

Supporting Information

**Organocatalytic Asymmetric Formal Oxidative Coupling for the
Construction of All-Aryl Quaternary Stereocenters**

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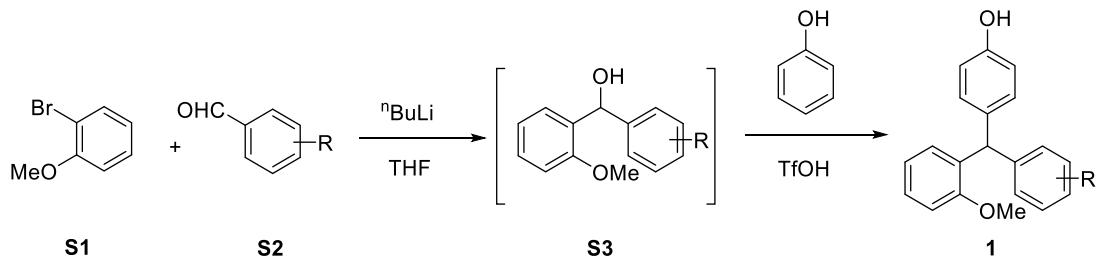
NMR Spectra and HPLC Traces

I. General Information

Flash column chromatography was performed over silica gel (200-300 mesh) purchased from Qindao Haiyang Co., China. All air or moisture sensitive reactions were conducted in oven-dried glassware under nitrogen atmosphere using anhydrous solvents. Anhydrous dichloromethane, toluene, diethyl ether, and tetrahydrofuran were purified by the Innovative® solvent purification system. Chemicals were purchased from commercial suppliers and used without further purification unless otherwise stated. ^1H , ^{13}C , ^{19}F NMR spectra were collected on a Bruker AV 400 MHz NMR spectrometer using residue solvent peaks as an internal standard (^1H NMR: CDCl_3 at 7.26 ppm, CD_2Cl_2 at 5.32 ppm, acetone- d_6 at 2.05 ppm; ^{13}C NMR: CDCl_3 at 77.23 ppm, CD_2Cl_2 at 53.84 ppm, acetone- d_6 at 29.84 ppm). Data for ^1H NMR are recorded as follows: chemical shift (δ , ppm), multiplicity (s = singlet; d = doublet; t = triplet; q = quarter; p = pentet; sept = septet; m = multiplet; br = broad), coupling constant (Hz), integration. Mass spectra were collected on an Agilent GC/MS 5975C system, a MALDI Micro MX mass spectrometer, or an API QSTAR XL System. IR spectra were recorded on Bruker TENSOR 27 spectrometer and reported in terms of frequency of absorption (cm^{-1}). Optical rotations were measured on JASCO P-2000 polarimeter with $[\alpha]_D$ values reported in degrees; concentration (c) is in 10 mg/mL. The enantiomeric excess values were determined by chiral HPLC using an Agilent 1200 LC instrument with Daicel CHIRALPAK® AD-H, IC, AS-H, or CHIRALCEL® OD-H columns.

II. Synthesis of the Triarylmethane Substrates

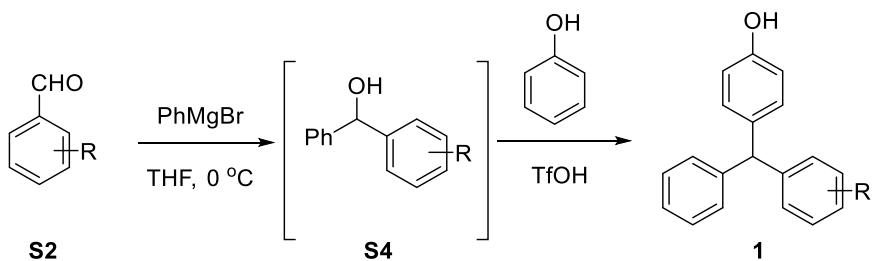
General Procedure A.



At -78°C , to a solution of 1-bromo-2-methoxybenzene **S1** (0.5 mL, 4.0 mmol) in THF (30 mL) was added dropwise n -butyllithium (2.5 M in pentane, 6.3 mL, 14.9 mmol). The mixture was kept stirring at the same temperature for 2 h. Then a solution of aldehyde **S2** (4.0 mmol) in THF (10 mL) was slowly added. The mixture was stirred at room temperature for 20 h before it was quenched with water (20 mL). The resulting mixture was extracted with DCM (30 mL \times 3). The combined organic layers were dried over anhydrous Na_2SO_4 , filtered, and concentrated. The crude mixture was purified by flash silica gel column chromatography (eluent: hexanes/EtOAc = 10:1) to give the diarylmethanol **S3**.

In a separate flask, to a solution of phenol (846 mg, 9.0 mmol,) and TfOH (0.05 mL, 0.57 mmol) in DCM (10 mL) was added a solution of the diarylmethanol **S3** in DCM (10 mL). The mixture was stirred for 10 min and concentrated. The residue was purified by silica gel flash column chromatography to afford the pure triarylmethane substrate **1**.

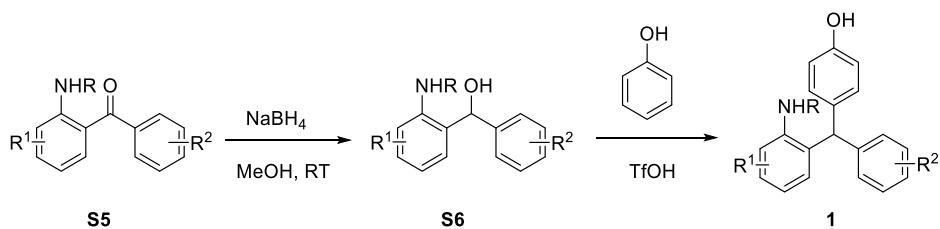
General Procedure B.



At 0°C , to a solution of **S2** (4.0 mmol) in THF (30 mL) was added PhMgBr (2.5 M in pentane, 6.3 mL, 14.9 mmol) dropwise. The mixture was kept stirring at the same temperature for 10 min. Next, the reaction was quenched with a saturated aqueous solution of NH₄Cl (20 mL). The mixture was extracted with DCM (30 mL \times 3). The combined organic layers were dried over anhydrous Na₂SO₄, filtered, and concentrated. The crude mixture was purified by flash silica gel column chromatography (eluent: hexanes/EtOAc = 10:1) to give alcohol **S4**.

In a separate flask, to a solution of phenol (846 mg, 9.0 mmol) and TfOH (0.05 mL, 0.57 mmol) in DCM (10 mL) was slowly added a solution of the alcohol **S4** in DCM (10 mL). The mixture was stirred for 10 min and concentrated. The residue was purified by silica gel flash column chromatography to afford the pure triarylmethane substrate **1**.

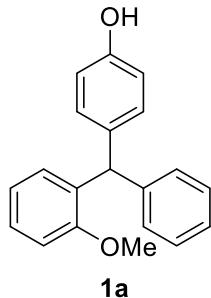
General Procedure C.



At 0°C , to a solution of **S5** (4.0 mmol) in MeOH (40 mL) was added NaBH₄ (908 mg, 24.0 mmol, 6 equiv) portionwise. The mixture was stirred at the same temperature for 1 h and concentrated. The residue was treated with a saturated

aqueous solution of NH₄Cl (30 mL), and extracted with DCM (40 mL × 3). The combined organic layers were dried over anhydrous Na₂SO₄, filtered, and concentrated to give the crude ketone **S6**.

In a separate flask, to a solution of phenol (846 mg, 9.0 mmol) and TfOH (0.05 mL, 0.57 mmol) in DCM (10 mL) was slowly added a solution of the crude ketone **S6** in DCM (10 mL). The mixture was stirred for 10 min and concentrated. The residue was purified by silica gel flash column chromatography to afford the pure triarylmethane substrate **1**.



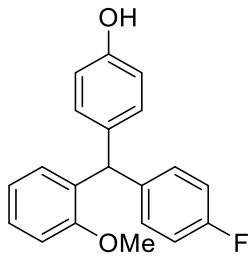
4-((2-Methoxyphenyl)(phenyl)methyl)phenol (1a) was prepared as a white solid from benzaldehyde (4.0 mmol) according to the General Procedure A (eluent: hexanes/EtOAc = 30:1 → 10:1) in 62% yield (over two steps)

¹H NMR (400 MHz, CDCl₃) δ 7.28 – 7.19 (m, 4H), 7.08 (d, *J* = 7.2 Hz, 2H), 6.95 (d, *J* = 8.4 Hz, 2H), 6.88 – 6.85 (m, 3H), 6.72 (d, *J* = 8.4 Hz, 2H), 5.86 (s, 1H), 3.72 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 157.2, 153.8, 144.3, 136.3, 133.0, 130.7, 130.4, 129.5, 128.2, 127.6, 126.1, 120.4, 115.1, 110.8, 55.7, 48.8.

IR (thin film) 3353, 3023, 2836, 1597, 1499, 1237, 1170, 748 cm⁻¹.

HRMS (CI+) Calcd for C₂₀H₁₈O₂ (M⁺): 290.1307, Found: 290.1308.



1b

4-((4-Fluorophenyl)(2-methoxyphenyl)methyl)phenol (1b) was prepared as a white solid from 4-fluorobenzaldehyde (4.0 mmol) according to the General Procedure A (eluent: hexanes/EtOAc = 30:1 → 10:1) in 27% yield (over two steps).

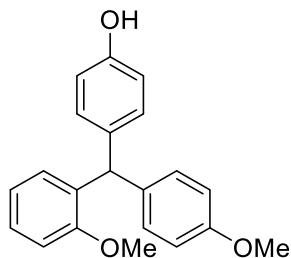
¹H NMR (400 MHz, CDCl₃) δ 7.22 (t, *J* = 7.4 Hz, 1H), 7.04 – 7.01 (m, 2H), 6.94 (t, *J* = 8.0 Hz, 4H), 6.88 – 6.81 (m, 3H), 6.73 (d, *J* = 8.4 Hz, 2H), 5.82 (s, 1H), 3.71 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 161.5 (d, *J* = 242.5 Hz), 157.2, 154.0, 140.1 (d, *J* = 3.2 Hz), 136.2, 132.9, 130.9 (d, *J* = 7.6 Hz), 130.7, 130.3, 127.8, 120.5, 115.3, 115.0 (d, *J* = 21.0 Hz), 110.9, 55.8, 48.2.

¹⁹F NMR (376 MHz, CDCl₃) δ -117.4.

IR (thin film) 3356, 3016, 2836, 1500, 1451, 1229, 1163, 752, 704 cm⁻¹.

HRMS (CI+) Calcd for C₂₀H₁₇FO₂ (M⁺): 308.1213, Found: 278.1223.



1c

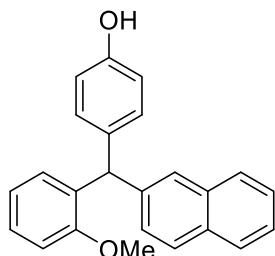
4-((2-Methoxyphenyl)(4-methoxyphenyl)methyl)phenol (1c) was prepared as a light yellow solid from 4-methoxybenzaldehyde (4.0 mmol) according to the General Procedure A (eluent: hexanes/EtOAc = 30:1 → 10:1) in 80% yield (over two steps).

¹H NMR (400 MHz, CDCl₃) δ 7.24 – 7.19 (m, 1H), 7.01 (d, *J* = 8.4, 2H), 6.94 (d, *J* = 8.4 Hz, 2H), 6.90 – 6.86 (m, 3H), 6.82 (d, *J* = 8.6 Hz, 2H), 6.72 (d, *J* = 8.4 Hz, 2H), 5.81 (s, 1H), 3.79 (s, 3H), 3.72 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 157.9, 157.2, 153.8, 136.7, 136.5, 133.4, 130.6, 130.4, 130.3, 127.6, 120.4, 115.1, 113.6, 110.9, 55.8, 55.4, 48.0.

IR (thin film) 3381, 2955, 2835, 1502, 1452, 1027, 824, 733, 574 cm⁻¹.

HRMS (CI+) Calcd for C₂₁H₂₀O₃ (M⁺): 320.1412, Found: 320.1407.



1d

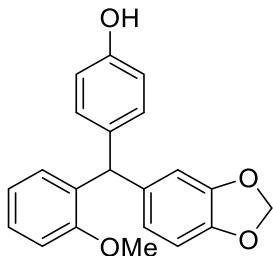
4-((2-Methoxyphenyl)(naphthalen-2-yl)methyl)phenol (1d) was prepared as a white solid from 2-naphthaldehyde (4.0 mmol) according to the General Procedure A (eluent: hexanes/EtOAc = 30:1 → 10:1) in 68% yield (over two steps).

¹H NMR (400 MHz, CDCl₃) δ 7.81 – 7.69 (m, 3H), 7.42 (d, *J* = 2.6 Hz, 3H), 7.30 – 7.22 (m, 2H), 7.00 (d, *J* = 7.6 Hz, 2H), 6.91– 6.86 (m, 3H), 6.74 (d, *J* = 8.5 Hz, 2H), 6.02 (s, 1H), 3.74 (s, 3H), 2.19 (s, 1H).

¹³C NMR (100 MHz, CDCl₃) δ 157.3, 154.0, 142.1, 136.1, 133.6, 132.9, 132.3, 130.9, 130.6, 128.6, 128.0, 127.8, 127.7, 127.6, 126.0, 125.6, 120.5, 115.2, 110.9, 55.8, 49.0.

IR (thin film) 3391, 3054, 2836, 1598, 1502, 1451, 1239, 1171, 822, 747 cm⁻¹.

HRMS (CI+) Calcd for C₂₄H₂₀O₂ (M⁺): 340.1463, Found: 340.1451.



1e

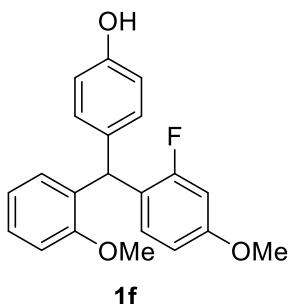
4-(Benzo[*d*][1,3]dioxol-5-yl)(2-methoxyphenyl)methylphenol (1e) was prepared as a yellow solid from benzo[*d*][1,3]dioxole-5-carbaldehyde (4.0 mmol) according to the General Procedure A (eluent: hexanes/EtOAc = 30:1 → 10:1) in 70% yield (over two steps).

¹H NMR (400 MHz, CDCl₃) δ 7.24 – 7.19 (m, 1H), 6.94 (d, *J* = 8.3 Hz, 2H), 6.88 – 6.87 (m, 3H), 6.74 – 6.70 (m, 3H), 6.60 – 6.53 (m, 2H), 5.91 (s, 2H), 5.78 (s, 1H), 4.81 (s, 1H), 3.73 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 157.2, 153.9, 147.6, 145.9, 138.4, 136.5, 133.1, 130.6, 130.3, 127.7, 122.5, 120.5, 115.2, 110.9, 110.2, 108.0, 101.0, 55.8, 48.5.

IR (thin film) 3383, 2891, 1598, 1485, 1437, 1234, 1033, 746 cm⁻¹.

HRMS (CI+) Calcd for C₂₁H₁₈O₄ (M⁺): 334.1205, Found: 334.1204



1f

4-((2-Fluoro-4-methoxyphenyl)(2-methoxyphenyl)methyl)phenol (1f) was prepared as a white solid from 2-fluoro-4-methoxybenzaldehyde (4.0 mmol) according to the General Procedure A (eluent: hexanes/EtOAc = 30:1 → 10:1) in 52% yield (over two steps).

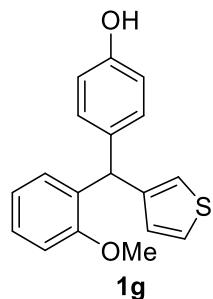
¹H NMR (400 MHz, CDCl₃) δ 7.26 – 7.22 (m, 1H), 6.95 (d, *J* = 8.4 Hz, 2H), 6.88 – 6.72 (m, 6H), 6.62 – 6.56 (m, 2H), 6.00 (s, 1H), 3.77 (s, 3H), 3.71 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 161.3 (d, *J* = 245.0 Hz), 159.5 (d, *J* = 10.5 Hz), 157.2, 154.0, 135.3, 132.1, 131.0 (d, *J* = 5.9 Hz), 130.5, 130.0, 127.8, 123.5 (d, *J* = 15.3 Hz), 120.4, 115.2, 110.9, 109.3 (d, *J* = 2.9 Hz), 101.8 (d, *J* = 26.0 Hz), 55.8, 55.7, 41.6 (d, *J* = 2.7 Hz).

¹⁹F NMR (376 MHz, CDCl₃) δ -114.2.

IR (thin film) 3353, 2836, 1617, 1587, 1500, 1451, 1241, 1098, 1025, 753 cm⁻¹.

HRMS (CI+) Calcd for C₂₁H₁₉FO₃ (M⁺): 338.1318, Found: 338.1321.



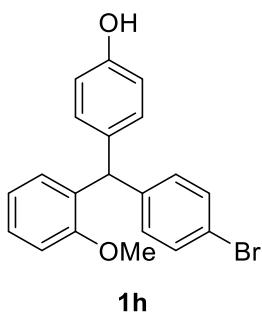
4-((2-Methoxyphenyl)(thiophen-3-yl)methyl)phenol (1g) was prepared as a yellow solid from thiophene-3-carbaldehyde (4.0 mmol) according to the General Procedure A (eluent: hexanes/EtOAc = 30:1 → 10:1) in 31% yield (over two steps).

¹H NMR (400 MHz, CDCl₃) δ 7.27 – 7.21 (m, 2H), 7.03 (d, *J* = 8.4 Hz, 2H), 6.96 (d, *J* = 7.0 Hz, 1H), 6.92 – 6.87 (m, 3H), 6.74 – 6.70 (m, 3H), 5.87 (s, 1H), 3.75 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 157.0, 153.9, 145.4, 136.3, 132.9, 130.3, 129.8, 129.1, 127.7, 125.3, 122.5, 120.6, 115.2, 110.9, 55.8, 44.5.

IR (thin film) 3395, 2834, 1597, 1497, 1447, 1232m 1169, 1100, 1021, 746 cm⁻¹.

HRMS (CI+) Calcd for C₁₈H₁₆O₂S (M⁺): 296.0871, Found: 296.0874.



1h

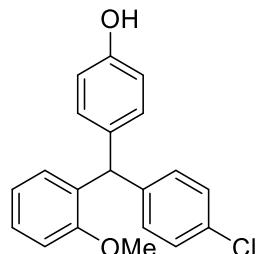
4-((4-Bromophenyl)(2-methoxyphenyl)methyl)phenol (1h) was prepared as a white solid from 4-bromobenzaldehyde (4.0 mmol) according to the General Procedure A (eluent: hexanes/EtOAc = 30:1 → 10:1) in 30% yield (over two steps).

¹H NMR (400 MHz, CDCl₃) δ 7.37 (d, *J* = 8.4 Hz, 2H), 7.03 (td, *J*₁ = 7.7 Hz, *J*₂ = 1.5 Hz, 1H), 6.96 – 6.82 (m, 7H), 6.73 (d, *J* = 8.4 Hz, 2H), 5.79 (s, 1H), 4.76 (br, 1H), 3.72 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 157.2, 154.1, 143.5, 135.7, 132.5, 131.4, 131.3, 130.7, 130.3, 127.9, 120.5, 120.0, 115.3, 110.9, 55.8, 48.5.

IR (thin film) 3360, 2835, 1597, 1510, 1485, 1240, 1173, 823, 754 cm⁻¹.

HRMS (CI+) Calcd for C₂₀H₁₇O₂Br (M⁺): 368.0412, Found: 368.0411.



1i

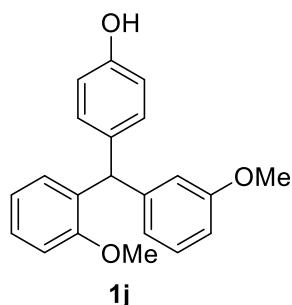
4-((4-Chlorophenyl)(2-methoxyphenyl)methyl)phenol (1i) was prepared as a white solid from 4-chlorobenzaldehyde (4.0 mmol) according to the General Procedure A (eluent: hexanes/EtOAc = 30:1 → 10:1) in 42% yield (over two steps).

¹H NMR (400 MHz, CDCl₃) δ 7.26 – 7.21 (m, 3H), 6.99 – 6.80 (m, 7H), 6.81 (d, *J* = 7.2 Hz, 2H), 5.80 (s, 1H), 3.71 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 157.2, 154.1, 143.0, 135.8, 132.6, 131.9, 130.8, 130.7, 130.3, 128.4, 127.9, 120.5, 115.3, 110.9, 55.8, 48.4.

IR (thin film) 3352, 3028, 1596, 1486, 1453, 1239, 1172, 822, 750 cm⁻¹.

HRMS (CI+) Calcd for C₂₀H₁₇O₂Cl (M⁺): 324.0917, Found: 324.0924.



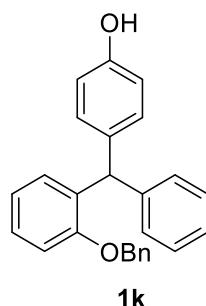
4-((2-Methoxyphenyl)(3-methoxyphenyl)methyl)phenol (1j) was prepared as a light yellow oil from 3-methoxybenzaldehyde (4.0 mmol) according to the General Procedure A (eluent: hexanes/EtOAc = 30:1 → 10:1) in 40% yield (over two steps).

¹H NMR (400 MHz, CDCl₃) δ 7.21 – 7.17 (m, 2H), 6.95 (d, J = 8.4 Hz, 2H), 6.88 – 6.86 (m, 3H), 6.76 – 6.66 (m, 5H), 5.84 (s, 1H), 3.74 (s, 3H), 3.72 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 159.6, 157.2, 153.9, 146.1, 136.2, 132.9, 130.7, 130.4, 129.2, 127.7, 122.2, 120.5, 115.7, 115.2, 111.2, 110.9, 55.8, 55.3, 48.8.

IR (thin film) 3383, 3026, 2889, 1690, 1603, 1506, 1452, 1249, 1053, 930, 703 cm⁻¹.

HRMS (CI+) Calcd for C₂₁H₂₀O₃ (M⁺): 320.1412, Found: 320.1407.



4-((2-(Benzylxy)phenyl)(phenyl)methyl)phenol (1k) was prepared as a white solid from 2-(benzyloxy)benzaldehyde (4.0 mmol) according to the

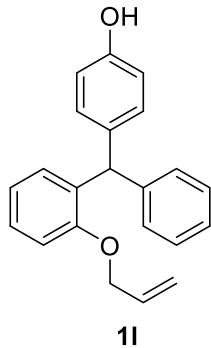
General Procedure B (eluent: hexanes/EtOAc = 30:1 → 10:1) in 55% yield (over two steps).

¹H NMR (400 MHz, CDCl₃) δ 7.29 – 7.17 (m, 7H), 7.09 – 7.03 (m, 4H), 6.97 – 6.83 (m, 5H), 6.76 – 6.72 (m, 2H), 5.88 (s, 1H), 4.96 (s, 2H).

¹³C NMR (100 MHz, CDCl₃) δ 156.3, 154.0, 144.3, 137.3, 136.3, 133.6, 130.9, 130.5, 129.6, 128.5, 128.3, 127.8, 127.7, 127.3, 126.2, 120.7, 115.2, 112.0, 70.1, 49.5.

IR (thin film) 3381, 3028, 1703, 1595, 1556, 1232, 1171, 743, 698 cm⁻¹.

HRMS (CI+) Calcd for C₂₆H₂₂O₂ (M⁺): 366.1620, Found: 366.1634.



1l

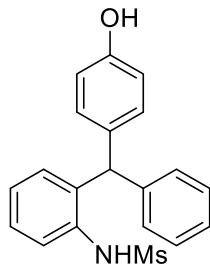
4-((2-(Allyloxy)phenyl)(phenyl)methyl)phenol (1l) was prepared as a white solid from 2-(allyloxy)benzaldehyde (4.0 mmol) according to the General Procedure B (eluent: hexanes/EtOAc = 30:1 → 10:1) in 30% yield (over two steps).

¹H NMR (400 MHz, CDCl₃) δ 7.31 – 7.20 (m, 4H), 7.13 (d, *J* = 7.3 Hz, 2H), 6.99 (d, *J* = 8.4 Hz, 2H), 6.90 – 6.87 (m, 3H), 6.74 (d, *J* = 8.4 Hz, 2H), 5.92 (s, 1H), 5.89 – 5.80 (m, 1H), 5.26 – 5.15 (m, 2H), 5.05 (s, 1H), 4.48 – 4.42 (m, 2H).

¹³C NMR (100 MHz, CDCl₃) δ 156.1, 153.8, 144.3, 136.1, 133.6, 133.4, 130.7, 130.4, 129.5, 128.3, 127.6, 126.1, 120.7, 116.9, 115.2, 112.3, 69.1, 49.3.

IR (thin film) 3524, 3025, 1597, 1497, 1447, 1229, 1170, 993, 738, 700 cm⁻¹.

HRMS (CI+) Calcd for C₂₂H₂₀O₂ (M⁺): 316.1463, Found: 316.1470.



1m

N-(2-(Hydroxy(4-hydroxyphenyl)(phenyl)methyl)phenyl)methane sulfonamide (1m)

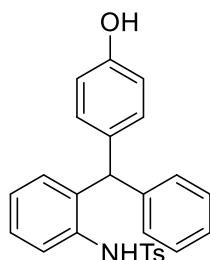
sulfonamide (1m) was prepared as a white solid from *N*-(2-benzoylphenyl)methanesulfonamide (4.0 mmol) according to the General Procedure C (eluent: hexanes/EtOAc = 20:1 → 3:1) in 40% yield (over two steps).

¹H NMR (400 MHz, CDCl₃) δ 7.54 – 7.52 (m, 1H), 7.34 – 7.26 (m, 4H), 7.15 – 7.07 (m, 3H), 6.96 – 6.94 (m, 2H), 6.88 – 6.85 (m, 1H), 6.80 – 6.78 (m, 2H), 6.10 (s, 1H), 5.58 (s, 1H), 4.98 (br, 1H), 2.73 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 154.9, 142.0, 136.9, 134.8, 133.6, 130.83, 130.77, 129.6, 129.1, 128.2, 127.3, 126.1, 123.0, 116.0, 51.7, 40.0.

IR (thin film) 3315, 1603, 1502, 1447, 1318, 1228, 1144, 972, 742, 701 cm⁻¹.

HRMS (CI+) Calcd for C₂₀H₁₉NO₃S (M⁺): 353.1086, Found: 353.1049.



1n

N-(2-(Hydroxy(4-hydroxyphenyl)(phenyl)methyl)phenyl)-4-methylbenzenesulfonamide (1n) was prepared as a white solid from *N*-(2-benzoylphenyl)-4-methylbenzenesulfonamide (4.0 mmol) according to the General Procedure C (eluent: hexanes/EtOAc = 20:1 → 3:1) in 90% yield (over two steps).

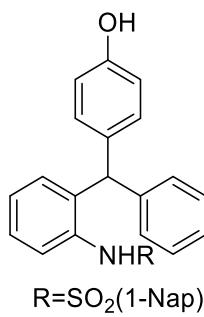
¹H NMR (400 MHz, CDCl₃) δ 7.58 (d, *J* = 8.2 Hz, 2H), 7.47 (dd, *J*₁ = 8.2 Hz, *J*₂ =

1.0 Hz, 1H), 7.30 – 7.20 (m, 5H), 7.09 – 7.05 (m, 1H), 6.74 – 6.71 (m, 4H), 6.63 – 6.61 (m, 3H), 6.09 (s, 1H), 4.91 (s, 1H), 4.82 (s, 1H), 2.45 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 154.8, 144.2, 141.8, 138.2, 137.2, 134.3, 133.5, 130.7, 130.1, 130.0, 129.4, 129.0, 127.9, 127.3, 127.2, 126.5, 126.2, 115.9, 51.2, 21.8.

IR (thin film) 3322, 3059, 3029, 1601, 1505, 1446, 1264, 1221, 1156, 1090, 752, 558 cm⁻¹.

HRMS (CI+) Calcd for C₂₆H₂₃NO₃S (M⁺): 429.1399, Found: 429.1286.



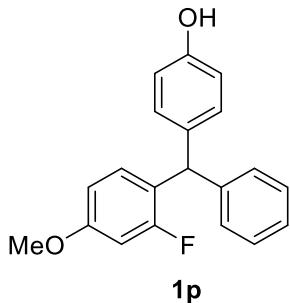
N-(2-(Hydroxy(4-hydroxyphenyl)(phenyl)methyl)phenyl)naphthalene-1-sulfonamide (1o) was prepared as a light yellow solid from *N*-(2-benzoylphenyl)naphthalene-1-sulfonamide (4.0 mmol) according to the General Procedure C (eluent: hexanes/EtOAc = 20:1 → 3:1) in 68% yield (over two steps).

¹H NMR (400 MHz, CDCl₃) δ 8.56 – 8.53 (m, 1H), 8.19 (d, *J* = 7.4 Hz, 1H), 8.10 (d, *J* = 8.1 Hz, 1H), 7.99 – 7.97 (m, 1H), 7.62 – 7.60 (m, 2H), 7.51 (t, *J* = 7.8 Hz, 1H), 7.41 (d, *J* = 8.0 Hz, 1H), 7.18 – 7.15 (m, 4H), 7.01 (t, *J* = 7.6 Hz, 1H), 6.66 (d, *J* = 8.4 Hz, 2H), 6.51 – 6.49 (m, 3H), 6.41 (d, *J* = 8.4 Hz, 2H), 6.30 (s, 1H), 4.78 (s, 1H).

¹³C NMR (100 MHz, CDCl₃) δ 154.7, 141.8, 138.3, 135.3, 134.8, 134.6, 134.3, 133.6, 130.6, 130.3, 130.0, 129.4, 129.3, 128.8, 128.4, 127.6, 127.3, 127.1, 126.5, 126.1, 124.7, 124.6, 115.8, 50.8.

IR (thin film) 3444, 3343, 1602, 1505, 1325, 1263, 1160, 758 cm⁻¹.

HRMS (CI+) Calcd for C₂₉H₂₃NO₃S (M⁺): 465.1399, Found: 465.1424.



4-((2-Fluoro-4-methoxyphenyl)(phenyl)methyl)phenol (1p) was prepared as a white solid from 2-fluoro-4-methoxybenzaldehyde (4.0 mmol) according to the General Procedure B (eluent: hexanes/EtOAc = 30:1 → 10:1) in 40% yield (over two steps).

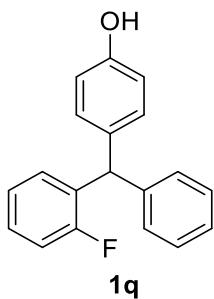
¹H NMR (400 MHz, CDCl₃) δ 7.31 – 7.20 (m, 3H), 7.10 (d, *J* = 7.5 Hz, 2H), 6.97 (d, *J* = 8.4 Hz, 2H), 6.82 (t, *J* = 8.7 Hz, 1H), 6.75 (d, *J* = 8.4 Hz, 2H), 6.64 – 6.60 (m, 2H), 5.69 (s, 1H), 3.78 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 161.2 (d, *J* = 245.0 Hz), 159.6 (d, *J* = 10.9 Hz), 154.2, 143.4, 135.4, 131.2 (d, *J* = 5.8 Hz), 130.6, 129.3, 128.5, 126.6, 123.5 (d, *J* = 15.0 Hz), 115.4, 109.6 (d, *J* = 3.0 Hz), 101.8 (d, *J* = 25.7 Hz), 55.7, 48.2 (d, *J* = 2.3 Hz).

¹⁹F NMR (376 MHz, CDCl₃) δ -114.4.

IR (thin film) 3377, 3021, 2837, 1617, 1586, 1502, 1441, 1252, 1156, 1096, 1027, 731, 699 cm⁻¹.

HRMS (CI+) Calcd for C₂₀H₁₇FO₂ (M⁺): 308.1213, Found: 308.1202.



4-((2-Fluorophenyl)(phenyl)methyl)phenol (1q) was prepared as a white solid from 2-fluorobenzaldehyde (4.0 mmol) according to the General Procedure B (eluent: hexanes/EtOAc = 30:1 → 10:1) in 77% yield (over two steps).

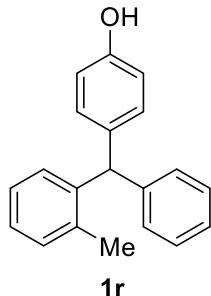
¹H NMR (400 MHz, CDCl₃) δ 7.30 – 7.18 (m, 4H), 7.10 – 6.91 (m, 7H), 6.75 (d, *J* = 8.5 Hz, 2H), 5.77 (s, 1H), 4.87 (s, 1H).

¹³C NMR (100 MHz, CDCl₃) δ 160.9 (d, *J* = 245.3 Hz), 154.3, 143.1, 135.1, 131.6 (d, *J* = 14.3 Hz), 131.0 (d, *J* = 3.9 Hz), 130.6, 129.4, 128.5, 128.3 (d, *J* = 8.4 Hz), 126.7, 124.0 (d, *J* = 3.5 Hz), 115.6, 115.4, 48.7 (d, *J* = 3.3 Hz).

¹⁹F NMR (376 MHz, CDCl₃) δ –116.5.

IR (thin film) 3298, 3026, 1600, 1485, 1446, 1223, 1172, 1095, 748, 698 cm⁻¹.

HRMS (CI+) Calcd for C₁₉H₁₅FO (M⁺): 278.1107, Found: 278.1103.



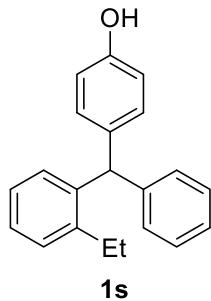
4-(Phenyl(*o*-tolyl)methyl)phenol (1r) was prepared as a white solid from 2-methylbenzaldehyde (3.3 mmol) according to the General Procedure B (eluent: hexanes/EtOAc = 30:1 → 10:1) in 35% yield (over two steps).

¹H NMR (400 MHz, CDCl₃) δ 7.30 – 7.28 (m, 1H), 7.23 – 7.05 (m, 6H), 6.94 – 6.91 (m, 2H), 6.83 – 6.80 (m, 1H), 6.76 – 6.74 (m, 2H), 5.61 (s, 1H), 4.74 (s, 1H), 2.22 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 154.1, 143.9, 142.8, 136.8, 135.9, 130.9, 130.6, 129.7, 129.5, 128.5, 126.5, 126.4, 125.9, 115.3, 52.9, 20.1.

IR (thin film) 3348, 3057, 3021, 2955, 1603, 1504, 1446, 1226, 1171, 828, 740, 701 cm⁻¹.

HRMS (CI+) Calcd for C₂₀H₁₈O (M⁺): 274.1358, Found: 274.1360.



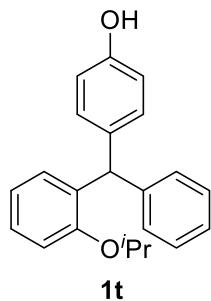
4-((2-Ethylphenyl)(phenyl)methyl)phenol (1s) was prepared as a colorless oil from 2-ethylbenzaldehyde (3.3 mmol) according to the General Procedure B (eluent: hexanes/EtOAc = 30:1 → 10:1) in 72% yield (over two steps).

¹H NMR (400 MHz, CDCl₃) δ 7.36 – 7.25 (m, 5H), 7.22 – 7.15 (m, 3H), 7.02 – 6.94 (m, 3H), 6.80 – 6.77 (m, 2H), 5.81 (s, 1H), 5.36 (br, 1H), 2.70 (q, *J* = 7.5 Hz, 2H), 1.23 (s, *J* = 7.5 Hz, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 153.8, 144.3, 142.4, 141.9, 136.3, 130.9, 129.9, 129.7, 128.7, 128.4, 126.7, 126.3, 125.8, 115.3, 51.9, 25.7, 15.2.

IR (thin film) 3357, 2362, 1604, 1507, 1446, 1257, 1233, 1171, 1101, 1029, 829, 734, 701 cm⁻¹.

HRMS (ES-) Calcd for C₂₁H₁₉O[·] (M-H⁺): 287.1441, Found: 287.1431.



4-((2-Isopropoxyphenyl)(phenyl)methyl)phenol (1t) was prepared as a colorless oil by the following procedure. To the solution of 4-(hydroxy(2-isopropoxy-phenyl)(phenyl)methyl)phenol (136.6 mg, 0.47 mmol) in DCM (3 mL) were added 2-(naphthalen-2-yl)-2,3-dihydrobenzo[d]thiazole (245.0 mg, 0.94 mmol, 2.0 equiv.) and 4-hydroxydinaphtho[2,1-d:1',2'-f][1,3,2]dioxaphosphepine 4-oxide (16.6 mg, 47 μmmol). The mixture was stirred for 16 h until

full conversion of the substrate. Then, the mixture was directly subjected to silica gel column chromatography (eluent: hexanes/ethyl acetate = 30:1 → 10:1) to afford the desired product **1t** in 85% yield.

¹H NMR (400 MHz, CDCl₃) δ 7.28 – 7.25 (m, 2H), 7.21 – 7.14 (m, 2H), 7.11 – 7.10 (m, 2H), 7.00 – 6.96 (m, 2H), 6.87 – 6.81 (m, 3H), 6.76 – 6.72 (m, 2H), 5.81 (s, 1H), 4.69 (s, 1H), 4.49 – 6.43 (m, 1H), 1.11 – 1.07 (m, 6H).

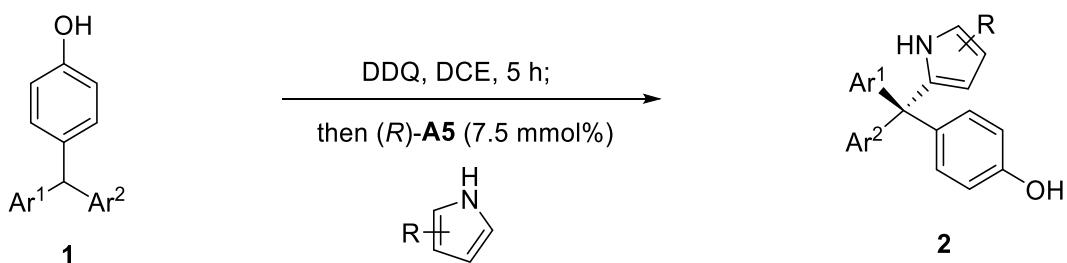
¹³C NMR (100 MHz, CDCl₃) δ 155.5, 153.8, 144.4, 136.4, 134.4, 130.9, 130.4, 129.6, 128.2, 127.4, 126.0, 120.0, 115.1, 113.1, 70.1, 49.5, 22.1, 22.0.

IR (thin film) 3358, 2976, 2360, 1707, 1508, 1483, 1446, 1336, 1233, 1171, 1111, 740, 700 cm⁻¹.

HRMS (ES-) Calcd for C₂₂H₂₁O₂⁻ (M-H⁺): 317.1547, Found: 317.1541.

III. Catalytic Asymmetric Oxidative Synthesis of Tetraarylmethanes

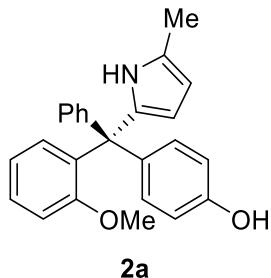
General Procedure D.



At room temperature, to an oven-dried 10-mL vial charged with a solution of the triarylmethane **1** (0.25 mmol) and DDQ (57.0 mg, 0.25 mmol, 1 equiv) in DCE (5 mL). The mixture was stirred for 5 h and then cooled to 0 °C. The catalyst (R)-**A5** (18.6 mg, 18.8 µmmol, 7.5 mol%) and the nucleophile pyrrole (0.5 mmol, 2 equiv) were added. The mixture was stirred at 0 °C for 3 h. Then, Na₂CO₃ (212 mg, 2.0 mmol) was added. The mixture was stirred for additional 10 min before it was concentrated under reduced pressure. The residue was purified by silica gel column chromatography to afford the desired product **2**.

General Procedure E.

At room temperature, to an oven-dried 10-mL vial charged with a solution of the triarylmethane **1** (0.25 mmol) and DDQ (57.0 mg, 0.25 mmol, 1 equiv) in DCE (5 mL). The mixture was stirred for 12 h and then cooled to -20 °C. The catalyst (R)-**A5** (18.6 mg, 18.8 µmmol, 7.5 mol%) and the nucleophile pyrrole (0.5 mmol, 2 equiv) were added. The mixture was stirred at -20 °C for 12 h. Then, Na₂CO₃ (212 mg, 2.0 mmol) was added. The mixture was stirred for additional 10 min before it was concentrated under reduced pressure. The residue was purified by silica gel column chromatography to afford the desired product **2**.



(R)-4-((2-Methoxyphenyl)(5-methyl-1H-pyrrol-2-yl)(phenyl)methyl)phenol

(2a) was prepared as a light yellow foam from 4-((2-methoxyphenyl)(phenyl)methyl)phenol (72.5 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/EtOAc = 20:1 → 10:1) in 95% yield (92.0 mg, 96% ee).

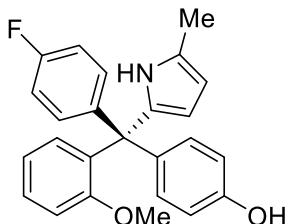
$[\alpha]_D^{22}$: +4.5 ($c = 1.0$, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK IC column; 3% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 10.8 min (major), 12.4 min (minor).

¹H NMR (400 MHz, CDCl₃) δ 7.59 (s, 1H), 7.30 – 7.13 (m, 7H), 7.00 – 6.87 (m, 5H), 6.68 (d, *J* = 8.6 Hz, 2H), 5.76 (s, 1H), 5.70 (t, *J* = 2.9 Hz, 1H), 3.20 (s, 3H), 2.17 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 158.1, 153.8, 145.7, 137.4, 136.8, 135.9, 131.5, 130.3, 129.9, 128.7, 127.4, 127.1, 126.0, 120.8, 114.2, 113.8, 110.4, 105.1, 58.5, 55.8, 13.4.

IR (thin film) 3446, 3384, 3056, 2983, 2839, 1588, 1247, 1039, 733, 700 cm⁻¹.

HRMS (CI+) Calcd for C₂₅H₂₃NO₂ (M⁺): 369.1729, Found: 369.1725.



(S)-4-((4-Fluorophenyl)(2-methoxyphenyl)(5-methyl-1H-pyrrol-2-yl)methyl)phenol **(2b)** was prepared as a light yellow foam from 4-((4-fluorophenyl)(2-

methoxyphenyl)methyl)phenol (77.7 mg, 0.25 mmol) and 2-methylpyrrole (40.6

mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/EtOAc = 20:1 → 10:1) in 61% yield (59.5 mg, 97% ee).

$[\alpha]_D^{22}$: +8.7 ($c = 1.0$, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK IC column; 3% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 9.0 min (major), 10.1 min (minor).

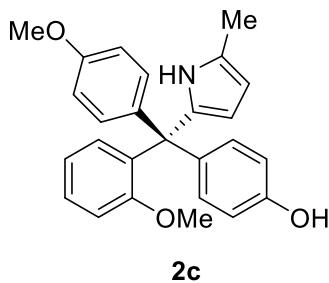
¹H NMR (400 MHz, acetone-*d*₆) δ 8.84 (s, 1H), 8.26 (s, 1H), 7.28 (t, *J* = 7.6 Hz, 1H), 7.15 – 7.12 (m, 2H), 6.98 – 6.87 (m, 7H), 6.70 (d, *J* = 8.6 Hz, 2H), 5.63 (s, 1H), 5.55 (t, *J* = 2.9 Hz, 1H), 3.22 (s, 3H), 2.13 (s, 3H).

¹³C NMR (100 MHz, acetone-*d*₆) δ 161.6 (d, *J* = 241.0 Hz), 158.7, 156.2, 142.9 (d, *J* = 3.0 Hz), 137.4, 136.4 (d, *J* = 6.0 Hz), 132.1 (d, *J* = 7.6 Hz), 131.9 (2C), 130.7, 129.3, 128.0, 121.0, 114.6, 114.1, 113.9 (d, *J* = 21.1 Hz), 110.4, 105.4, 58.6, 55.4, 13.1.

¹⁹F NMR (376 MHz, acetone-*d*₆) δ -119.8.

IR (thin film) 3447, 3053, 2980, 2936, 1593, 1476, 1261, 1020, 733 cm⁻¹.

HRMS (CI+) Calcd for C₂₅H₂₂NFO₂ (M⁺): 387.1635, Found: 387.1638.



(S)-4-((2-Methoxyphenyl)(4-methoxyphenyl)(5-methyl-1*H*-pyrrol-2-yl)methyl)phenol (2c) was prepared as a light yellow foam from 4-((2-methoxyphenyl)(4-methoxyphenyl)methyl)phenol (80.0 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/EtOAc = 20:1 → 10:1) in 95% yield (94.4 mg, 90% ee).

$[\alpha]_D^{22}$: +5.6 ($c = 1.0$, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK IC column; 5% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 8.9 min (major), 10.3 min (minor).

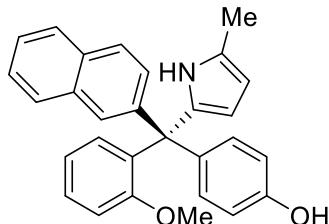
¹H NMR (400 MHz, CDCl₃) δ 7.60 (s, 1H), 7.29 – 7.26 (m, 1H), 7.05 (d, *J* = 8.8 Hz,

2H), 6.99 (d, J = 8.6 Hz, 2H), 6.94 – 6.87 (m, 3H), 6.77 (d, J = 8.8 Hz, 2H), 6.68 (d, J = 8.6 Hz, 2H), 5.76 (s, 1H), 5.69 (t, J = 2.9 Hz, 1H), 5.08 (s, 1H), 3.79 (s, 3H), 3.23 (s, 3H), 2.17 (s, 3H).

^{13}C NMR (100 MHz, CDCl_3) δ 158.1, 157.7, 153.7, 137.9, 137.8, 137.0, 136.2, 131.2, 131.1, 130.3, 128.6, 126.9, 120.8, 114.1, 113.8, 112.6, 110.2, 105.1, 57.8, 55.9, 55.4, 13.4.

IR (thin film) 3445, 3389, 3050, 2838, 1590, 1249, 1176, 1031, 821, 735 cm^{-1} .

HRMS (LD+) Calcd for $\text{C}_{26}\text{H}_{25}\text{NO}_3$ (M^+): 399.1834, Found: 399.1827.



2d

(S)-4-((2-Methoxyphenyl)(5-methyl-1*H*-pyrrol-2-yl)(naphthalen-2-yl)methyl)phenol (2d) was prepared as a light yellow foam from 4-((2-methoxyphenyl)(naphthalen-2-yl)methyl)phenol (85.1 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/EtOAc = 20:1 → 10:1) in 95% yield (99.1 mg, 94% ee).

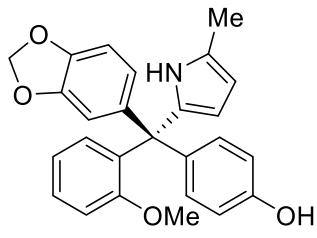
$[\alpha]_D^{22}$: +22.3 (c = 1.0, CHCl_3). HPLC analysis of the product: Daicel CHIRALPAK IC column; 3% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 11.4 min (major), 12.9 min (minor).

^1H NMR (400 MHz, acetone- d_6) δ 8.89 (s, 1H), 8.27 (s, 1H), 7.83 (d, J = 7.4 Hz, 1H), 7.72 – 7.60 (m, 3H), 7.45 – 7.28 (m, 4H), 7.09 – 6.90 (m, 5H), 6.73 (d, J = 8.4 Hz, 2H), 5.69 (s, 2H), 3.16 (s, 3H), 2.13 (s, 3H).

^{13}C NMR (100 MHz, acetone- d_6) δ 159.0, 156.1, 144.1, 137.4, 136.8, 136.3, 133.9, 132.8, 132.0, 130.9, 130.4, 129.3, 128.9, 128.4, 128.1, 128.0, 126.4, 126.3, 126.2, 121.1, 114.5, 114.0, 110.5, 105.5, 59.2, 55.5, 13.1.

IR (thin film) 3441, 3380, 3052, 2836, 1588, 1243, 734, 698 cm^{-1} .

HRMS (LD+) Calcd for C₂₉H₂₅NO₂ (M⁺): 419.1885, Found: 419.1883.



2e

(S)-4-(Benzo[d][1,3]dioxol-5-yl(2-methoxyphenyl)(5-methyl-1*H*-pyrrol-2-yl)methylphenol (2e) was prepared as a light yellow foam from 4-(benzo[d][1,3]dioxol-5-yl(2-methoxyphenyl)methyl)phenol (83.5 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/EtOAc = 20:1 → 10:1) in 90% yield (92.7 mg, 92% ee).

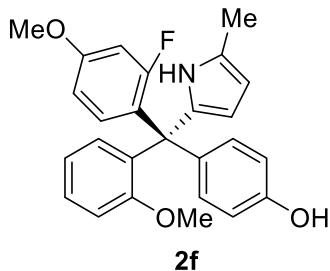
[α]_D²²: +9.2 (c = 1.0, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK IC column; 5% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 10.1 min (major), 14.1 min (minor).

¹H NMR (400 MHz, CDCl₃) δ 7.61 (s, 1H), 7.30 – 7.28 (m, 1H), 7.00 (d, J = 8.6 Hz, 2H), 6.95 – 6.86 (m, 3H), 6.70 – 6.67 (m, 4H), 6.60 (d, J = 8.2 Hz, 1H), 5.92 (s, 2H), 5.76 (s, 1H), 5.70 (t, J = 2.7 Hz, 1H), 3.27 (s, 3H), 2.18 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 158.0, 153.8, 146.9, 145.6, 139.9, 137.6, 136.7, 135.8, 131.4, 130.3, 128.7, 127.0, 123.1, 120.8, 114.1, 113.7, 111.0, 110.3, 107.0, 105.2, 101.0, 58.3, 55.8, 13.4.

IR (thin film) 3442, 3403, 3057, 2837, 1587, 1236, 734 cm⁻¹.

HRMS (LD+) Calcd for C₂₆H₂₃NO₄ (M⁺): 413.1627, Found: 413.1637.



(R)-4-((2-Fluoro-4-methoxyphenyl)(2-methoxyphenyl)(5-methyl-1*H*-pyrrol-2-yl)methyl)phenol (2f) was prepared as a light yellow foam from 4-((2-fluoro-4-methoxyphenyl)(2-methoxyphenyl)methyl)phenol (84.6 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/EtOAc = 20:1 → 10:1) in 87% yield (90.5 mg, 89% ee).

$[\alpha]_D^{22}$: -5.1 ($c = 1.0$, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK IC column; 10% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 6.6 min (major), 9.2 min (minor).

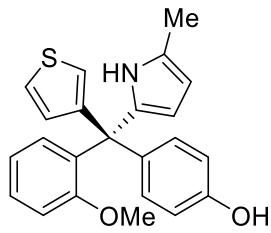
¹H NMR (400 MHz, CDCl₃) δ 7.76 (s, 1H), 7.30 – 7.26 (m, 1H), 7.03 – 6.85 (m, 6H), 6.67 (d, $J = 8.7$ Hz, 2H), 6.58 – 6.55 (m, 2H), 5.77 (s, 1H), 5.66 (s, 1H), 3.78 (s, 3H), 3.25 (s, 3H), 2.18 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 161.5 (d, $J = 247.8$ Hz), 159.8 (d, $J = 11.1$ Hz), 157.9, 153.7, 136.6, 135.3, 134.2, 131.5 (d, $J = 5.4$ Hz), 131.0, 130.2, 128.7, 126.4, 125.7 (d, $J = 11.5$ Hz), 120.9, 114.0 (d, $J = 37.4$ Hz), 109.6, 108.7, 105.4, 102.6, 102.3, 55.93, 55.90, 55.7, 13.5.

¹⁹F NMR (376 MHz, CDCl₃) δ -99.6.

IR (thin film) 3446, 3391, 3053, 2839, 1616, 1256, 734 cm⁻¹.

HRMS (CI+) Calcd for C₂₆H₂₄FNO₃ (M⁺): 417.1740, Found: 417.1727.



2g

(R)-4-((2-Methoxyphenyl)(5-methyl-1H-pyrrol-2-yl)(thiophen-3-yl)methyl)phenol (2g) was prepared as a light yellow foam from 4-((2-methoxyphenyl)(thiophen-3-yl)methyl)phenol (74.2 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/EtOAc = 20:1 → 10:1) in 94% yield (88.4 mg, 84% ee).

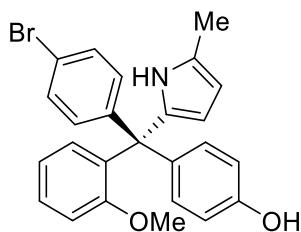
$[\alpha]_D^{22}$: +8.7 ($c = 1.0$, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK AD-H column; 5% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 13.3 min (major), 15.0 min (minor).

¹H NMR (400 MHz, CDCl₃) δ 7.84 (s, 1H), 7.28 – 7.26 (m, 1H), 7.21 – 7.19 (m, 1H), 6.99 (d, *J* = 8.6 Hz, 2H), 6.93 – 6.87 (m, 4H), 6.78 – 6.77 (m, 1H), 6.69 – 6.67 (m, 2H), 5.78 (s, 1H), 5.70 – 5.69 (m, 1H), 3.29 (s, 3H), 2.20 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 157.9, 153.8, 147.0, 138.4, 136.6, 135.2, 130.8, 130.2, 130.0, 128.7, 126.9, 123.7, 123.4, 120.9, 114.3, 113.9, 109.8, 105.3, 56.0, 55.5, 13.4.

IR (thin film) 3443, 3393, 3049, 2837, 1587, 1251, 1175, 733, 703 cm⁻¹.

HRMS (LD+) Calcd for C₂₃H₂₁NO₂S (M⁺): 375.1298, Found: 375.1311.



2h

(S)-4-((4-Bromophenyl)(2-methoxyphenyl)(5-methyl-1H-pyrrol-2-yl)methyl)phenol (2h) was prepared as a light yellow foam from 4-((4-bromophenyl)(2-methoxyphenyl)methyl)phenol (92.0 mg, 0.25 mmol) and 2-

methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/EtOAc = 20:1 → 10:1) in 90% yield (100.3 mg, 94% ee).

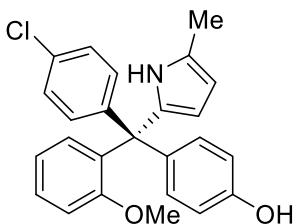
$[\alpha]_D^{22}$: +7.7 ($c = 1.0$, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK IC column; 3% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 11.0 min (major), 12.2 min (minor).

¹H NMR (400 MHz, CDCl₃) δ 7.53 (s, 1H), 7.35 – 7.26 (m, 3H), 7.05 – 6.86 (m, 7H), 6.70 (d, $J = 8.6$ Hz, 2H), 5.76 (s, 1H), 5.68 (t, $J = 2.9$ Hz, 1H), 3.24 (s, 3H), 2.17 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 157.8, 153.9, 144.9, 136.8, 136.1, 135.3, 131.6, 131.5, 130.3, 130.2, 128.9, 127.4, 120.9, 119.9, 114.3, 113.5, 110.5, 105.3, 58.1, 55.6, 13.4.

IR (thin film) 3436, 1589, 1483, 1435, 1245, 1177, 1015, 820, 747 cm⁻¹.

HRMS (LD+) Calcd for C₂₅H₂₂BrNO₂ (M⁺): 447.0834, Found: 447.0829.



2i

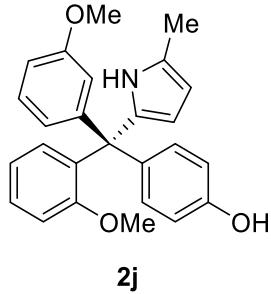
(S)-4-((4-Chlorophenyl)(2-methoxyphenyl)(5-methyl-1*H*-pyrrol-2-yl)methyl)phenol (2i)

was prepared as a light yellow foam from 4-((4-chlorophenyl)(2-methoxyphenyl)methyl)phenol (81.0 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/EtOAc = 15:1 → 10:1) in 89% yield (89.6 mg, 95% ee).

$[\alpha]_D^{22}$: +2.6 ($c = 1.0$, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK IC column; 3% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 8.8 min (major), 9.7 min (minor).

¹H NMR (400 MHz, CDCl₃) δ 7.51 (s, 1H), 7.31–7.27 (m, 1H), 7.19 (d, $J = 8.6$ Hz, 2H), 7.08 (d, $J = 8.6$ Hz, 2H), 6.94 – 6.86 (m, 5H), 6.66 (d, $J = 8.6$ Hz, 2H), 5.76 (s, 1H), 5.68 – 5.67 (m, 1H), 3.24 (s, 2H), 2.18 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 157.8, 153.9, 144.4, 136.8, 136.2, 135.4, 131.6, 131.5, 131.2, 130.2, 128.9, 127.4, 120.9, 114.3, 113.5, 110.5, 105.3, 58.1, 55.6, 13.4.
IR (thin film) 3400, 3019, 2941, 1597, 1486, 1453, 1240, 1172, 1098, 823, 753 cm⁻¹.
HRMS (LD+) Calcd for C₂₅H₂₂ClNO₂ (M⁺): 403.1339, Found: 403.1336.



(S)-4-((2-Methoxyphenyl)(3-methoxyphenyl)(5-methyl-1*H*-pyrrol-2-yl)methyl)phenol (2j) was prepared as a light yellow foam from 4-((2-methoxyphenyl)(3-methoxyphenyl)methyl)phenol (80.0 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/EtOAc = 20:1 → 10:1) in 97% yield (96.8 mg, 96% ee).

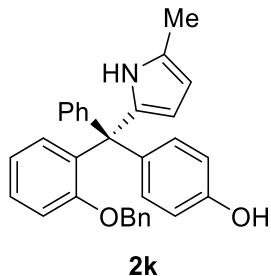
[α]_D²²: +10.1 (*c* = 1.0, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK IC column; 5% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 8.4 min (major), 10.4 min (minor).

¹H NMR (400 MHz, acetone-*d*₆) δ 8.82 (s, 1H), 8.18 (s, 1H), 7.27 (t, *J* = 7.8 Hz, 1H), 7.12 (t, *J* = 7.8 Hz, 1H), 6.97 – 6.93 (m, 4H), 6.87 (t, *J* = 7.6 Hz, 1H), 6.74 – 6.68 (m, 5H), 5.64 – 5.62 (m, 1H), 5.58 (t, *J* = 2.9 Hz, 1H), 3.66 (s, 3H), 3.21 (s, 3H), 2.13 (s, 3H).

¹³C NMR (100 MHz, acetone-*d*₆) δ 159.7, 159.0, 156.2, 148.7, 137.7, 136.8, 136.4, 132.1, 130.9, 129.2, 128.3, 127.8, 123.4, 121.1, 117.4, 114.5, 114.1, 110.8, 110.4, 105.4, 59.2, 55.6, 55.2, 13.1.

IR (thin film) 3445, 3384, 3053, 2931, 2839, 1590, 1375, 1042, 735, 700 cm⁻¹.

HRMS (CI+) Calcd for C₂₆H₂₅NO₃ (M⁺): 399.1834, Found: 399.1828.



(R)-4-((2-(Benzyl)phenyl)(5-methyl-1H-pyrrol-2-yl)(phenyl)methyl)phenol (2k) was prepared as a yellow foam from 4-((2-(benzyloxy)phenyl)(phenyl)methyl)phenol (91.6 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/EtOAc = 20:1 → 10:1) in 91% yield (101.1 mg, 97% ee).

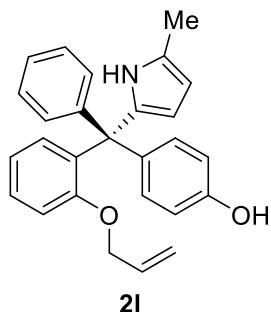
$[\alpha]_D^{22}$: +5.2 ($c = 1.0$, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK AD-H column; 5% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 13.4 min (major), 14.4 min (minor).

¹H NMR (400 MHz, CDCl₃) δ 7.94 (s, 1H), 7.30–7.17 (m, 9H), 7.08 – 7.03 (m, 3H), 6.95 – 6.91 (m, 2H), 6.79–6.77 (m, 2H), 6.66 (d, *J* = 8.6 Hz, 2H), 5.77 – 5.73 (m, 2H), 4.64 (s, 2H), 2.07 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 156.7, 153.7, 145.6, 137.4, 136.9, 136.3, 135.6, 131.6, 130.7, 130.0, 128.6, 128.2, 127.6, 127.4, 127.2, 127.0, 126.0, 120.8, 114.1, 113.6, 110.7, 104.9, 70.1, 58.6, 13.2.

IR (thin film) 3445, 3386, 3056, 2856, 1588, 1253, 1230, 731, 696 cm⁻¹.

HRMS (CI+) Calcd for C₃₁H₂₇NO₂ (M⁺): 445.2042, Found: 445.2058.



(R)-4-((2-(Allyloxy)phenyl)(5-methyl-1H-pyrrol-2-yl)(phenyl)methyl)phenol

(2l) was prepared as a red foam from 4-((2-(allyloxy)phenyl)(phenyl)methyl)phenol (79.0 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/EtOAc = 15:1 → 10:1) in 93% yield (91.9 mg, 93% ee).

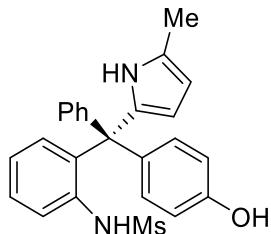
$[\alpha]_D^{22}$: -1.5 ($c = 1.0$, CHCl₃). HPLC analysis of the product derivative¹ after deallylation: Daicel CHIRALPAK IC column; 3% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 14.6 min (minor), 20.0 min (major).

¹H NMR (400 MHz, acetone-*d*₆) δ 8.83 (s, 1H), 8.19 (s, 1H), 7.27 – 7.12 (m, 6H), 7.00 – 6.85 (m, 5H), 6.71 – 6.67 (m, 2H), 5.64 – 5.62 (m, 1H), 5.58 – 5.56 (m, 1H), 5.41 – 5.32 (m, 1H), 4.91 – 4.84 (m, 2H), 4.08 – 3.98 (m, 2H), 2.12 (s, 3H).

¹³C NMR (100 MHz, acetone-*d*₆) δ 157.8, 156.1, 146.9, 137.5, 136.6, 136.5, 134.5, 132.2, 131.0, 130.6, 129.1, 127.7, 127.6, 126.2, 121.1, 116.3, 114.6, 114.4, 110.5, 105.4, 69.6, 59.2, 13.1.

IR (thin film) 3400, 3019, 2941, 1597, 1486, 1453, 1240, 1172, 1098, 823, 753 cm⁻¹.

HRMS (CI+) Calcd for C₂₇H₂₅NO₂ (M⁺): 395.1885, Found: 395.1871.



2m

(R)-N-(2-((4-Hydroxyphenyl)(phenyl)methyl)phenyl)methanesulfon-amide (2m) was prepared as a light yellow foam from *N*-(2-((4-hydroxyphenyl)(phenyl)methyl)phenyl)methanesulfonamide (88.3 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure E (eluent: hexanes/EtOAc = 15:1 → 10:1) in 85% yield (91.7 mg, 87%

(1) D. R. Vutukuri, P. Bharathi, Z. Yu, K. Rajasekaran, M.-H. Tran and S. Thayumanavan, *J. Org. Chem.*, 2003, **68**, 1146–1149.

ee).

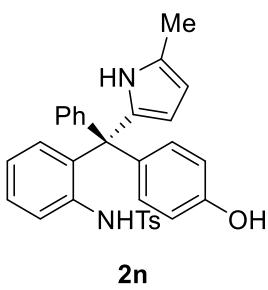
$[\alpha]_D^{22}$: +3.8 ($c = 1.0$, acetone). HPLC analysis of the product: Daicel CHIRALPAK IC column; 30% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 11.7 min (minor), 13.0 min (major).

$^1\text{H NMR}$ (400 MHz, acetone- d_6) δ 9.38 (s, 1H), 8.48 (s, 1H), 7.73 (d, $J = 8.1$ Hz, 1H), 7.35 – 7.25 (m, 4H), 7.10 (d, $J = 7.4$ Hz, 2H), 7.02 – 6.90 (m, 4H), 6.84 (s, 1H), 6.77 (d, $J = 8.6$ Hz, 2H), 6.03 – 6.02 (m, 1H), 5.82 (s, 1H), 2.41 (s, 3H), 2.16 (s, 3H).

$^{13}\text{C NMR}$ (100 MHz, acetone- d_6) δ 157.0, 146.1, 138.8, 135.9, 134.6, 133.0, 132.2, 131.9, 130.9, 130.4, 129.3, 128.7, 127.5, 123.7, 117.9, 115.5, 110.8, 106.4, 58.9, 38.1, 13.0.

IR (thin film) 3433, 3361, 3021, 2926, 2852, 1588, 1265, 1149, 750, 701 cm⁻¹.

HRMS (LD+) Calcd for C₂₅H₂₄N₂O₃S (M⁺): 432.1508, Found: 432.1507.



2n

(R)-N-(2-((4-Hydroxyphenyl)(5-methyl-1*H*-pyrrol-2-yl)(phenyl)methyl)-phenyl)-4-methylbenzene-sulfonamide (2n) was prepared as a light yellow foam from *N*-(2-((4-hydroxyphenyl)(phenyl)methyl)phenyl)-4-methylbenzenesulfonamide (107.3 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure E (eluent: hexanes/EtOAc = 20:1 → 10:1) in 82% yield (105.1 mg, 88% ee).

$[\alpha]_D^{22}$: +2.1 ($c = 1.0$, acetone). HPLC analysis of the product: Daicel CHIRALPAK IC column; 15% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 18.6 min (minor), 20.9 min (major).

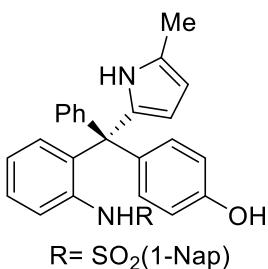
$^1\text{H NMR}$ (400 MHz, acetone- d_6) δ 9.35 (s, 1H), 8.48 (s, 1H), 7.45 – 7.40 (m, 3H), 7.37 – 7.26 (m, 6H), 7.21 – 7.16 (m, 1H), 7.08 – 7.06 (m, 2H), 6.90 – 6.86 (m, 4H),

6.76 – 6.74 (m, 2H), 6.06 (s, 1H), 5.87 (m, 1H), 2.39 (s, 3H), 2.19 (s, 3H).

¹³C NMR (100 MHz, acetone-*d*₆) δ 157.0, 146.0, 144.6, 138.3, 138.2, 136.1, 133.4, 133.0, 132.2, 131.8, 131.0, 130.8, 130.4, 128.8, 128.5, 128.3, 127.4, 122.8, 117.0, 115.4, 111.0, 106.3, 58.9, 21.4, 13.2.

IR (thin film) 3435, 3341, 3054, 2924, 2853, 1590, 1153, 739, 698 cm⁻¹.

HRMS (LD+) Calcd for C₃₁H₂₈N₂O₃S (M⁺): 508.1821, Found: 508.1802.



(R)-N-(2-((4-Hydroxyphenyl)(phenyl)methyl)phenyl)-naphthalene-1-sulfonamide (2o) was prepared as a light yellow foam from *N*-(2-((4-hydroxyphenyl)(phenyl)methyl)phenyl)-naphthalene-1-sulfonamide (116.3mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure E (eluent: hexanes/EtOAc = 20:1 → 10:1) in 49% yield (66.9 mg, 98% ee).

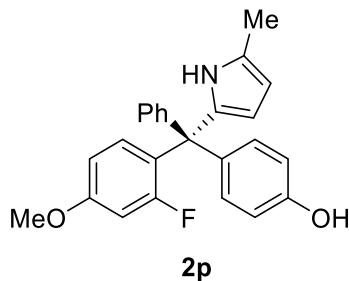
$[\alpha]_D^{22}$: +1.4 (*c* = 1.0, acetone). HPLC analysis of the product: Daicel CHIRALPAK IC column; 30% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 12.3 min (major), 15.0 min (minor).

¹H NMR (400 MHz, acetone-*d*₆) δ 9.44 (s, 1H), 8.48 (s, 1H), 8.24 – 8.21 (m, 2H), 8.05 – 7.99 (m, 2H), 7.69 – 7.62 (m, 3H), 7.45 (s, 1H), 7.29 – 7.25 (m, 3H), 7.14 – 7.01 (m, 4H), 6.95 – 6.87 (m, 4H), 6.76 (d, *J* = 8.6 Hz, 2H), 6.14 – 6.13 (m, 1H), 5.98 – 5.93 (m, 1H), 2.12 (s, 3H).

¹³C NMR (100 MHz, acetone-*d*₆) δ 157.0, 146.0, 139.0, 138.3, 136.3, 135.3, 135.2, 134.5, 133.3, 132.3, 131.8, 131.2, 130.9, 129.7, 129.3, 129.1, 128.8, 128.7, 128.5, 127.8, 127.5, 125.39, 125.30, 123.3, 119.0, 115.4, 111.0, 106.4, 59.0, 13.2.

IR (thin film) 3435, 3376, 3057, 2925, 2851, 1589, 1264, 743, 700 cm^{-1} .

HRMS (LD+) Calcd for $\text{C}_{34}\text{H}_{28}\text{N}_2\text{O}_3\text{S} (\text{M}^+)$: 544.1821, Found: 544.1840.



2p

(R)-4-((2-Fluoro-4-methoxyphenyl)(5-methyl-1*H*-pyrrol-2-yl)(phenyl)methyl)phenol (2p) was prepared as a light yellow foam from 4-((2-fluoro-4-methoxyphenyl)(phenyl)methyl)phenol (77.0 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/EtOAc = 20:1 → 10:1) in 94% yield (90.6 mg, 94% ee).

$[\alpha]_D^{22}$: +1.4 ($c = 1.0$, CHCl_3). HPLC analysis of the product: Daicel CHIRALCEL OD-H column; 5% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 19.5 min (major), 24.4 min (minor).

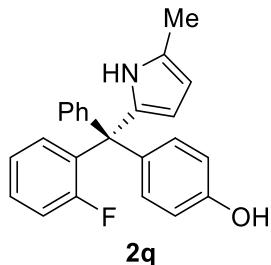
$^1\text{H NMR}$ (400 MHz, CDCl_3) δ 7.37 (s, 1H), 7.29 – 7.15 (m, 6H), 7.01 (d, $J = 8.5$ Hz, 2H), 6.89 – 6.84 (m, 1H), 6.73 (d, $J = 8.7$ Hz, 2H), 6.60 – 6.55 (m, 2H), 5.79 – 5.77 (m, 2H), 3.79 (s, 3H), 2.18 (s, 3H).

$^{13}\text{C NMR}$ (100 MHz, CDCl_3) δ 161.4 (d, $J = 247.8$ Hz), 160.3 (d, $J = 11.0$ Hz), 154.2, 145.1, 137.0, 134.6, 131.2, 131.0 (d, $J = 5.5$ Hz), 129.7, 127.8, 127.4, 127.0 (d, $J = 12.1$ Hz), 126.6, 114.6, 110.3, 108.9 (d, $J = 2.7$ Hz), 105.5, 102.9 (d, $J = 26.5$ Hz), 57.2, 55.7, 13.4.

$^{19}\text{F NMR}$ (376 MHz, CDCl_3) δ -99.3.

IR (thin film) 3446, 3054, 2932, 1616, 1261, 824, 733 cm^{-1} .

HRMS (CI+) Calcd for $\text{C}_{25}\text{H}_{22}\text{FNO}_2 (\text{M}^+)$: 387.1635, Found: 387.1624.



(R)-4-((2-Fluorophenyl)(5-methyl-1H-pyrrol-2-yl)(phenyl)methyl)phenol (2q)

was prepared as a light yellow foam from 4-((2-fluorophenyl)(phenyl)methyl)phenol (69.5 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure E (eluent: hexanes/EtOAc = 20:1 → 10:1) in 74% yield (66.2 mg, 83% ee).

$[\alpha]_D^{22}$: +6.9 ($c = 1.0$, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK AD-H column; 5% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 12.4 min (major), 13.4 min (minor).

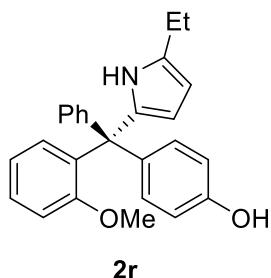
¹H NMR (400 MHz, CDCl₃) δ 7.38 (s, 1H), 7.32 – 7.17 (m, 6H), 7.09 – 6.97 (m, 5H), 6.75 – 6.71 (m, 2H), 5.81 – 5.79 (m, 2H), 4.95 (s, 1H), 2.19 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 161.0 (d, $J = 248.0$ Hz), 154.2, 144.7, 136.7, 134.9 (d, $J = 11.5$ Hz), 134.3, 131.3, 130.6 (d, $J = 3.4$ Hz), 129.8, 129.3 (d, $J = 8.4$ Hz), 127.9, 127.6, 126.7, 124.0 (d, $J = 3.0$ Hz), 116.6 (d, $J = 22.8$ Hz), 114.7, 110.5, 105.5, 57.8, 13.4.

¹⁹F NMR (376 MHz, CDCl₃) δ -101.3.

IR (thin film) 3440, 3057, 2854, 1586, 1484, 1443, 1262, 1220, 1177, 753 cm⁻¹.

HRMS (CI+) Calcd for C₂₄H₂₀FNO (M⁺): 357.1529, Found: 357.1527.



(R)-4-((5-Ethyl-1H-pyrrol-2-yl)(2-methoxyphenyl)(phenyl)methyl)phenol (2r)

was prepared as a light yellow foam from 4-((2-methoxyphenyl)(phenyl)methyl)phenol (72.5 mg, 0.25 mmol) and 2-ethylpyrrole (47.5 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/EtOAc = 20:1 → 10:1) in 98% yield (94.1 mg, 81% ee).

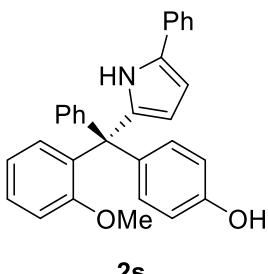
$[\alpha]_D^{22}$: -10.6 ($c = 1.0$, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK IC column; 3% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 9.3 min (major), 10.1 min (minor).

¹H NMR (400 MHz, CDCl₃) δ 7.65 (s, 1H), 7.31 – 7.19 (m, 6H), 7.03 – 7.00 (m, 2H), 6.97 – 6.94 (m, 1H), 6.92 – 6.88 (m, 2H), 6.70 – 6.68 (m, 2H), 5.81 – 5.79 (m, 1H), 5.74 – 5.72 (m, 1H), 3.21 (s, 3H), 2.52 (q, $J = 7.5$ Hz, 2H), 1.16 (t, $J = 7.5$ Hz, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 158.1, 153.7, 145.8, 137.5, 136.8, 135.7, 133.8, 131.6, 130.4, 129.9, 128.7, 127.4, 126.1, 120.9, 114.2, 113.8, 110.2, 103.3, 58.6, 55.8, 21.1, 13.8.

IR (thin film) 3444, 3384, 3055, 2836, 1587, 1241, 1117, 1024, 734, 697 cm⁻¹.

HRMS (LD+) Calcd for C₂₆H₂₅NO₂ (M⁺): 383.1885, Found: 383.1898.



(R)-4-((2-Methoxyphenyl)(phenyl)(5-phenyl-1*H*-pyrrol-2-yl)methyl)phenol

(2s) was prepared as a colorless foam from 4-((2-methoxyphenyl)(phenyl)methyl)phenol (72.5 mg, 0.25 mmol) and 2-phenylpyrrole (71.5 mg, 0.5 mmol) according to the General Procedure E (eluent: hexanes/EtOAc = 20:1 → 10:1) in 94% yield (101.5 mg, 82% ee).

$[\alpha]_D^{22}$: +2.7 ($c = 1.0$, CHCl₃). HPLC analysis of the product: Daicel CHIRALCEL OD-H column; 5% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 11.2 min

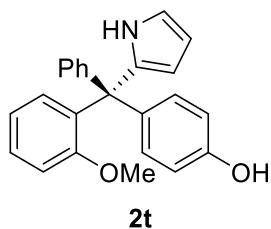
(minor), 12.2 min (major).

¹H NMR (400 MHz, acetone-*d*₆) δ 9.48 (s, 1H), 8.26 (s, 1H), 7.48 (d, *J* = 7.7 Hz, 2H), 7.31 (t, *J* = 6.9 Hz, 1H), 7.27 – 7.16 (m, 7H), 7.10 – 7.06 (m, 2H), 7.02 – 6.96 (m, 3H), 6.90 (t, *J* = 7.5 Hz, 1H), 6.74 (d, *J* = 8.7 Hz, 2H), 6.43 (t, *J* = 3.1 Hz, 1H), 5.82 (t, *J* = 3.0 Hz, 1H), 3.26 (s, 3H).

¹³C NMR (100 MHz, acetone-*d*₆) δ 158.9, 156.4, 146.8, 139.9, 137.0, 136.5, 134.2, 132.4, 132.2, 131.1, 130.6, 129.4, 129.3, 127.8, 126.5, 126.2, 124.4, 121.3, 114.7, 114.4, 112.6, 105.8, 59.4, 55.7.

IR (thin film) 3449, 3366, 3054, 2836, 1600, 1254, 732, 697 cm⁻¹.

HRMS (LD+) Calcd for C₃₀H₂₅NO₂ (M⁺): 431.1885, Found: 431.1876.



(R)-4-((2-Methoxyphenyl)(phenyl)(1*H*-pyrrol-2-yl)methyl)phenol (2t) was prepared as a light yellow foam from 4-((2-methoxyphenyl)(phenyl)methyl)phenol (72.5 mg, 0.25 mmol) and pyrrole (83.75 mg, 1.25 mmol, 5 equiv) according to the General Procedure E (eluent: hexanes/EtOAc = 20:1 → 10:1) in 99% yield (89.0 mg, 99% ee).

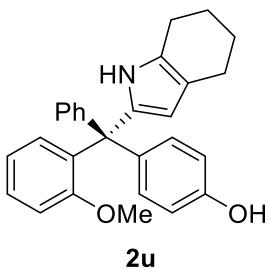
[α]_D²²: +1.1 (*c* = 1.0, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK IC column; 2% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 13.6 min (major >99%).

¹H NMR (400 MHz, acetone-*d*₆) δ 9.04 (s, 1H), 8.27 (s, 1H), 7.30 – 7.14 (m, 6H), 6.96 – 6.85 (m, 5H), 6.74 – 6.69 (m, 3H), 6.01 – 5.99 (m, 1H), 5.78 (s, 1H), 3.19 (s, 3H).

¹³C NMR (100 MHz, acetone-*d*₆) δ 158.8, 156.1, 146.9, 137.8, 137.4, 136.7, 132.0, 130.7, 130.4, 129.2, 127.6, 126.3, 121.0, 118.0, 114.5, 114.0, 110.1, 107.3, 59.0, 55.5.

IR (thin film) 3447, 3055, 2835, 1598, 1241, 1176, 731 cm⁻¹.

HRMS (LD+) Calcd for C₂₄H₂₁NO₂ (M⁺): 355.1572, Found: 355.1556.



(R)-4-((2-Methoxyphenyl)(phenyl)(4,5,6,7-tetrahydro-1*H*-indol-2-yl)methyl)phenol (3u) was prepared as a colorless foam from 4-((2-methoxyphenyl)(phenyl)methyl)phenol (72.5 mg, 0.25 mmol) and 4,5,6,7-tetrahydro-1*H*-indole (60.5 mg, 0.5 mmol) according to the General Procedure E (eluent: hexanes/EtOAc = 15:1 → 10:1) in 95% yield (97.1 mg, 95% ee).

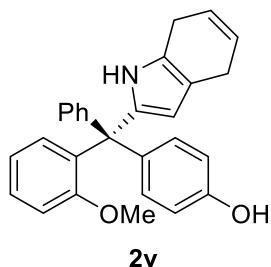
[α]_D²²: -0.5 (c = 1.0, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK IC column; 3% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 9.0 min (major), 10.1 min (minor).

¹H NMR (400 MHz, acetone-*d*₆) δ 8.48 (s, 1H), 8.20 (s, 1H), 7.28 – 7.21 (m, 6H), 7.02 – 6.85 (m, 5H), 6.69 (d, *J* = 8.6 Hz, 2H), 5.44 – 5.43 (m, 1H), 3.16 (s, 3H), 2.45 (t, *J* = 5.7 Hz, 2H), 2.39 (t, *J* = 5.5 Hz, 2H), 1.72 – 1.67 (m, 4H).

¹³C NMR (100 MHz, acetone-*d*₆) δ 158.9, 156.1, 147.0, 137.9, 136.7, 136.1, 132.1, 130.9, 130.5, 129.1, 127.5, 127.1, 126.2, 121.0, 115.5, 114.4, 114.0, 109.7, 59.2, 55.5, 24.8, 24.4, 23.7, 23.4.

IR (thin film) 3447, 3402, 3054, 2844, 1598, 1241, 1174, 733, 700 cm⁻¹.

HRMS (LD+) Calcd for C₂₈H₂₇NO₂ (M⁺): 409.2042, Found: 409.2054.



(R)-4-((4,7-Dihydro-1H-indol-2-yl)(2-methoxyphenyl)(phenyl)methyl)phenol (2v)

was prepared as a colorless foam from 4-((2-methoxyphenyl)(phenyl)methyl)phenol (72.5 mg, 0.25 mmol) and 4,7-dihydro-1*H*-indole (59.5 mg, 0.5 mmol) according to the General Procedure E (eluent: hexanes/EtOAc = 15:1 → 10:1) in 80% yield (81.4 mg, 86% ee).

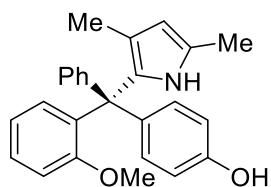
$[\alpha]_D^{22}$: -0.5 ($c = 1.0$, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK IC column; 3% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 9.8 min (major), 11.9 min (minor).

¹H NMR (400 MHz, acetone-*d*₆) δ 8.65 (s, 1H), 8.25 (s, 1H), 7.29 – 7.12 (m, 6H), 7.02 – 6.86 (m, 5H), 6.72 – 6.69 (m, 2H), 5.85 – 5.83 (m, 1H), 5.79 – 5.76 (m, 1H), 5.53 – 5.52 (m, 1H), 3.17 (s, 3H), 3.15 – 3.10 (m, 4H).

¹³C NMR (100 MHz, acetone-*d*₆) δ 158.9, 156.1, 146.9, 137.7, 137.0, 136.6, 132.1, 130.8, 130.5, 129.2, 127.6, 126.5, 126.2, 124.31, 124.27, 121.0, 114.5, 114.0, 112.7, 108.9, 59.2, 55.5, 25.6, 24.7.

IR (thin film) 3447, 3405, 2831, 1696, 1599, 1255, 1175, 1108, 820, 731 cm⁻¹.

HRMS (CI+) Calcd for C₂₈H₂₅NO₂ (M⁺): 407.1885, Found: 407.1888.



4-((3,5-Dimethyl-1H-pyrrol-2-yl)(2-methoxyphenyl)(phenyl)methyl)phenol (2x)

was prepared as a colorless foam from 4-((2-methoxyphenyl)(phenyl)methyl)phenol (72.5 mg, 0.25 mmol) and 2,4-dimethyl-1*H*-pyrrole

(47.6 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/EtOAc = 10:1 → 3:1) in 76% yield (73.2 mg, 94% ee).

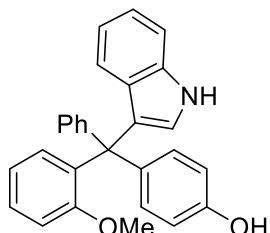
$[\alpha]_D^{22}$: +3.2 (c = 1.0, CH₂Cl₂). HPLC analysis of the product: Daicel CHIRALPAK AD-H column; 2% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 16.0 min (minor), 17.9 min (major).

¹H NMR (400 MHz, CDCl₃) δ 7.49 (s, 1H), 7.34 – 7.30 (m, 1H), 7.22 (d, J = 8.7 Hz, 3H), 7.12 (d, J = 8.6 Hz, 2H), 7.01 (d, J = 8.6 Hz, 2H), 6.96 – 6.90 (m, 2H), 6.70 (d, J = 8.7 Hz, 2H), 5.721–5.715 (m, 1H), 5.17 (br, 1H), 3.22 (s, 3H), 2.18 (s, 3H), 1.30 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 158.1, 153.6, 143.8, 137.1, 134.4, 131.6, 131.5, 131.4, 130.7, 129.6, 128.8, 127.5, 124.0, 120.7, 116.9, 114.4, 113.6, 109.9, 57.9, 55.6, 13.2, 13.0.

IR (thin film) 3451, 1589, 1484, 1432, 1259, 1177, 1099, 824, 734 cm⁻¹.

