Li, K.-G. Ji, H.-X. Wang, S. Ali and Y.-M. Liang, *J. Org. Chem.*, 2011, **76**, 744–747.
4. P. Sharma and N. Jain, *J. Org. Chem.*, 2019, **84**, 13045–13052.

**6. NMR spectra [^1H , and $^{13}\text{C}\{^1\text{H}\}$] of
synthesized products:**



¹H NMR: 400 MHz; Solvent: CDCl₃



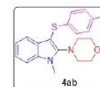
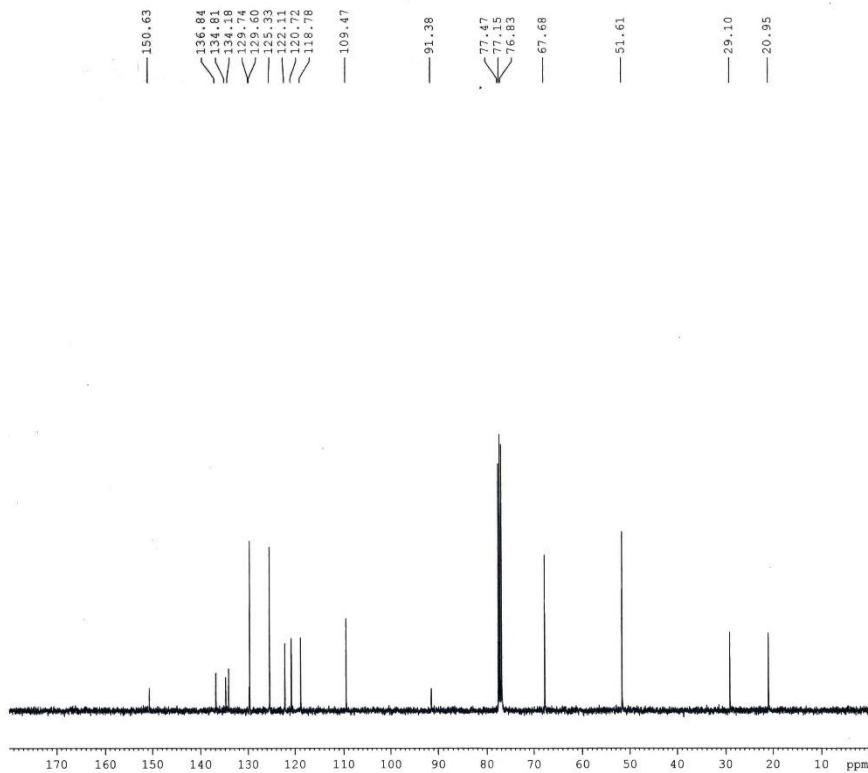
Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 360
PROCNO 1

F2 - Acquisition Parameters
Date_ 20230603
Time_ 12.51
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 8
DS 1
SWH 8223.685 Hz
FIDRES 0.250967 Hz
AQ 1.992294 sec
RG 87.66
DW 60.800 usec
DE 6.50 usec
TE 299.0 K
D1 1.00000000 sec
TD0 1

==== CHANNEL f1 =====
SF01 400.1524711 MHz
NUC1 1H
P1 14.75 usec
PLW1 12.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1500097 MHz
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 2.00

¹³C{¹H} NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 361
PROCNO 1

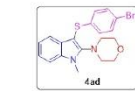
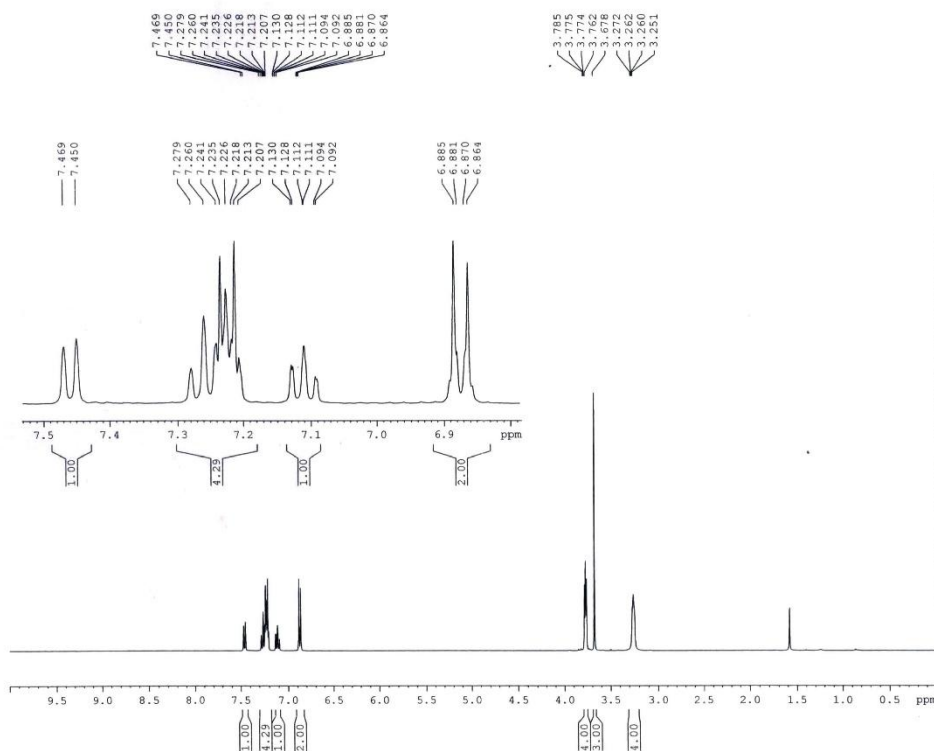
F2 - Acquisition Parameters
Date_ 20230603
Time_ 13.08
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 240
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 87.66
DW 20.800 usec
DE 6.50 usec
TE 299.1 K
D1 2.00000000 sec
D11 0.03000000 sec
TD0 1

==== CHANNEL f1 =====
SF01 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

==== CHANNEL f2 =====
SF02 400.1516006 MHz
NUC2 1H
CDEPRG2 waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6179973 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

¹H NMR: 400 MHz; Solvent: CDCl₃



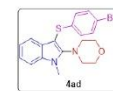
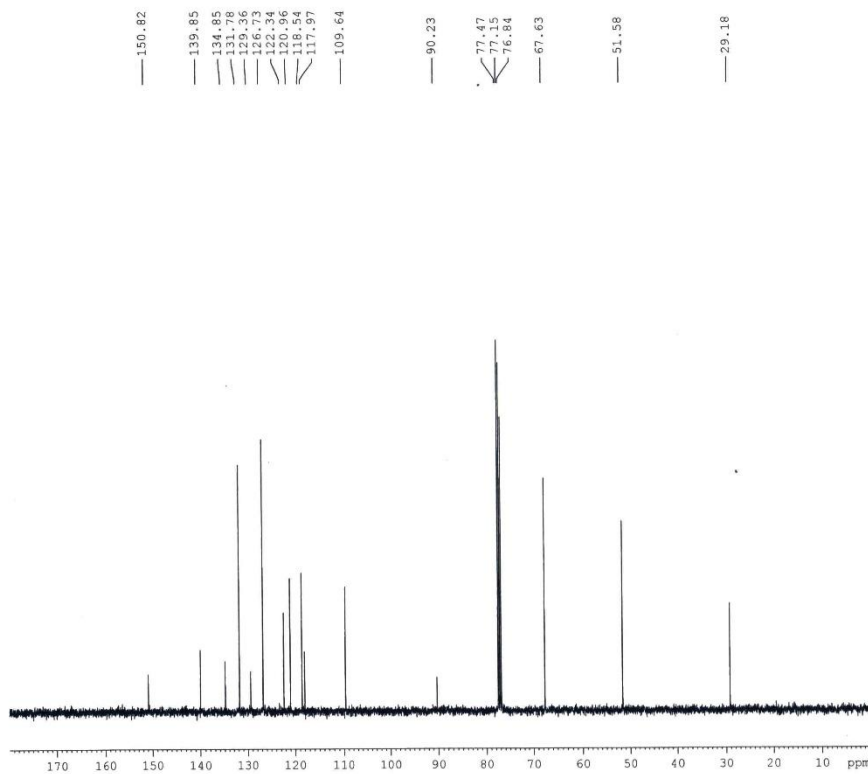
Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 536
PROCNO 1

F2 - Acquisition Parameters
Date 20230727
Time 13.28
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 8
DS 1
SWH 8223.685 Hz
FIDRES 0.250967 Hz
AQ 1.922944 sec
RG 67.81
DW 60.800 usec
DE 6.50 usec
TE 296.8 K
D1 1.00000000 sec
TDO 1

----- CHANNEL f1 -----
SFO1 400.1524711 MHz
NUC1 1H
P1 14.75 usec
PLW1 12.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1500228 MHz
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 2.00

¹³C{¹H} NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 537
PROCNO 1

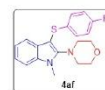
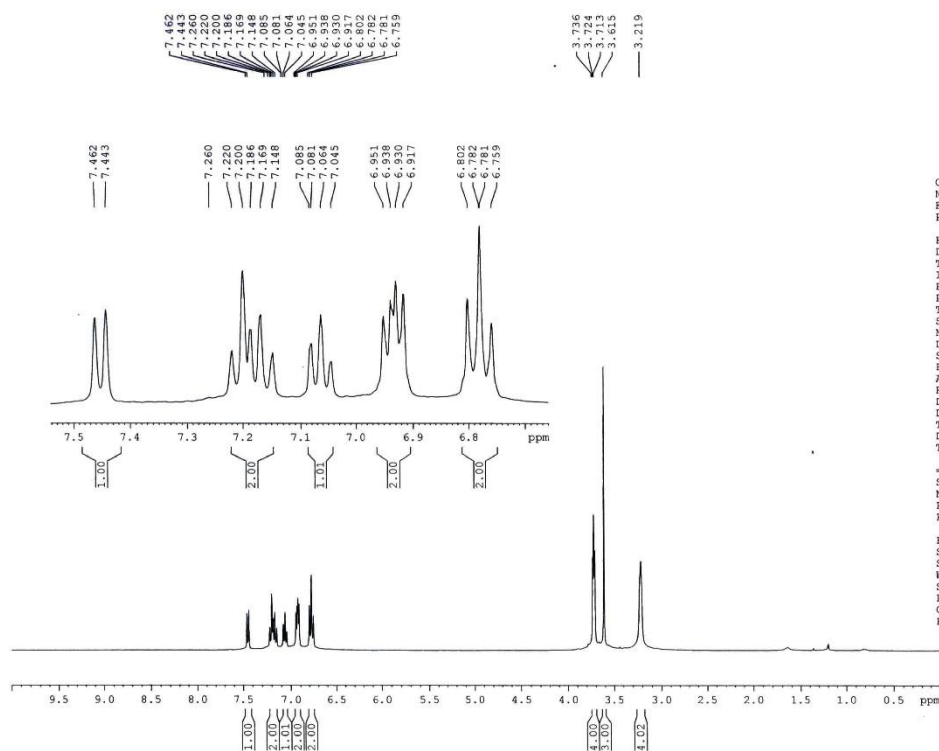
F2 - Acquisition Parameters
Date 20230727
Time 13.39
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 160
DS 2
SWH 24638.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 67.81
DW 20.800 usec
DE 6.50 usec
TE 297.1 K
D1 2.00000000 sec
D11 0.03000000 sec
TDO 1

----- CHANNEL f1 -----
SFO1 100.6276588 MHz
NUC1 13C
P1 6.90 usec
PLW1 54.00000000 W

----- CHANNEL f2 -----
SFO2 400.1516006 MHz
NUC2 1H
CFDPRG12 waltz16
PCFD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32210000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177873 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

¹H NMR: 400 MHz; Solvent: CDCl₃



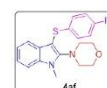
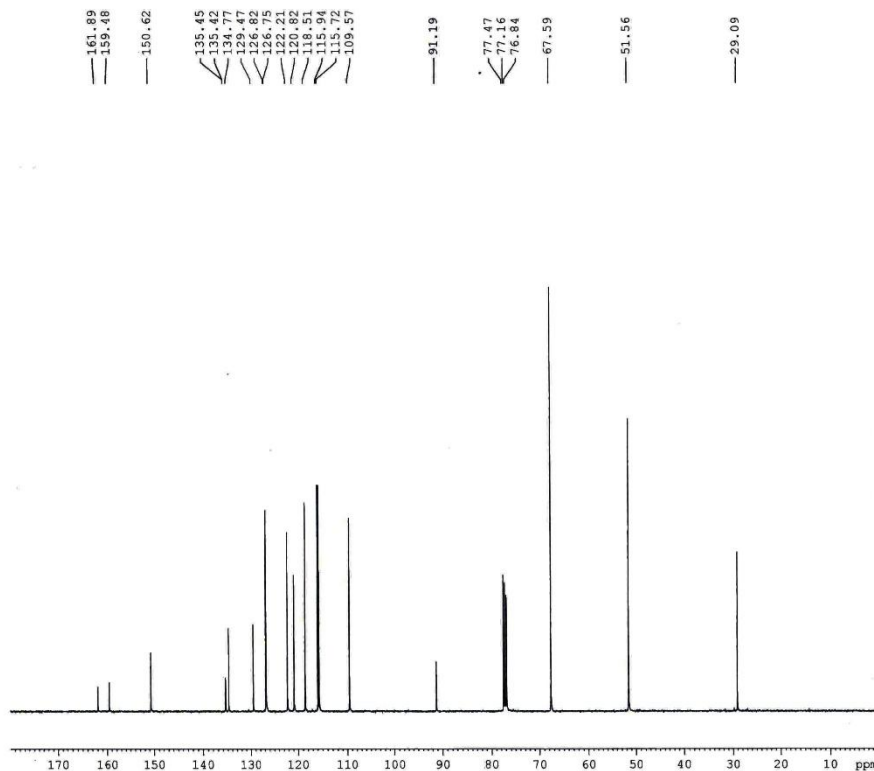
Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 523
PROCNO 1

F2 - Acquisition Parameters
Date_ 20230726
Time 10.45
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 8
DS 1
SWH 8223.685 Hz
FIDRES 0.250967 Hz
AQ 1.9922944 sec
RG 26.53
DW 60.800 usec
DE 6.50 usec
TE 294.0 K
D1 1.0000000 sec
TDC 1

----- CHANNEL f1 -----
SFO1 400.1524711 MHz
NUC1 1H
P1 14.75 usec
PLW1 12.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1500348 MHz
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 2.00

¹³C{¹H} NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 523
PROCNO 1

F2 - Acquisition Parameters
Date_ 20230726
Time 12.32
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 1024
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 186.42
DW 20.800 usec
DE 6.50 usec
TE 296.0 K
D1 2.0000000 sec
D11 0.03000000 sec
TDO 0