HRMS (ES-) Calcd for C₂₆H₂₄NO₂⁻ (M-H⁺): 382.1813, Found: 382.1810.



2y

4-((1*H*-indol-3-yl)(2-methoxyphenyl)(phenyl)methyl)phenol (2y) was prepared as a colorless foam from 4-((2-methoxyphenyl)(phenyl)methyl)phenol (72.5 mg, 0.25 mmol) and indole (58.6 mg, 0.5 mmol) according to the General Procedure D (24 h for step 2 at room temperature, eluent: hexanes/EtOAc = 10:1 → 3:1) in 99% yield (101.1 mg, 21% ee).

HPLC analysis of the product: Daicel CHIRALPAK IC column; 3% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 29.9 min (minor), 38.2 min (major).

¹H NMR (400 MHz, acetone-*d*₆) δ 10.02 (s, 1H), 8.15 (s, 1H), 7.36 – 7.34 (m, 1H),

7.29 – 7.10 (m, 7H), 7.04 – 7.01 (m, 2H), 6.98 – 6.93 (m, 2H), 6.82 – 6.78 (m, 2H), 6.70 – 6.62 (m, 4H), 3.15 (s, 3H).

^{13}C NMR (100 MHz, acetone- d_6) δ 159.3, 155.8, 147.4, 138.3, 137.2, 137.0, 132.2, 131.2, 130.8, 129.0, 128.6, 127.5, 126.5, 126.0, 123.8, 122.8, 121.6, 120.7, 119.1, 114.4, 114.1, 112.1, 58.2, 55.4.

IR (thin film) 3411, 3054, 1706, 1594, 1484, 1449, 1334, 1238, 1175, 1104, 821, 735, 703 cm⁻¹.

HRMS (ES+) Calcd for C₂₈H₂₃NO₂ (M⁺): 405.1729, Found: 405.1717.

IV. Mechanistic Study



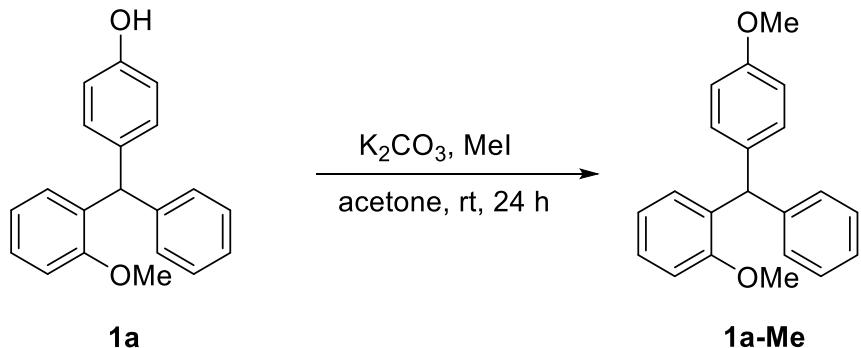
4-((2-Methoxyphenyl)(phenyl)methylene)cyclohexa-2,5-dien-1-one (QM). At room temperature, to a solution of triarylmethane **1a** (72.5 mg, 0.25 mmol) in DCM (5 mL) were added DDQ (57 mg, 0.25 mmol, 1 equiv) and 4 Å molecular sieves (50 mg). Then the mixture was stirred for 5 h and concentrated. The residue was purified by silica gel flash column chromatography to afford the quinone methide as a yellow foam in 49% yield (35.0 mg). The reaction is clean. The low yield was due to instability of this quinone methide.

1H NMR (400 MHz, CDCl₃) δ 7.47 – 7.35 (m, 5H), 7.26 – 7.25 (m, 2H), 7.19 (dd, J₁ = 10 Hz, J₂ = 2.4 Hz, 1H), 7.12 (d, J = 7.6 Hz, 1H), 7.03 (t, J = 7.4 Hz, 1H), 6.95 (d, J = 8.0 Hz, 1H), 6.44 (dd, J₁ = 10 Hz, J₂ = 1.6 Hz, 1H), 6.37 (dd, J₁ = 10 Hz, J₂ = 1.6 Hz, 1H), 3.59 (s, 3H).

13C NMR (100 MHz, CDCl₃) δ 187.6, 158.1, 157.6, 140.1, 139.8, 139.1, 132.6, 131.4, 131.1, 130.4, 129.6, 129.3, 128.9, 128.5, 128.2, 120.6, 112.0, 55.8.

IR (thin film) 3513, 2940, 1620, 1487, 1232, 1165, 1017, 699 cm⁻¹.

HRMS (CI+) Calcd for C₂₀H₁₆O₂ (M⁺): 288.1150, Found: 288.1152.



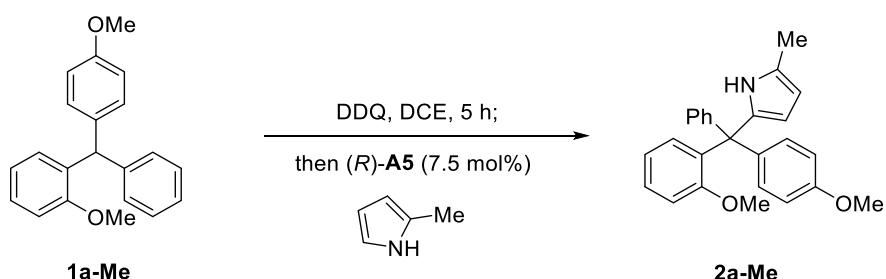
1-Methoxy-2-((4-methoxyphenyl)(phenyl)methyl)benzene (1a-Me). At room temperature, to a solution of **1a** (290 mg, 1 mmol) in acetone (10 mL) were sequentially added MeI (0.3 mL, 3 mmol) and K₂CO₃ (649 mg, 5 mmol). The reaction mixture was stirred for 24 h before it was filtered and concentrated. The residue was purified by silica gel flash chromatography to afford the desired product (eluent: hexanes/EtOAc = 30:1 → 20:1) as white foam in 69% yield (211 mg).

¹H NMR (400 MHz, CDCl₃) δ 7.28 – 7.19 (m, 4H), 7.09 (d, *J* = 7.2 Hz, 2H), 7.02 – 7.00 (m, 2H), 6.89 – 6.80 (m, 5H), 5.88 (s, 1H), 3.79 (s, 3H), 3.72 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 158.0, 157.2, 144.5, 136.2, 133.1, 130.5, 130.4, 129.5, 128.3, 127.6, 126.1, 120.4, 113.7, 110.8, 55.8, 55.3, 48.9.

IR (thin film) 3027, 2835, 1596, 1333, 1293, 1240, 1029, 734, 701 cm⁻¹.

HRMS (CI+) Calcd for C₂₁H₂₀O₂ (M⁺): 304.1463, Found: 304.1469.



2-((2-Methoxyphenyl)(4-methoxyphenyl)(phenyl)methyl)-5-methyl-1*H*-pyrrole (2a-Me) was prepared as a colorless foam from 1-methoxy-2-((4-methoxyphenyl)(phenyl)methyl)benzene (72.5 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure D

(eluent: hexanes/DCM = 30:1 → 10:1) in 58% yield (55.6 mg).

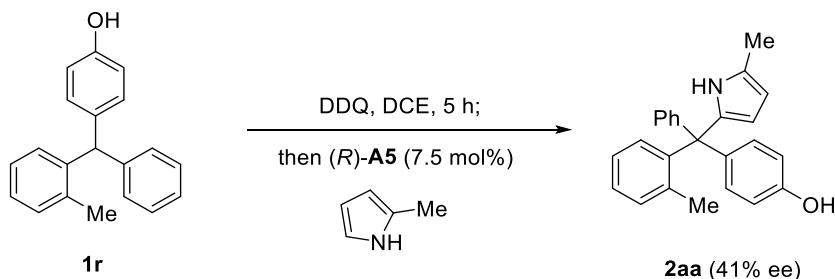
HPLC analysis of the product: Daicel CHIRALCEL OD-H column; 0.7% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 9.0 min, 11.6 min.

1H NMR (400 MHz, CDCl₃) δ 7.59 (s, 1H), 7.31 – 7.15 (m, 6H), 7.10 – 7.06 (m, 2H), 6.98 – 6.96 (m, 1H), 6.92 – 6.88 (m, 2H), 6.81 – 6.77 (m, 2H), 5.78 – 5.76 (m, 1H), 5.72 – 5.70 (m, 1H), 3.80 (s, 3H), 3.21 (s, 3H), 2.18 (s, 3H).

13C NMR (100 MHz, CDCl₃) δ 158.1, 157.8, 145.8, 137.3, 136.9, 135.9, 131.4, 130.3, 129.9, 128.7, 127.3, 127.0, 125.9, 120.8, 113.8, 112.6, 110.4, 105.2, 58.5, 55.8, 55.4, 13.4.

IR (thin film) 3460, 239, 1589, 1501, 1246, 1180, 1113, 1031, 907, 729 cm⁻¹.

HRMS (CI+) Calcd for C₂₆H₂₅NO₂ (M⁺): 383.1885, Found: 383.1880.



4-((5-Methyl-1*H*-pyrrol-2-yl)(phenyl)(*o*-tolyl)methyl)phenol (2aa) was prepared as a yellow foam from 4-(phenyl(*o*-tolyl)methyl)phenol (68.5 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/DCM = 30:1 → 10:1) in 99% yield (87 mg, 41% ee).

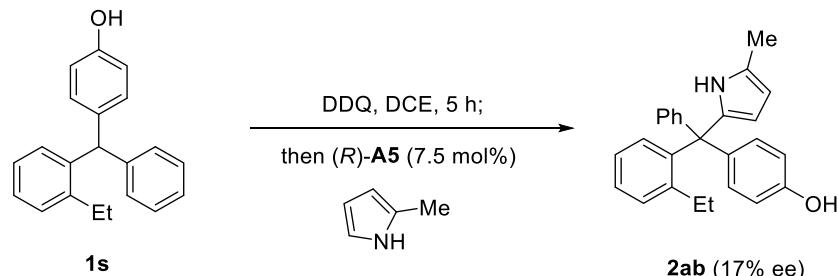
HPLC analysis of the product: Daicel CHIRALPAK AD-H column; 2% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 20.5 min (minor), 24.6 min (major).

1H NMR (400 MHz, CDCl₃) δ 7.27 – 7.08 (m, 9H), 6.98 – 6.94 (m, 2H), 6.84 – 6.82 (m, 1H), 6.69 – 6.65 (m, 2H), 5.83 – 5.80 (m, 2H), 5.03 (br, 1H), 2.19 (s, 3H), 1.66 (s, 3H).

13C NMR (100 MHz, CDCl₃) δ 153.9, 145.8, 145.3, 139.5, 137.9, 135.5, 132.4, 131.8, 130.3, 130.0, 127.8, 127.23, 127.22, 126.4, 125.7, 114.6, 110.2, 105.4, 59.8, 22.7, 13.4.

IR (thin film) 3441, 3056, 2921, 2857, 1590, 1504, 1441, 1259, 1176, 1040, 824, 740, 700 cm⁻¹.

HRMS (Cl+) Calcd for C₂₅H₂₃NO₂ (M⁺): 353.1780, Found: 353.1744.



4-((2-Ethylphenyl)(5-methyl-1H-pyrrol-2-yl)(phenyl)methyl)phenol (2ab)

was prepared as a yellow foam from 4-((2-ethylphenyl)(phenyl)methyl)phenol (72.0 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/DCM = 10:1 → 5:1) in 99% yield (90.5 mg, 17% ee).

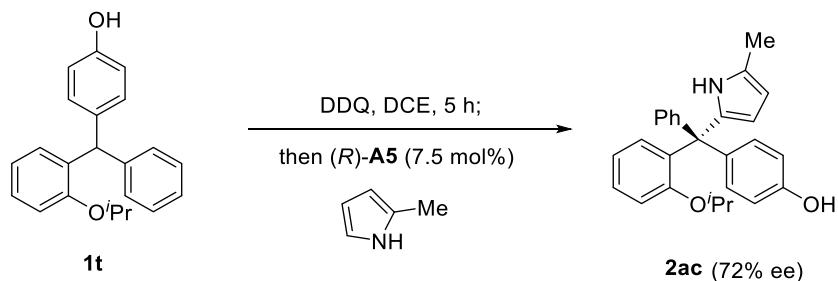
HPLC analysis of the product: Daicel CHIRALPAK AD-H column; 3% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 14.7 min (minor), 16.0 min (major).

1H NMR (400 MHz, CDCl₃) δ 7.33 – 7.24 (m, 8H), 7.17 – 7.07 (m, 3H), 6.91 (d, *J* = 7.9 Hz, 1H), 6.73 (d, *J* = 8.6 Hz, 2H), 5.90 – 5.86 (m, 2H), 5.28 (br, 1H), 2.28 – 2.21 (m, 5H), 0.52 (t, *J* = 7.4 Hz, 3H).

13C NMR (100 MHz, CDCl₃) δ 153.8, 146.0, 145.0, 144.6, 138.0, 136.0, 131.8, 130.45, 130.37, 129.6, 127.7, 127.4, 127.3, 126.3, 125.5, 114.5, 109.9, 105.4, 59.6, 27.8, 14.2, 13.3.

IR (thin film) 3446, 2361, 1601, 1506, 1438, 1263, 1176, 825, 734 cm⁻¹.

HRMS (ES-) Calcd for C₂₆H₂₄NO⁻ (M-H⁺): 366.1863, Found: 366.1861.



4-((2-Isopropoxypyhenyl)(5-methyl-1H-pyrrol-2-yl)(phenyl)methyl)phenol

(2ac) was prepared as a white foam from 4-((2-isopropoxypyhenyl)(phenyl)methyl)phenol (79.5 mg, 0.25 mmol) and 2-methylpyrrole (40.6 mg, 0.5 mmol) according to the General Procedure D (eluent: hexanes/DCM = 10:1 → 5:1) in 75% yield (74 mg, 72% ee).

HPLC analysis of the product: Daicel CHIRALPAK IC column; 2% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 9.2 min (major), 10.4 min (minor).

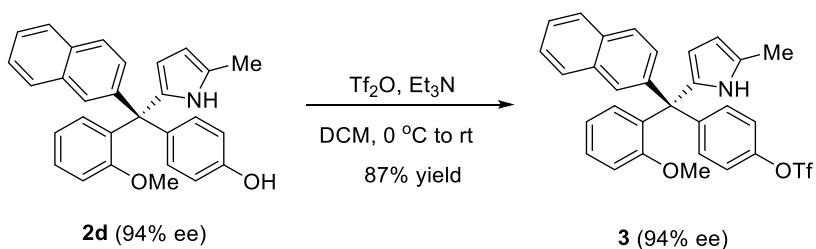
1H NMR (400 MHz, CDCl₃) δ 7.88 (s, 1H), 7.30 – 7.18 (m, 6H), 7.06 – 7.02 (m, 3H), 6.87 – 6.82 (m, 2H), 6.69 (d, *J* = 8.7 Hz, 2H), 5.81 – 5.79 (m, 1H), 5.72 – 5.70 (m, 1H), 5.06 (br, 1H), 4.40 – 4.34 (m, 1H), 2.21 (s, 3H), 0.80 – 0.77 (m, 6H).

13C NMR (100 MHz, CDCl₃) δ 155.2, 153.6, 145.7, 137.3, 136.2, 135.9, 131.9, 130.9, 130.2, 128.4, 127.1, 126.8, 125.8, 119.4, 113.9, 112.1, 110.6, 104.9, 68.3, 58.6, 21.13, 21.07, 13.4.

IR (thin film) 3446, 3055, 2978, 2928, 1589, 1508, 1443, 1254, 1176, 1116, 732, 699 cm⁻¹.

HRMS (ES-) Calcd for C₂₇H₂₆NO₂⁻ (M-H⁺): 396.1969, Found: 396.1968.

V. Product Derivatizations



(S)-4-((2-Methoxyphenyl)(5-methyl-1*H*-pyrrol-2-yl)(naphthalen-2-yl)methyl)phenyl trifluoromethanesulfonate (3). At 0 °C, to a solution of **2d** (210 mg, 0.5 mmol) and Et₃N (0.16 mL, 1.0 mmol) in DCM (12 mL) was slowly added Tf₂O (0.1 mL, 0.6 mmol) over 1 min. The reaction mixture was stirred at 0 °C for 30 min and then at room temperature for 1 h. Next, water (8 mL) was added, and the reaction mixture was extracted with DCM (16 mL × 2). The combined organic layers were dried over anhydrous Na₂SO₄ and concentrated. The residue was purified by silica gel flash chromatography to afford the desired product (eluent: hexanes/EtOAc = 30:1 → 20:1) as a colorless foam in 87% yield (239 mg, 94% ee).

$[\alpha]_D^{22}$: -7.4 ($c = 1.0$, CHCl_3). HPLC analysis of the product: Daicel CHIRALPAK AD-H column; 1% *i*-PrOH in hexanes; 0.6 mL/min; retention times: 5.5 min (major), 6.4 min (minor).

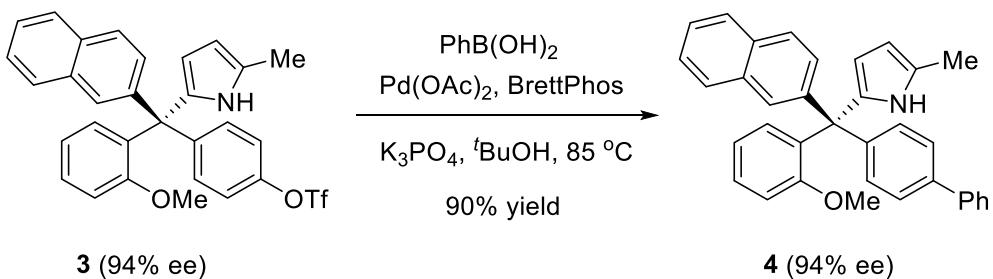
¹H NMR (400 MHz, CDCl₃) δ 7.84 (d, *J* = 7.8 Hz, 1H), 7.76 – 7.70 (m, 2H), 7.55 – 7.43 (m, 4H), 7.38 – 7.31 (m, 4H), 7.15 (d, *J* = 8.8 Hz, 2H), 7.05 (d, *J* = 7.2 Hz, 1H), 6.97 – 6.91 (m, 2H), 5.80 (s, 1H), 5.75 – 5.74 (s, 1H), 3.20 (s, 3H), 2.17 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 157.8, 147.7, 146.5, 141.4, 135.6, 134.4, 133.1, 132.3, 131.6, 130.3, 129.6, 129.3, 128.5, 128.3, 127.9, 127.5, 126.7, 126.2, 126.1, 121.1, 120.0, 119.0 (q, *J* = 332 Hz), 113.6, 110.9, 105.5, 58.9, 55.4, 13.4.

¹⁹F NMR (376 MHz, CDCl₃) δ -72.9.

IR (thin film) 3442, 3057, 1590, 1491, 1420, 1212, 1139, 1021, 889, 756, 571 cm⁻¹.

HRMS (CI+) Calcd for C₃₀H₂₄F₃NO₄S (M⁺): 551.1378, Found: 551.1354.



(S)-2-([1,1'-Biphenyl]-4-yl(2-methoxyphenyl)(naphthalen-2-yl)methyl)-5-methyl-1*H*-pyrrole (4) was prepared by a modified literature procedure.² Under N₂, to an oven-dried 4-mL vial were added triflate **3a** (110.2 mg, 0.2 mmol), phenylboronic acid (73.2 mg, 0.6 mmol), Pd(OAc)₂ (1.1 mg, 10 µmol, 5 mol%), BrettPhos (5.7 mg, 20 µmol, 10 mol%), K₃PO₄ (127.2 mg, 0.6 mmol, 3.0 equiv), and *tert*-butanol (2 mL). The mixture was stirring at 85 °C for 12 h. After that, the mixture was cooled to room temperature, concentrated and then purified by silica gel flash chromatography (eluent: hexanes/EtOAc = 40:1 → 30:1) to afford the desired product as a colorless foam in 90% yield (86.3 mg, 94% ee).

$[\alpha]_D^{22}$: +4.4 (*c* = 1.0, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK AD-H column; 1% *i*-PrOH in hexanes; 0.6 mL/min; retention times: 7.1 min (minor), 8.8 min (major).

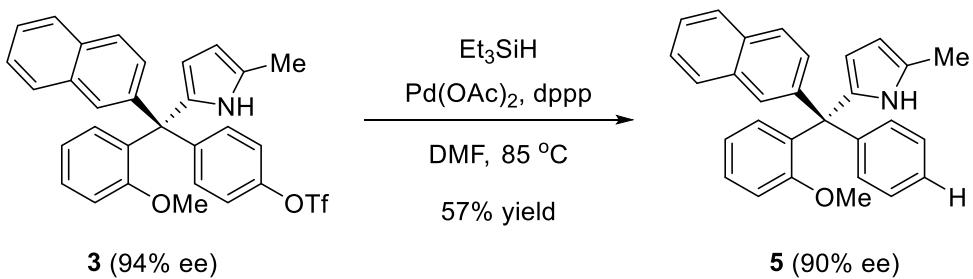
¹H NMR (400 MHz, CDCl₃) δ 7.82 (d, *J* = 7.8 Hz, 1H), 7.32 (d, *J* = 7.8 Hz, 2H), 7.64 – 7.62 (m, 4H), 7.52 – 7.28 (m, 11H), 7.08 – 7.07 (m, 1H), 6.96 – 6.91 (m, 2H), 5.81 (s, 2H), 3.19 (s, 3H), 2.18 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 158.2, 144.5, 142.7, 141.0, 138.6, 136.3, 135.2, 133.2, 132.2, 130.6, 130.5, 129.7, 128.94, 128.92, 128.6, 128.0, 127.5, 127.32, 127.30, 127.1, 126.5, 126.0, 125.9, 125.8, 120.9, 113.8, 110.7, 105.4, 59.0, 55.7, 13.5.

IR (thin film) 3446, 1589, 1512, 1255, 1111, 1021, 735 cm⁻¹.

HRMS (CI+) Calcd for C₃₅H₂₉NO (M⁺): 479.2249, Found: 479.2256.

(2) B. Bhayana, B. P. Fors and S. L. Buchwald, *Org. Lett.*, 2009, **11**, 3954–3957.



(S)-2-(2-Methoxyphenyl)(naphthalen-2-yl)(phenyl)methyl-5-methyl-1*H*-pyrrole (5) was prepared by a modified literature procedure.³ Under N₂, an oven-dried 4-mL vial equipped with magnetic stir bar was charged with triflate **3a** (110.2 mg, 0.2 mmol), Pd(OAc)₂ (1.1 mg, 5 µmol, 2.5 mol%), dppp (2.1 mg, 5 µmol, 2.5 mol%), and DMF (1 mL). The vial was sealed and kept stirring at 60 °C for 15 min. Then Et₃SiH (80 µL, 0.5 mmol, 2.5 equiv) was added and kept stirring for 48 h. After that, the mixture was cooled to room temperature and Et₂O was added. The solution was washed by water, saturated aqueous NaHCO₃ solution and NaCl solution, respectively. Next, the organic layer was dried over Na₂SO₄ and concentrated. The residue was purified by silica gel flash chromatography to afford the desired product (eluent: hexanes → hexanes/EtOAc = 50:1) as a colorless foam in 57% yield (45.5 mg, 90% ee).

[α]_D²²: +13.8 (*c* = 1.0, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK AD-H column; 0.7% *i*-PrOH in hexanes; 0.6 mL/min; retention times: 8.0 min (minor), 9.1 min (major).

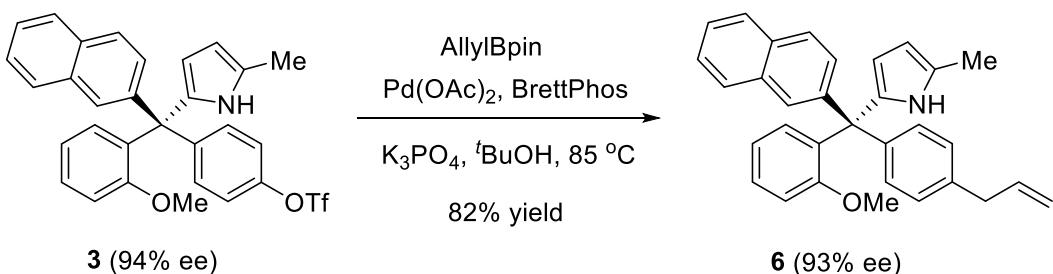
¹H NMR (400 MHz, CDCl₃) δ 7.85 (d, *J* = 7.6 Hz, 1H), 7.75 (d, *J* = 8.4 Hz, 2H), 7.63 (s, 2H), 7.53 – 7.23 (m, 9H), 7.10 (d, *J* = 7.6 Hz, 1H), 6.99 – 6.93 (m, 2H), 5.84 (s, 2H), 3.20 (s, 3H), 2.20 (s, 3H).

¹³C NMR (100 MHz, CDCl₃) δ 158.2, 145.3, 142.6, 136.3, 135.2, 133.1, 132.1, 130.5, 130.1, 129.8, 128.8, 128.5, 128.0, 127.42, 127.35, 127.2, 126.3, 126.1, 125.8, 125.7, 120.8, 113.7, 110.6, 105.3, 59.2, 55.6, 13.4.

(3) G. L. Larson and J. L. Fry, *Org. React.*, 2008, 1–737.

IR (thin film) 3447, 3055, 2928, 1589, 1487, 1247, 1024, 700 cm^{-1} .

HRMS (CI+) Calcd for $\text{C}_{29}\text{H}_{25}\text{NO}$ (M^+): 403.1936, Found: 403.1949.



(S)-2-([1,1'-Biphenyl]-4-yl(2-methoxyphenyl)(naphthalen-2-yl)methyl)-5-methyl-1*H*-pyrrole (6) was prepared by a modified literature procedure.² Under N_2 , to an oven-dried 4-mL vial were added triflate **3a** (110.2 mg, 0.2 mmol), allylboronic acid pinacol ester (34.0 mg, 0.4 mmol), $\text{Pd}(\text{OAc})_2$ (1.1 mg, 5 μmol , 2.5 mol%), BrettPhos (5.3 mg, 10 μmol , 5 mol%), K_3PO_4 (127.2 mg, 0.6 mmol, 3.0 equiv), and *tert*-butanol (2 mL). The mixture was stirring at 85 $^\circ\text{C}$ for 12 h. After that, the mixture was cooled to room temperature, concentrated and then purified by silica gel flash chromatography (eluent: hexanes/EtOAc = 40:1 \rightarrow 30:1) to afford the desired product as a colorless foam in 82% yield (72.5 mg, 93% ee).

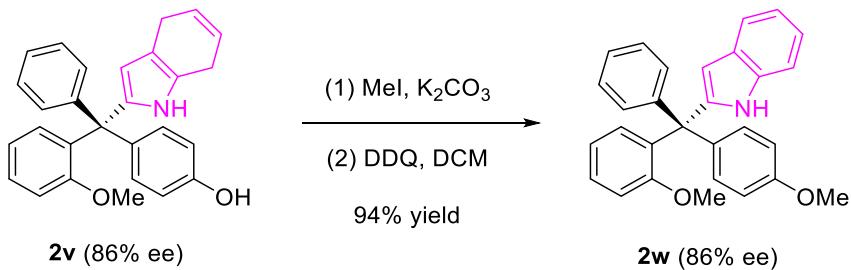
$[\alpha]_{\text{D}}^{22}: +14.7$ ($c = 1.0$, CHCl_3). HPLC analysis of the product: Daicel CHIRALPAK AD-H column; 0.7% *i*-PrOH in hexanes; 0.6 mL/min; retention times: 9.9 min (major), 10.9 min (minor).

$^1\text{H NMR}$ (400 MHz, CDCl_3) δ 7.83 (d, $J = 7.6$ Hz, 1H), 7.73 – 7.71 (m, 2H), 7.62 – 7.61 (m, 2H), 7.48 – 7.41 (m, 2H), 7.37 – 7.31 (m, 2H), 7.19 – 7.16 (m, 2H), 7.11 – 7.06 (m, 3H), 6.97 – 6.91 (m, 2H), 6.08 – 5.98 (m, 1H), 5.81 – 5.79 (m, 2H), 5.13 – 5.08 (m, 2H), 3.41 (d, $J = 6.7$ Hz, 2H), 3.19 (s, 3H), 2.19 (s, 3H).

$^{13}\text{C NMR}$ (100 MHz, CDCl_3) δ 158.2, 143.0, 142.9, 137.8, 137.7, 136.5, 135.3, 133.2, 132.1, 130.5, 130.3, 129.6, 128.8, 128.5, 127.9, 127.6, 127.5, 127.1, 125.77, 125.73, 120.9, 115.81, 115.79, 113.7, 110.6, 105.3, 59.0, 55.7, 40.0, 13.4.

IR (thin film) 3447, 1636, 1588, 1490, 1253, 1113, 1025, 818, 739 cm^{-1} .

HRMS (Cl⁺) Calcd for C₃₂H₂₉NO (M⁺): 443.2249, Found: 443.2267.



(R)-2-((2-Methoxyphenyl)(4-methoxyphenyl)(phenyl)methyl)-1*H*-indole

(2w).⁴ At room temperature, to a solution of **2v** (50.0 mg, 0.12 mmol) in acetone (4 mL) were sequentially added MeI (52.7 mg, 0.37 mmol) and K₂CO₃ (85.1 mg, 0.61 mmol). The reaction mixture was stirred for 24 h and filtered through a short pad of silica gel, which was washed with Et₂O. The filtrate was concentrated under reduced pressure. The crude product was dissolved in anhydrous DCM (8 mL). Next, DDQ (31.0 mg, 0.143 mmol) was added in one portion. The reaction mixture was stirred at room temperature for 3 h and diluted with DCM (15 mL). Then it was washed with an aqueous solution of NaOH (15 mL, 10 wt%) and water (15 mL × 2), dried over Na₂SO₄, and filtered. The filtrate was concentrated. The residue was purified by silica gel flash chromatography to afford the desired product (eluent: hexanes/EtOAc = 15:1 → 10:1) in 96% yield for two steps (48.5 mg, 86% ee).

[α]_D²²: -0.3 (c = 1.0, CHCl₃). HPLC analysis of the product: Daicel CHIRALPAK AD-H column; 3% *i*-PrOH in hexanes; 1.0 mL/min; retention times: 10.5 min (minor), 11.3 min (major).

¹H NMR (400 MHz, acetone-*d*₆) δ 9.46 (s, 1H), 7.46 (d, *J* = 7.7 Hz, 1H), 7.34 – 7.31 (m, 1H), 7.28 – 7.20 (m, 6H), 7.09 (d, *J* = 8.7 Hz, 2H), 7.01 – 6.90 (m, 5H), 6.82 (d, *J* = 8.8 Hz, 2H), 6.19 (s, 1H), 3.78 (s, 3H), 3.20 (s, 3H).

¹³C NMR (100 MHz, acetone-*d*₆) δ 159.0, 158.8, 146.0, 145.5, 137.7, 137.2, 136.1,

(4) H. Çavdar and N. Saracoğlu, *Tetrahedron*, 2005, **61**, 2401–2405.

132.2, 130.81, 130.75, 129.7, 128.7, 127.9, 126.7, 121.7, 121.2, 120.7, 119.8, 114.0,
113.2, 111.8, 104.3, 59.6, 55.5, 55.4.

IR (thin film) 3448, 3053, 2836, 1592, 1290, 1105, 799, 731 cm⁻¹.

HRMS (CI+) Calcd for C₂₉H₂₅NO₂ (M)⁺: 419.1885, Found: 419.1886.

VI. Biological Study

Virus culture:⁵

Human rhabdomyosarcoma (RD) cells (ATCC #CCL-136) were cultured in Dulbecco's modified Eagle's medium (DMEM) with 10% fetal bovine serum (FBS), supplemented with 100 U/mL of penicillin and streptomycin. The cells were incubated at 37 °C in a tissue culture incubator supplied with 5% CO₂. Cells were cultured in 10-cm dishes and infected with enterovirus A71 (EV-A71, SHZ98 strain, GenBank accession number AF302996.1) at multiplicity of infection (MOI) of 1. EV-A71 virus was propagated on a 90% confluent cell monolayer in DMEM with 2% FBS. After incubation for 24 hours, the culture media were collected for preparation of virus stocks. The virus titre was determined by 50% tissue culture infective dose (TCID₅₀) assay or plaque assays (pfu). The virus stocks were aliquoted and stored in -80 °C until use.

Cytotoxicity assay:^{6,7}

The cytotoxicity of compounds was measured by determining the cell viability using MTT [3-(4,5-dimethylthiazol-2-yl)-2,5-diphenyltetrazolium bromide] assays. Briefly, RD cells were set in 96-well plates (10⁴ cells per well) overnight. The testing compound diluted in a 10-fold series was added in the culture media and incubated for another 48 hours. The media were then replaced with 0.5 mg/mL MTT medium and incubated for another 4 hours. The MTT solution was removed from the wells and the formazan crystals were dissolved in DMSO. Absorbances of the formazan products were measured at 550 nm with

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- (5) L. Lu, L. Yi, J. Zhao., J. Yu, Y. Chen, M. C. Lin, H. F. Kung and M. L. He, *J. Virol.*, 2012, **86**, 3767–3776.
 - (6) Q. Dong, R. Men, X. Dan, Y. Chen, H. Li, G. Chen, B. Zee and M. H. T. Wang, M. L. He, *Antiviral Res.*, 2018, **150**, 39–46.
 - (7) F. Zhou, Q. Wan, J. Lu, Y. Chen, G. Lu and M. L. He, *iScience*, 2019, **19**, 715–727.

the reference wavelength at 690 nm. The cell viability assay was carried out in triplicates by three independent experiments. The CC₅₀ (50% cytotoxic concentration) value was calculated.

Cytopathic effects (CPE) assay:^{6,7}

RD cells were set in 24-well plates and incubated overnight. The culture media were replaced with DMEM containing 2% FBS, and then infected with nil (mock) or EV-A71 at MOI of 0.1. After adsorption for 1 hour, 0.4 mL of DMEM medium containing 10% FBS was added to each well. The cells were cultured for 24 hours. The cell morphology was monitored and recorded using a phase-contrast microscope associated with a CCD camera and computer at different time points.

Viral RNA isolation:^{5,7}

RD cells were cultured in 48-well plates overnight. The cells were treated with serial diluted testing compounds for 2 hours, then infected with EV-A71 at MOI of 0.1 and cultured for 24 hours. Four duplicates were set for each concentration of the testing compound. 100 µL of culture media were collected to isolate viral RNA. The viral RNA was isolated with TRIzol reagent (Ambion, Life Technologies), and 2 µL of total RNA was used to synthesize cDNA using ImProm-II™ Reverse Transcriptase (Promega) according to the manufacturer's instructions. To set up the standard curve of infectious viruses, the viral titres were first determined by TCID₅₀ or plaque assay; then the viral RNA was extracted from the tittered infectious EV-A71 viruses and used as the standard.

Quantitative reverse-transcription polymerase chain reaction (RT-qPCR):^{5,7}

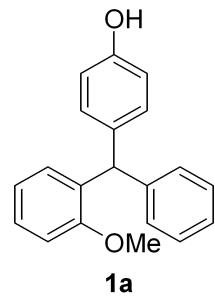
Quantitative Real-time PCR (RT-qPCR) was performed using SYBR Green Mix (Life Technologies) on Applied Biosystems QuantStudio™ 3 Real-Time PCR Systems. The RT-qPCR was performed by using the following primer pairs:

EV-A71, 5'-GCAGCCCAAAAGAACTTCAC-3' (forward) and 5'-ATTCAGCAGCTTGGAGTGC-3'(reverse). The target fragment amplification was carried out as follows: reverse transcription at 50 °C for 30 min; initial activation of HotStar Taq DNA Polymerase at 95 °C for 15 min; 45 cycles in four steps: 94 °C for 10 s, 56 °C for 30 s, and 72 °C for 30 s. At the end of the amplification cycles, melting temperature analysis was carried out by a slow increase in temperature (0.1 °C/s) up to 95 °C. The copy number of virions was used to calculate the testing compound's IC₅₀ (half maximal inhibitory concentration).

V. Determination of the Product Structure

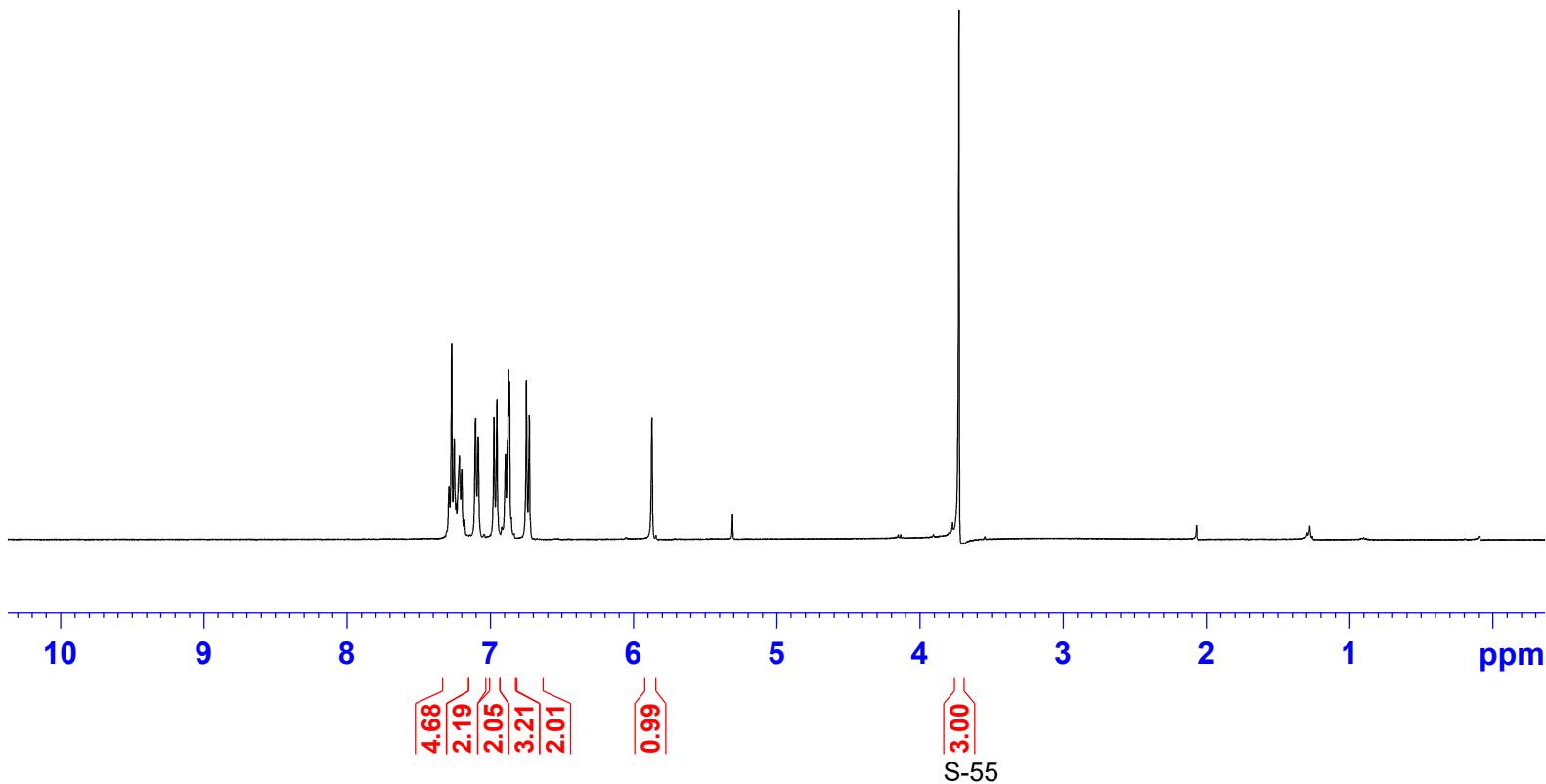
The structure and absolute stereochemistry of the products were determined by comparison with the literature report.⁸

(8) X. Li, M. Duan, Z. Deng, Q. Shao, M. Chen, G. Zhu, K. N. Houk and J. Sun, *Nat. Catal.*, 2020, **3**, 1010–1019.



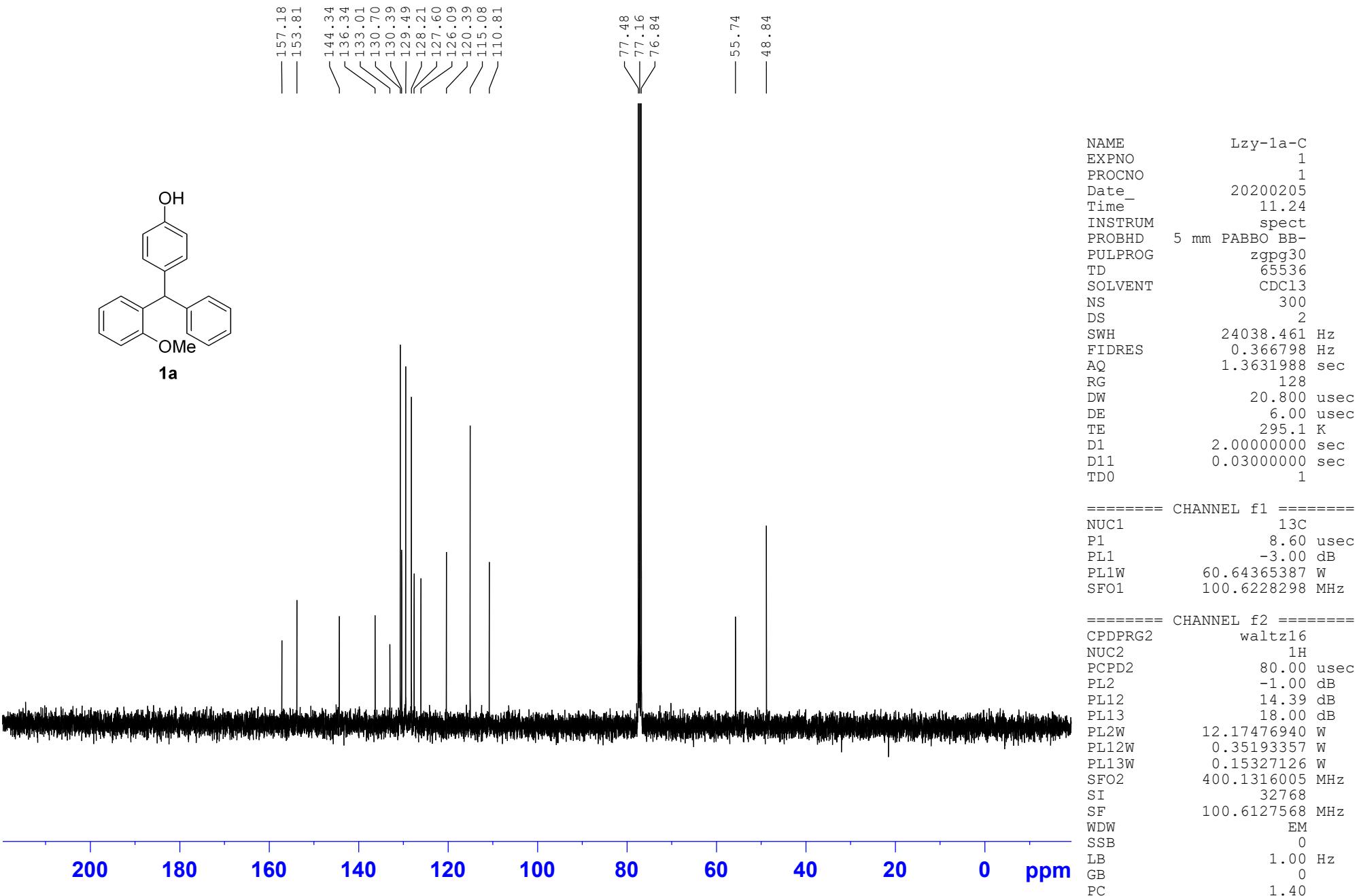
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6.86
6.85
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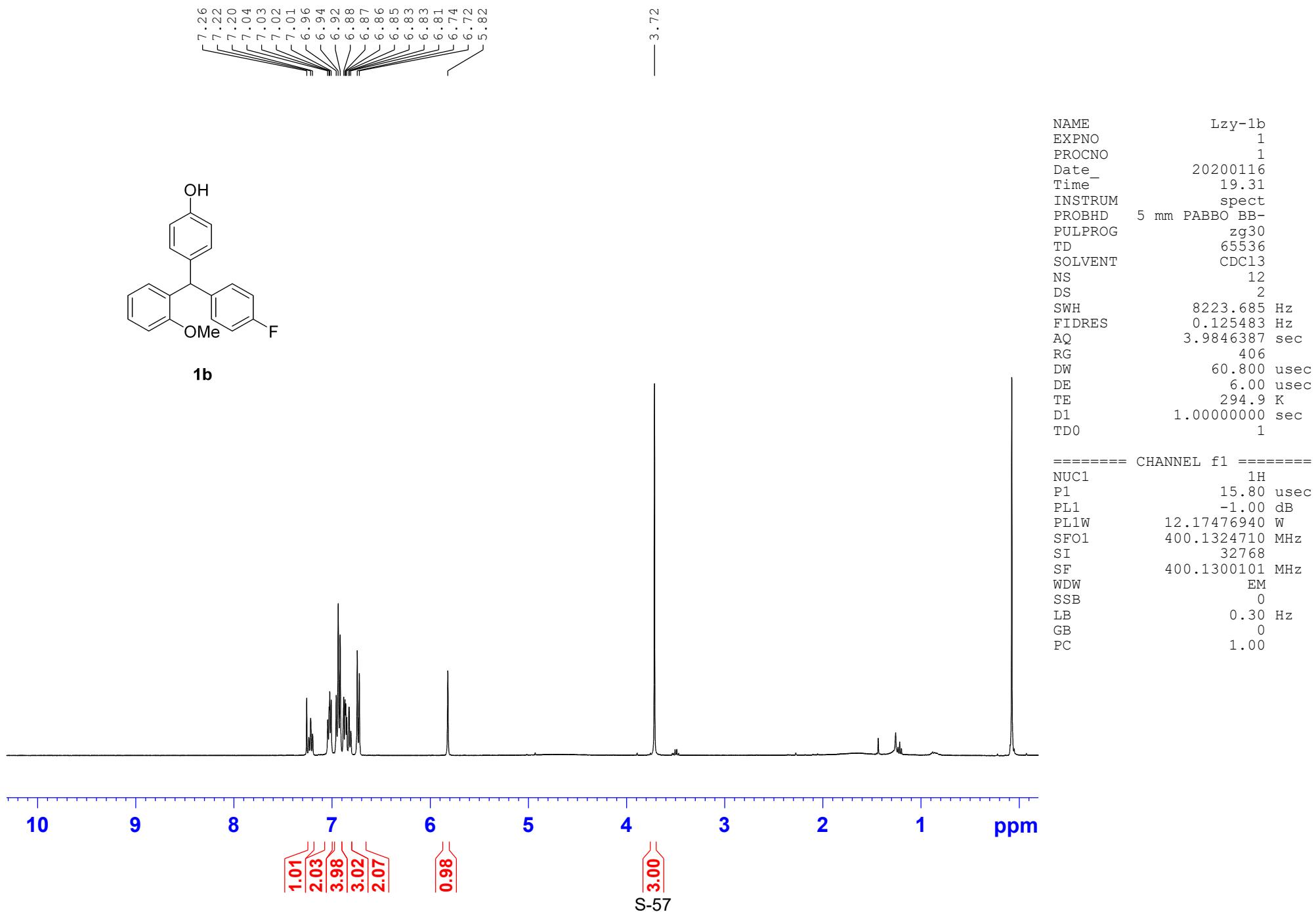
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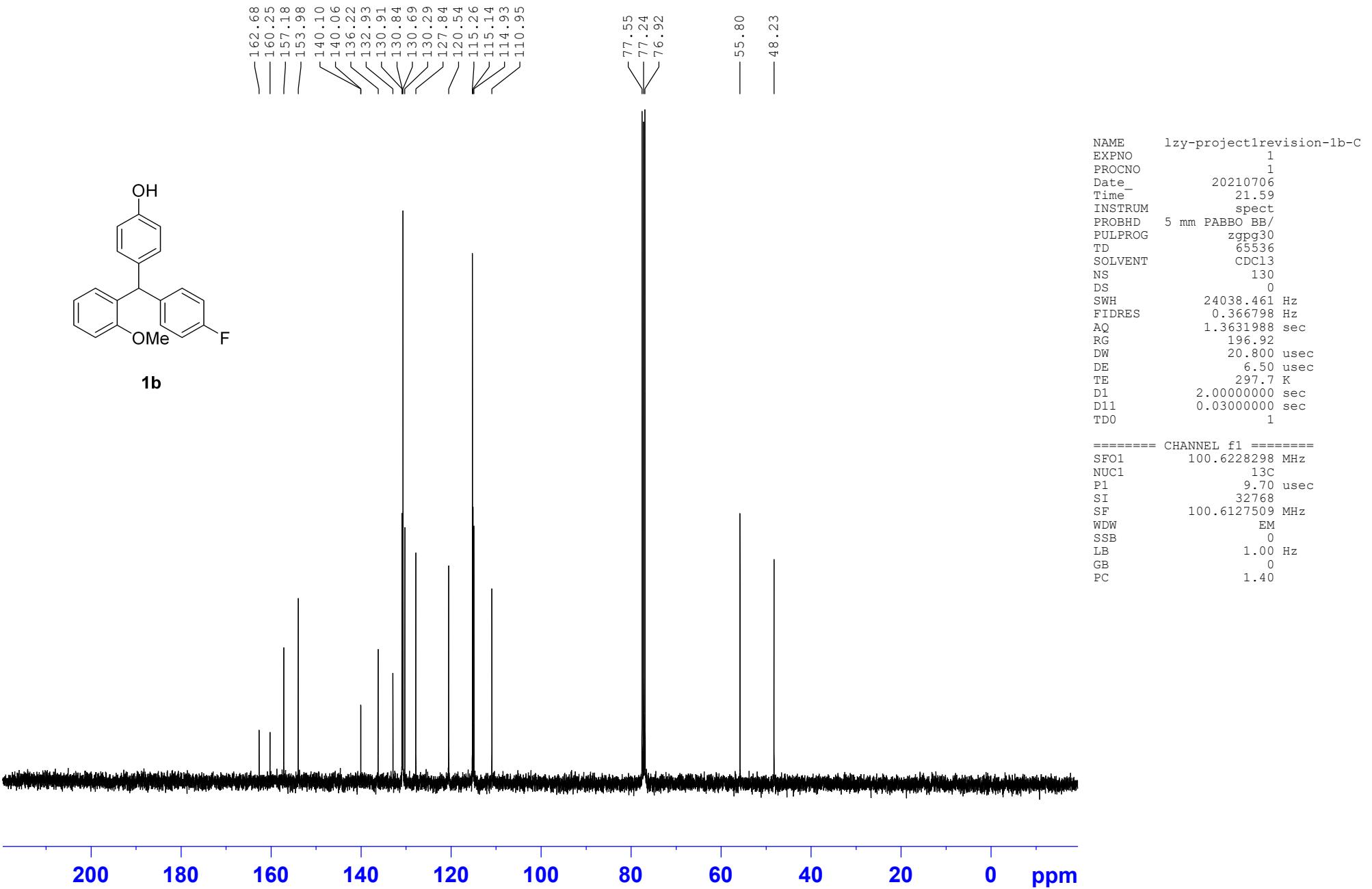


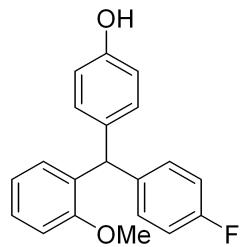
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DW 60.800 usec
DE 6.00 usec
TE 294.9 K
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TD0 1

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1b

-117.45

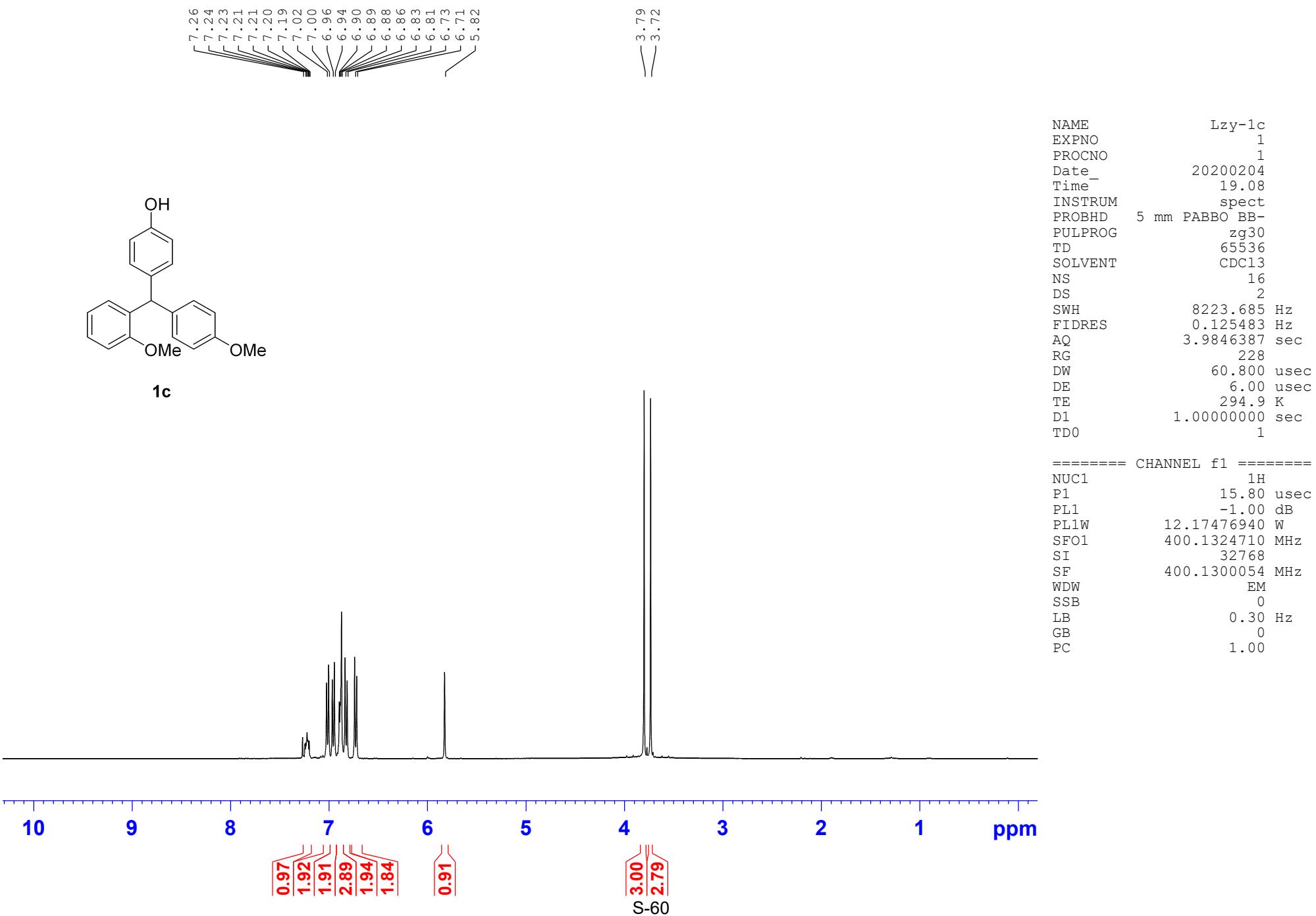
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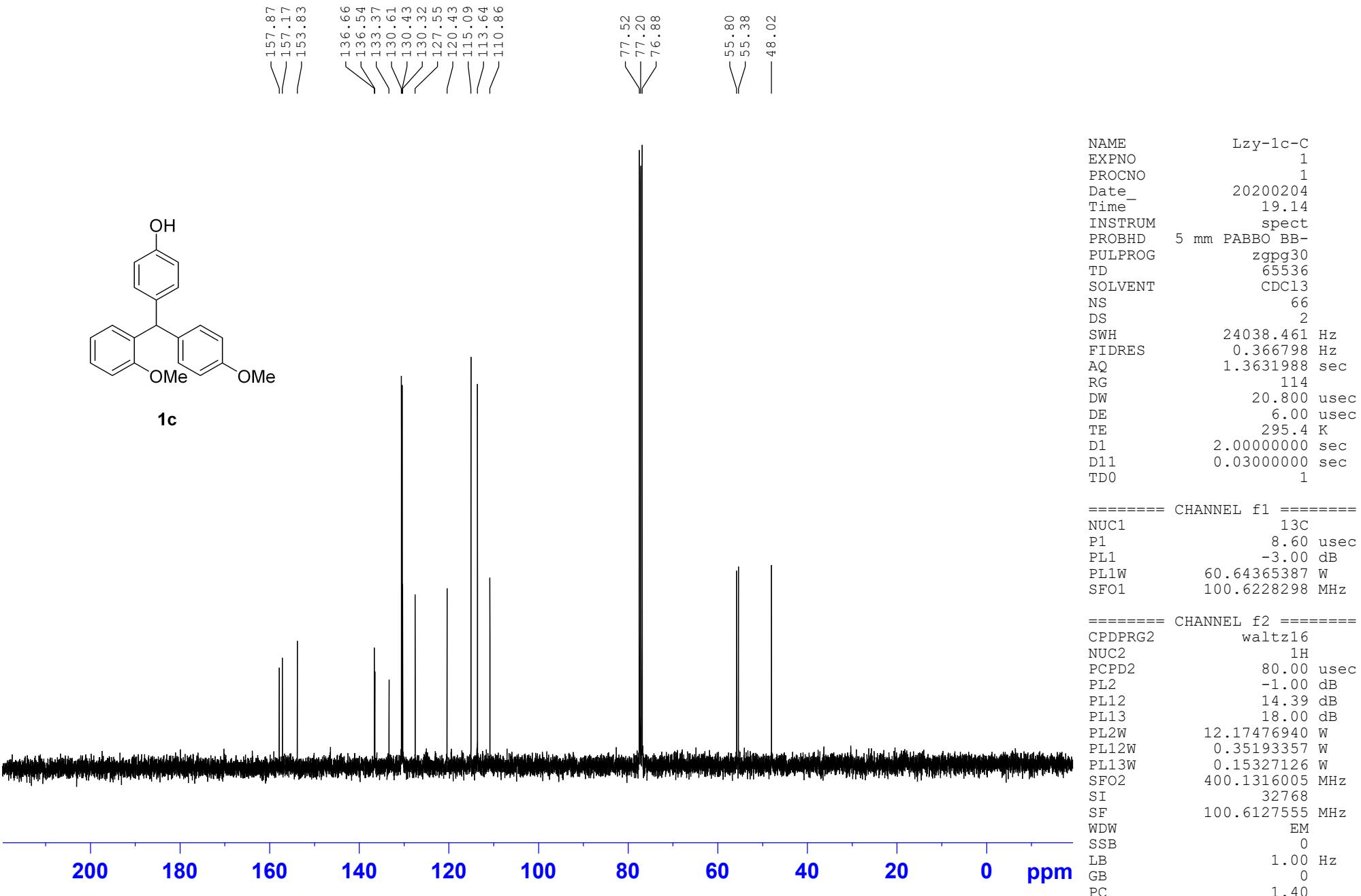
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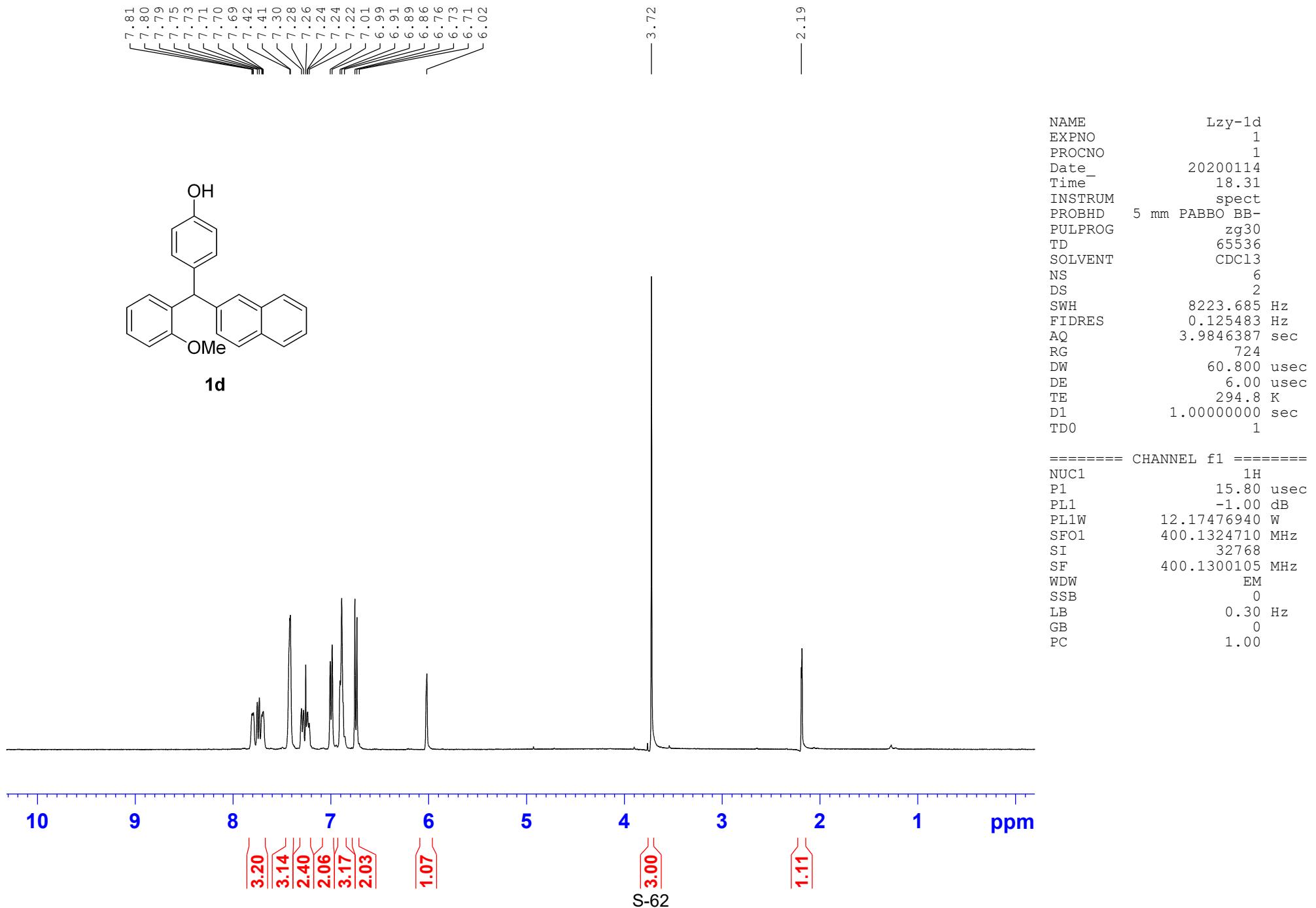
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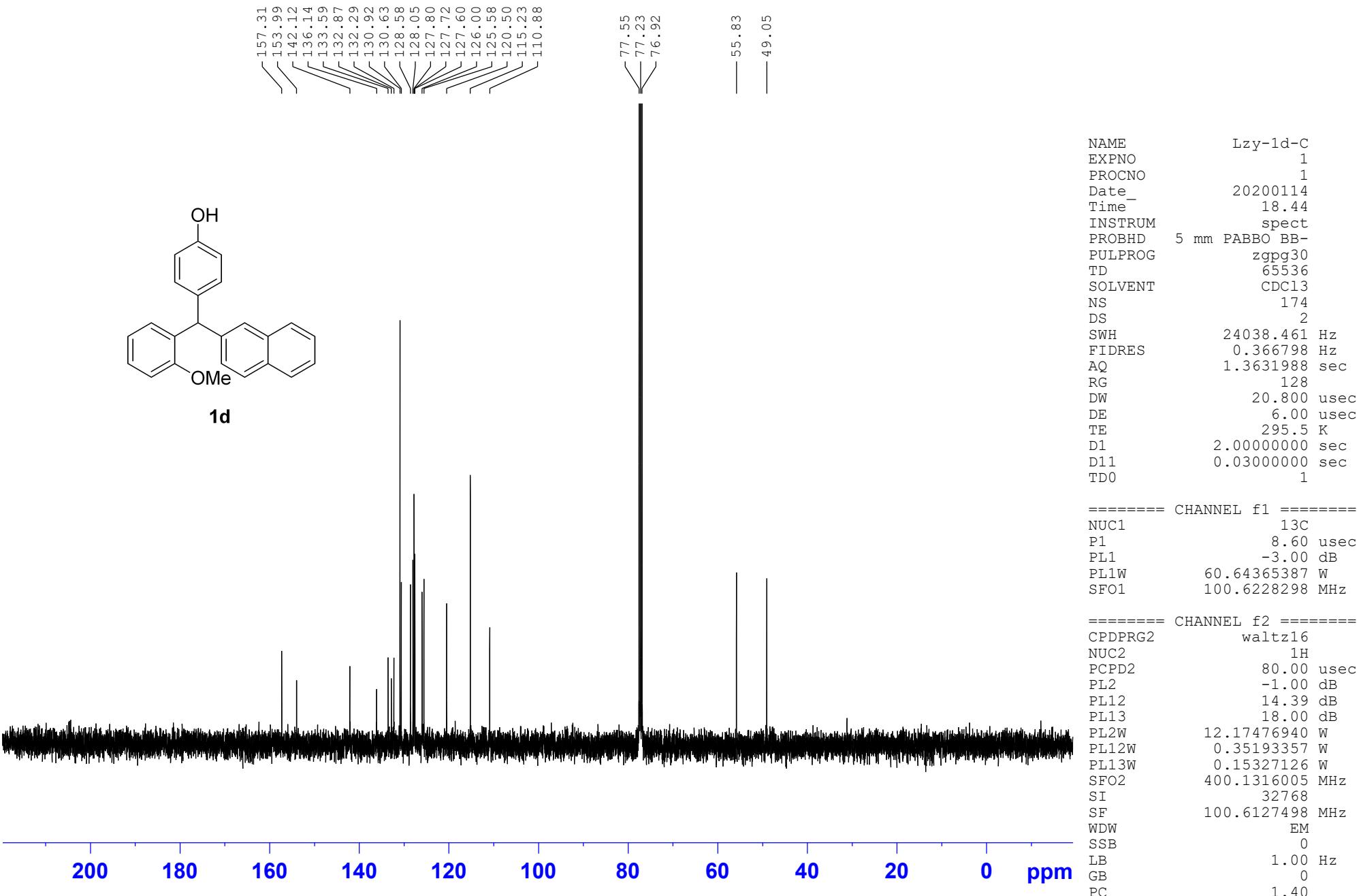
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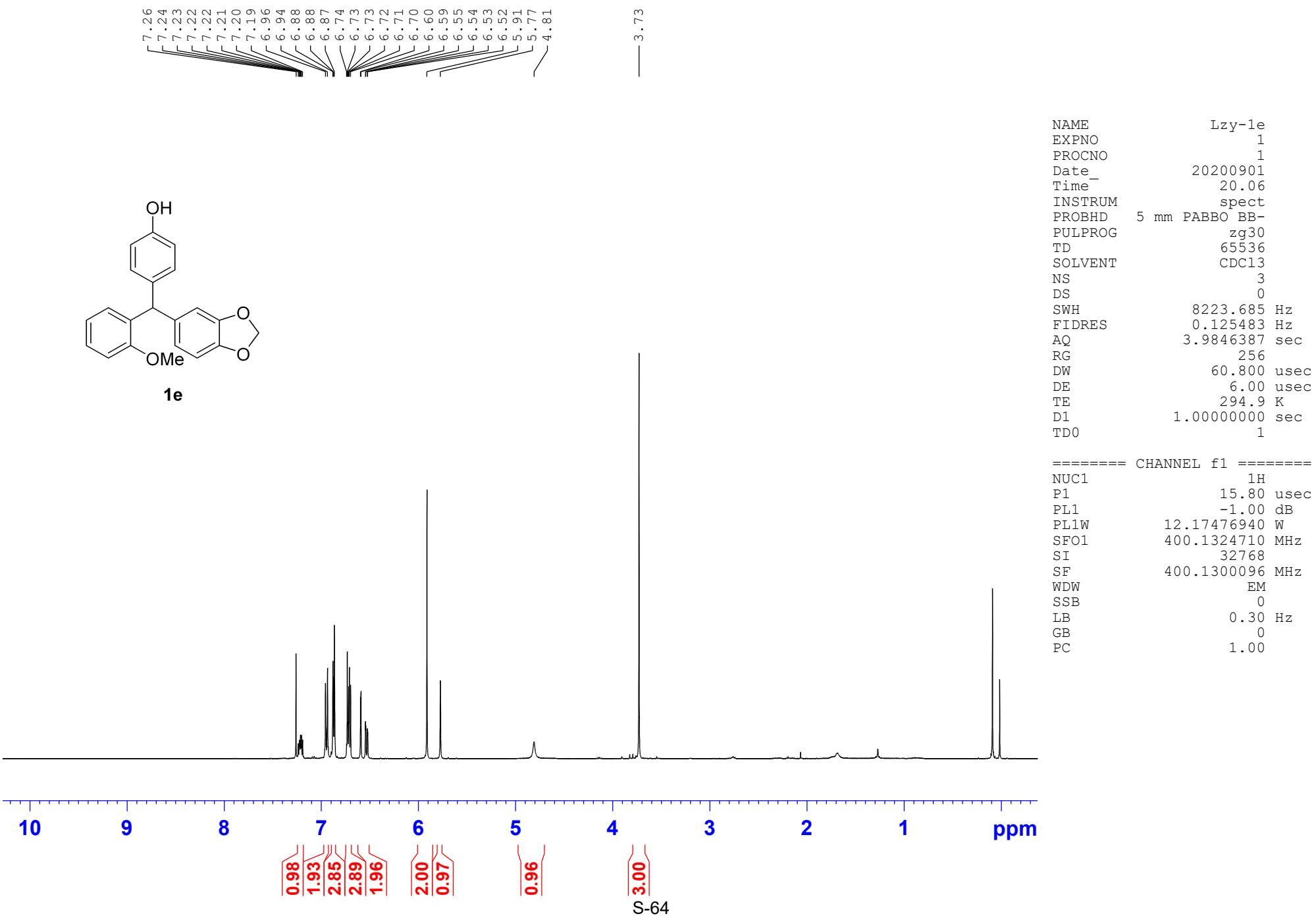
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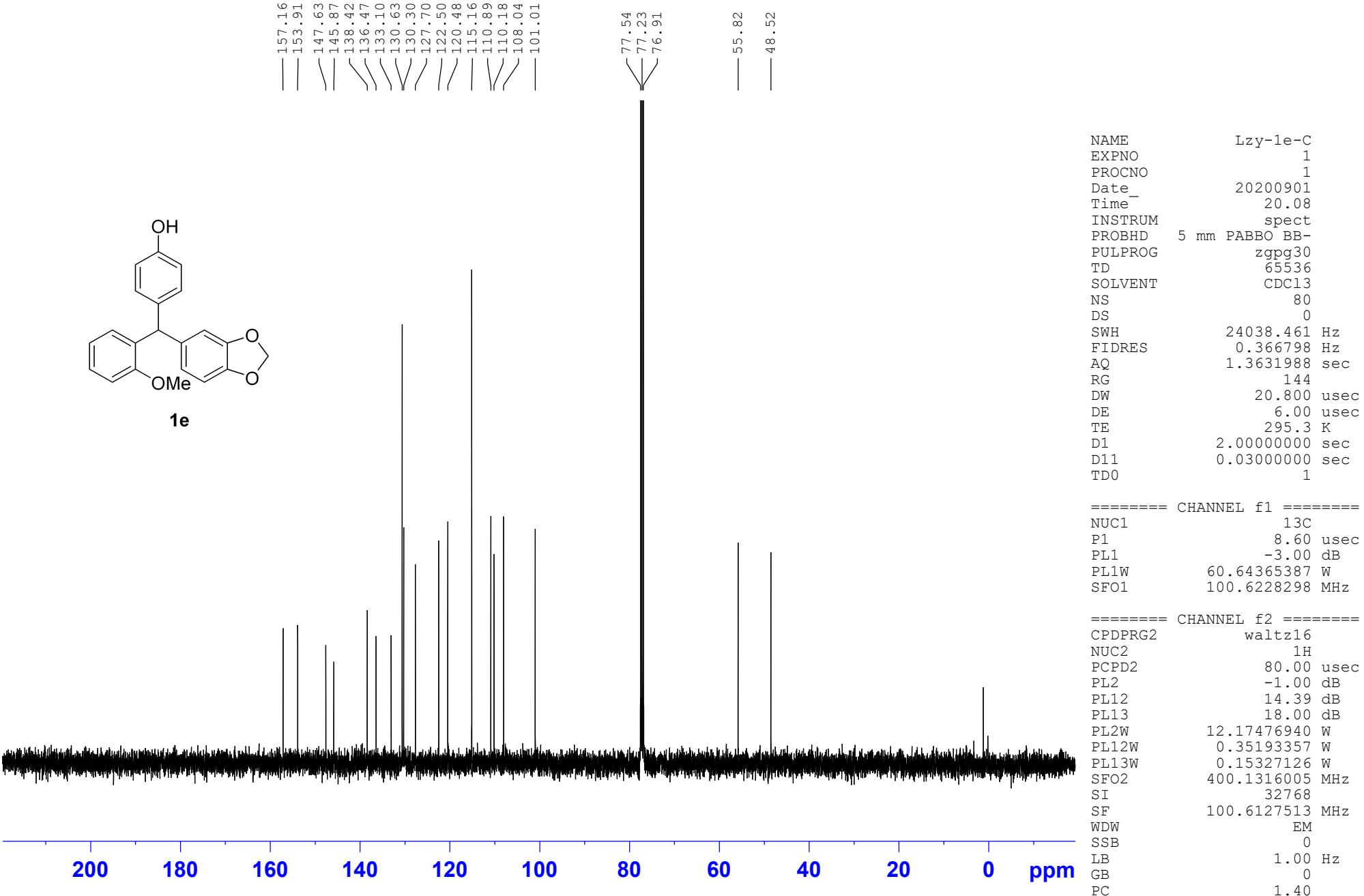


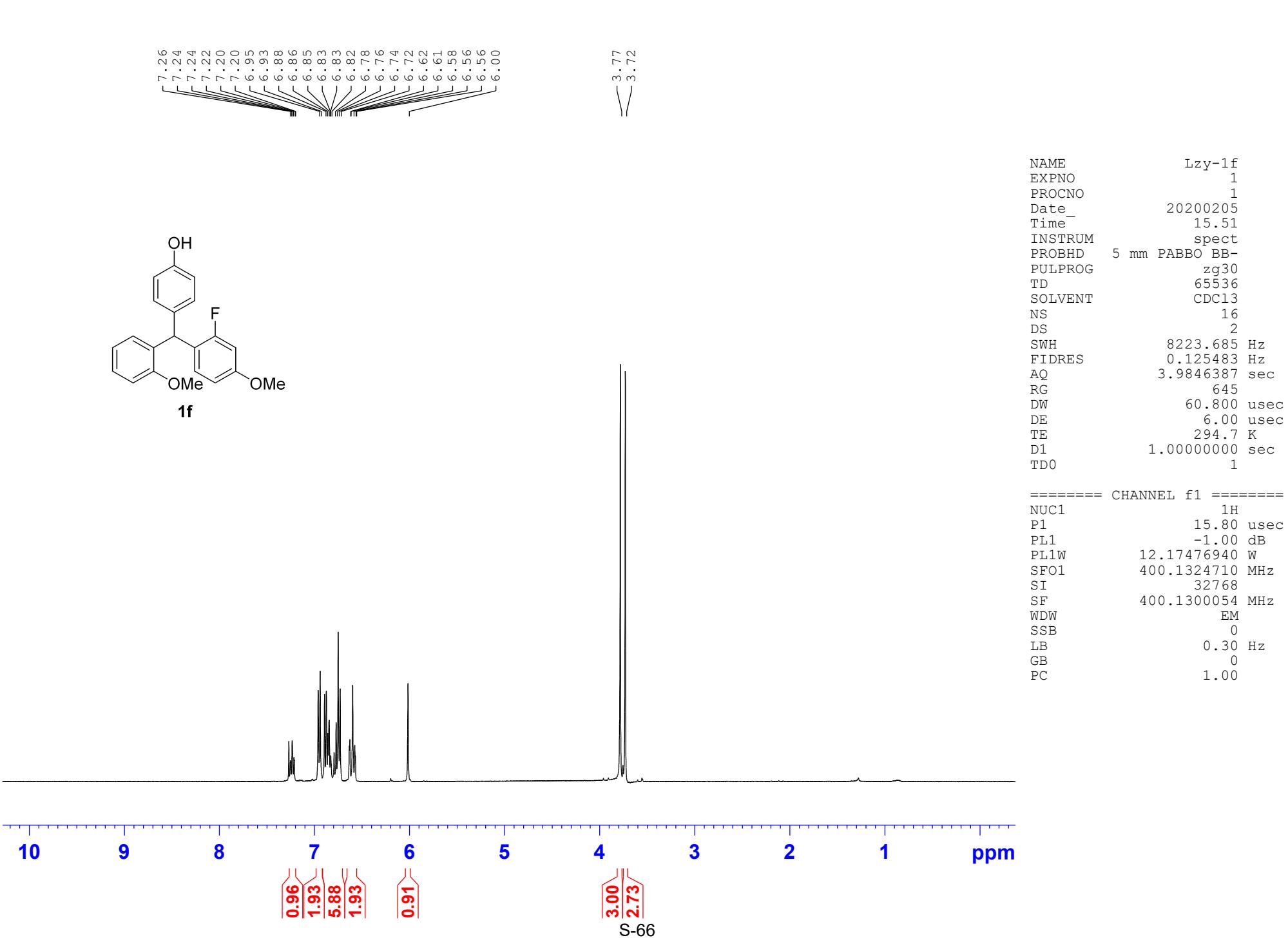
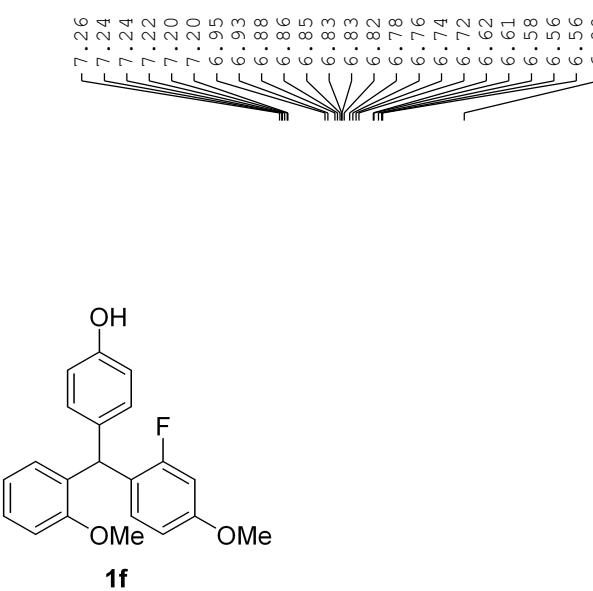


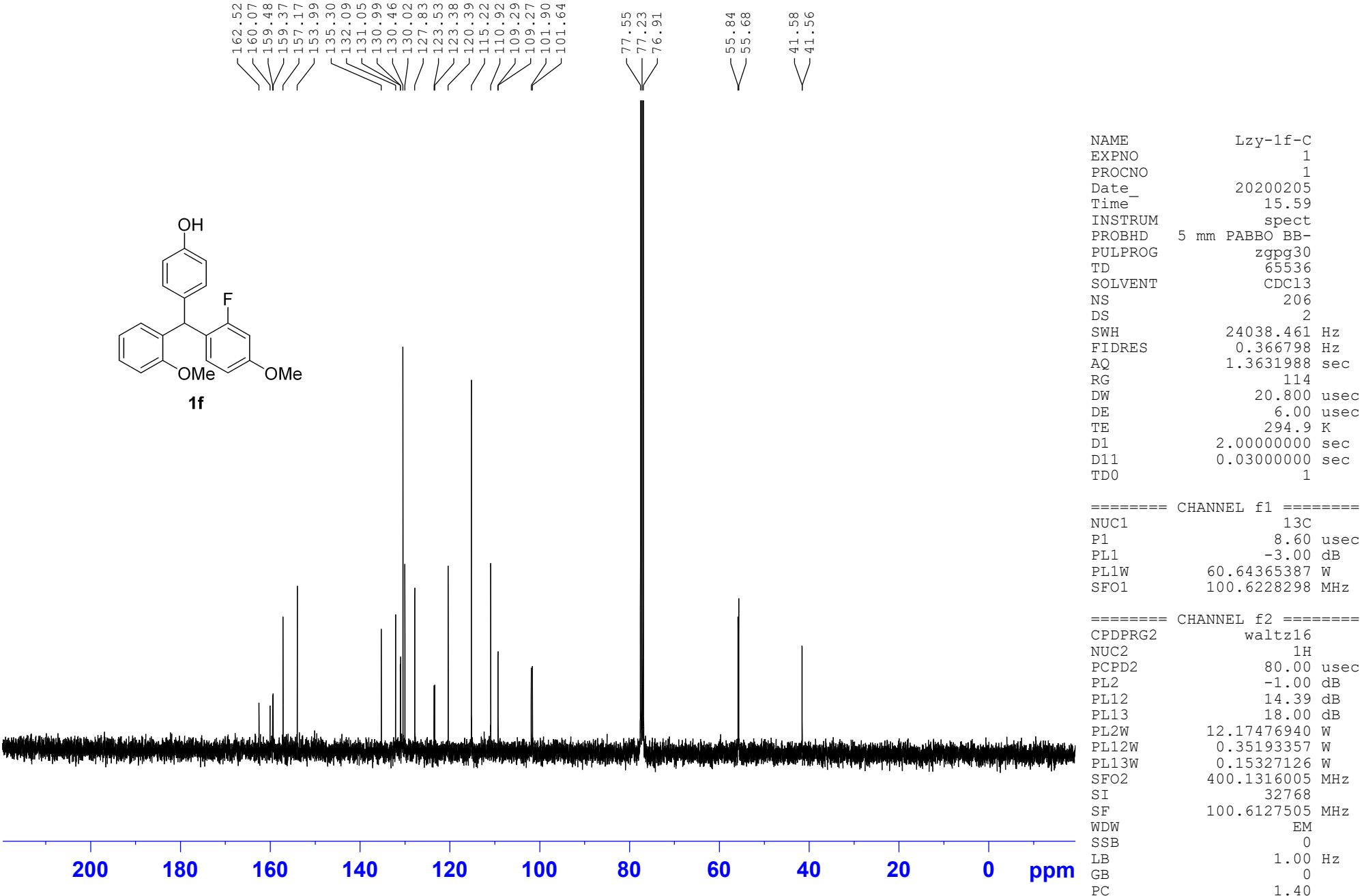


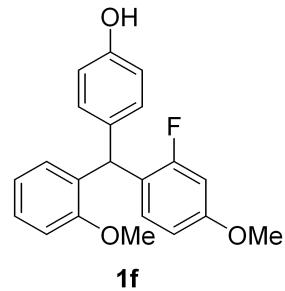










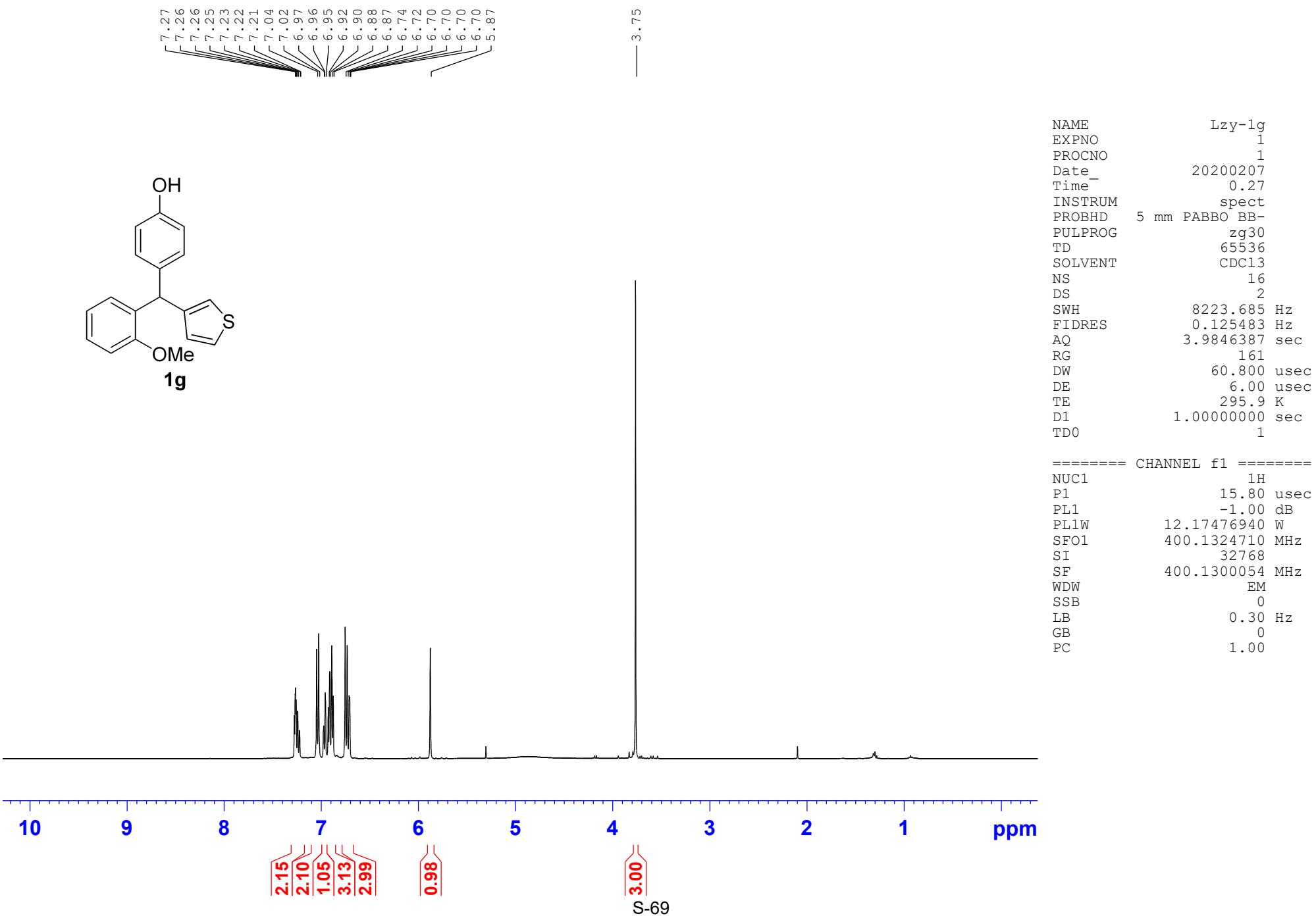


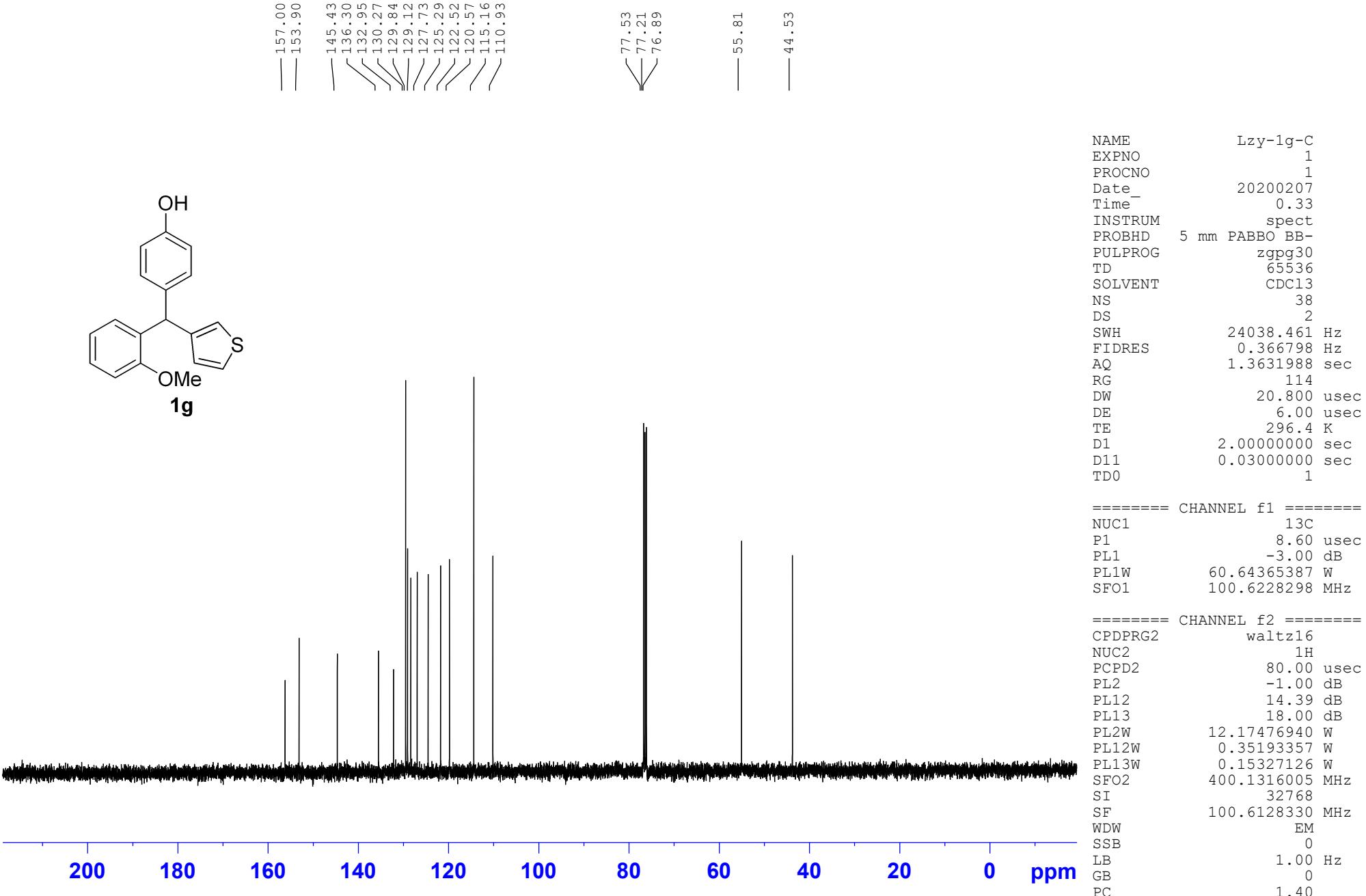
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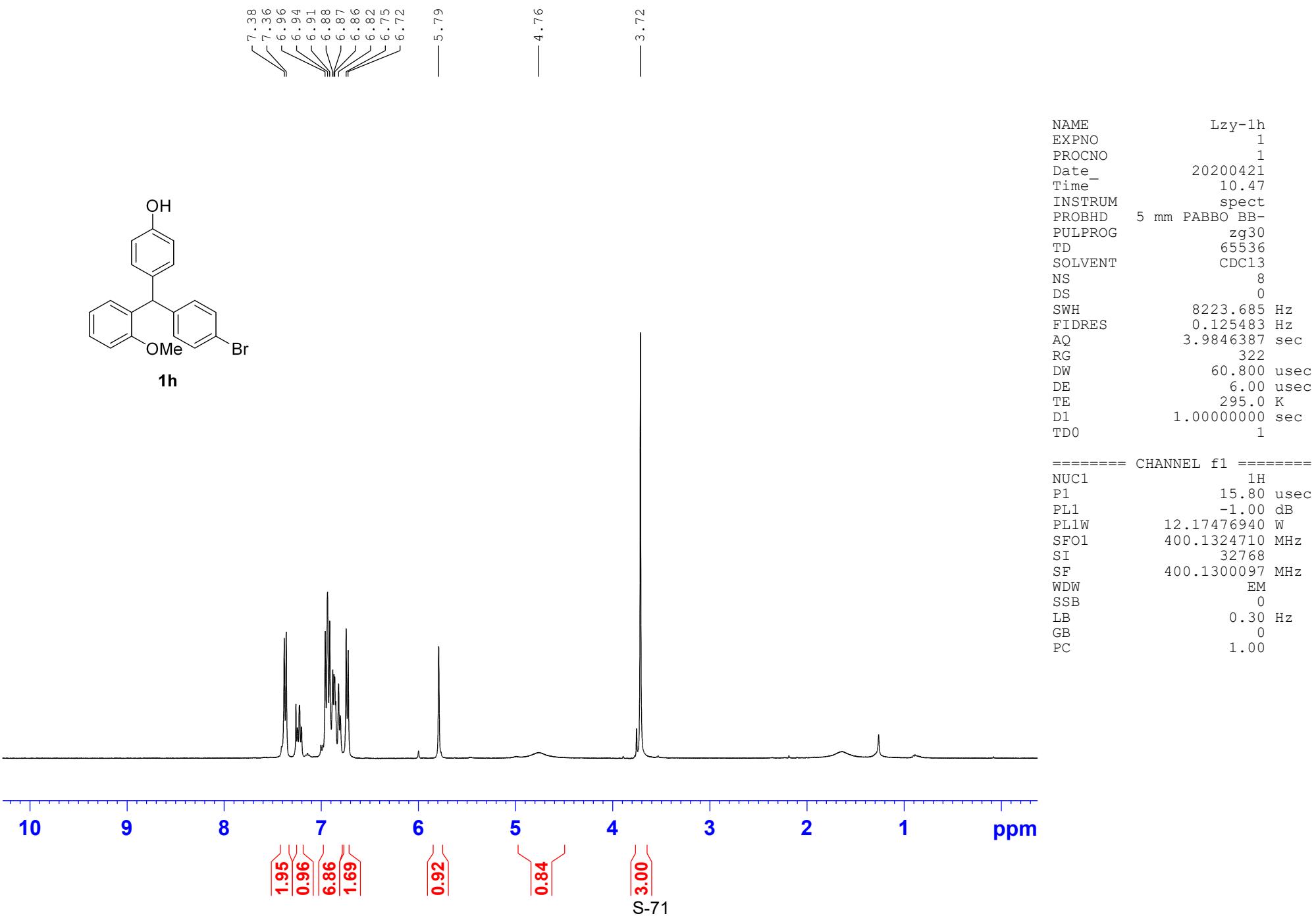
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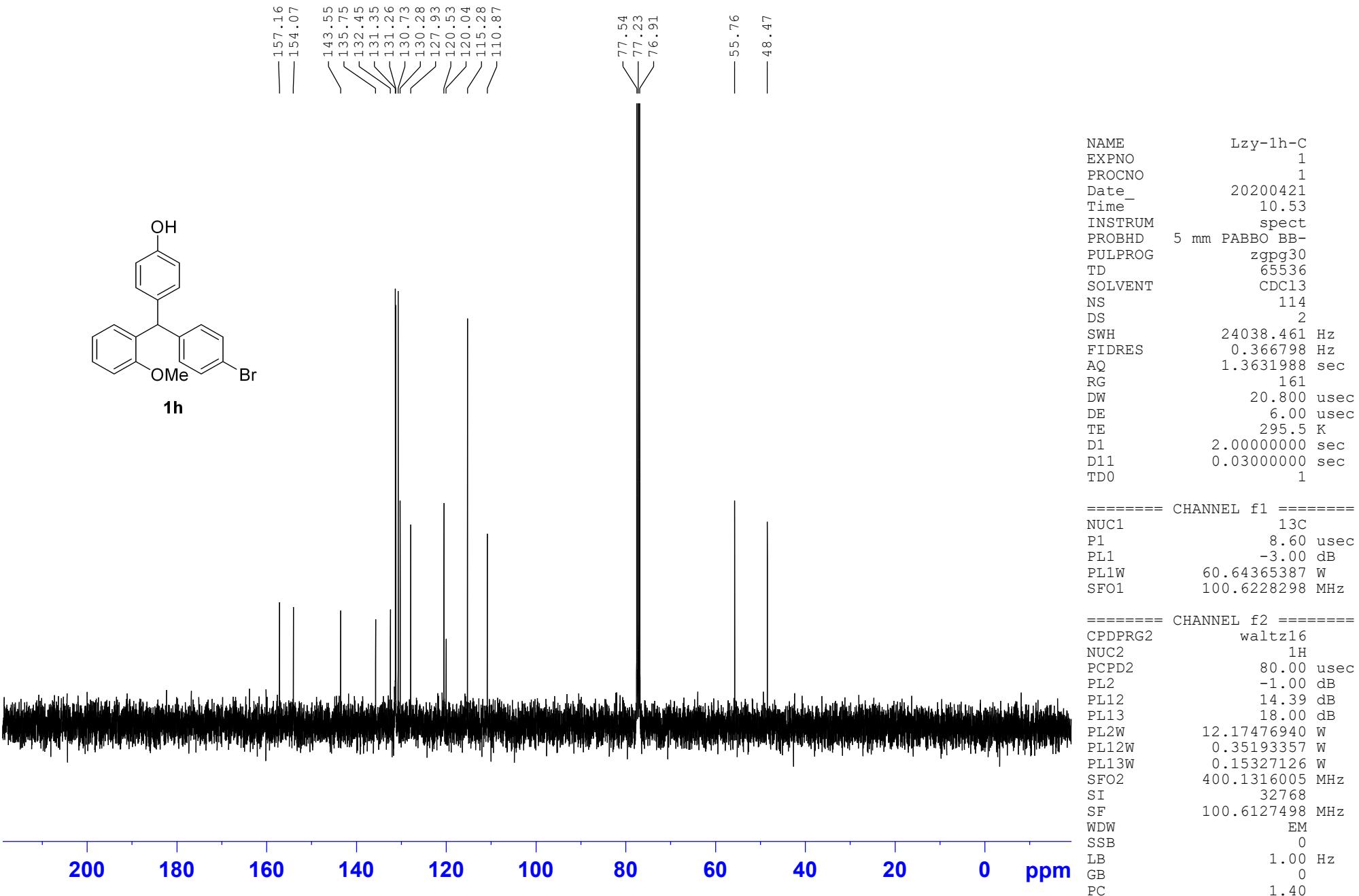
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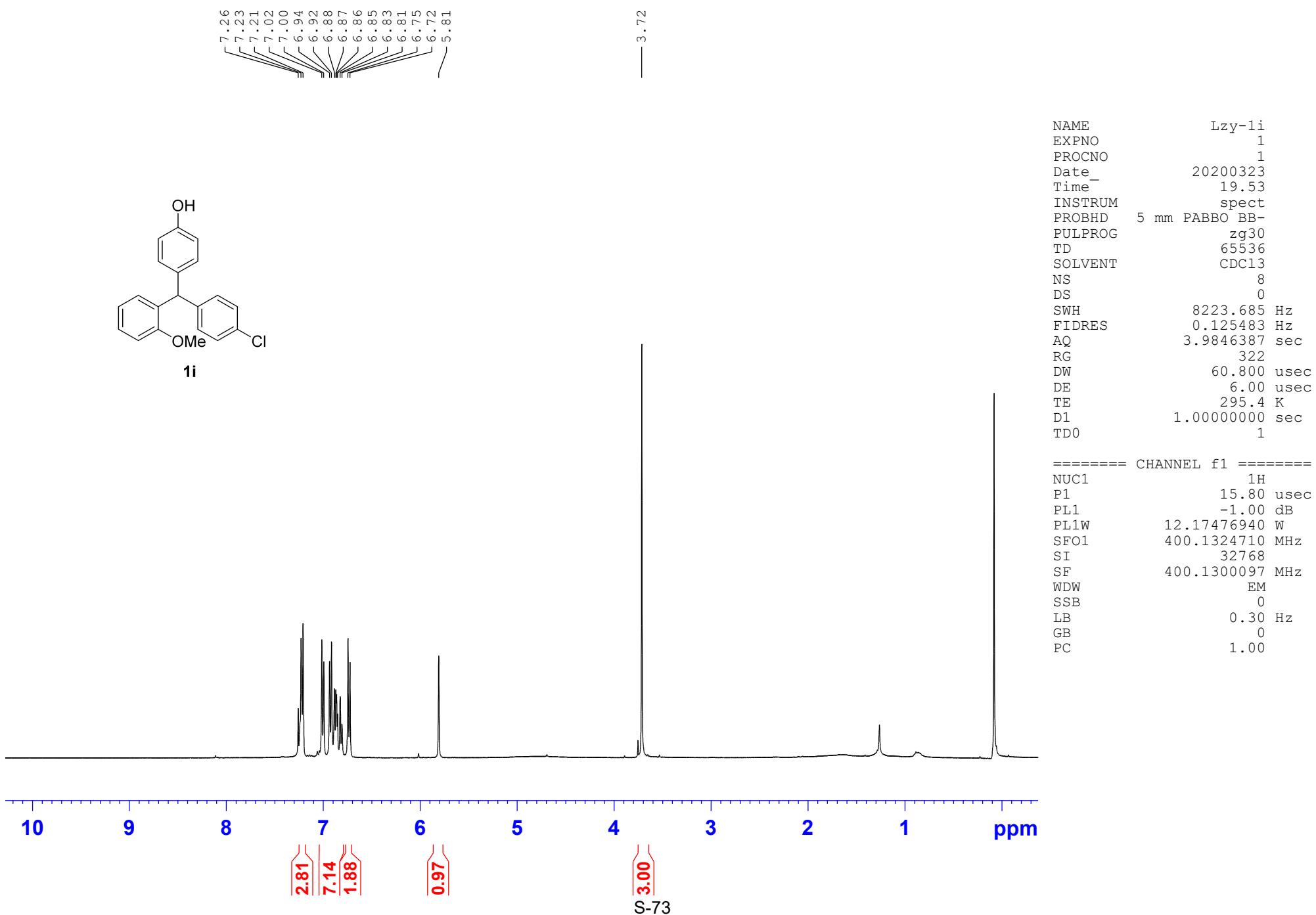
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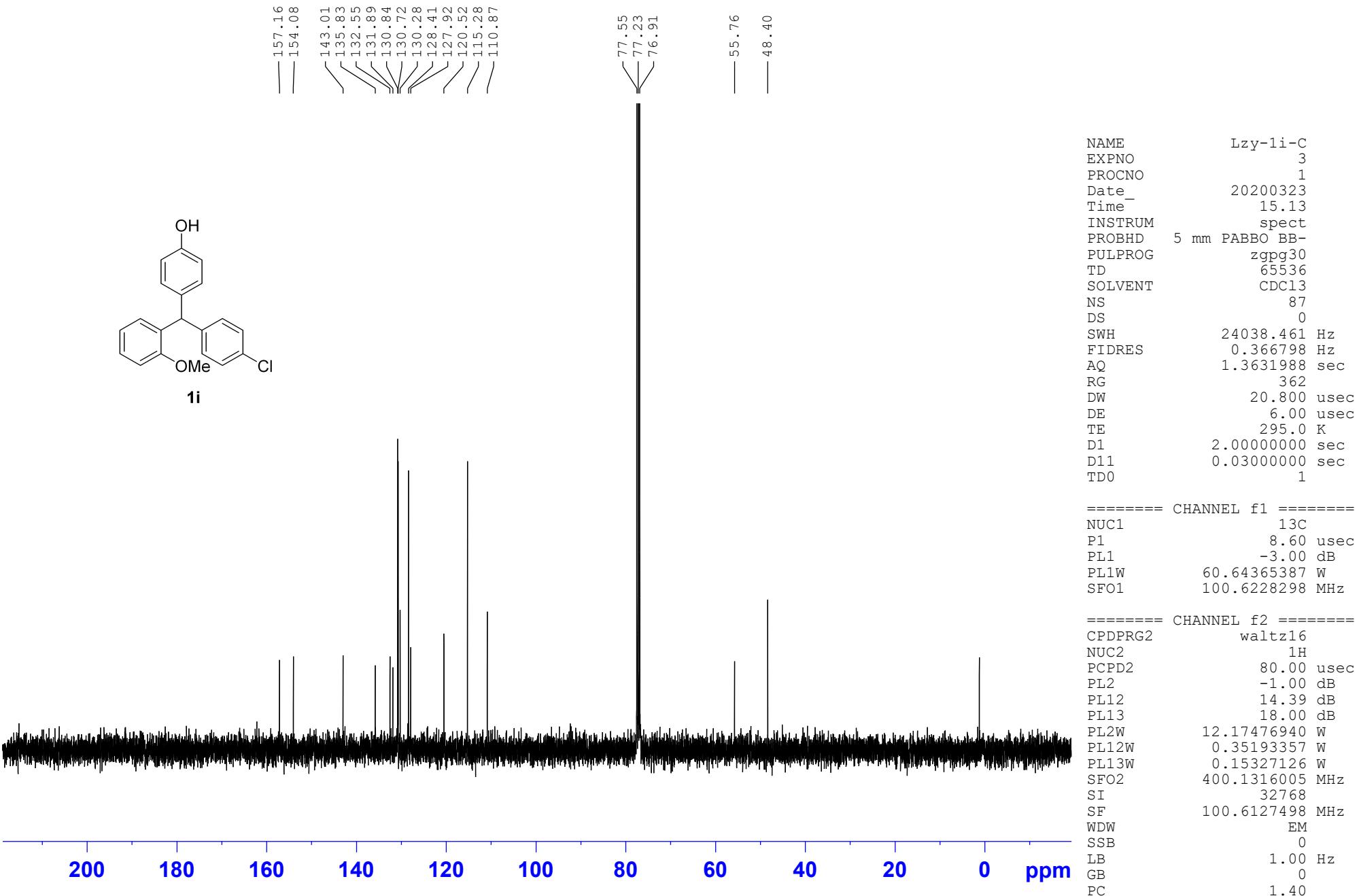


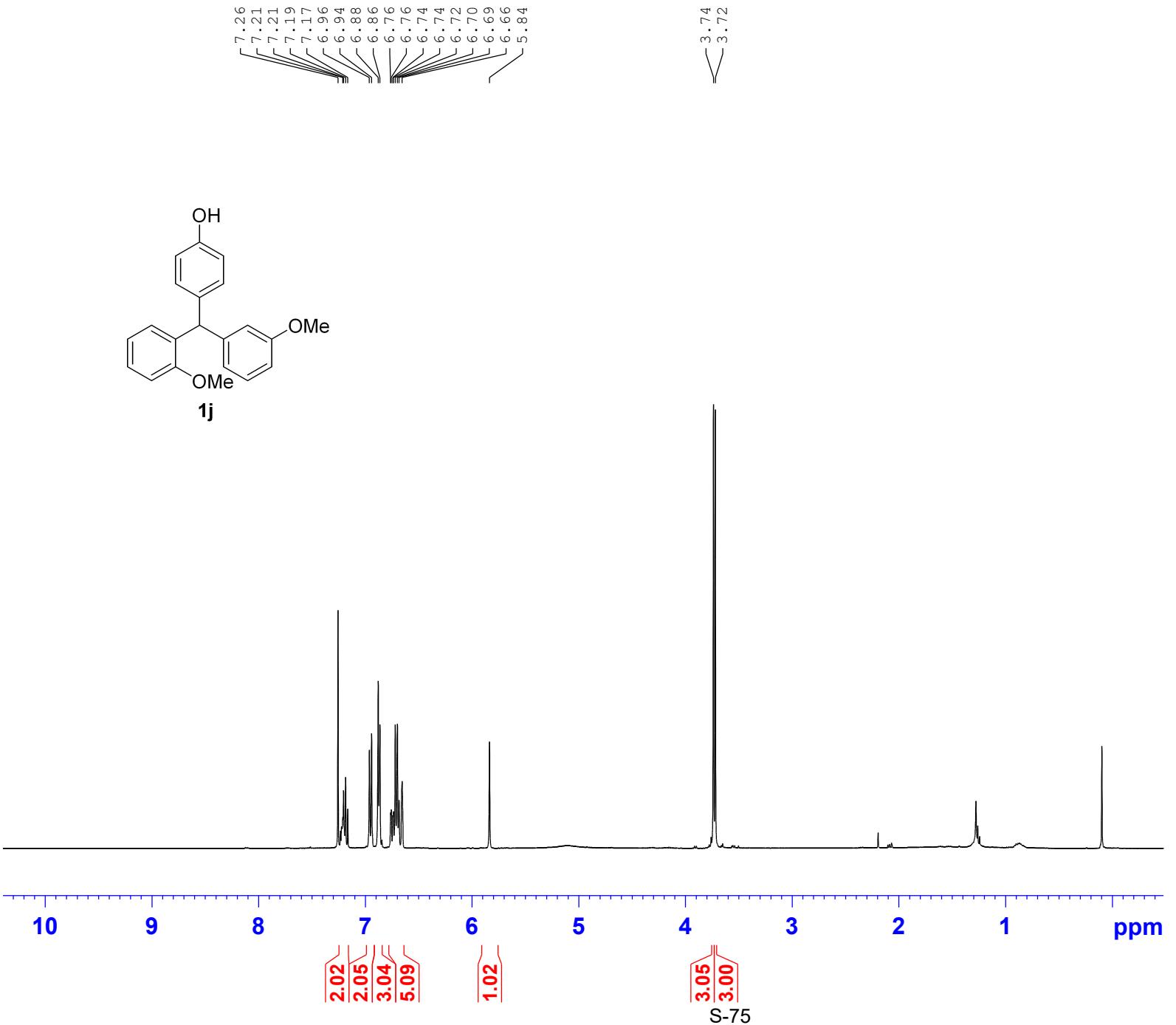
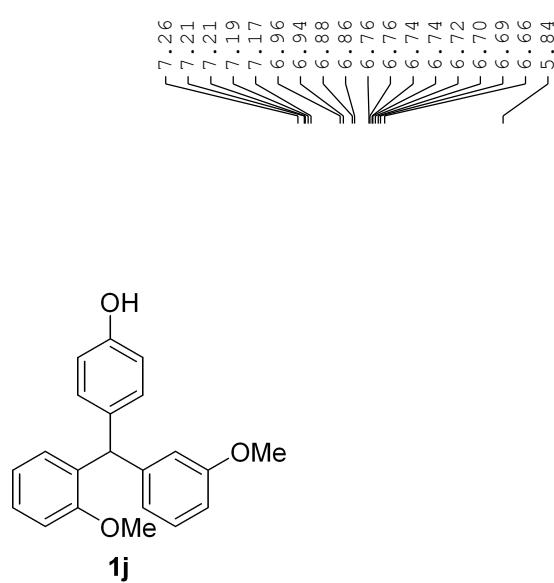








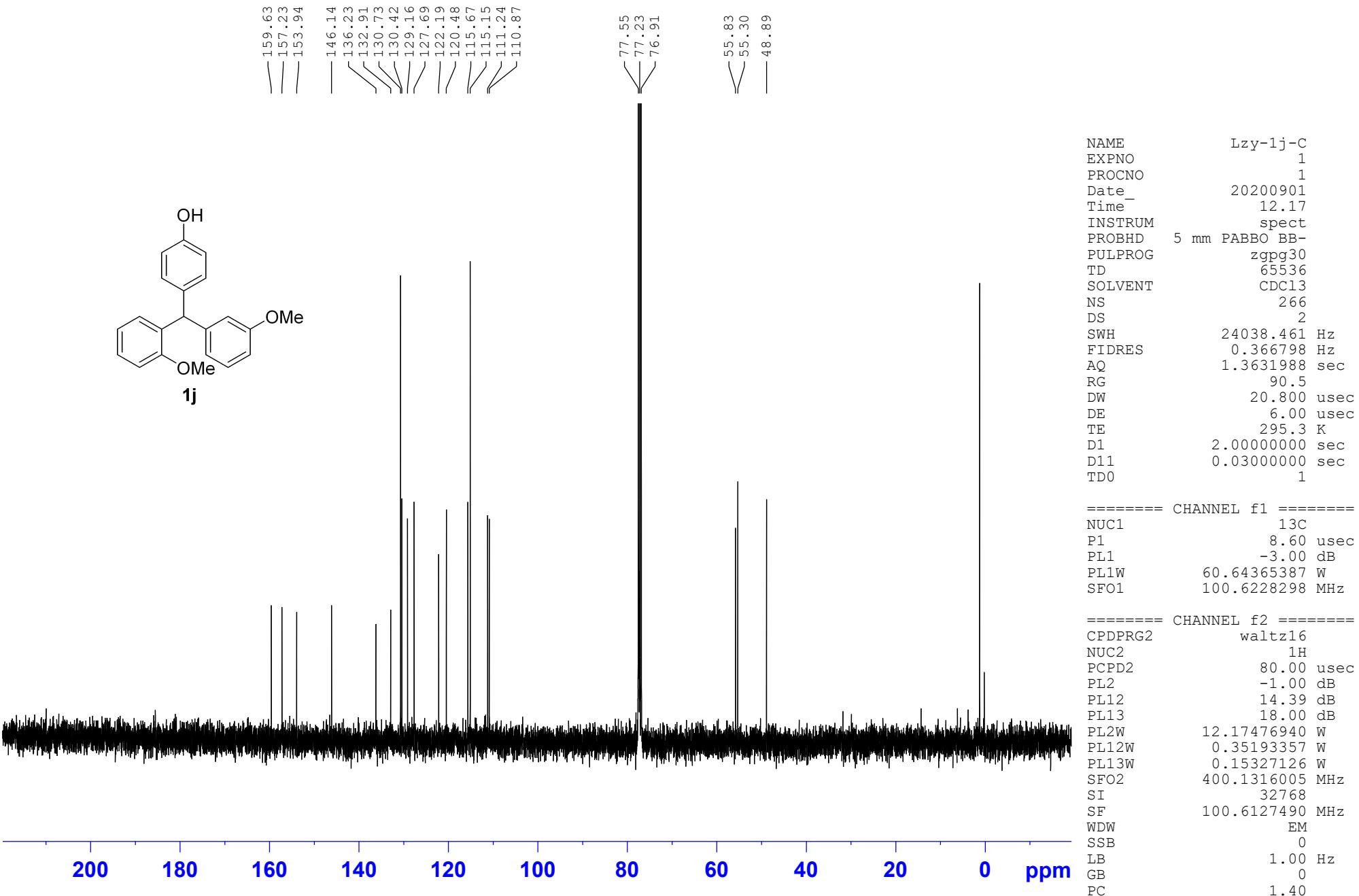


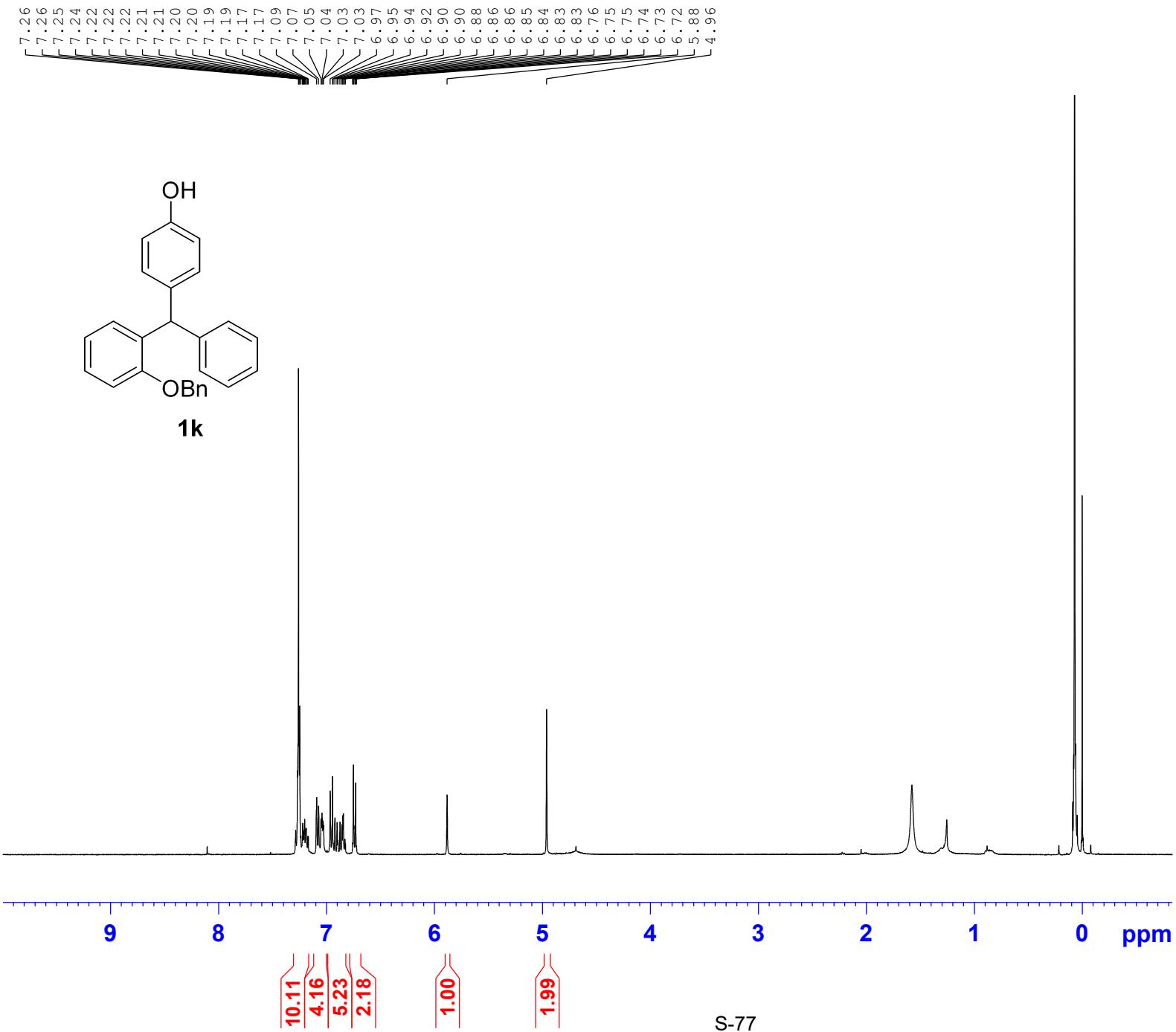


3.74
3.72

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TD0 1

===== CHANNEL f1 =====
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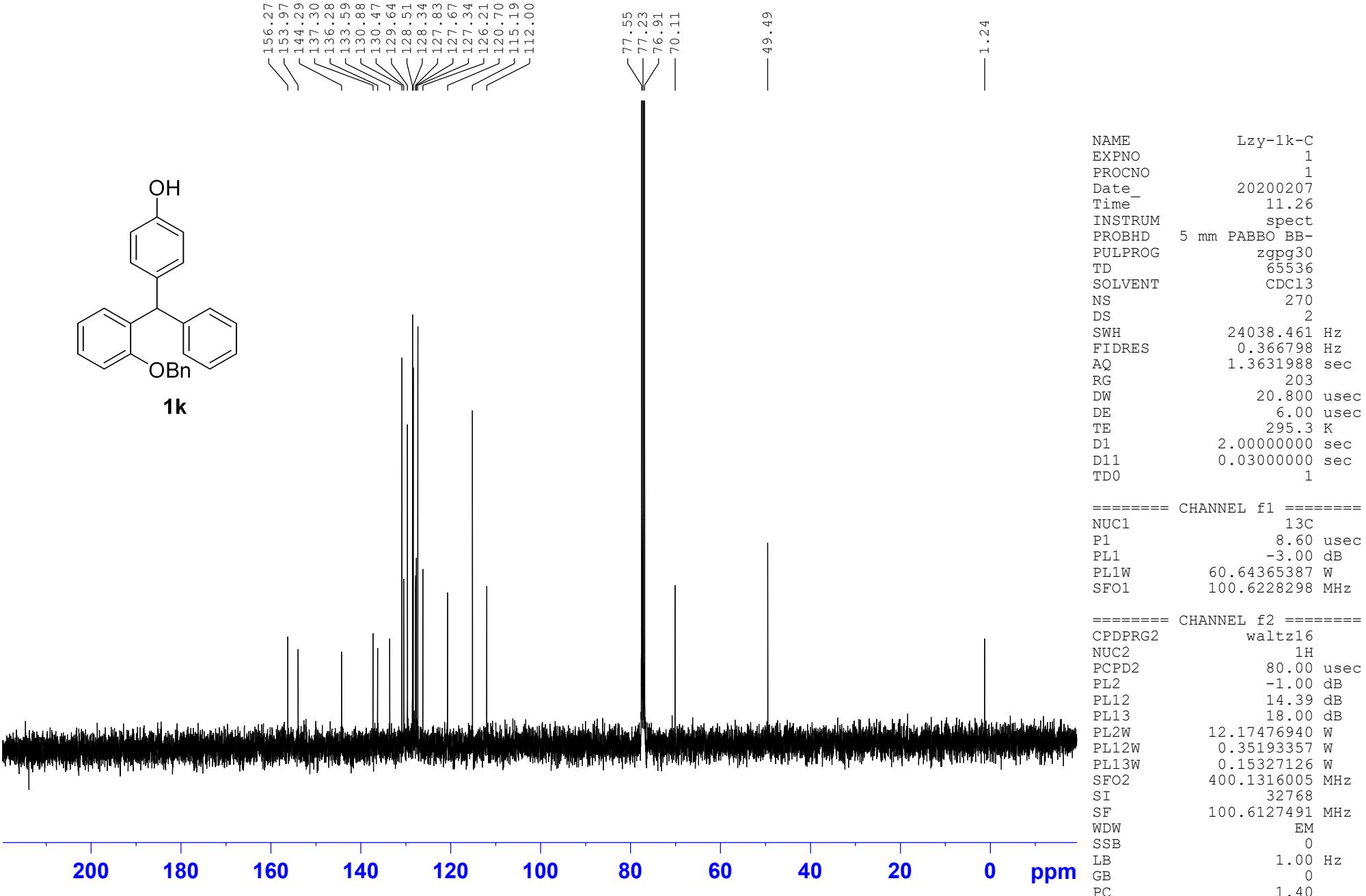


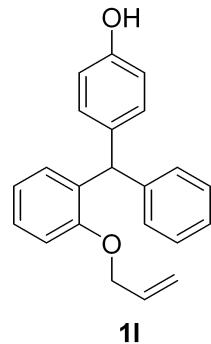
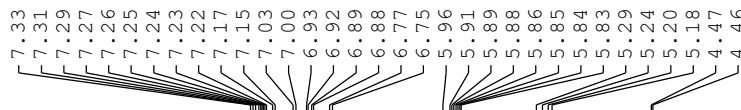


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 NS 16
 DS 0
 SWH 8223.685 Hz
 FIDRES 0.125483 Hz
 AQ 3.9846387 sec
 RG 645
 DW 60.800 usec
 DE 6.00 usec
 TE 294.2 K
 D1 1.0000000 sec
 TDO 1

===== CHANNEL f1 =====

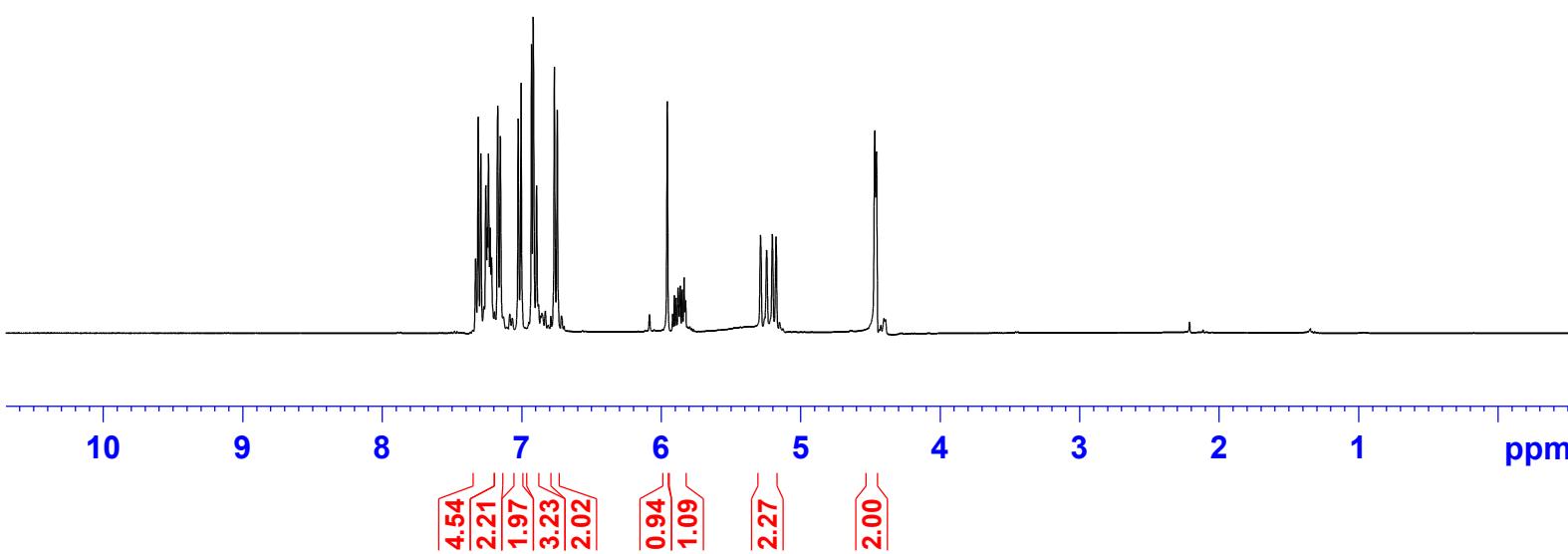
NUC1 ¹H
 P1 15.80 usec
 PL1 -1.00 dB
 PL1W 12.17476940 W
 SFO1 400.1324710 MHz
 SI 32768
 SF 400.1300099 MHz
 WDW EM
 SSB 0
 LB 0.30 Hz
 GB 0
 PC 1.00

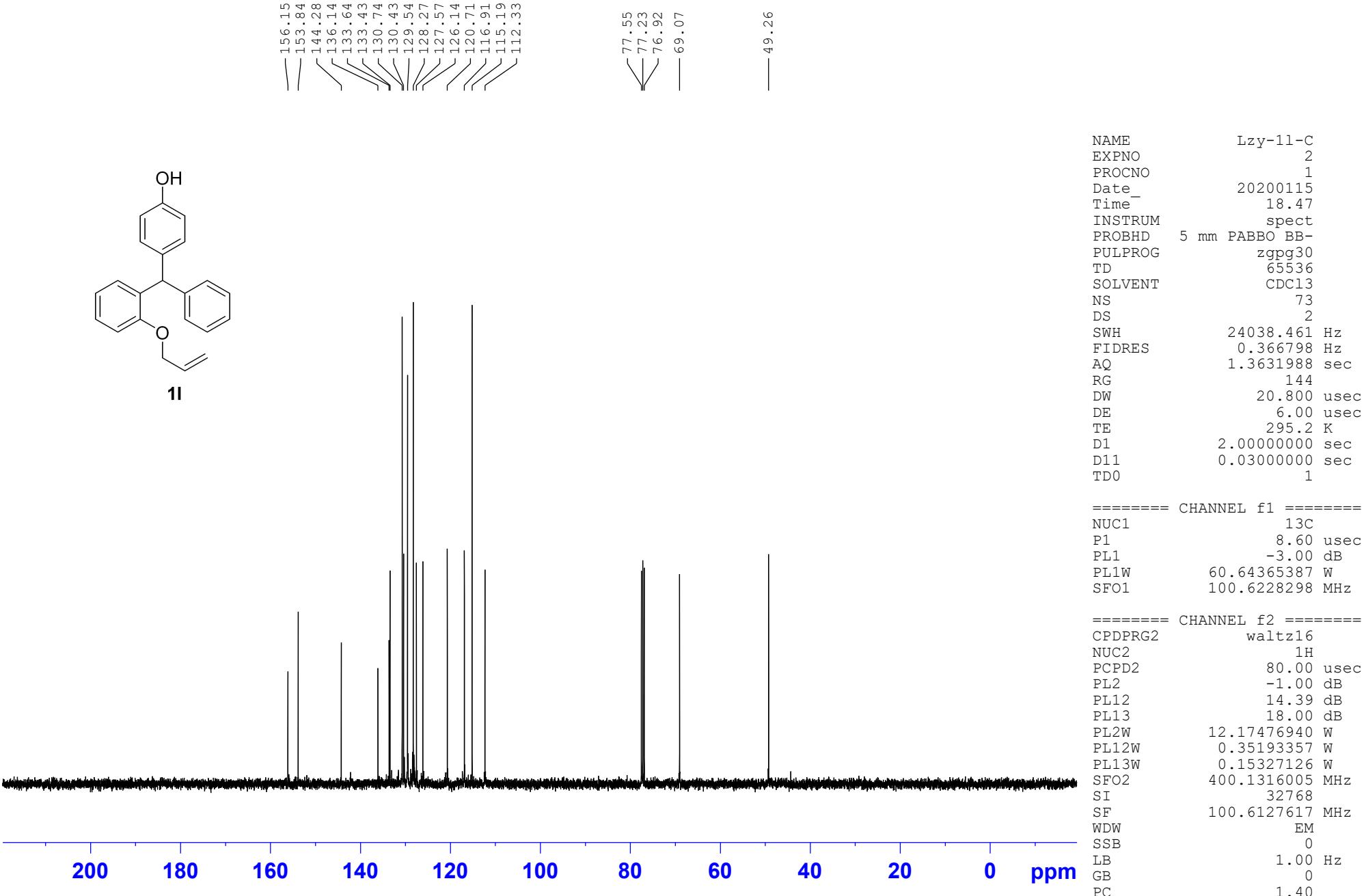


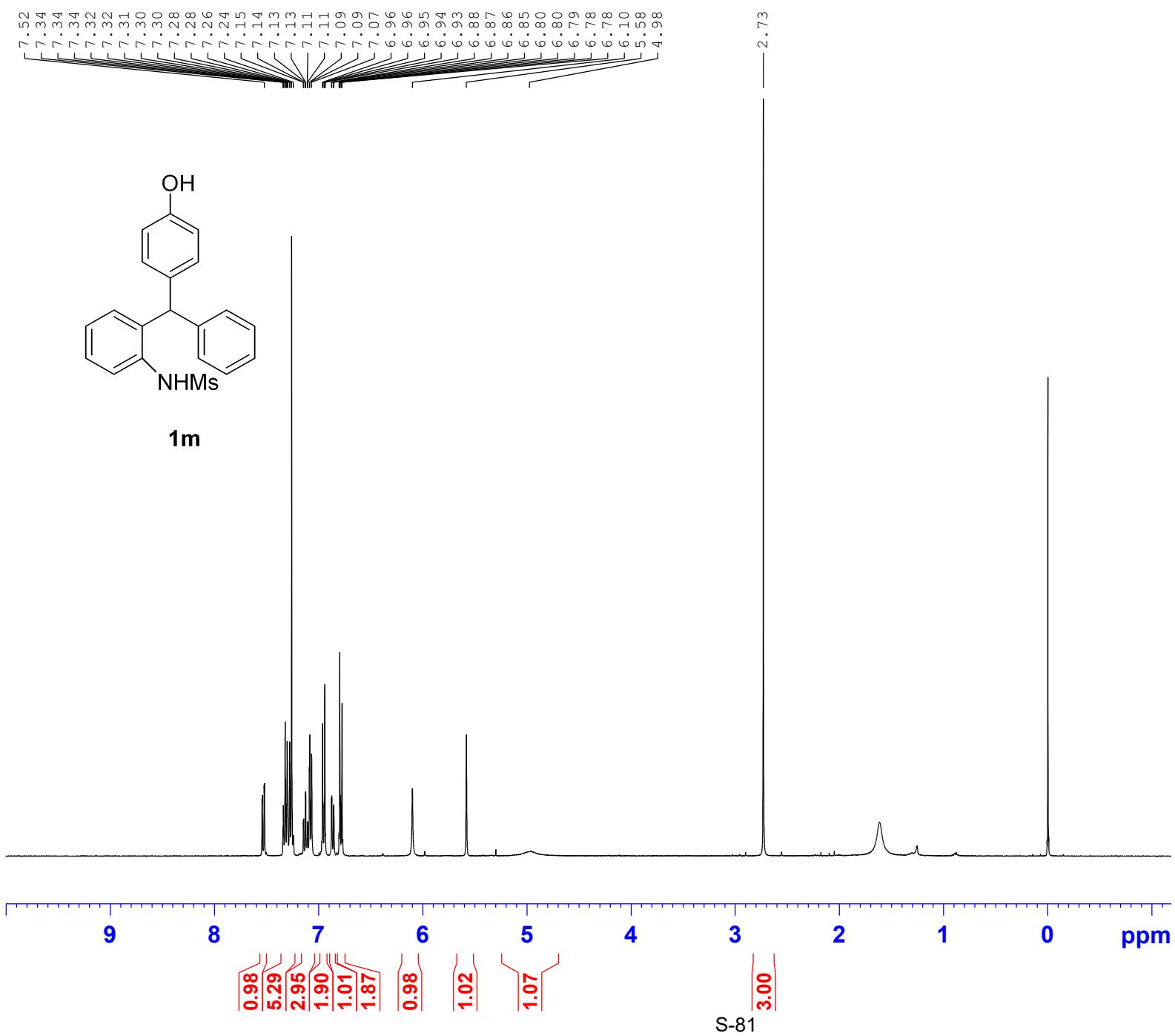


NAME Lzy-11
 EXPNO 2
 PROCNO 1
 Date 20200115
 Time 18.41
 INSTRUM spect
 PROBHD 5 mm PABBO BB-
 PULPROG zg30
 TD 65536
 SOLVENT CDCl3
 NS 16
 DS 2
 SWH 8223.685 Hz
 FIDRES 0.125483 Hz
 AQ 3.9846387 sec
 RG 64
 DW 60.800 usec
 DE 6.00 usec
 TE 294.7 K
 D1 1.00000000 sec
 TD0 1

===== CHANNEL f1 =====
 NUC1 1H
 P1 15.80 usec
 PL1 -1.00 dB
 PL1W 12.17476940 W
 SFO1 400.1324710 MHz
 SI 32768
 SF 400.1300091 MHz
 WDW EM
 SSB 0
 LB 0.30 Hz
 GB 0
 PC 1.00