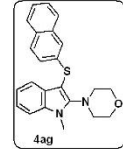
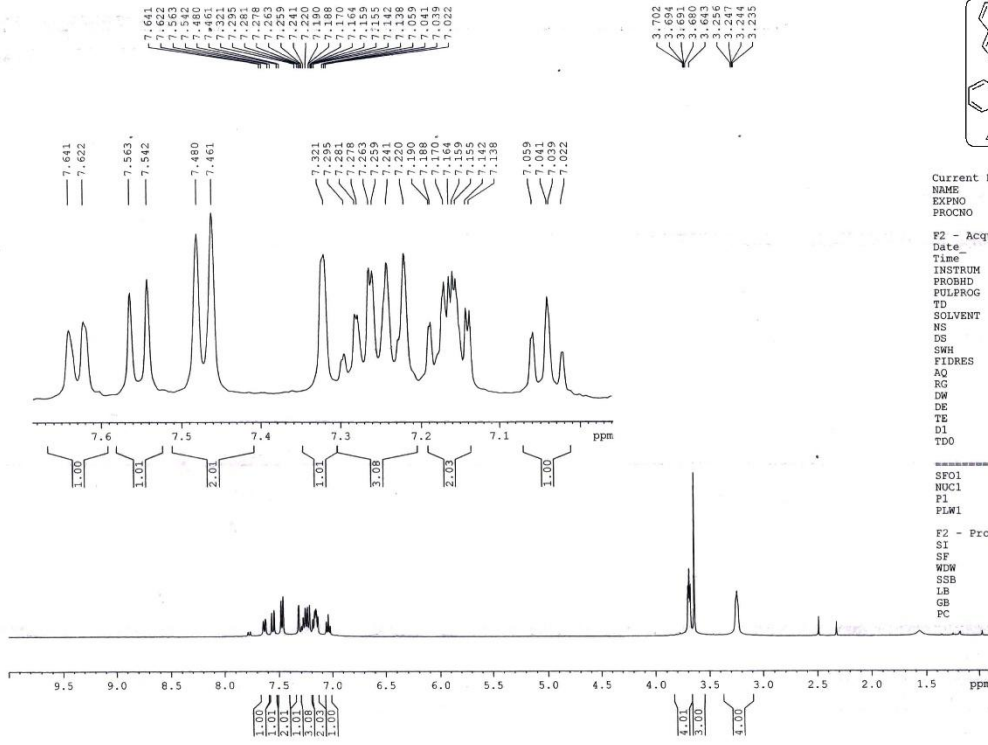
----- CHANNEL f1 -----
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

----- CHANNEL f2 -----
SFO2 400.1516006 MHz
NUC2 1H
CPCPRG[2] waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.3231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177961 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

¹H of VBSP-283/12

¹H NMR: 400 MHz; Solvent: CDCl₃



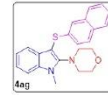
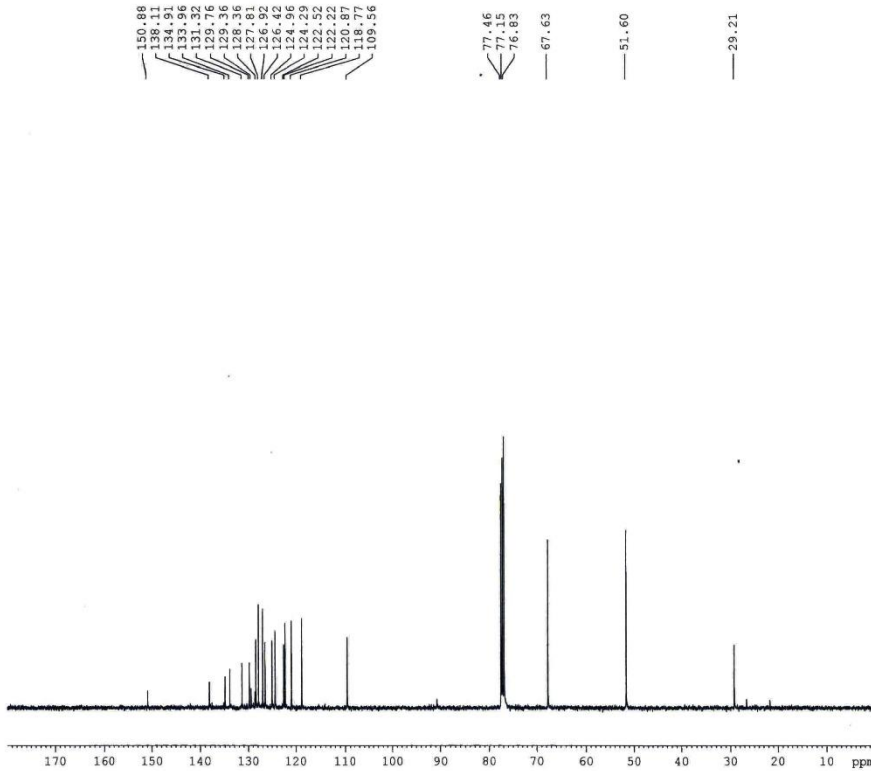
Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 543
PROCNO 1

F2 - Acquisition Parameters
Date_ 20230729
Time 10.21
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 8
DS 1
SWH 8223.685 Hz
FIDRES 0.250867 Hz
AQ 1.9922944 sec
RG 106.66
DW 60.800 use
DE 6.50 use
TE 299.4 K
D1 1.0000000 sec
TDO 1

CHANNEL f1
SFO1 400.1524711 MHz
NUC1 1H
P1 14.75 use
PLW1 12.0000000 W

F2 - Processing parameters
SI 16384
SF 400.1500514 MHz
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 2.00

¹³C{¹H} NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 544
PROCNO 1

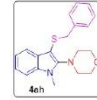
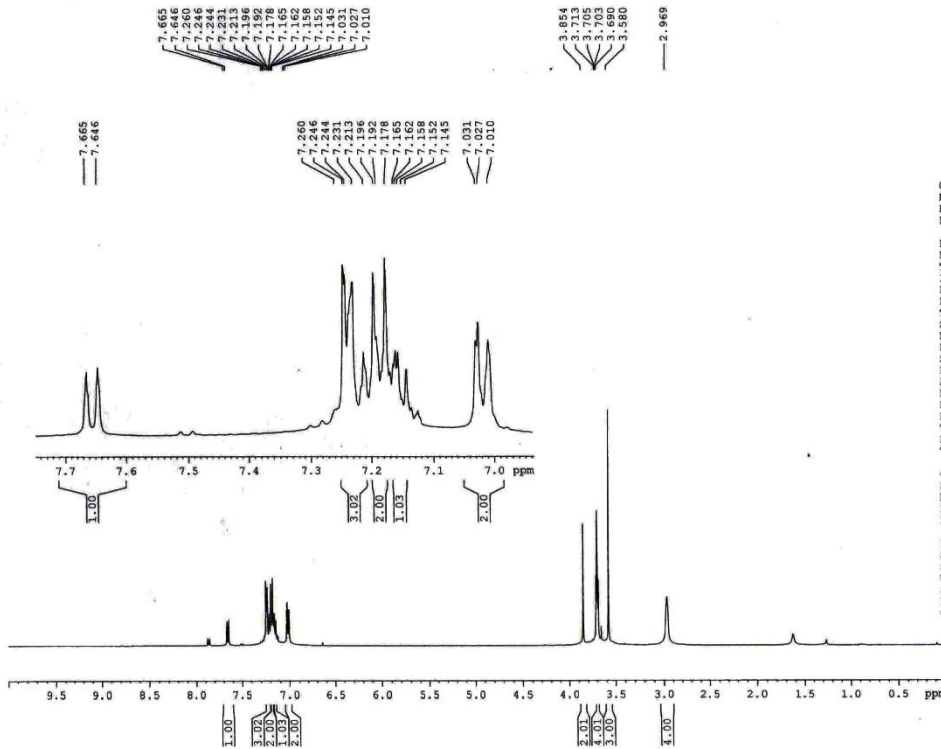
F2 - Acquisition Parameters
Date_ 20230729
Time 11.10
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 720
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6915744 sec
RG 196.66
DW 20.800 use
DE 6.50 use
TE 299.6 K
D1 2.0000000 sec
D11 0.0300000 sec
TDO 1

CHANNEL f1
SFO1 100.6278568 MHz
NUC1 13C
P1 8.90 use
PLW1 54.0000000 W

CHANNEL f2
SFO2 400.1516006 MHz
NUC2 1H
CFPRG2 waltz16
PCPD2 90.00 use
PLW2 12.0000000 W
PLW3 0.12231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177887 MHz
WDW DM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

¹H NMR: 400 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJER 2023
EXPNO 534
PROCNO 1

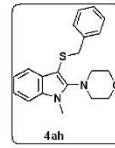
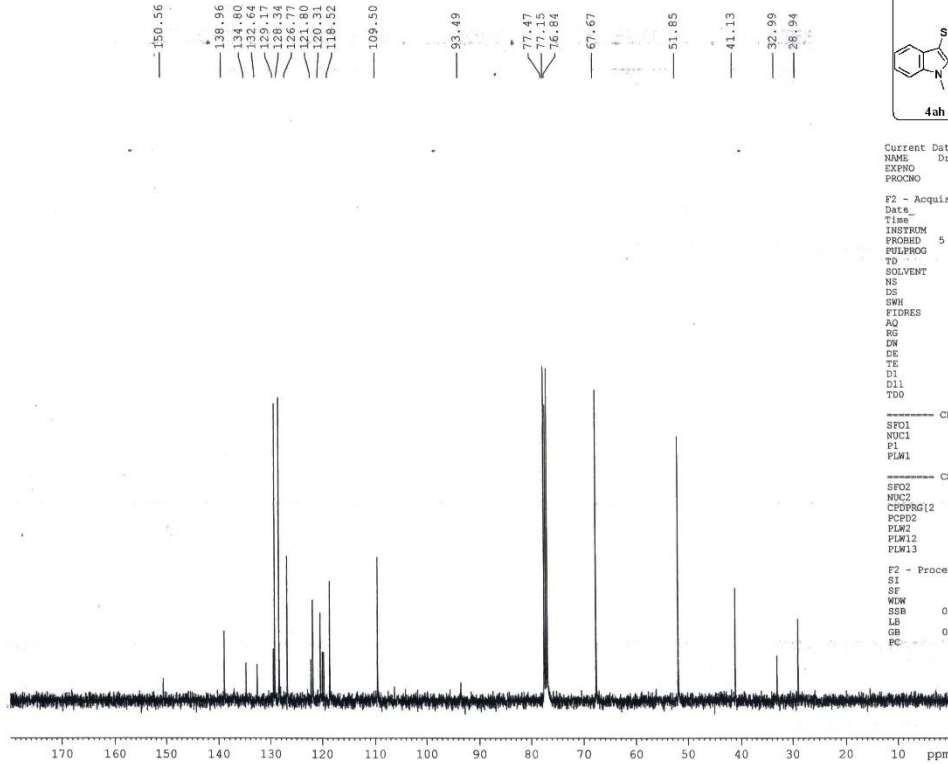
F2 - Acquisition Parameter
Date 20230726
Time 21.42
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl₃
NS 8
DS 1
SWH 8223.685 Hz
FIDRES 0.250967 Hz
AQ 1.9922944 se
RG 67.61
DW 60.800 us
DE 6.50 us
TE 296.1 K
SI 1.00000000 se
TDO 1

CHANNEL f1
SFO1 400.1524711 MH
NUC1 1H
PI 14.75 us
PLW1 12.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1500147 MH
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 2.00

¹³C of VBSP-283/3

¹³C{¹H} NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJER 2023
EXPNO 535
PROCNO 1

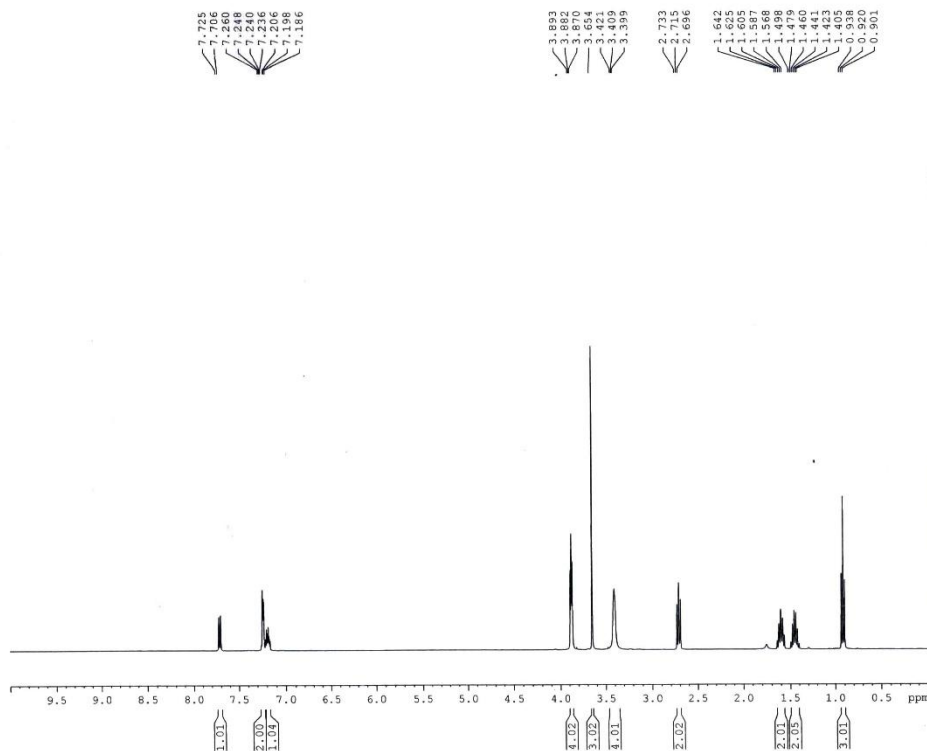
F2 - Acquisition Parameters
Date 20230726
Time 22.03
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl₃
NS 120
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 67.61
DW 20.800 usec
DE 6.50 usec
TE 296.8 K
SI 2.00000000 sec
D11 0.03000000 sec
TDO 1

CHANNEL f1
SFO1 100.6278568 MHz
NUC1 13C
PI 8.90 usec
PLW1 54.00000000 W

CHANNEL f2
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG2 waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177894 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