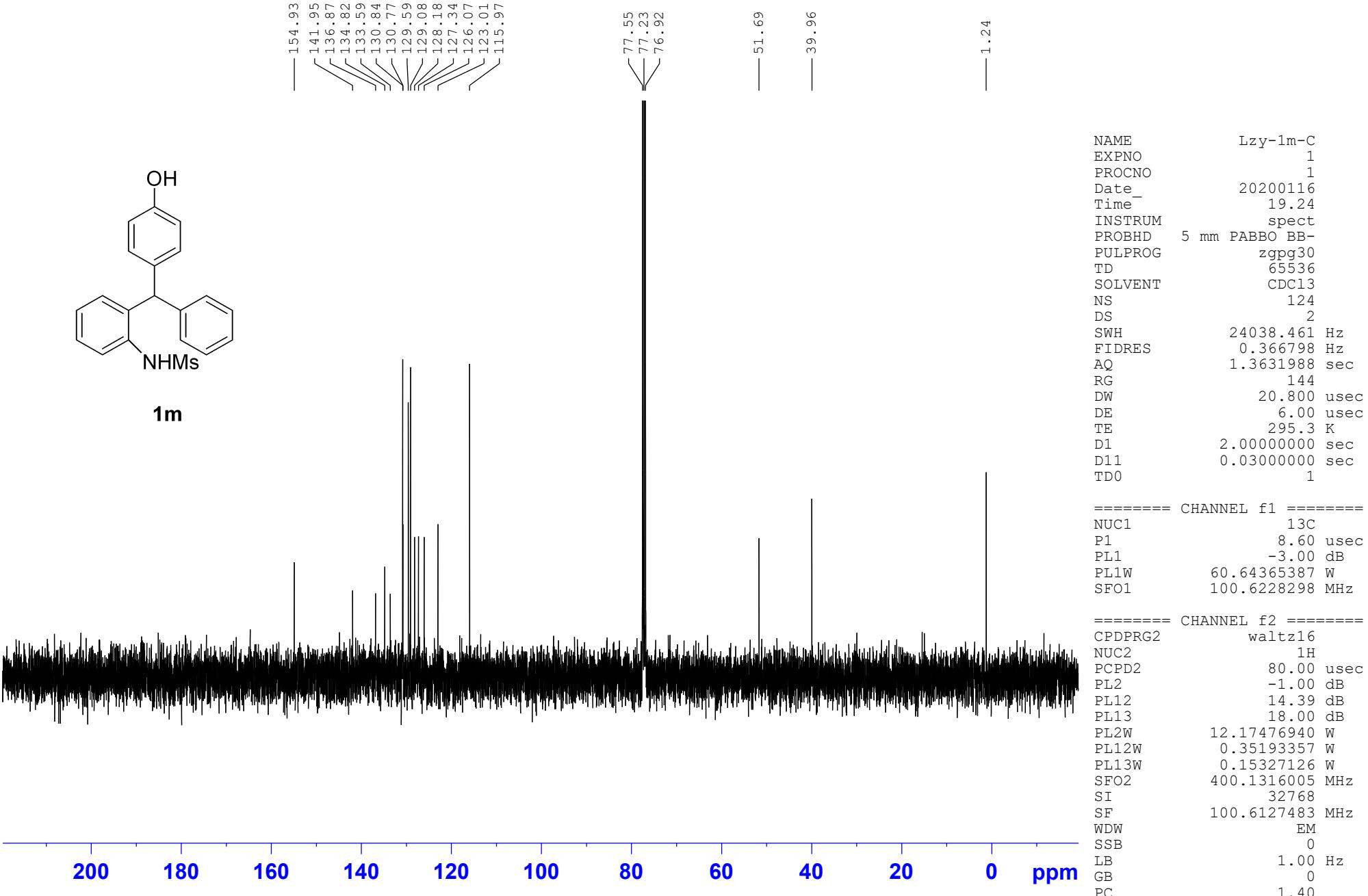


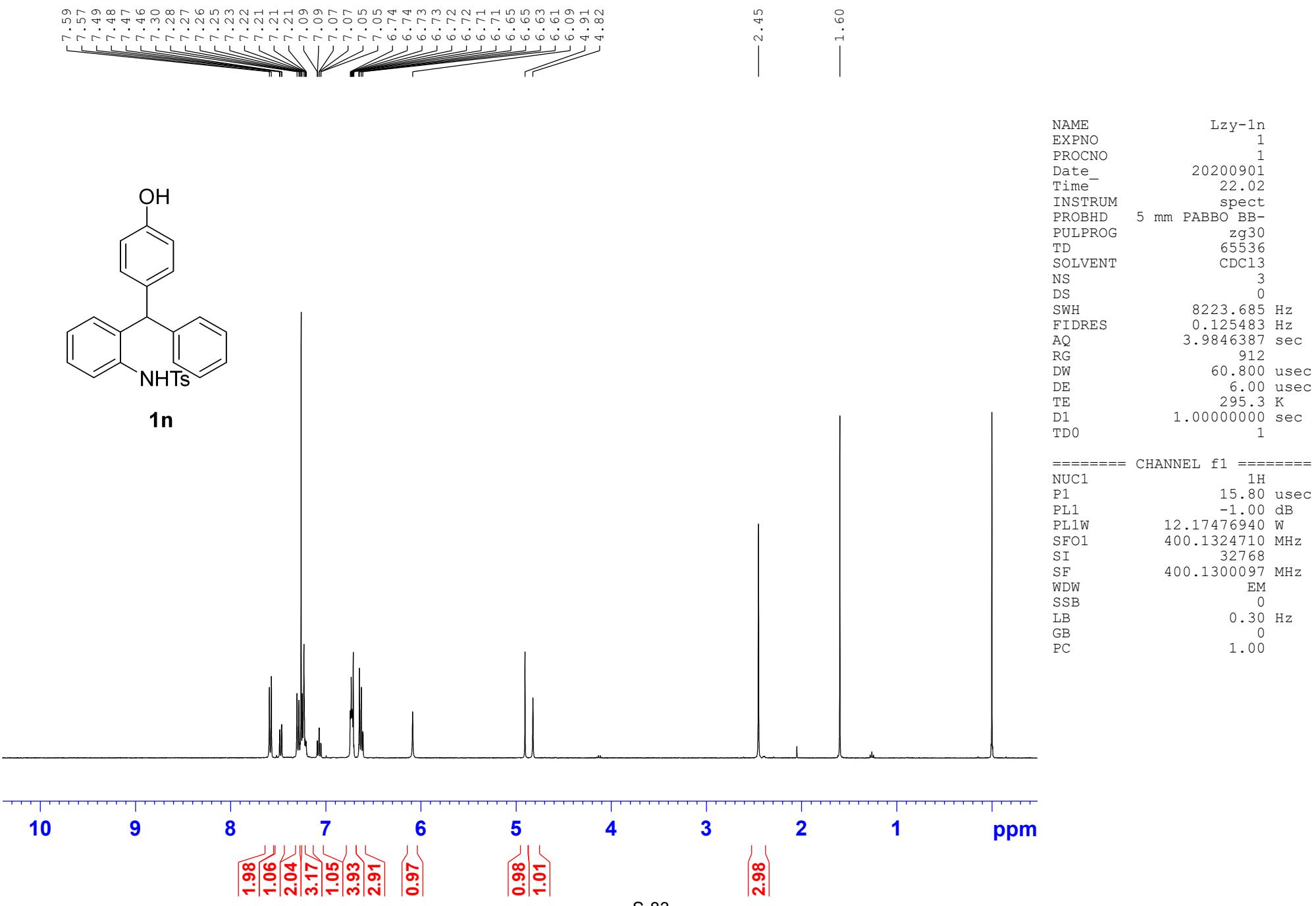


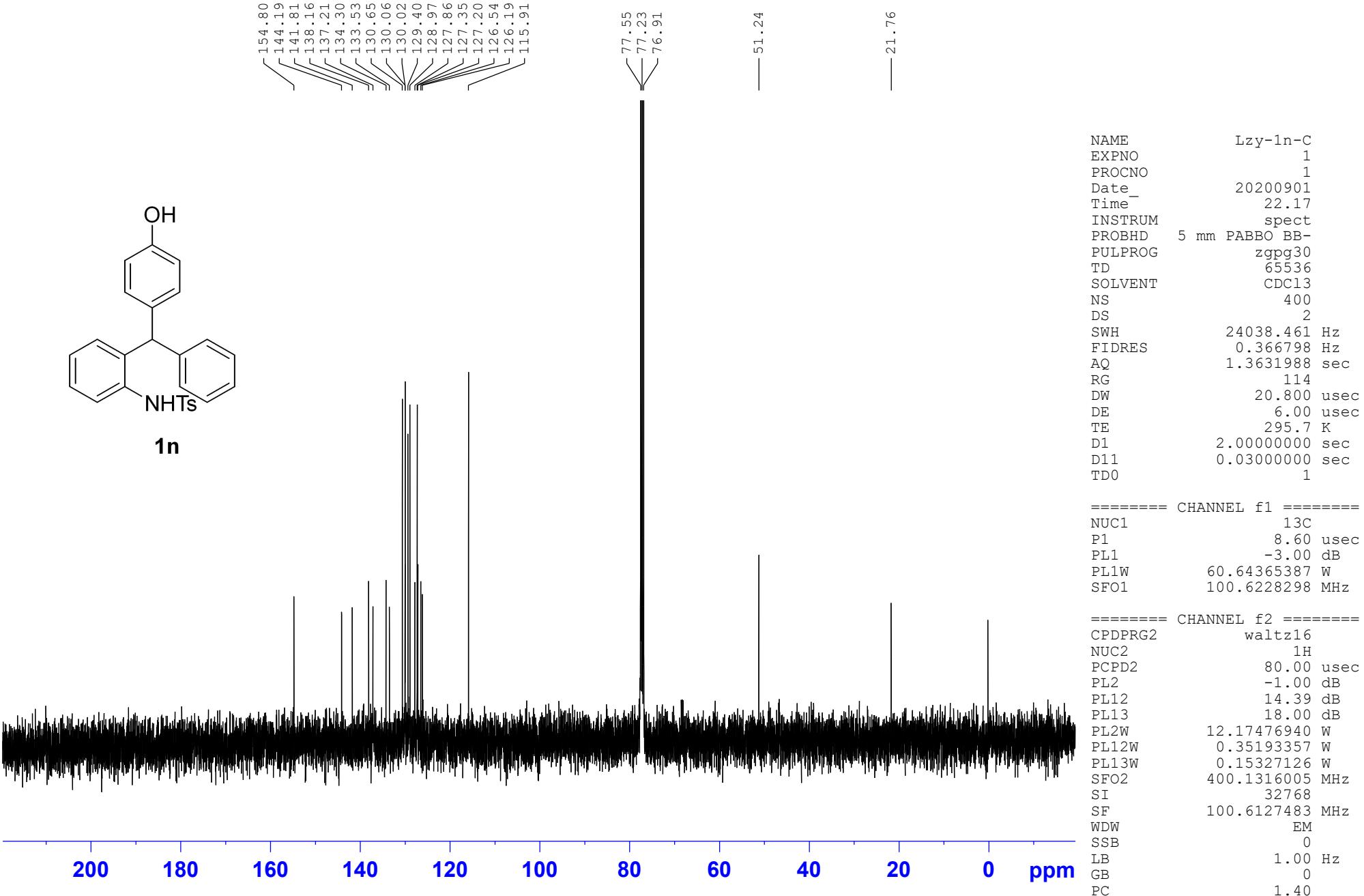
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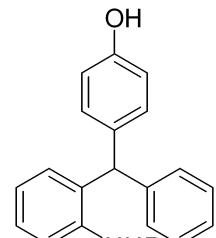
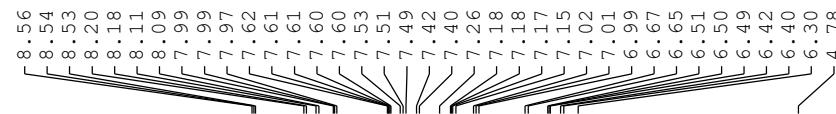
NAME      lzy-project1-1m-old
EXPNO      1
PROCNO     1
Date_      20210512
Time_      13.57
INSTRUM   spect
PROBHD   5 mm DUL 13C-1
PULPROG  zg30
TD        65536
SOLVENT   CDC13
NS         8
DS         0
SWH       8223.685 Hz
FIDRES   0.125483 Hz
AQ        3.9846387 sec
RG        575
DW        60.800 usec
DE        6.00  usec
TE        293.6 K
D1
TD0      1.00000000 sec

```



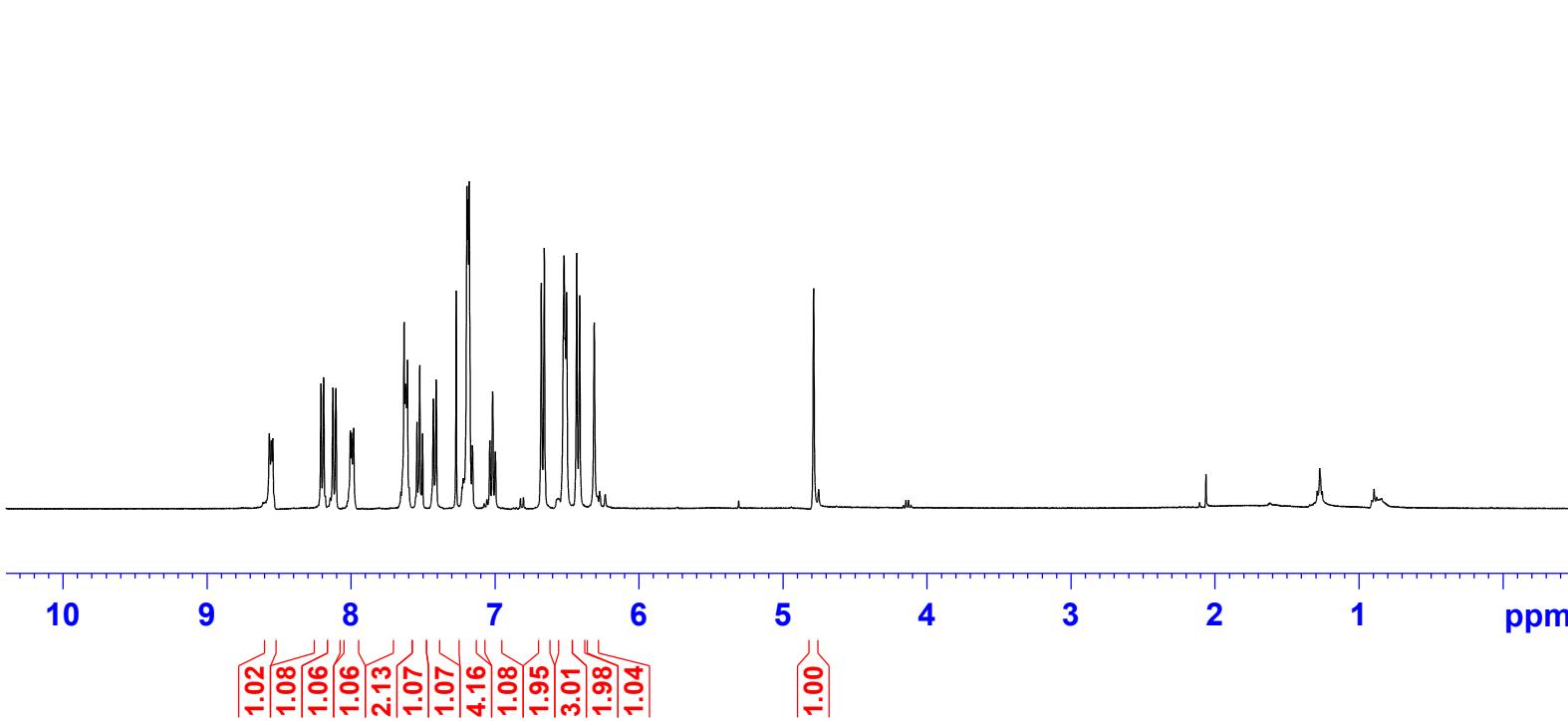






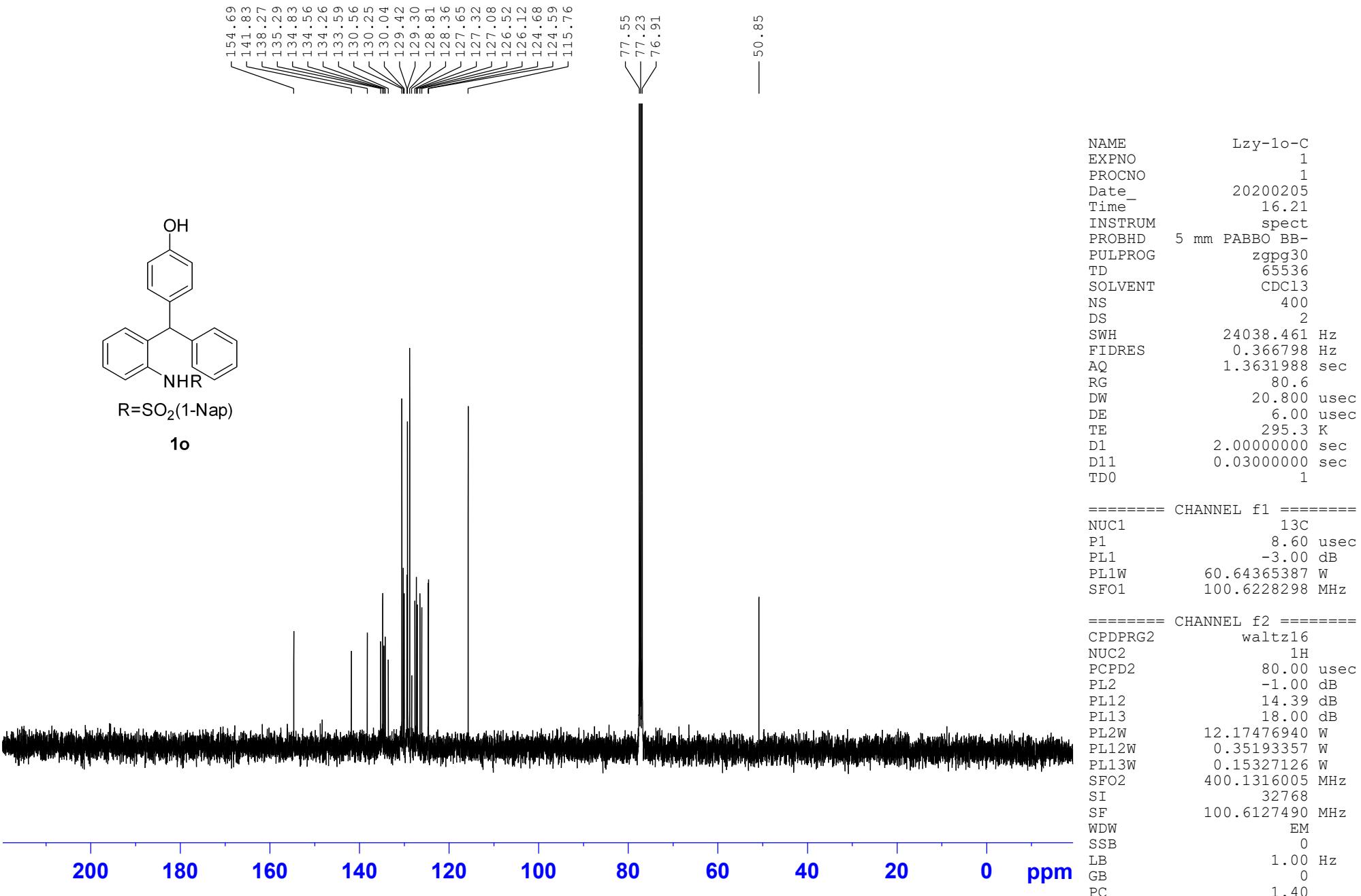
R=SO₂(1-Nap)

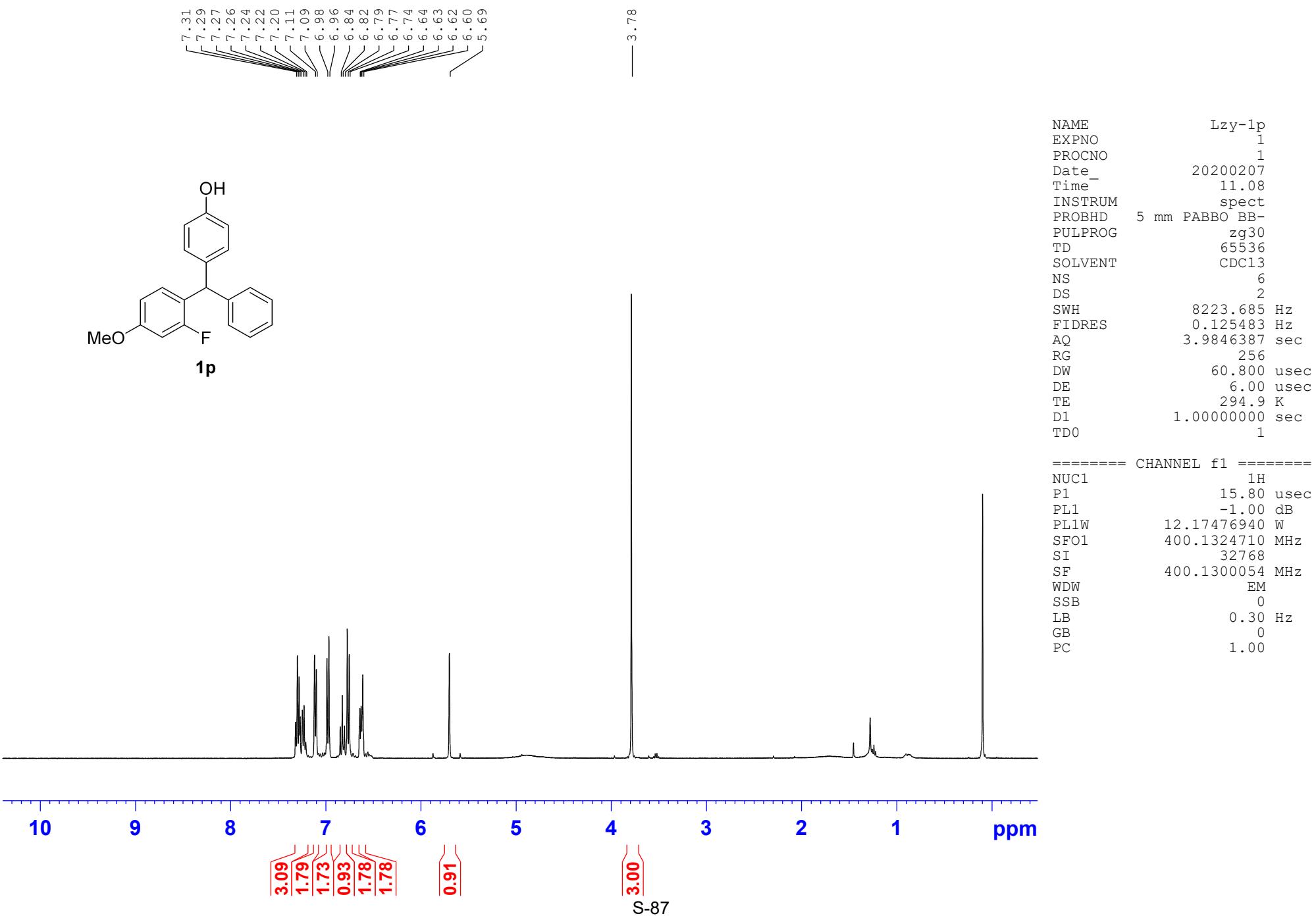
1o

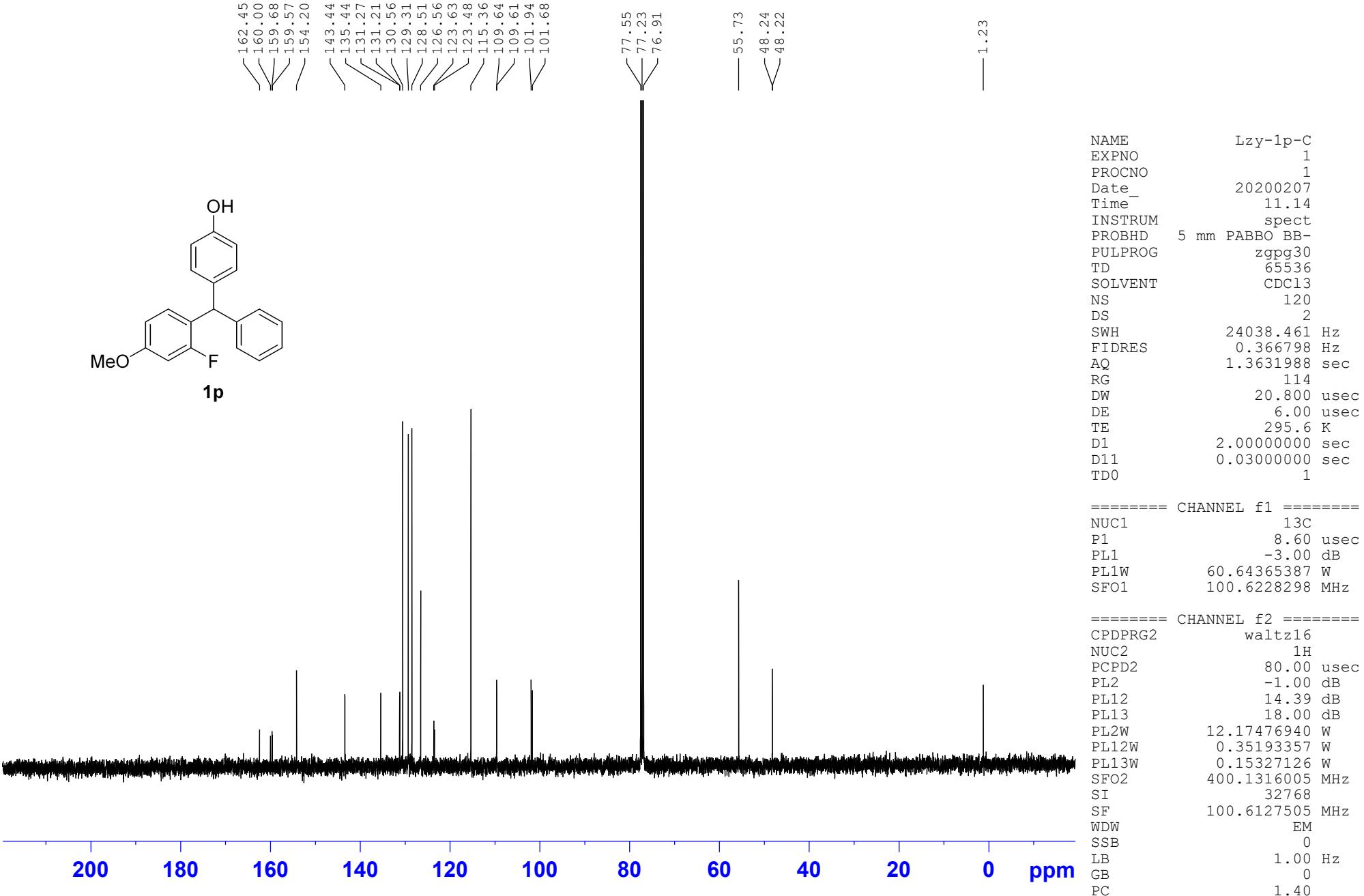


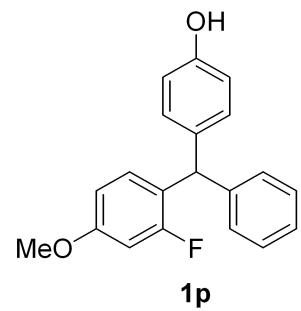
NAME	Lzy-lo
EXPNO	1
PROCNO	1
Date	20200205
Time	16.11
INSTRUM	spect
PROBHD	5 mm PABBO BB-
PULPROG	zg30
TD	65536
SOLVENT	CDCl ₃
NS	12
DS	2
SWH	8223.685 Hz
FIDRES	0.125483 Hz
AQ	3.9846387 sec
RG	512
DW	60.800 usec
DE	6.00 usec
TE	294.7 K
D1	1.00000000 sec
TD0	1

===== CHANNEL f1 =====	
NUC1	1H
P1	15.80 usec
PL1	-1.00 dB
PL1W	12.17476940 W
SFO1	400.1324710 MHz
SI	32768
SF	400.1300054 MHz
WDW	EM
SSB	0
LB	0.30 Hz
GB	0
PC	1.00







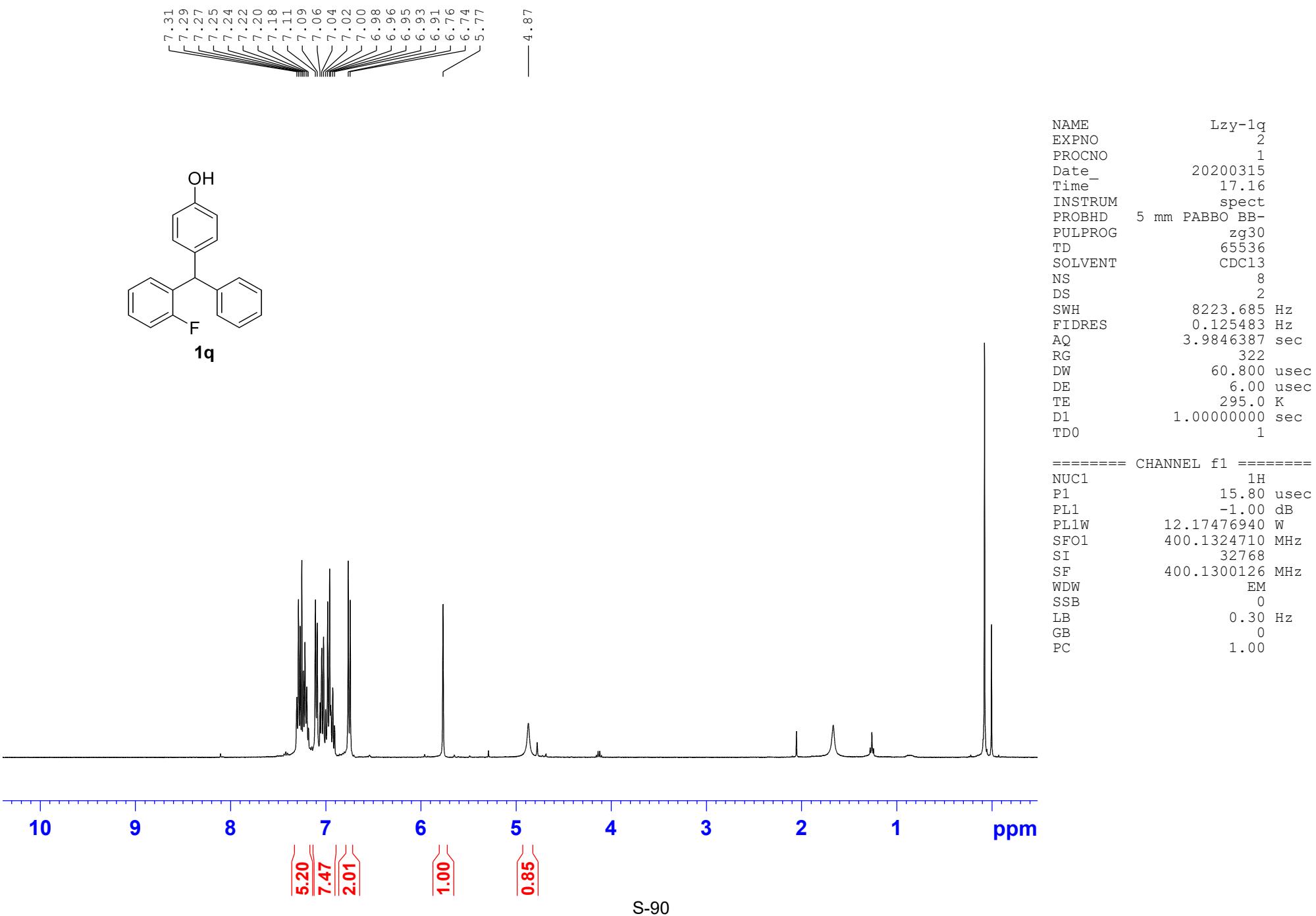


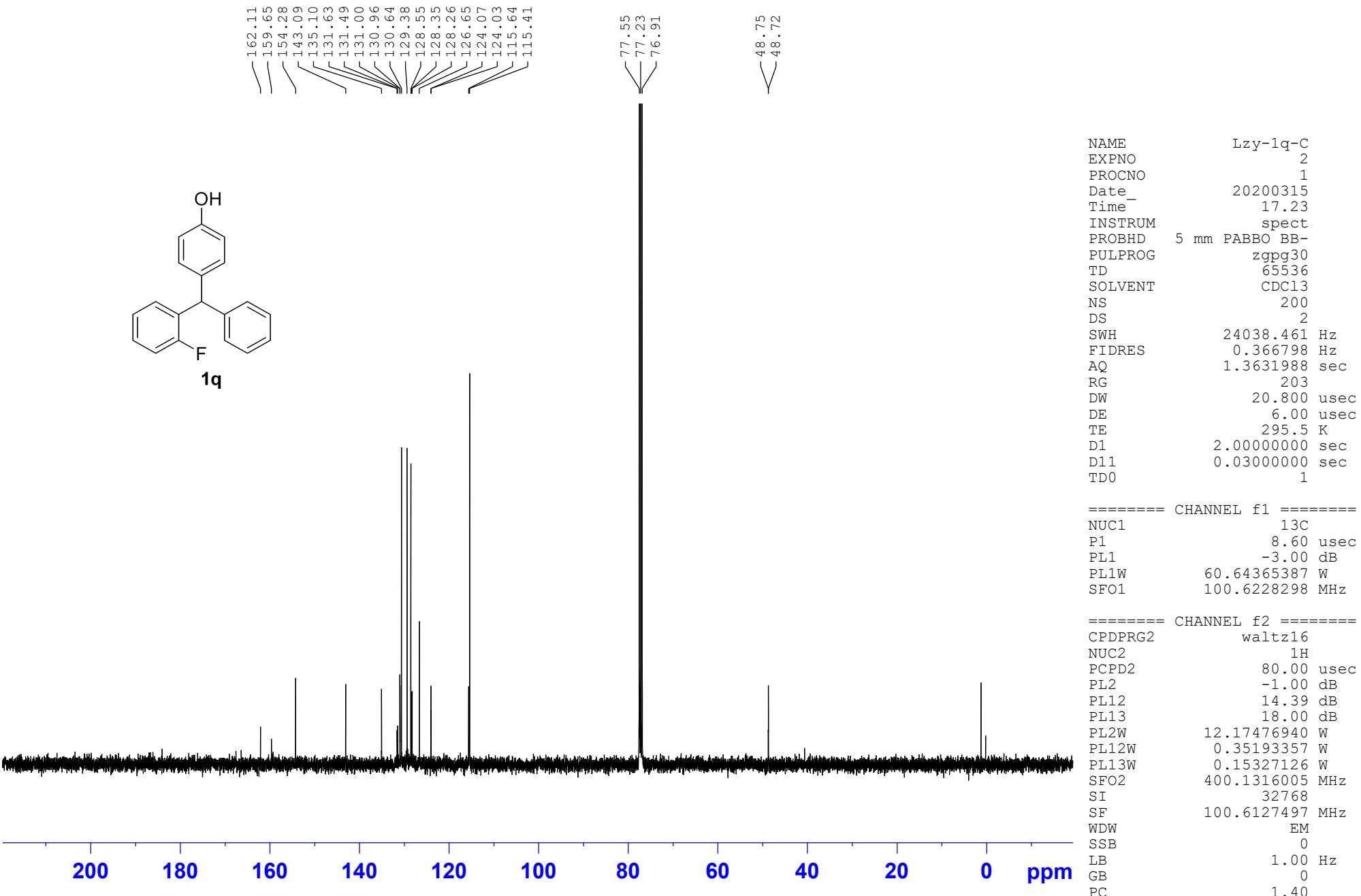
-114.41

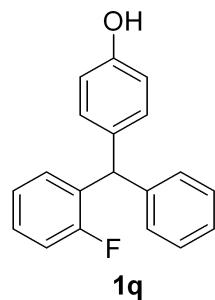
NAME Lzy-1p-F-new
 EXPNO 1
 PROCNO 1
 Date 20200207
 Time 14.00
 INSTRUM spect
 PROBHD 5 mm PABBO BB-
 PULPROG zgflqn
 TD 131072
 SOLVENT CDC13
 NS 16
 DS 4
 SWH 89285.711 Hz
 FIDRES 0.681196 Hz
 AQ 0.7340532 sec
 RG 2050
 DW 5.600 usec
 DE 6.00 usec
 TE 294.9 K
 D1 1.0000000 sec
 TDO 1

===== CHANNEL f1 =====
 NUC1 19F
 P1 19.50 usec
 PL1 -4.00 dB
 PL1W 16.97275162 W
 SFO1 376.4607164 MHz
 SI 65536
 SF 376.4983660 MHz
 WDW EM
 SSB 0
 LB 0.30 Hz
 GB 0
 PC 1.00

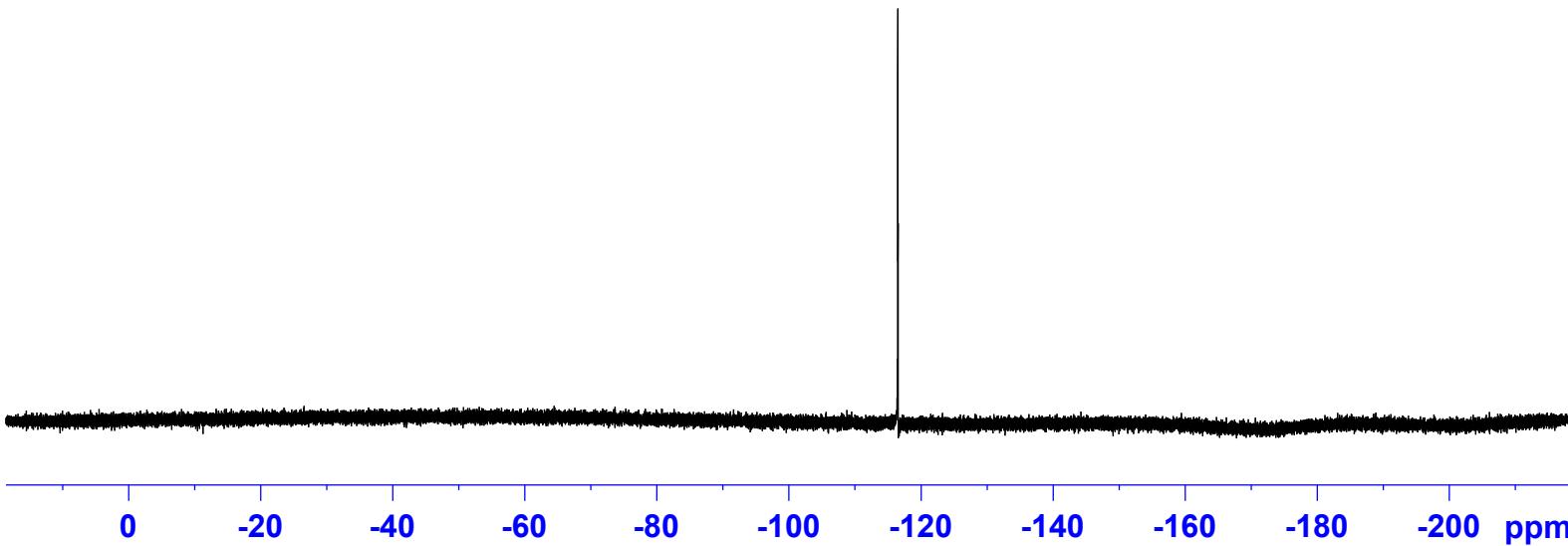
0 -20 -40 -60 -80 -100 -120 -140 -160 -180 -200 ppm



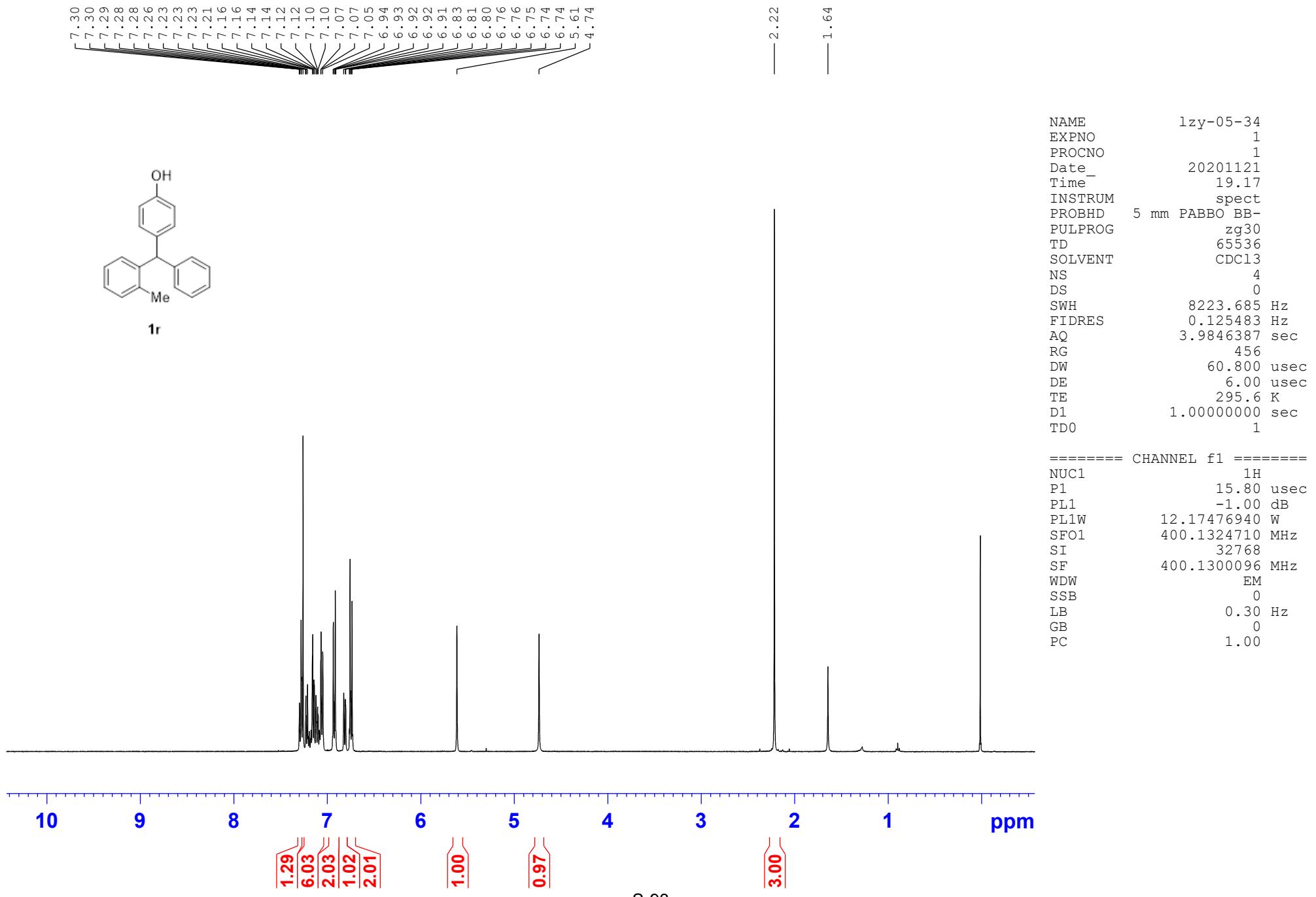




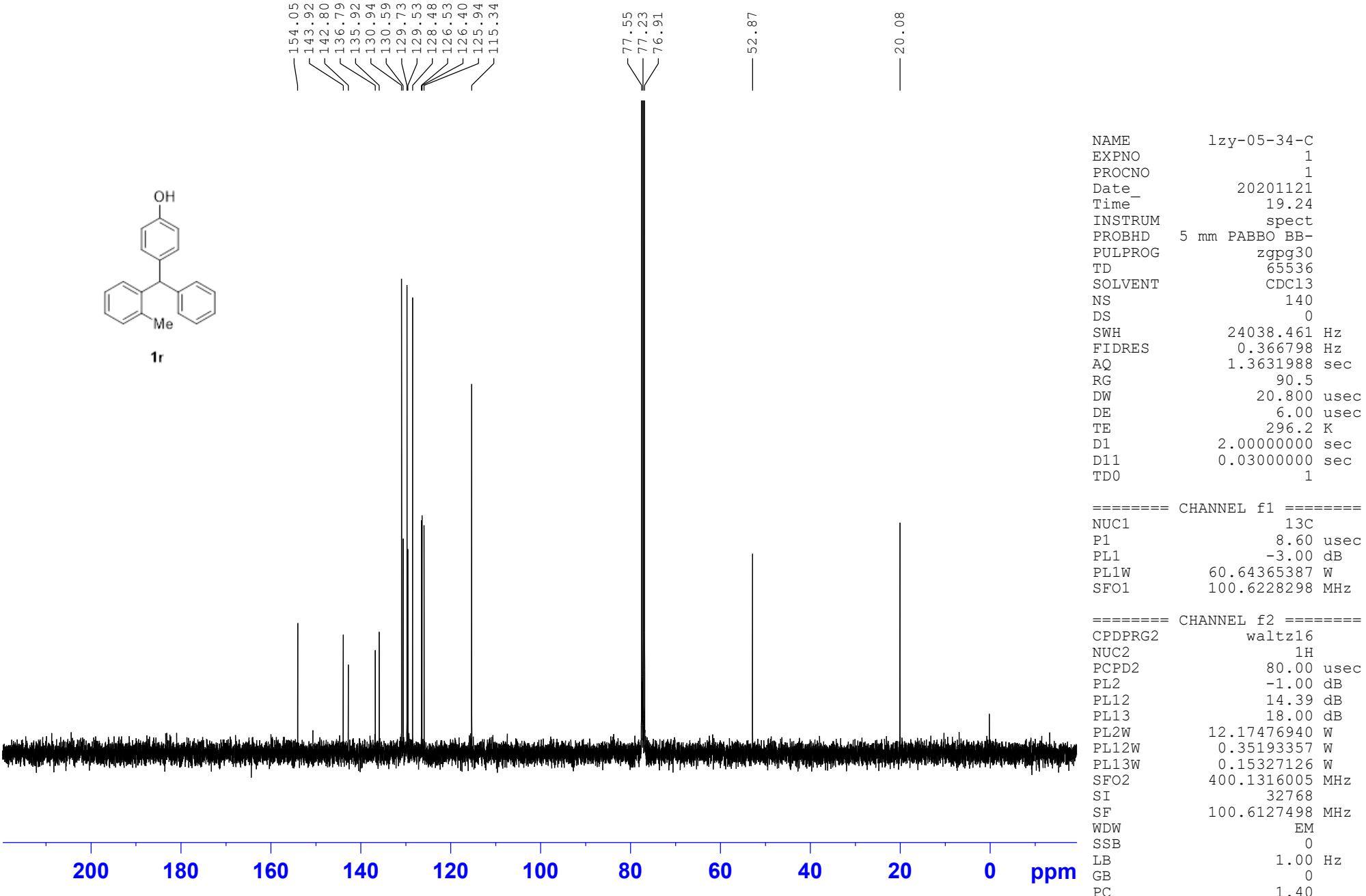
-116.48

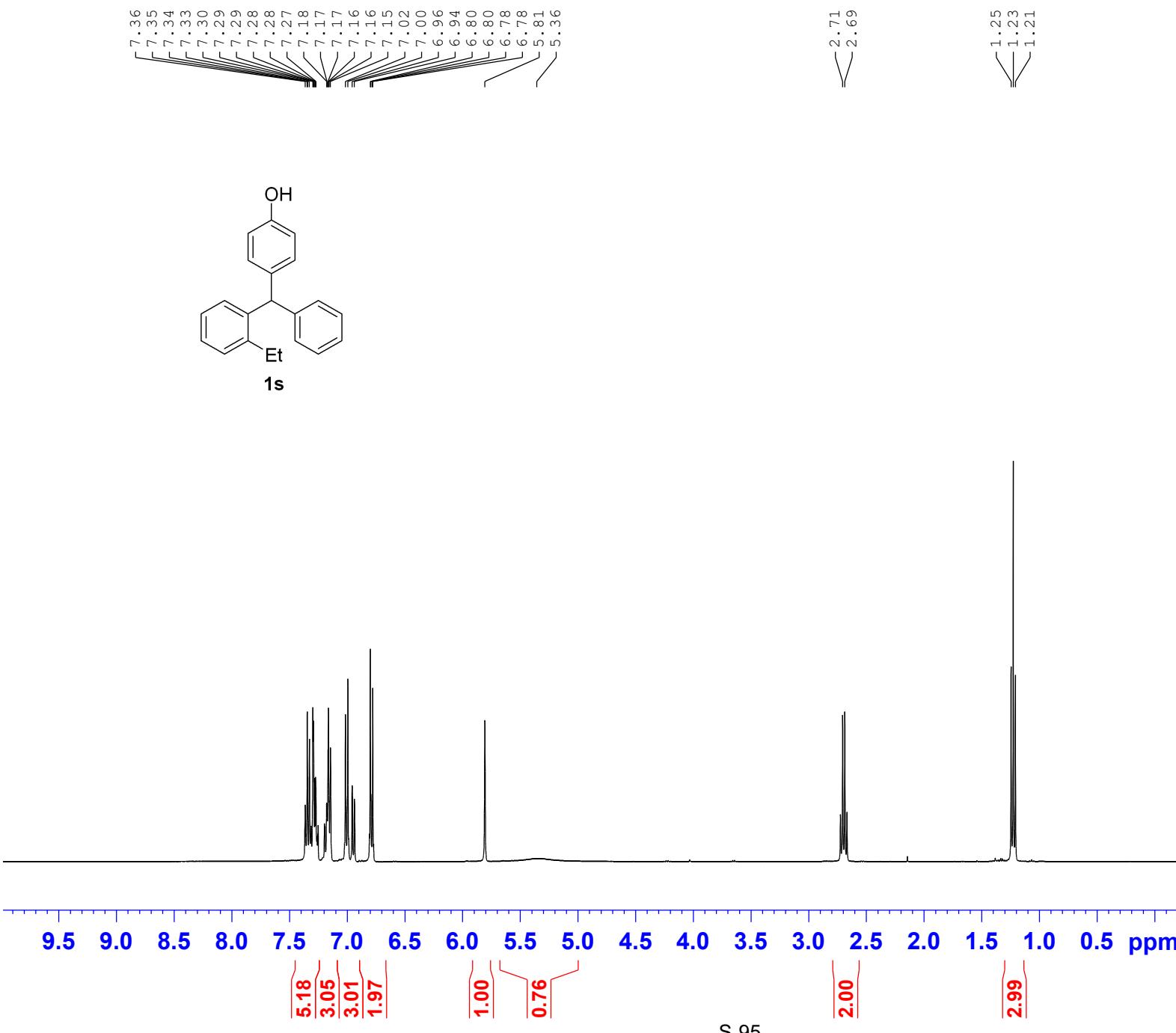


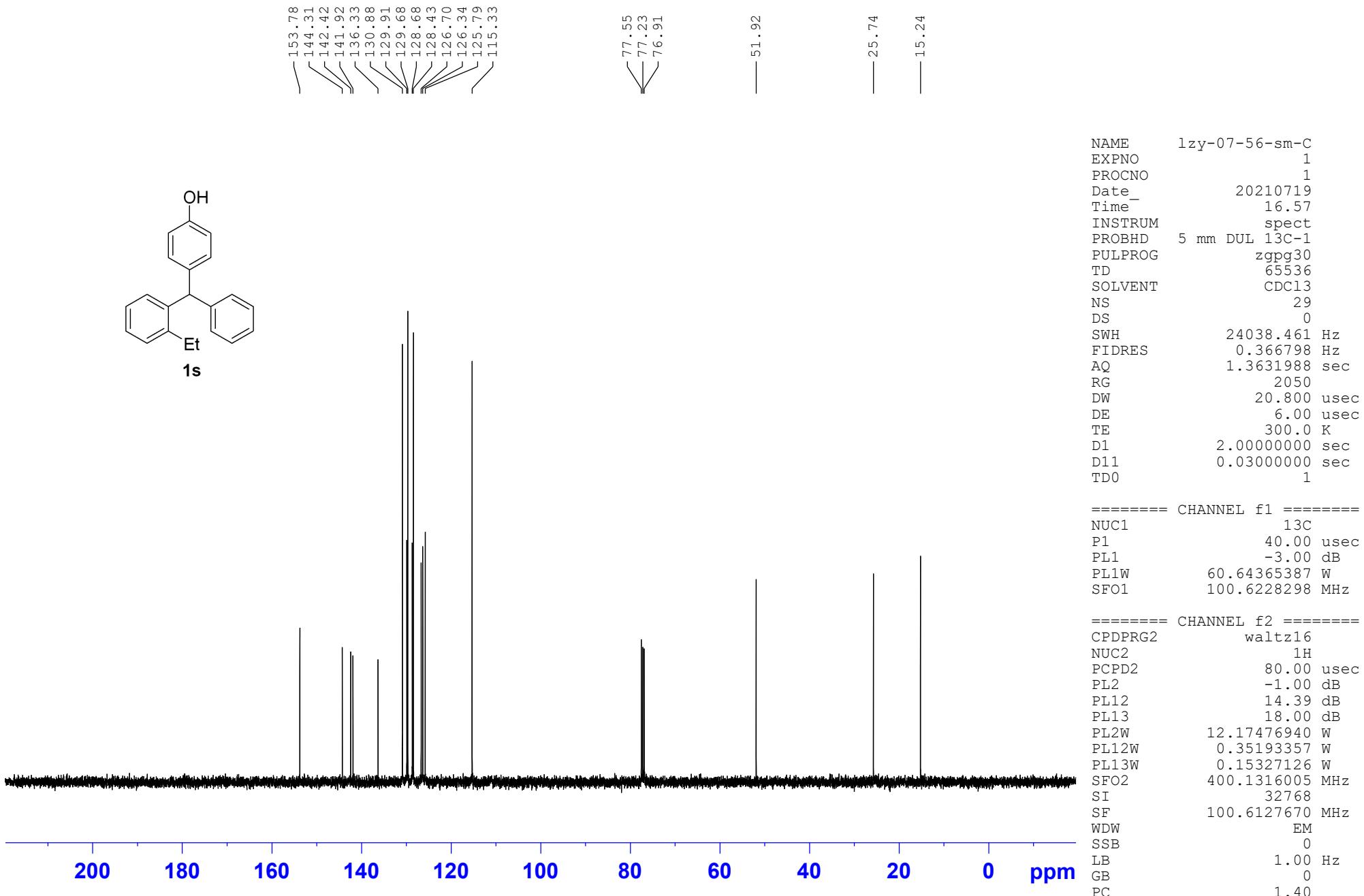
NAME	lzy-project1-1q-F
EXPNO	3
PROCNO	1
Date	20210717
Time	20.57
INSTRUM	spect
PROBHD	5 mm PABBO BB/
PULPROG	zgflqn
TD	131072
SOLVENT	CDC13
NS	5
DS	4
SWH	89285.711 Hz
FIDRES	0.681196 Hz
AQ	0.7340532 sec
RG	196.92
DW	5.600 usec
DE	6.50 usec
TE	296.7 K
D1	1.00000000 sec
TD0	1
===== CHANNEL f1 =====	
SFO1	376.4607164 MHz
NUC1	19F
P1	14.70 usec
SI	65536
SF	376.4983660 MHz
WDW	EM
SSB	0
LB	0.30 Hz
GB	0
PC	1.00

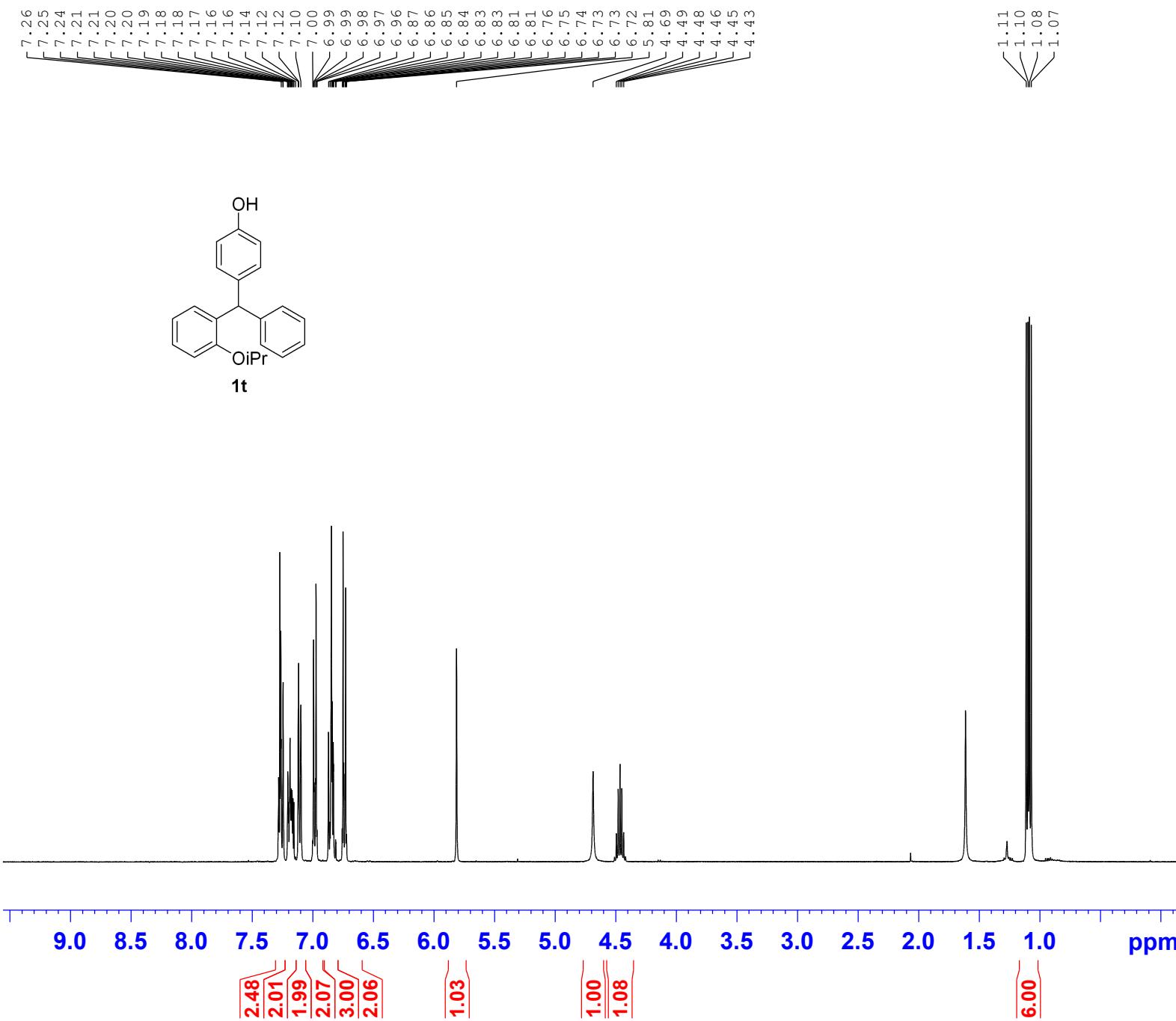


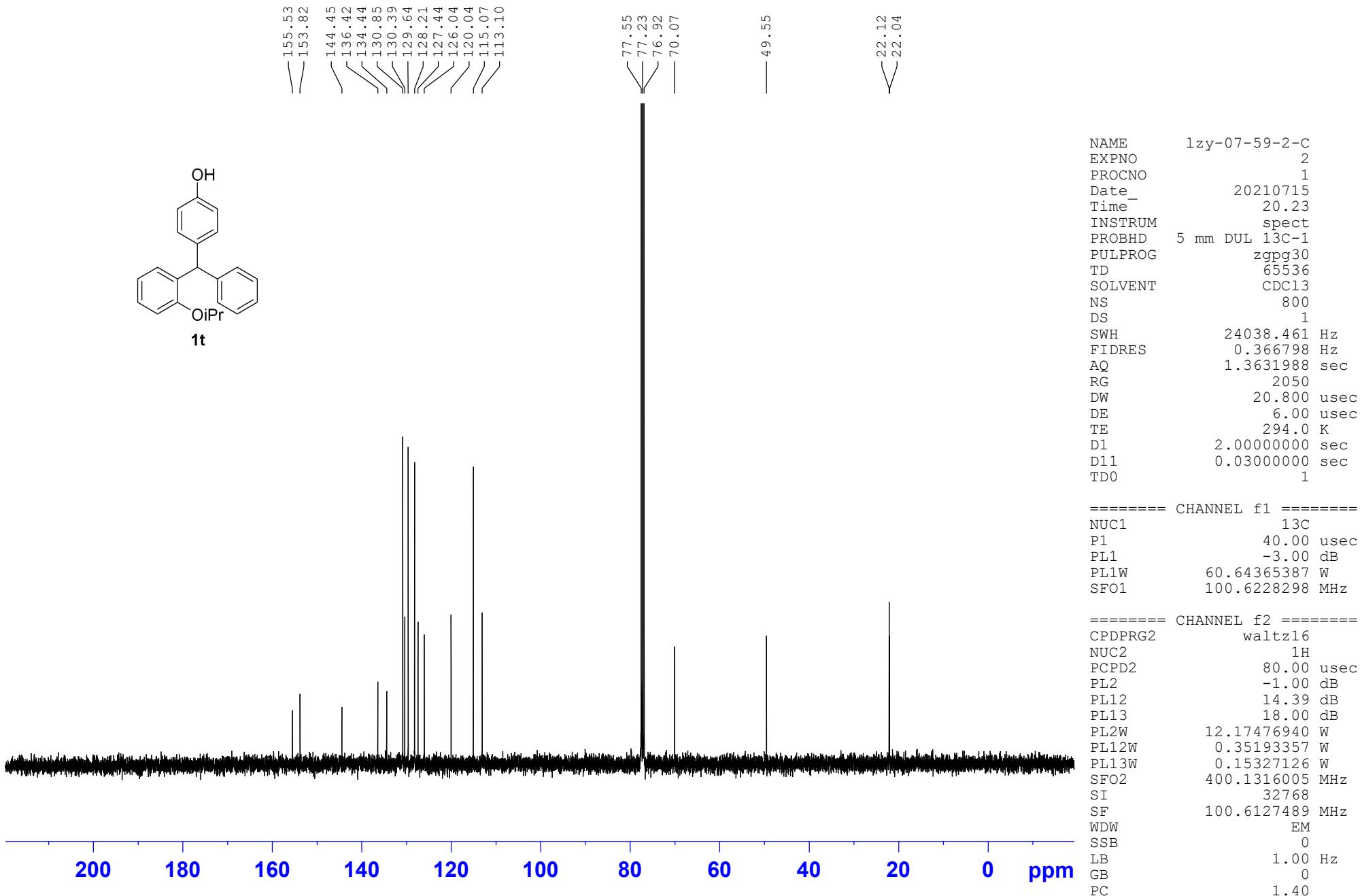
NAME	lzy-05-34
EXPNO	1
PROCNO	1
Date_	20201121
Time	19.17
INSTRUM	spect
PROBHD	5 mm PABBO BB-
PULPROG	zg30
TD	65536
SOLVENT	CDC13
NS	4
DS	0
SWH	8223.685 Hz
FIDRES	0.125483 Hz
AQ	3.9846387 sec
RG	456
DW	60.800 usec
DE	6.00 usec
TE	295.6 K
D1	1.00000000 sec
TDO	1

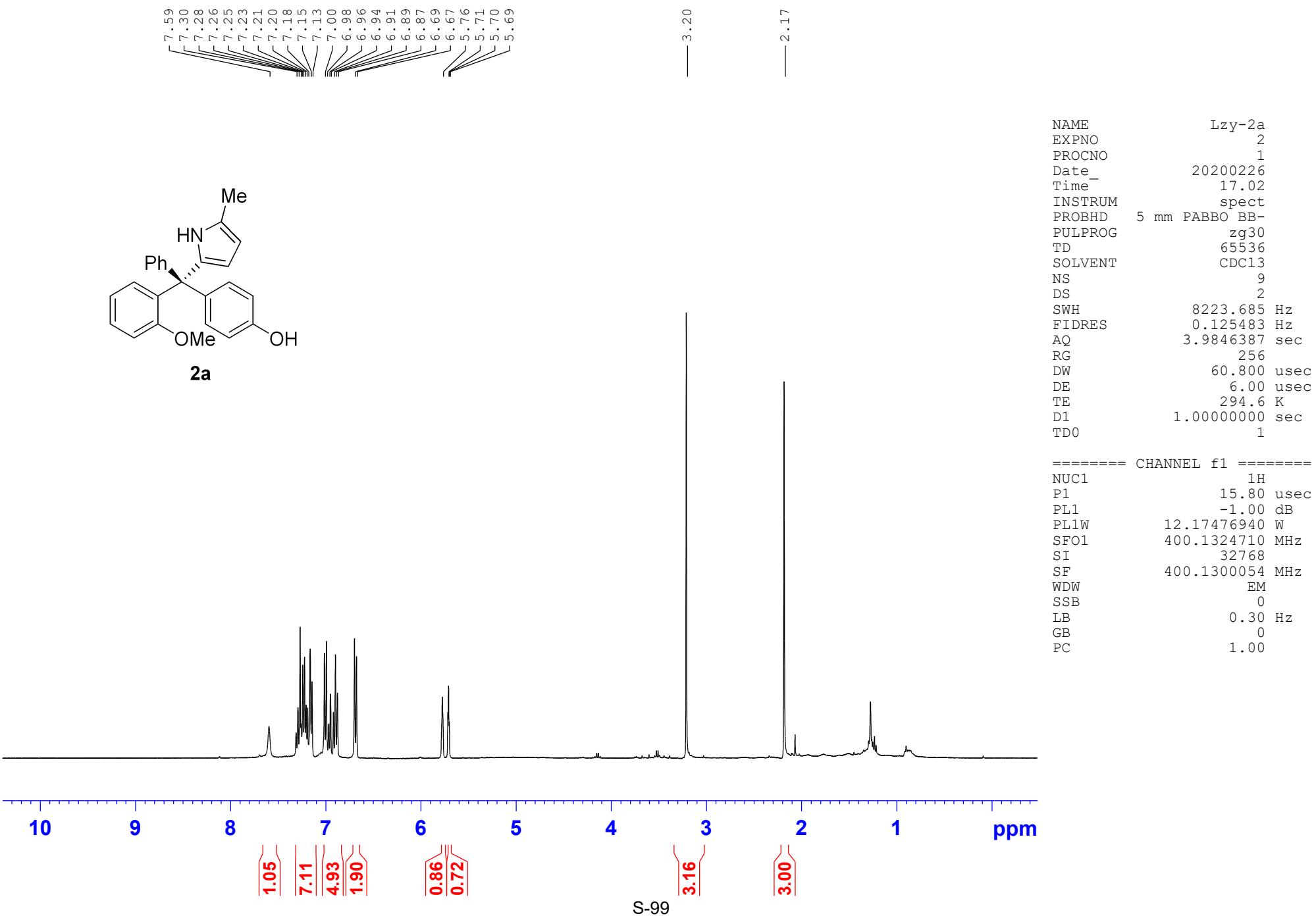


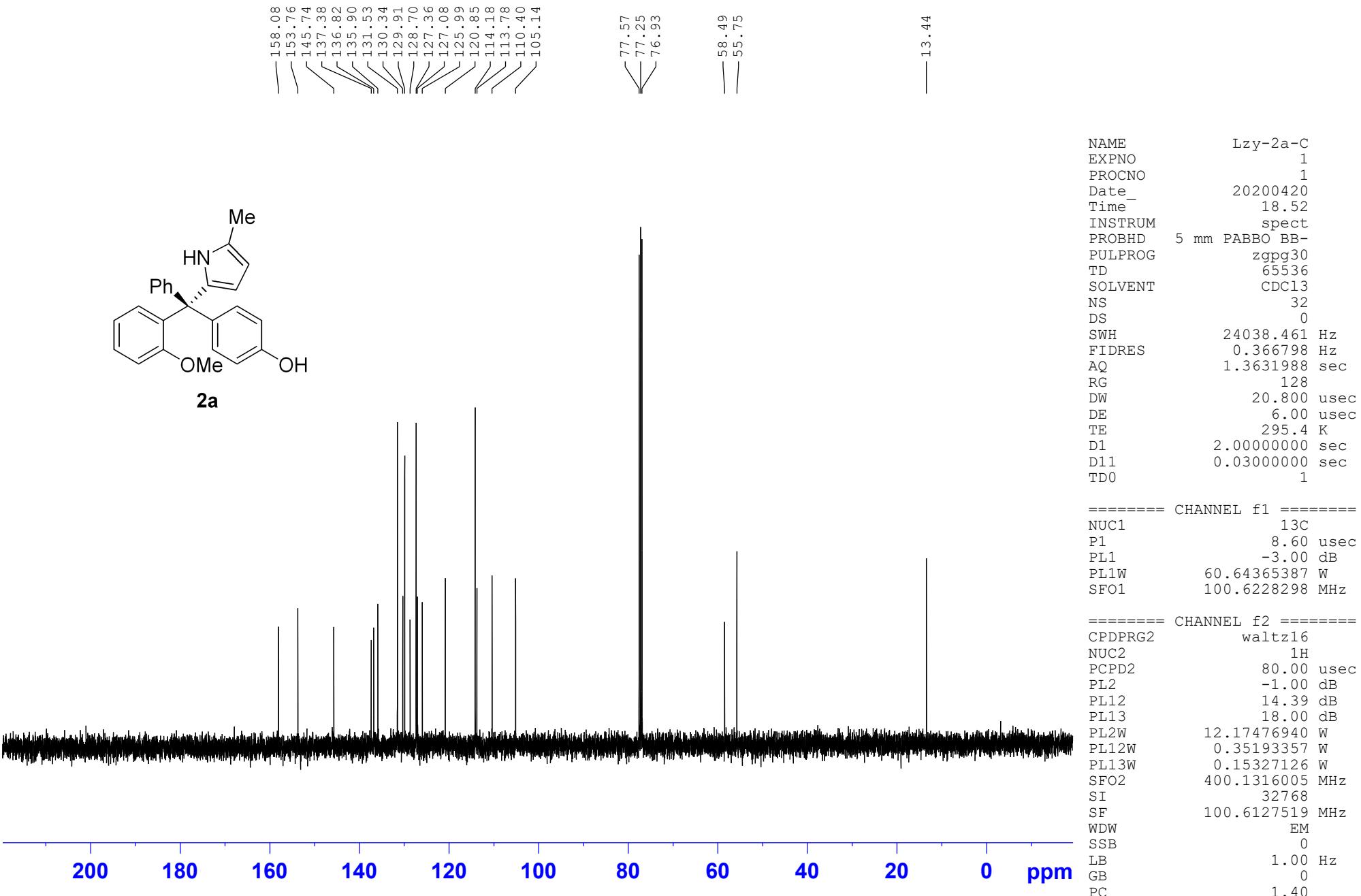


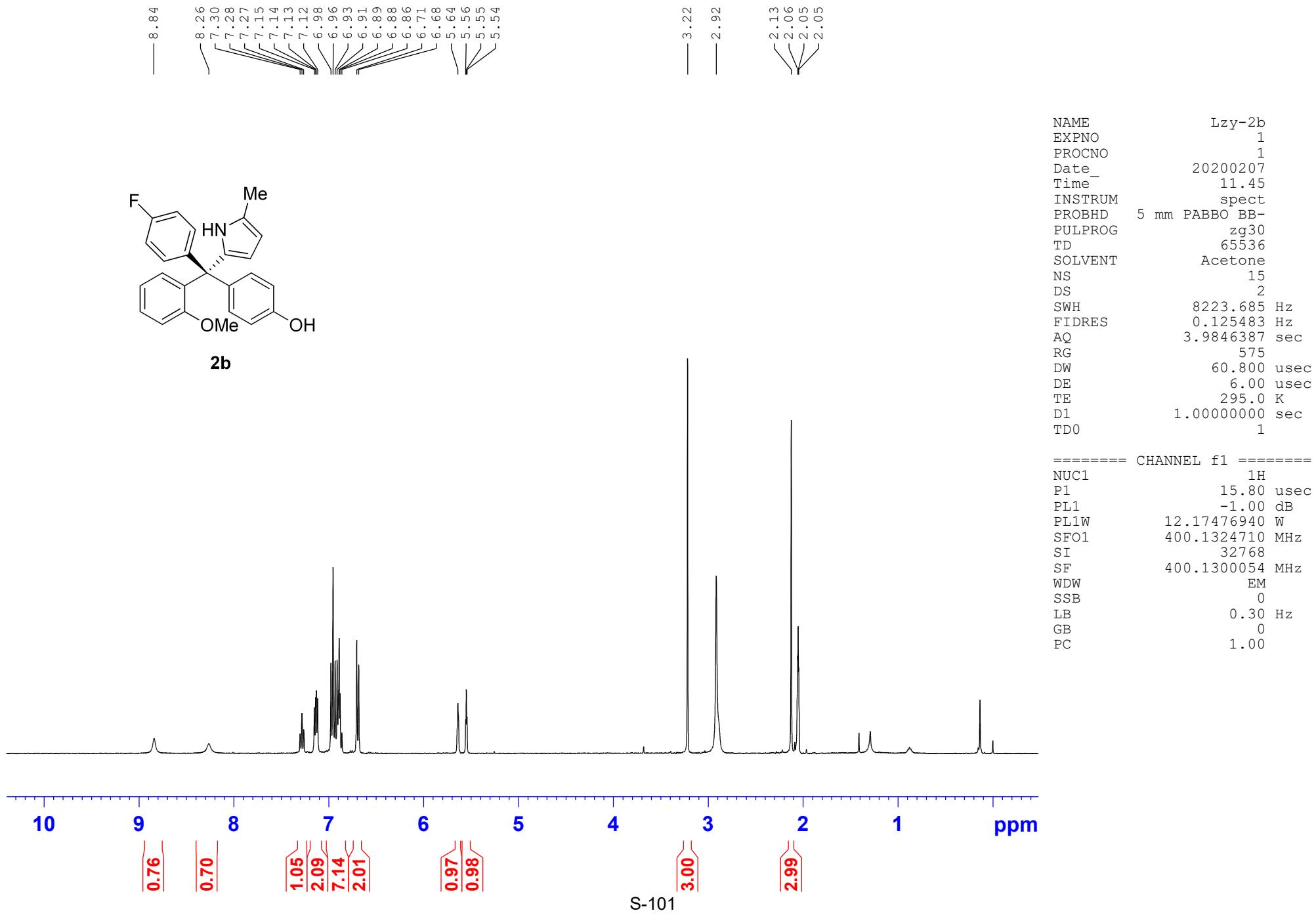


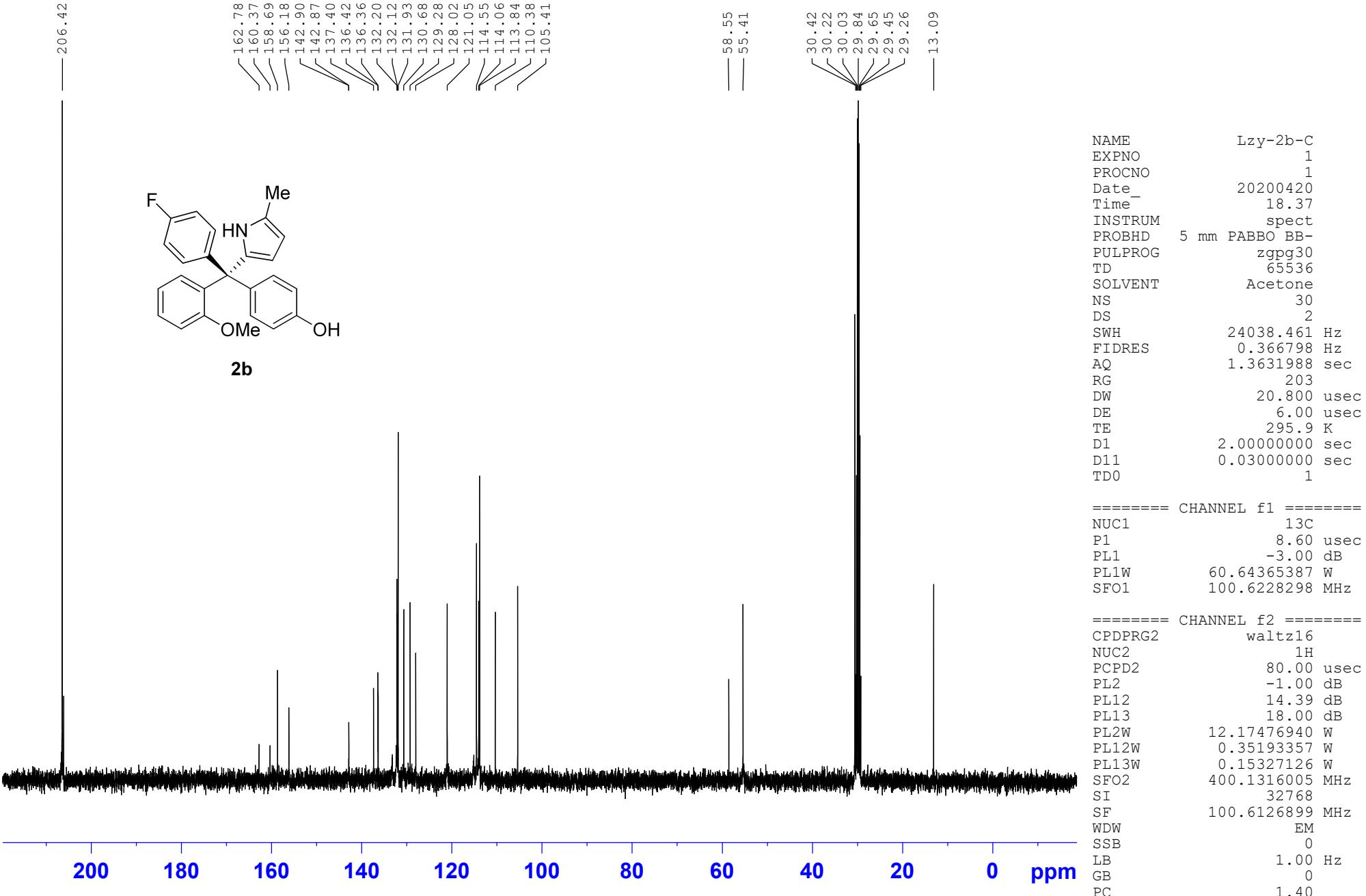


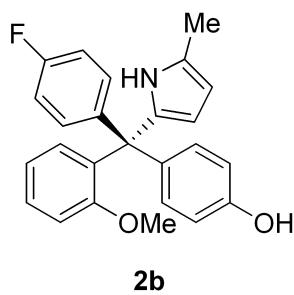








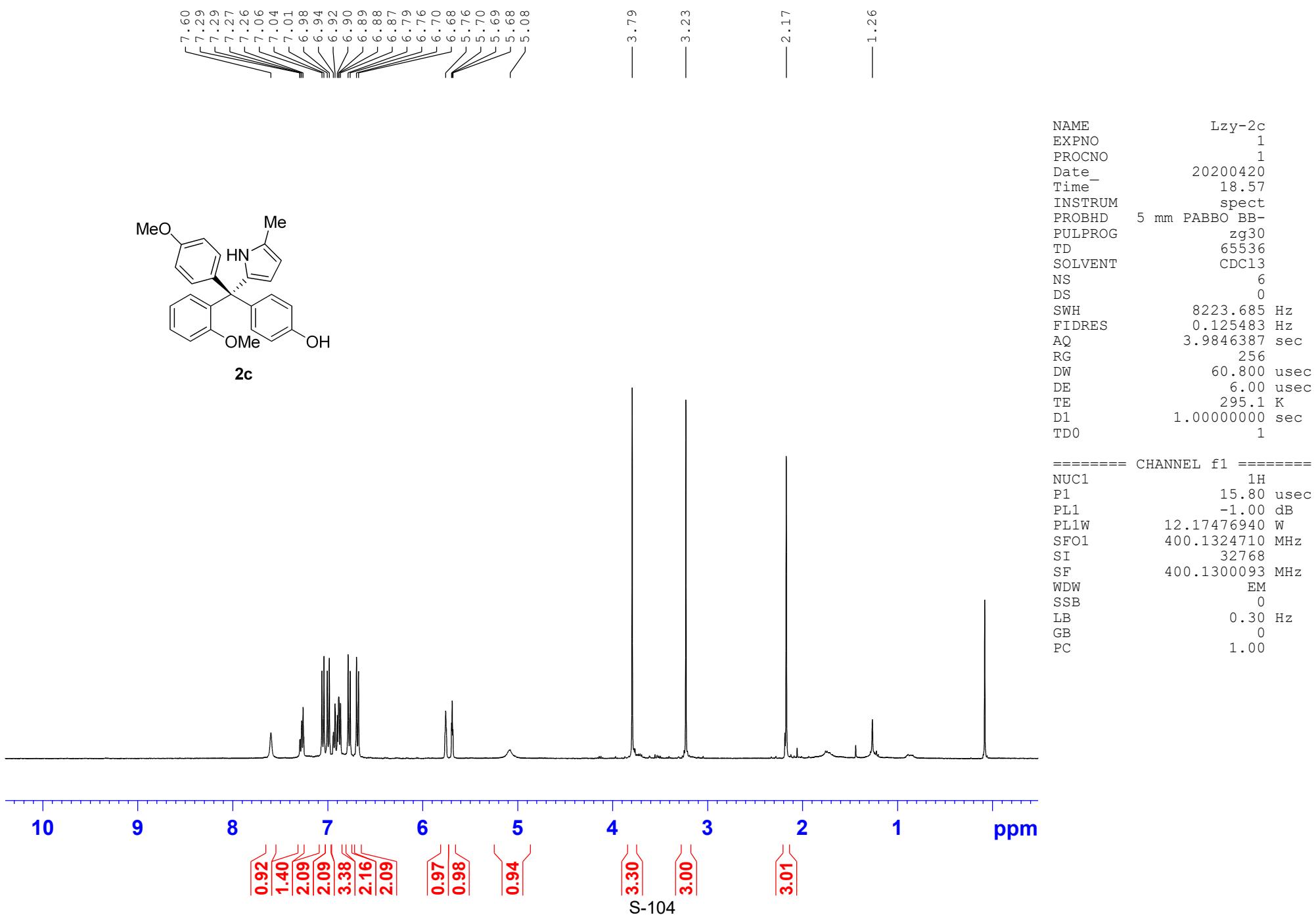


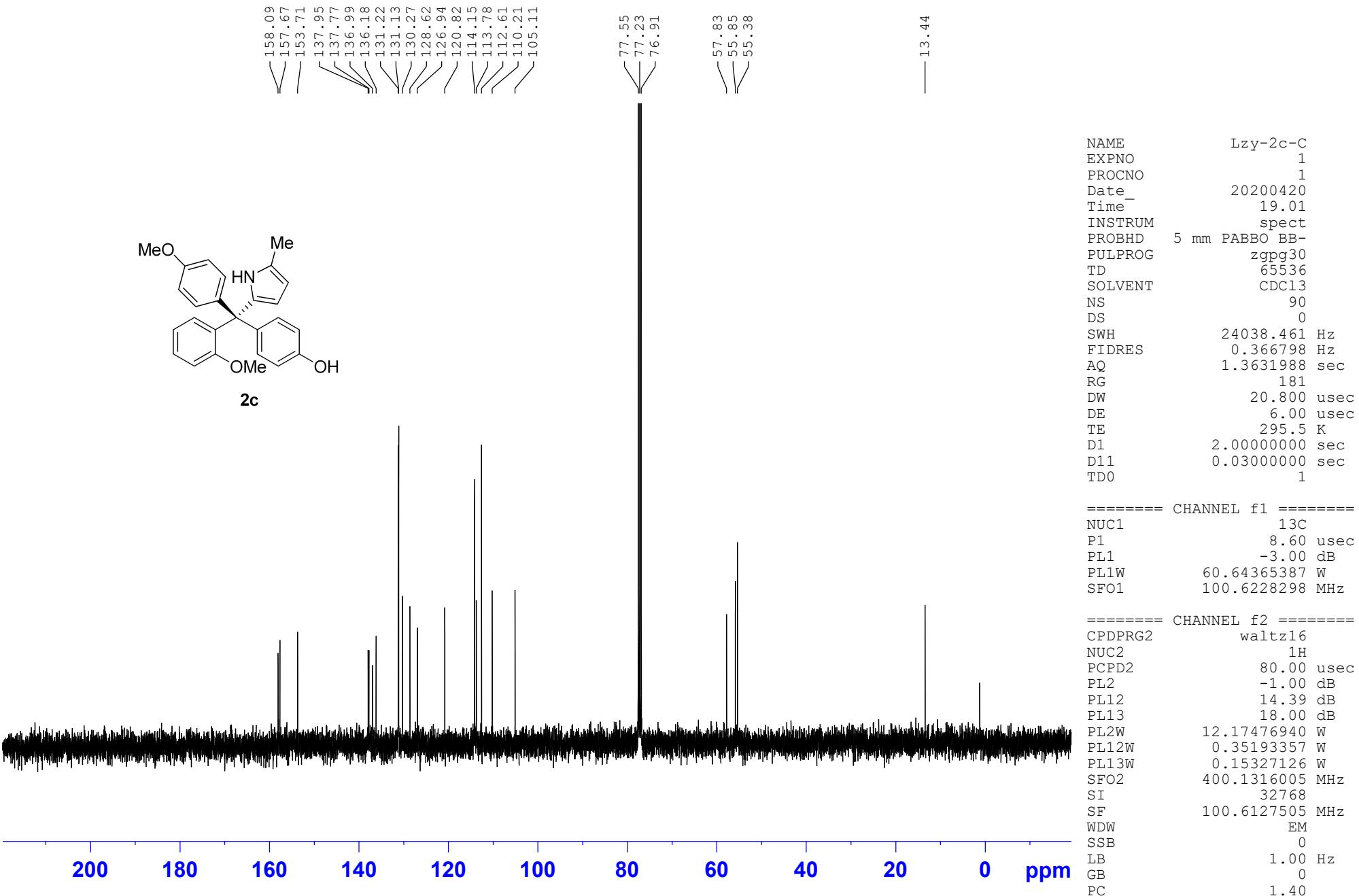


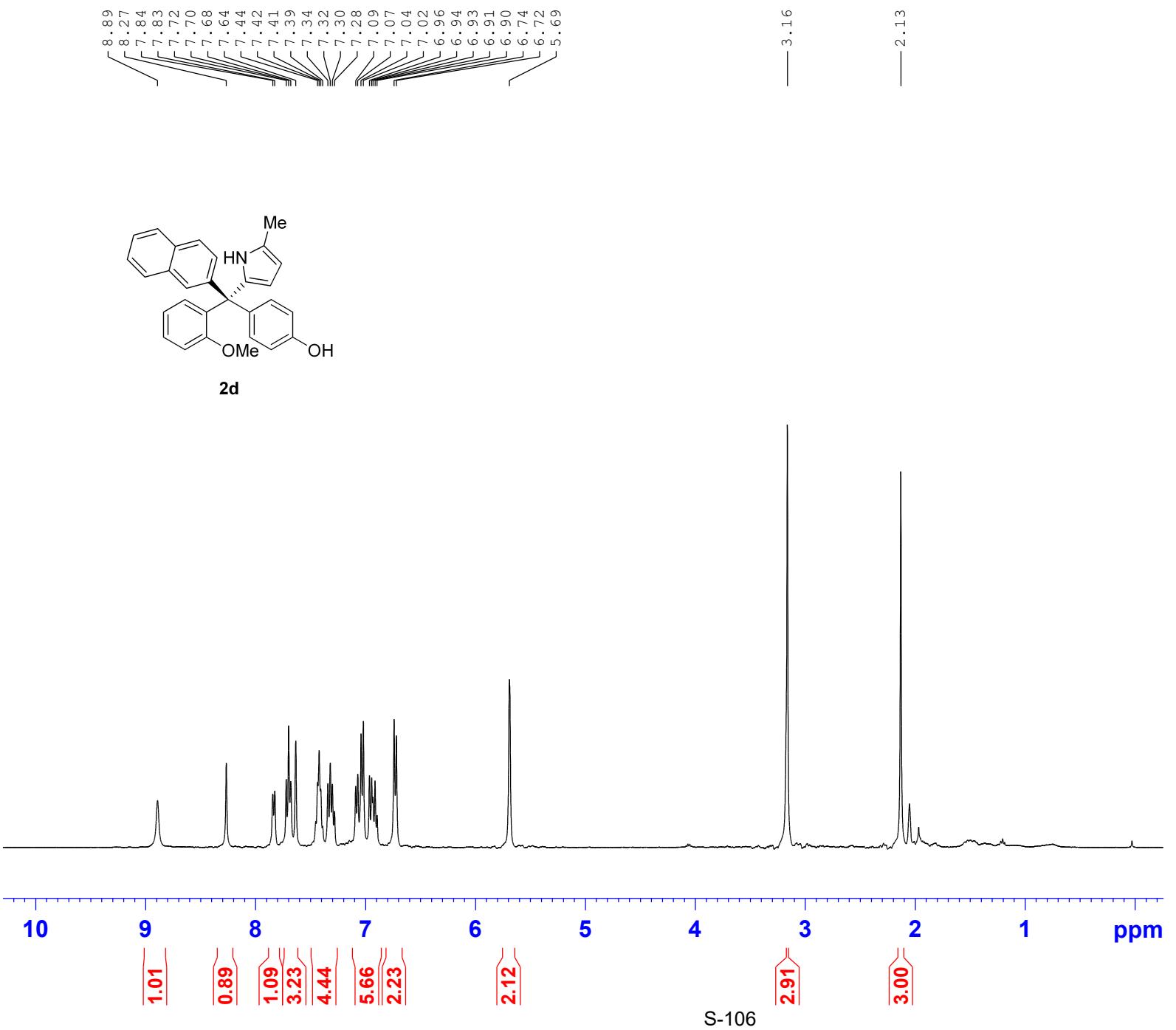
-119.82

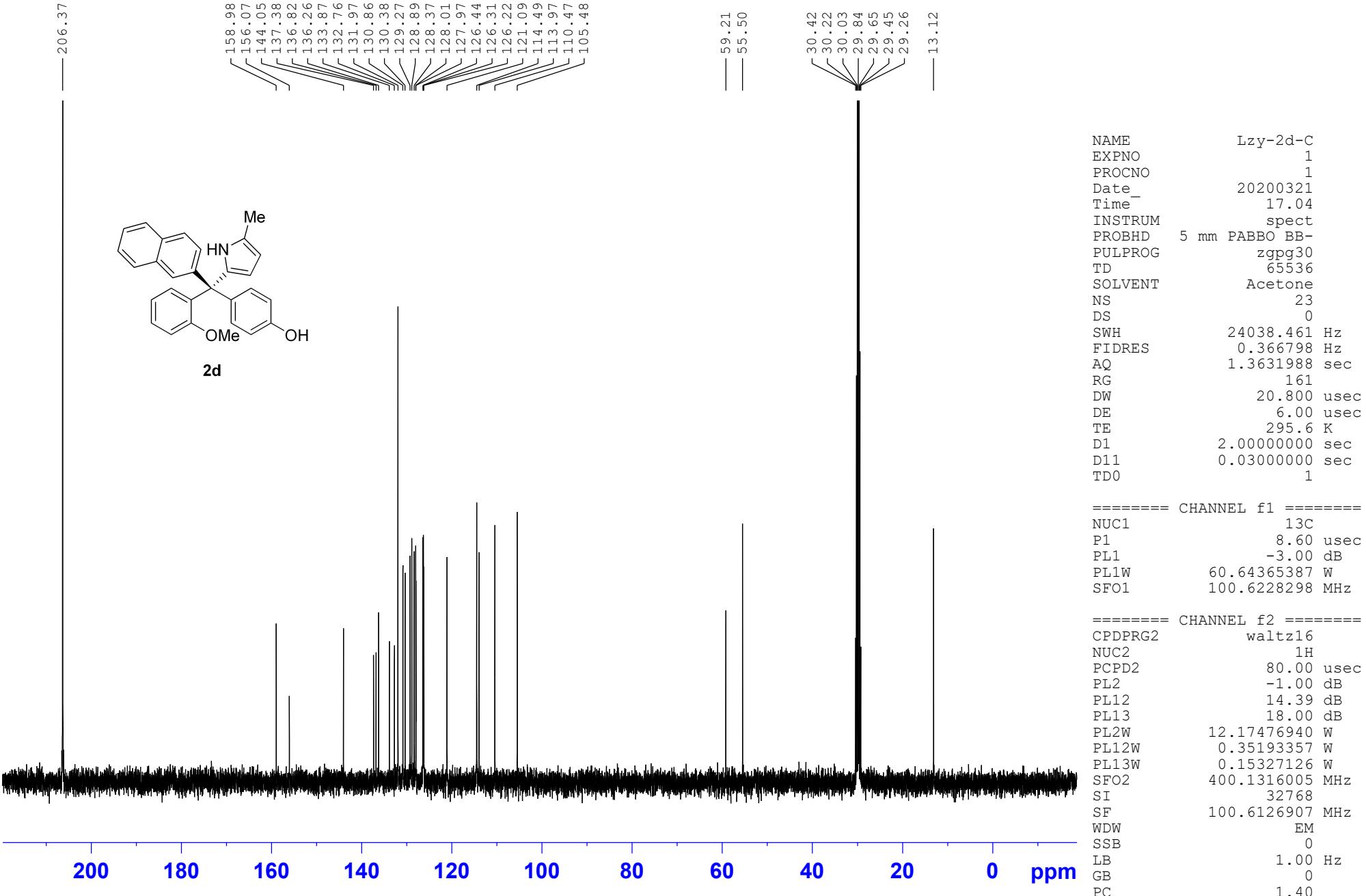
NAME	Lzy-2b-F
EXPNO	1
PROCNO	1
Date	20200207
Time	11.59
INSTRUM	spect
PROBHD	5 mm PABBO BB-
PULPROG	zgflqn
TD	131072
SOLVENT	Acetone
NS	7
DS	4
SWH	89285.711 Hz
FIDRES	0.681196 Hz
AQ	0.7340532 sec
RG	1290
DW	5.600 usec
DE	6.00 usec
TE	294.9 K
D1	1.00000000 sec
TD0	1
===== CHANNEL f1 =====	
NUC1	19F
P1	19.50 usec
PL1	-4.00 dB
PL1W	16.97275162 W
SFO1	376.4607164 MHz
SI	65536
SF	376.4983660 MHz
WDW	EM
SSB	5
LB	0.30 Hz
GB	0.05
PC	1.00

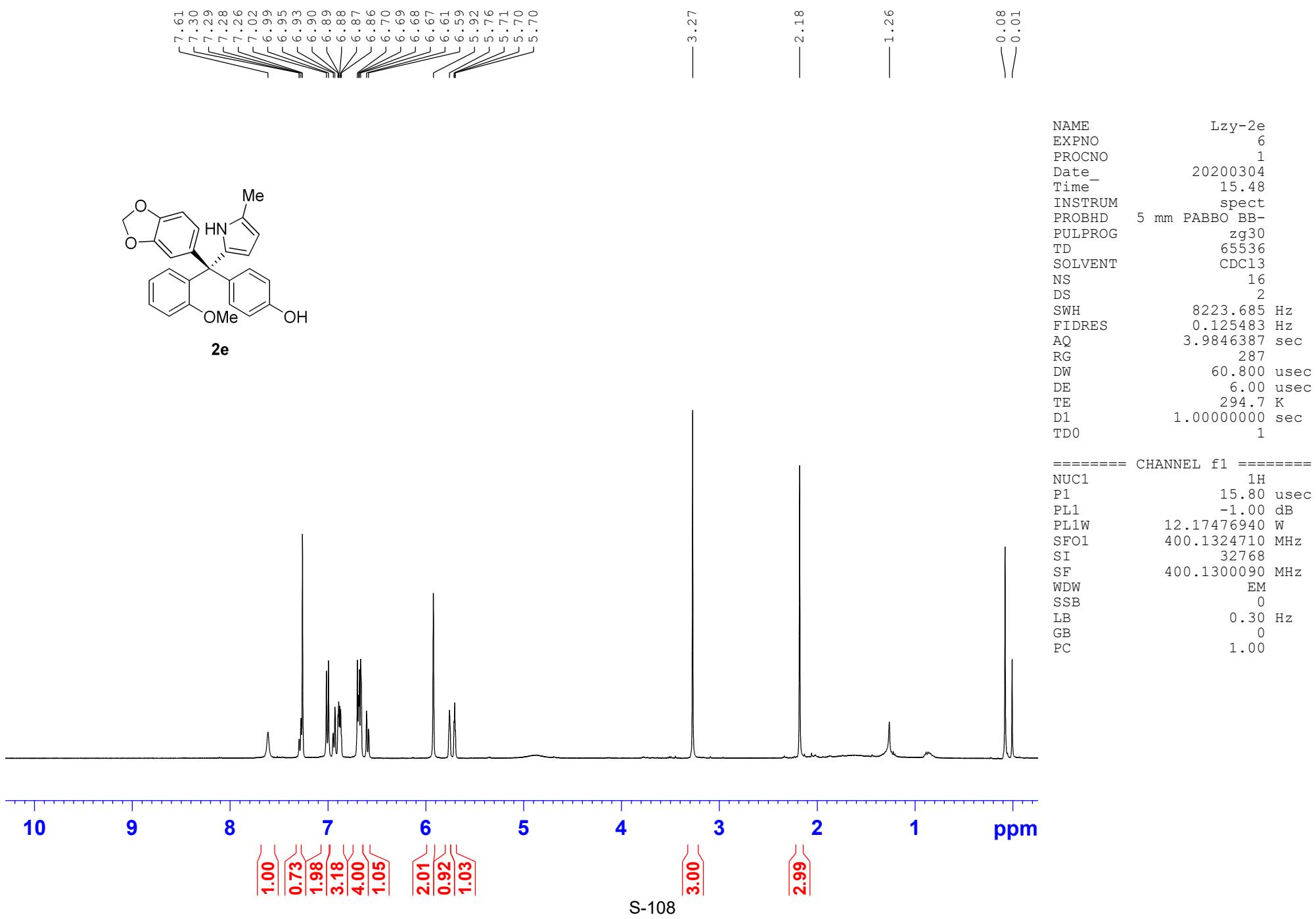
0 -20 -40 -60 -80 -100 -120 -140 -160 -180 -200 ppm

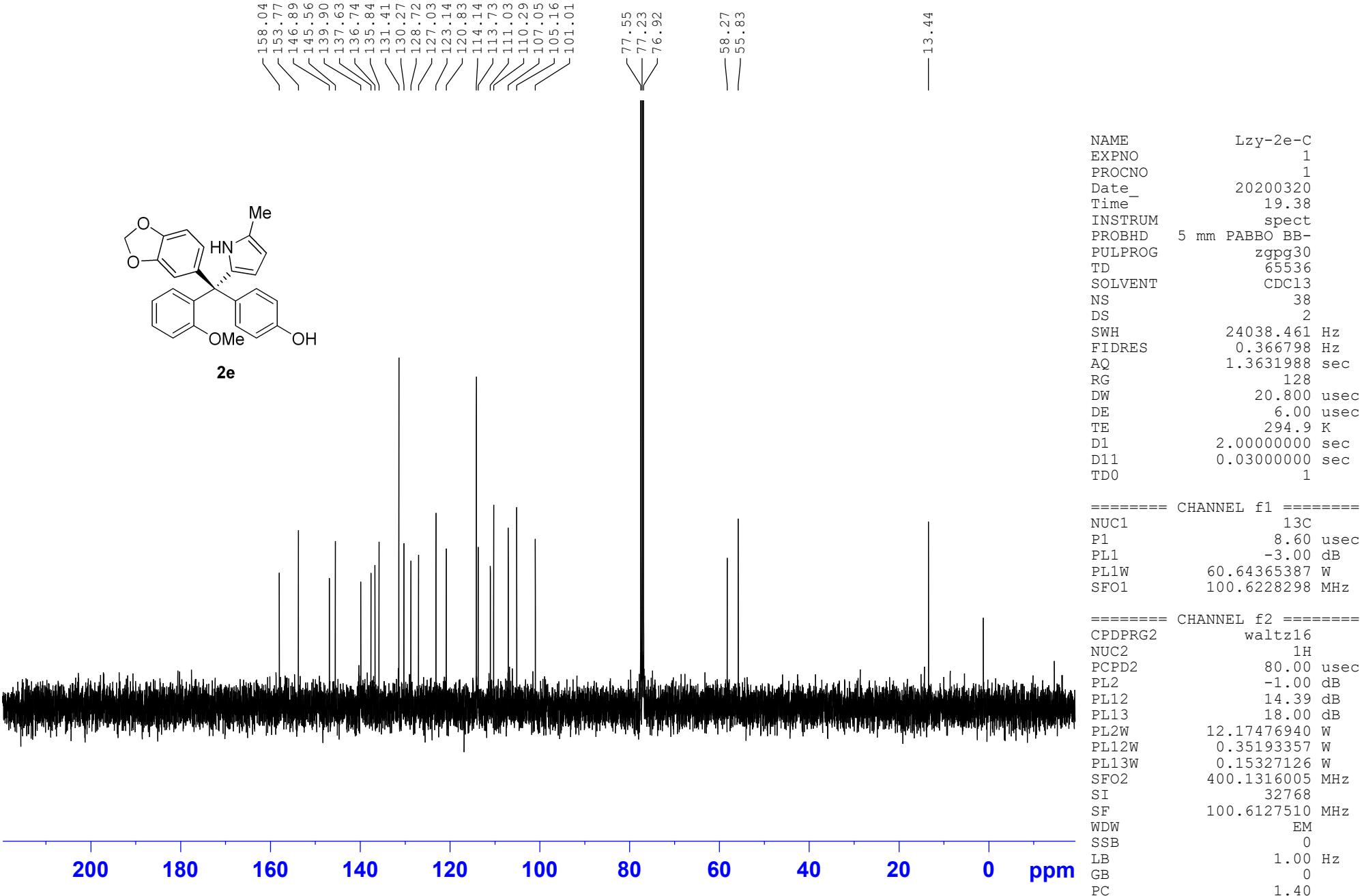


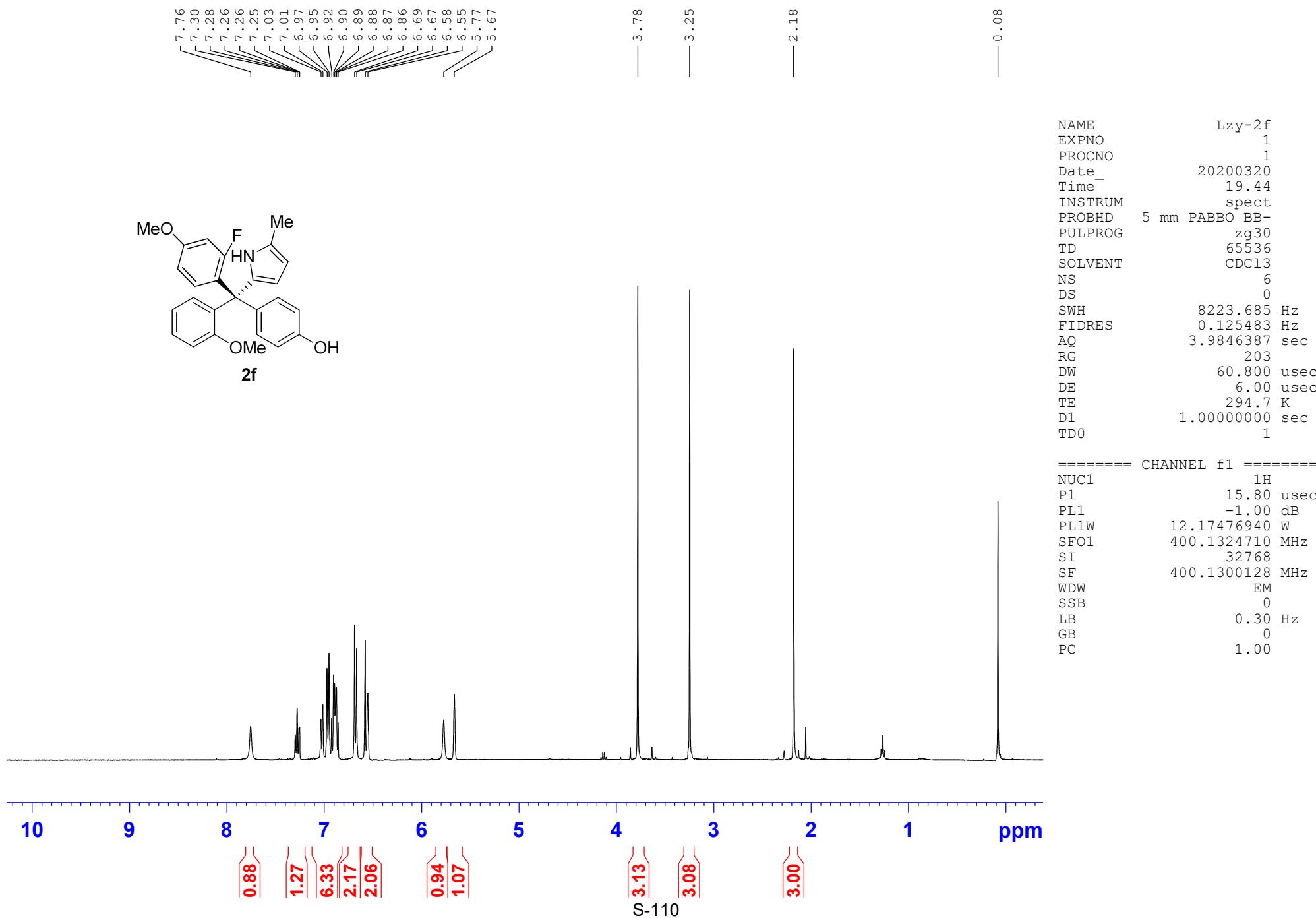


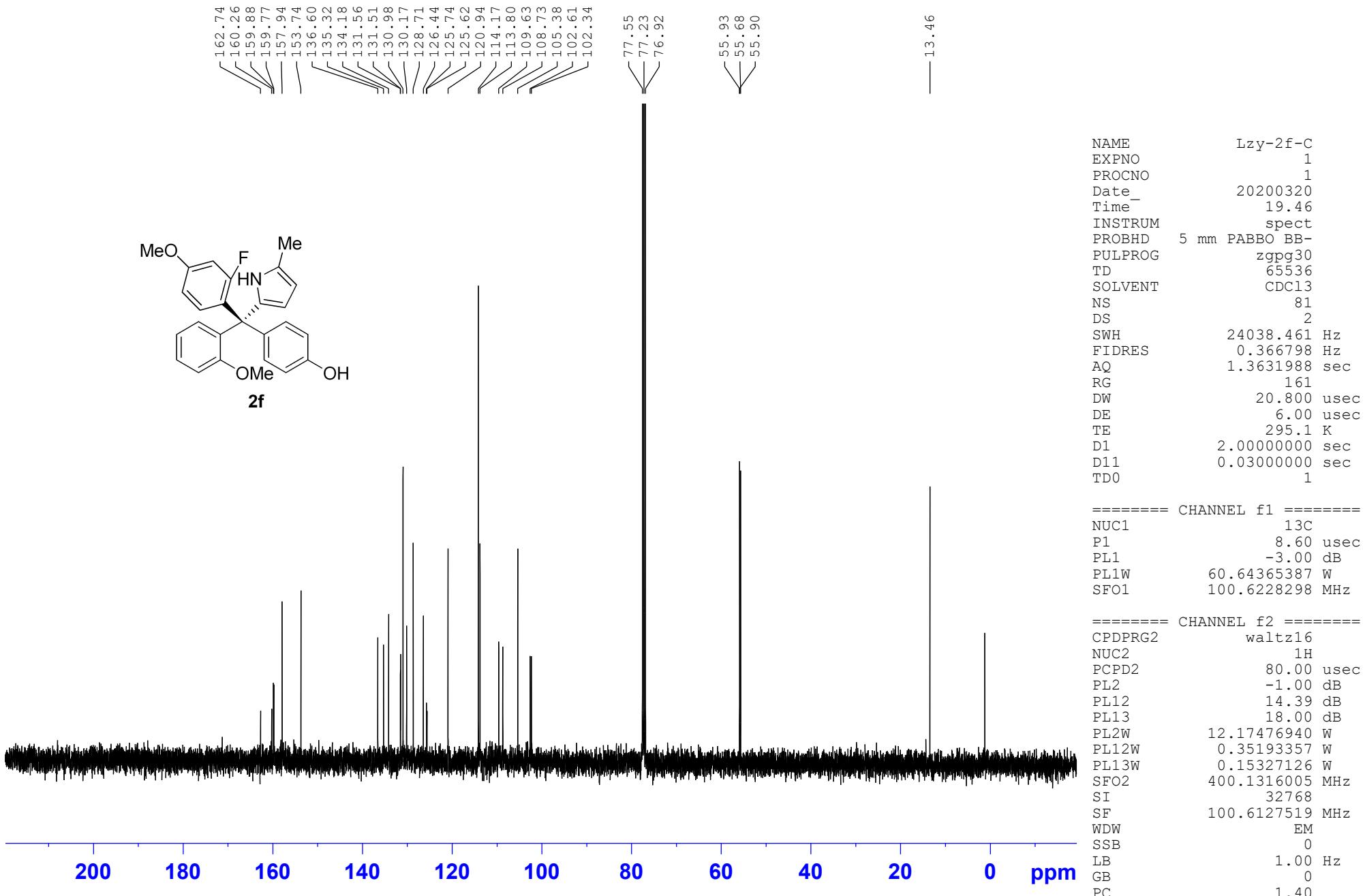


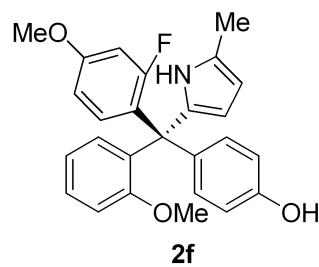




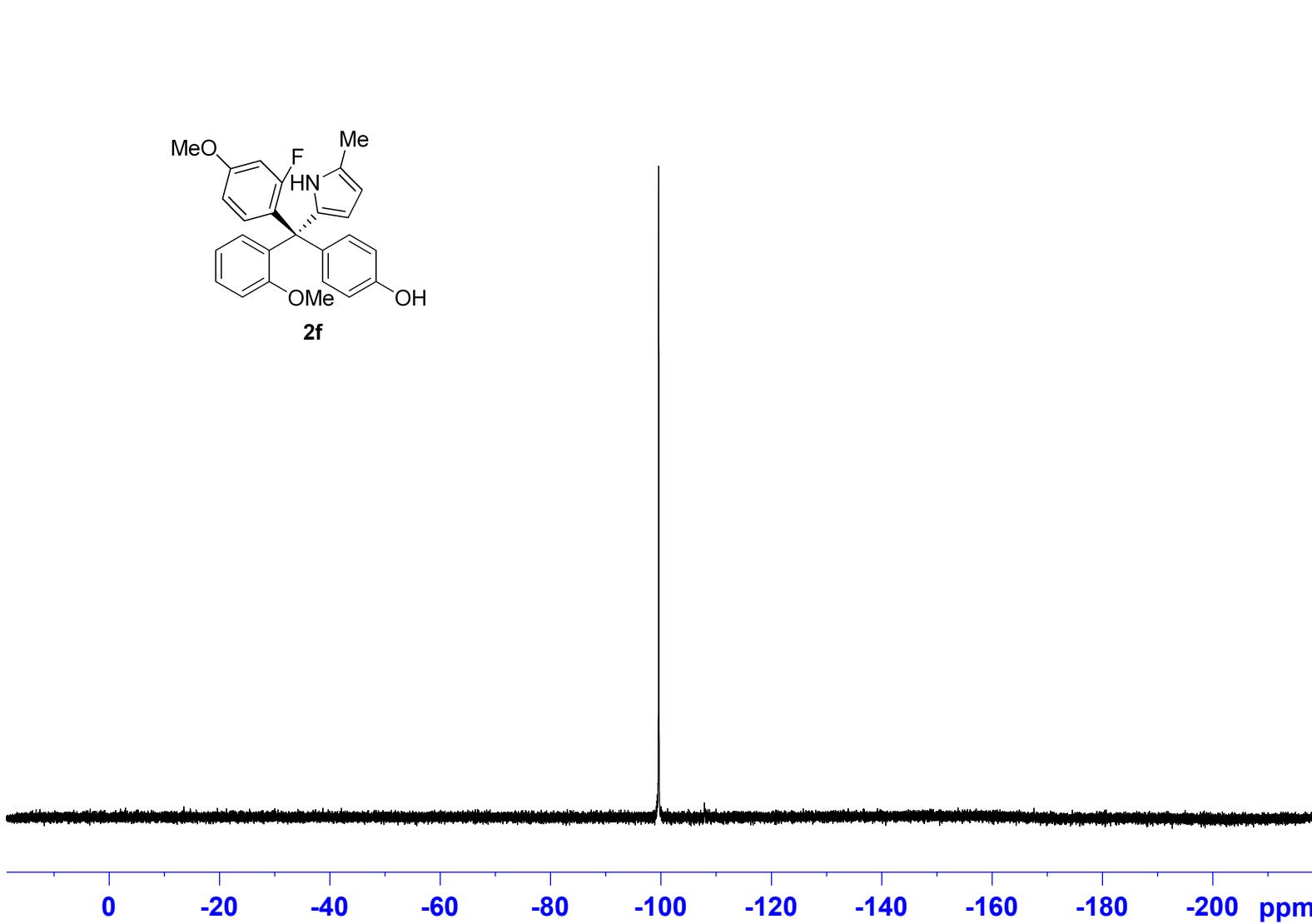






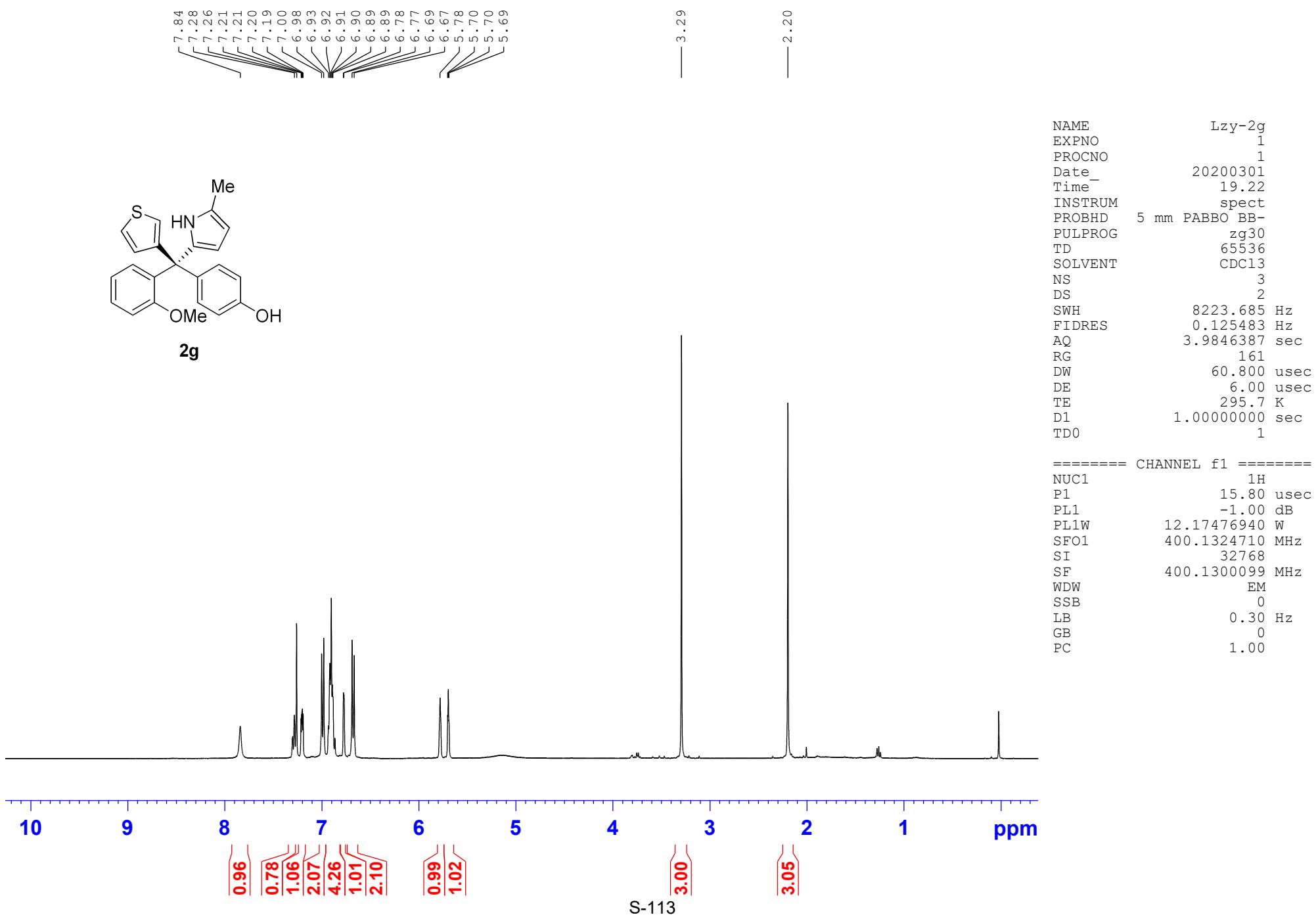


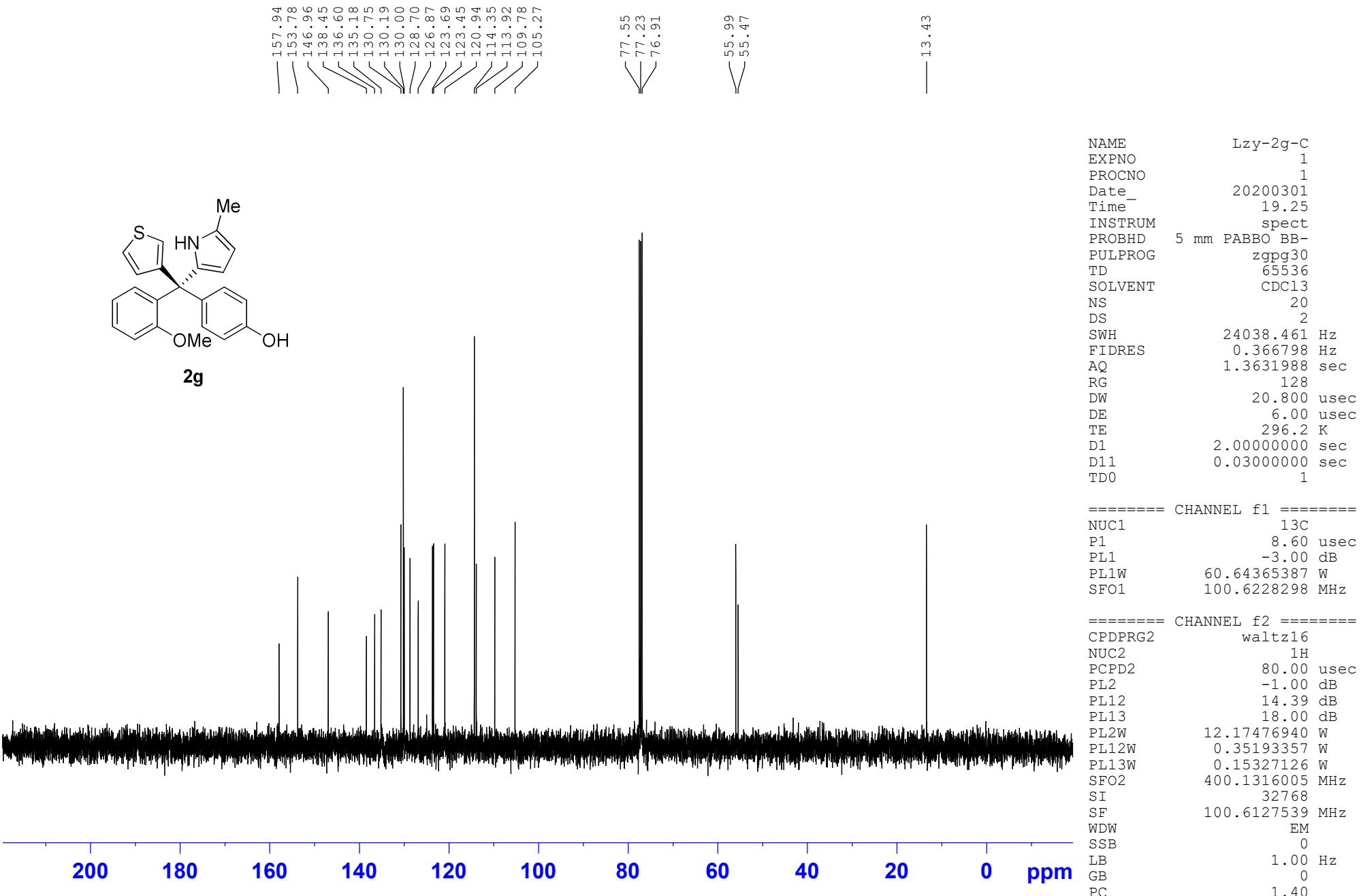
-99.59

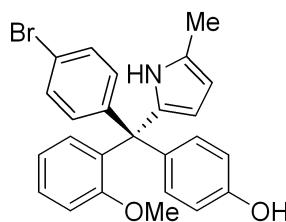
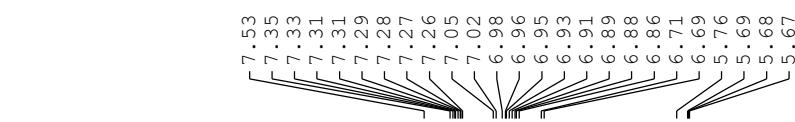


NAME Lzy-2f-F
 EXPNO 1
 PROCNO 1
 Date 20200320
 Time 19.53
 INSTRUM spect
 PROBHD 5 mm PABBO BB-
 PULPROG zgflqn
 TD 131072
 SOLVENT CDC13
 NS 2
 DS 0
 SWH 89285.711 Hz
 FIDRES 0.681196 Hz
 AQ 0.7340532 sec
 RG 1620
 DW 5.600 usec
 DE 6.00 usec
 TE 294.8 K
 D1 1.00000000 sec
 TD0 1

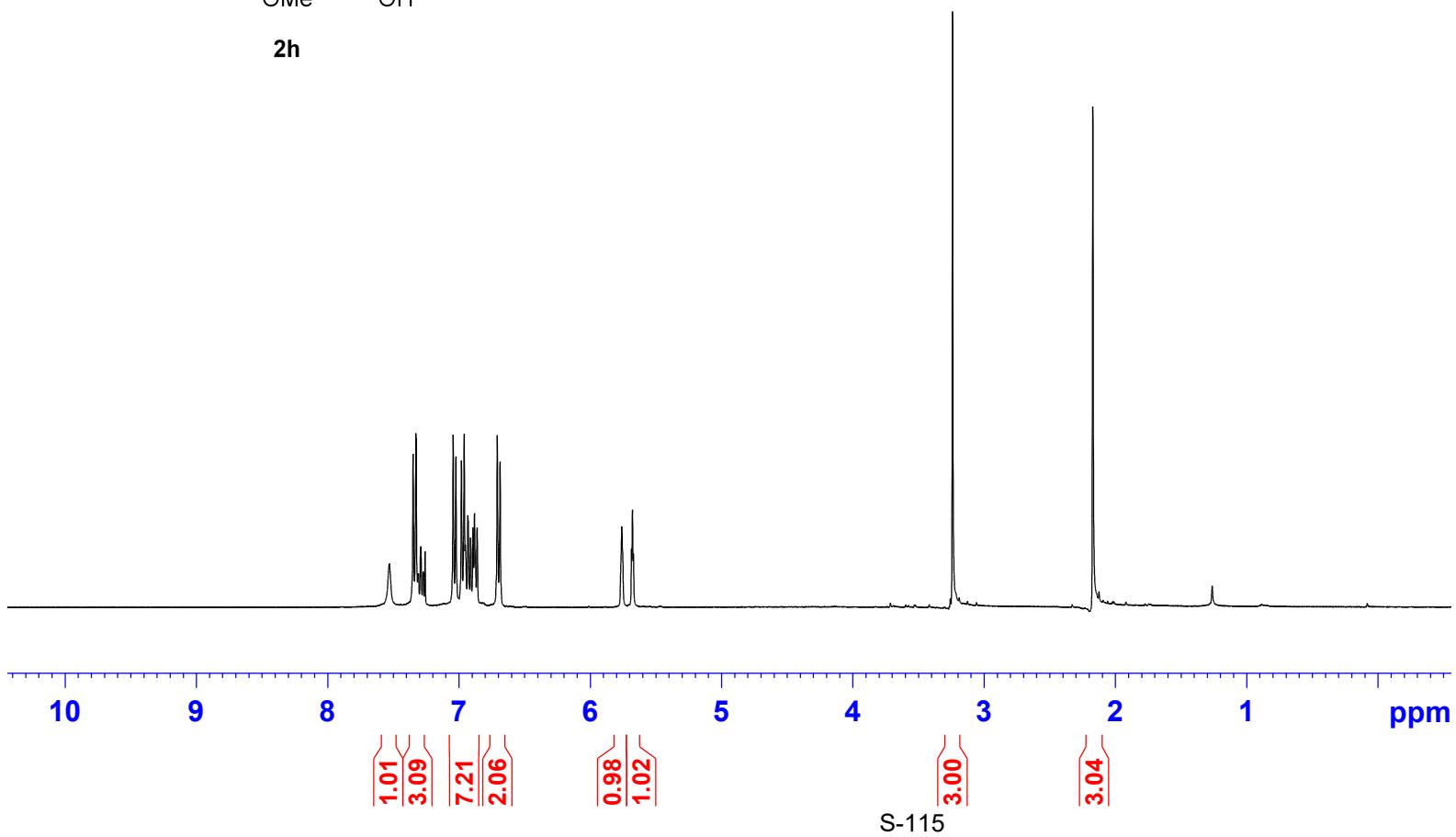
===== CHANNEL f1 ======
 NUC1 19F
 P1 19.50 usec
 PL1 -4.00 dB
 PL1W 16.97275162 W
 SFO1 376.4607164 MHz
 SI 65536
 SF 376.4983660 MHz
 WDW EM
 SSB 0
 LB 0.30 Hz
 GB 0
 PC 1.00

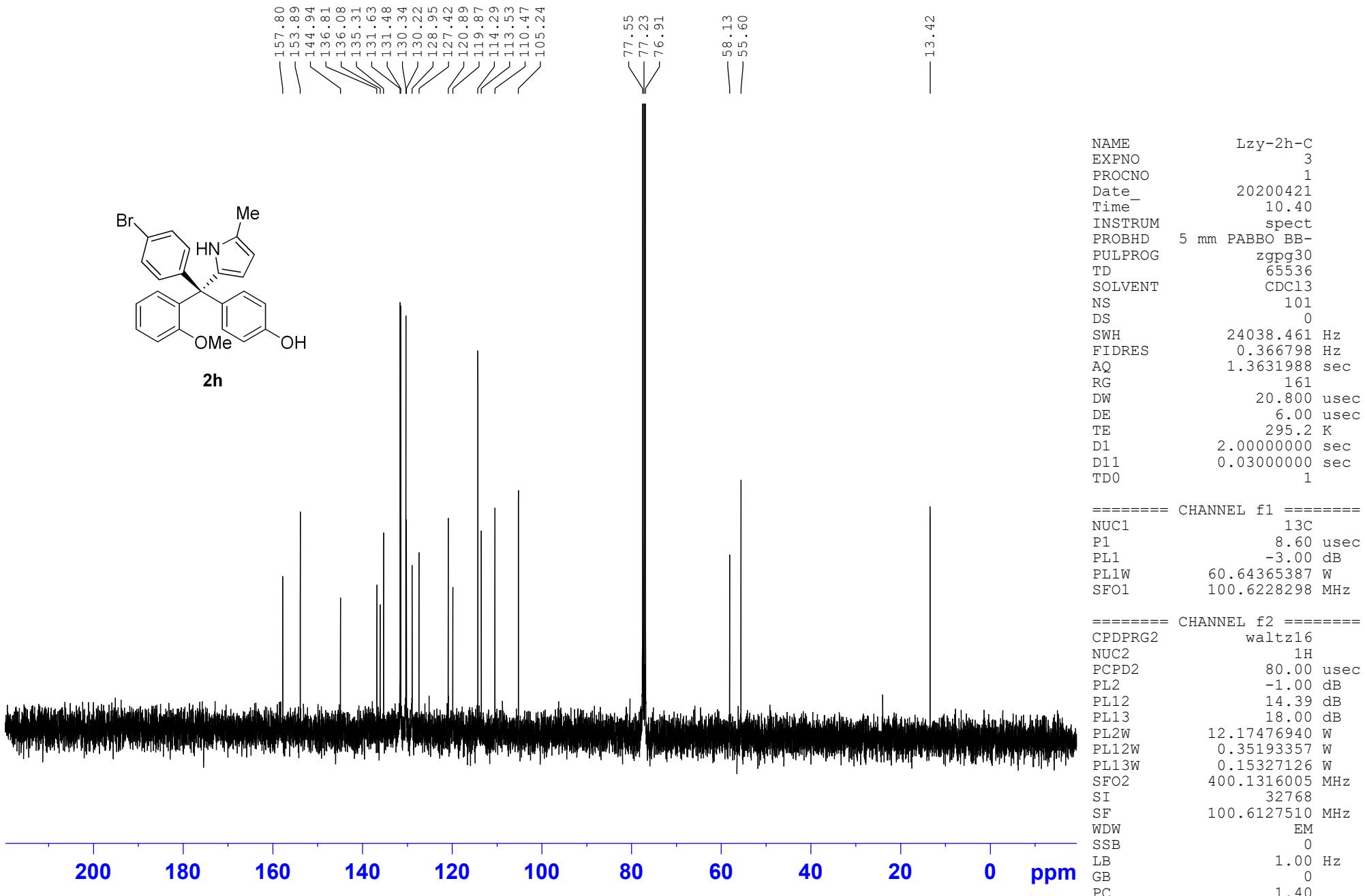


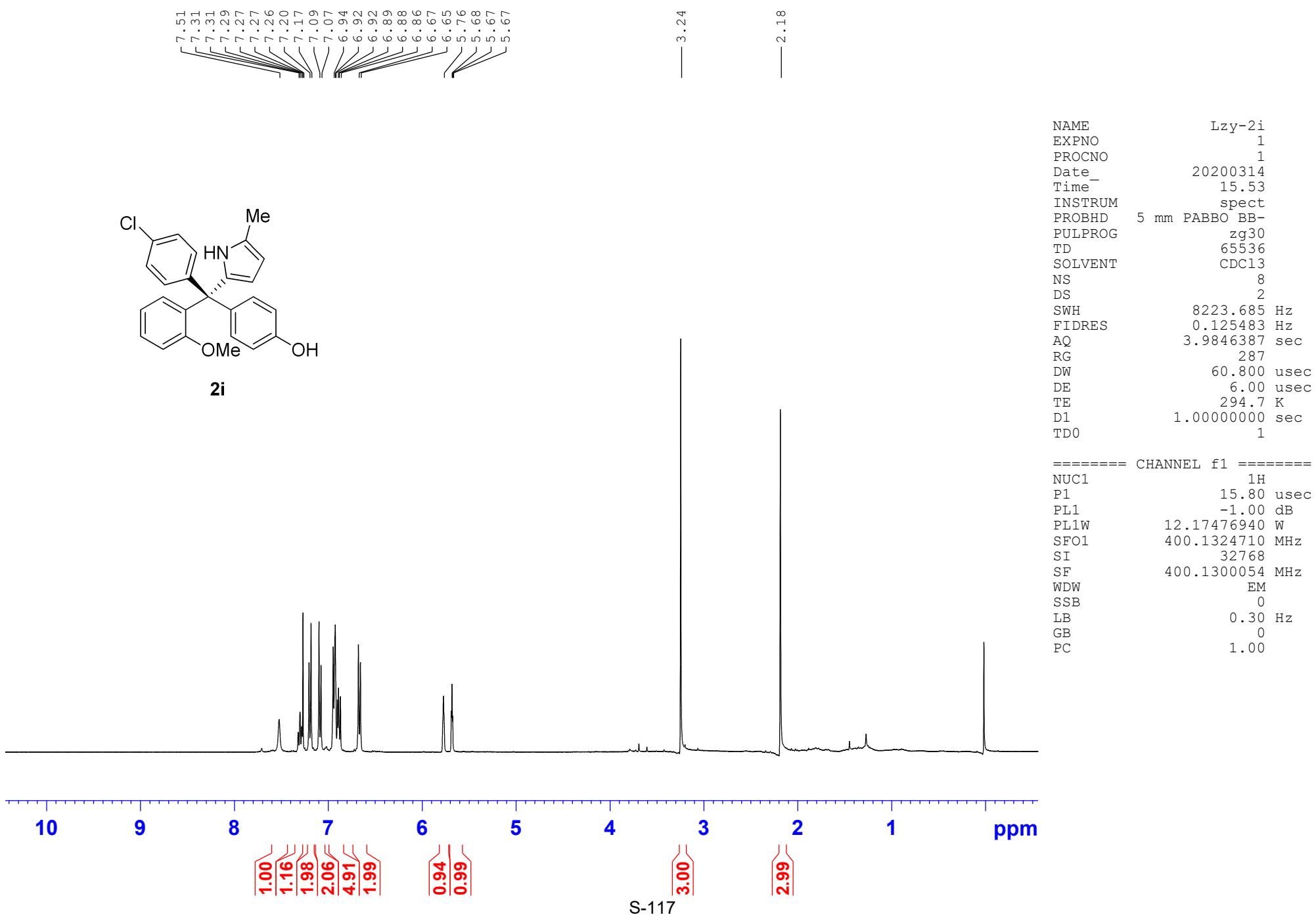


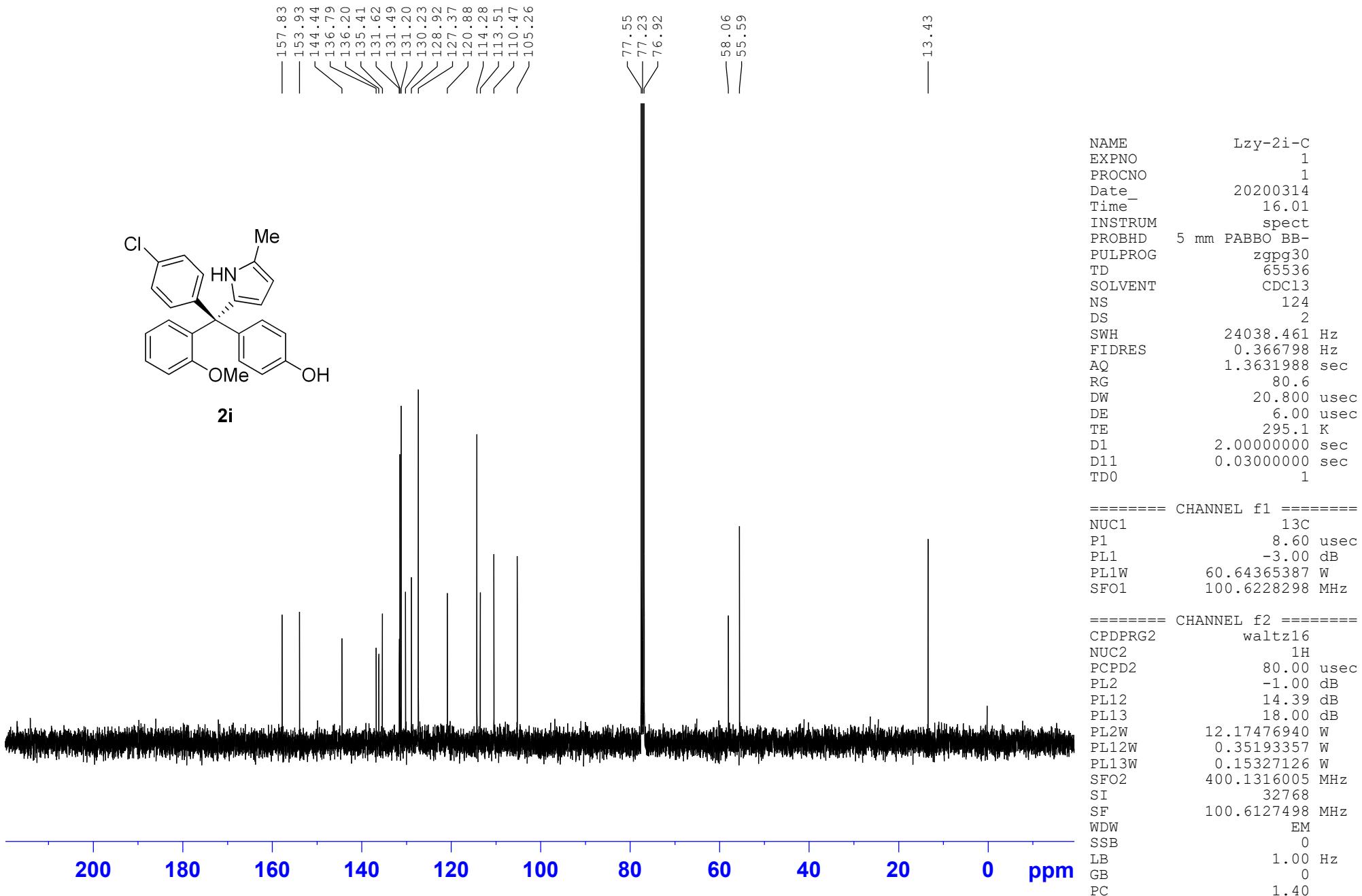


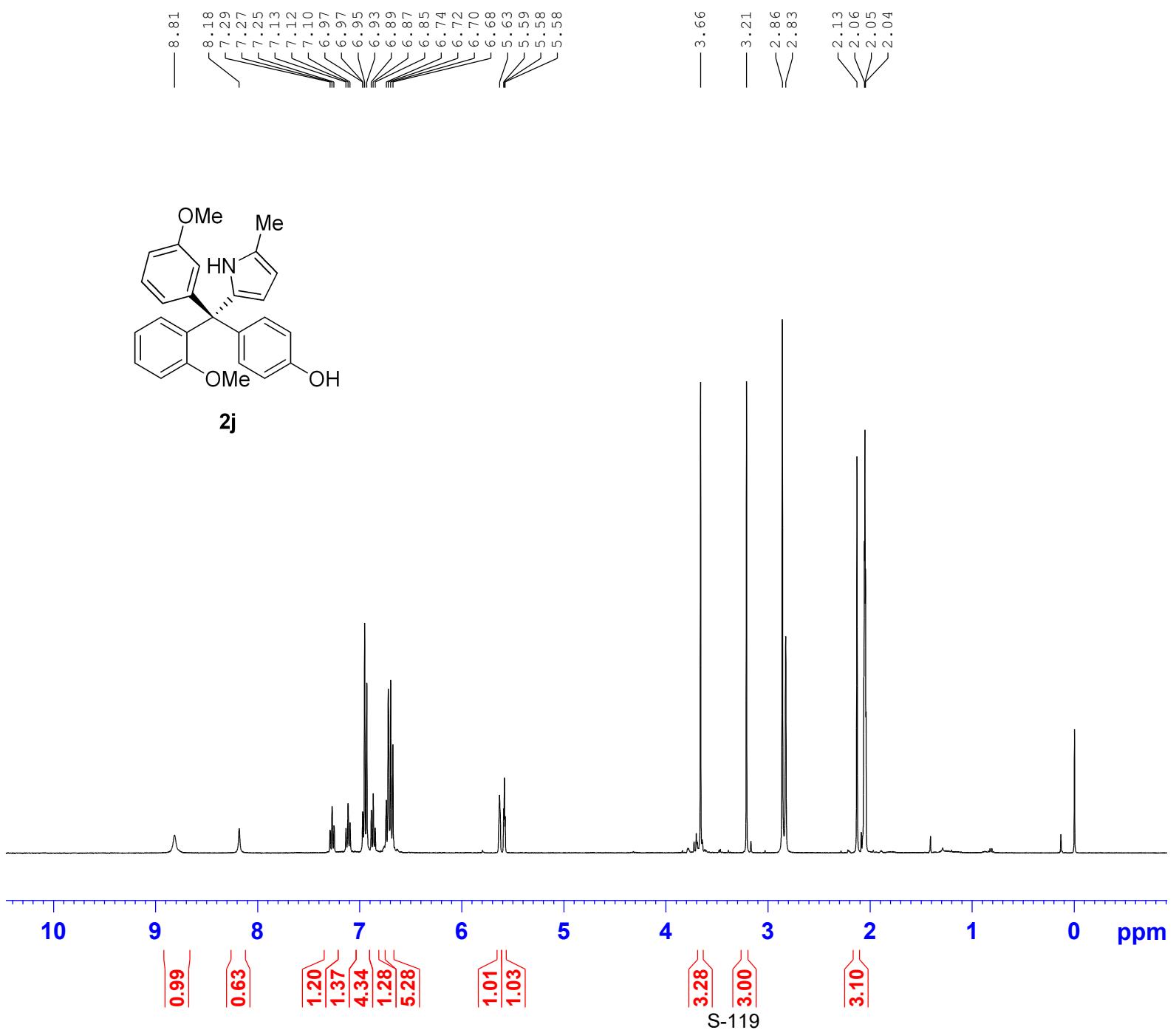
2h









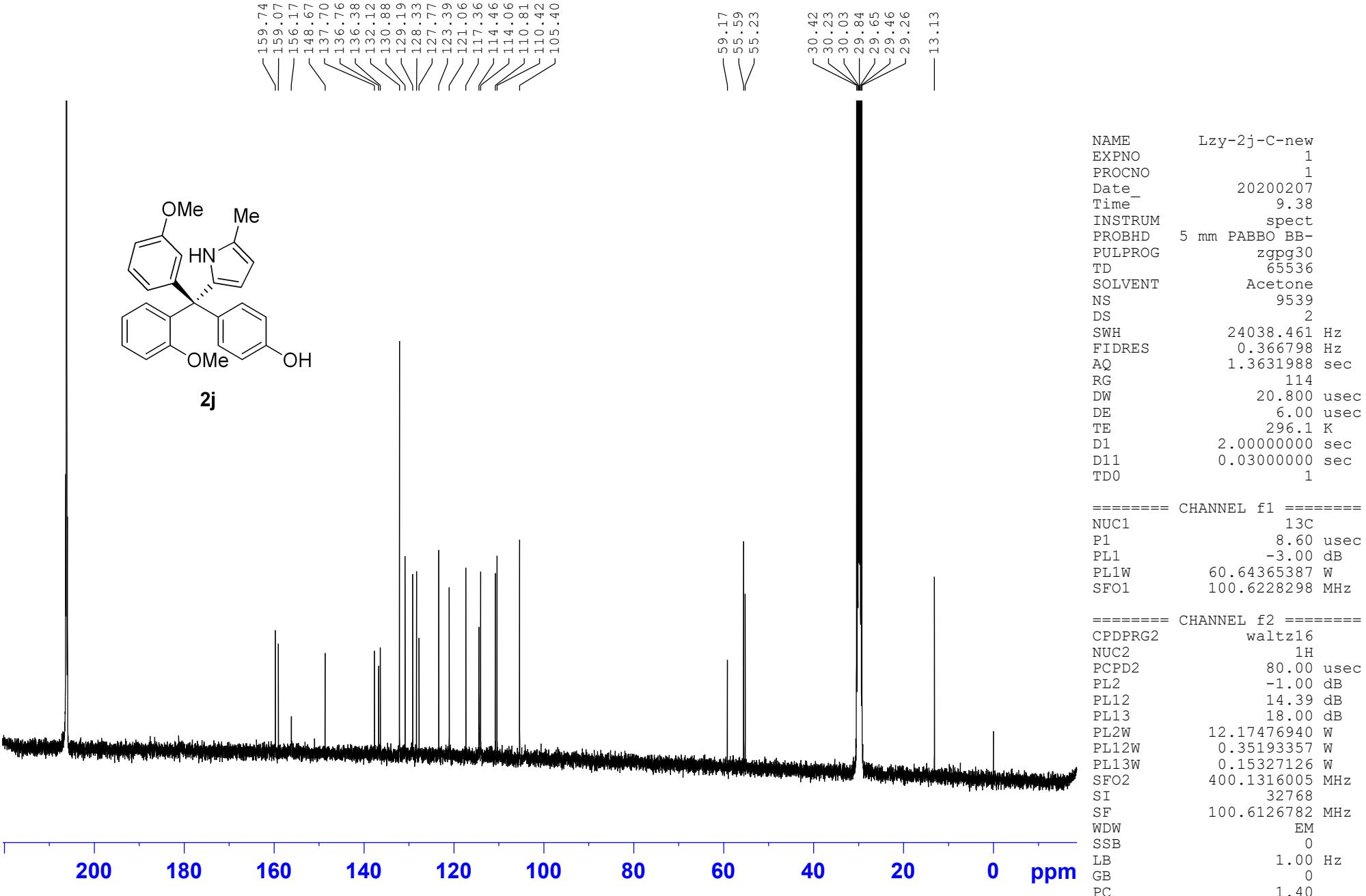


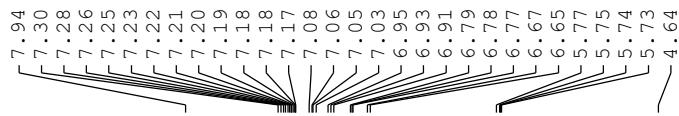
```