¹H NMR: 400 MHz; Solvent: CDCl₃



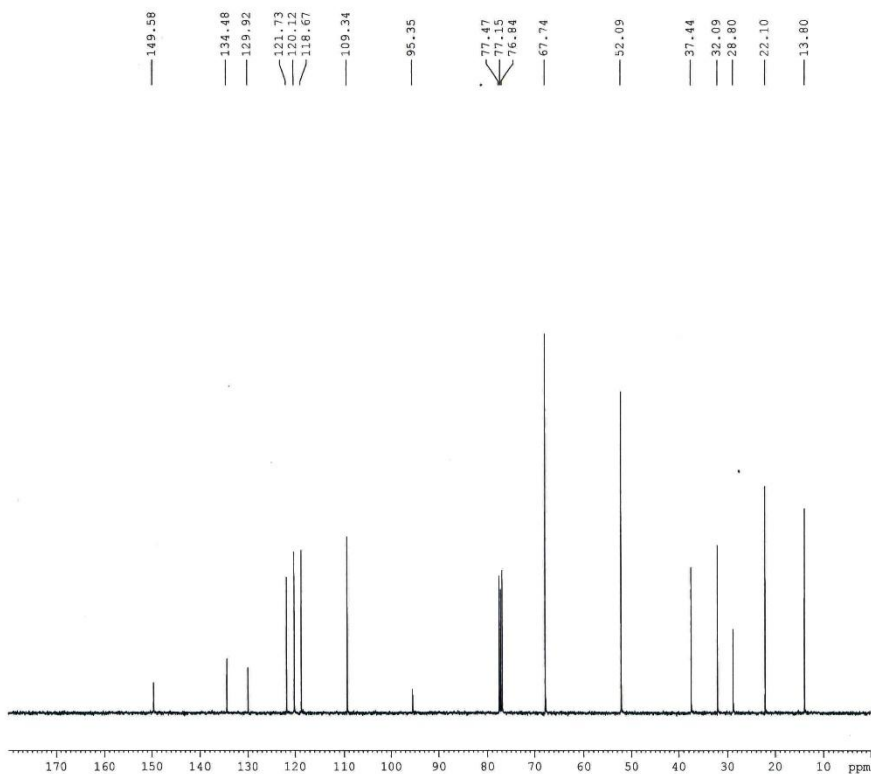
Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 487
PROCNO 1

F2 - Acquisition Parameters
Date_ 20230712
Time 0.25
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 8
DS 1
SWH 8223.685 Hz
FIDRES 0.250967 Hz
AQ 1.992344 sec
RG 32.25
DW 60.800 usec
DE 6.50 usec
TE 298.2 K
D1 1.00000000 sec
TDO 1

==== CHANNEL f1 =====
SF01 400.1524711 MHz
NUC1 1H
P1 14.75 usec
PLW1 12.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1500122 MHz
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 2.00

¹³C{¹H} NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 488
PROCNO 1

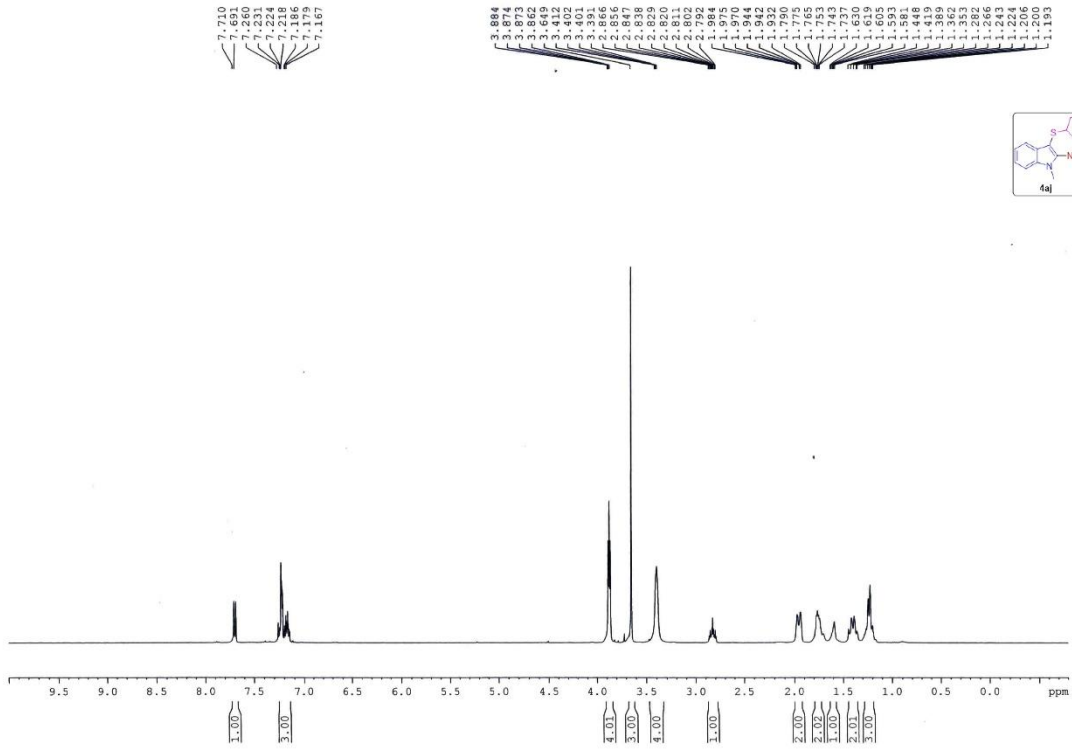
F2 - Acquisition Parameters
Date_ 20230712
Time 0.39
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 256
DS 2
SWH 24038.461 Hz
FIDRES 0.733896 Hz
AQ 0.6815744 sec
RG 32.25
DW 20.800 usec
DE 6.50 usec
TE 298.2 K
D1 2.00000000 sec
D11 0.03000000 sec
TDC 1

==== CHANNEL f1 =====
SF01 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

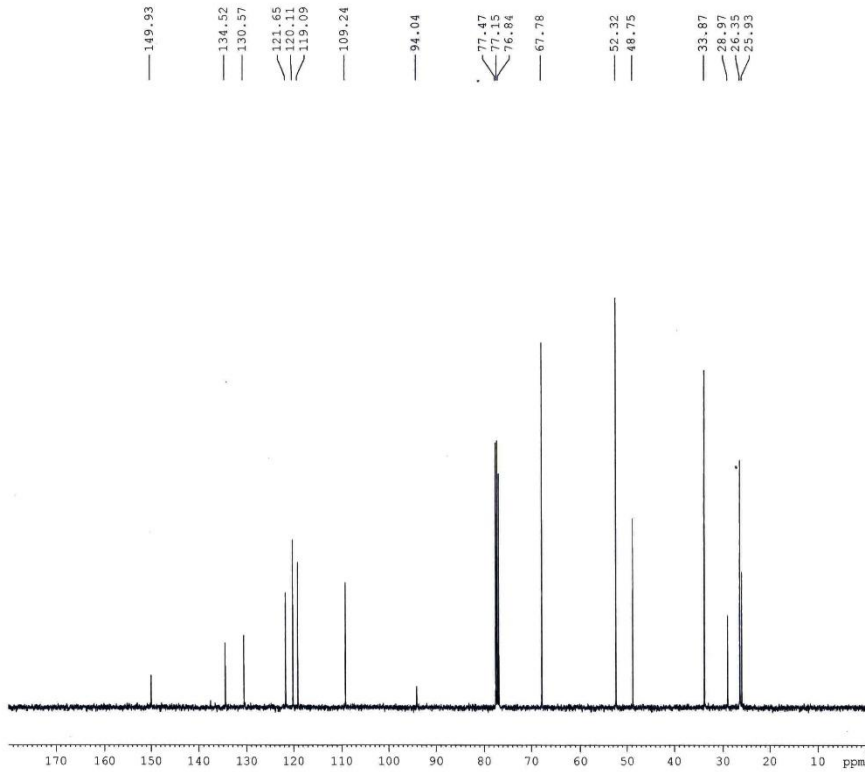
==== CHANNEL f2 =====
SF02 400.1516006 MHz
NUC2 1H
CPDPRG2 waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.3221000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177955 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

¹H NMR: 400 MHz; Solvent: CDCl₃



¹³C{¹H} NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 482
PROCNO 1

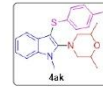
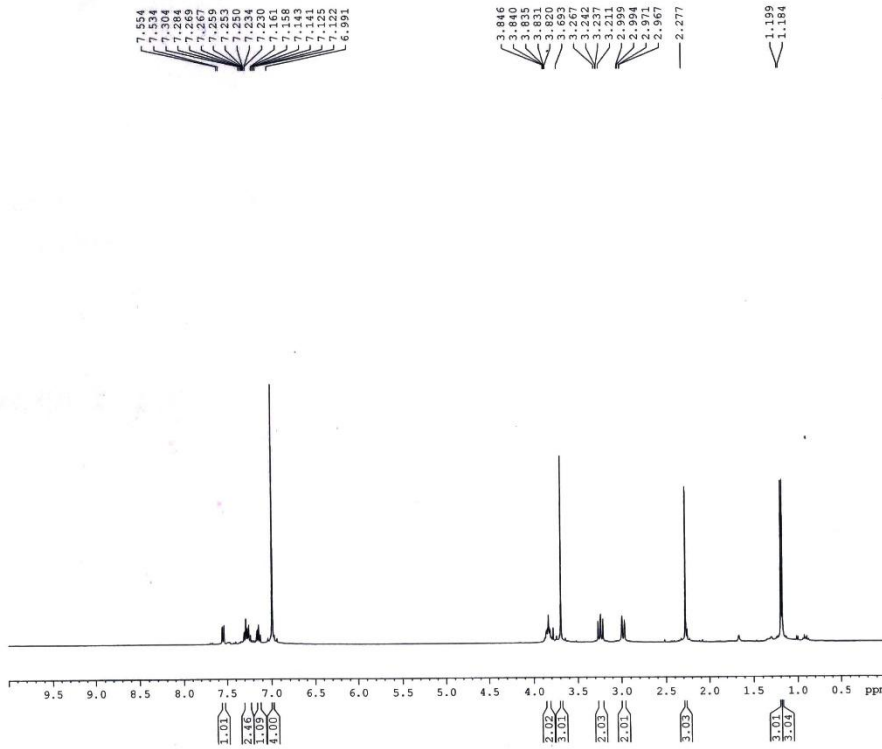
F2 - Acquisition Parameters
Date_ 20230709
Time 16.03
INSTRUM spect
PROBHD 5-mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl₃
NS 512
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 5.13
DW 20.800 usec
DE 6.50 usec
TE 297.0 K
D1 2.00000000 sec
D11 0.03000000 sec
YD0 1

----- CHANNEL f1 -----
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

----- CHANNEL f2 -----
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG2 waltz16
PCPD2 90.00 usec
PLM2 12.00000000 W
PLW2 0.32231000 W
PLM13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177902 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

¹H NMR: 400 MHz; Solvent: CDCl₃



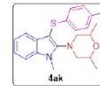
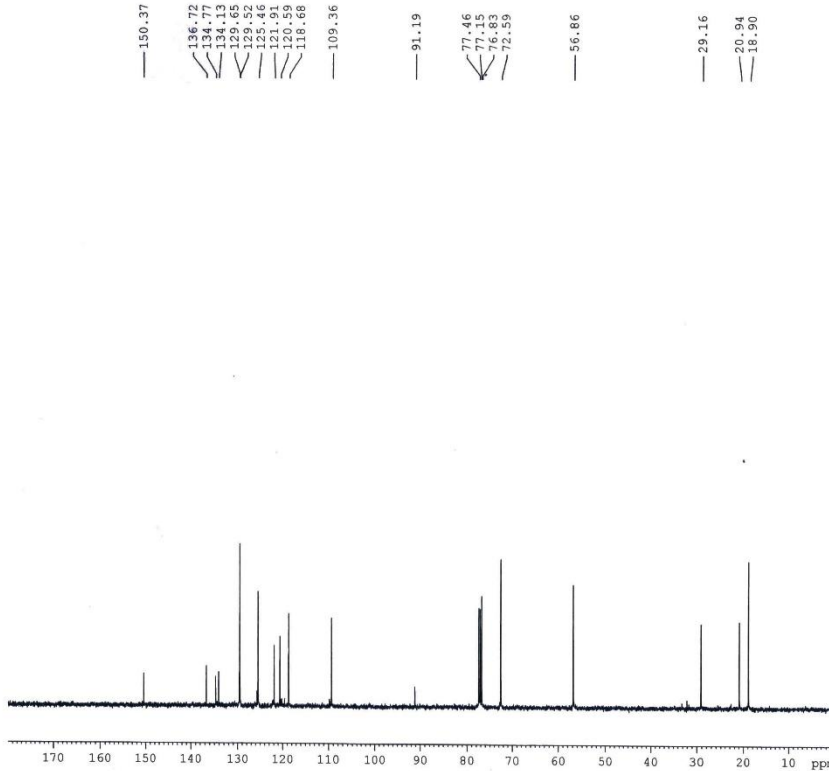
Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 870
PROCNO 1

F2 - Acquisition Parameters
Date_ 20231203
Time 0.58
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT cdcl3
NS 8
DS 1
SWH 9223.685 Hz
FIDRES 0.250967 Hz
AQ 1.9922944 sec
RG 40.87
DW 60.800 usec
DE 6.50 usec
TE 292.6 K
D1 1.0000000 sec
TDO 1

----- CHANNEL f1 -----
SF01 400.1524711 MHz
NUC1 1H
P1 14.75 usec
PLW1 12.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1500097 MHz
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 2.00

¹³C{¹H} NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 871
PROCNO 1

F2 - Acquisition Parameters
Date_ 20231203
Time 1.10
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 120
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6813746 sec
RG 40.87
DW 20.800 usec
DE 6.50 usec
TE 293.6 K
D1 2.0000000 sec
D11 0.0300000 sec
TDO 1

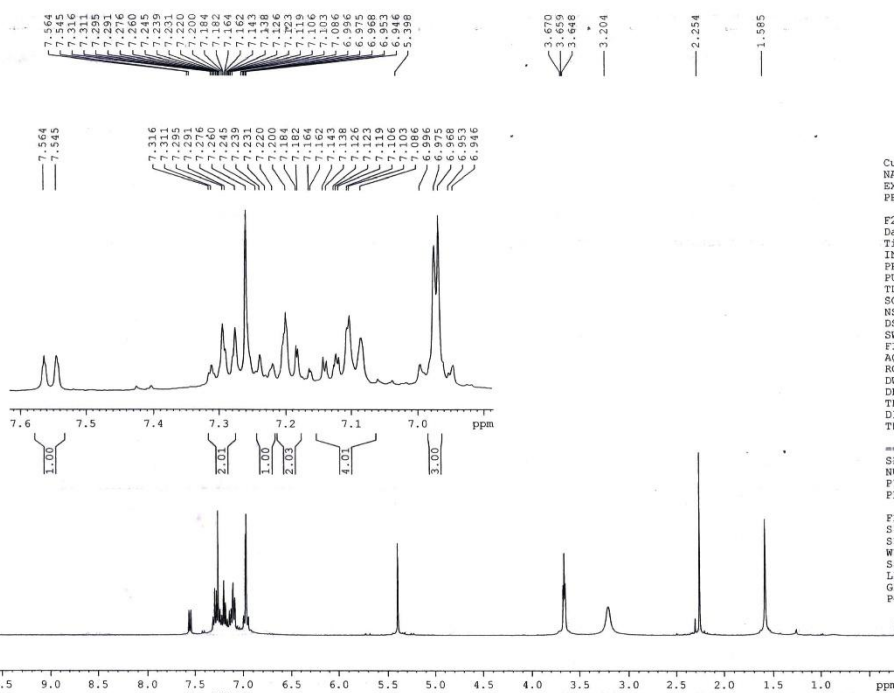
----- CHANNEL f1 -----
SF01 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