NAME          Lzy-2j
EXPNO         1
PROCNO        1
Date_ 20200203
Time_ 19.48
INSTRUM      spect
PROBHD      5 mm PABBO BB-
PULPROG zg30
TD       65536
SOLVENT      Acetone
NS          16
DS           2
SWH        8223.685 Hz
FIDRES      0.125483 Hz
AQ        3.9846387 sec
RG          456
DW        60.800 usec
DE          6.00 usec
TE        294.7 K
D1    1.00000000 sec
TD0             1

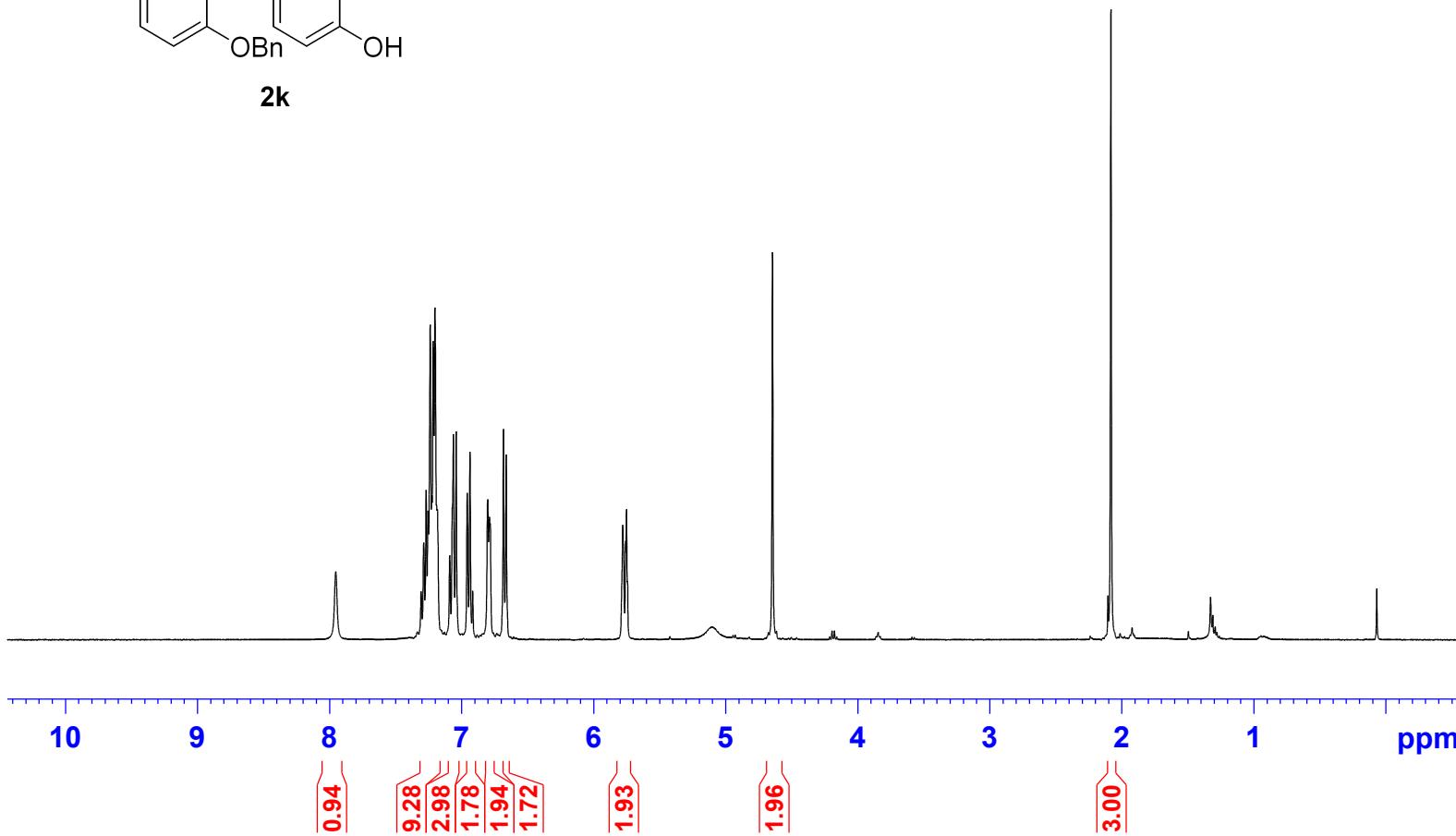
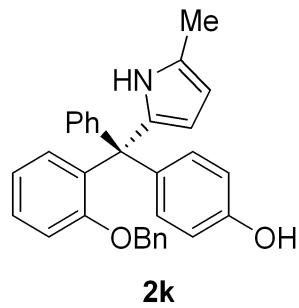
===== CHANNEL f1 =====
NUC1            1H
P1        15.80 usec
PL1           -1.00 dB
PL1W        12.17476940 W
SFO1        400.1324710 MHz
SI            32768
SF        400.1300062 MHz
WDW
SSB            0
LB          0.30 Hz
GB            0
PC          1.00

```





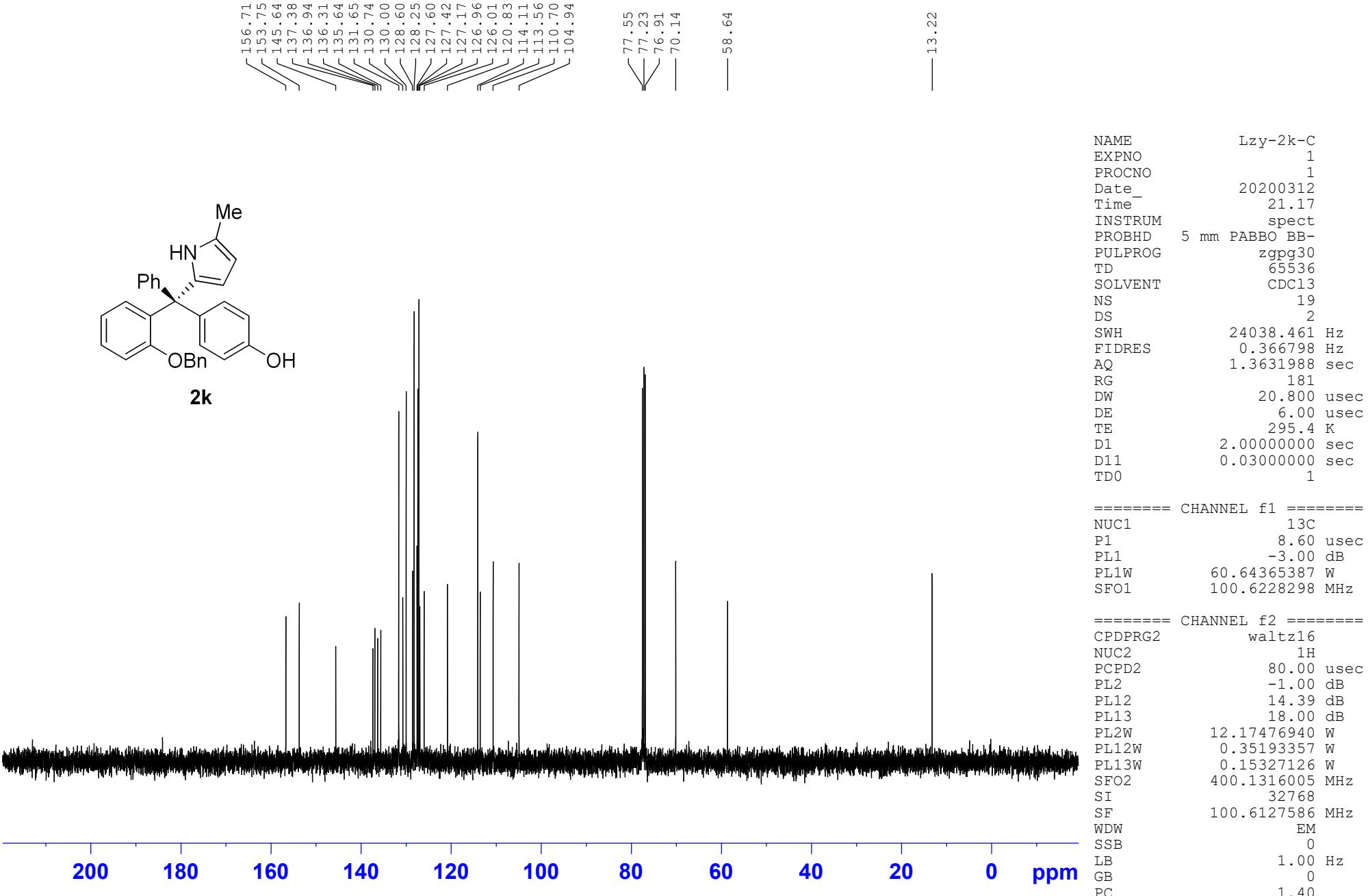
— 2.07 —

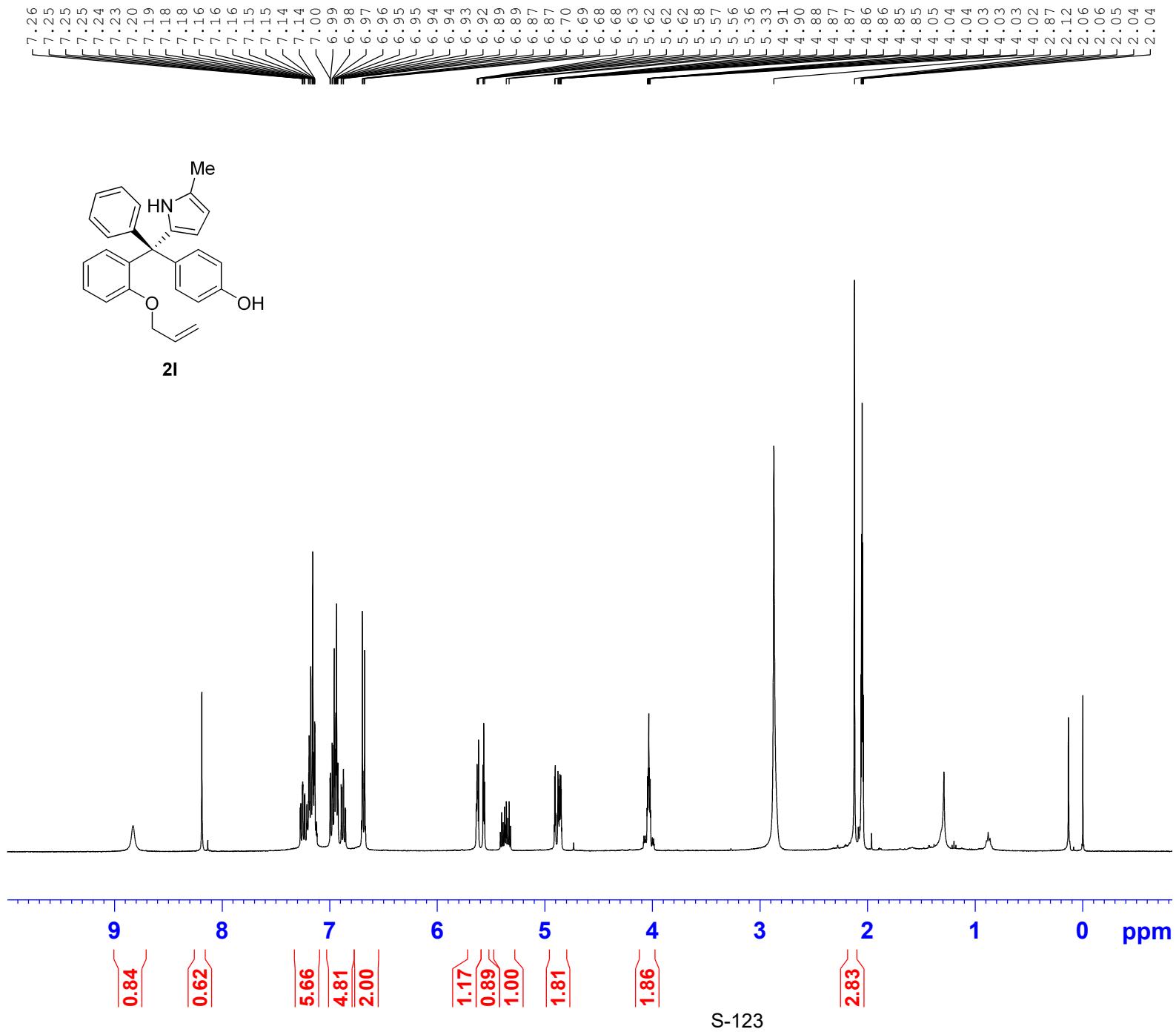


NAME	Lzy-2k
EXPNO	2
PROCNO	1
Date	20200312
Time	21.20
INSTRUM	spect
PROBHD	5 mm PABBO BB-
PULPROG	zg30
TD	65536
SOLVENT	CDCl ₃
NS	2
DS	2
SWH	8223.685 Hz
FIDRES	0.125483 Hz
AQ	3.9846387 sec
RG	101
DW	60.800 usec
DE	6.00 usec
TE	295.2 K
D1	1.00000000 sec
TD0	1

===== CHANNEL f1 =====

NUC1	1H
P1	15.80 usec
PL1	-1.00 dB
PL1W	12.17476940 W
SFO1	400.1324710 MHz
SI	32768
SF	400.1300054 MHz
WDW	EM
SSB	0
LB	0.30 Hz
GB	0
PC	1.00



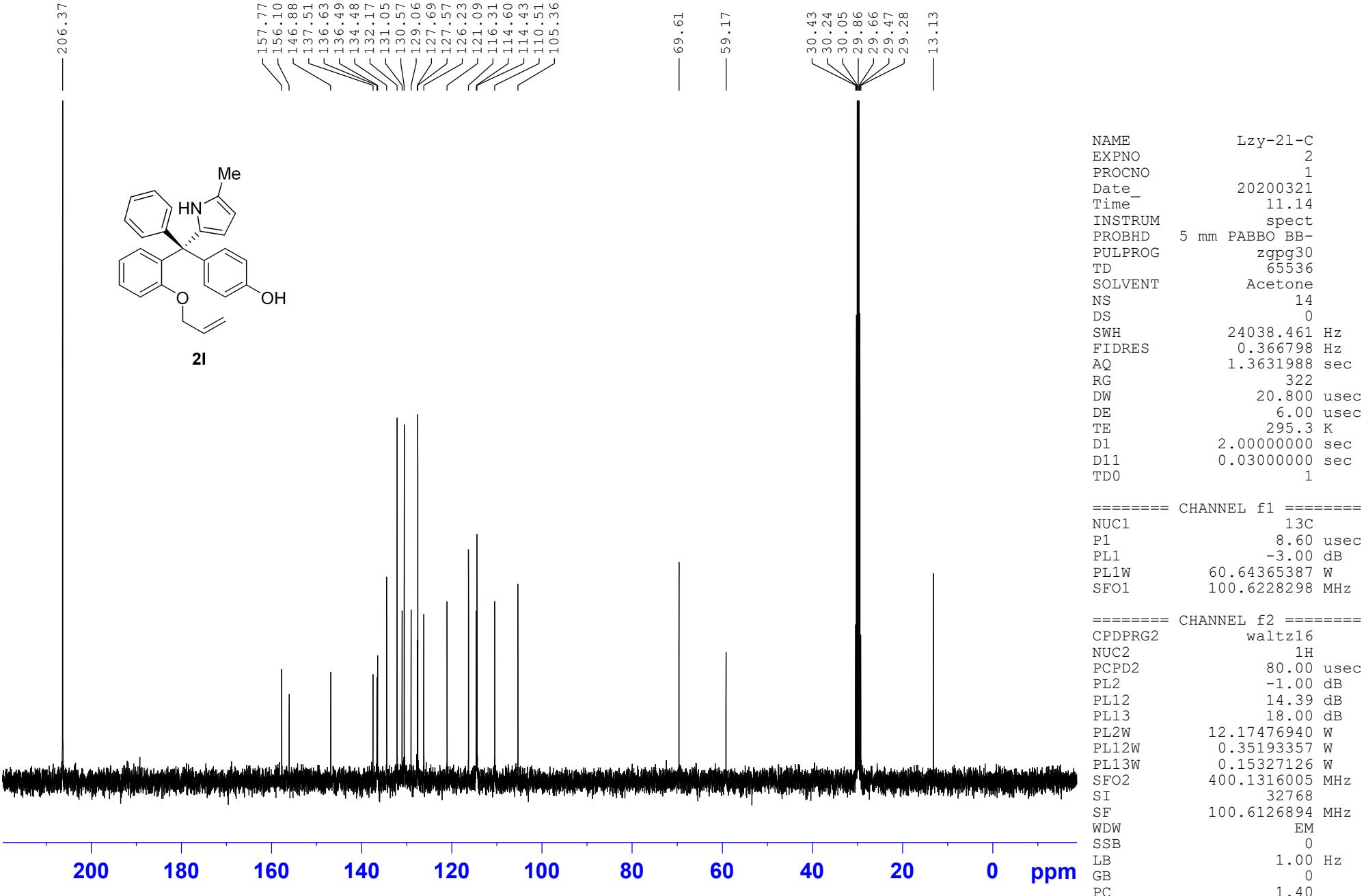


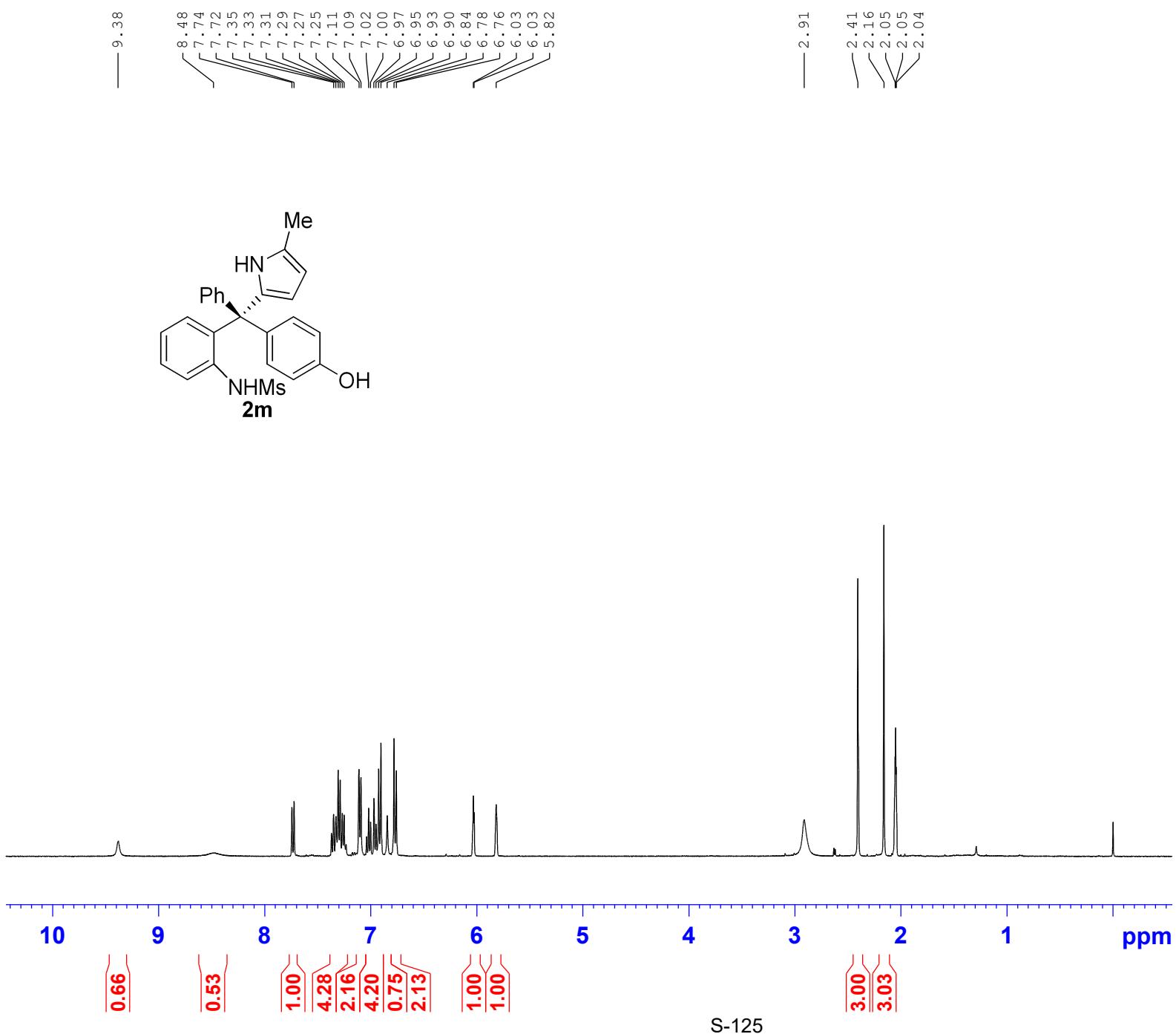
```

NAME      lzy-project1-11-H-1
EXPNO
PROCNO
Date_     20210509
Time_     19.32
INSTRUM  spect
PROBHD   5 mm PABBO BB/
PULPROG  zg30
TD        65536
SOLVENT   Acetone
NS
DS
SWH      8012.820 Hz
FIDRES   0.122266 Hz
AQ        4.0894966 sec
RG        126.97
DW        62.400 usec
DE        6.50  usec
TE        298.2 K
D1        1.00000000 sec
TDO0

```

```
===== CHANNEL f1 ======  
SFO1          400.1324710 MHz  
NUC1           1H  
P1             14.50 usec  
SI              65536  
SF          400.1300069 MHz  
WDW            EM  
SSB              0  
LB             0.30 Hz  
GB              0  
PC             1.00
```

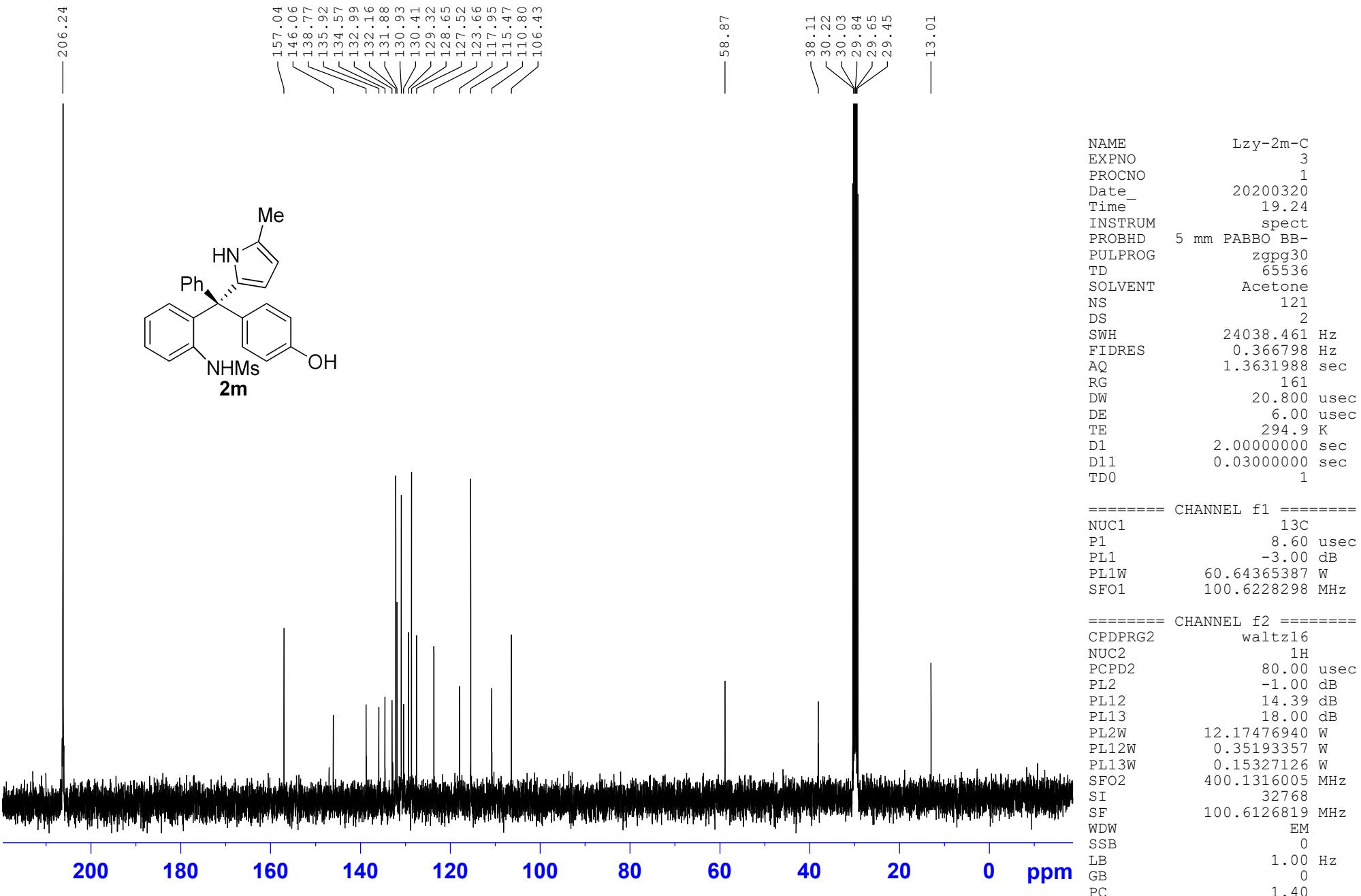


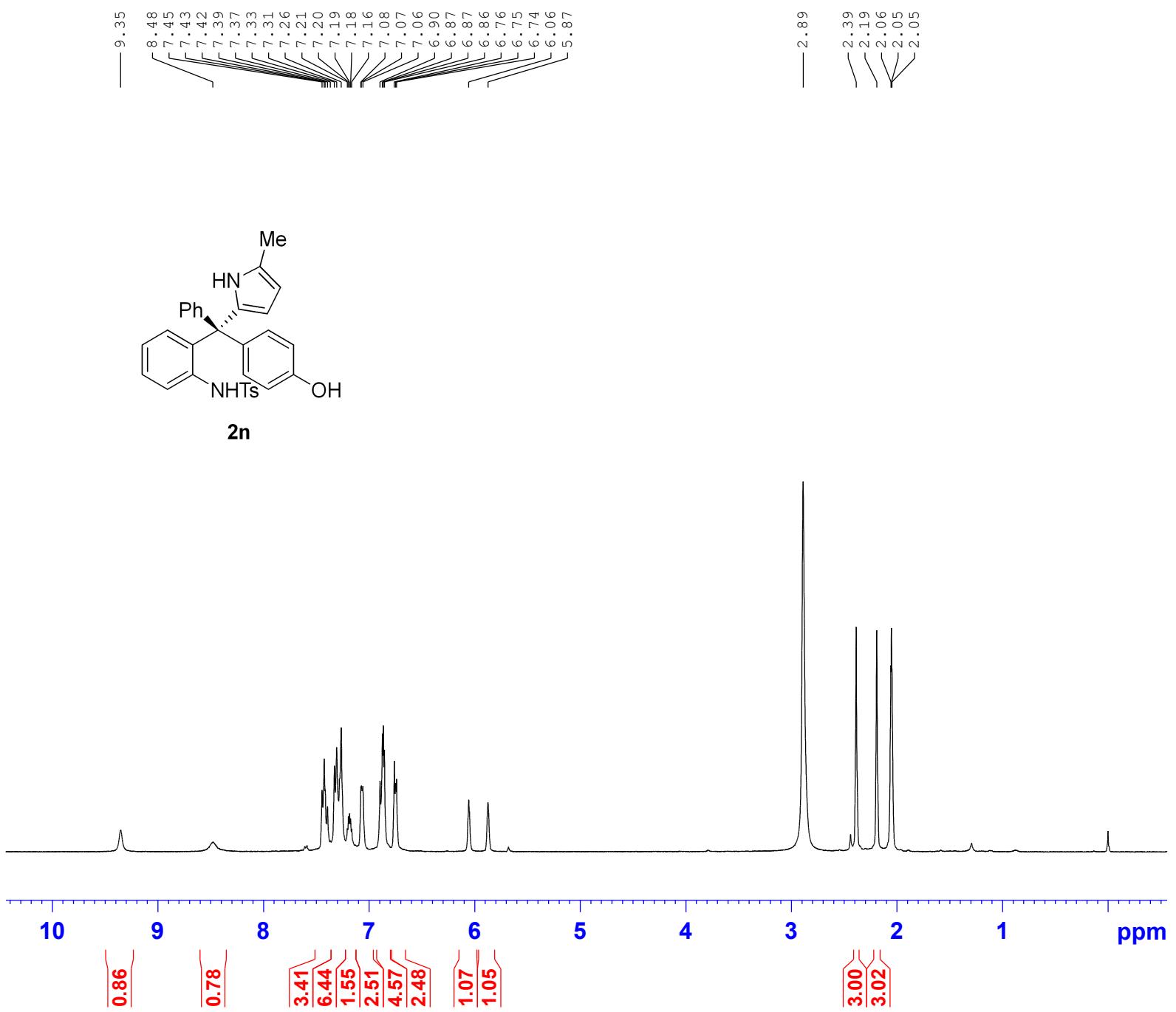


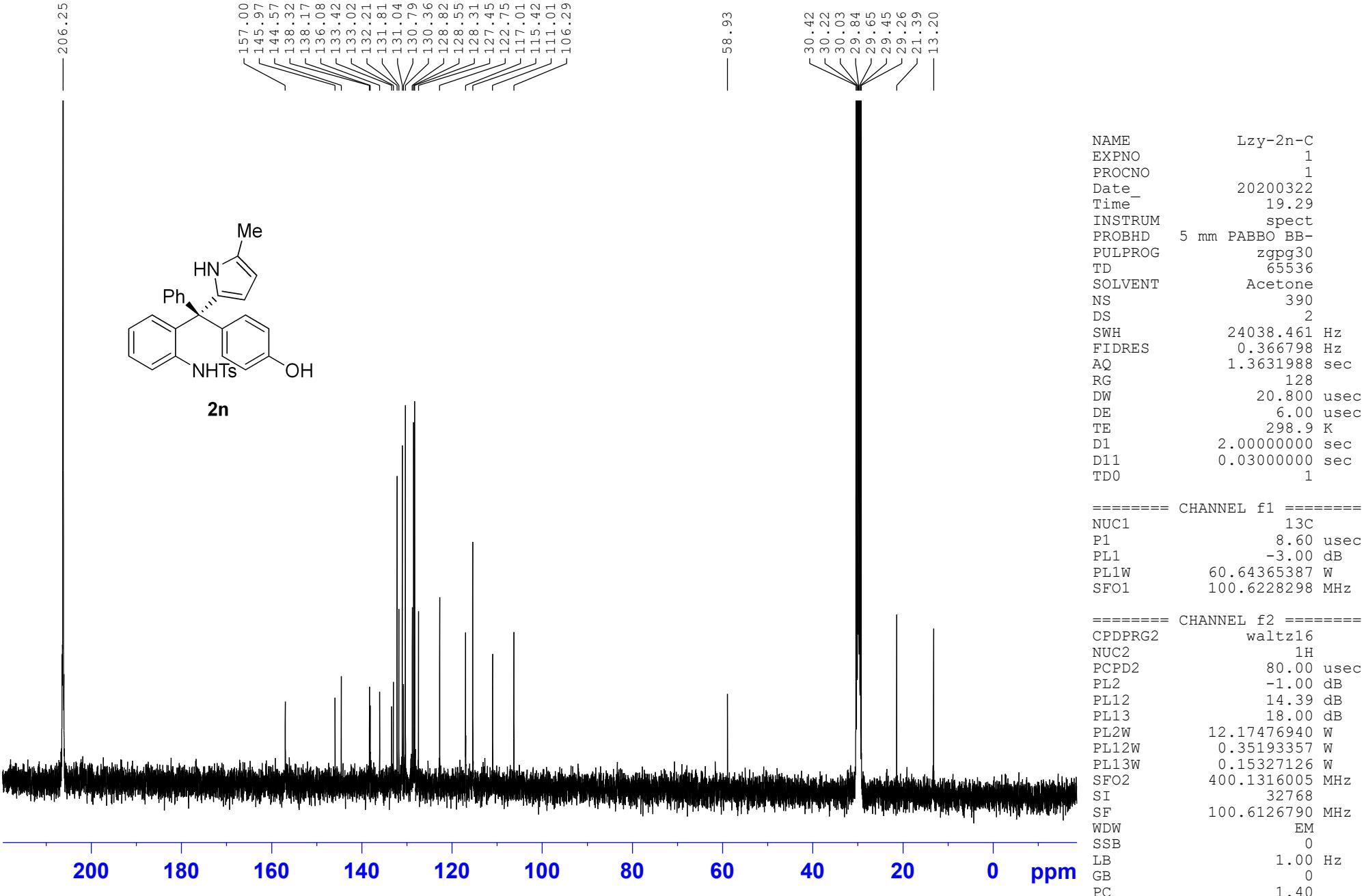
NAME	Lzy-2m
EXPNO	3
PROCNO	1
Date	20200320
Time	19.22
INSTRUM	spect
PROBHD	5 mm PABBO BB-
PULPROG	zg30
TD	65536
SOLVENT	Acetone
NS	6
DS	0
SWH	8223.685 Hz
FIDRES	0.125483 Hz
AQ	3.9846387 sec
RG	575
DW	60.800 usec
DE	6.00 usec
TE	294.7 K
D1	1.00000000 sec
TD0	1

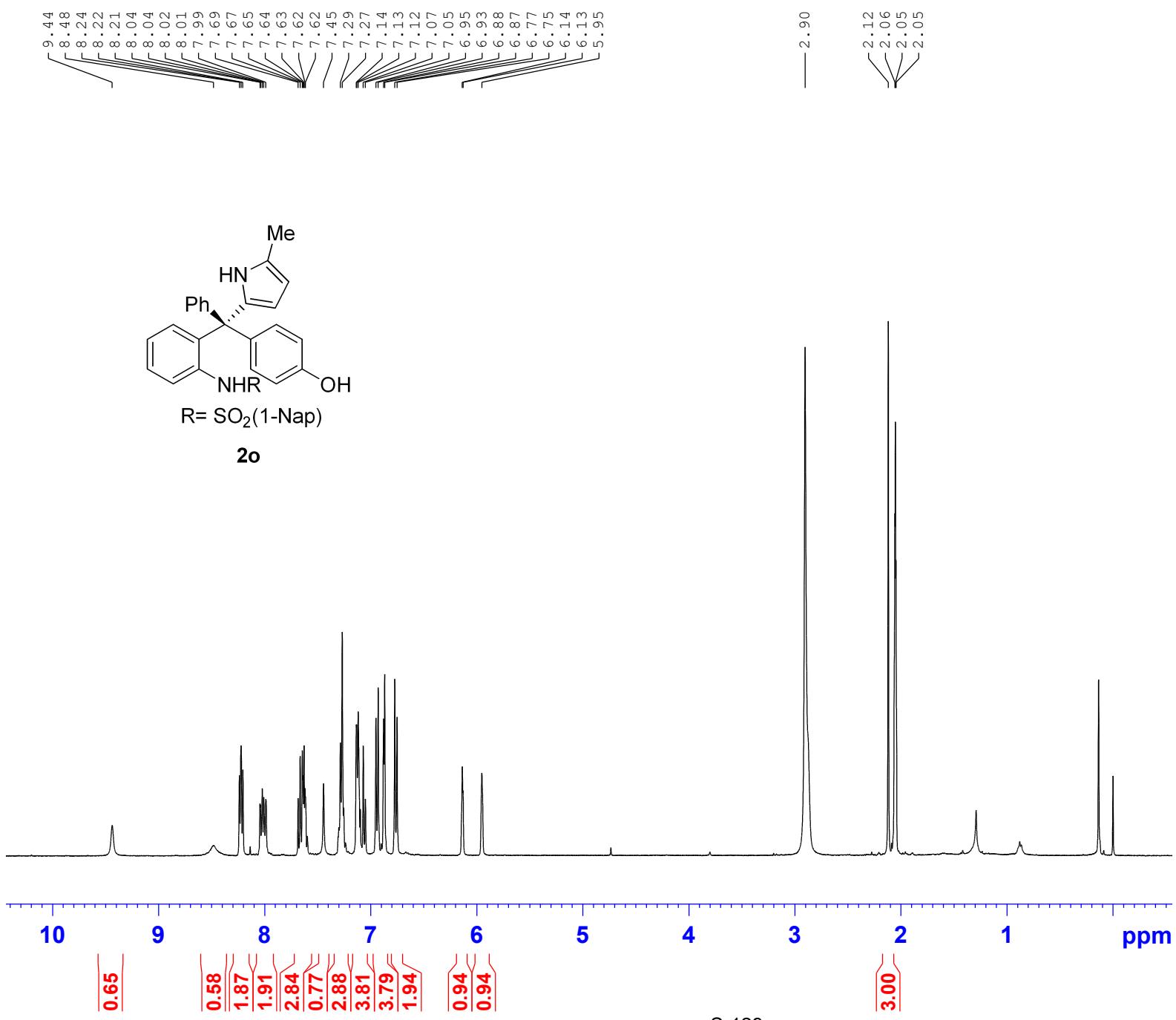
===== CHANNEL f1 =====

NUC1	1H
P1	15.80 usec
PL1	-1.00 dB
PL1W	12.17476940 W
SFO1	400.1324710 MHz
SI	32768
SF	400.1300063 MHz
WDW	EM
SSB	0
LB	0.30 Hz
GB	0
PC	1.00





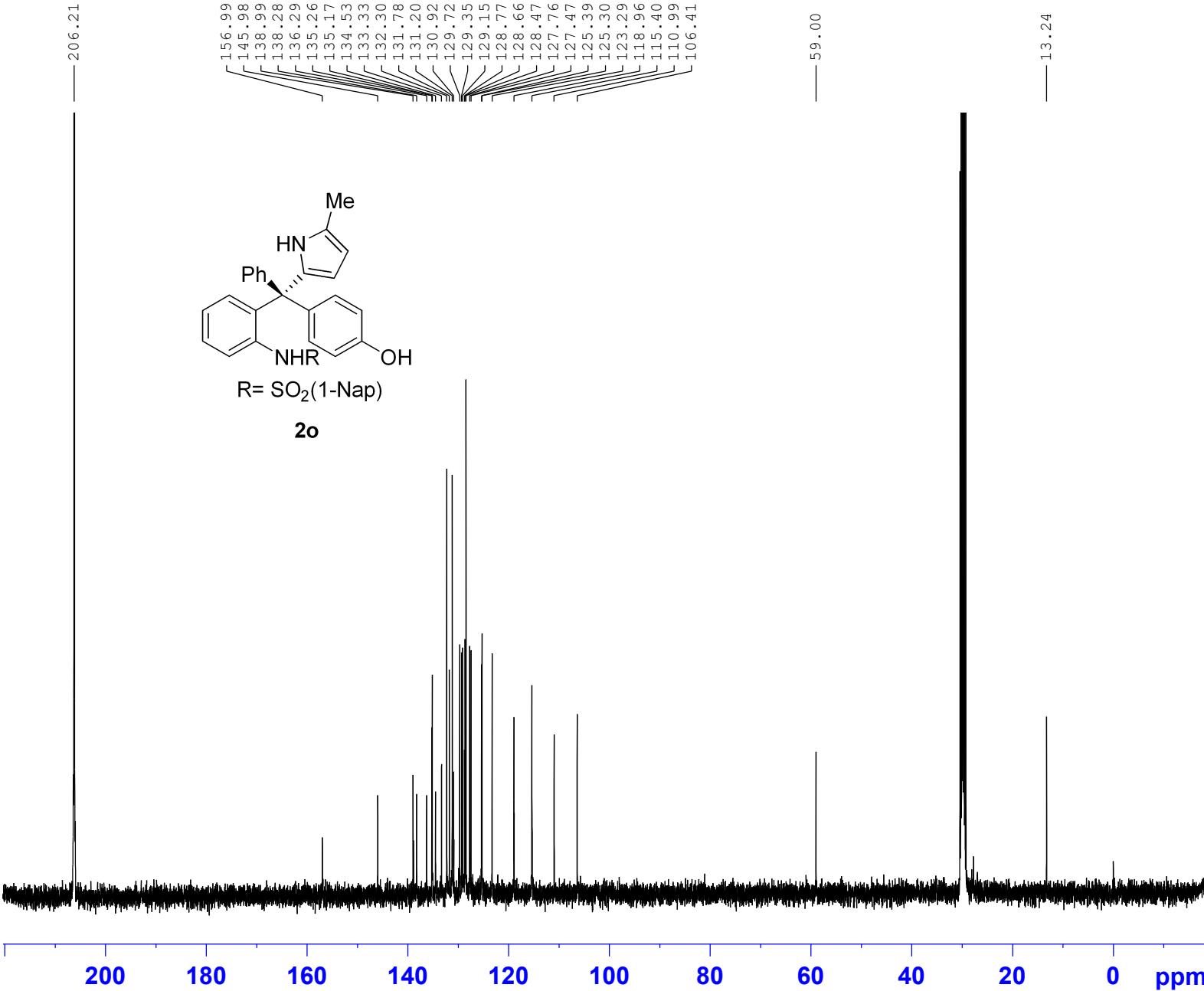




NAME Lzy-2o
 EXPNO 2
 PROCNO 1
 Date 20200319
 Time 19.45
 INSTRUM spect
 PROBHD 5 mm PABBO BB-
 PULPROG zg30
 TD 65536
 SOLVENT Acetone
 NS 6
 DS 0
 SWH 8223.685 Hz
 FIDRES 0.125483 Hz
 AQ 3.9846387 sec
 RG 322
 DW 60.800 usec
 DE 6.00 usec
 TE 294.8 K
 D1 1.00000000 sec
 TDO 1

===== CHANNEL f1 =====

NUC1 1H
 P1 15.80 usec
 PL1 -1.00 dB
 PL1W 12.17476940 W
 SFO1 400.1324710 MHz
 SI 32768
 SF 400.1300054 MHz
 WDW EM
 SSB 0
 LB 0.30 Hz
 GB 0
 PC 1.00



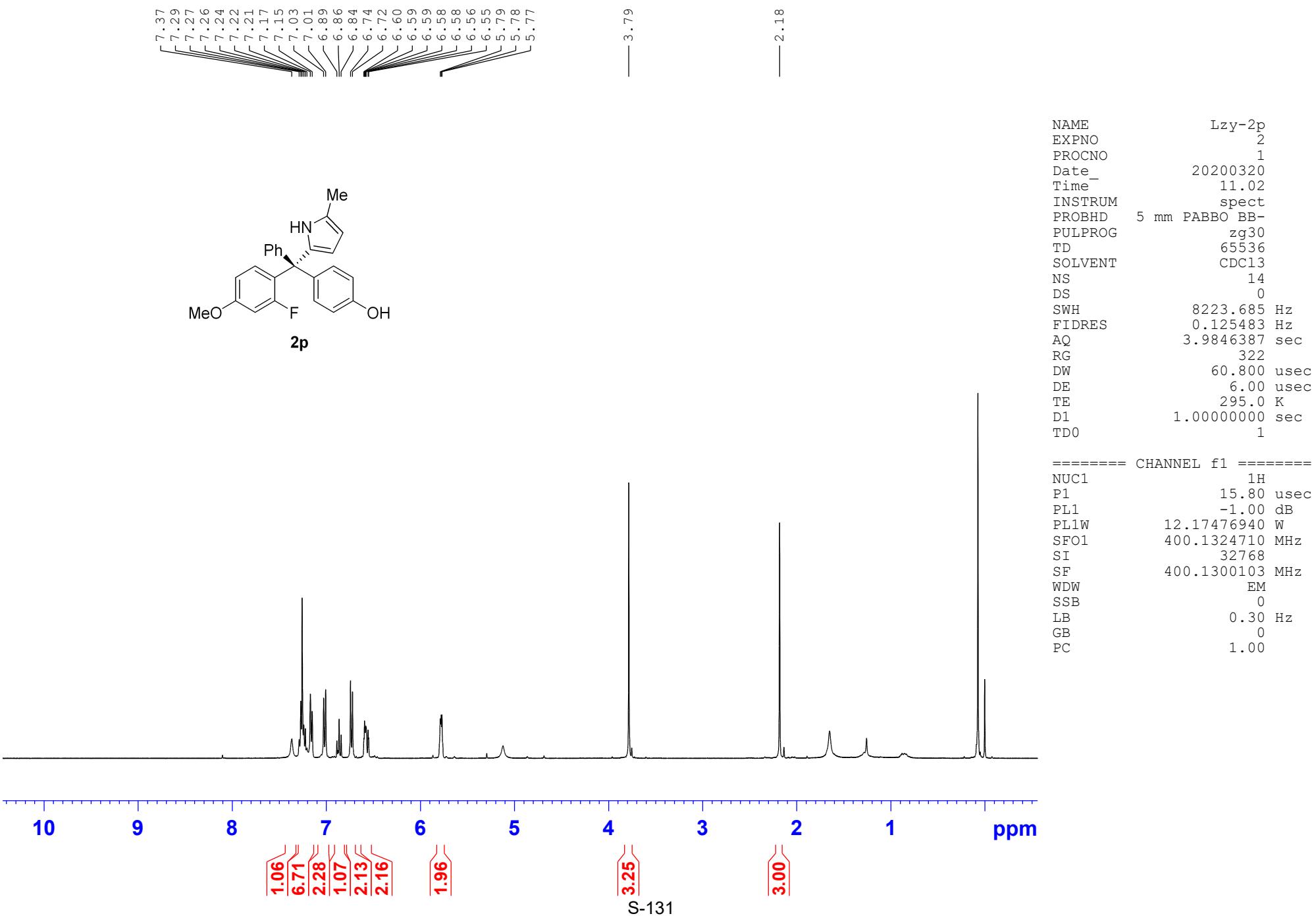
NAME lzy-project1-2o-C-0711
 EXPNO 2
 PROCNO 1
 Date 20210711
 Time 17.58
 INSTRUM spect
 PROBHD 5 mm DUL 13C-1
 PULPROG zgpg30
 TD 65536
 SOLVENT Acetone
 NS 2951
 DS 0
 SWH 24038.461 Hz
 FIDRES 0.366798 Hz
 AQ 1.3631988 sec
 RG 2050
 DW 20.800 usec
 DE 6.00 usec
 TE 293.6 K
 D1 2.0000000 sec
 D11 0.03000000 sec
 TDO 1

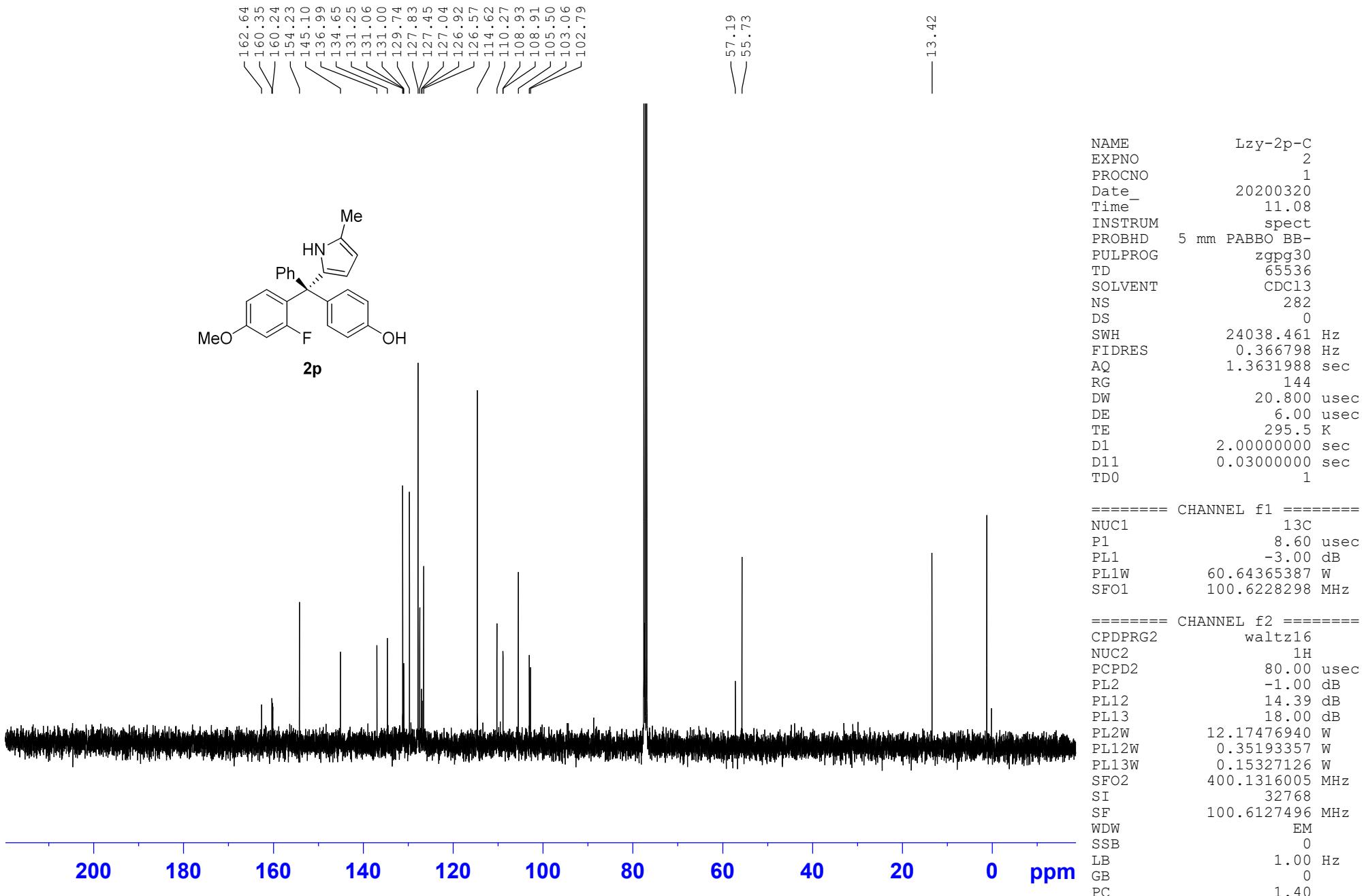
===== CHANNEL f1 =====

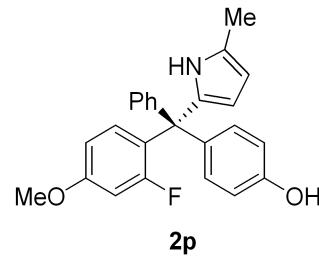
NUC1 13C
 P1 40.00 usec
 PL1 -3.00 dB
 PL1W 60.64365387 W
 SFO1 100.6228298 MHz

===== CHANNEL f2 =====

CPDPRG2 waltz16
 NUC2 1H
 PCPD2 80.00 usec
 PL2 -1.00 dB
 PL12 14.39 dB
 PL13 18.00 dB
 PL2W 12.17476940 W
 PL12W 0.35193357 W
 PL13W 0.15327126 W
 SFO2 400.1316005 MHz
 SI 32768
 SF 100.6126807 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40

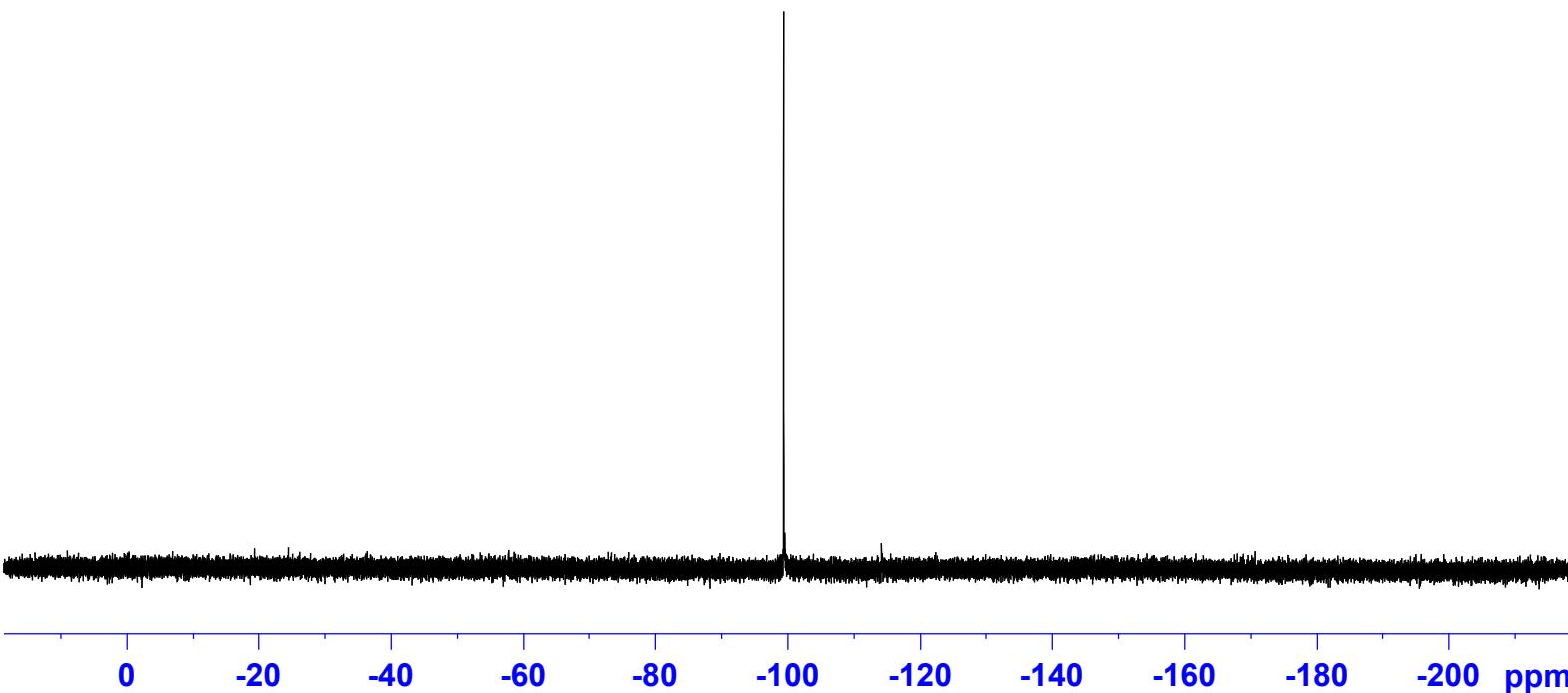


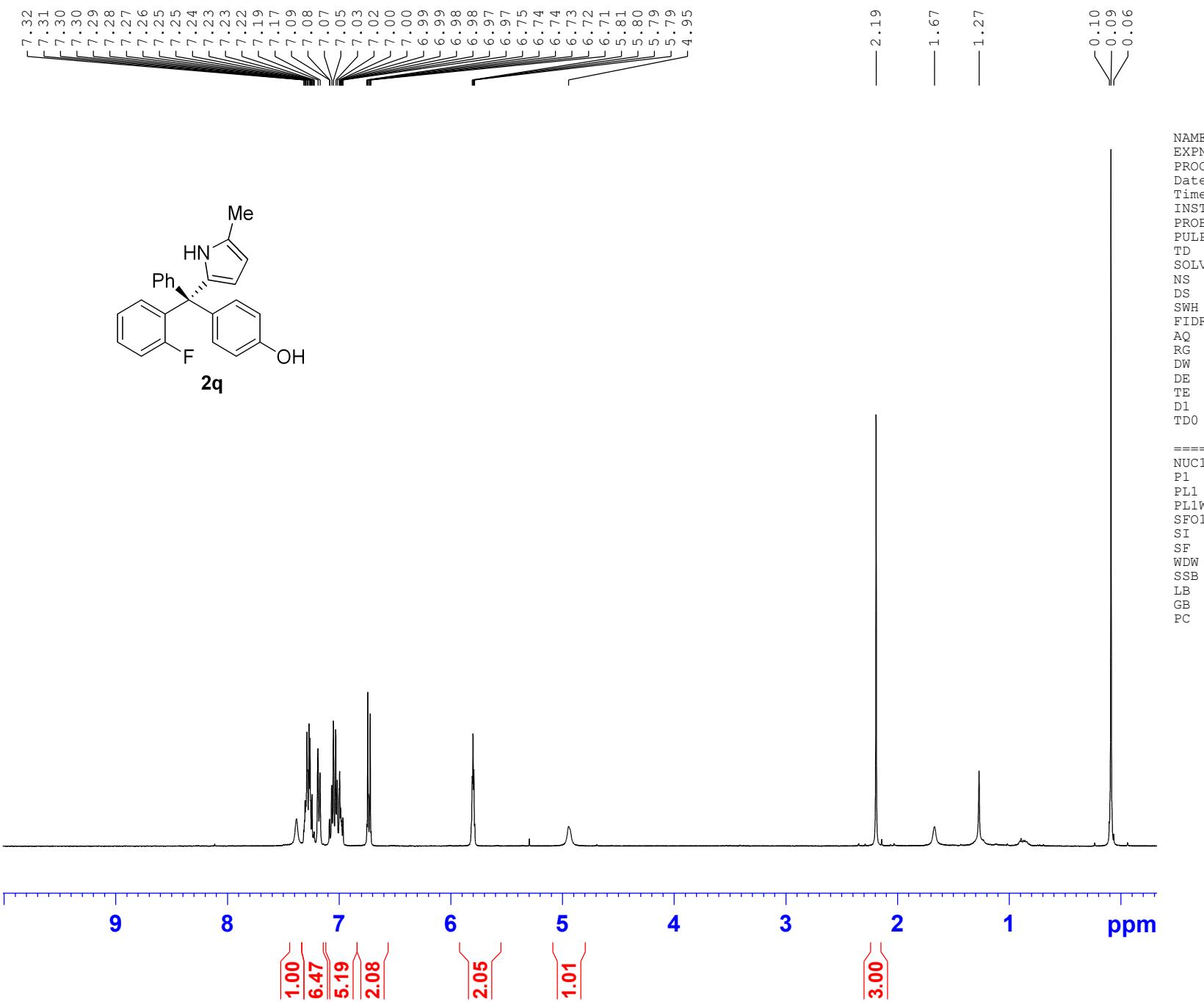


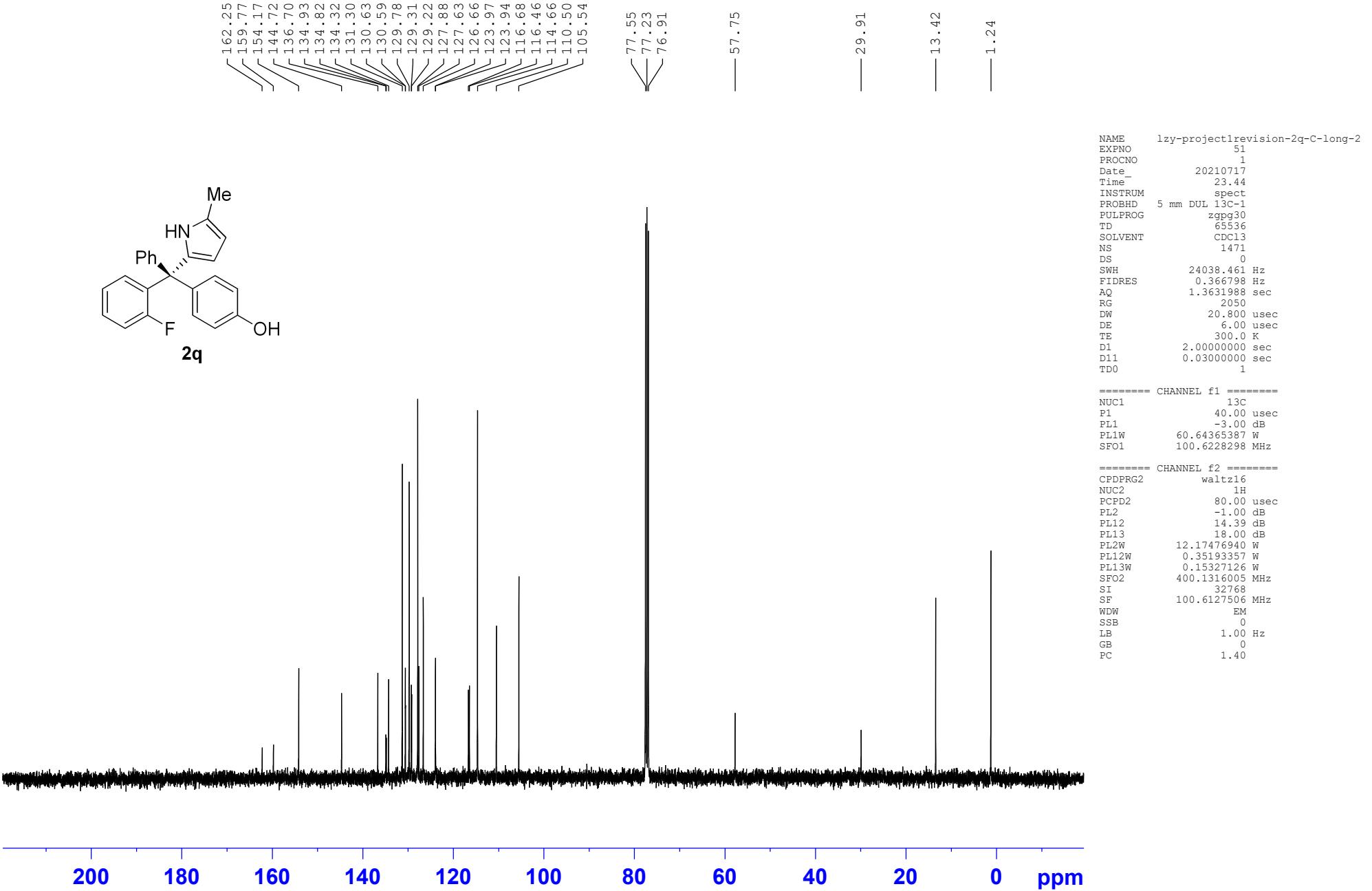


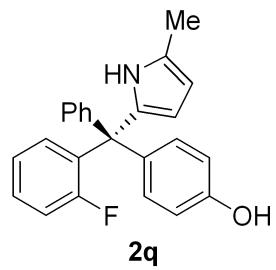
NAME Lzy-2p-F
 EXPNO 1
 PROCNO 1
 Date 20200322
 Time 18.57
 INSTRUM spect
 PROBHD 5 mm PABBO BB-
 PULPROG zgflqn
 TD 131072
 SOLVENT CDC13
 NS 1
 DS 4
 SWH 89285.711 Hz
 FIDRES 0.681196 Hz
 AQ 0.7340532 sec
 RG 1440
 DW 5.600 usec
 DE 6.00 usec
 TE 296.5 K
 D1 1.0000000 sec
 TD0 1

===== CHANNEL f1 =====
 NUC1 19F
 P1 19.50 usec
 PL1 -4.00 dB
 PL1W 16.97275162 W
 SFO1 376.4607164 MHz
 SI 65536
 SF 376.4983660 MHz
 WDW EM
 SSB 0
 LB 0.30 Hz
 GB 0
 PC 1.00







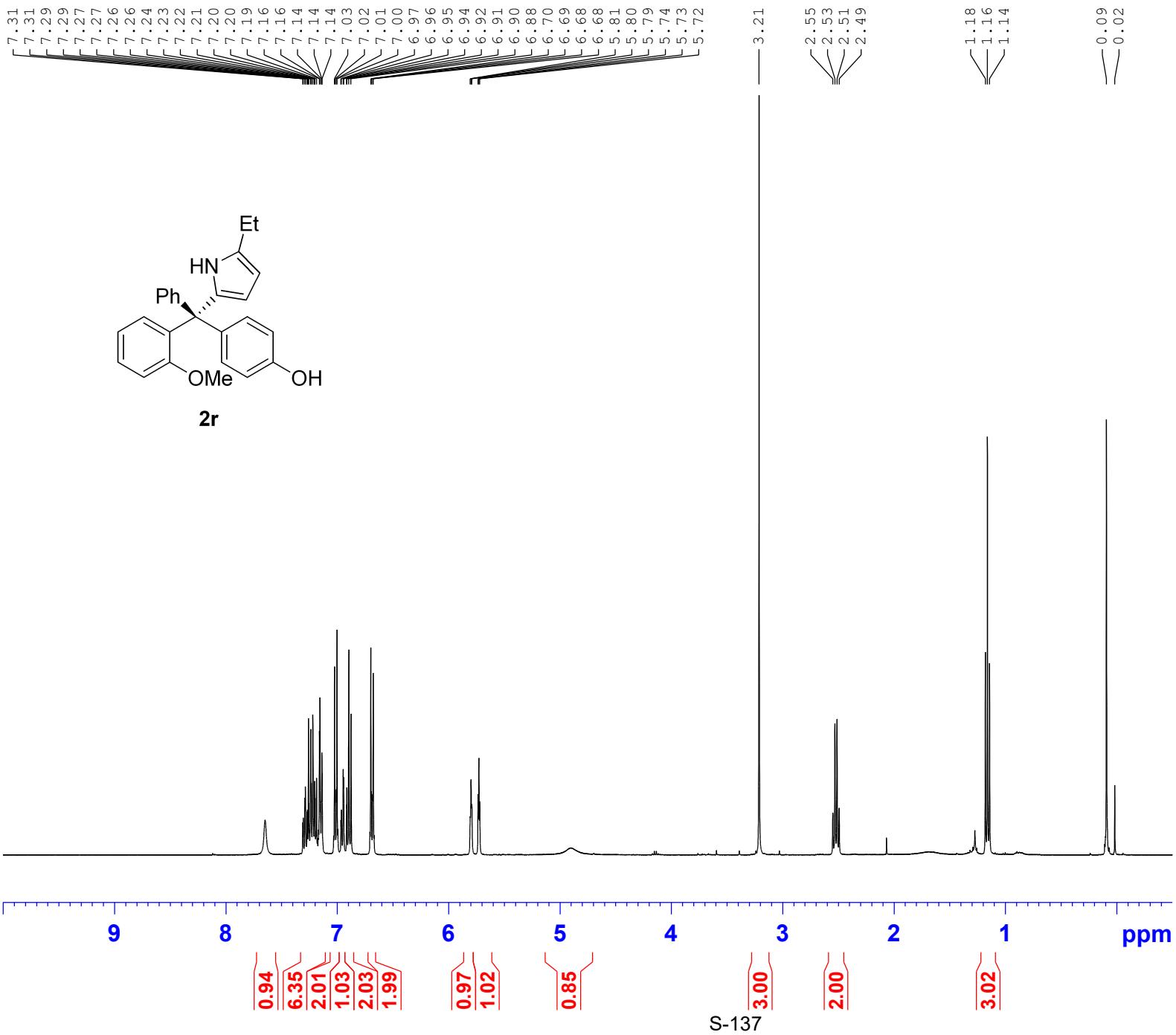


-101.32

NAME lzy-07-project1revision-2q-F
 EXPNO 1
 PROCNO 1
 Date 20210707
 Time 18.05
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgflqn
 TD 131072
 SOLVENT CDCl3
 NS 2
 DS 4
 SWH 89285.711 Hz
 FIDRES 0.681196 Hz
 AQ 0.7340532 sec
 RG 196.92
 DW 5.600 usec
 DE 6.50 usec
 TE 296.8 K
 D1 1.0000000 sec
 TDO 1

===== CHANNEL f1 =====
 SFO1 376.4607164 MHz
 NUC1 19F
 P1 14.70 usec
 SI 65536
 SF 376.4983660 MHz
 WDW EM
 SSB 0
 LB 0.30 Hz
 GB 0
 PC 1.00

0 -20 -40 -60 -80 -100 -120 -140 -160 -180 -200 ppm

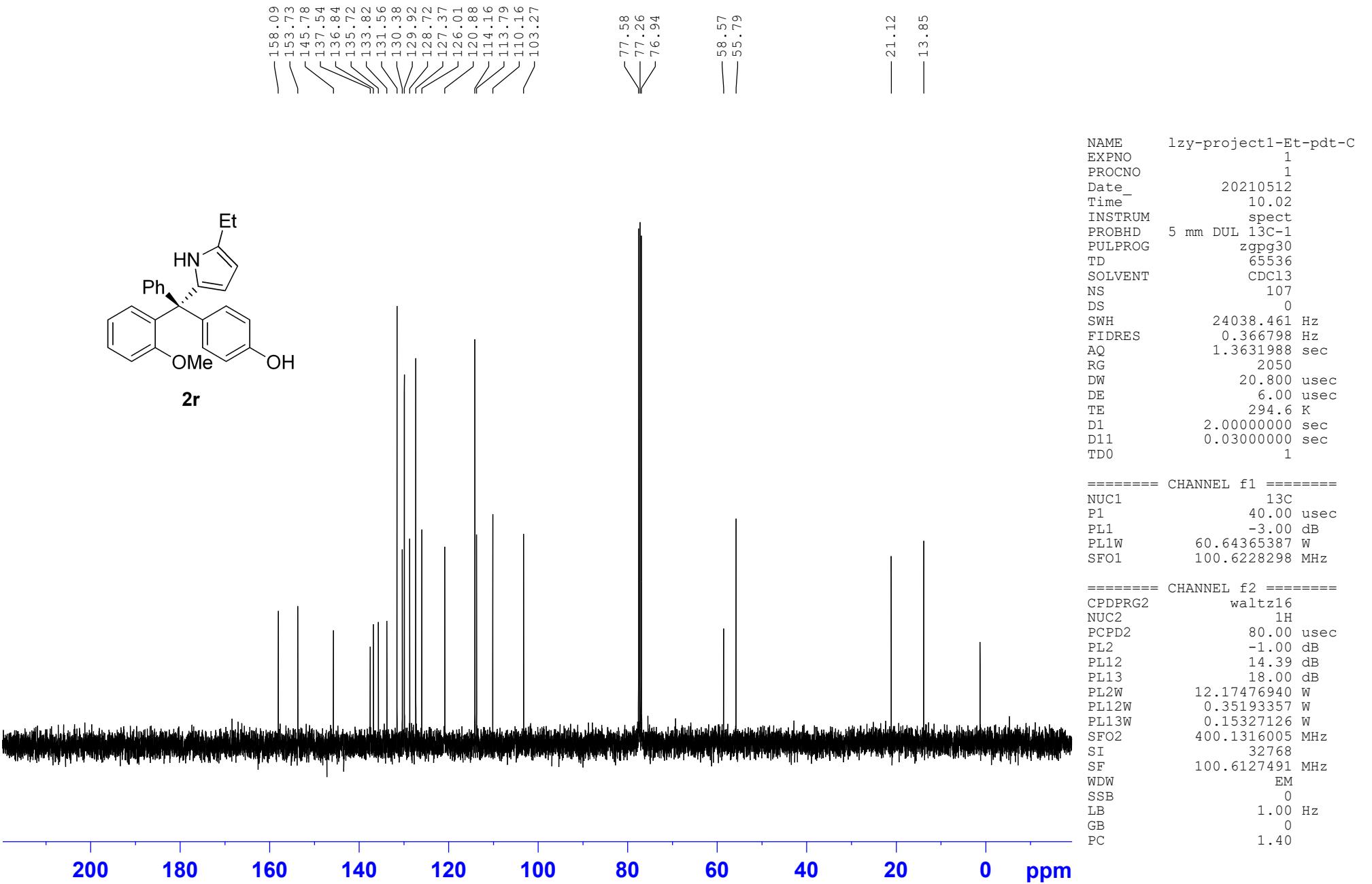


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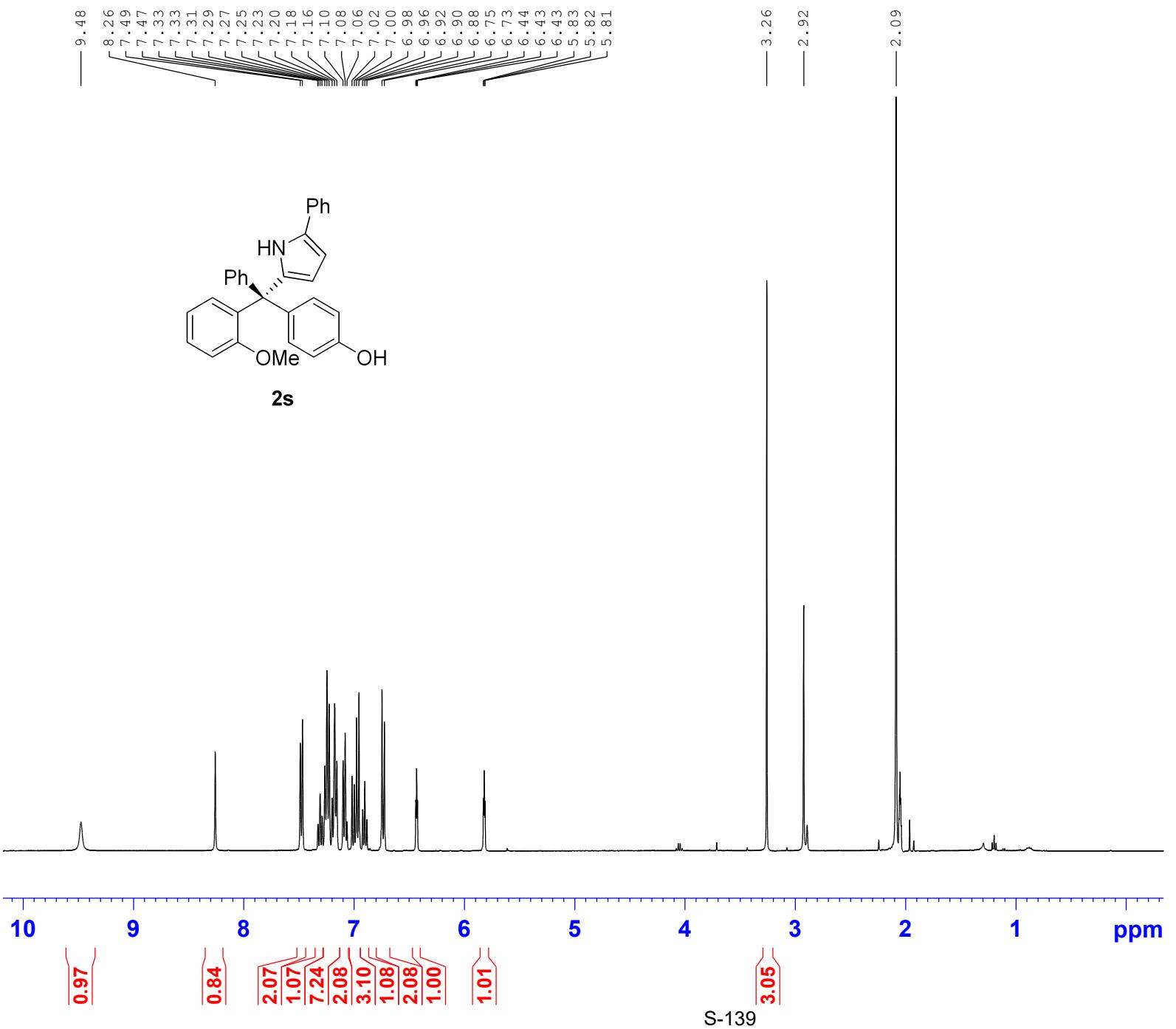
NAME      lzy-project1-Et-pdt
EXPNO     1
PROCNO    1
Date      20210512
Time      10.00
INSTRUM   spect
PROBHD   5 mm DUL 13C-1
PULPROG  zg30
TD       65536
SOLVENT   CDC13
NS        4
DS         0
SWH      8223.685 Hz
FIDRES   0.125483 Hz
AQ       3.9846387 sec
RG        203
DW       60.800 usec
DE        6.00 usec
TE       294.4 K
D1      1.00000000 sec
TD0          1

===== CHANNEL f1 =====
NUC1      1H
P1        15.80 usec
PL1      -1.00 dB
PL1W    12.17476940 W
SFO1    400.1324710 MHz
SI        32768
SF      400.1300099 MHz
WDW        EM
SSB        0
LB        0.30 Hz
GB        0
PC        1.00