----- CHANNEL f2 -----
SF02 400.1516006 MHz
NUC2 1H
CPDPRG12 waltz16
FCDPR 90.000 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

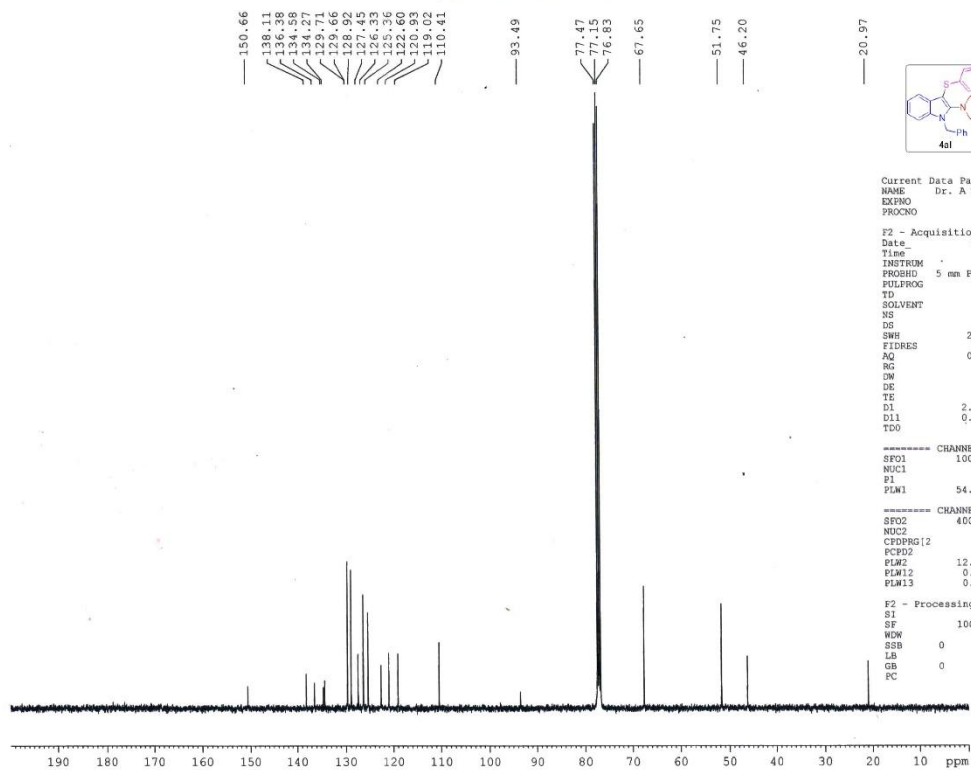
F2 - Processing parameters
SI 16384
SF 100.6177946 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

¹H of VBSP-1

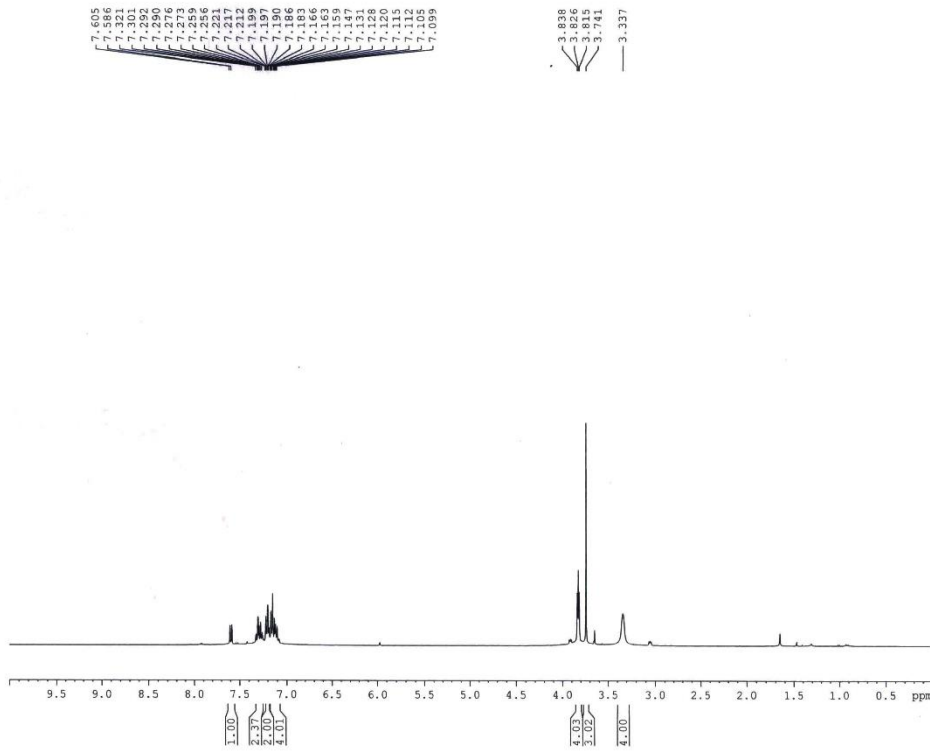
¹H NMR: 400 MHz; Solvent: CDCl₃



¹³C NMR: 100 MHz; Solvent: CDCl₃



¹H NMR: 400 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE 2023
EXFNO 494
PROCNO 1

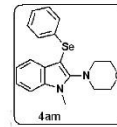
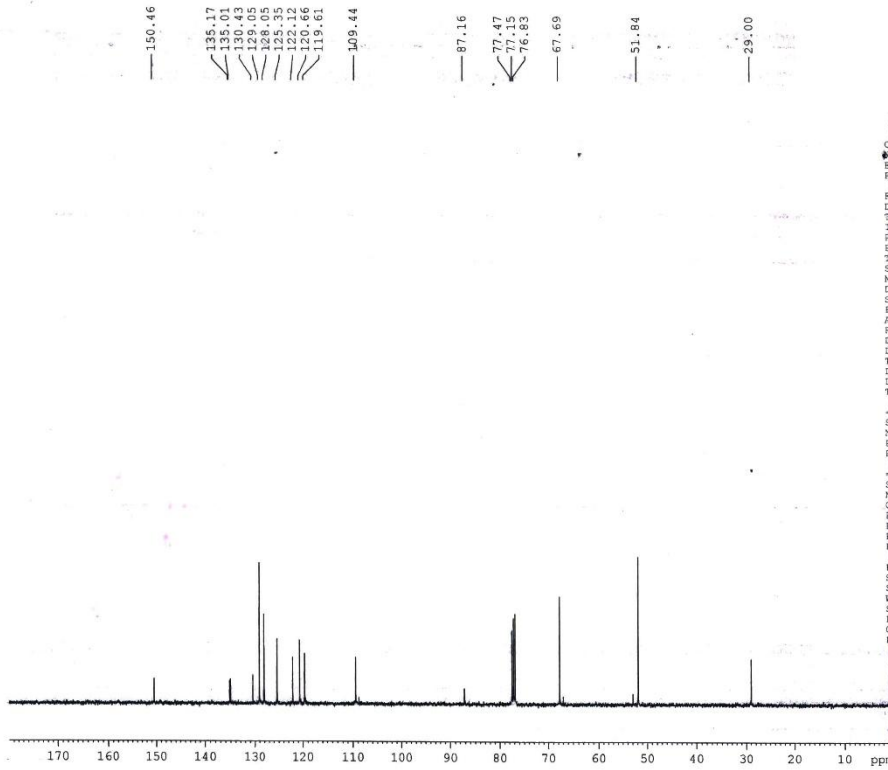
F2 - Acquisition Parameters:
Date 20230718
Time 14.33
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl₃
NS 8
DS 1
SWH 8223.685 Hz
FIDRES 0.250967 Hz
AQ 1.5922944 sec
RG 54.07
DW 60.800 usec
DE 6.50 usec
TE 299.4 K
D1 1.00000000 sec
TDO 1

----- CHANNEL f1 -----
SF01 400.1524711 MHz
NUC1 1H
P1 14.75 usec
PLW1 12.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1500097 MHz
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 2.00

13C of VBSP-283/8

¹³C NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE 2023
EXFNO 496
PROCNO 1

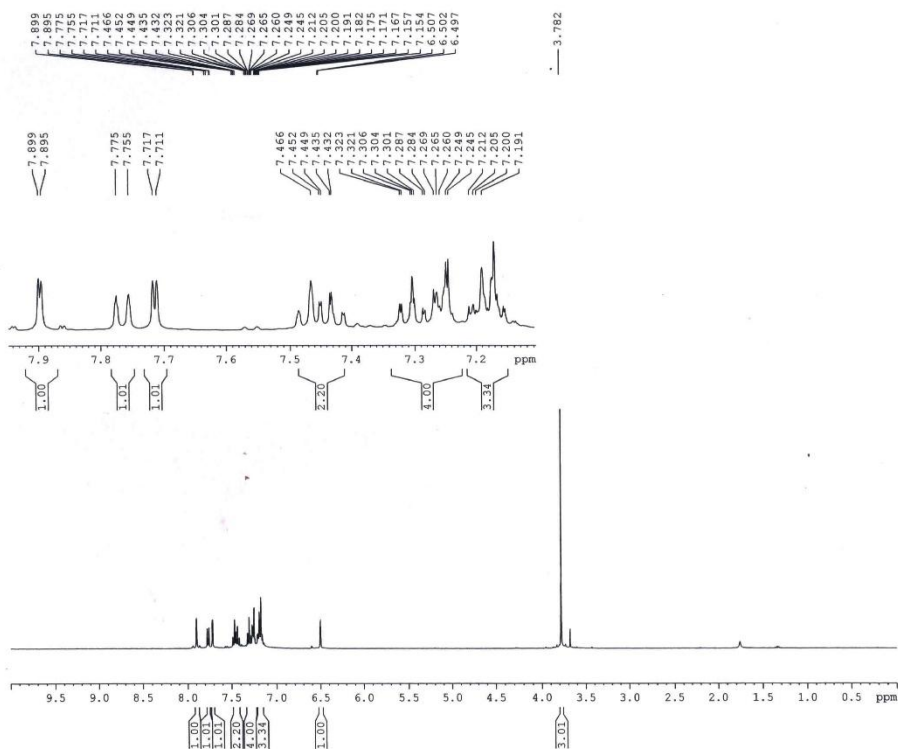
F2 - Acquisition Parameters:
Date 20230718
Time 14.56
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl₃
NS 120
DS 2
SWH 24039.461 Hz
FIDRES 0.733596 Hz
AQ 0.6915744 sec
RG 54.07
DW 20.800 usec
DE 6.50 usec
TE 299.9 K
D1 2.00000000 sec
D11 0.03000000 sec
TDO 1

----- CHANNEL f1 -----
SF01 100.6278588 MHz
NUC1 13C
P1 8.50 usec
PLW1 54.00000000 W

----- CHANNEL f2 -----
SF02 400.1516006 MHz
NUC2 1H
CDEPRG2 waltz16
PCPD2 80.00 usec
PLW2 12.00000000 W
PLW12 0.32210000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177932 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

¹H NMR: 400 MHz; Solvent: CDCl₃



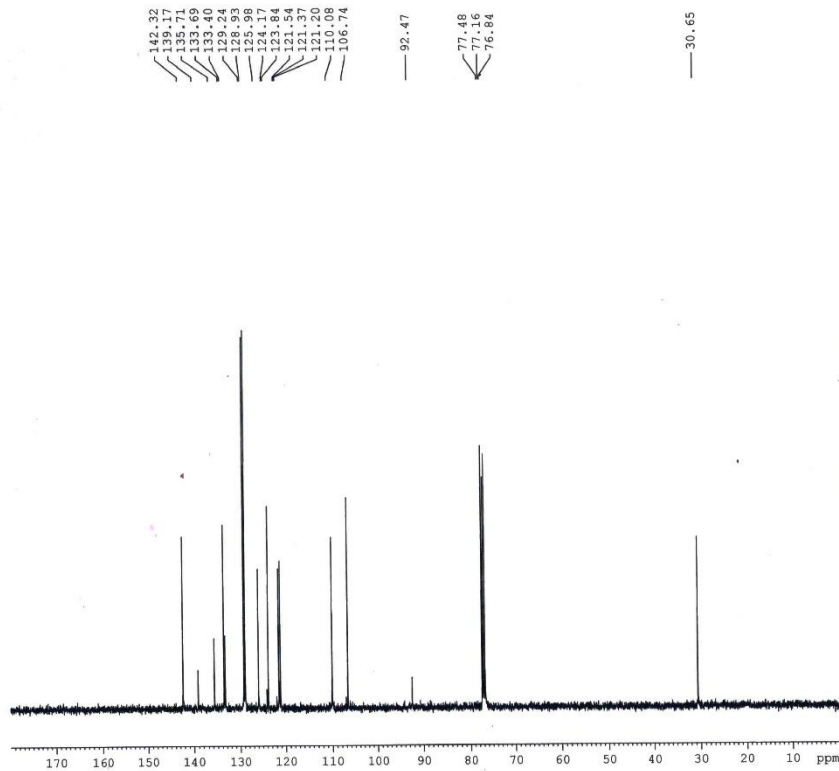
Current Data Parameters
NAME Dr. A MAJEE 2023
EXPRO 761
PROCNO 1

F2 - Acquisition Parameters
Date 20231102
Time 17.23
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 8
DS 1
SWH 8223.685 Hz
FIDRES 0.250967 Hz
AQ 1.9922944 sec
RG 93.46
DW 60.800 usec
DE 6.50 usec
TE 293.6 K
D1 1.00000000 sec
TDO 1

===== CHANNEL f1 =====
SFO1 400.1524711 MHz
NUC1 1H
P1 14.75 usec
PLW1 12.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1499916 MHz
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 2.00

¹³C [H] NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE 2023
EXPRO 762
PROCNO 1

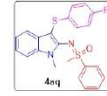
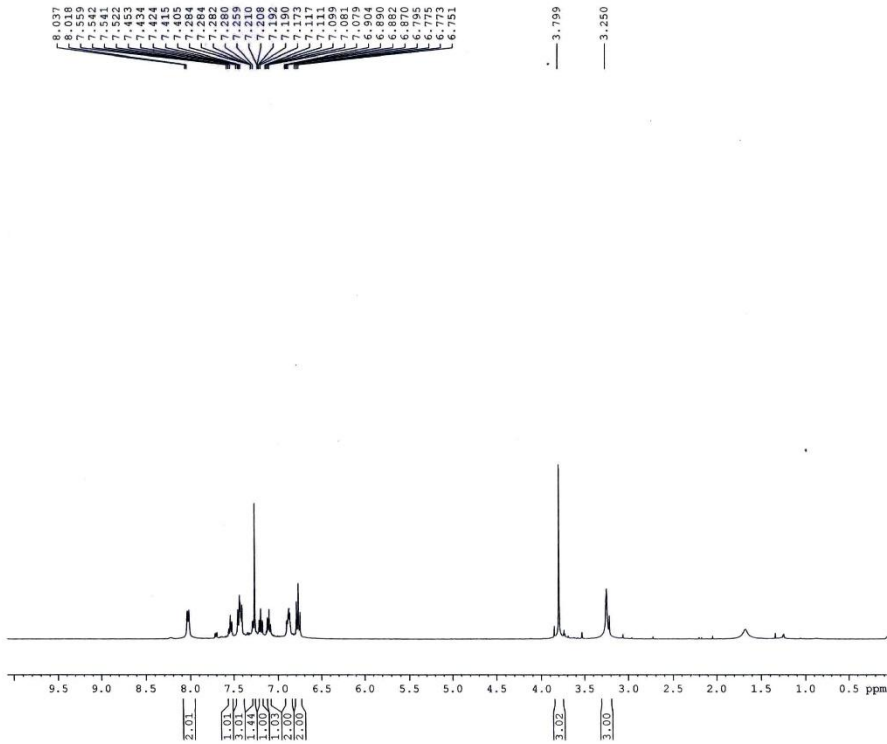
F2 - Acquisition Parameters
Date 20231102
Time 17.39
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 256
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 93.46
DW 20.800 usec
DE 6.50 usec
TE 293.7 K
D1 2.00000000 sec
D11 0.03000000 sec
TDO 1

===== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG2 waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177916 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

¹H NMR: 400 MHz; Solvent: CDCl₃



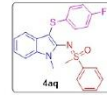
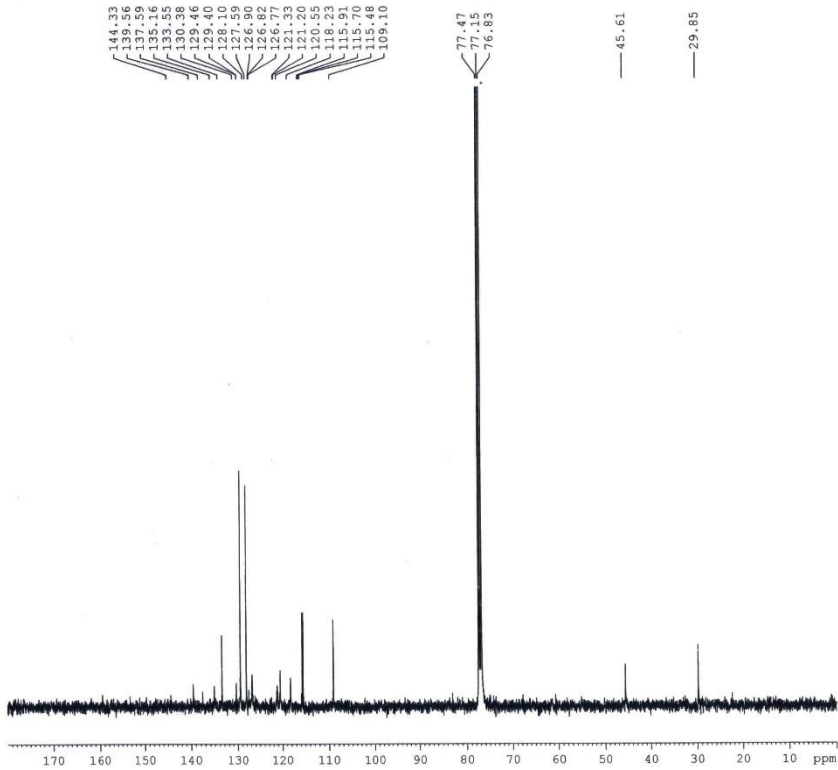
Current Data Parameters
NAME Dr. A MAJEE 2024
EXPNO 38
PROCNO 1

F2 - Acquisition Parameters
Date_ 20240203
Time 20.59
INSTRUM spect
PROBHD 5 mm PABBO BB/
FULPROG zg30
TD 32768
SOLVENT CDCl₃
NS 8
DS 1
SWH 8223.685 Hz
FIDRES 0.250967 Hz
AQ 1.9922944 sec
RG 135.7
DW 60.800 usec
DE 6.50 usec
TE 291.0 K
D1 1.00000000 sec
TDO 1

----- CHANNEL f1 -----
SF01 400.1524711 MHz
NUC1 1H
F1 14.75 usec
PLW1 12.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1500097 MHz
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 2.00

¹³C{¹H} NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE 2024
EXPNO 40
PROCNO 1

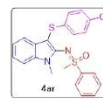
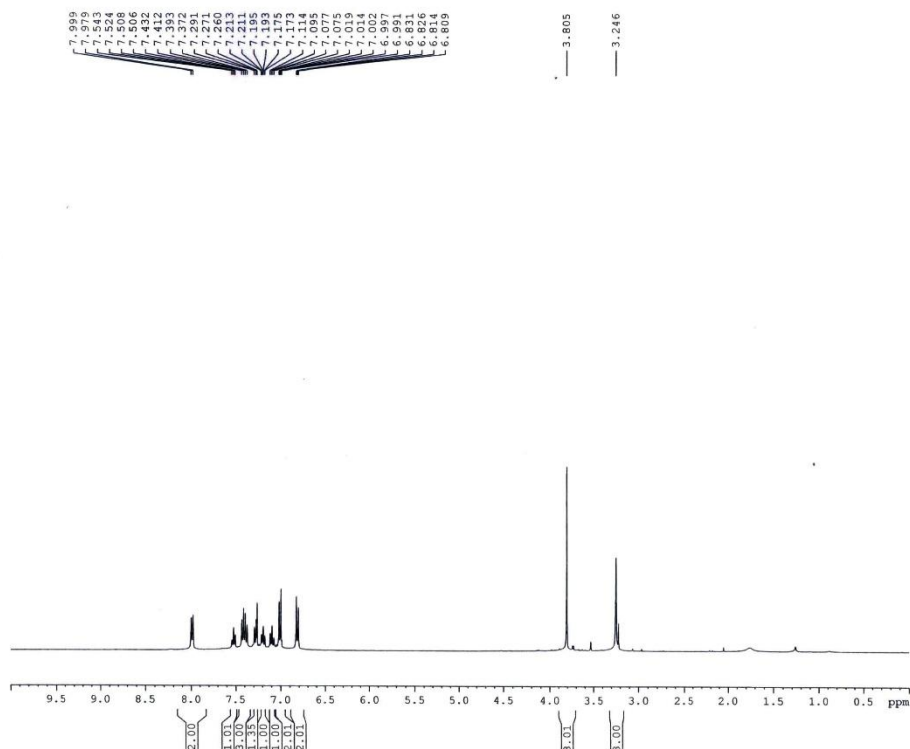
F2 - Acquisition Parameters
Date_ 20240203
Time 22.19
INSTRUM spect
PROBHD 5 mm PABBO BB/
FULPROG zgpg30
TD 32768
SOLVENT CDCl₃
NS 1500
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 135.7
DW 20.800 usec
DE 6.50 usec
TE 292.2 K
D1 2.00000000 sec
D11 0.03000000 sec
TDO 1

----- CHANNEL f1 -----
SF01 100.6278588 MHz
NUC1 13C
F1 8.90 usec
PLW1 54.00000000 W

----- CHANNEL f2 -----
SF02 400.1516006 MHz
NUC2 1H
CPRPRG12 waltz16
PCD2 90.00 usec
FLM2 12.00000000 W
FLM12 0.32231000 W
FLM13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6178783 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

¹H NMR: 400 MHz; Solvent: CDCl₃



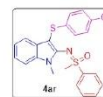
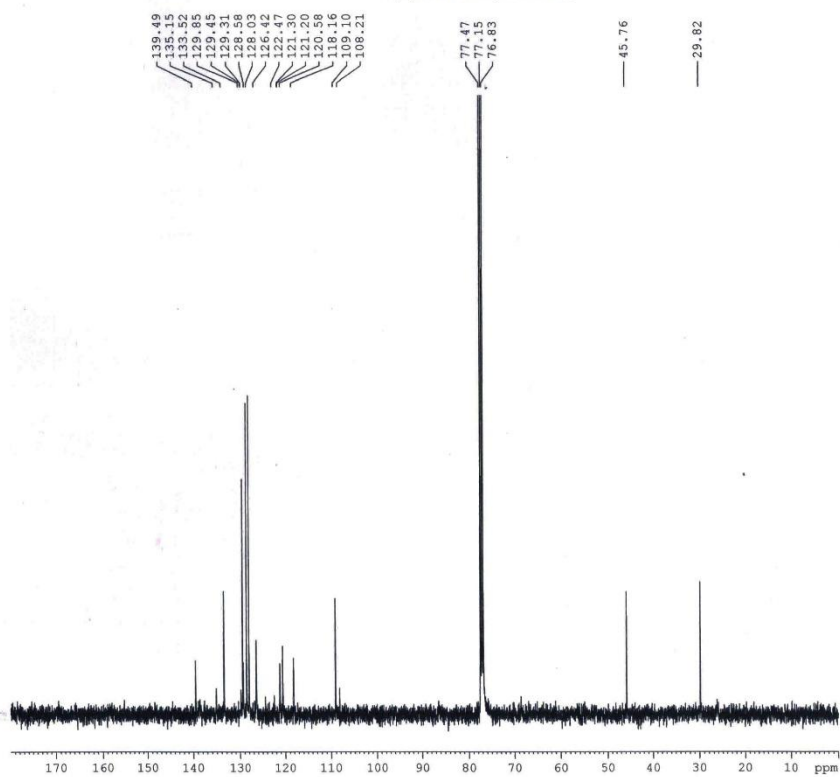
Current Data Parameters
NAME Dr. A MAJEE 2024
EXPNO 44
PROCNO 1

F2 - Acquisition Parameters
Date 20240204
Time 18.57
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl₃
NS 8
DS 1
SWH 8223.685 Hz
FIDRES 0.250967 Hz
AQ 1.9922944 sec
RG 106.66
DW 60.800 usec
DE 6.50 usec
TE 290.8 K
D1 1.00000000 sec
TDO 1

==== CHANNEL f1 =====
SF01 400.1524711 MHz
NUC1 ¹H
P1 14.75 usec
PLW1 12.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1500092 MHz
WDW EM
SSB 0
LB 0 0.30 Hz
GB 0
PC 2.00

¹³C NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE 2024
EXPNO 45
PROCNO 1

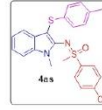
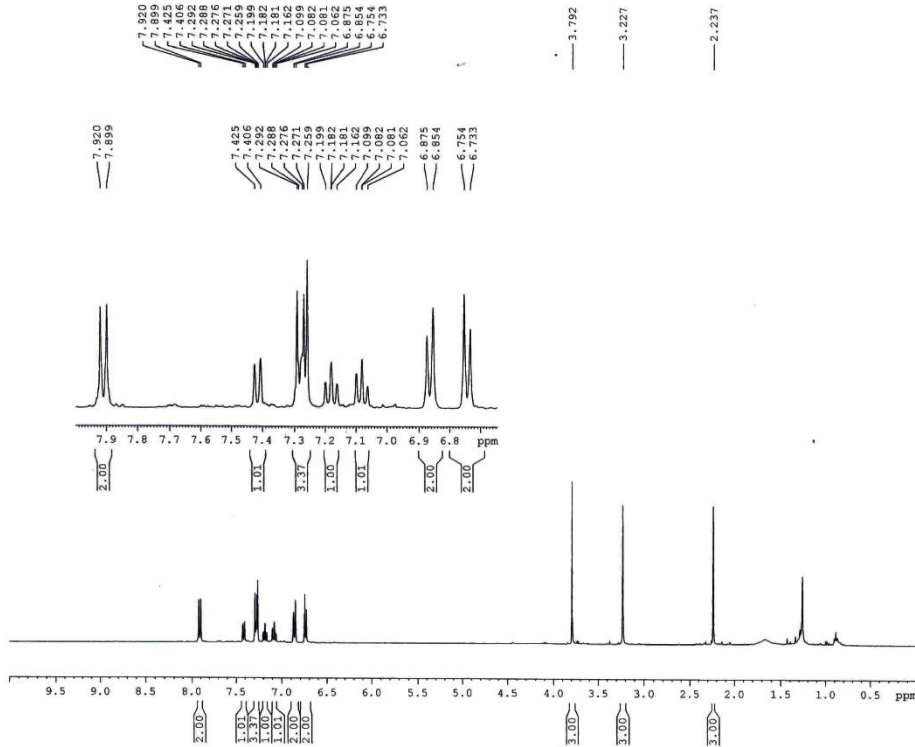
F2 - Acquisition Parameters
Date 20240204
Time 19.18
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl₃
NS 400
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 106.66
DW 20.800 usec
DE 6.50 usec
TE 292.1 K
D1 2.00000000 sec
D11 0.03000000 sec
TDO 1

==== CHANNEL f1 =====
SF01 100.6278568 MHz
NUC1 ¹³C
P1 8.90 usec
PLW1 54.00000000 W

==== CHANNEL f2 =====
SF02 400.1516006 MHz
NUC2 ¹H
CPCPRG12 waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLM12 0.32231000 W
PLM13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177887 MHz
WDW DM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