```

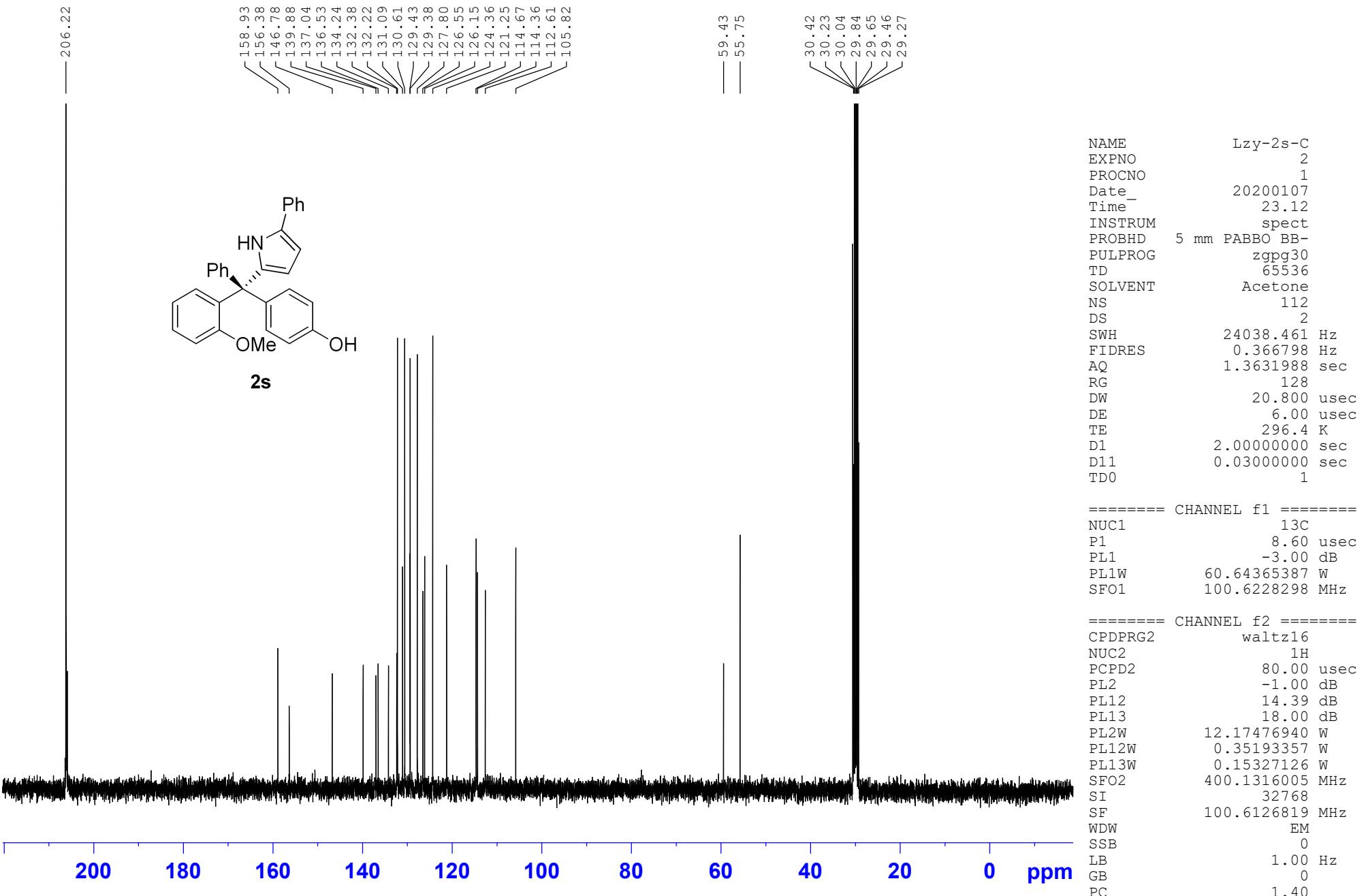


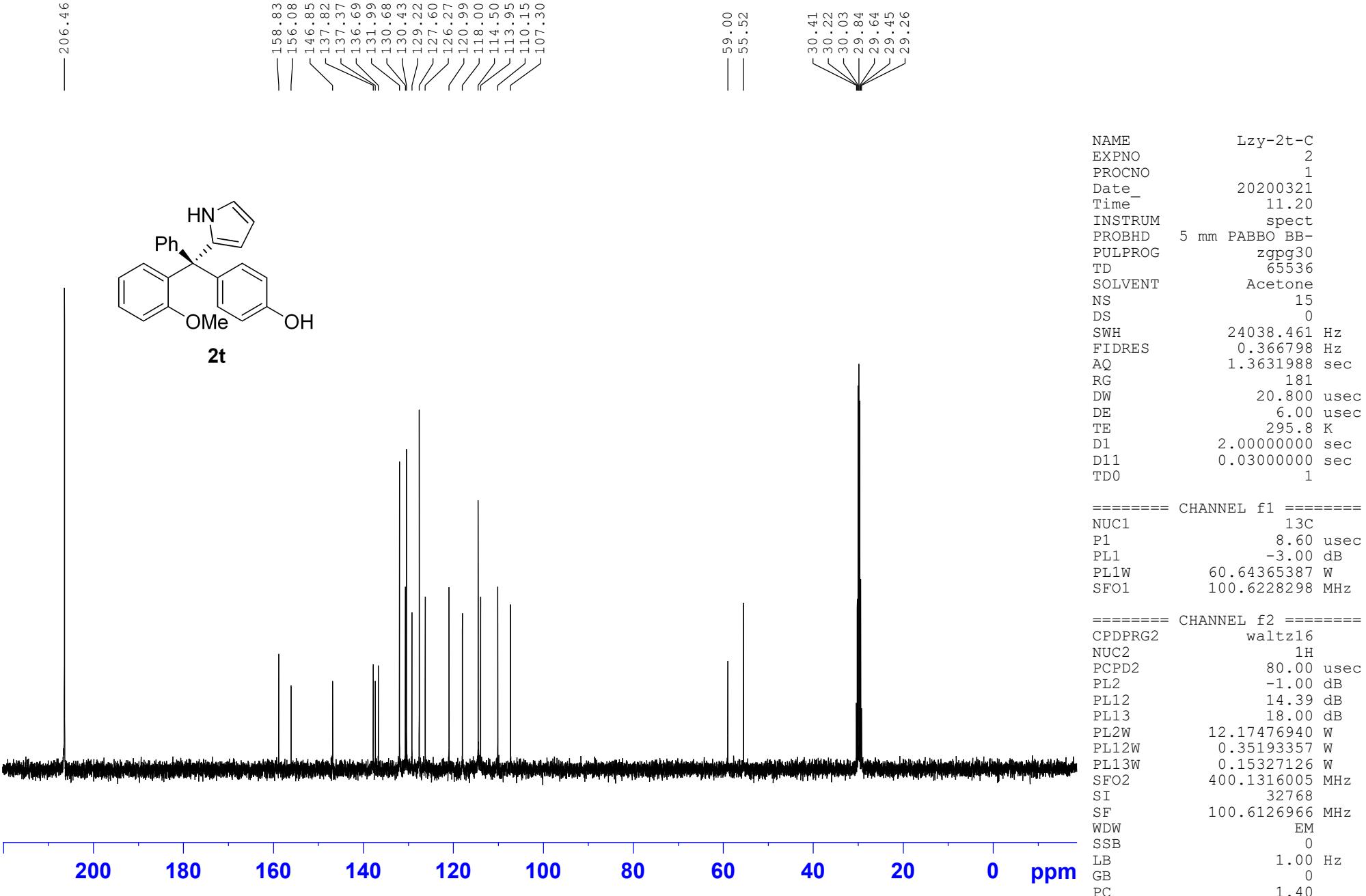
200 180 160 140 120 100 80 60 40 20 0 ppm

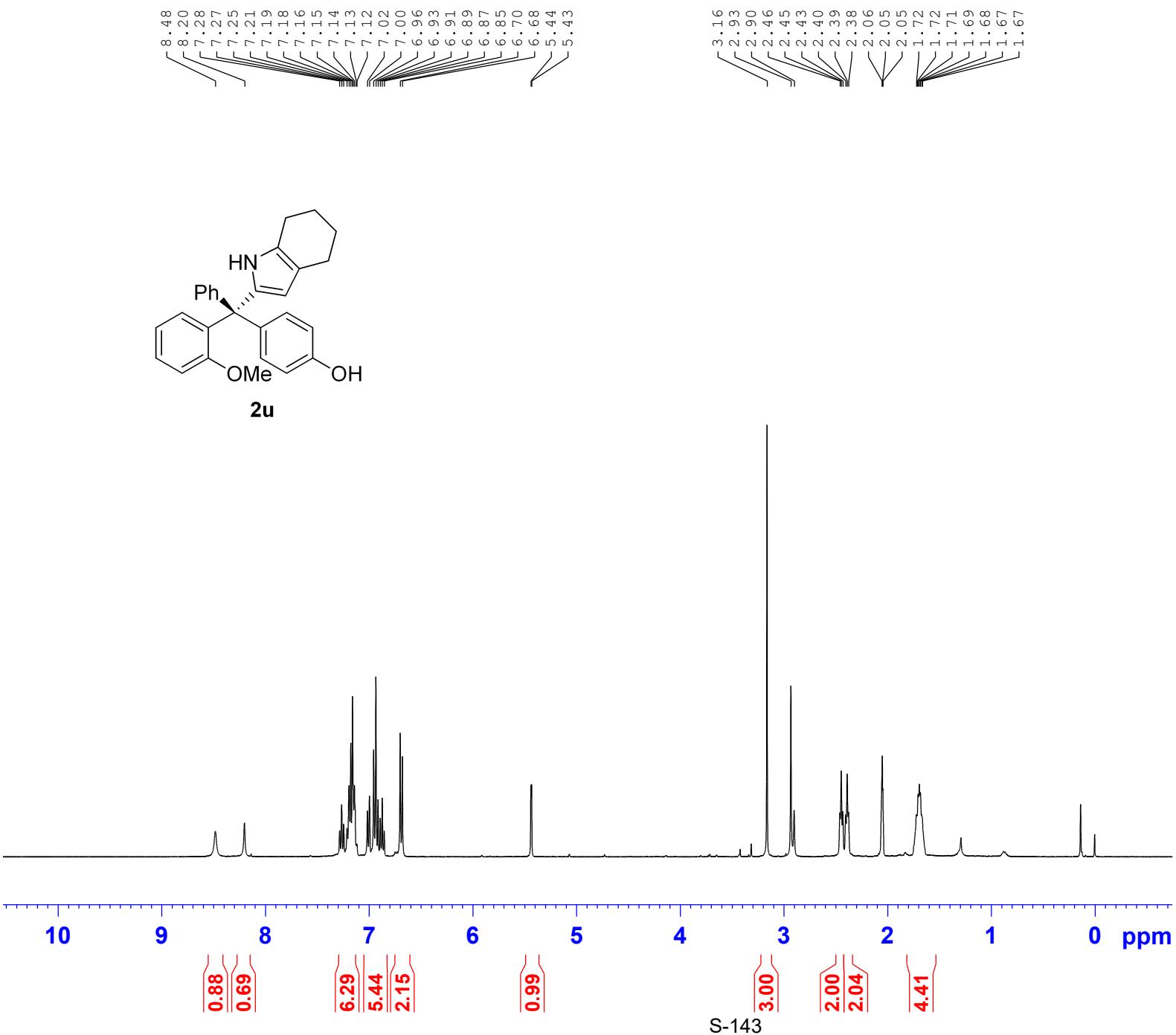


NAME Lzy-2s
 EXPNO 2
 PROCNO 1
 Date 20200107
 Time 23.06
 INSTRUM spect
 PROBHD 5 mm PABBO BB-
 PULPROG zg30
 TD 65536
 SOLVENT Acetone
 NS 16
 DS 2
 SWH 8223.685 Hz
 FIDRES 0.125483 Hz
 AQ 3.9846387 sec
 RG 161
 DW 60.800 usec
 DE 6.00 usec
 TE 295.8 K
 D1 1.00000000 sec
 TD0 1

===== CHANNEL f1 =====
 NUC1 1H
 P1 15.80 usec
 PL1 -1.00 dB
 PL1W 12.17476940 W
 SFO1 400.1324710 MHz
 SI 32768
 SF 400.1300065 MHz
 WDW EM
 SSB 0
 LB 0.30 Hz
 GB 0
 PC 1.00





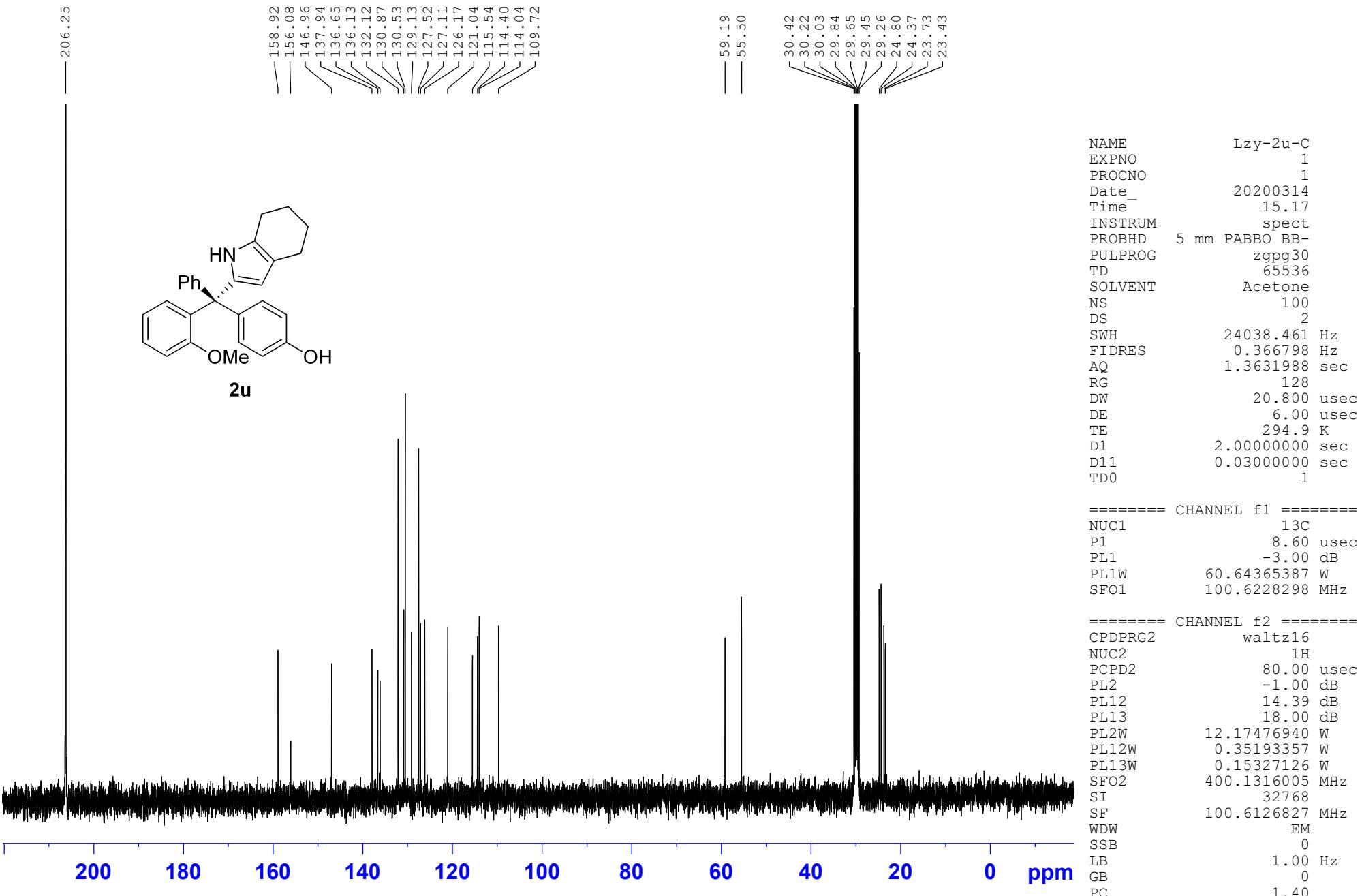


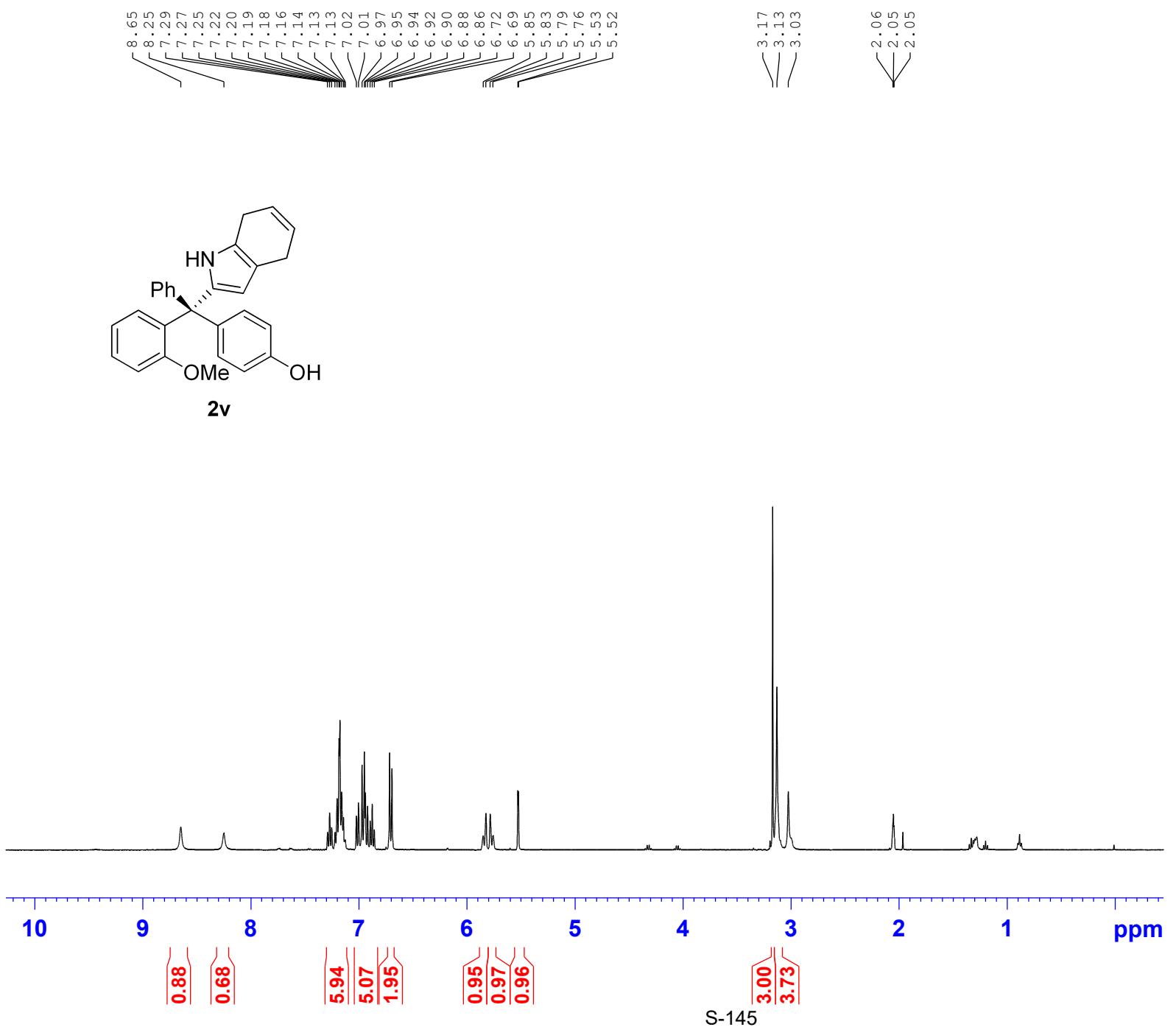
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NAME          Lzy-2u
EXPNO         1
PROCNO        1
Date 20200314
Time 15.15
INSTRUM spect
PROBHD 5 mm PABBO BB-
PULPROG zg30
TD 65536
SOLVENT Acetone
NS 12
DS 0
SWH 8223.685 Hz
FIDRES 0.125483 Hz
AQ 3.9846387 sec
RG 406
DW 60.800 usec
DE 6.00 usec
TE 294.7 K
D1 1.00000000 sec
TD0 1

===== CHANNEL f1 =====
NUC1           1H
P1 15.80 usec
PL1 -1.00 dB
PL1W 12.17476940 W
SFO1 400.1324710 MHz
SI 32768
SF 400.1300054 MHz
WDW EM
SSB 0
LB 0.30 Hz
GB 0
PC 1.00

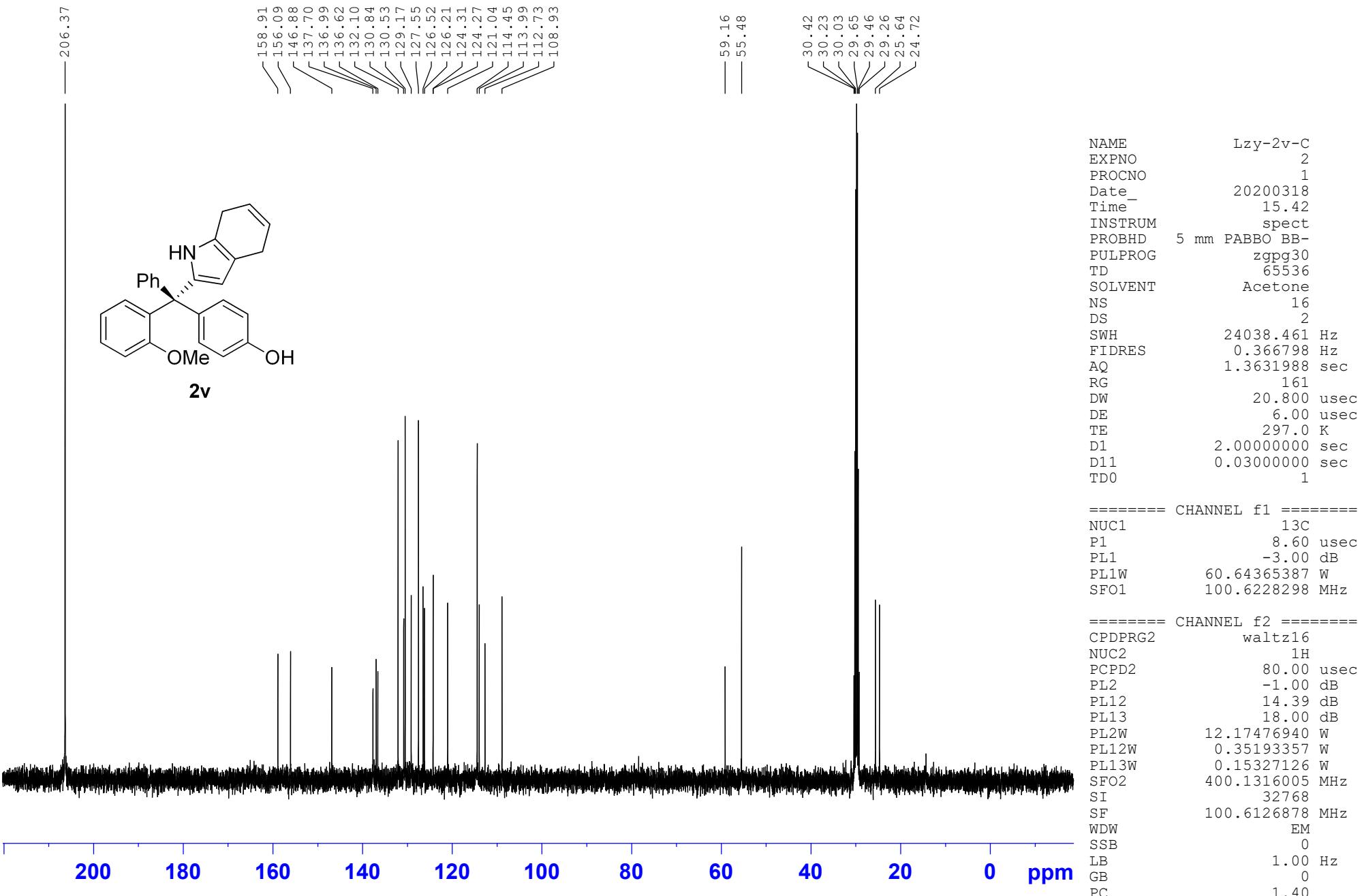
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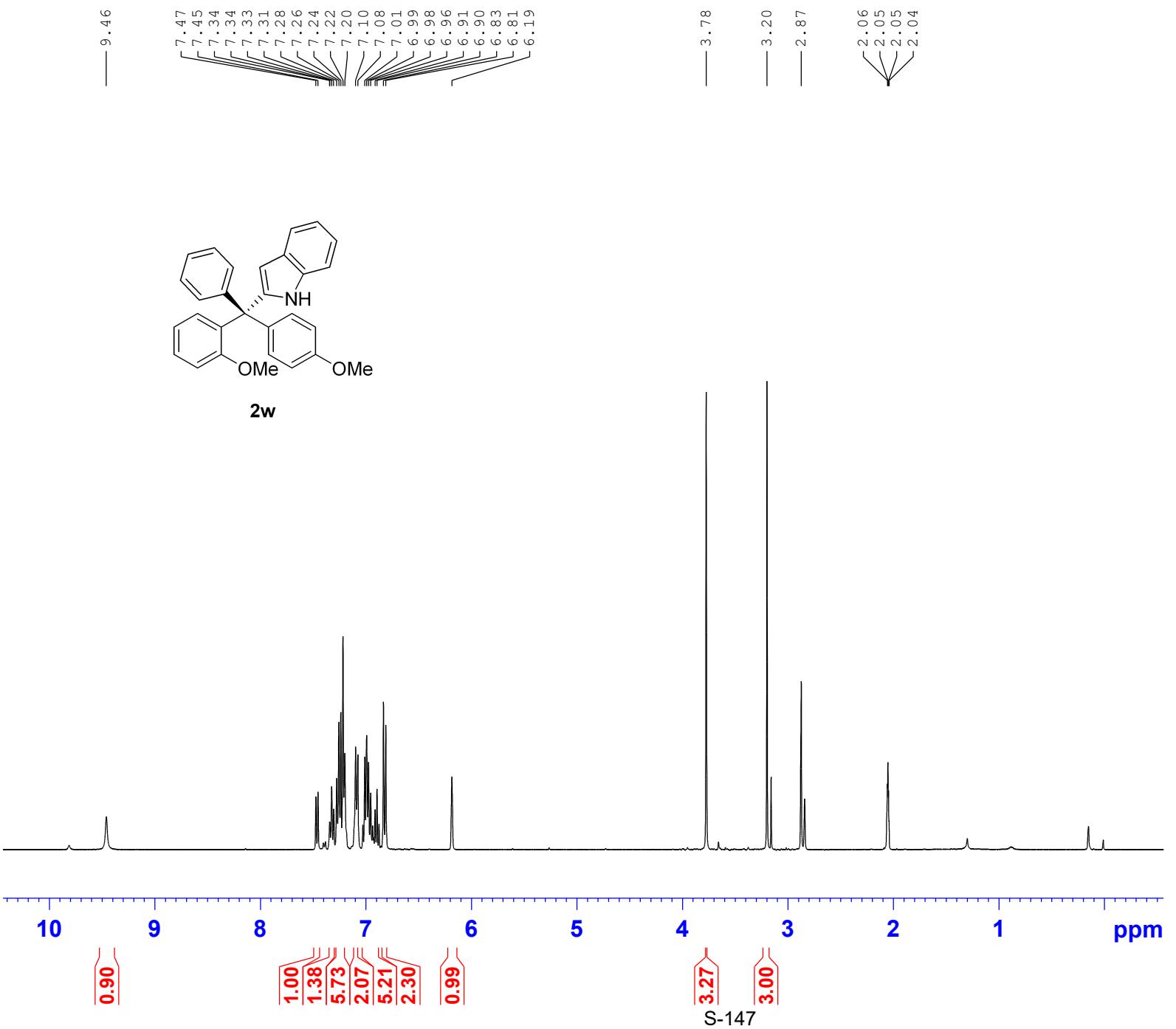




NAME Lzy-2v
 EXPNO 4
 PROCNO 1
 Date 20200318
 Time 20.08
 INSTRUM spect
 PROBHD 5 mm PABBO BB-
 PULPROG zg30
 TD 65536
 SOLVENT Acetone
 NS 2
 DS 0
 SWH 8223.685 Hz
 FIDRES 0.125483 Hz
 AQ 3.9846387 sec
 RG 114
 DW 60.800 usec
 DE 6.00 usec
 TE 295.0 K
 D1 1.00000000 sec
 TD0 1

===== CHANNEL f1 =====
 NUC1 1H
 P1 15.80 usec
 PL1 -1.00 dB
 PL1W 12.17476940 W
 SFO1 400.1324710 MHz
 SI 32768
 SF 400.1300054 MHz
 WDW EM
 SSB 0
 LB 0.30 Hz
 GB 0
 PC 1.00





```

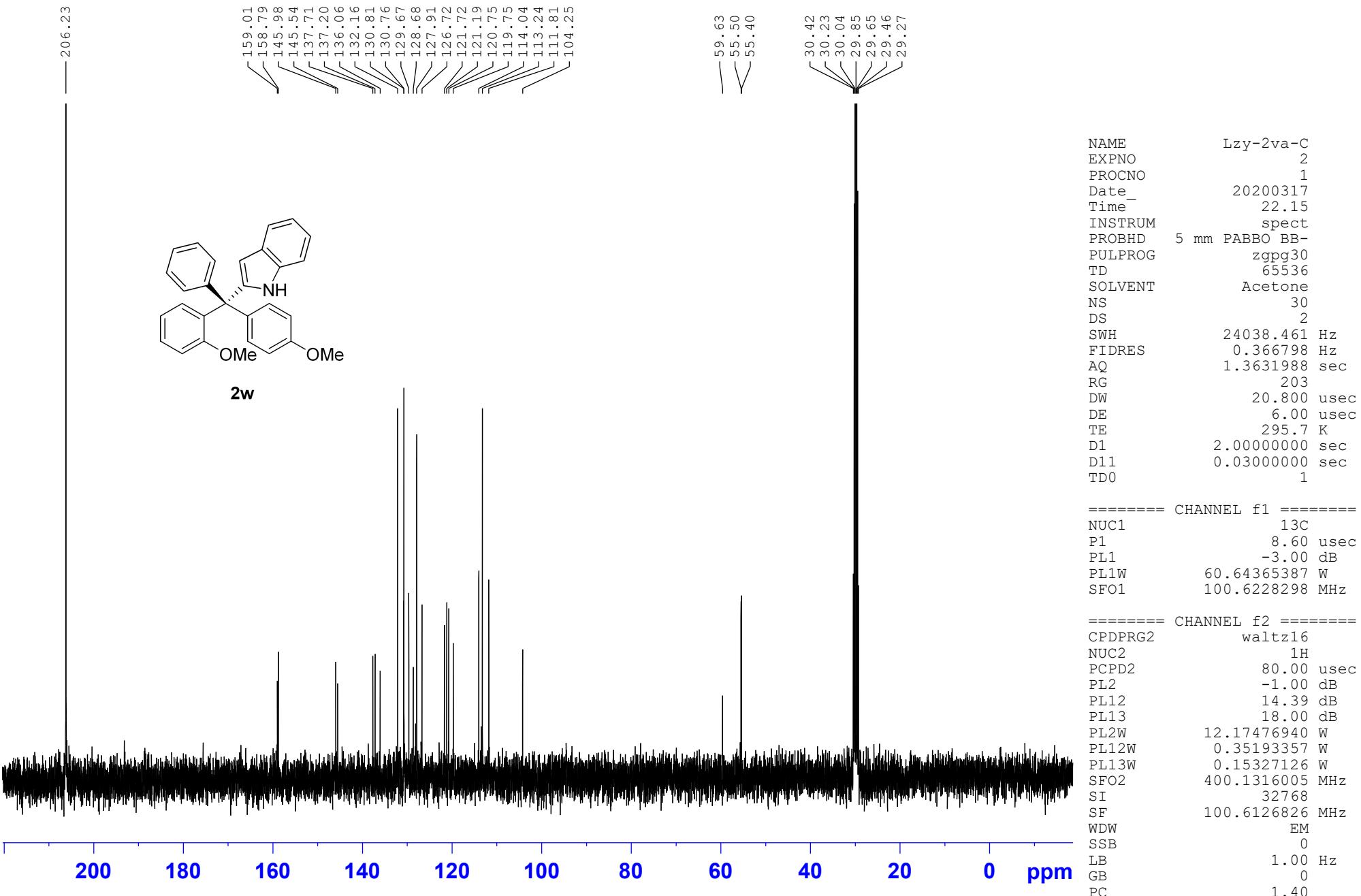
NAME          Lzy-2va
EXPNO         2
PROCNO        1
Date_ 20200317
Time_ 22.12
INSTRUM      spect
PROBHD 5 mm PABBO BB-
PULPROG zg30
TD        65536
SOLVENT      Acetone
NS          4
DS          0
SWH        8223.685 Hz
FIDRES       0.125483 Hz
AQ        3.9846387 sec
RG          161
DW        60.800 usec
DE          6.00 usec
TE        295.3 K
D1        1.00000000 sec
TD0          1

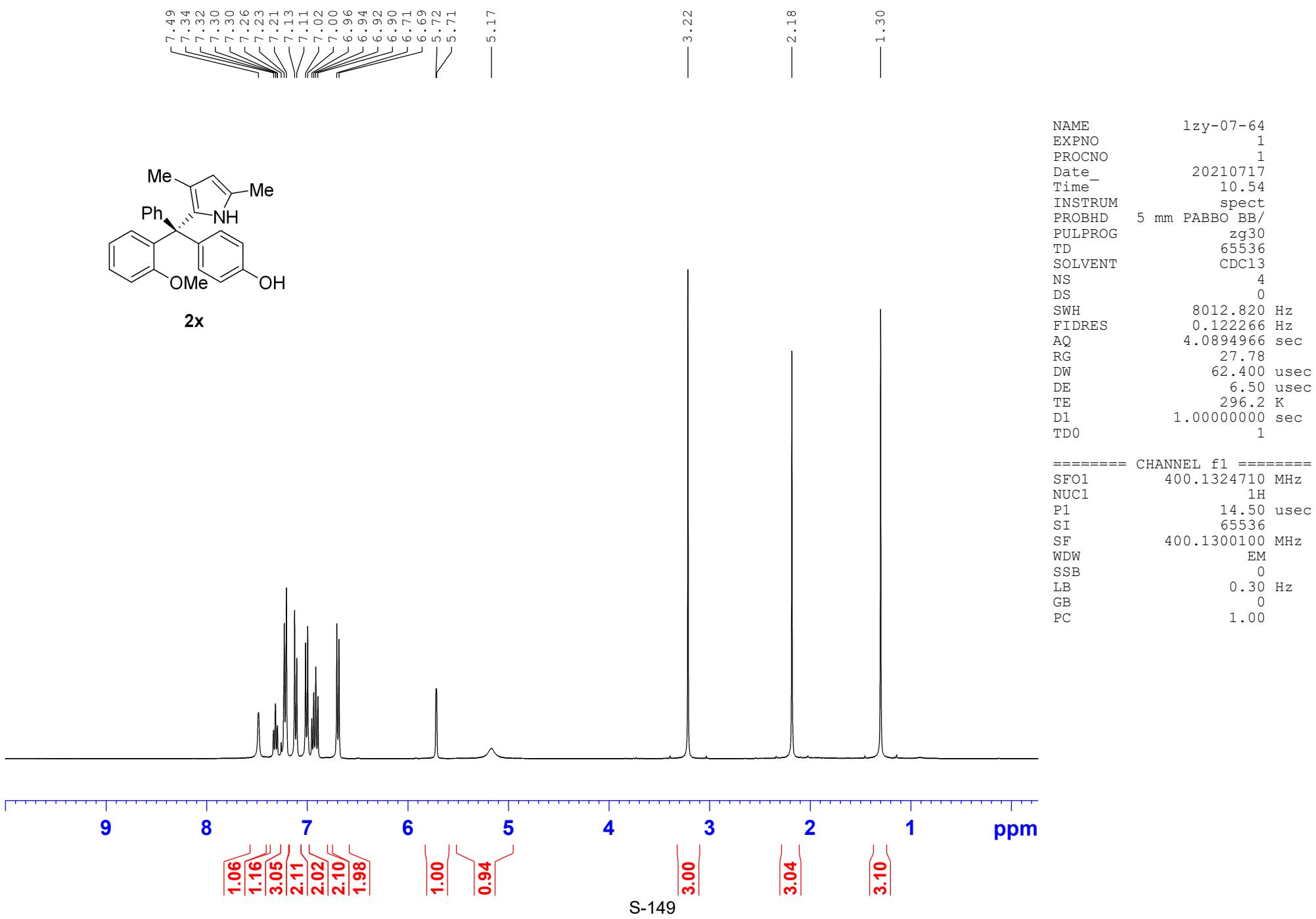
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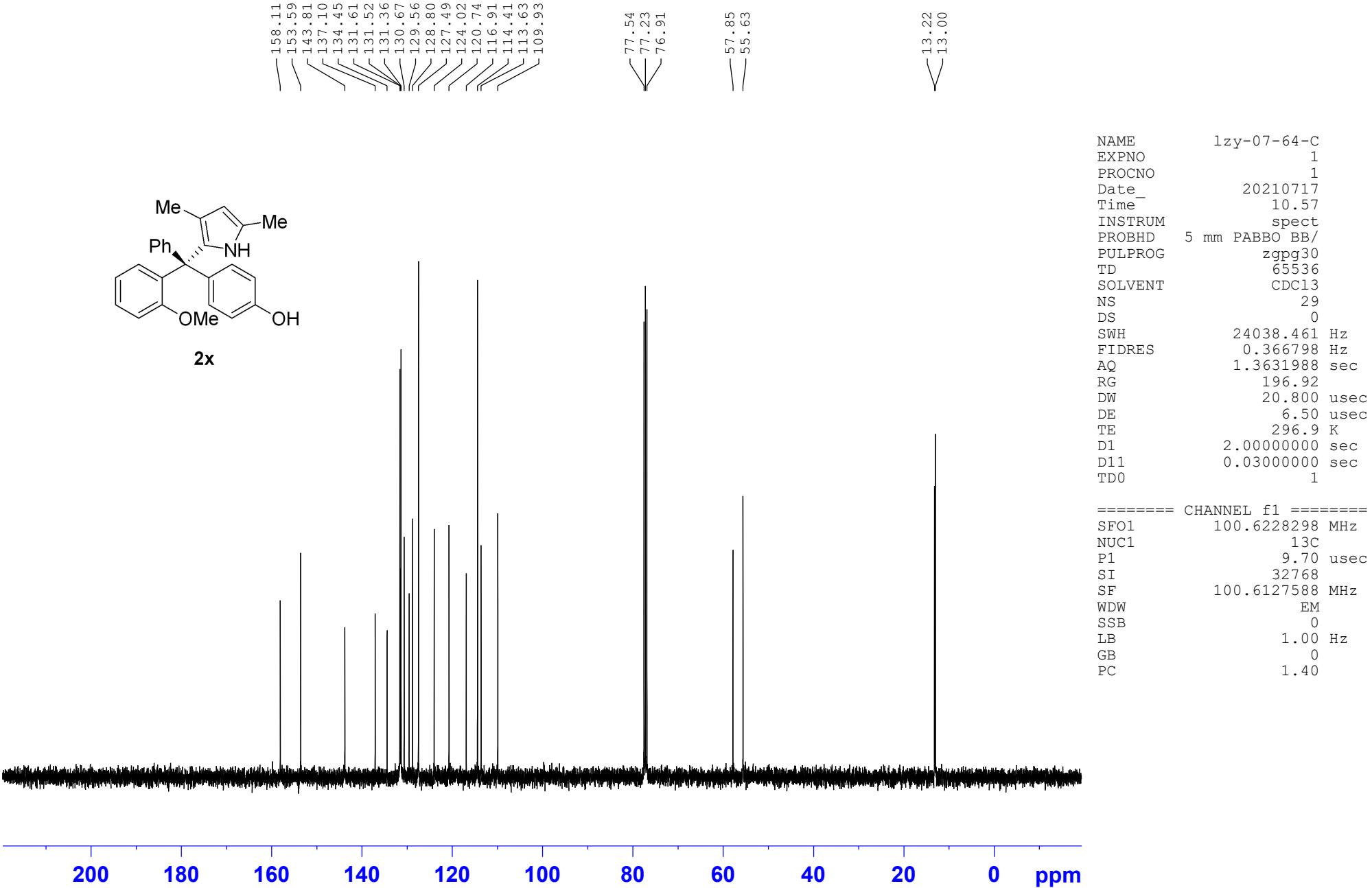
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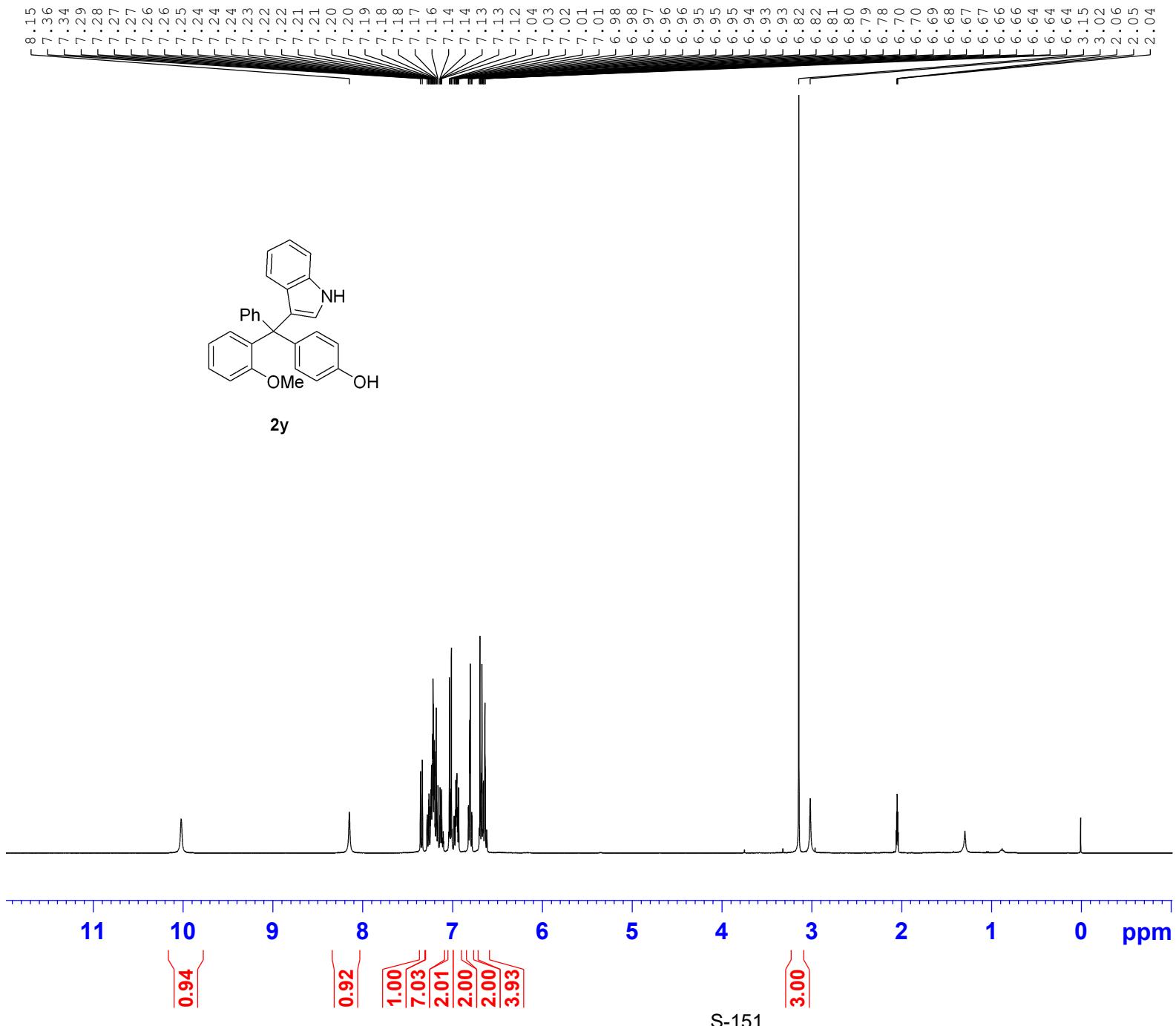
===== CHANNEL f1 =====
NUC1           1H
P1        15.80 usec
PL1        -1.00 dB
PL1W      12.17476940 W
SFO1      400.1324710 MHz
SI          32768
SF      400.1300054 MHz
WDW            EM
SSB            0
LB        0.30 Hz
GB            0
PC          1.00

```





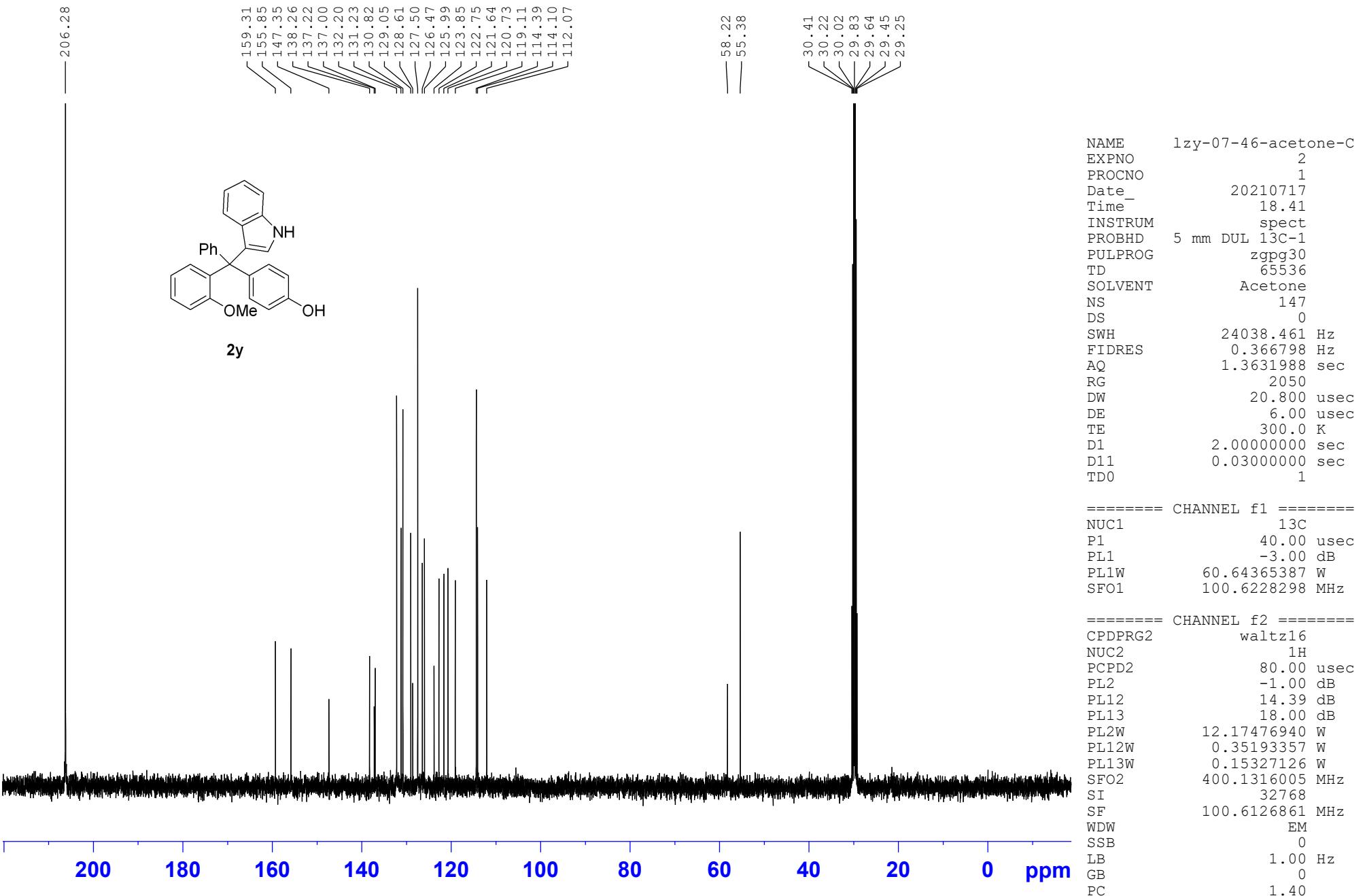


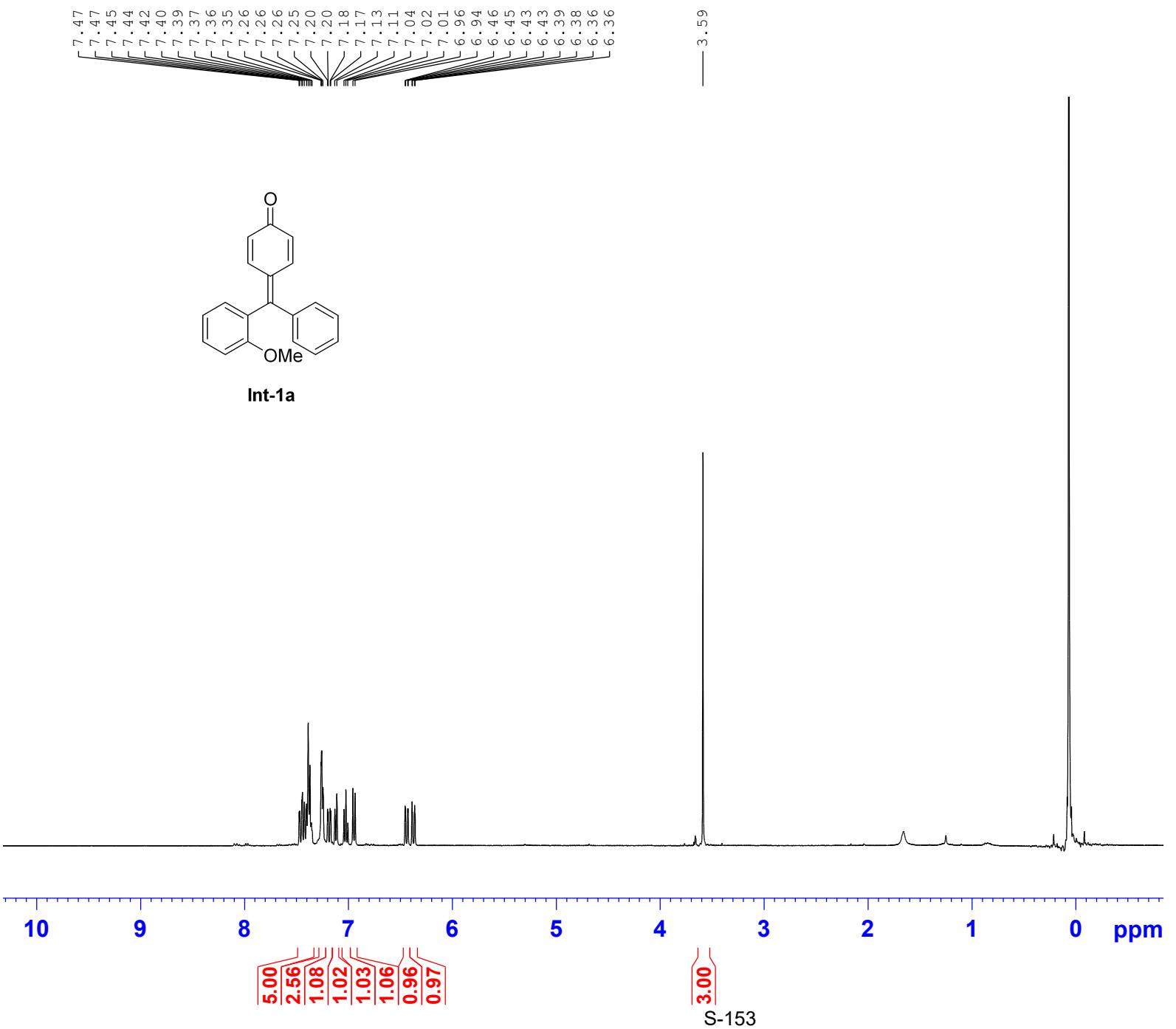


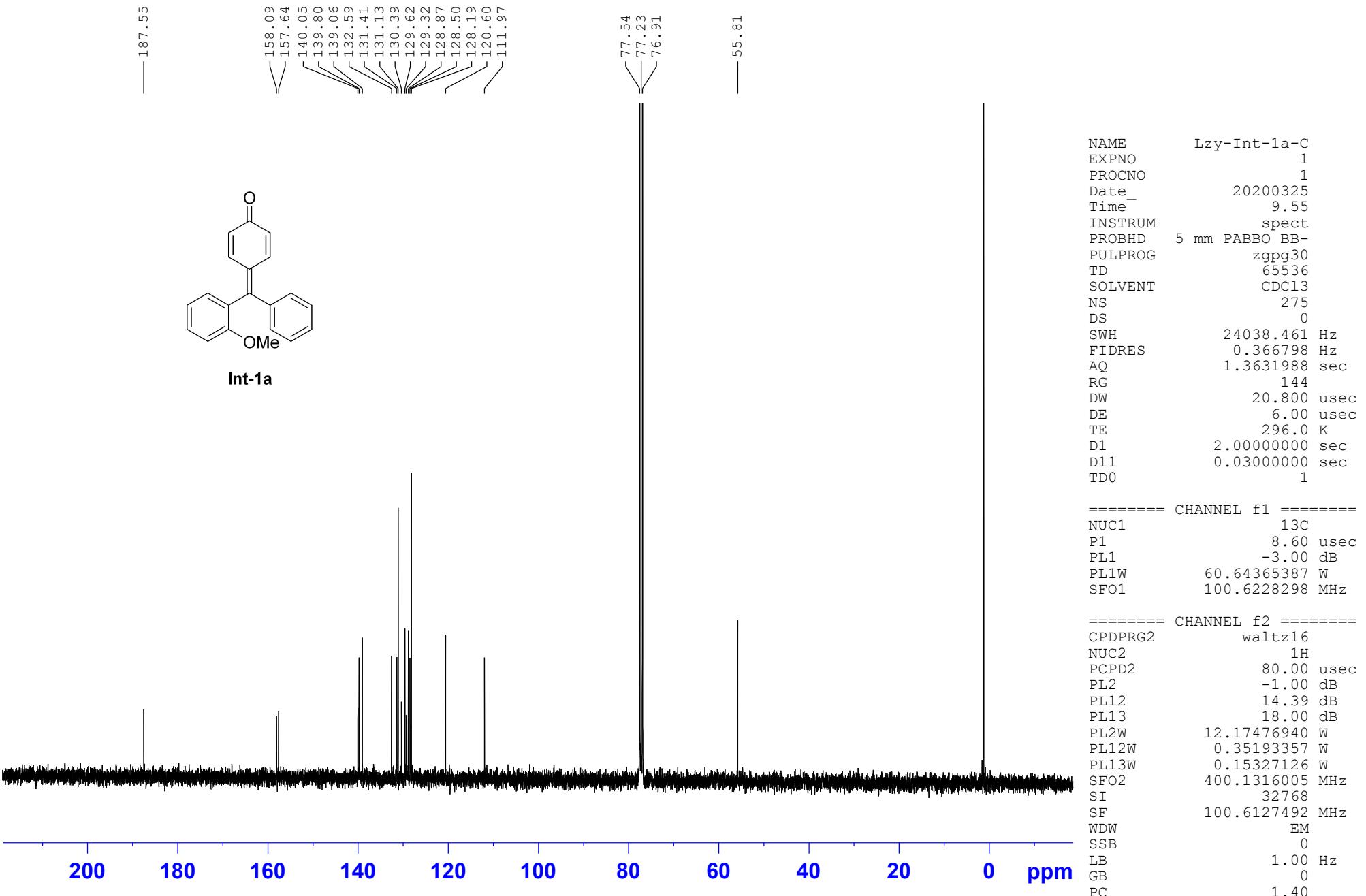
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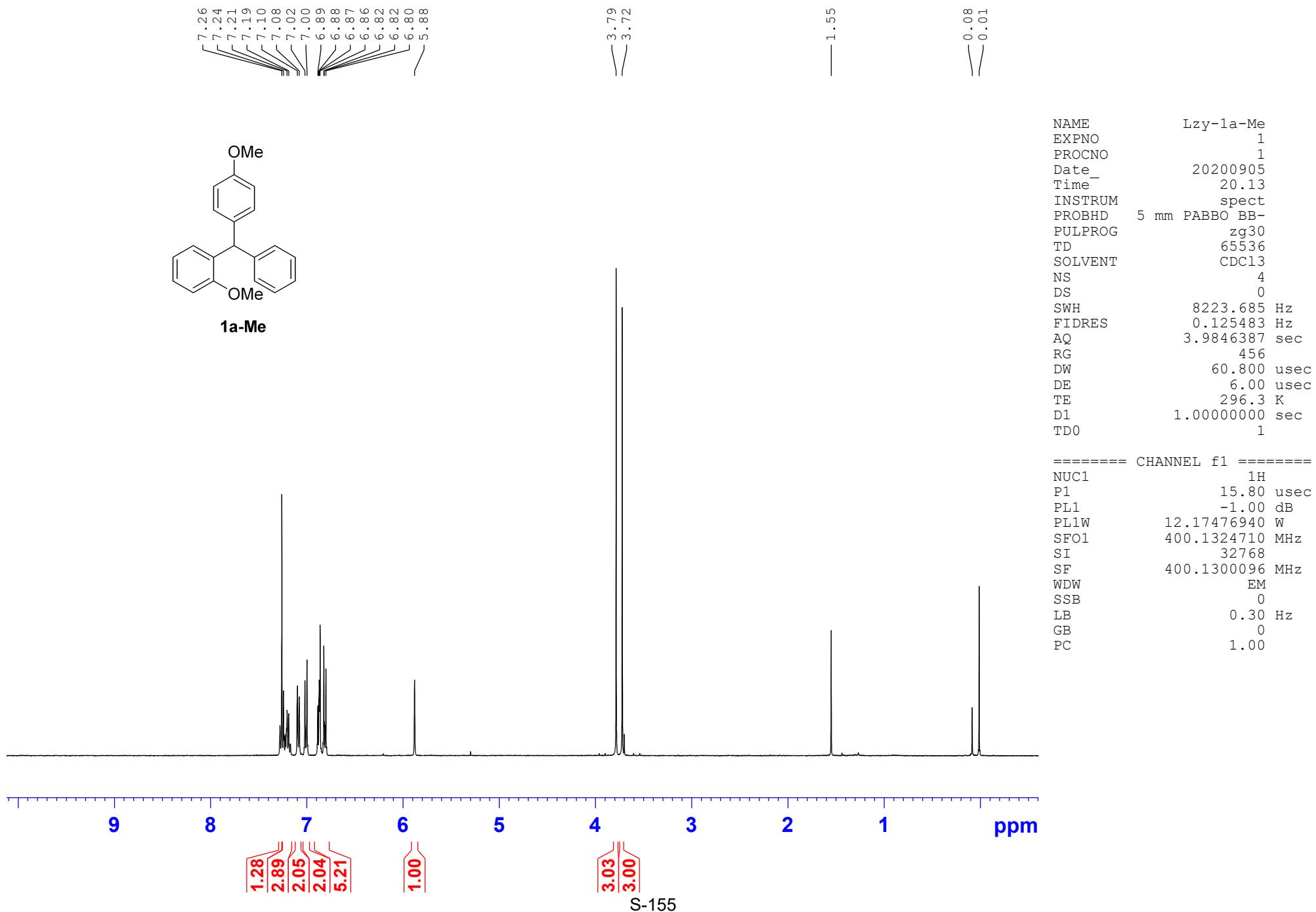
NAME      lzy-07-46-acetone-H
EXPNO          1
PROCNO         1
Date_        20210717
Time         18.39
INSTRUM      spect
PROBHD      5 mm DUL 13C-1
PULPROG     zg30
TD           65536
SOLVENT      Acetone
NS            3
DS            0
SWH         8223.685 Hz
FIDRES     0.125483 Hz
AQ        3.9846387 sec
RG           181
DW         60.800 usec
DE           6.00 usec
TE         300.0 K
D1      1.000000000 sec
TD0             1

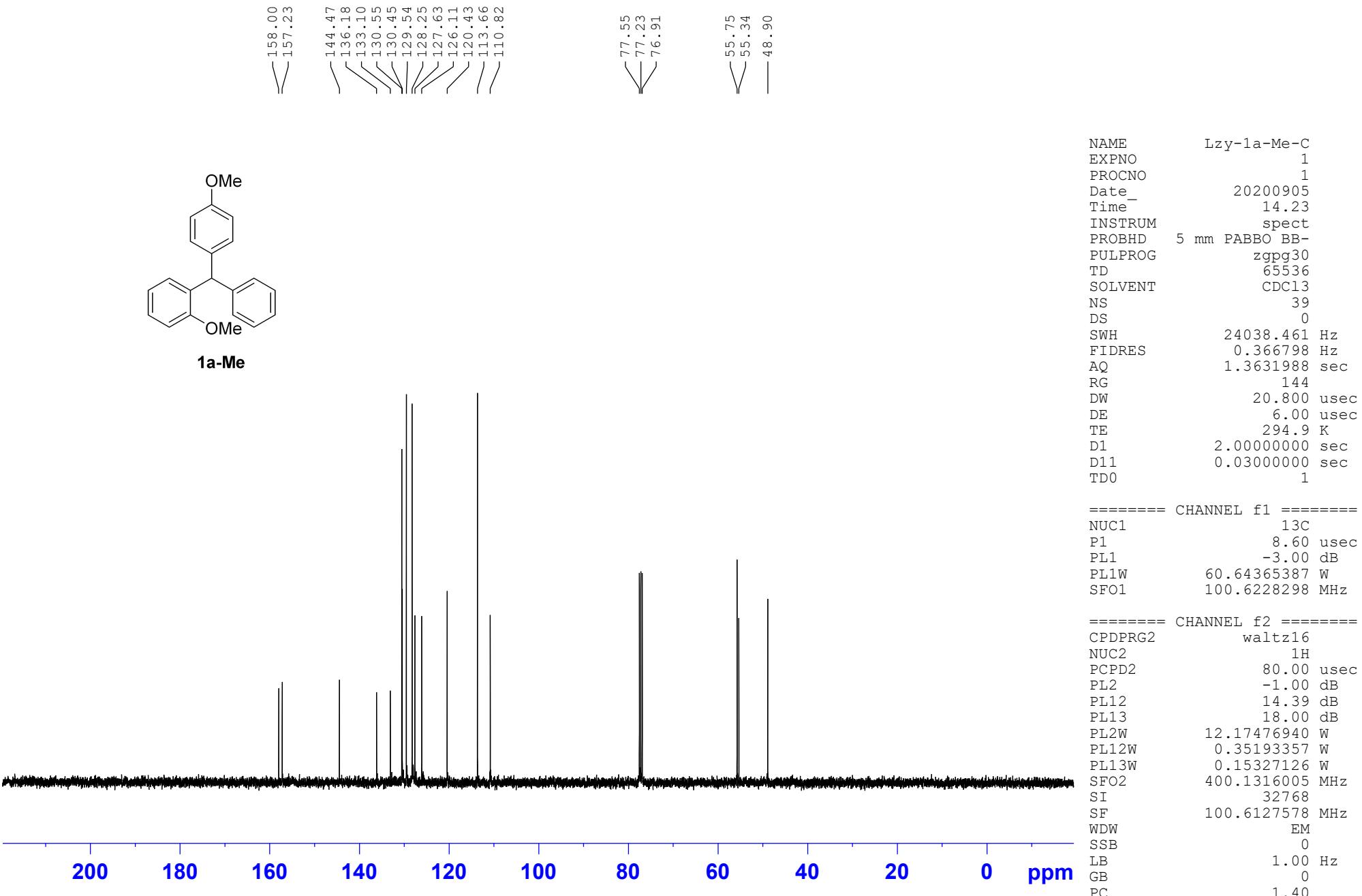
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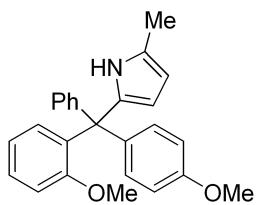
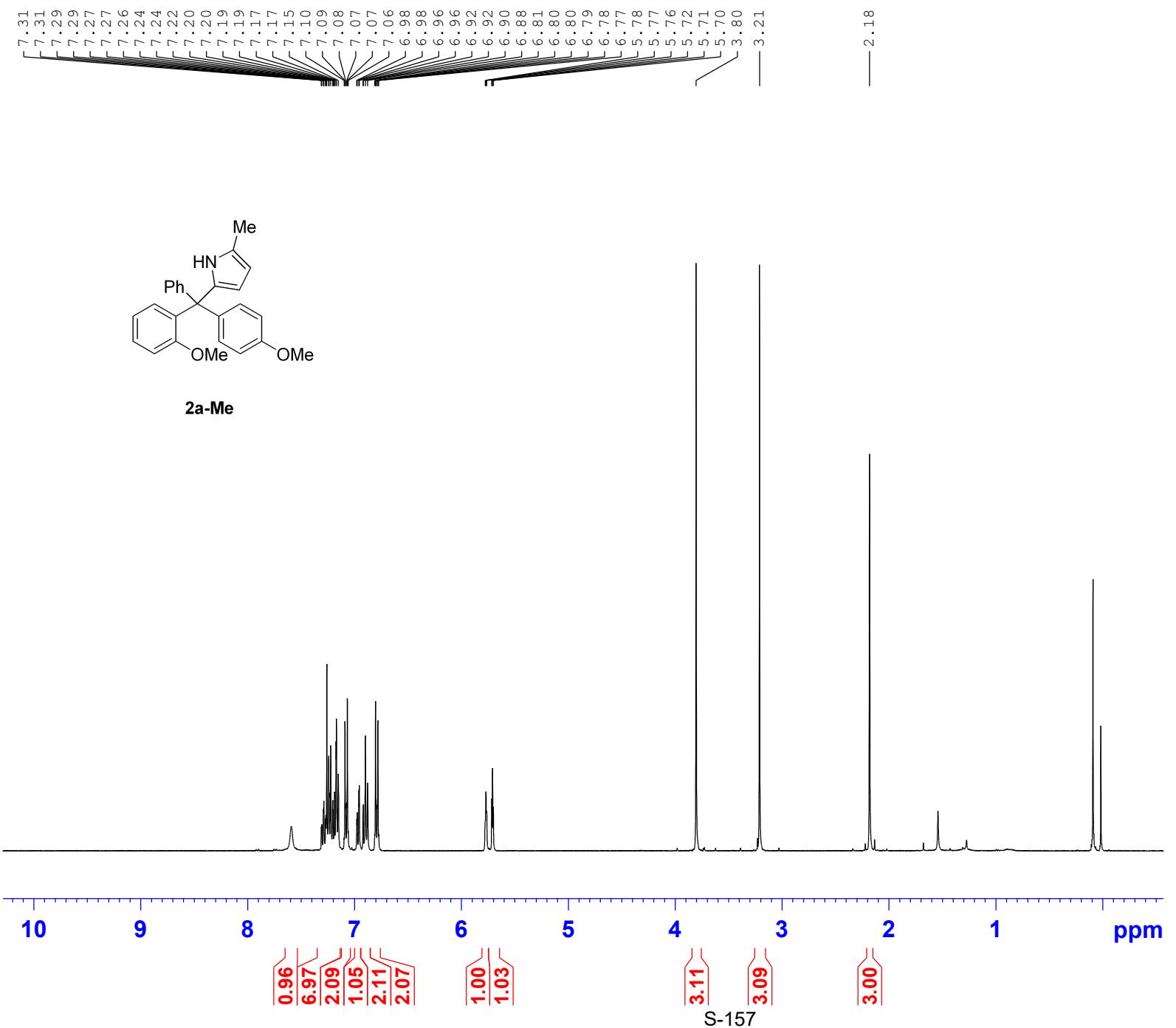