¹H NMR: 400 MHz; Solvent: CDCl₃



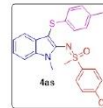
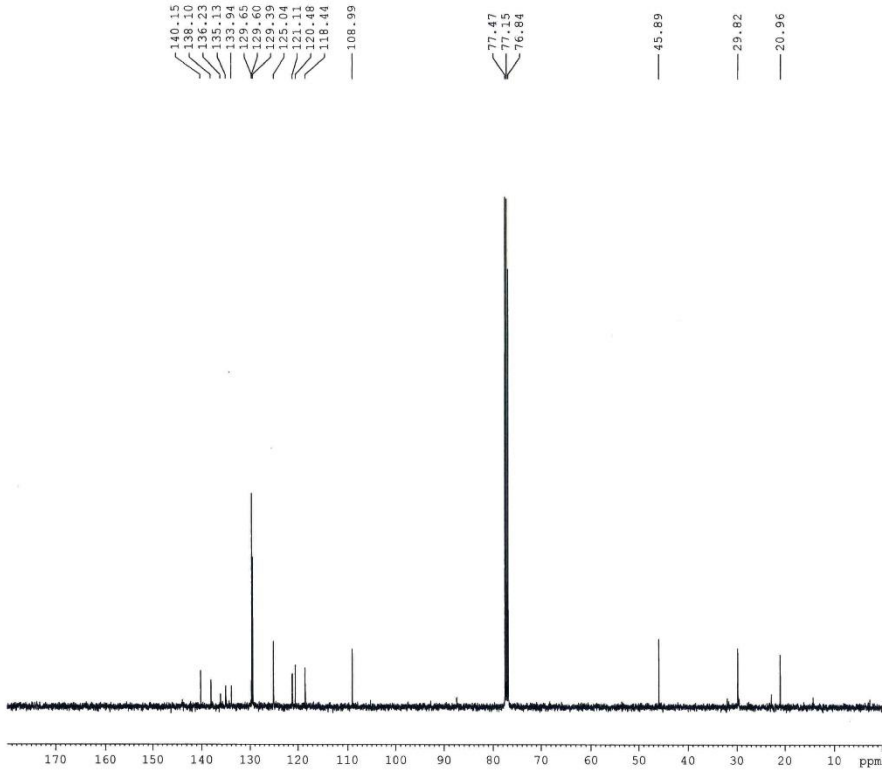
Current Data Parameters
NAME Dr. A MAJEE 2023
EXNO 506
PROCNO 1

F2 - Acquisition Parameters
Date_ 20230719
Time 17.18
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 8
DS 1
SWH 8223.685 Hz
FIDRES 0.250967 Hz
AQ 1.992944 sec
RG 93.46
DW 60.800 usec
DE 6.50 usec
TE 297.3 K
D1 1.0000000 sec
TDO 1

CHANNEL f1
SFO1 400.1524711 MHz
NUC1 ¹H
P1 14.75 usec
PLW1 12.0000000 W

F2 - Processing parameters
SI 16384
SF 400.1500097 MHz
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 2.00

¹³C{¹H} NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE 2023
EXNO 507
PROCNO 1

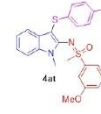
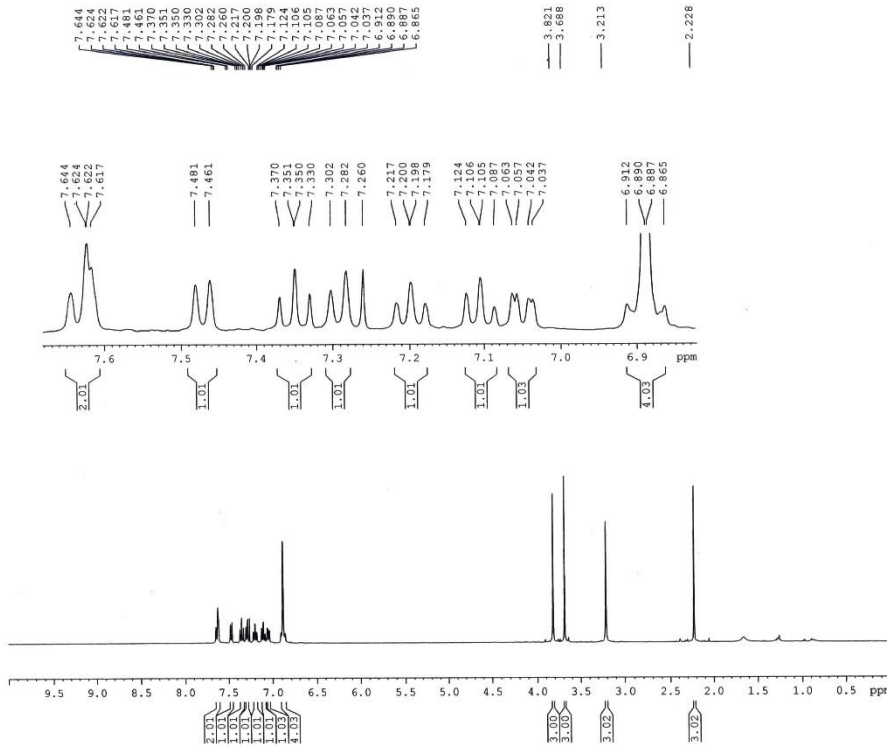
F2 - Acquisition Parameter
Date_ 20230719
Time 17.53
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 512
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 93.46
DW 20.800 usec
DE 6.50 usec
TE 297.7 K
D1 2.0000000 sec
D11 0.03000000 sec
TDO 1

CHANNEL f1
SFO1 100.6278588 MHz
NUC1 ¹³C
P1 8.90 usec
PLW1 54.0000000 W

CHANNEL f2
SFO2 400.1516006 MHz
NUC2 ¹H
CPCPD2 waltz16
PCPD2 90.00 usec
PLW2 12.0000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177858 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

¹H NMR: 400 MHz; Solvent: CDCl₃



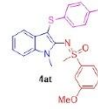
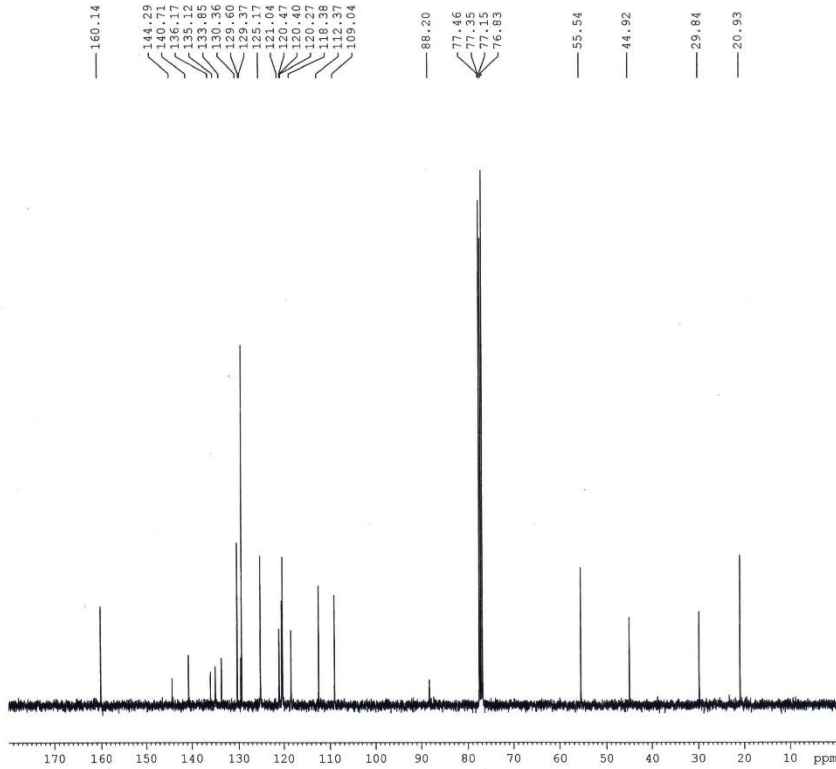
Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 874
PROCNO 1

F2 - Acquisition Parameters
Date_ 20231203
Time 20.06
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 8
DS 1
SWH 8223.685 Hz
FIDRES 0.250967 Hz
AQ 1.9922944 sec
RG 93.46
DW 60.800 usec
DE 6.50 usec
TE 293.3 K
D1 1.00000000 sec
TDO 1

CHANNEL f1
SF01 400.1524711 MHz
NUC1 1H
P1 14.75 usec
PLW1 12.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1500092 MHz
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 2.00

¹³C NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 875
PROCNO 1

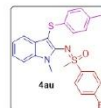
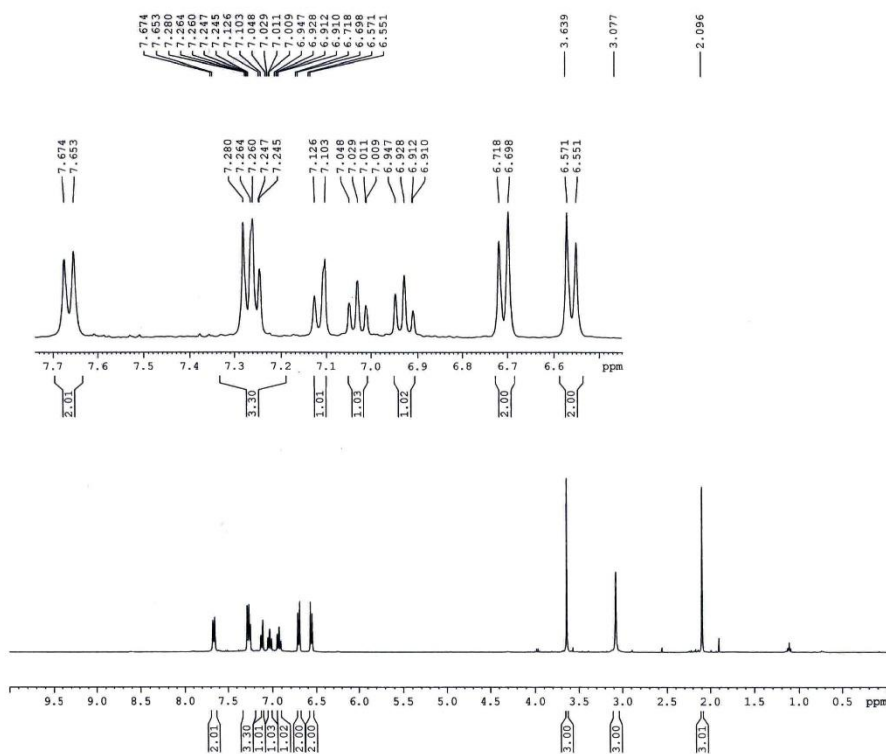
F2 - Acquisition Parameters
Date_ 20231203
Time 20.25
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 320
DS 2
SWH 24938.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 93.46
DW 20.800 usec
DE 6.50 usec
TE 292.9 K
D1 2.00000000 sec
D11 0.03000000 sec
TDO 1

CHANNEL f1
SF01 100.6278588 MHz
NUC1 13C
P1 8.30 usec
PLW1 54.00000000 W

CHANNEL f2
SF02 400.1516006 MHz
NUC2 1H
CFPRG(2) waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177902 MHz
WDW DM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

¹H NMR: 400 MHz; Solvent: CDCl₃



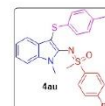
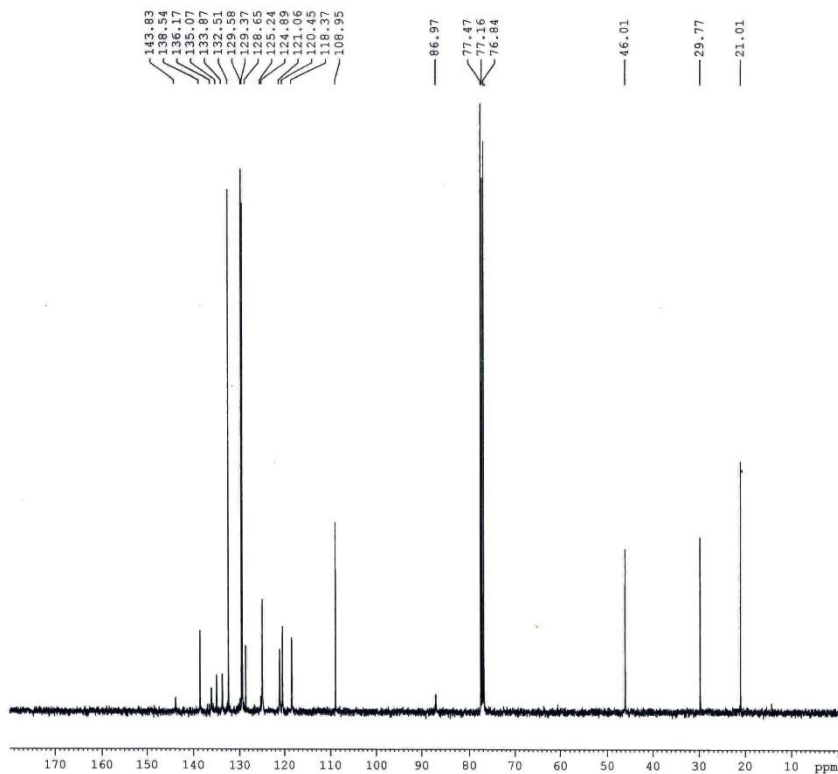
Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 882
PROCNO 1

F2 - Acquisition Parameters
Date_ 20231204
Time 13.08
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 8
DS 1
SWH 8223.685 Hz
FIDRES 0.250967 Hz
AQ 1.9922944 sec
RG 93.46
DW 60.800 usec
DE 6.50 usec
TE 291.4 K
D1 1.00000000 sec
TDO 1

CHANNEL f1
SFO1 400.1524711 MHz
NUC1 1H
P1 14.75 usec
PLW1 12.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1500724 MHz
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 2.00

¹³C{¹H} NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 883
PROCNO 1

F2 - Acquisition Parameters
Date_ 20231204
Time 13.45
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 720
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 93.46
DW 20.800 usec
DE 6.50 usec
TE 292.0 K
D1 2.00000000 sec
D11 0.03000000 sec
TDO 1

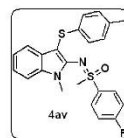
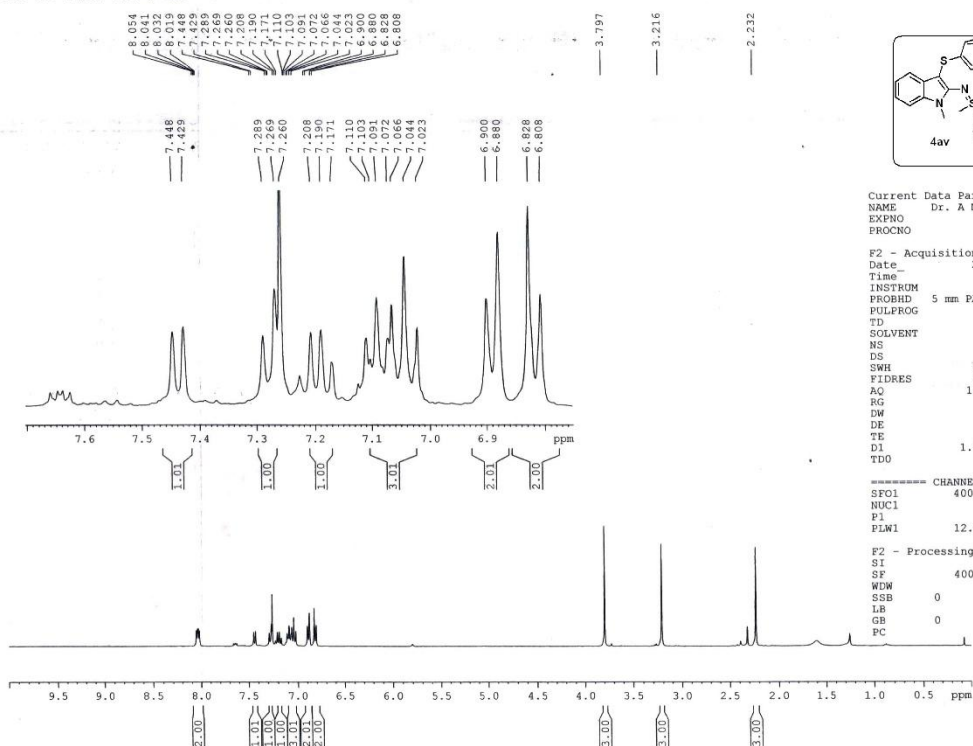
CHANNEL f1
SFO1 100.6278588 MHz
NUC1 13C
P1 8.50 usec
PLW1 54.00000000 W

CHANNEL f2
SFO2 400.1516006 MHz
NUC2 1H
CPDPRG2 waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.617917 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

¹H of VBSP-RR-8 UP

¹H NMR: 400 MHz; Solvent: CDCl₃



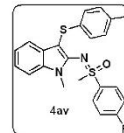
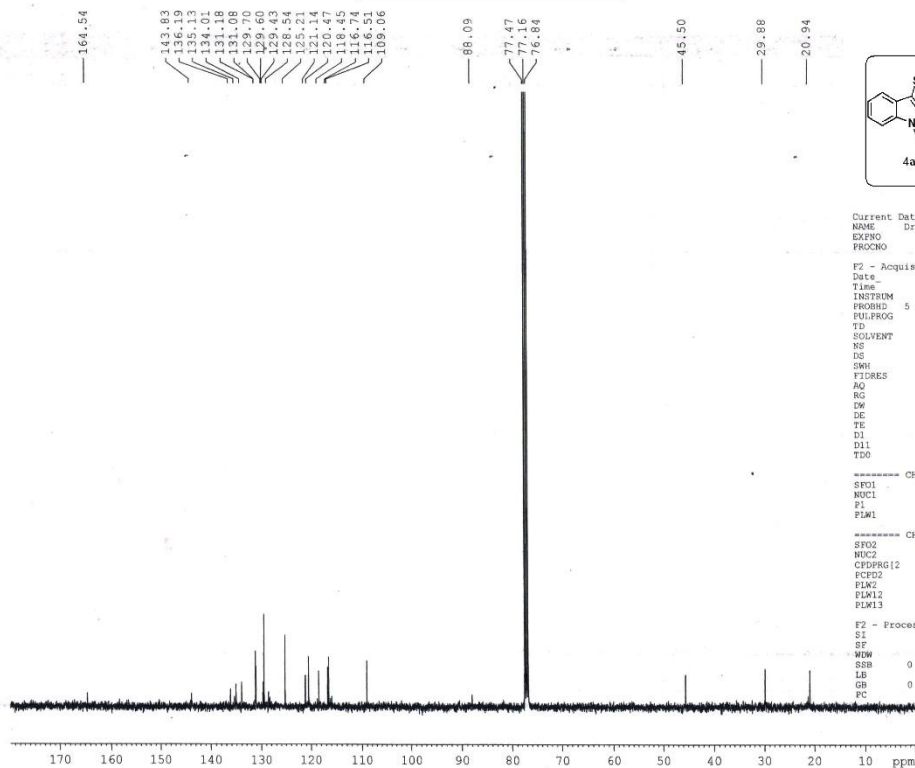
Current Data Parameters
NAME Dr. A MAJEE 2026-1H
EXPNO 123
PROCNO 1
F2 - Acquisition Parameters
Date 20260412
Time 11.00
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl₃
NS 8
DS 1
SWH 9223.695 Hz
FIDRES 0.250967 Hz
AQ 1.9922944 sec
RG 135.7
DW 60.800 usec
DE 6.50 usec
TE 293.1 K
D1 1.00000000 sec
TDO 1

CHANNEL f1
SF01 400.1524711 MHz
NUC1 1H
P1 14.75 usec
PLW1 12.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1500097 MHz
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 2.00

¹³C of VBSP-RR-8 UP

¹³C{¹H} NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE-2026-13C
EXPNO 81
PROCNO 1

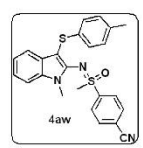
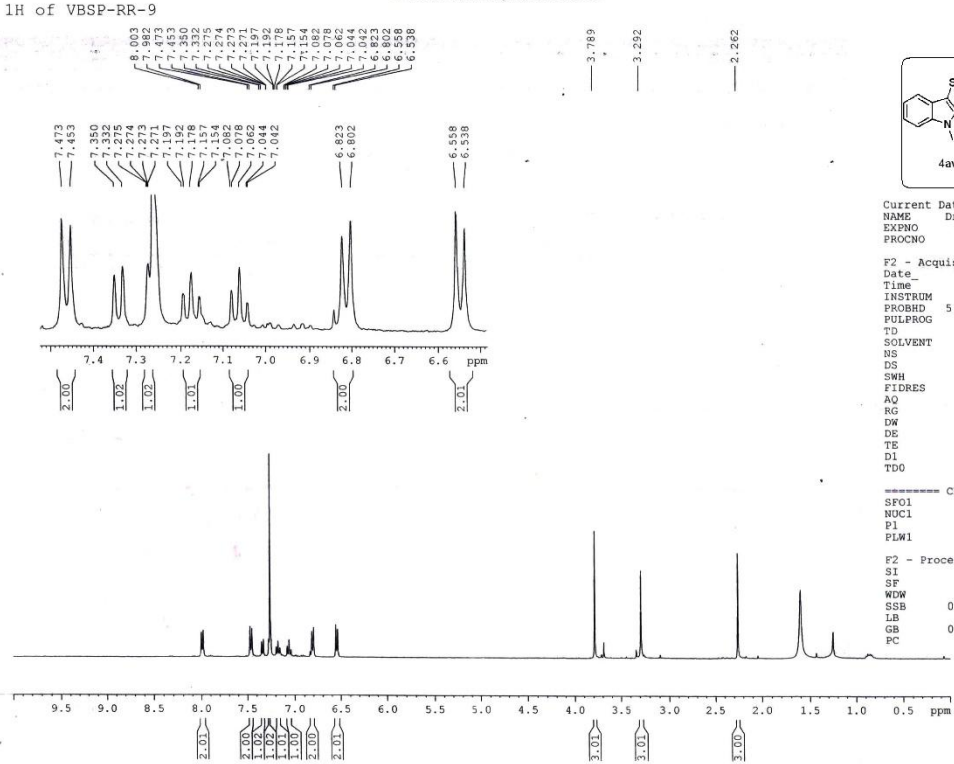
F2 - Acquisition Parameters
Date 20260412
Time 11.50
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl₃
NS 90
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 135.7
DW 20.800 usec
DE 6.50 usec
TE 292.5 K
D1 2.00000000 sec
D11 0.03000000 sec
TDO 1

CHANNEL f1
SF01 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

CHANNEL f2
SF02 400.1516006 MHz
NUC2 1H
CPDPRG12 waltz16
PCPD 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16215000 W

F2 - Processing parameters
SI 16384
SF 100.6177859 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40

¹H NMR: 400 MHz; Solvent: CDCl₃



Current Data Parameters
 NAME Dr. A MAJEE 2026-1H
 EXPNO 130
 PROCNO 1

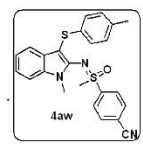
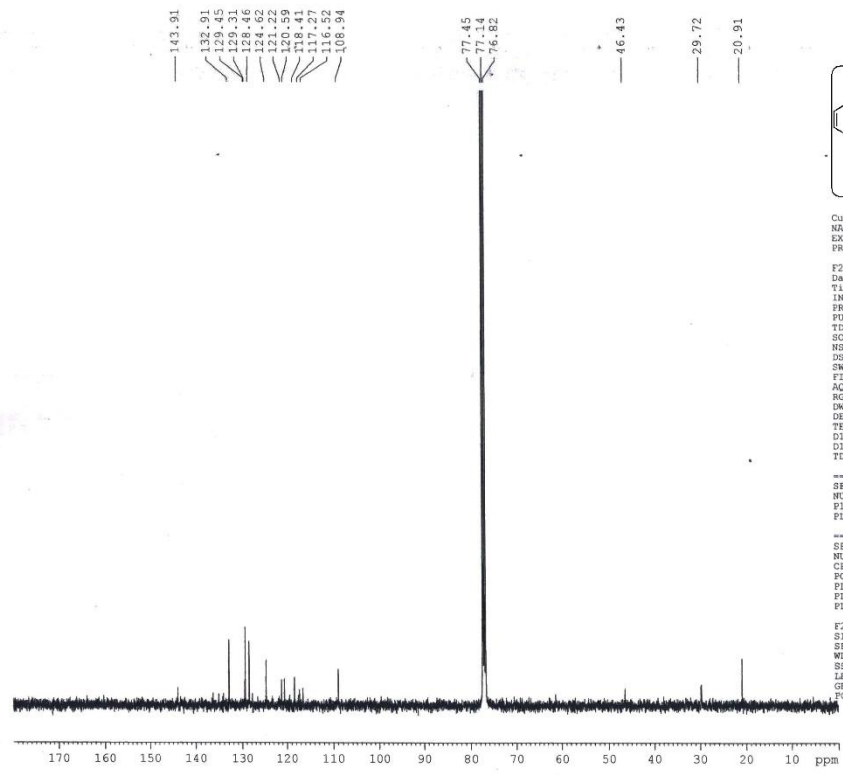
F2 - Acquisition Parameters
 Date 20260413
 Time 15.22
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zg30
 TD 32768
 SOLVENT CDCl3
 NS 8
 DS 1
 SWH 8223.685 Hz
 FIDRES 0.250967 Hz
 AQ 1.992944 sec
 RG 186.42
 DW 60.800 usec
 DE 6.50 usec
 TE 292.8 K
 D1 1.00000000 sec
 TDO 1

==== CHANNEL f1 =====
 SF01 400.1524711 MHz
 NUC1 1H
 P1 14.75 usec
 PLW1 12.00000000 W

F2 - Processing parameters
 SI 16384
 SF 400.1500097 MHz
 WDW EM
 SSB 0
 LB 0.30 Hz
 GB 0
 PC 2.00

¹³C NMR of VBSP-RR-9 rpt

¹³C[¹H] NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
 NAME Dr. A MAJEE-2026-13C
 EXPNO 91
 PROCNO 1

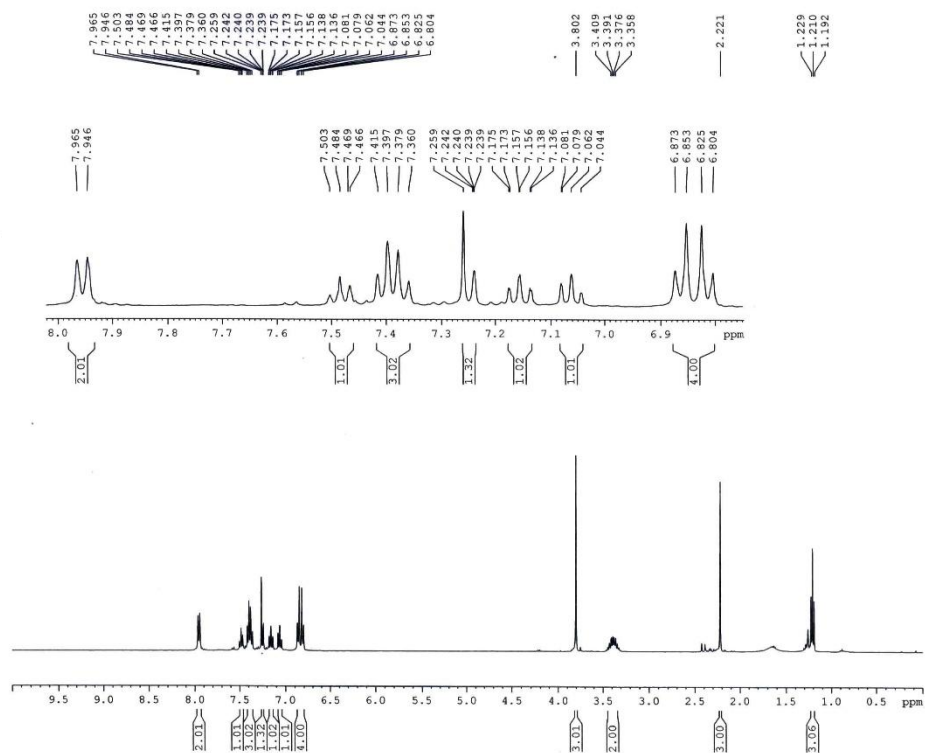
F2 - Acquisition Parameters
 Date 20260413
 Time 23.52
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 32768
 SOLVENT CDCl3
 NS 2400
 DS 2
 SWH 24039.461 Hz
 FIDRES 0.733596 Hz
 AQ 0.6815744 sec
 RG 186.42
 DW 20.800 usec
 DE 6.50 usec
 TE 296.1 K
 D1 2.00000000 sec
 D11 0.03000000 sec
 TDO 1

==== CHANNEL f1 =====
 SF01 100.6278588 MHz
 NUC1 13C
 P1 8.90 usec
 PLW1 54.00000000 W

==== CHANNEL f2 =====
 SF02 400.1516006 MHz
 NUC2 1H
 CPDPRG(2) waltz16
 PCPD 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.32231000 W
 PLW13 0.16215000 W

F2 - Processing parameters
 SI 16384
 SF 100.6177868 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40