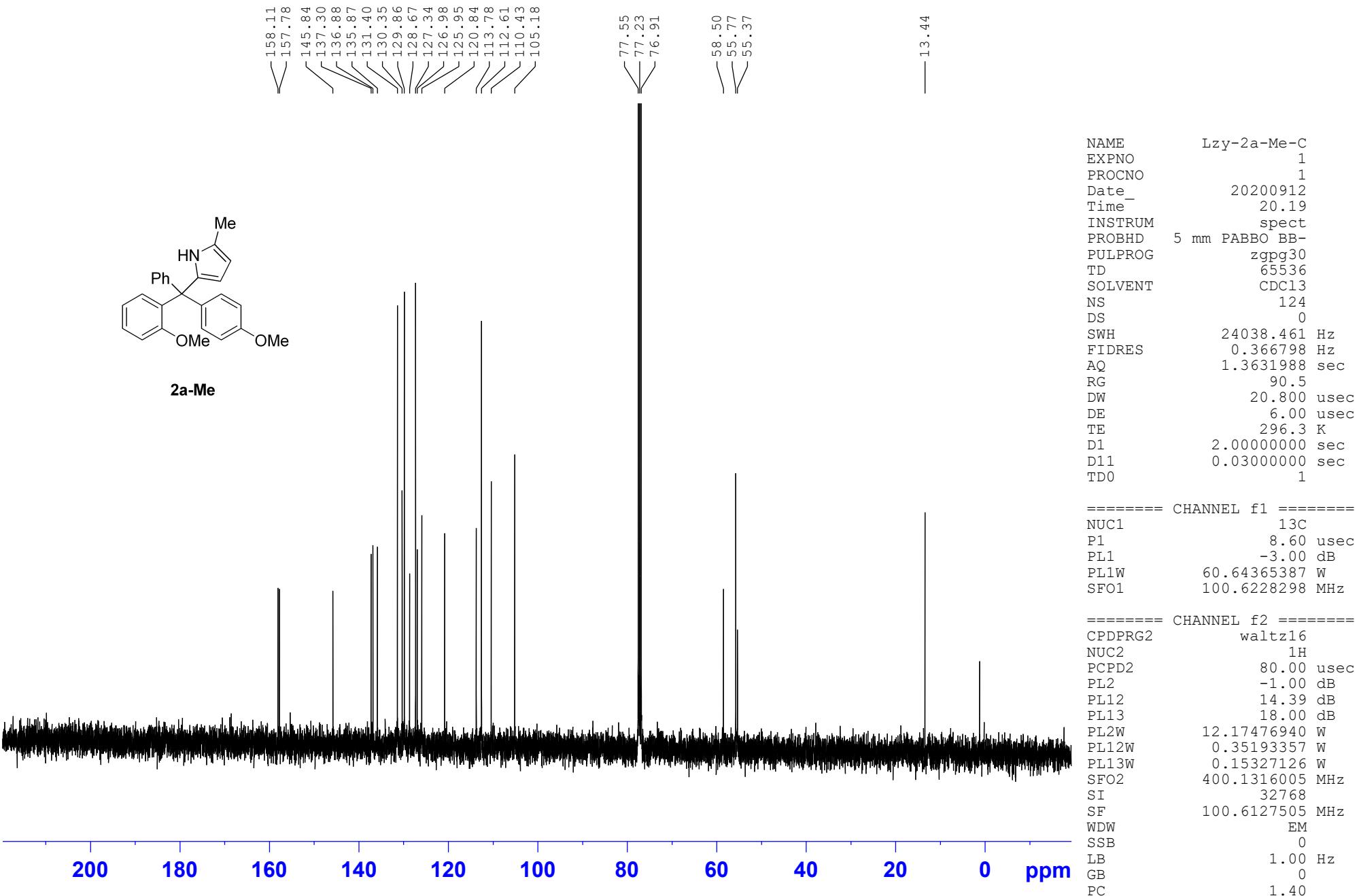


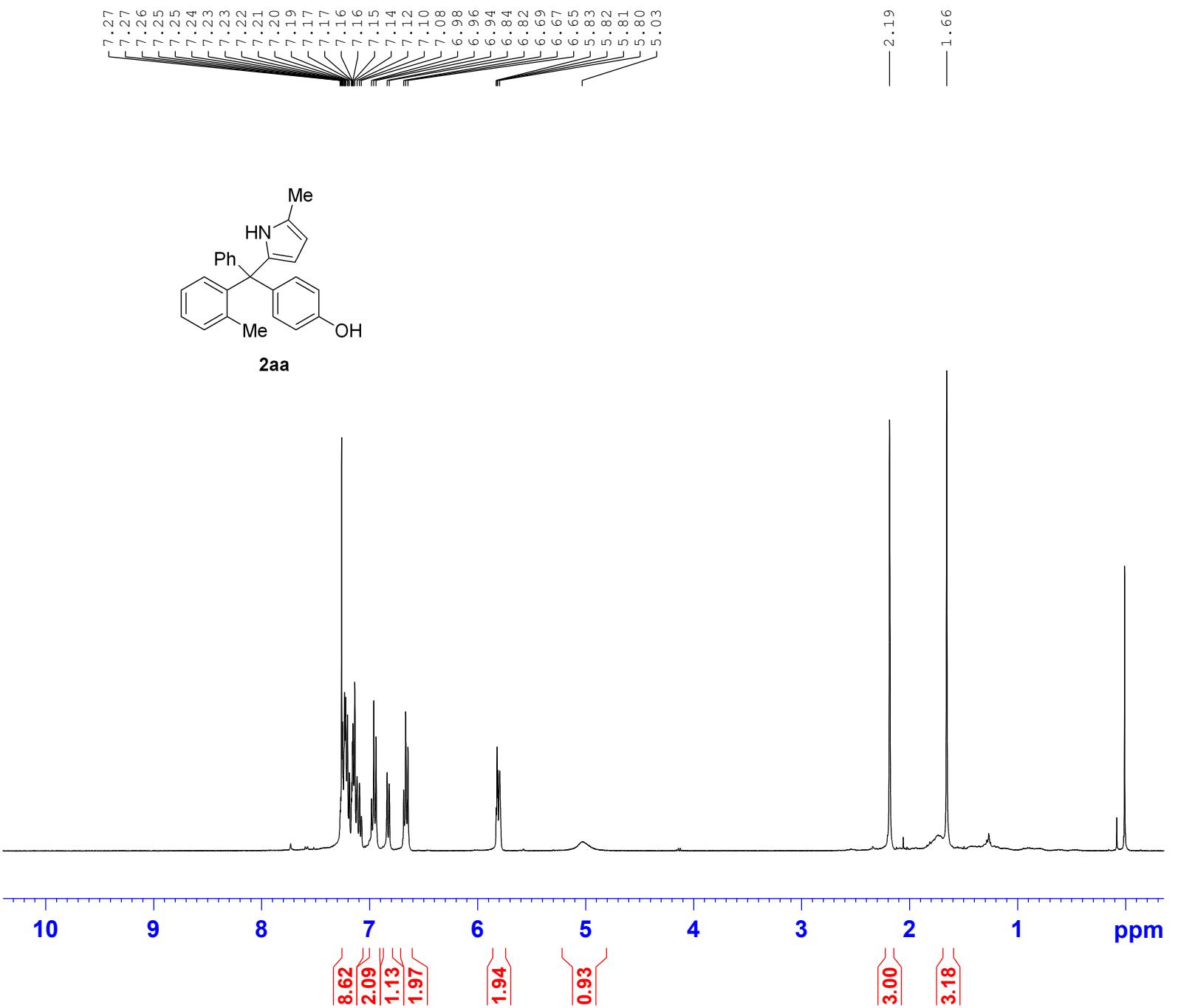


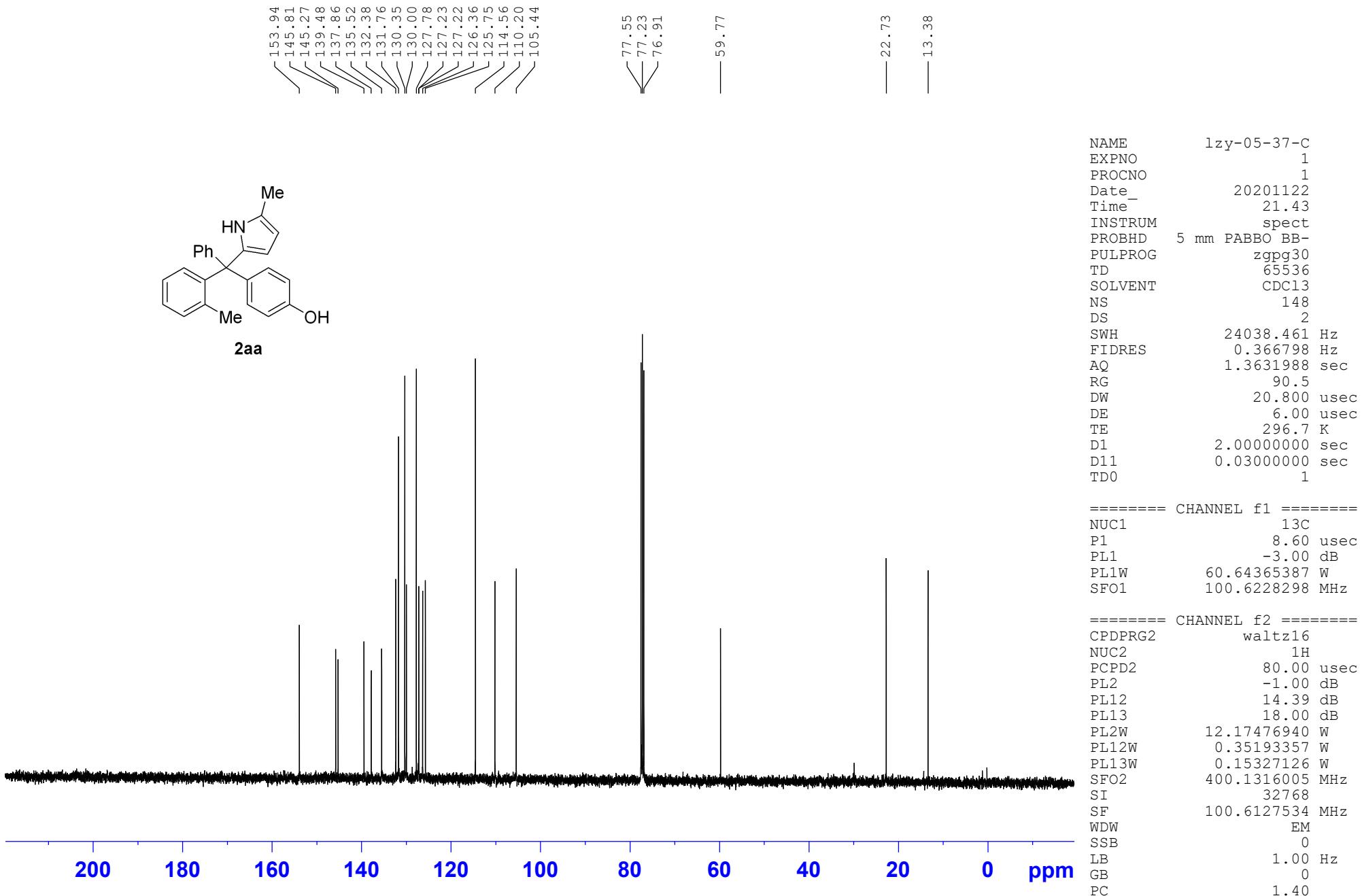


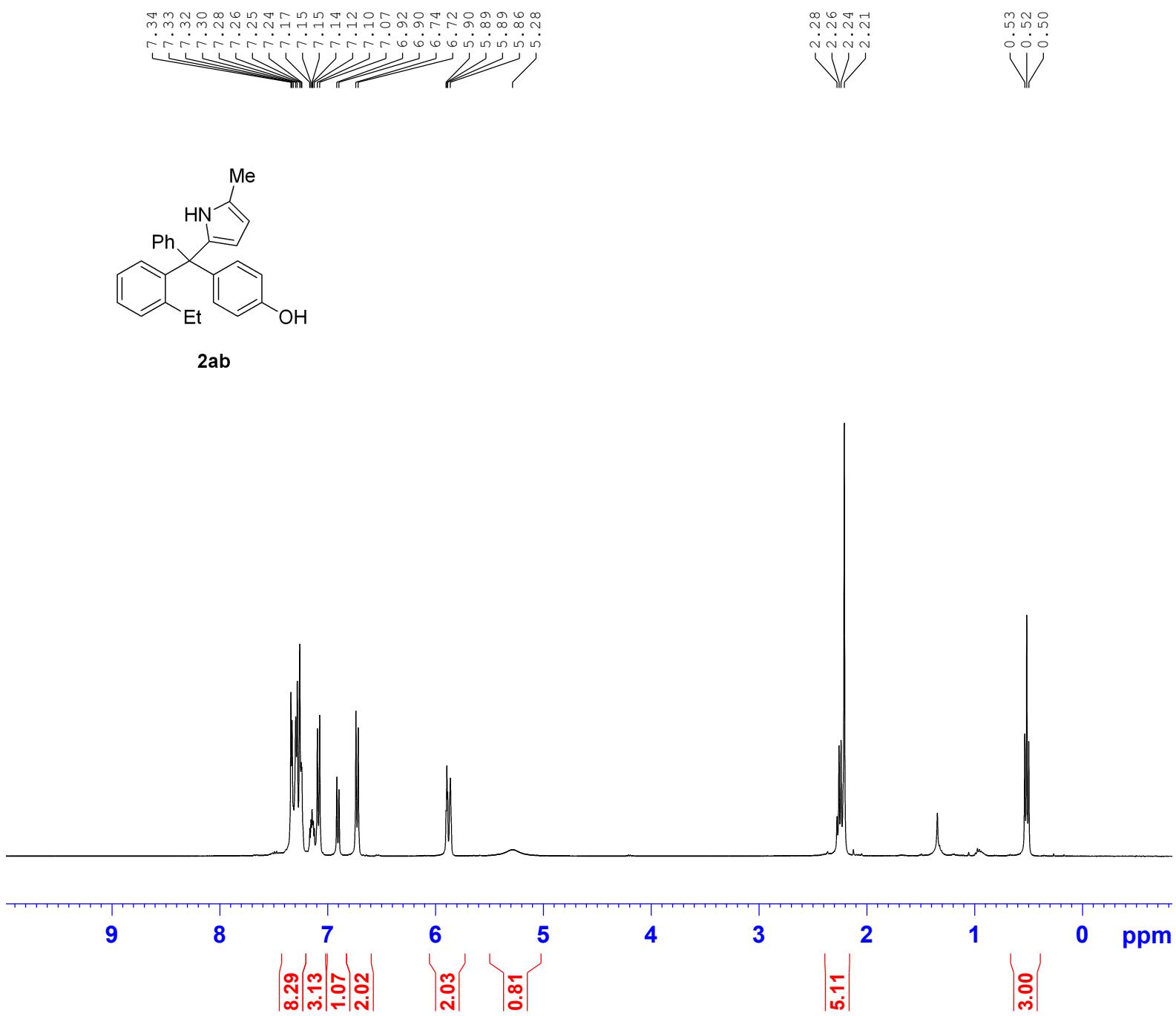


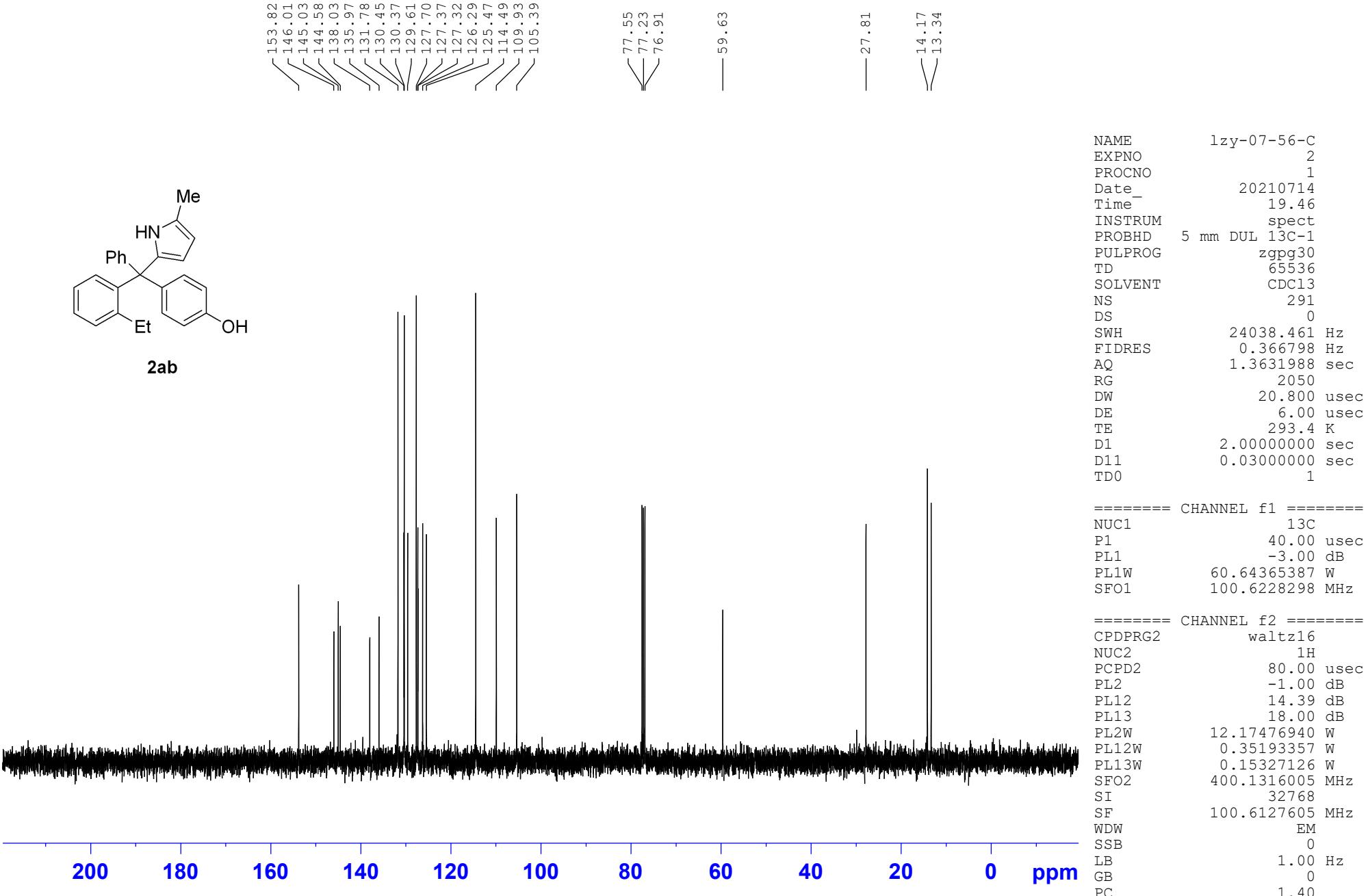
2a-Me

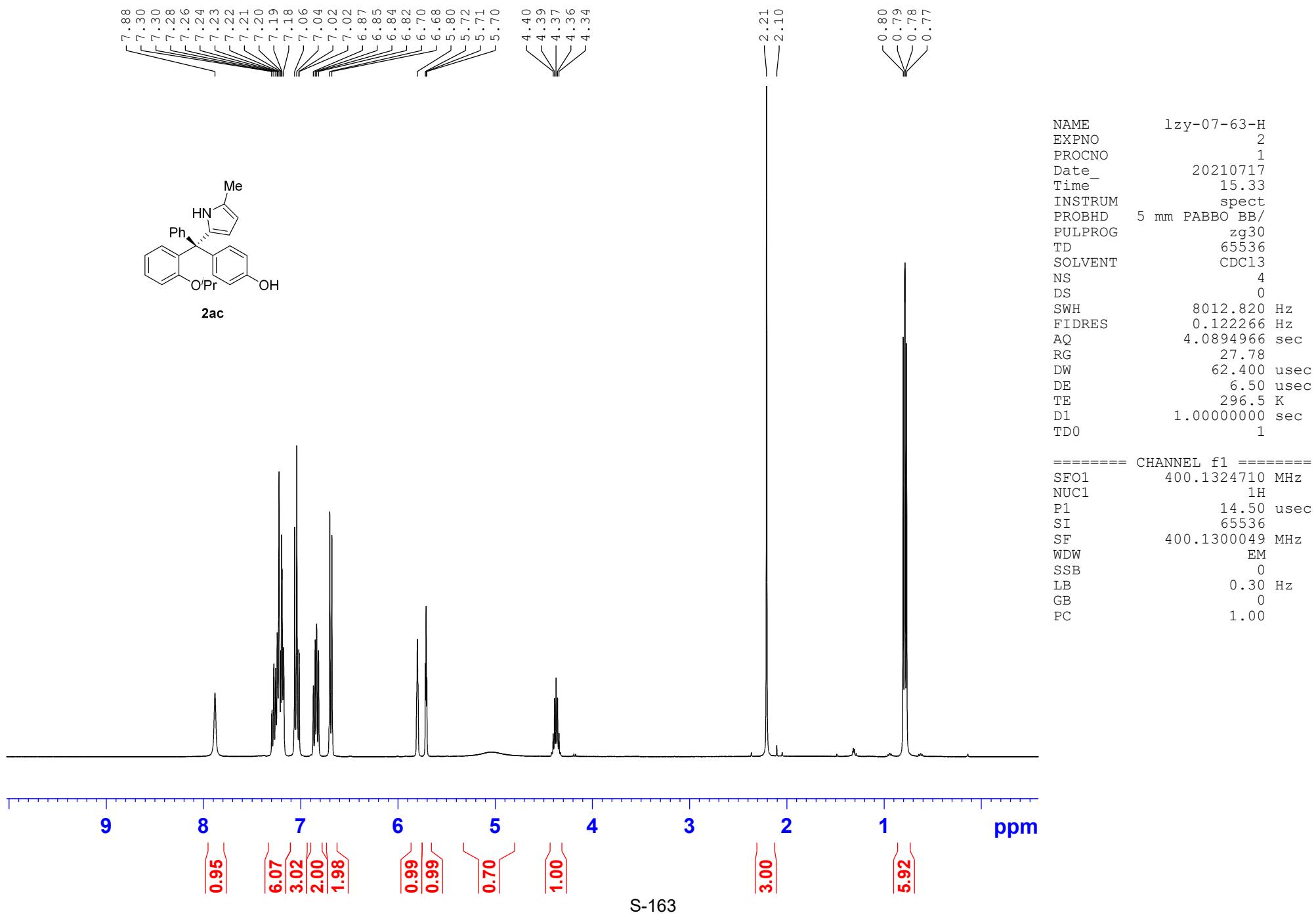


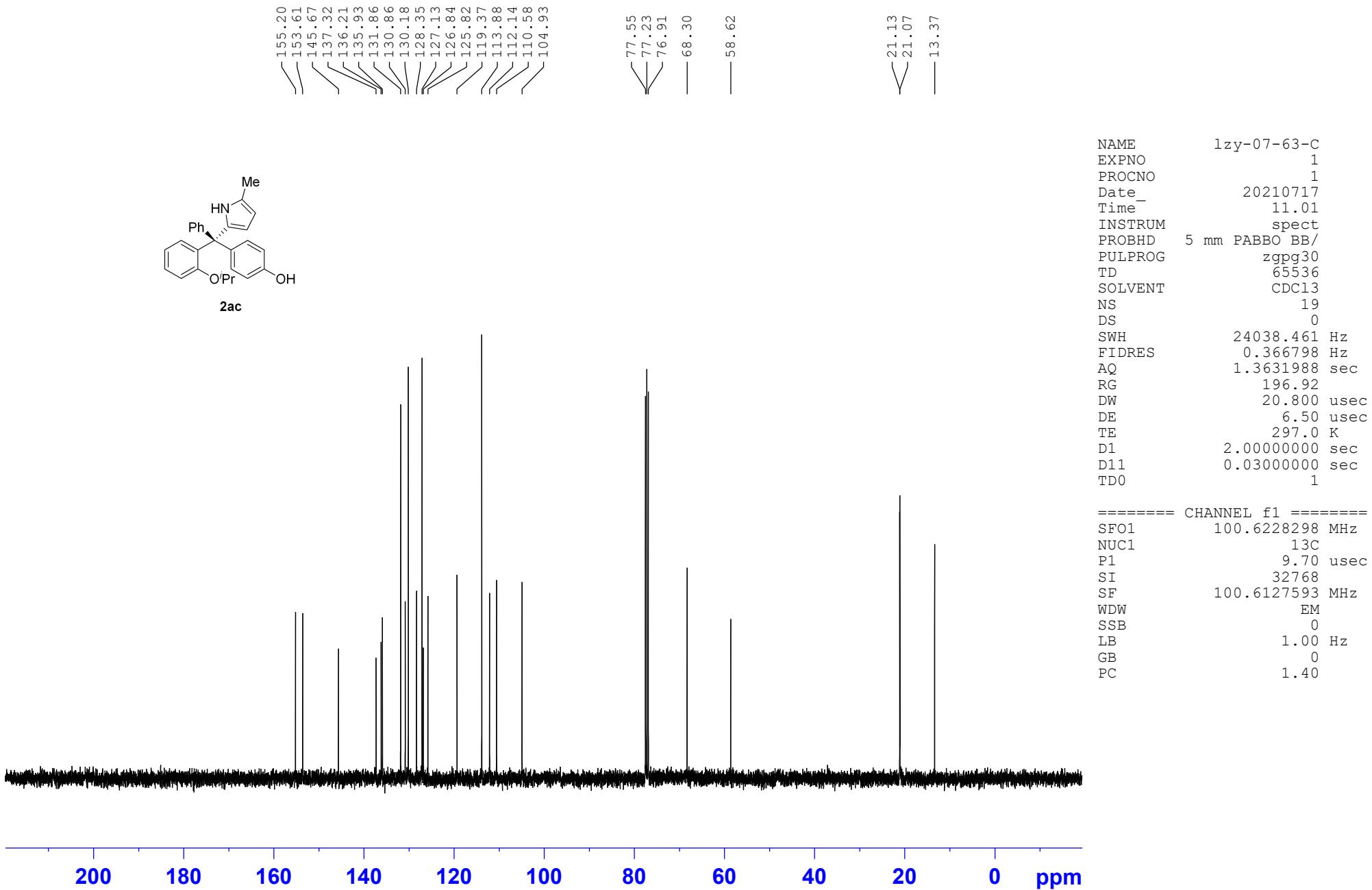


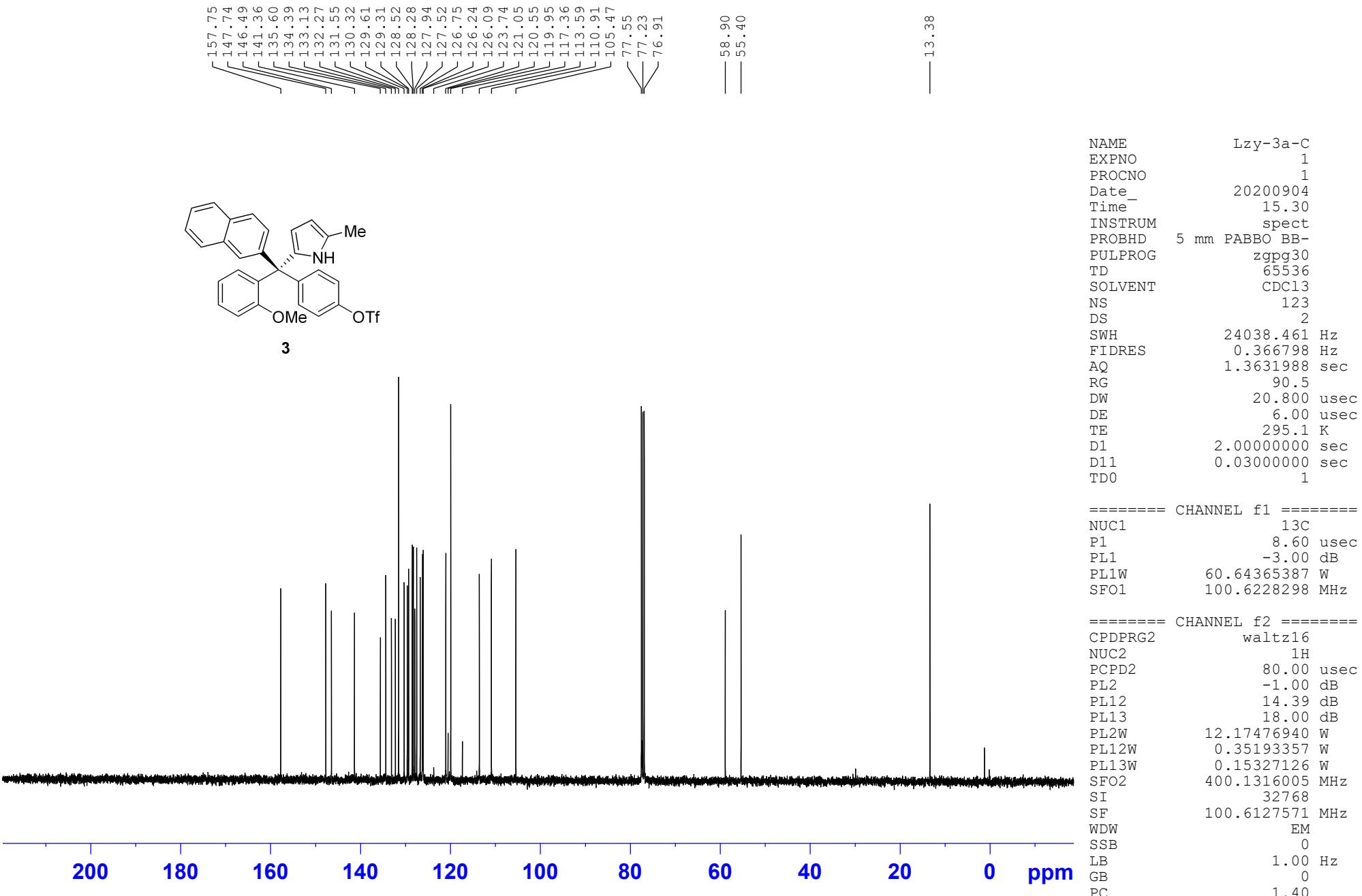


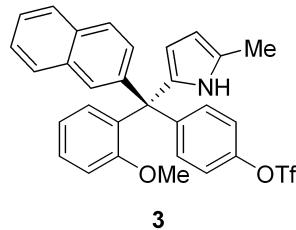








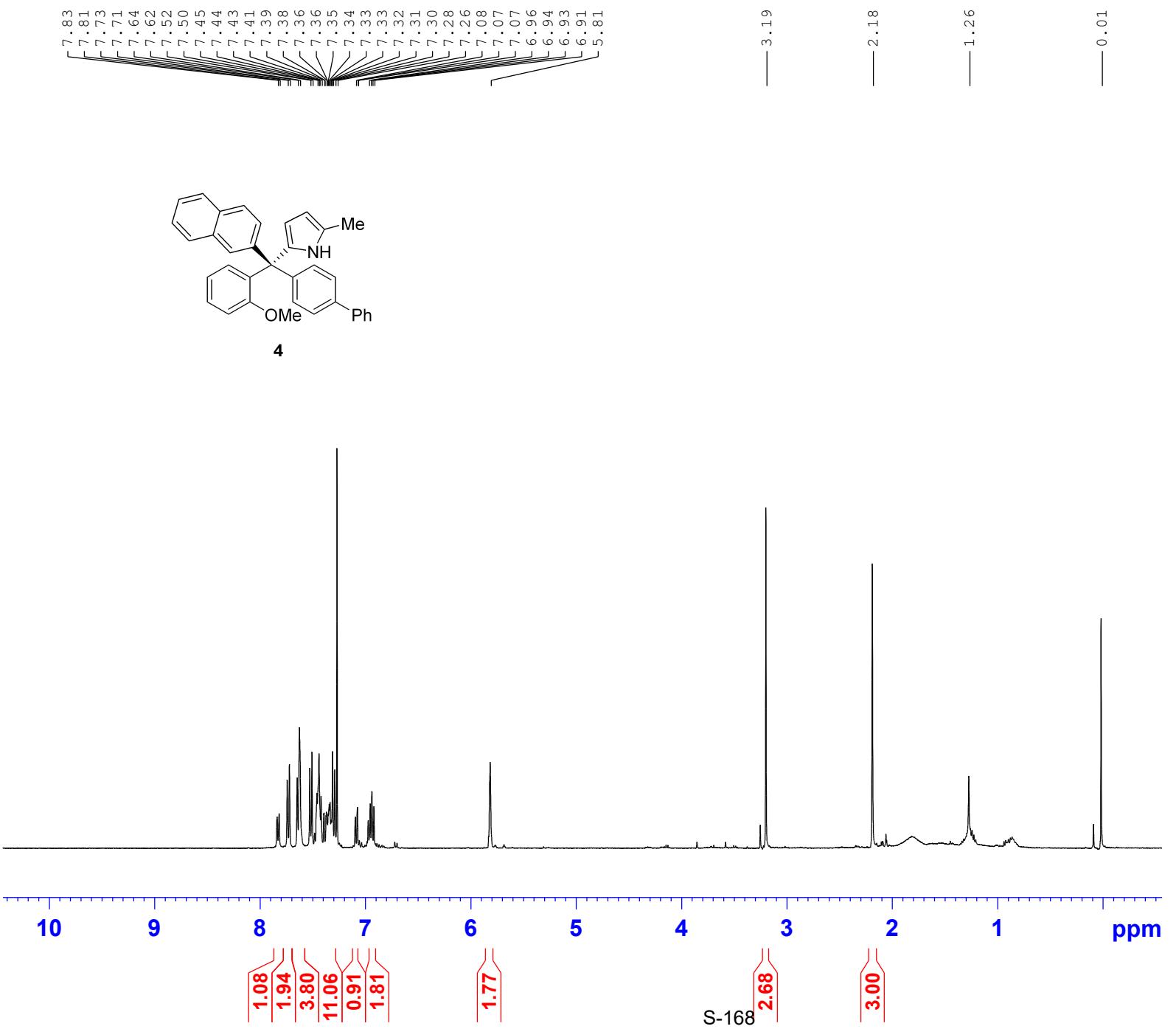




-72.90

NAME	Lzy-3a-F
EXPNO	1
PROCNO	1
Date	20200304
Time	17.36
INSTRUM	spect
PROBHD	5 mm PABBO BB-
PULPROG	zgflqn
TD	131072
SOLVENT	CDCl ₃
NS	6
DS	4
SWH	89285.711 Hz
FIDRES	0.681196 Hz
AQ	0.7340532 sec
RG	1820
DW	5.600 usec
DE	6.00 usec
TE	294.7 K
D1	1.00000000 sec
TD0	1
===== CHANNEL f1 =====	
NUC1	19F
P1	19.50 usec
PL1	-4.00 dB
PL1W	16.97275162 W
SFO1	376.4607164 MHz
SI	65536
SF	376.4983660 MHz
WDW	EM
SSB	5
LB	0.30 Hz
GB	0.05
PC	1.00

0 -20 -40 -60 -80 -100 -120 -140 -160 -180 -200 ppm



```

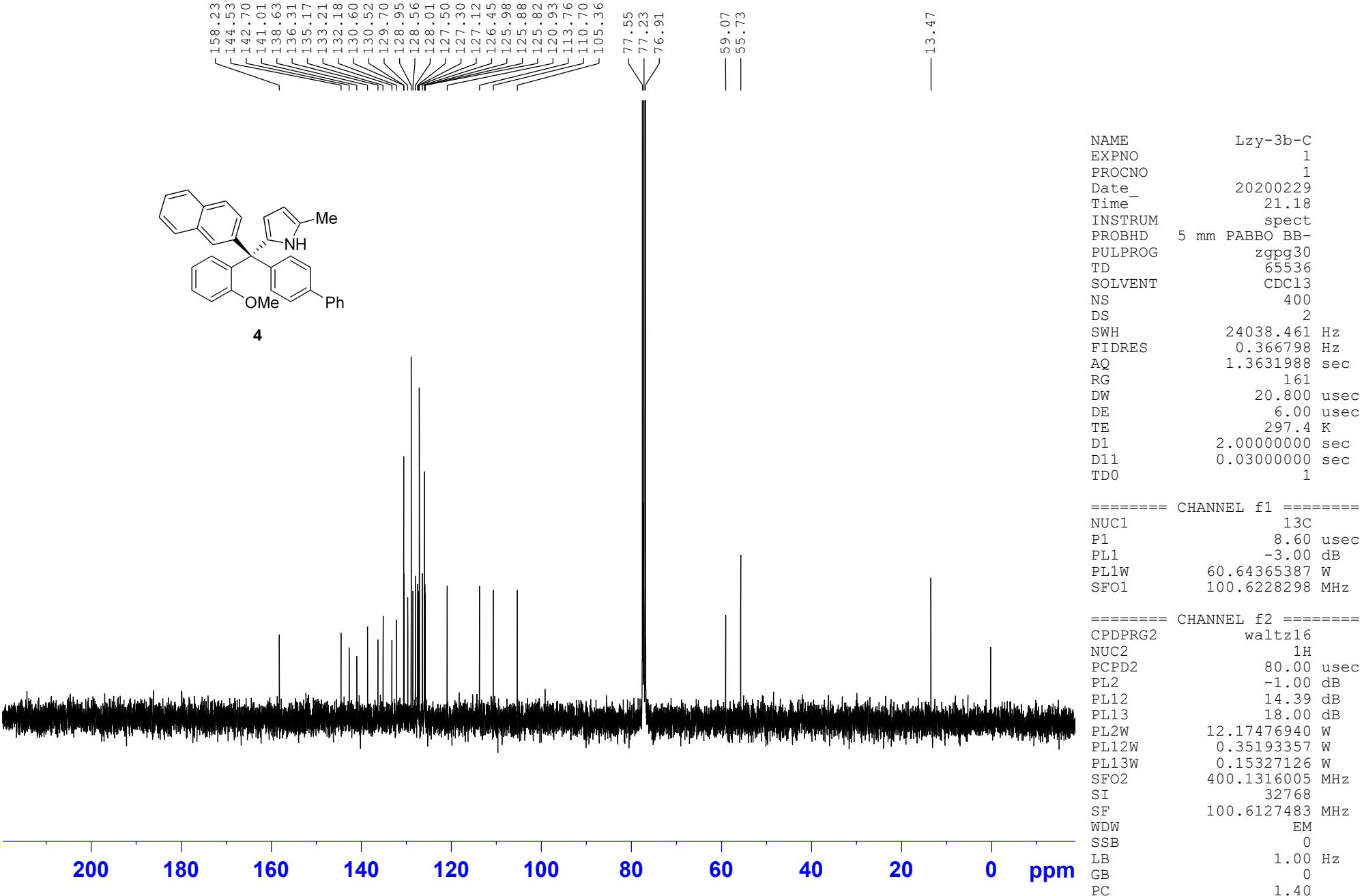
NAME          Lzy-3b
EXPNO         1
PROCNO        1
Date_ 20200229
Time_ 21.12
INSTRUM      spect
PROBHD 5 mm PABBO BB-
PULPROG zg30
TD        65536
SOLVENT       CDC13
NS           6
DS            2
SWH        8223.685 Hz
FIDRES      0.125483 Hz
AQ        3.9846387 sec
RG           406
DW        60.800 usec
DE          6.00 usec
TE        296.9 K
D1        1.00000000 sec
TD0             1

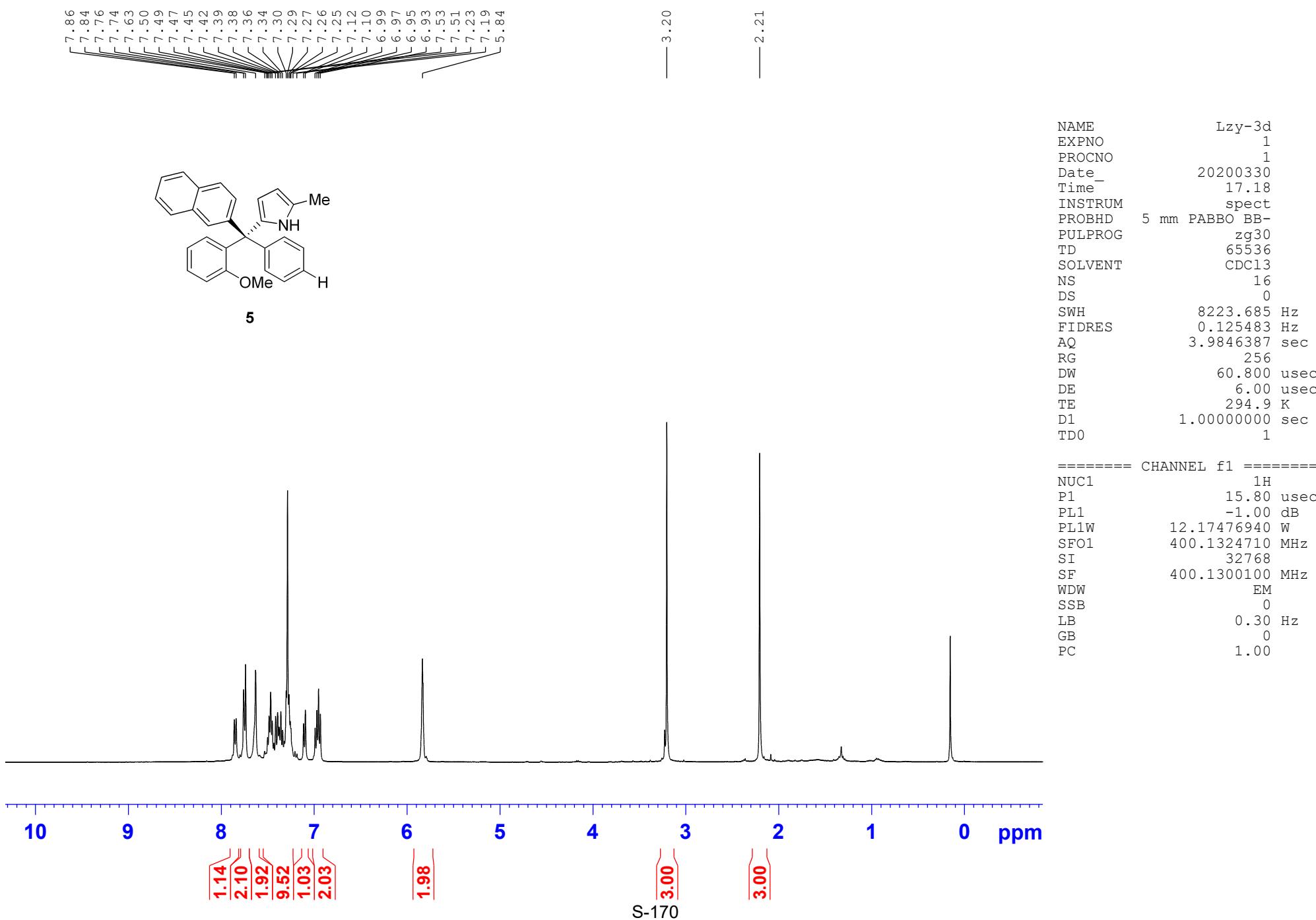
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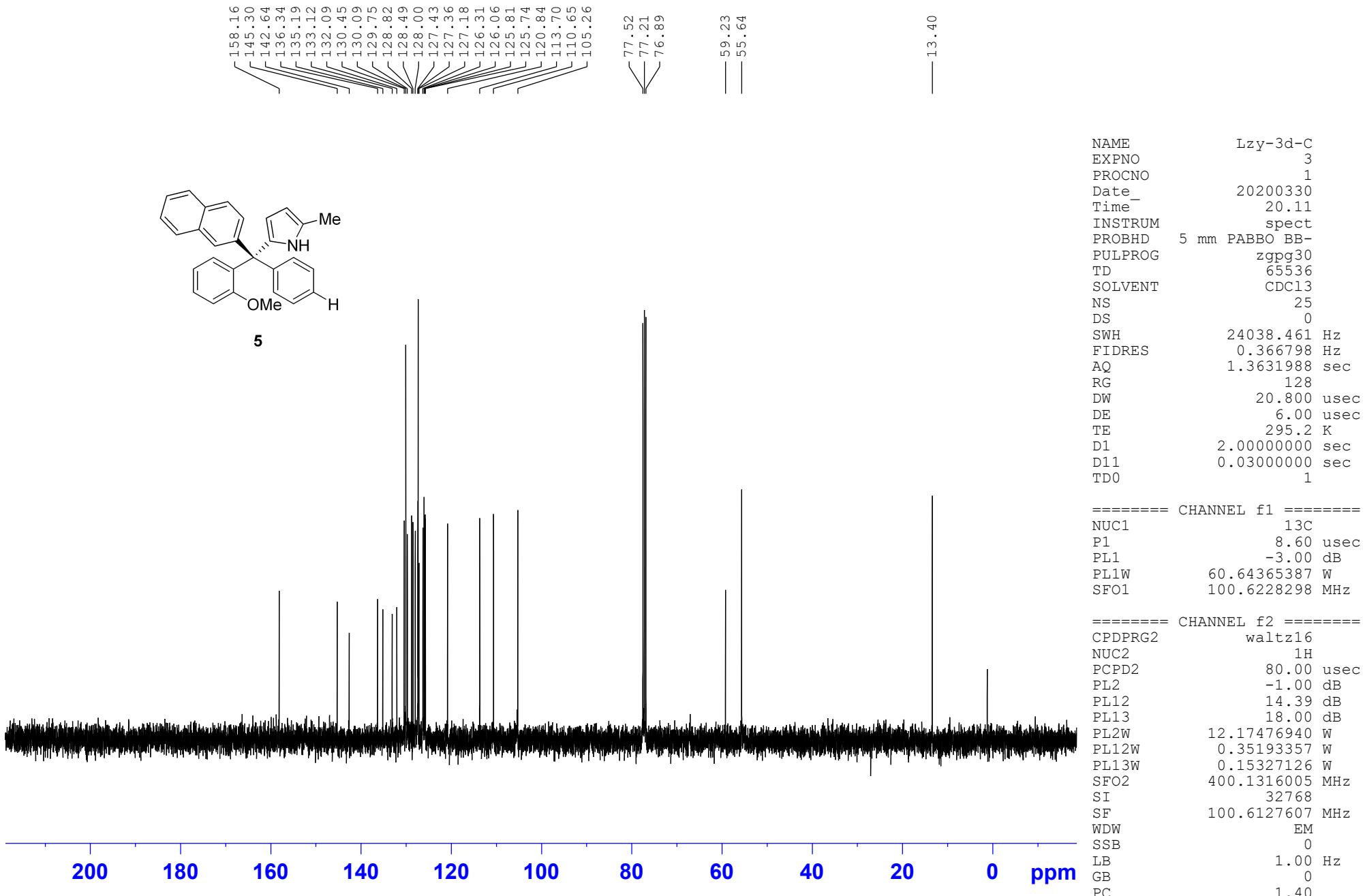
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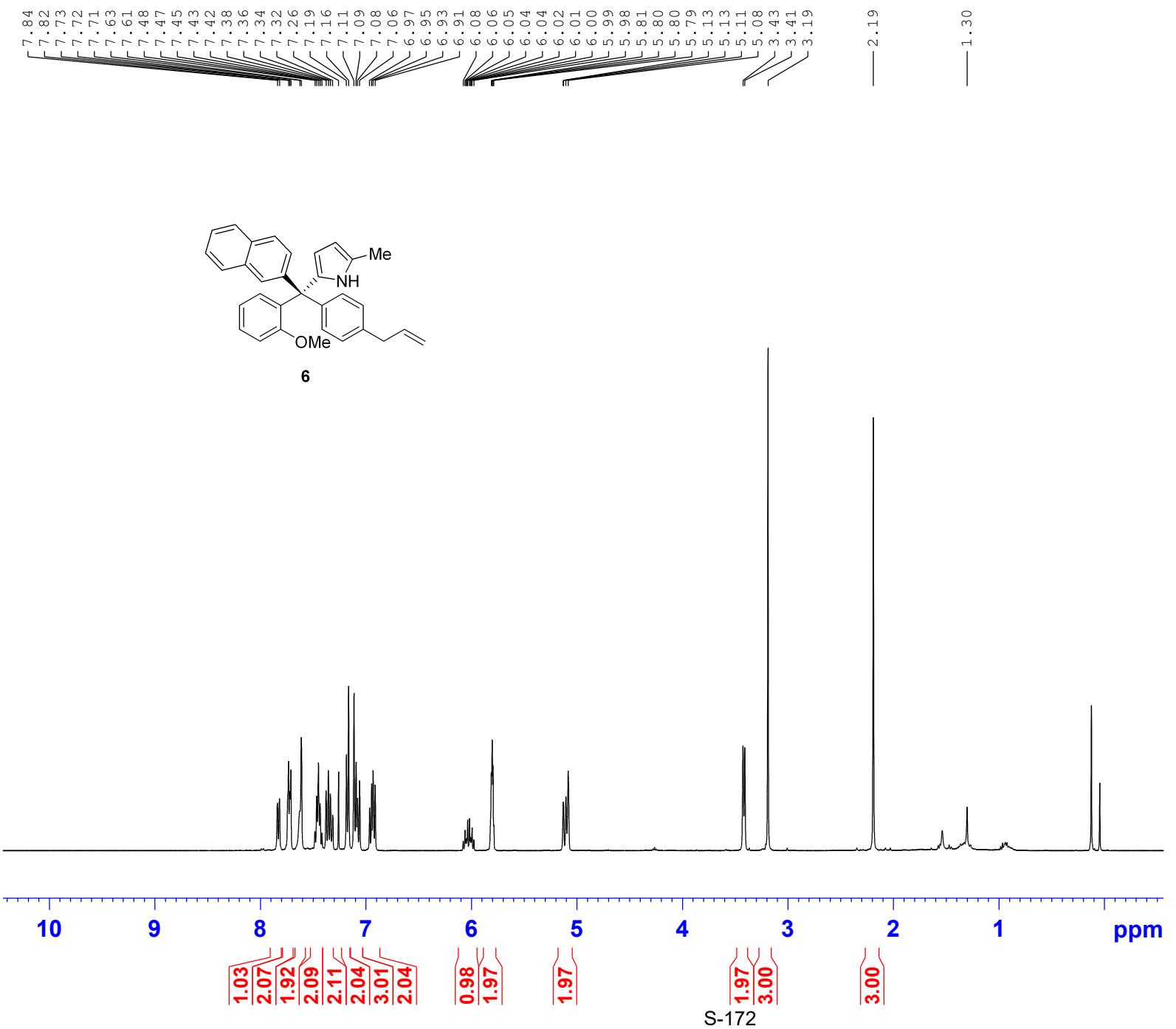
===== CHANNEL f1 =====
NUC1           1H
P1        15.80 usec
PL1        -1.00 dB
PL1W      12.17476940 W
SFO1      400.1324710 MHz
SI           32768
SF      400.1300054 MHz
WDW           EM
SSB            0
LB        0.30 Hz
GB            0
PC        1.00

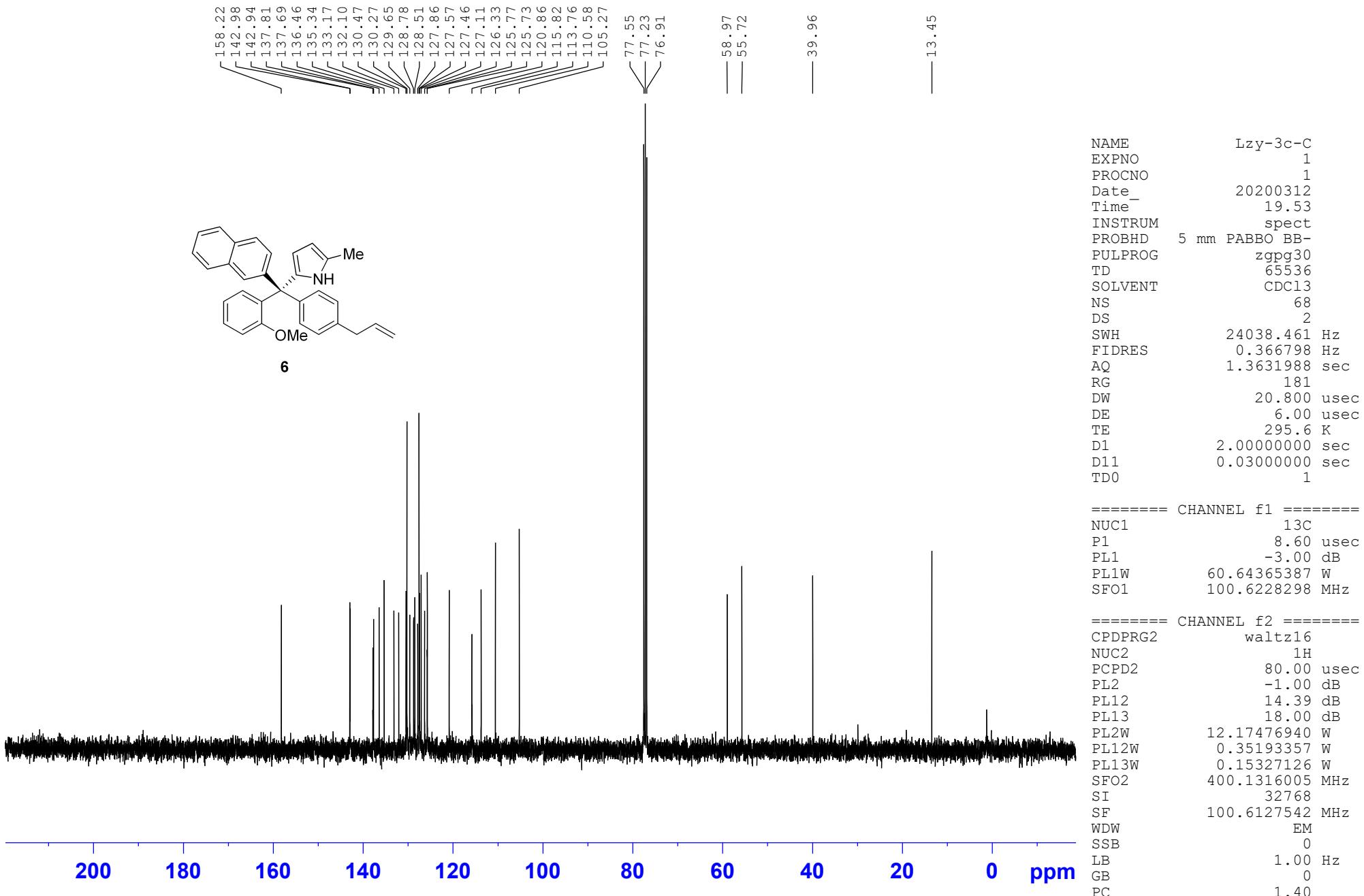
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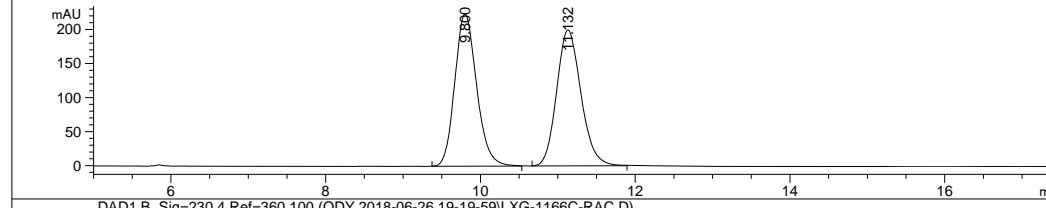


Sample Name:

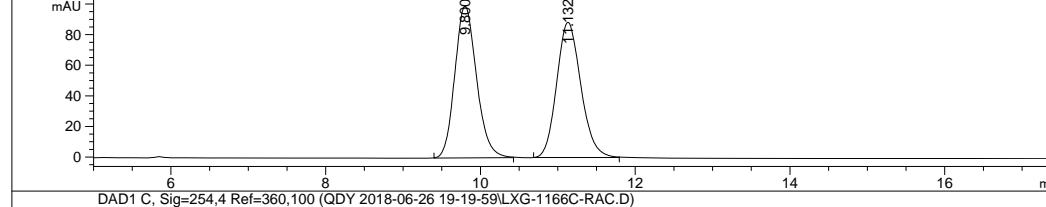
=====
 Acq. Operator : Seq. Line : 5
 Acq. Instrument : Instrument 1 Location : Vial 62
 Injection Date : 6/26/2018 8:47:00 PM Inj : 1
 Inj Volume : 5.000 μ l
 Different Inj Volume from Sequence ! Actual Inj Volume : 2.000 μ l
 Acq. Method : C:\CHEM32\1\DATA\QDY 2018-06-26 19-19-59\IC-03-20.M
 Last changed : 6/26/2018 8:46:07 PM
 (modified after loading)
 Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
 Last changed : 11/24/2020 3:16:16 PM
 (modified after loading)

Additional Info : Peak(s) manually integrated

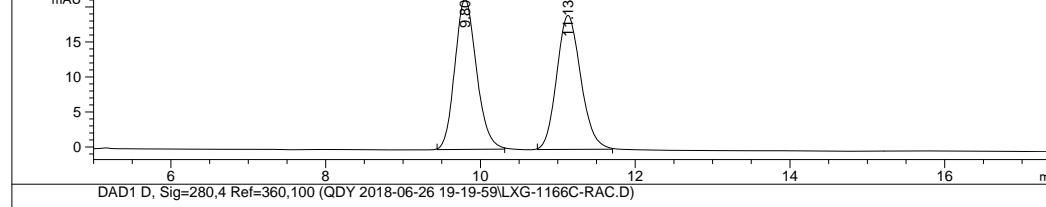
DAD1 A, Sig=210,4 Ref=360,100 (QDY 2018-06-26 19-19-59\LXG-1166C-RAC.D)



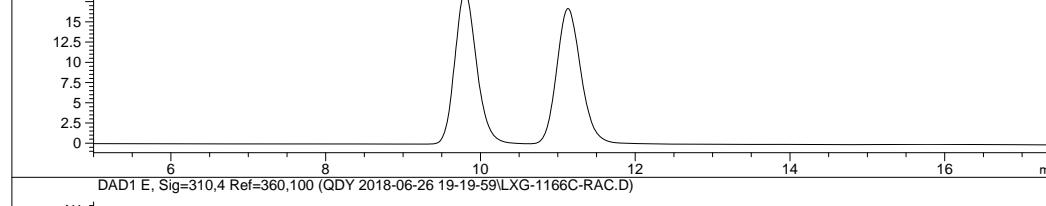
DAD1 B, Sig=230,4 Ref=360,100 (QDY 2018-06-26 19-19-59\LXG-1166C-RAC.D)



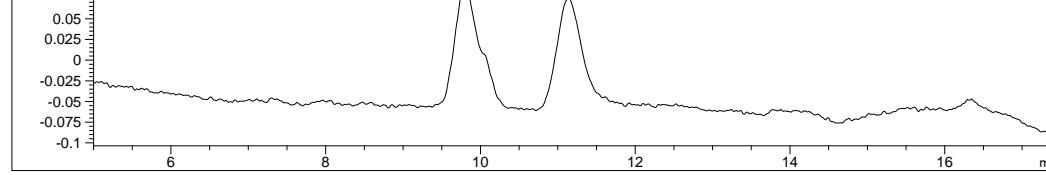
DAD1 C, Sig=254,4 Ref=360,100 (QDY 2018-06-26 19-19-59\LXG-1166C-RAC.D)



DAD1 D, Sig=280,4 Ref=360,100 (QDY 2018-06-26 19-19-59\LXG-1166C-RAC.D)



DAD1 E, Sig=310,4 Ref=360,100 (QDY 2018-06-26 19-19-59\LXG-1166C-RAC.D)



Data File C:\CHEM32\1\DATA\QDY 2018-06-26 19-19-59\LXG-1166C-RAC.D
Sample Name:

===== Area Percent Report =====

Sorted By : Signal
 Multiplier : 1.0000
 Dilution : 1.0000
 Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.800	BB	0.3023	4374.23193	223.54536	50.1048
2	11.132	BB	0.3385	4355.92871	199.66124	49.8952

Totals : 8730.16064 423.20660

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.800	BB	0.3012	1931.31421	99.15396	50.2023
2	11.132	BB	0.3369	1915.74731	88.36066	49.7977

Totals : 3847.06152 187.51462

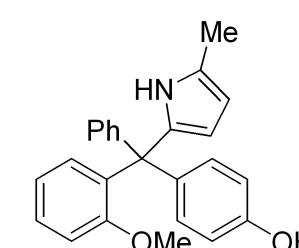
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.800	BB	0.3003	416.10309	21.45372	50.1396
2	11.132	BB	0.3341	413.78525	19.14697	49.8604

Totals : 829.88834 40.60069

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

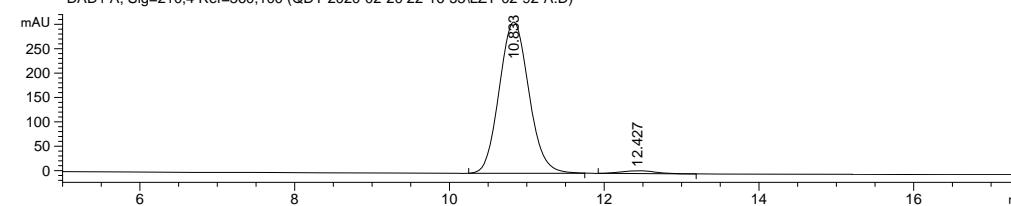
Signal 5: DAD1 E, Sig=310,4 Ref=360,100



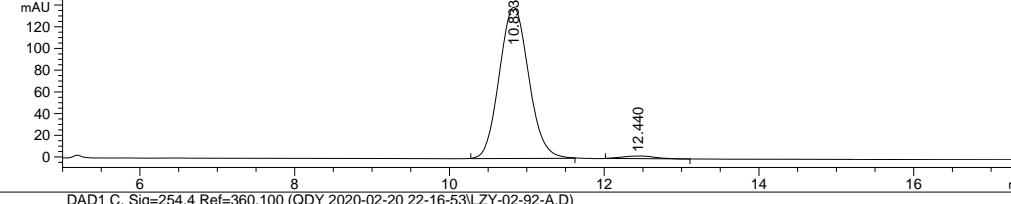
Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 2
Acq. Instrument : Instrument 1             Location : Vial 81
Injection Date : 2/20/2020 10:30:17 PM       Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-02-20 22-16-53\IC-03-20.M
Last changed : 2/20/2020 10:29:24 PM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 3:16:16 PM        (modified after loading)
Additional Info : Peak(s) manually integrated
```

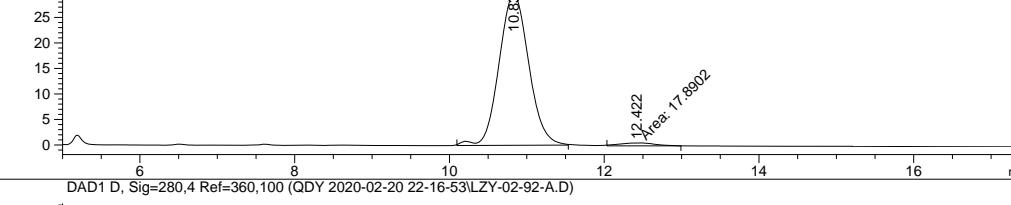
DAD1 A, Sig=210,4 Ref=360,100 (QDY 2020-02-20 22-16-53\LZY-02-92-A.D)



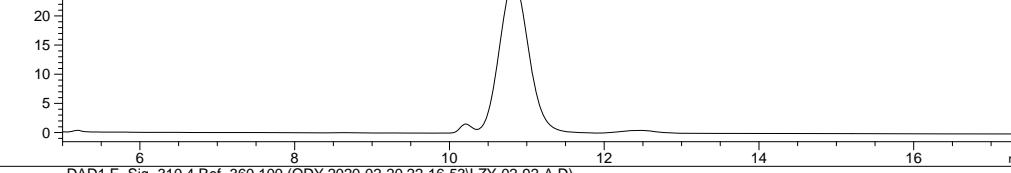
DAD1 B, Sig=230,4 Ref=360,100 (QDY 2020-02-20 22-16-53\LZY-02-92-A.D)



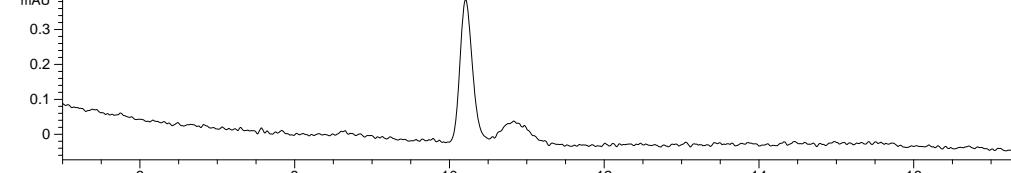
DAD1 C, Sig=254,4 Ref=360,100 (QDY 2020-02-20 22-16-53\LZY-02-92-A.D)



DAD1 D, Sig=280,4 Ref=360,100 (QDY 2020-02-20 22-16-53\LZY-02-92-A.D)



DAD1 E, Sig=310,4 Ref=360,100 (QDY 2020-02-20 22-16-53\LZY-02-92-A.D)



Sample Name:

===== Area Percent Report =====

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.833	BB	0.4149	8212.09082	310.16901	97.9796
2	12.427	BB	0.4284	169.33453	5.87353	2.0204

Totals : 8381.42535 316.04254

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.833	BB	0.4110	3668.44141	139.42035	98.1103
2	12.440	BB	0.4167	70.65849	2.55482	1.8897

Totals : 3739.09990 141.97517

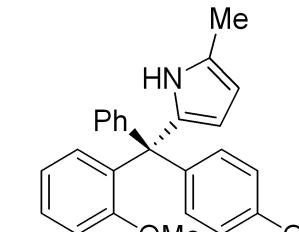
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.833	BB	0.4135	785.15674	29.59425	97.7722
2	12.422	MM	0.5004	17.89020	5.95917e-1	2.2278

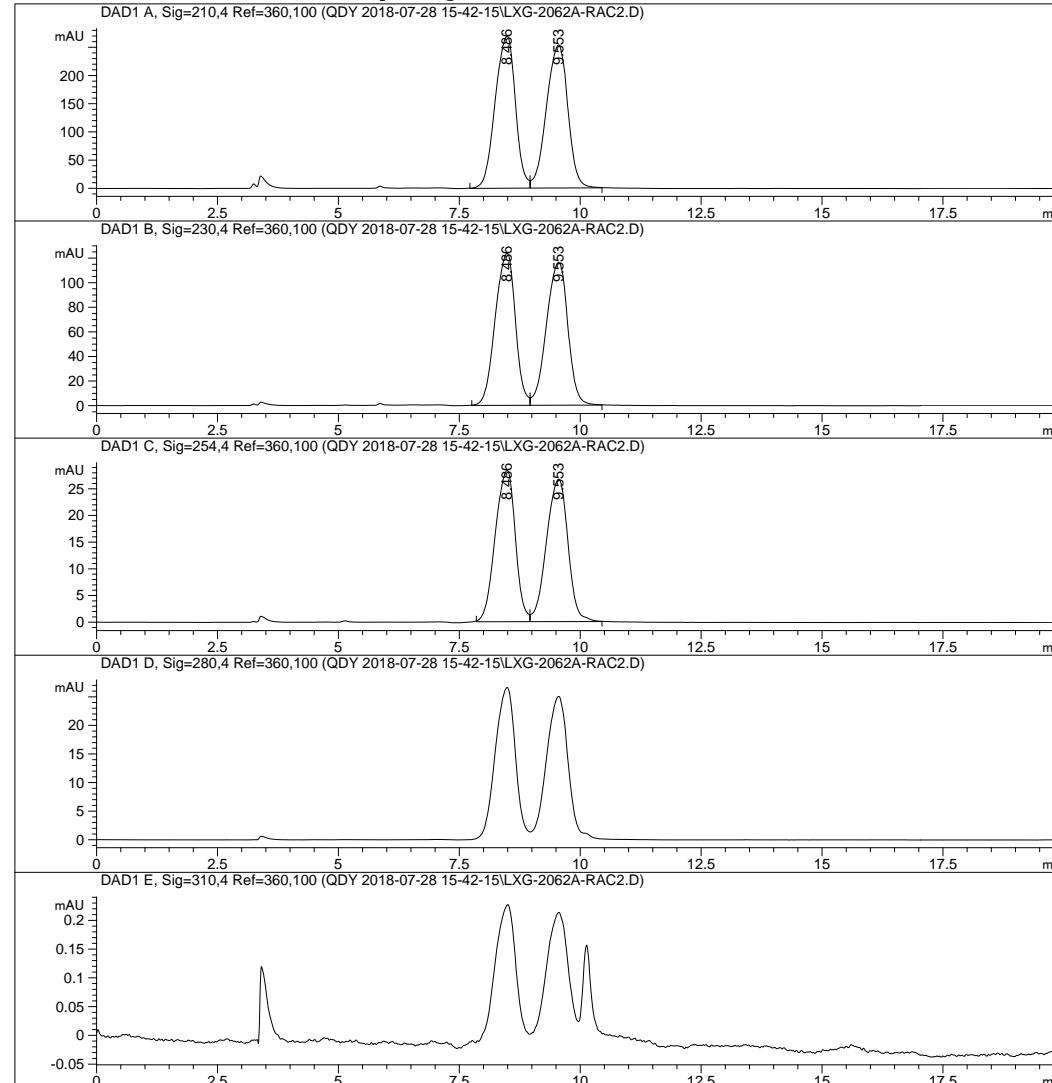
Totals : 803.04694 30.19017

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



```
=====
Acq. Operator :                               Seq. Line :   6
Acq. Instrument : Instrument 1             Location : Vial 67
Injection Date : 7/28/2018 5:12:39 PM        Inj : 1
                                                Inj Volume : 5.000 µl
Acq. Method :  C:\CHEM32\1\DATA\QDY 2018-07-28 15-42-15\IC-03-20.M
Last changed : 7/28/2018 5:11:47 PM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 10/29/2018 7:27:37 PM       (modified after loading)
Additional Info : Peak(s) manually integrated
```



```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.486	BV	0.4663	7820.67480	269.73096	50.0394
2	9.553	VB	0.4933	7808.35107	253.52272	49.9606
Totals :					1.56290e4	523.25368

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

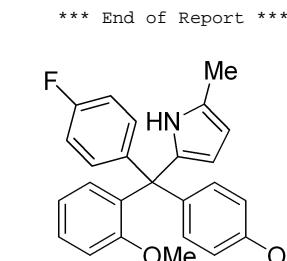
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.486	BV	0.4654	3579.20850	123.75388	50.0570
2	9.553	VB	0.4926	3571.05396	116.18679	49.9430
Totals :					7150.26245	239.94067

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.486	BV	0.4638	816.48468	28.36868	49.8220
2	9.553	VB	0.4936	822.31885	26.67982	50.1780
Totals :					1638.80353	55.04851

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

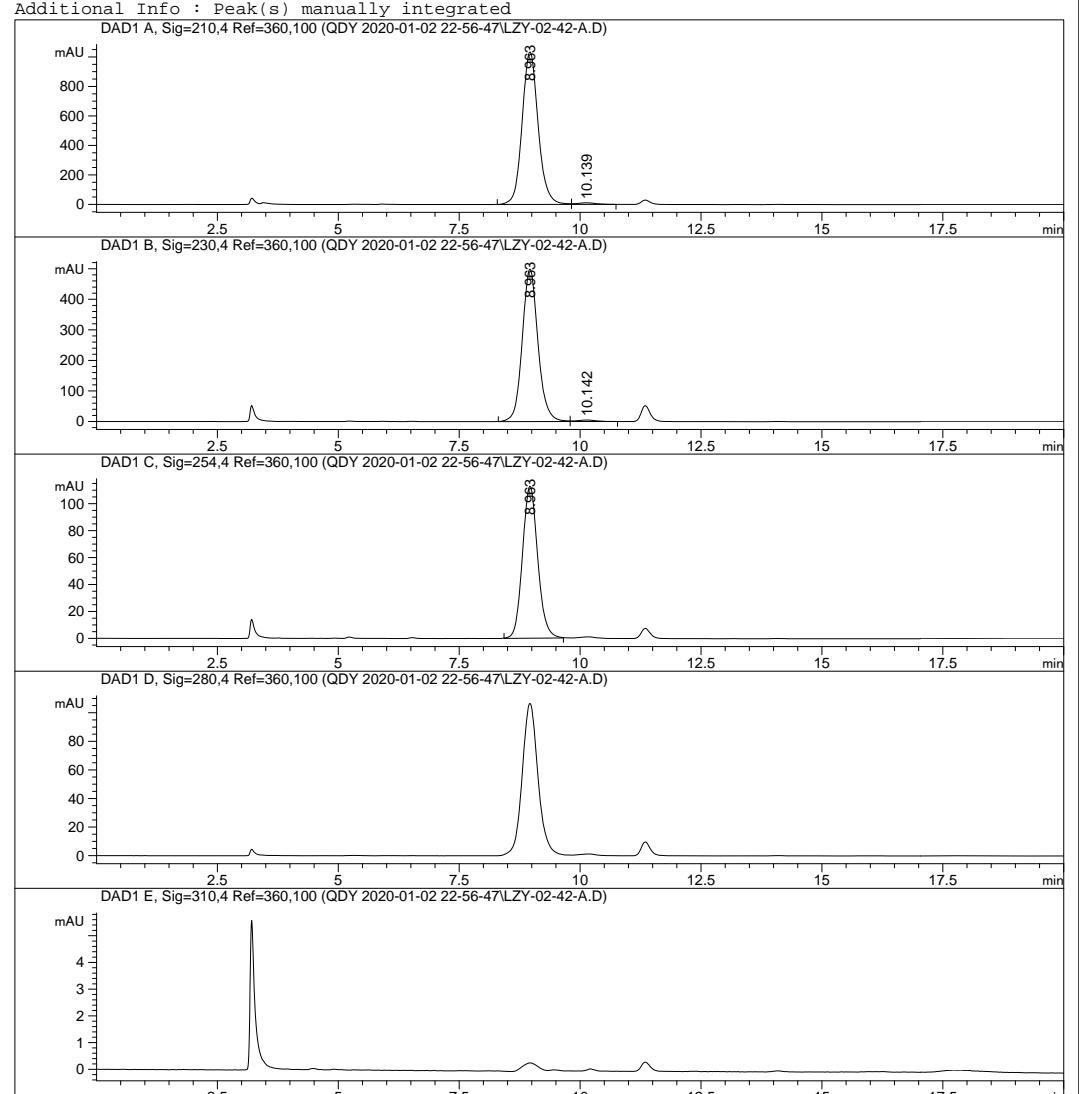
Signal 5: DAD1 E, Sig=310,4 Ref=360,100



2b-Rac

Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 17
Acq. Instrument : Instrument 1             Location : Vial 91
Injection Date : 1/3/2020 6:37:36 AM        Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-01-02 22-56-47\IC-03-20.M
Last changed : 1/3/2020 6:36:43 AM
                                                (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-10-10.M
Last changed : 9/9/2020 9:55:03 AM
                                                (modified after loading)
Additional Info : Peak(s) manually integrated
```



Data File C:\CHEM32\1\DATA\QDY 2020-01-02 22-56-47\LZY-02-42-A.D
Sample Name:

```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.963	BB	0.3580	2.39175e4	1033.63550	98.7632
2	10.139	BB	0.4029	299.51294	11.16648	1.2368

Totals : 2.42170e4 1044.80198

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.963	BB	0.3507	1.13165e4	498.77319	98.8202
2	10.142	BB	0.4064	135.10916	4.98238	1.1798

Totals : 1.14516e4 503.75557

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

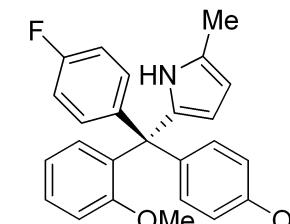
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.963	BB	0.3423	2478.50781	112.81509	100.0000

Totals : 2478.50781 112.81509

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

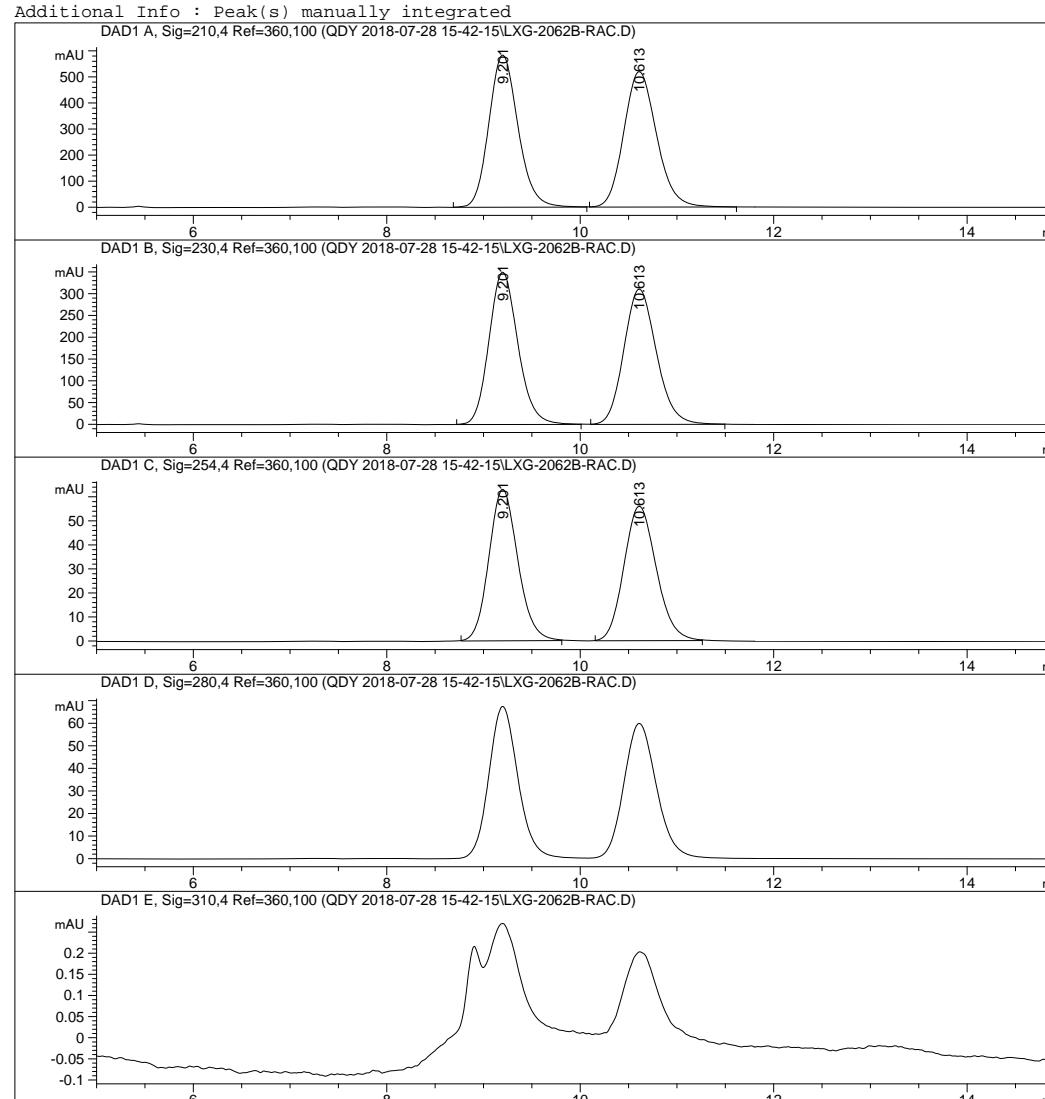
Signal 5: DAD1 E, Sig=310,4 Ref=360,100

===== *** End of Report *** =====

**2b**

Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 3
Acq. Instrument : Instrument 1             Location : Vial 68
Injection Date : 7/28/2018 4:17:04 PM        Inj : 1
                                                Inj Volume : 5.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2018-07-28 15-42-15\IC-05-20.M
Last changed : 6/25/2018 1:35:20 AM
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 3:21:46 PM
                                                (modified after loading)
Additional Info : Peak(s) manually integrated
```