¹H NMR: 400 MHz; Solvent: CDCl₃



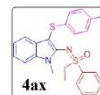
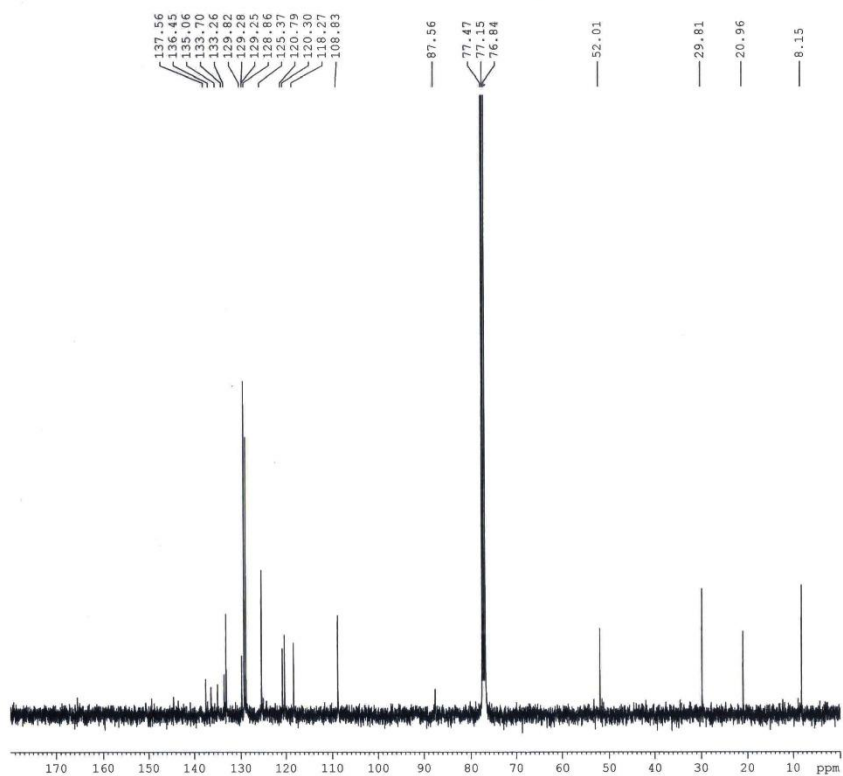
Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 751
PROCNO 1

F2 - Acquisition Parameters
Date 20231018
Time 14.36
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl₃
NS 8
DS 1
SWH 8223.685 Hz
FIDRES 0.250967 Hz
AQ 1.9922944 sec
RG 135.7
DW 60.800 usec
DE 6.50 usec
TE 297.3 K
D1 1.00000000 sec
TDO 1

===== CHANNEL f1 =====
SF01 400.1524711 MHz
NUC1 1H
P1 14.75 usec
PLW1 12.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1500097 MHz
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 2.00

¹³C{¹H} NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE 2023
EXPNO 752
PROCNO 1

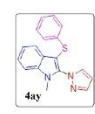
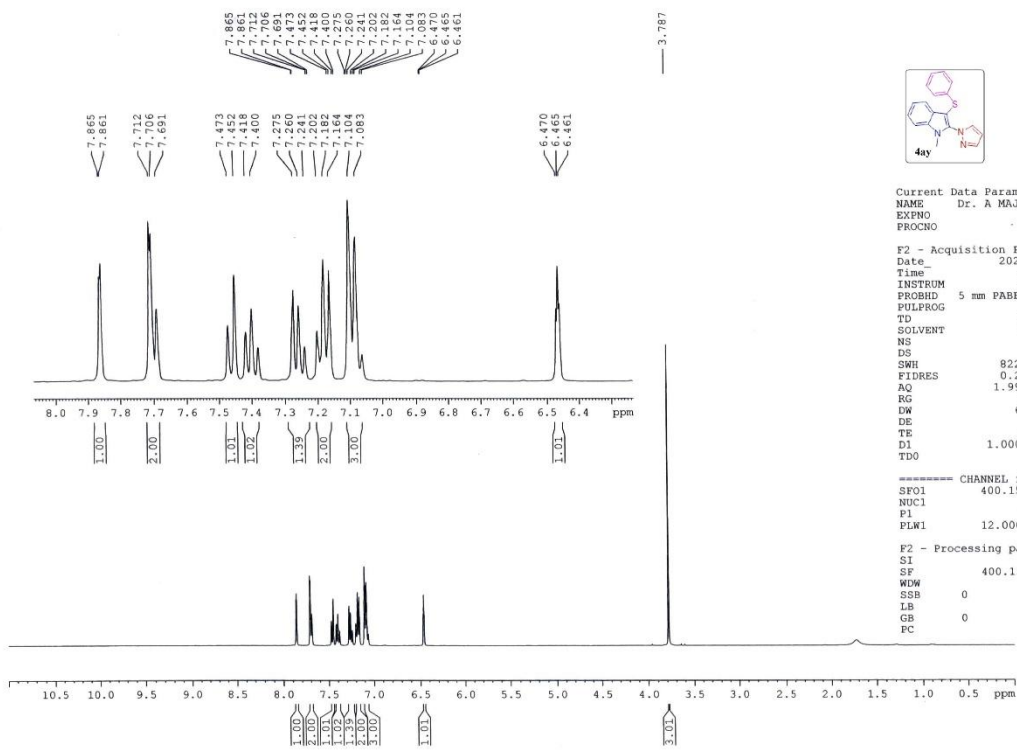
F2 - Acquisition Parameters
Date 20231018
Time 15.09
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl₃
NS 2
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 135.7
DW 20.800 usec
DE 6.50 usec
TE 297.4 K
D1 2.00000000 sec
D11 0.03000000 sec
TDO 1

===== CHANNEL f1 =====
SF01 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

===== CHANNEL f2 =====
SF02 400.1516006 MHz
NUC2 1H
CPDPRG2 waltz16
PCPD2 90.00 usec
PLW2 12.00000000 W
PLW12 0.32231000 W
PLW13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177858 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.00

¹H NMR: 400 MHz; Solvent: CDCl₃



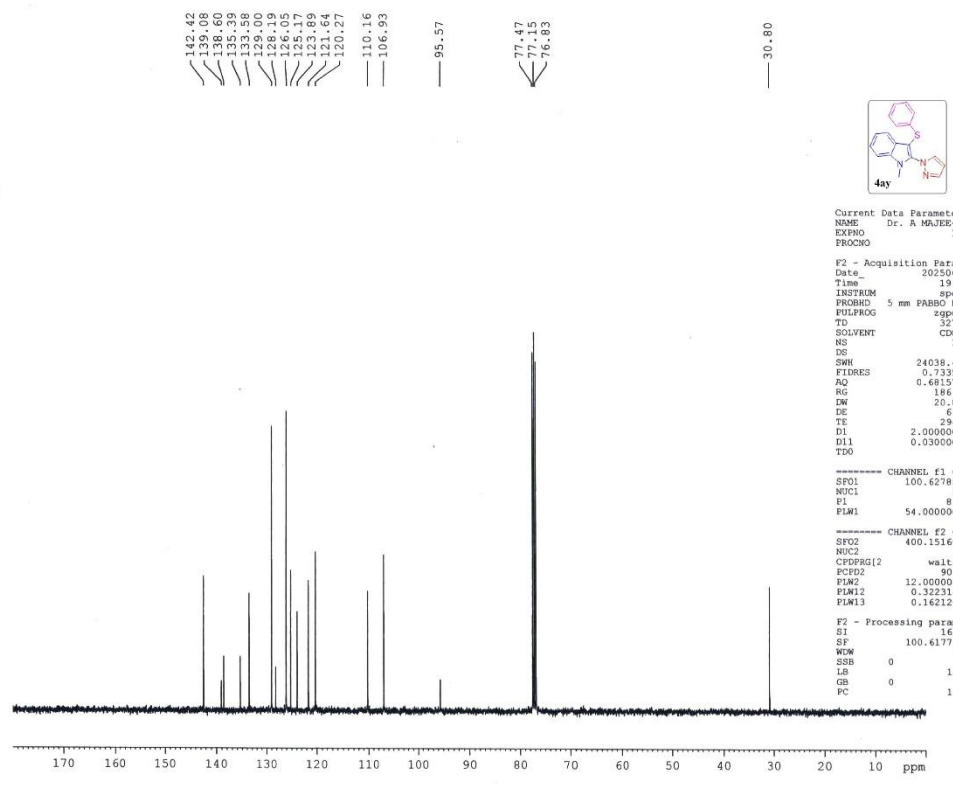
Current Data Parameters
NAME Dr. A MAJEE 2025-1H
EXPNO 259
PROCNO 1

F2 - Acquisition Parameters
Date 20250620
Time 16.18
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zg30
TD 32768
SOLVENT CDCl3
NS 8
DS 1
SWH 8223.685 Hz
FIDRES 0.250967 Hz
AQ 1.992944 sec
RG 77.59
DW 60.800 usec
DE 6.50 usec
TE 293.5 K
D1 1.0000000 sec
TDO 1

==== CHANNEL f1 =====
SFO1 400.1524711 MHz
NUC1 1H
P1 14.75 usec
PLW1 12.00000000 W

F2 - Processing parameters
SI 16384
SF 400.1500037 MHz
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 2.00

¹³C NMR: 100 MHz; Solvent: CDCl₃



Current Data Parameters
NAME Dr. A MAJEE-2025-13C
EXPNO 179
PROCNO 1

F2 - Acquisition Parameters
Date 20250620
Time 19.03
INSTRUM spect
PROBHD 5 mm PABBO BB/
PULPROG zgpg30
TD 32768
SOLVENT CDCl3
NS 300
DS 2
SWH 24038.461 Hz
FIDRES 0.733596 Hz
AQ 0.6815744 sec
RG 186.42
DW 20.800 usec
DE 6.50 usec
TE 294.5 K
D1 2.0000000 sec
D11 0.0500000 sec
TDO 1

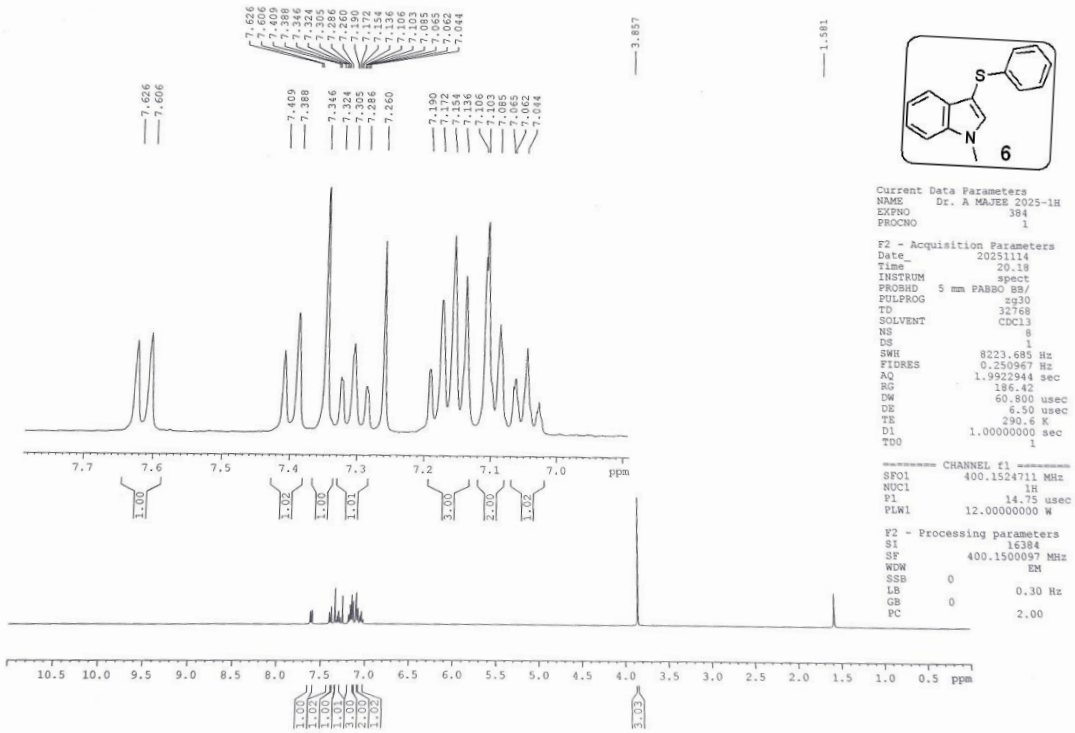
==== CHANNEL f1 =====
SFO1 100.6278588 MHz
NUC1 13C
P1 8.90 usec
PLW1 54.00000000 W

==== CHANNEL f2 =====
SFO2 400.1516006 MHz
NUC2 1H
CPCPRG2 waitz16
RG2 90.00 usec
P1A2 12.00000000 W
P1A2 0.32231000 W
P1A13 0.16212000 W

F2 - Processing parameters
SI 16384
SF 100.6177867 MHz
WDW EM
SSB 0
LB 1.00 Hz
GB 0
PC 1.40

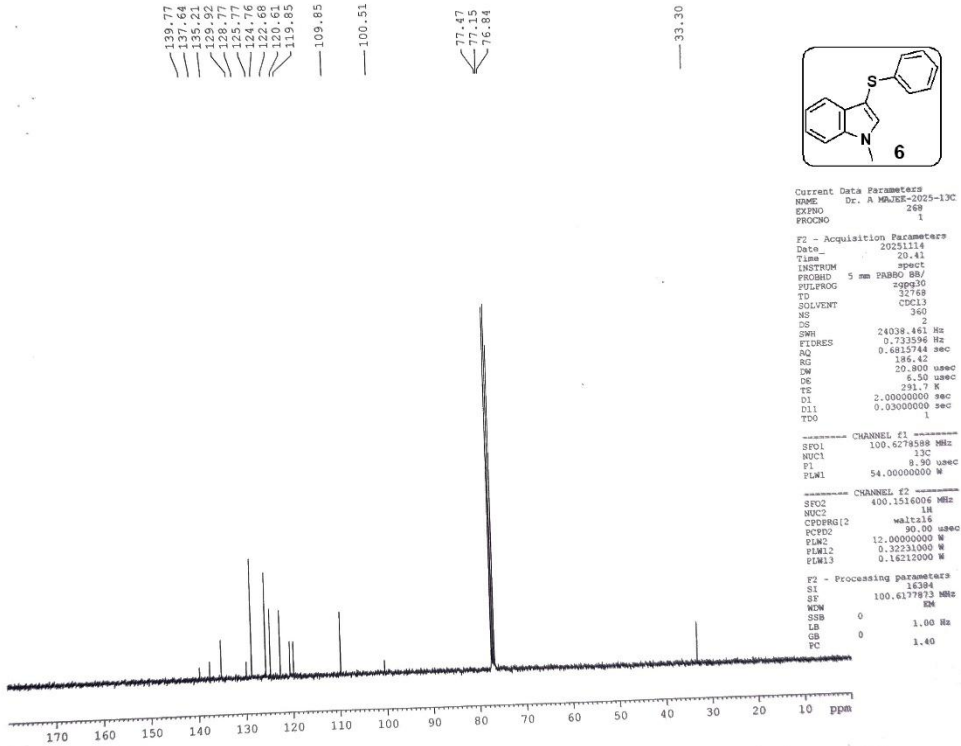
¹H of VBSPTP- Sph

¹H NMR: 400 MHz; Solvent: CDCl₃



¹³C of VBSPTP- Sph

¹³C{¹H} NMR: 100 MHz; Solvent: CDCl₃



Prof. A. Maiee
SP-21 9 (0.169) Cm (2:31)

Visva-Bharati University

Ph-Se-I

TOF MS ES+
306.8696 6.44e3