Data File C:\CHEM32\1\DATA\QDY 2018-07-28 15-42-15\LXG-2062B-RAC.D
Sample Name:

```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.201	BB	0.3245	1.22368e4	583.89270	49.9836
2	10.613	BB	0.3668	1.22448e4	519.77472	50.0164

Totals : 2.44816e4 1103.66742

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.201	BB	0.3226	7273.56543	349.71793	50.0562
2	10.613	BB	0.3606	7257.22803	310.55371	49.9438

Totals : 1.45308e4 660.27164

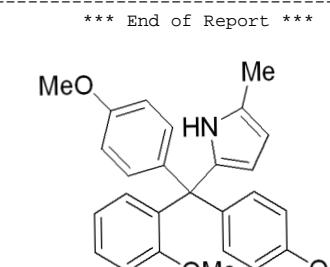
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.201	BB	0.3225	1310.82092	63.05332	50.2282
2	10.613	BB	0.3610	1298.91028	55.91644	49.7718

Totals : 2609.73120 118.96976

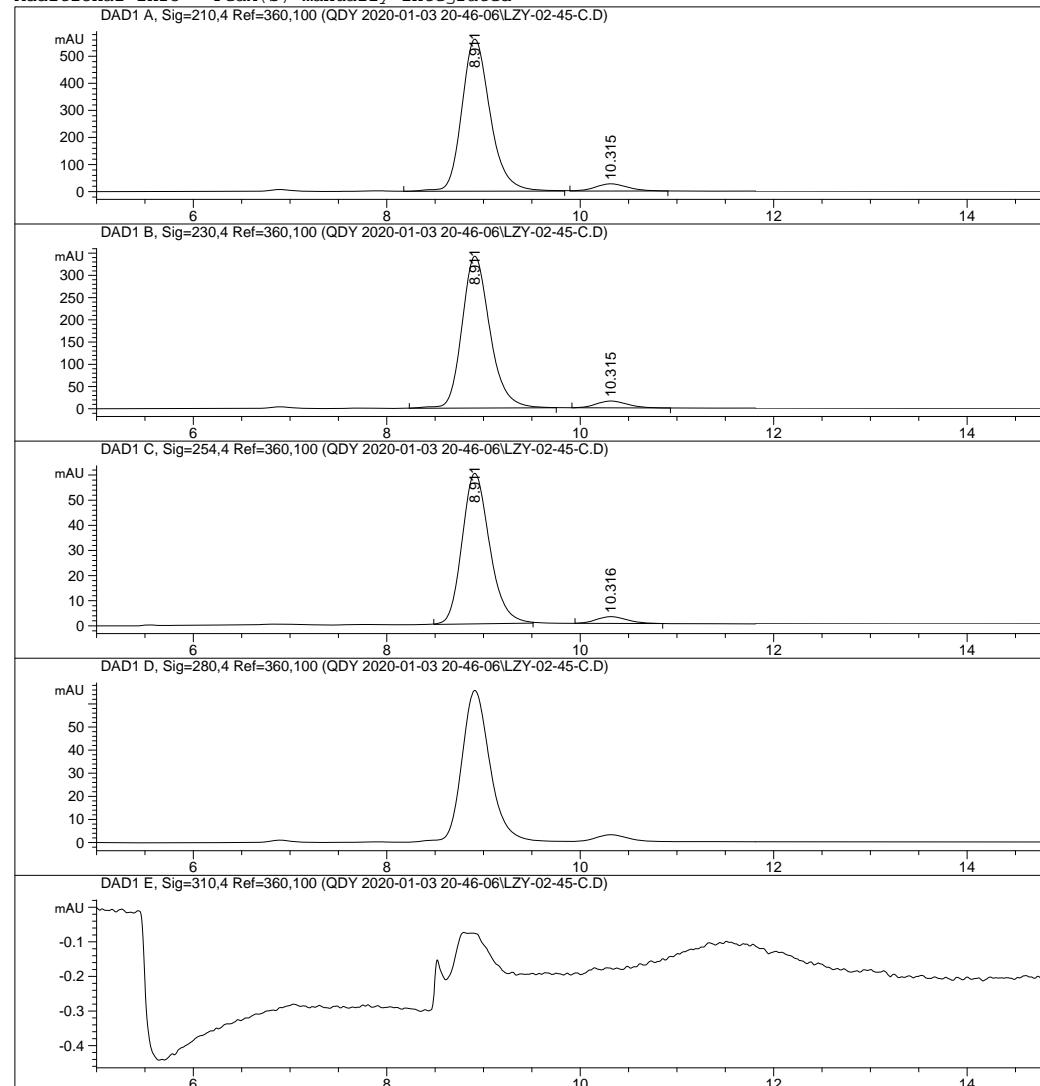
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

=====
 Acq. Operator : Seq. Line : 18
 Acq. Instrument : Instrument 1 Location : Vial 84
 Injection Date : 1/4/2020 2:37:46 AM Inj : 1
 Inj Volume : 5.000 μ l
 Different Inj Volume from Sequence ! Actual Inj Volume : 4.000 μ l
 Acq. Method : C:\CHEM32\1\DATA\QDY 2020-01-03 20-46-06\IC-05-20.M
 Last changed : 6/25/2018 1:35:20 AM
 Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
 Last changed : 11/24/2020 3:21:46 PM
 (modified after loading)
 Additional Info : Peak(s) manually integrated



=====
 Area Percent Report
 =====

Sorted By : Signal
 Multiplier : 1.0000
 Dilution : 1.0000
 Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.911	VB	0.3109	1.1411e4	561.88049	94.7741
2	10.315	BB	0.3607	629.21075	26.91560	5.2259
Totals :				1.2040e4	588.79610	

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

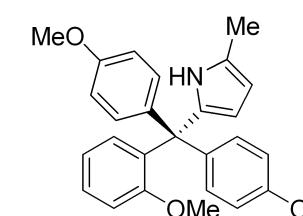
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.911	BB	0.3075	6835.92041	341.57336	95.2597
2	10.315	BB	0.3438	340.16577	15.39702	4.7403
Totals :				7176.08618	356.97038	

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.911	BB	0.3044	1183.26990	59.91114	95.2574
2	10.316	BB	0.3351	58.91190	2.69437	4.7426
Totals :				1242.18180	62.60551	

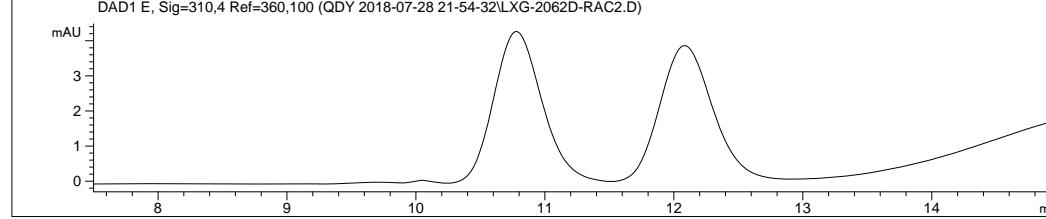
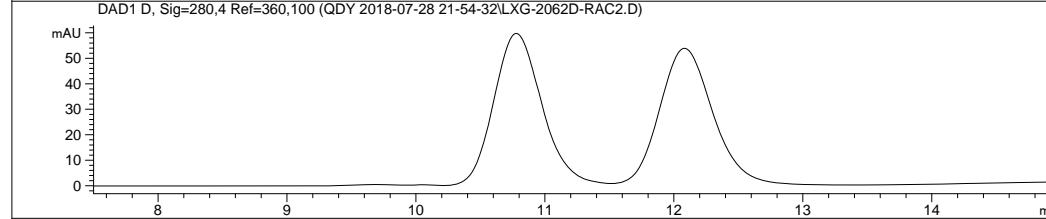
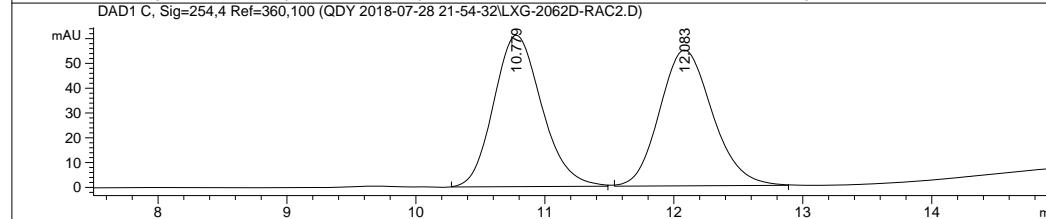
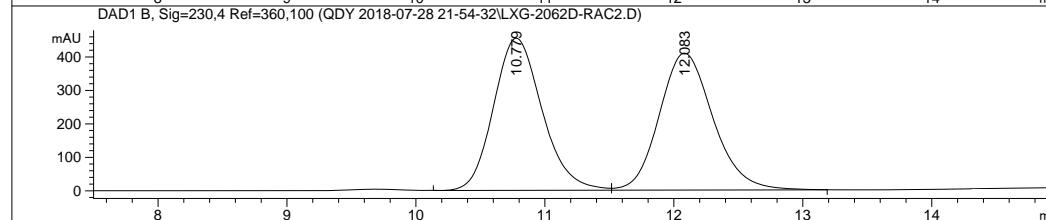
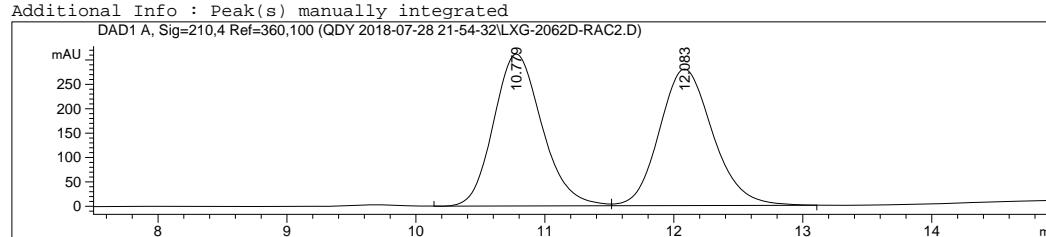
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 4
Acq. Instrument : Instrument 1             Location : Vial 70
Injection Date : 7/28/2018 10:40:28 PM       Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2018-07-28 21-54-32\IC-03-30.M
Last changed : 7/28/2018 10:39:35 PM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 4:51:53 PM        (modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.779	VV	0.4014	8103.85645	311.56830	50.1858
2	12.083	VB	0.4451	8043.84814	280.02228	49.8142

Totals : 1.61477e4 591.59058

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.779	VV	0.4023	1.18681e4	454.95987	50.0875
2	12.083	VB	0.4469	1.18267e4	409.59912	49.9125

Totals : 2.36948e4 864.55899

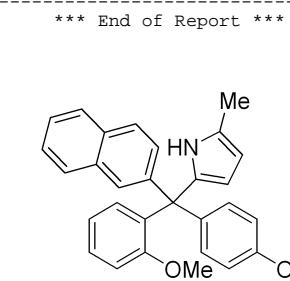
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.779	BB	0.3996	1578.74353	61.07904	50.3709
2	12.083	BB	0.4413	1555.49561	54.78296	49.6291

Totals : 3134.23914 115.86200

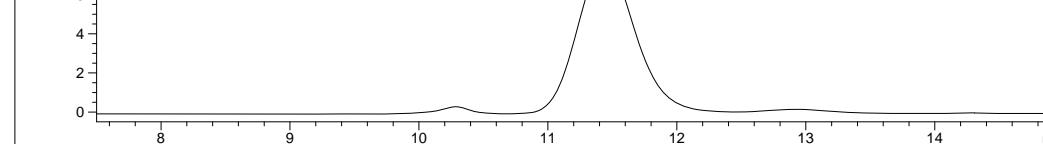
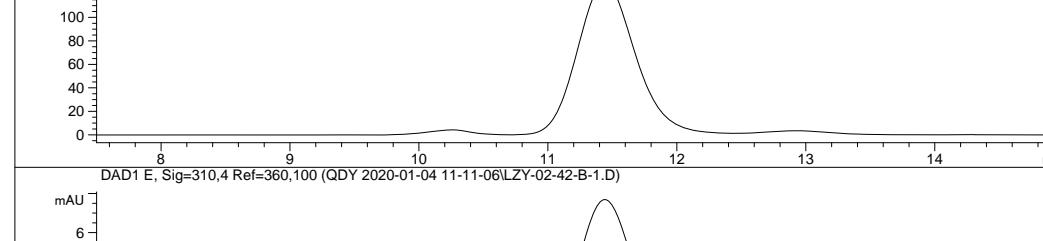
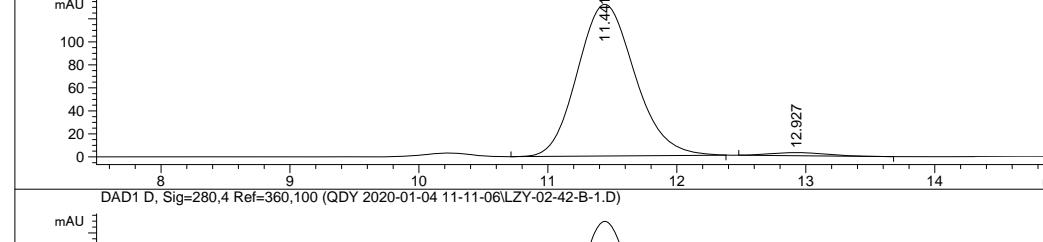
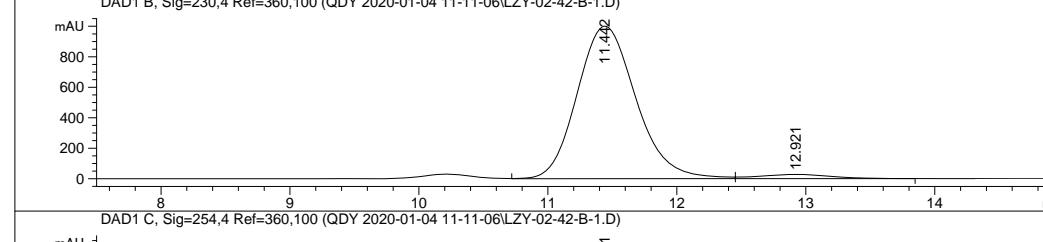
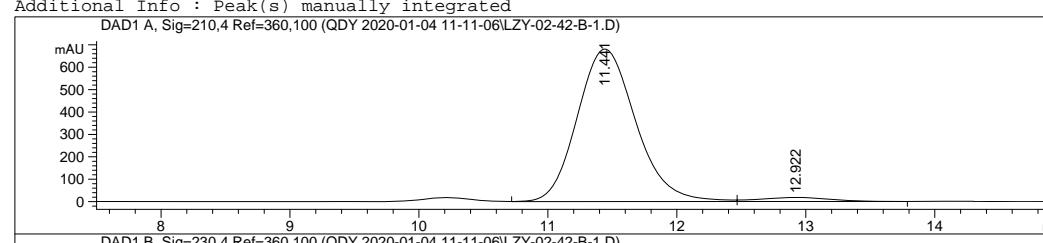
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

```
=====
Acq. Operator : Seq. Line : 2
Acq. Instrument : Instrument 1 Location : Vial 81
Injection Date : 1/4/2020 11:24:02 AM Inj : 1
Inj Volume : 5.000 µl
Different Inj Volume from Sequence ! Actual Inj Volume : 4.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-01-04 11-11-06\IC-03-30.M
Last changed : 1/4/2020 11:23:09 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 4:58:53 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.441	VB	0.4889	2.14076e4	680.77356	96.9799
2	12.922	BB	0.5514	666.66748	18.11265	3.0201
Totals :					2.20743e4	698.88621

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

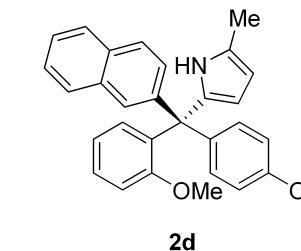
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.442	VV	0.4930	3.16293e4	1000.15277	96.7740
2	12.921	VB	0.5688	1054.37415	27.65328	3.2260
Totals :					3.26837e4	1027.80605

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.441	VB	0.4829	4081.43433	131.92607	98.0331
2	12.927	BB	0.4493	81.88832	2.71925	1.9669
Totals :					4163.32265	134.64532

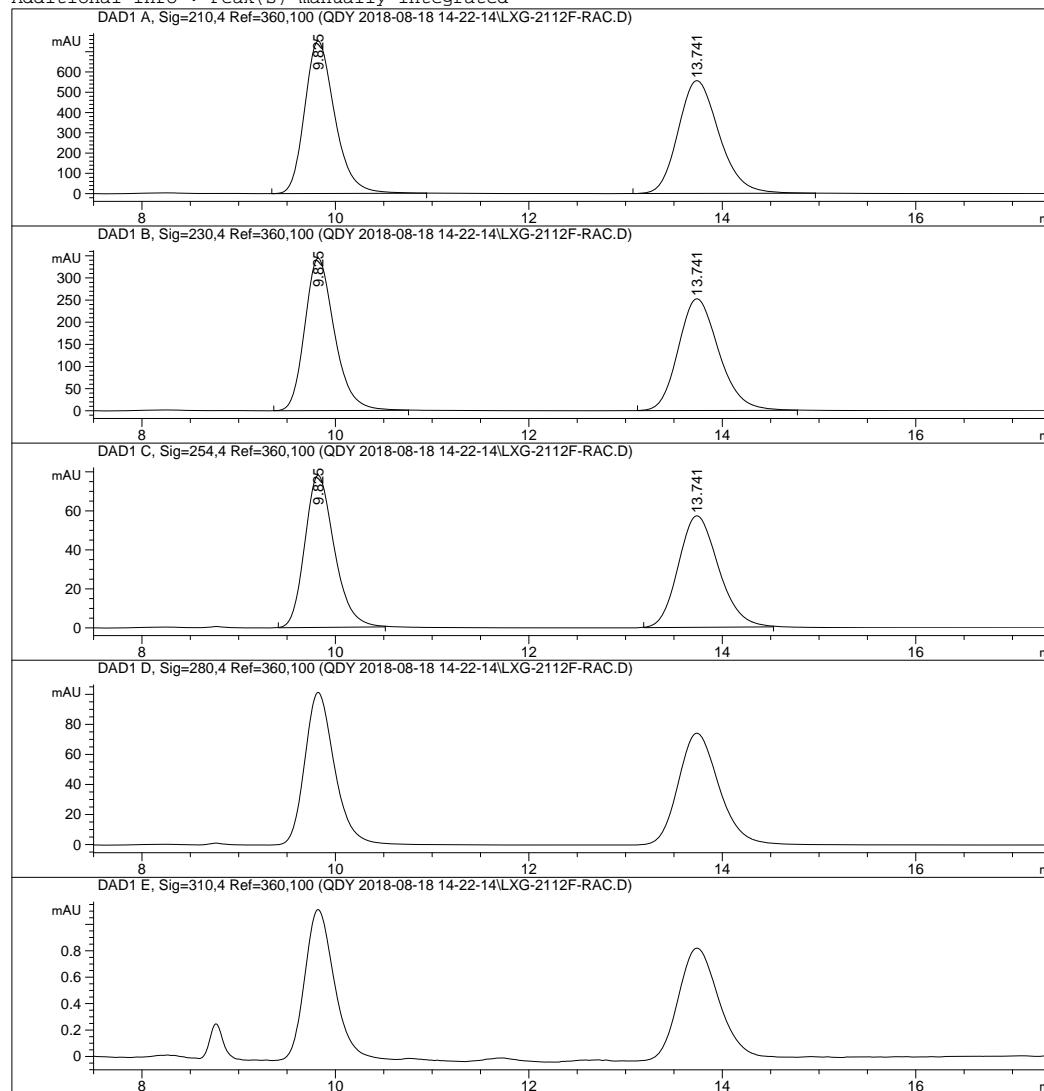
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 18
Acq. Instrument : Instrument 1             Location : Vial 67
Injection Date : 8/18/2018 8:15:50 PM        Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !       Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2018-08-18 14-22-14\IC-05-20.M
Last changed : 6/25/2018 1:35:20 AM
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 5:00:30 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



Data File C:\CHEM32\1\DATA\QDY 2018-08-18 14-22-14\LXG-2112F-RAC.D
Sample Name:

===== Area Percent Report =====

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.825	BB	0.3304	1.62644e4	751.89240	49.8544
2	13.741	BB	0.4551	1.63594e4	556.24500	50.1456
Totals :					3.26238e4	1308.13739

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

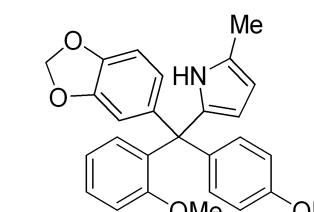
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.825	BB	0.3268	7339.66406	344.20276	49.9939
2	13.741	BB	0.4495	7341.46680	252.26933	50.0061
Totals :					1.46811e4	596.47209

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.825	BB	0.3253	1652.97595	77.99519	49.9956
2	13.741	BB	0.4472	1653.26514	57.20220	50.0044
Totals :					3306.24109	135.19738

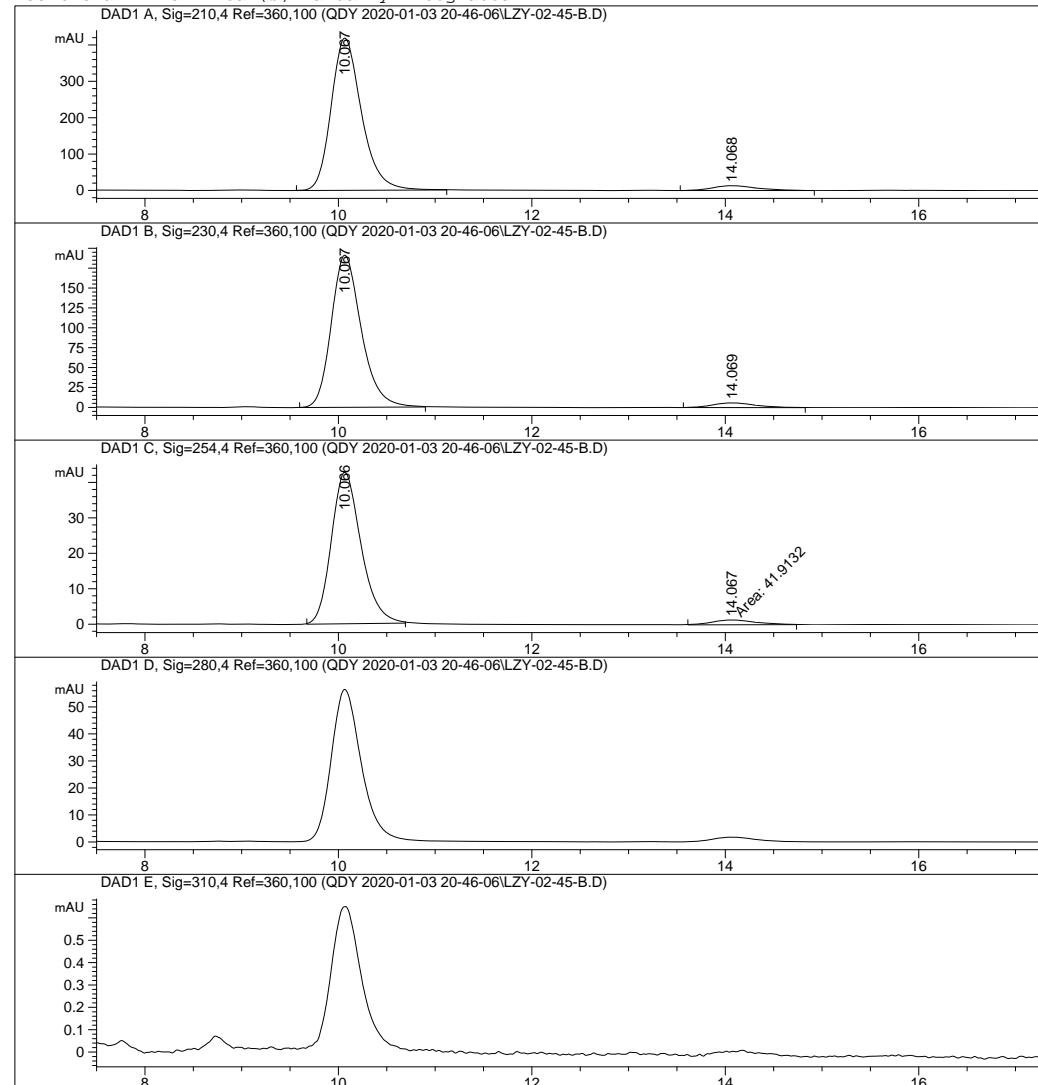
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100

**2e-Rac**

Sample Name:

```
=====
Acq. Operator : Seq. Line : 17
Acq. Instrument : Instrument 1 Location : Vial 83
Injection Date : 1/4/2020 2:16:31 AM Inj : 1
Inj Volume : 5.000 µl
Different Inj Volume from Sequence ! Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-01-03 20-46-06\IC-05-20.M
Last changed : 6/25/2018 1:35:20 AM
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 5:00:30 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.067	BB	0.3325	9129.89355	418.59555	95.7750
2	14.068	BB	0.4503	402.75327	13.33718	4.2250

Totals : 9532.64682 431.93273

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.067	BB	0.3298	4127.84521	191.22440	95.8873
2	14.069	BB	0.4608	177.04517	5.88723	4.1127

Totals : 4304.89038 197.11163

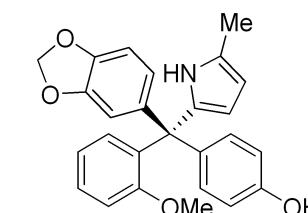
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.066	BB	0.3270	912.89362	42.76598	95.6103
2	14.067	MM	0.5100	41.91323	1.36984	4.3897

Totals : 954.80684 44.13582

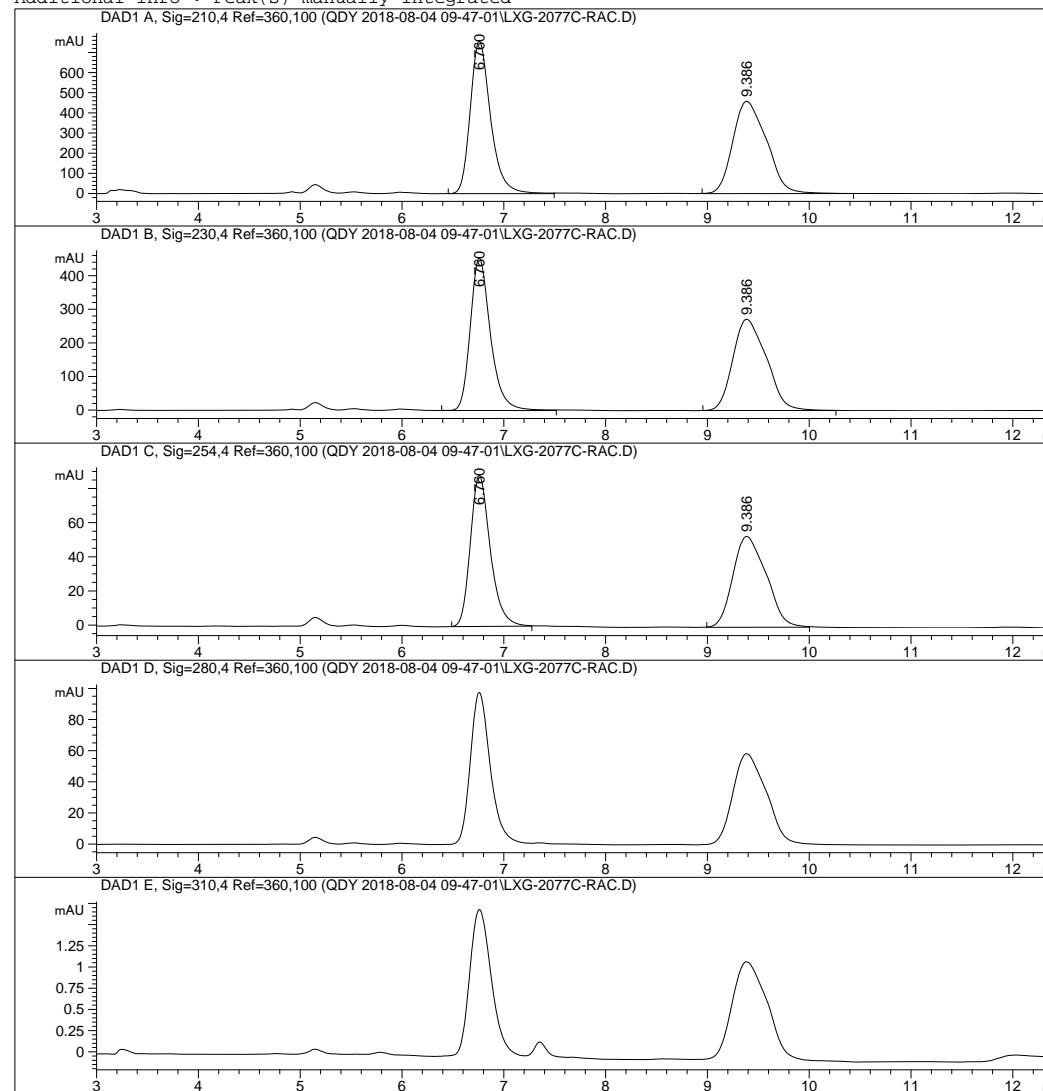
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

=====
 Acq. Operator : Seq. Line : 2
 Acq. Instrument : Instrument 1 Location : Vial 65
 Injection Date : 8/4/2018 10:00:32 AM Inj : 1
 Inj Volume : 5.000 μ l
 Different Inj Volume from Sequence ! Actual Inj Volume : 3.000 μ l
 Acq. Method : C:\CHEM32\1\DATA\QDY 2018-08-04 09-47-01\IC-10-20.M
 Last changed : 6/25/2018 2:45:05 PM
 Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
 Last changed : 11/24/2020 5:05:20 PM
 (modified after loading)
 Additional Info : Peak(s) manually integrated



===== Area Percent Report =====

Sorted By : Signal
 Multiplier : 1.0000
 Dilution : 1.0000
 Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.760	BB	0.2119	1.05081e4	757.44958	49.9888
2	9.386	BB	0.3400	1.05128e4	457.72897	50.0112

Totals : 2.10209e4 1215.17856

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.760	VB	0.2090	6253.97754	453.15485	50.1874
2	9.386	BB	0.3386	6207.27100	271.59845	49.8126

Totals : 1.24612e4 724.75330

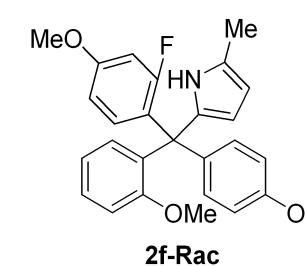
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.760	BB	0.2068	1206.07825	88.58540	50.0098
2	9.386	BB	0.3364	1205.60706	53.18979	49.9902

Totals : 2411.68530 141.77519

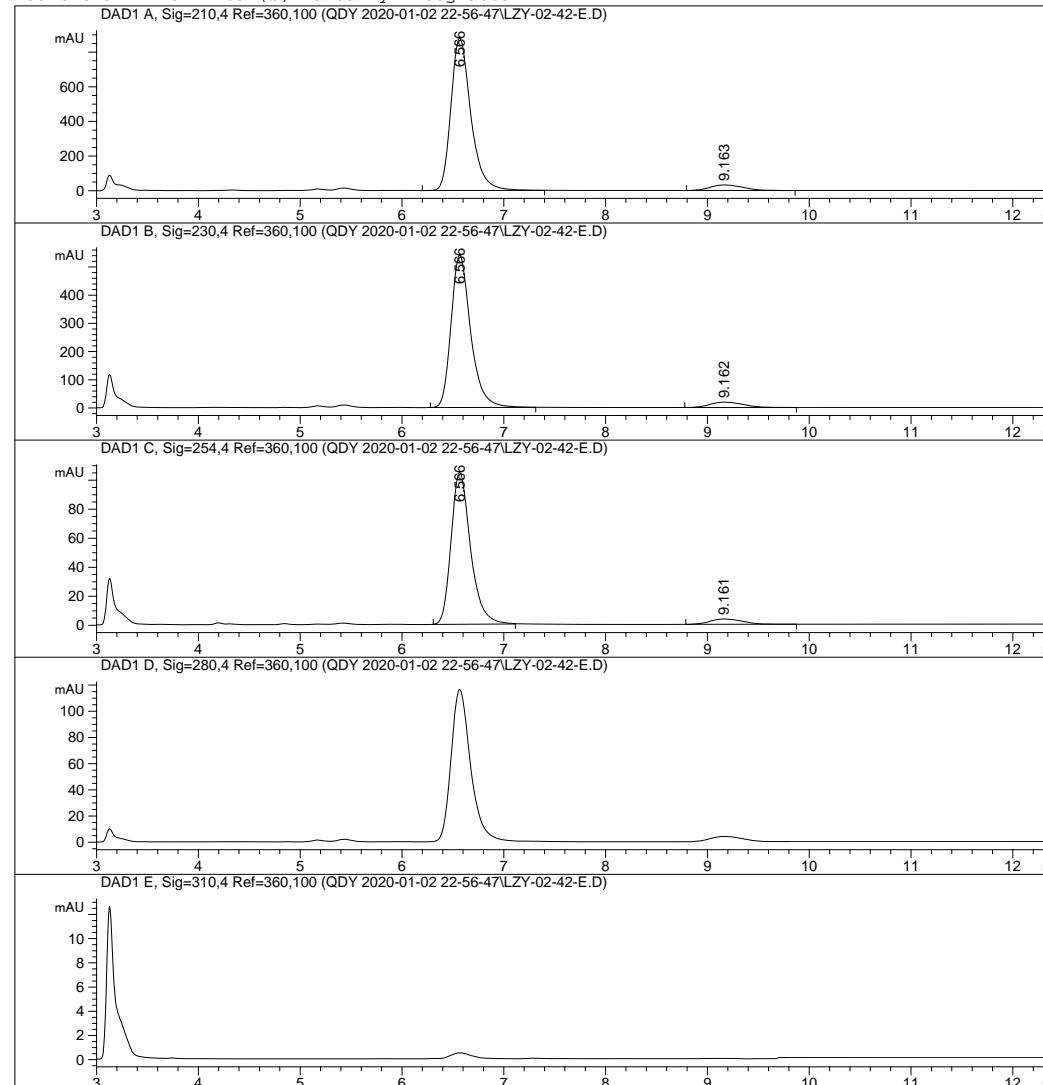
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

=====
 Acq. Operator : Seq. Line : 19
 Acq. Instrument : Instrument 1 Location : Vial 93
 Injection Date : 1/3/2020 7:09:51 AM Inj : 1
 Inj Volume : 5.000 μ l
 Different Inj Volume from Sequence ! Actual Inj Volume : 3.000 μ l
 Acq. Method : C:\CHEM32\1\DATA\QDY 2020-01-02 22-56-47\IC-10-20.M
 Last changed : 6/25/2018 2:45:05 PM
 Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
 Last changed : 11/24/2020 5:05:20 PM
 (modified after loading)
 Additional Info : Peak(s) manually integrated



Sample Name:

=====
 Area Percent Report
 =====

Sorted By : Signal
 Multiplier : 1.0000
 Dilution : 1.0000
 Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.566	VB	0.2083	1.19365e4	879.94104	94.2529
2	9.163	BB	0.3637	727.83667	32.69982	5.7471
Totals :					1.26644e4	912.64086

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

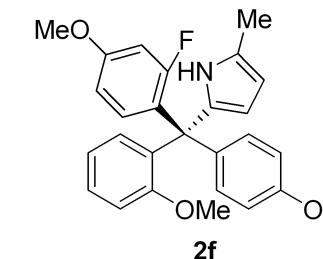
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.566	BB	0.2040	7221.55176	540.07068	94.3915
2	9.162	BB	0.3651	429.08340	19.31783	5.6085
Totals :					7650.63516	559.38850

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

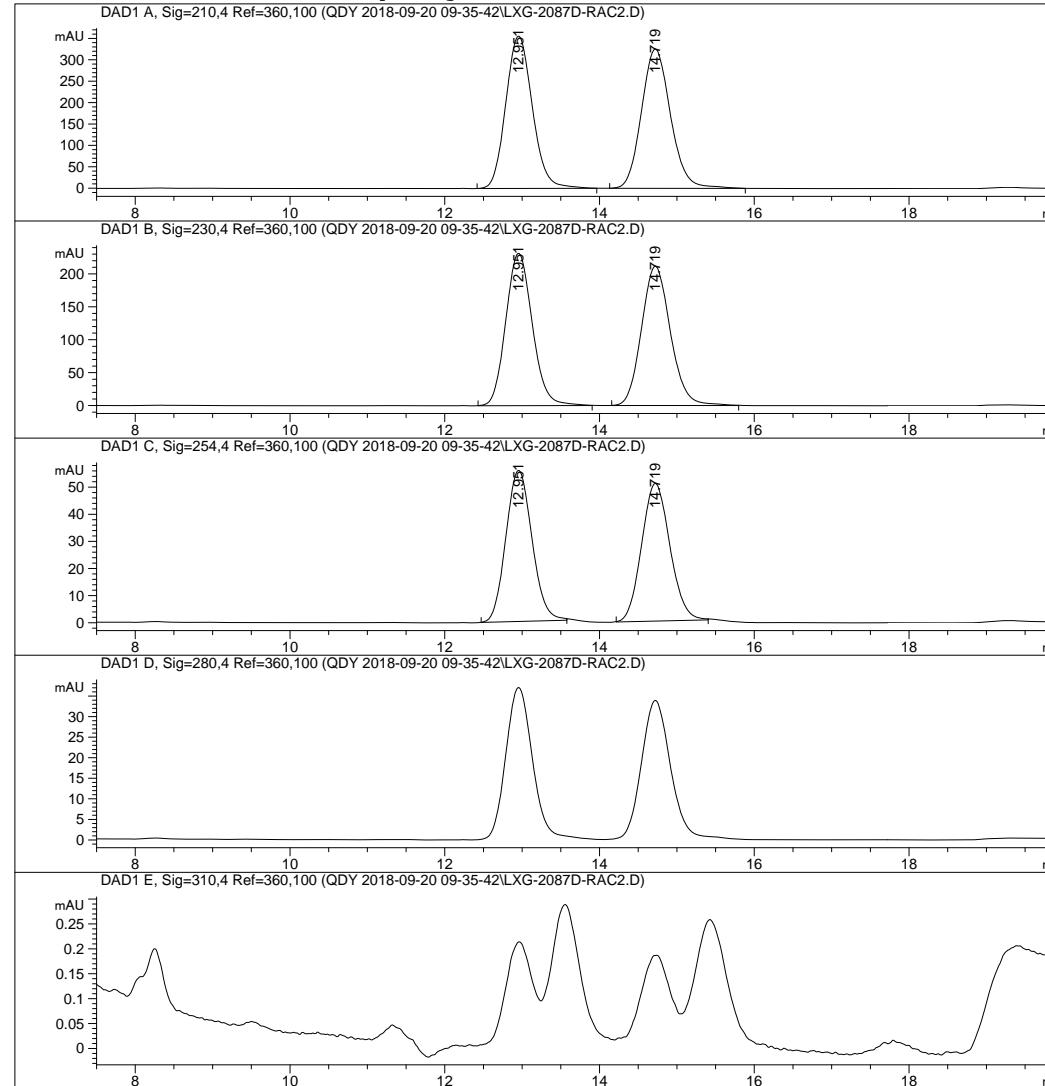
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	6.566	BB	0.2029	1390.37463	104.74338	94.4932
2	9.161	BB	0.3611	81.02735	3.70638	5.5068
Totals :					1471.40199	108.44976

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



```
=====
Acq. Operator :                               Seq. Line :    7
Acq. Instrument : Instrument 1             Location : Vial 63
Injection Date : 9/20/2018 12:26:06 PM       Inj : 1
                                                Inj Volume : 5.000 µl
Acq. Method :  C:\CHEM32\1\DATA\QDY 2018-09-20 09-35-42\AD-05-20.M
Last changed : 9/20/2018 12:25:13 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 5:10:49 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.951	BB	0.3743	8495.08398	356.08966	49.9916
2	14.719	BB	0.4046	8497.92188	325.53262	50.0084

Totals : 1.69930e4 681.62228

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.951	BB	0.3692	5503.25586	231.59538	50.0357
2	14.719	BB	0.4031	5495.40234	211.55399	49.9643

Totals : 1.09987e4 443.14937

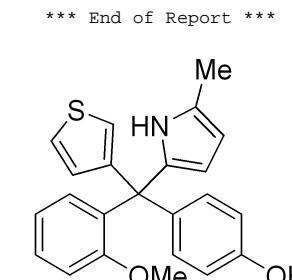
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.951	BB	0.3671	1304.67407	55.73872	50.2787
2	14.719	BB	0.3966	1290.21057	50.75733	49.7213

Totals : 2594.88464 106.49606

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

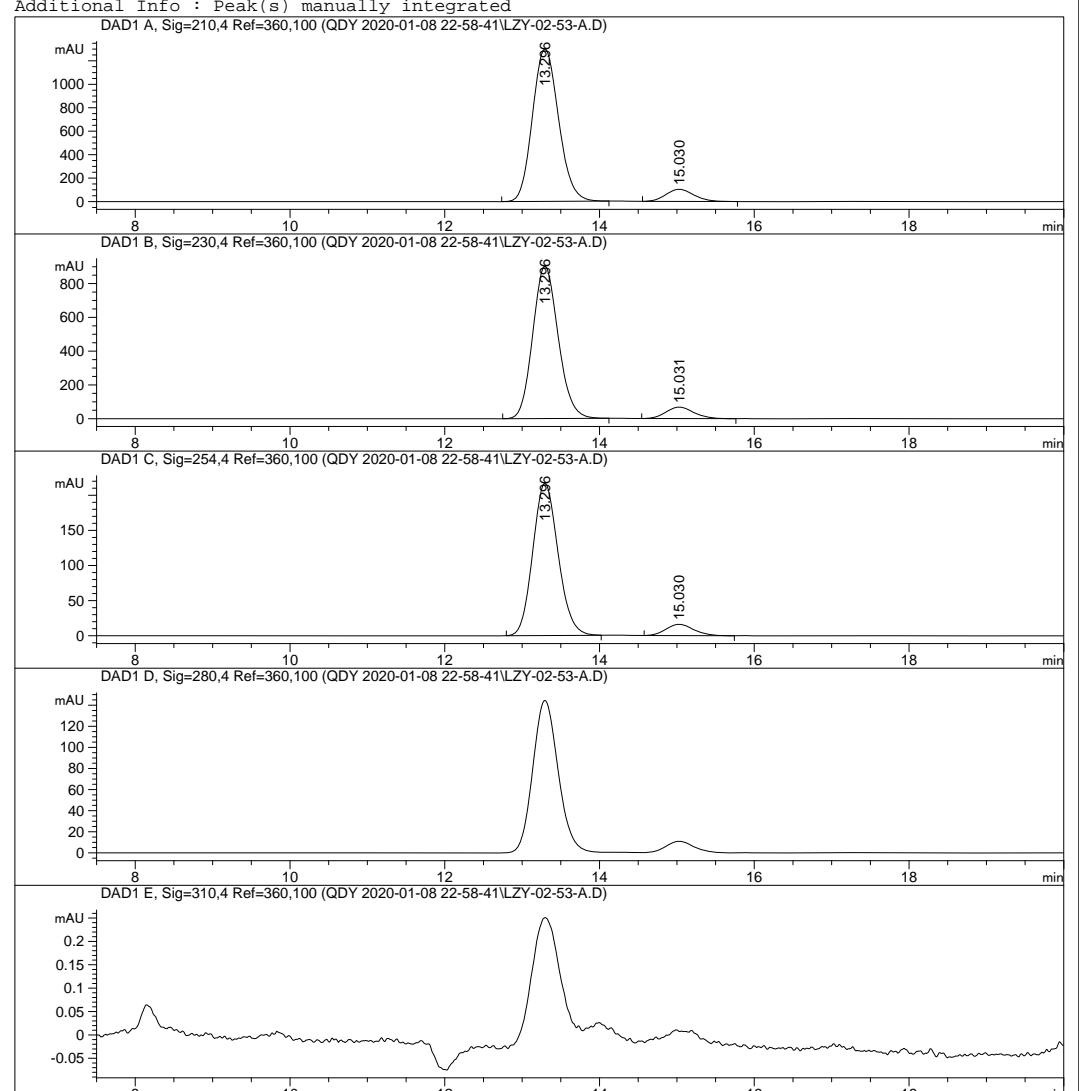
Signal 5: DAD1 E, Sig=310,4 Ref=360,100



2g-Rac

Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 9
Acq. Instrument : Instrument 1             Location : Vial 94
Injection Date : 1/9/2020 1:09:11 AM        Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 4.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-01-08 22-58-41\AD-05-20.M
Last changed : 1/9/2020 1:08:18 AM          (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 5:17:00 PM         (modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	13.296	BB	0.3672	3.04301e4	1299.27332	92.1366
2	15.030	BB	0.3860	2597.06372	103.75465	7.8634

Totals : 3.30272e4 1403.02796

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	13.296	BB	0.3541	2.05805e4	902.40430	92.3592
2	15.031	BB	0.3883	1702.60925	67.95928	7.6408

Totals : 2.22831e4 970.36358

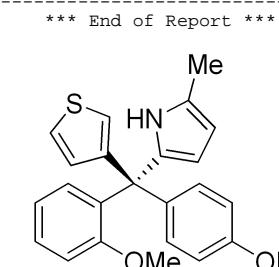
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	13.296	BB	0.3530	4924.81836	216.78227	92.4727
2	15.030	BB	0.3889	400.88251	16.07955	7.5273

Totals : 5325.70087 232.86182

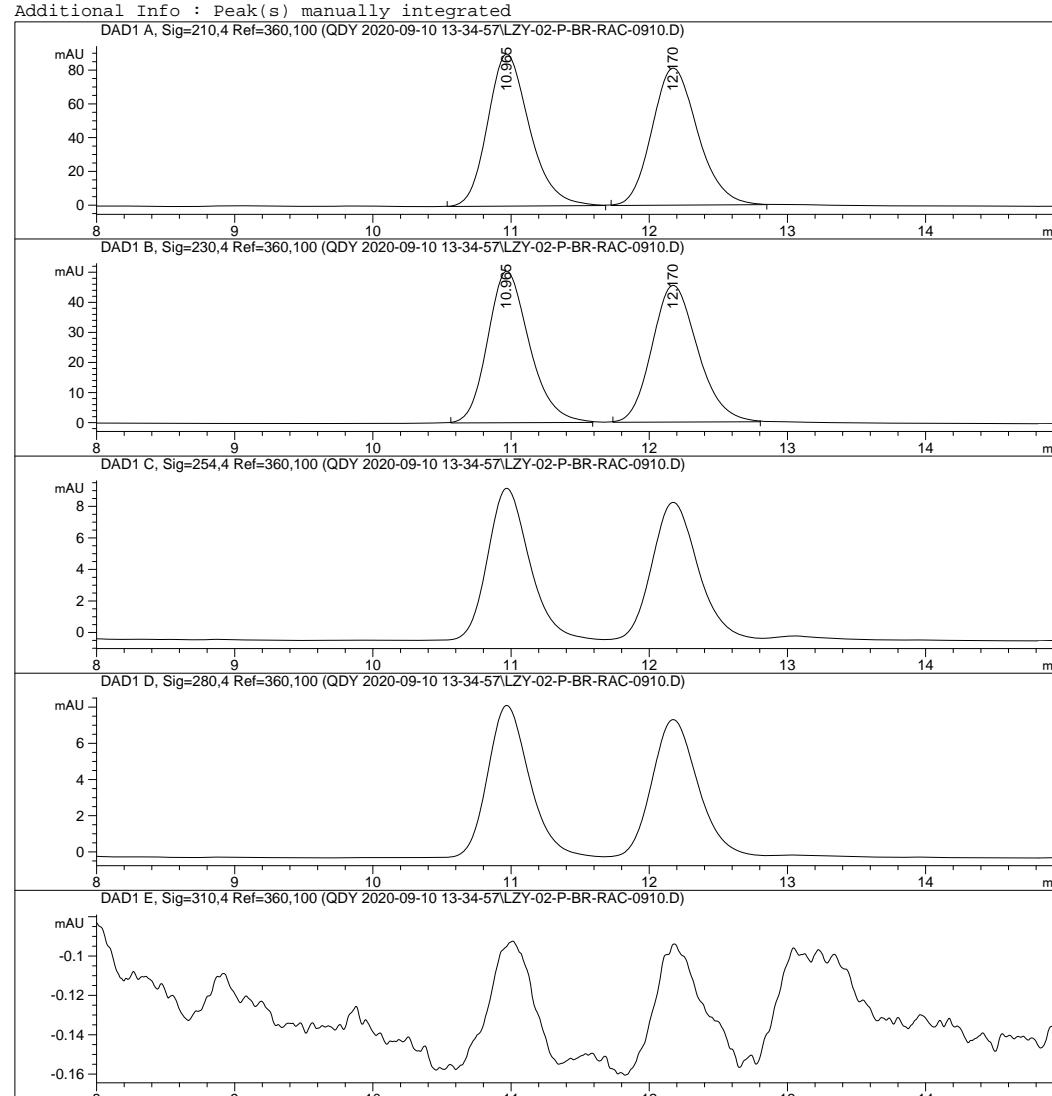
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

```
=====
Acq. Operator : Seq. Line : 30
Acq. Instrument : Instrument 1 Location : Vial 92
Injection Date : 9/10/2020 10:47:37 PM Inj : 1
Inj Volume : 5.000 µl
Different Inj Volume from Sequence ! Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-09-10 13-34-57\IC-02-20.M
Last changed : 9/10/2020 10:46:45 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 2:43:12 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.965	BB	0.3233	1893.90601	90.08872	50.3472
2	12.170	BB	0.3542	1867.78394	81.24993	49.6528

Totals : 3761.68994 171.33865

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.965	BB	0.3202	1054.73108	50.38411	50.2401
2	12.170	BB	0.3540	1044.65027	45.48300	49.7599

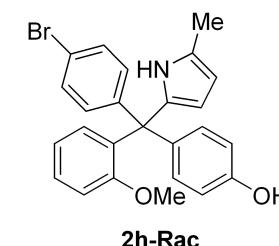
Totals : 2099.38135 95.86711

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100

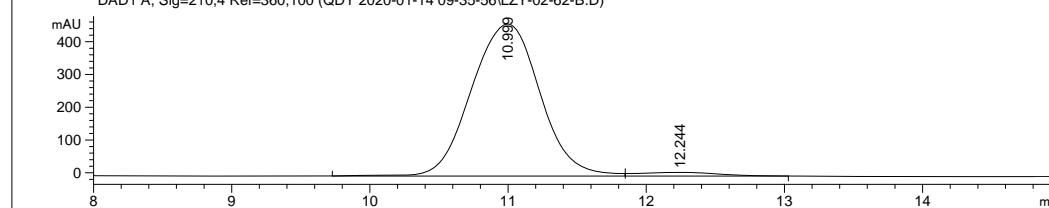
===== *** End of Report ***



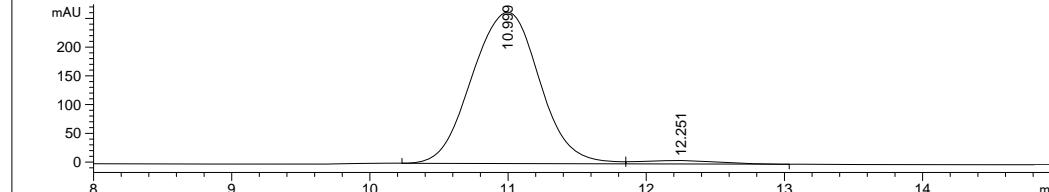
Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 3
Acq. Instrument : Instrument 1             Location : Vial 92
Injection Date : 1/14/2020 10:20:16 AM       Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 4.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-01-14 09-35-56\IC-02-30.M
Last changed : 1/14/2020 9:48:23 AM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 2:43:12 PM        (modified after loading)
Additional Info : Peak(s) manually integrated
```

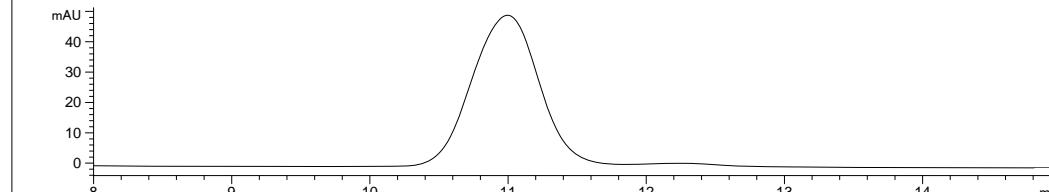
DAD1 A, Sig=210,4 Ref=360,100 (QDY 2020-01-14 09-35-56\LZY-02-62-B.D)



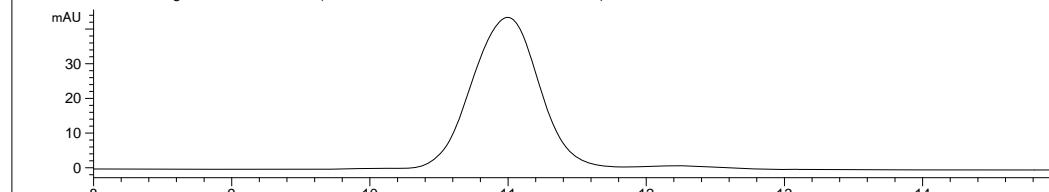
DAD1 B, Sig=230,4 Ref=360,100 (QDY 2020-01-14 09-35-56\LZY-02-62-B.D)



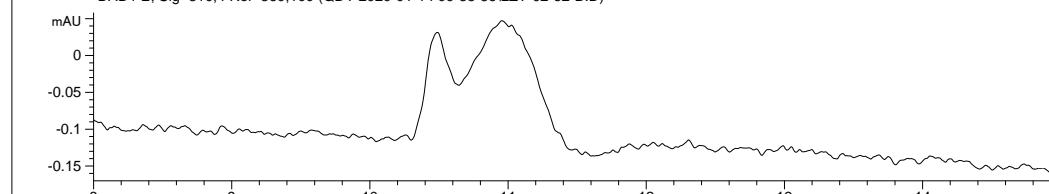
DAD1 C, Sig=254,4 Ref=360,100 (QDY 2020-01-14 09-35-56\LZY-02-62-B.D)



DAD1 D, Sig=280,4 Ref=360,100 (QDY 2020-01-14 09-35-56\LZY-02-62-B.D)



DAD1 E, Sig=310,4 Ref=360,100 (QDY 2020-01-14 09-35-56\LZY-02-62-B.D)



Sample Name:

```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.999	BV	0.5506	1.6009e4	463.92984	97.2556
2	12.244	VB	0.5797	451.74054	11.26800	2.7444
Totals :				1.64607e4	475.19784	

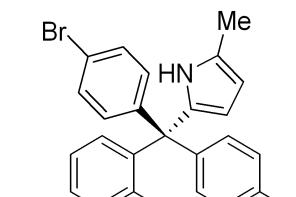
Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.999	BV	0.5418	8935.40234	263.35187	97.4918
2	12.251	VB	0.5462	229.88318	5.93145	2.5082
Totals :				9165.28552	269.28331	

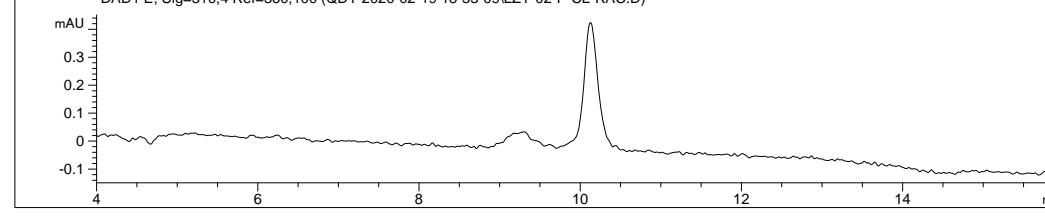
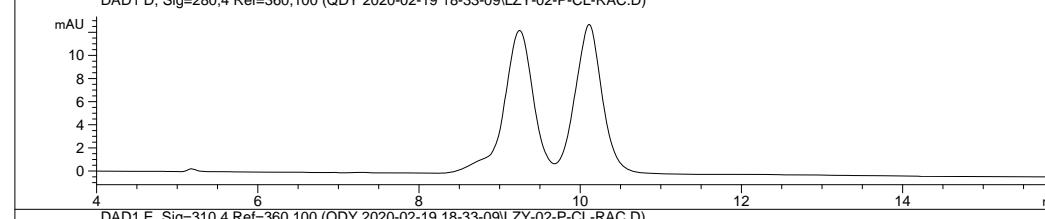
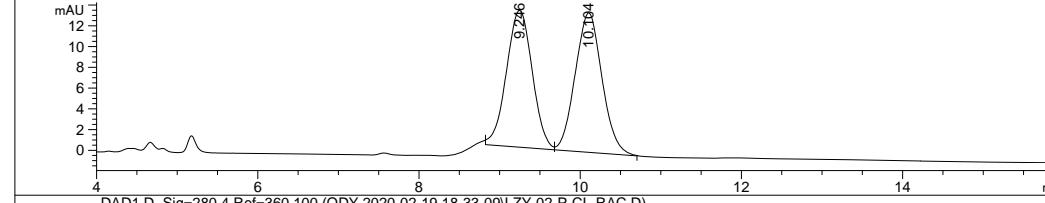
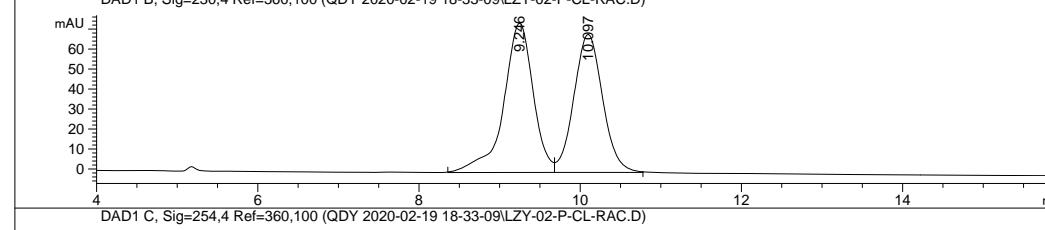
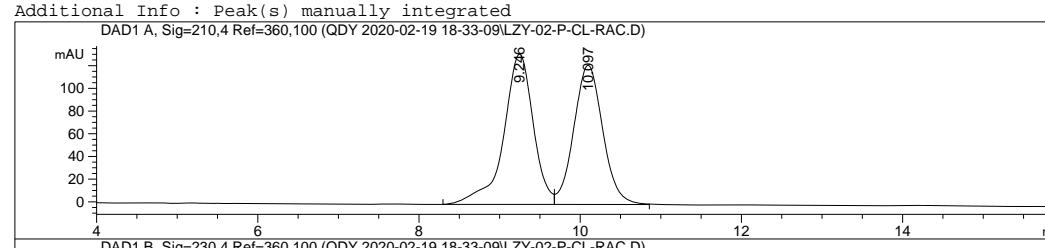
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100

**2h**

=====
 Acq. Operator : Seq. Line : 17
 Acq. Instrument : Instrument 1 Location : Vial 98
 Injection Date : 2/20/2020 12:03:49 AM Inj : 1
 Inj Volume : 5.000 μ l
 Different Inj Volume from Sequence ! Actual Inj Volume : 3.000 μ l
 Acq. Method : C:\CHEM32\1\DATA\QDY 2020-02-19 18-33-09\IC-03-20.M
 Last changed : 2/19/2020 11:41:56 PM (modified after loading)
 Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
 Last changed : 11/24/2020 2:48:22 PM (modified after loading)
 Additional Info : Peak(s) manually integrated



=====
 Area Percent Report
 =====

Sorted By : Signal
 Multiplier : 1.0000
 Dilution : 1.0000
 Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.246	BV	0.3879	3384.03345	132.54121	52.5468
2	10.097	VB	0.3846	3056.00171	122.70458	47.4532

Totals : 6440.03516 255.24580

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.246	BV	0.3848	1900.14783	74.68499	52.6616
2	10.097	VB	0.3847	1708.07288	69.03443	47.3384

Totals : 3608.22070 143.71942

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.246	BV	0.3505	292.29184	13.19219	48.5612
2	10.104	VB	0.3568	309.61221	13.43825	51.4388

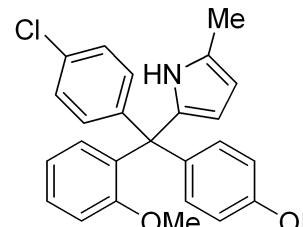
Totals : 601.90405 26.63044

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100

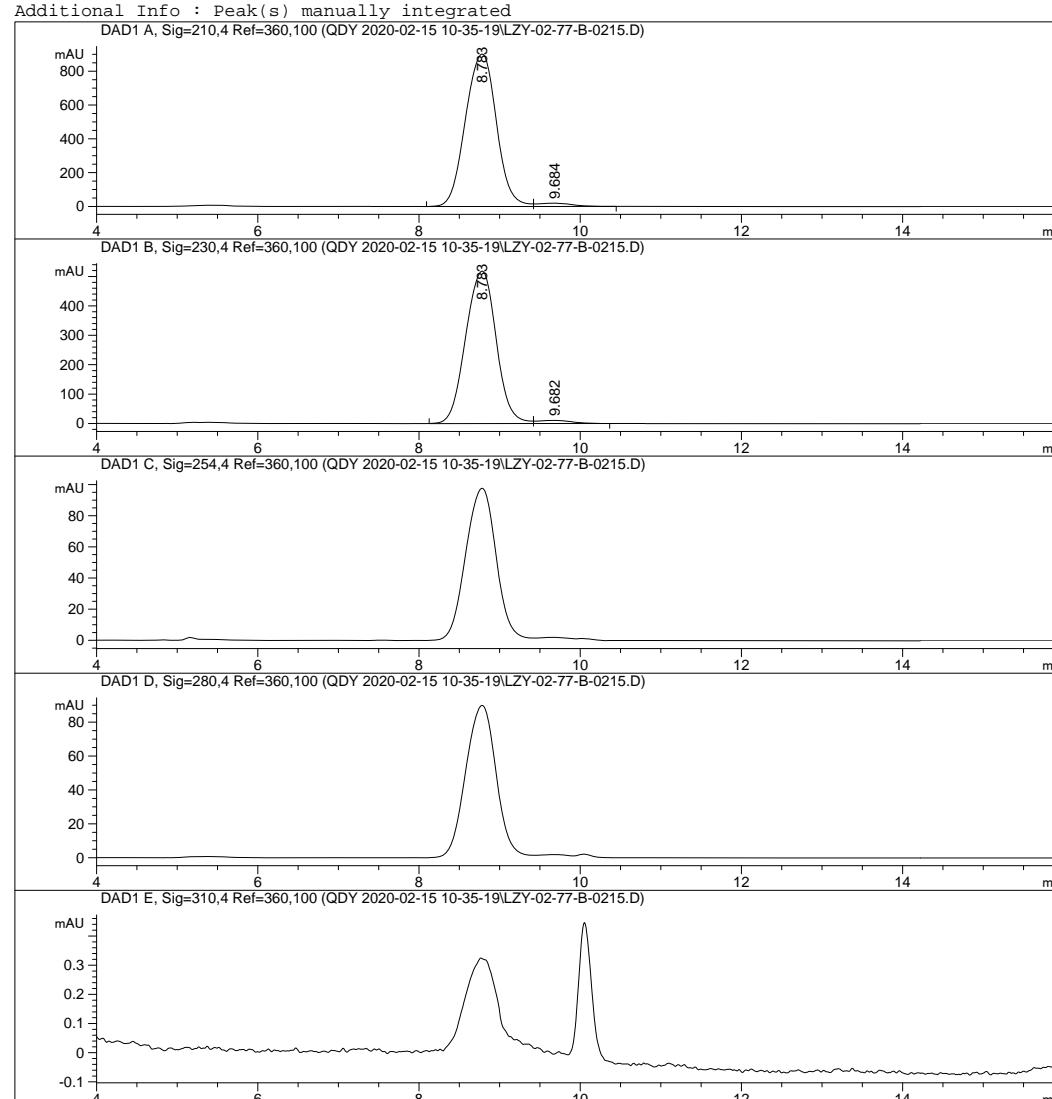
=====

*** End of Report ***



Sample Name:

```
=====
Acq. Operator :                               Seq. Line :   3
Acq. Instrument : Instrument 1             Location : Vial 92
Injection Date : 2/15/2020 11:21:29 AM       Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 4.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-02-15 10-35-19\IC-03-20.M
Last changed : 2/15/2020 11:20:39 AM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 8:38:09 PM        (modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

```
=====
Area Percent Report
=====
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100
```

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.783	BV	0.4250	2.40417e4	901.99292	97.6987
2	9.684	VB	0.4423	566.31244	19.41345	2.3013

Totals : 2.46080e4 921.40637

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

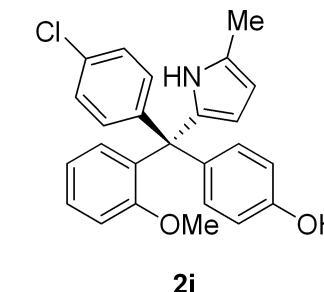
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.783	BV	0.4207	1.36287e4	518.34625	97.7838
2	9.682	VB	0.4400	308.88916	10.66089	2.2162

Totals : 1.39376e4 529.00714

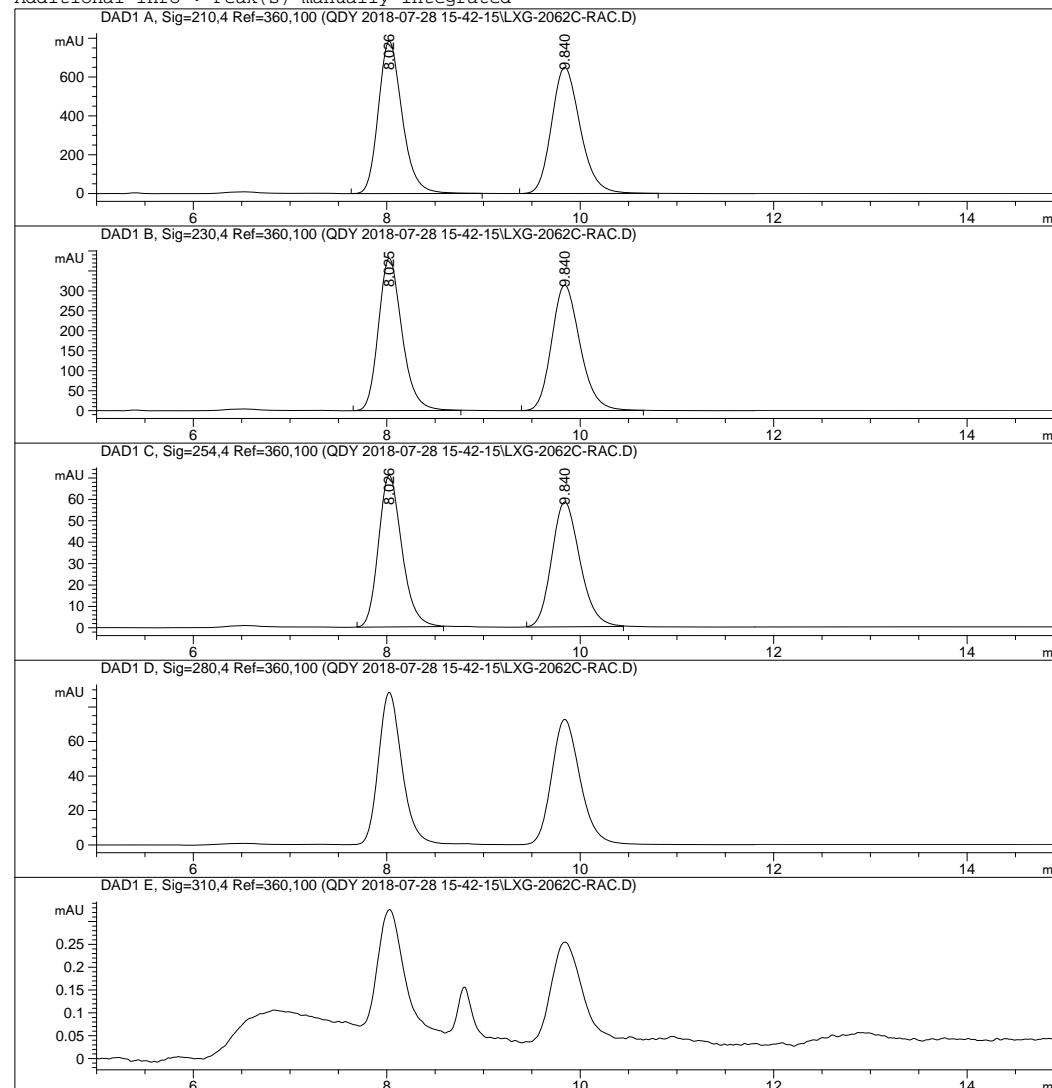
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



```
=====
Acq. Operator :                               Seq. Line : 4
Acq. Instrument : Instrument 1             Location : Vial 69
Injection Date : 7/28/2018 4:38:09 PM        Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !       Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2018-07-28 15-42-15\IC-05-20.M
Last changed : 6/25/2018 1:35:20 AM
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 3:21:46 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.026	BB	0.2630	1.33934e4	785.38489	50.0026
2	9.840	BB	0.3185	1.33920e4	649.64520	49.9974

Totals : 2.67854e4 1435.03009

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.025	BB	0.2608	6438.29688	381.67386	50.0475
2	9.840	BB	0.3165	6426.07471	314.29382	49.9525

Totals : 1.28644e4 695.96768

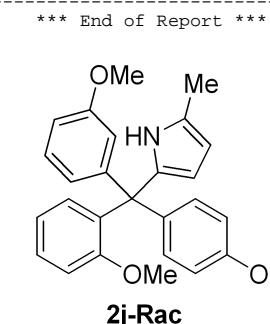
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.026	BB	0.2597	1188.85413	70.86761	50.0373
2	9.840	BB	0.3153	1187.08093	58.36302	49.9627

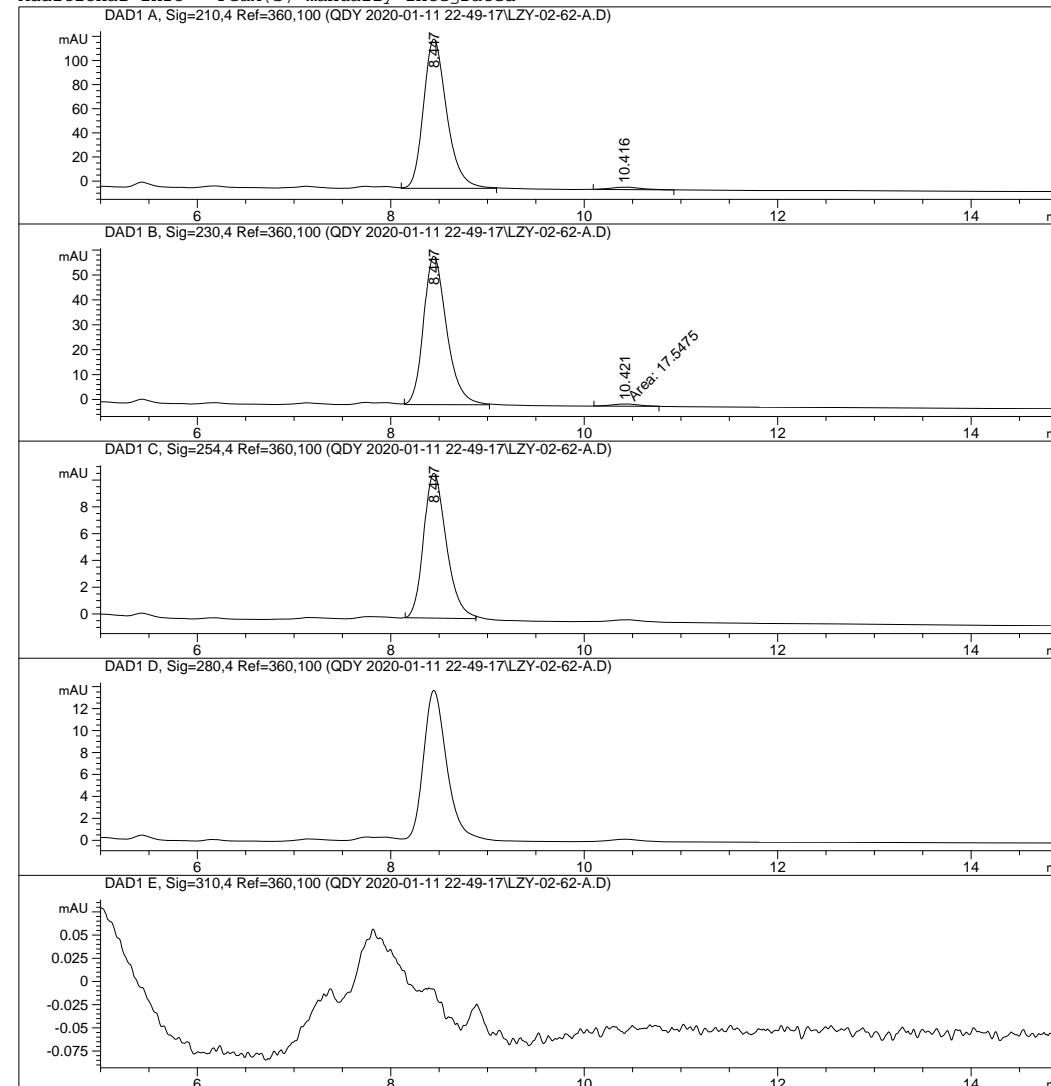
Totals : 2375.93506 129.23063

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



```
=====
Acq. Operator :                               Seq. Line : 2
Acq. Instrument : Instrument 1             Location : Vial 81
Injection Date : 1/11/2020 11:02:30 PM       Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 1.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-01-11 22-49-17\IC-05-20.M
Last changed : 6/25/2018 1:35:20 AM
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 4:45:25 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



```
=====
Area Percent Report
=====

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100
Peak RetTime Type Width Area Height Area
# [min] [min] [mAU*s] [mAU] %
-----|-----|-----|-----|-----|
1 8.447 VB 0.2545 2062.70898 123.73442 97.9110
2 10.416 BB 0.2953 44.00969 2.04281 2.0890

Totals : 2106.71867 125.77723

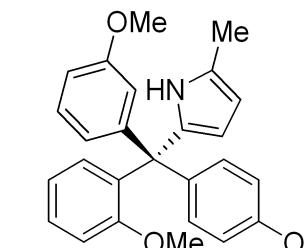
Signal 2: DAD1 B, Sig=230,4 Ref=360,100
Peak RetTime Type Width Area Height Area
# [min] [min] [mAU*s] [mAU] %
-----|-----|-----|-----|-----|
1 8.447 BB 0.2537 989.99640 59.61902 98.2584
2 10.421 MM 0.3280 17.54748 8.91600e-1 1.7416

Totals : 1007.54388 60.51062

Signal 3: DAD1 C, Sig=254,4 Ref=360,100
Peak RetTime Type Width Area Height Area
# [min] [min] [mAU*s] [mAU] %
-----|-----|-----|-----|-----|
1 8.447 BB 0.2520 176.08235 10.81169 100.0000

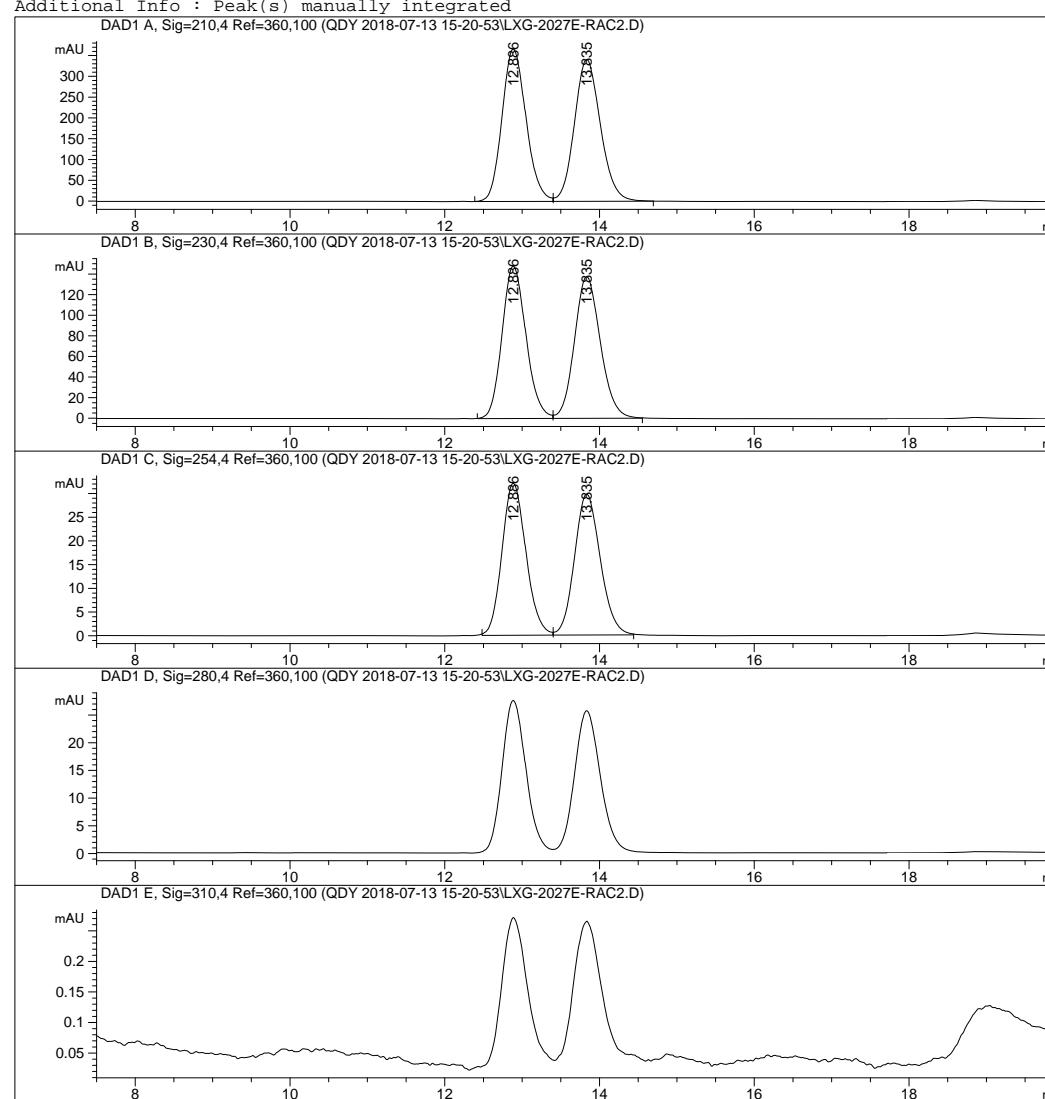
Totals : 176.08235 10.81169

Signal 4: DAD1 D, Sig=280,4 Ref=360,100
Signal 5: DAD1 E, Sig=310,4 Ref=360,100
=====
```



Sample Name:

```
=====
Acq. Operator :                               Seq. Line :    7
Acq. Instrument : Instrument 1             Location : Vial 67
Injection Date : 7/13/2018 5:00:32 PM       Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2018-07-13 15-20-53\AD-05-30.M
Last changed : 7/17/2016 12:02:24 PM
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 5:17:00 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.886	BV	0.3340	7914.43408	366.33969	49.7959
2	13.835	VB	0.3628	7979.30078	341.21024	50.2041

Totals : 1.58937e4 707.54993

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.886	BV	0.3330	3183.48584	147.96762	49.8640
2	13.835	VB	0.3592	3200.85205	137.67828	50.1360

Totals : 6384.33789 285.64590

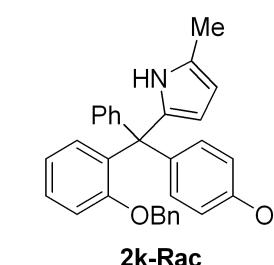
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.886	BV	0.3330	689.41772	32.04320	50.0482
2	13.835	VB	0.3580	688.09100	29.73436	49.9518

Totals : 1377.50873 61.77756

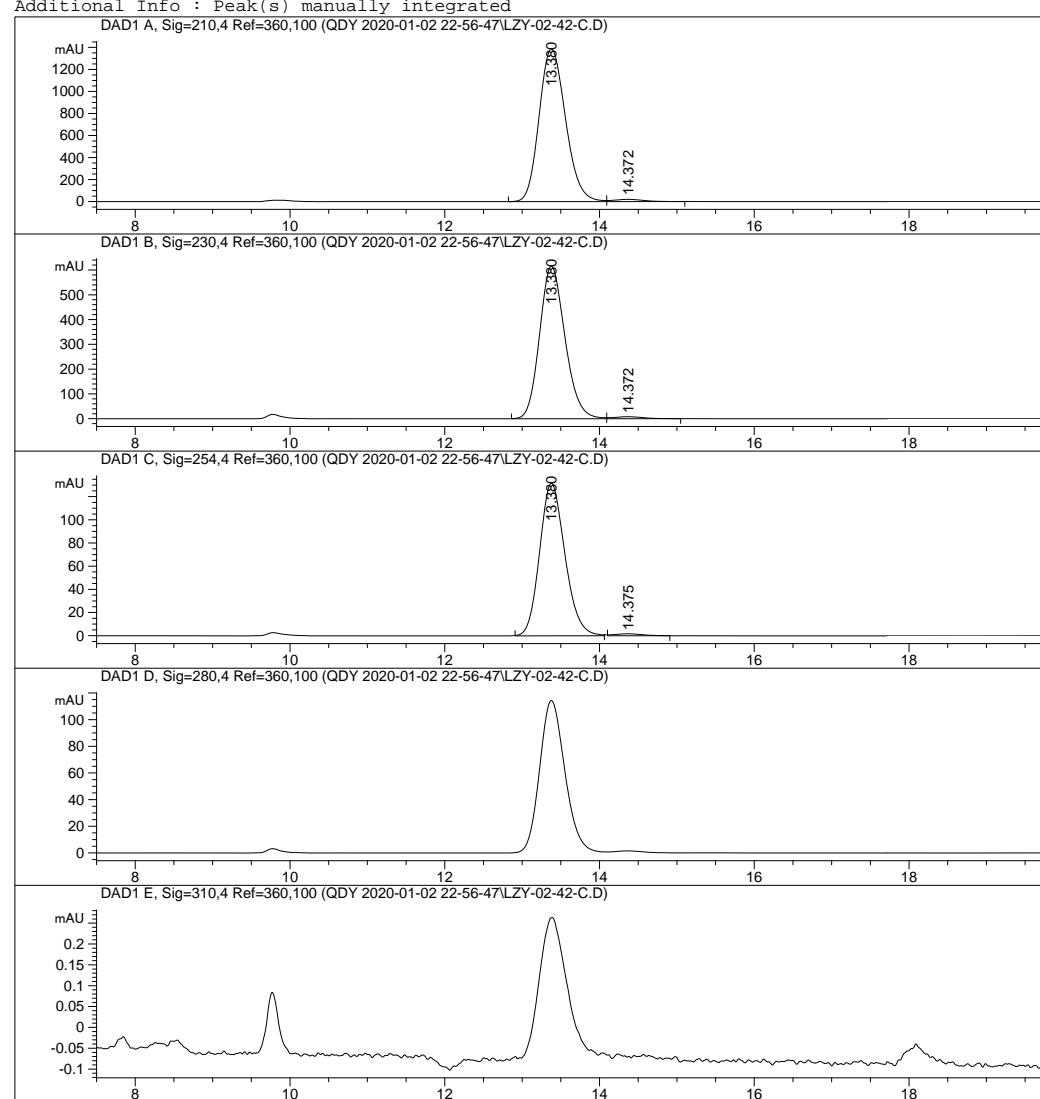
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

```
=====
Acq. Operator : Seq. Line : 12
Acq. Instrument : Instrument 1 Location : Vial 92
Injection Date : 1/3/2020 4:21:26 AM Inj : 1
Inj Volume : 5.000 µl
Different Inj Volume from Sequence ! Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-01-02 22-56-47\AD-05-30.M
Last changed : 7/17/2016 12:02:24 PM
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 8:39:23 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	13.380	BV	0.3707	3.28630e4	1385.41077	98.4422
2	14.372	VB	0.3881	520.05066	20.35078	1.5578
Totals :					3.33830e4	1405.76155

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

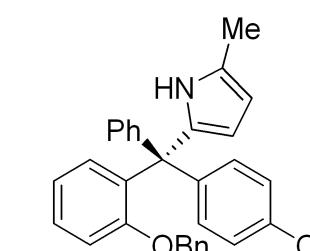
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	13.380	BV	0.3536	1.40167e4	615.71234	98.5424
2	14.372	VB	0.3918	207.32739	8.12356	1.4576
Totals :					1.422240e4	623.83590

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	13.380	BB	0.3508	2988.84521	131.66220	98.5785
2	14.375	BB	0.3519	43.10011	1.72278	1.4215
Totals :					3031.94532	133.38498

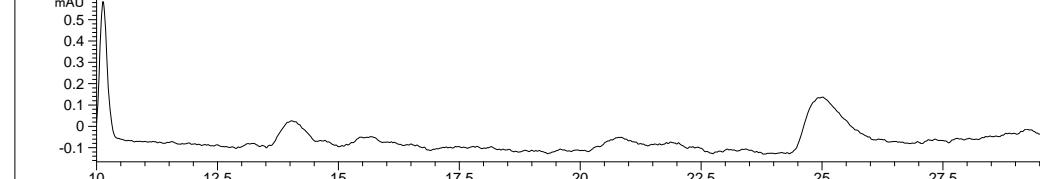
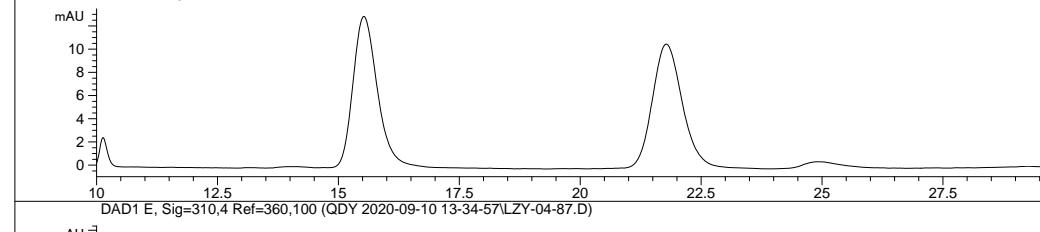
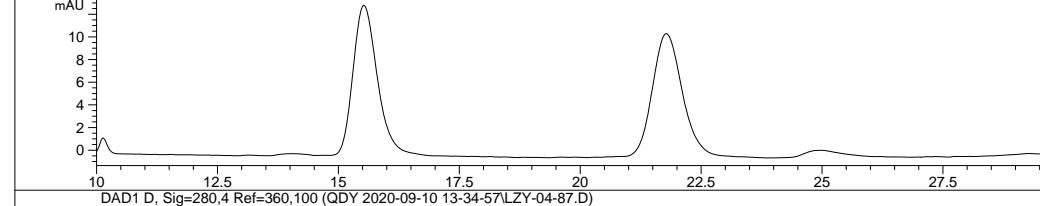
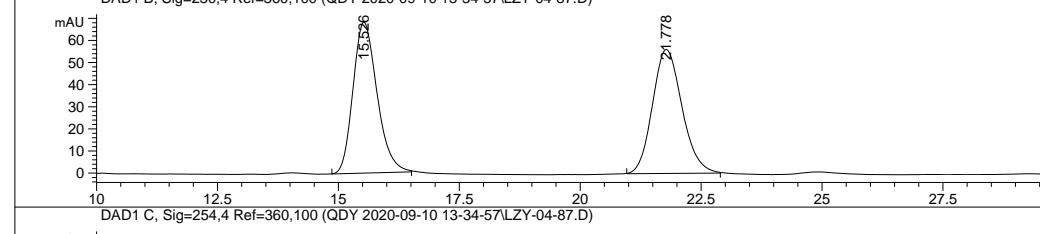
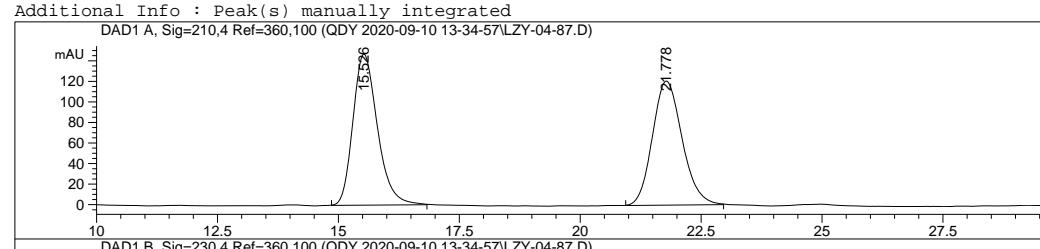
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 11
Acq. Instrument : Instrument 1             Location : Vial 91
Injection Date : 9/10/2020 5:12:31 PM       Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 2.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-09-10 13-34-57\IC-03-40.M
Last changed : 9/10/2020 5:11:38 PM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 2:52:38 PM        (modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

===== Area Percent Report =====

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	15.526	BB	0.5420	5167.96240	147.07161	50.1571
2	21.778	BB	0.6683	5135.58691	120.71682	49.8429
Totals :					1.03035e4	267.78843

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

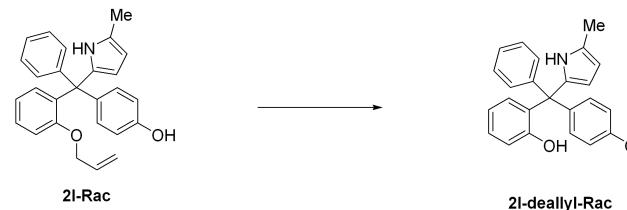
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	15.526	BB	0.5343	2353.96167	68.27108	49.7489
2	21.778	BB	0.6587	2377.72070	56.07371	50.2511
Totals :					4731.68237	124.34479

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

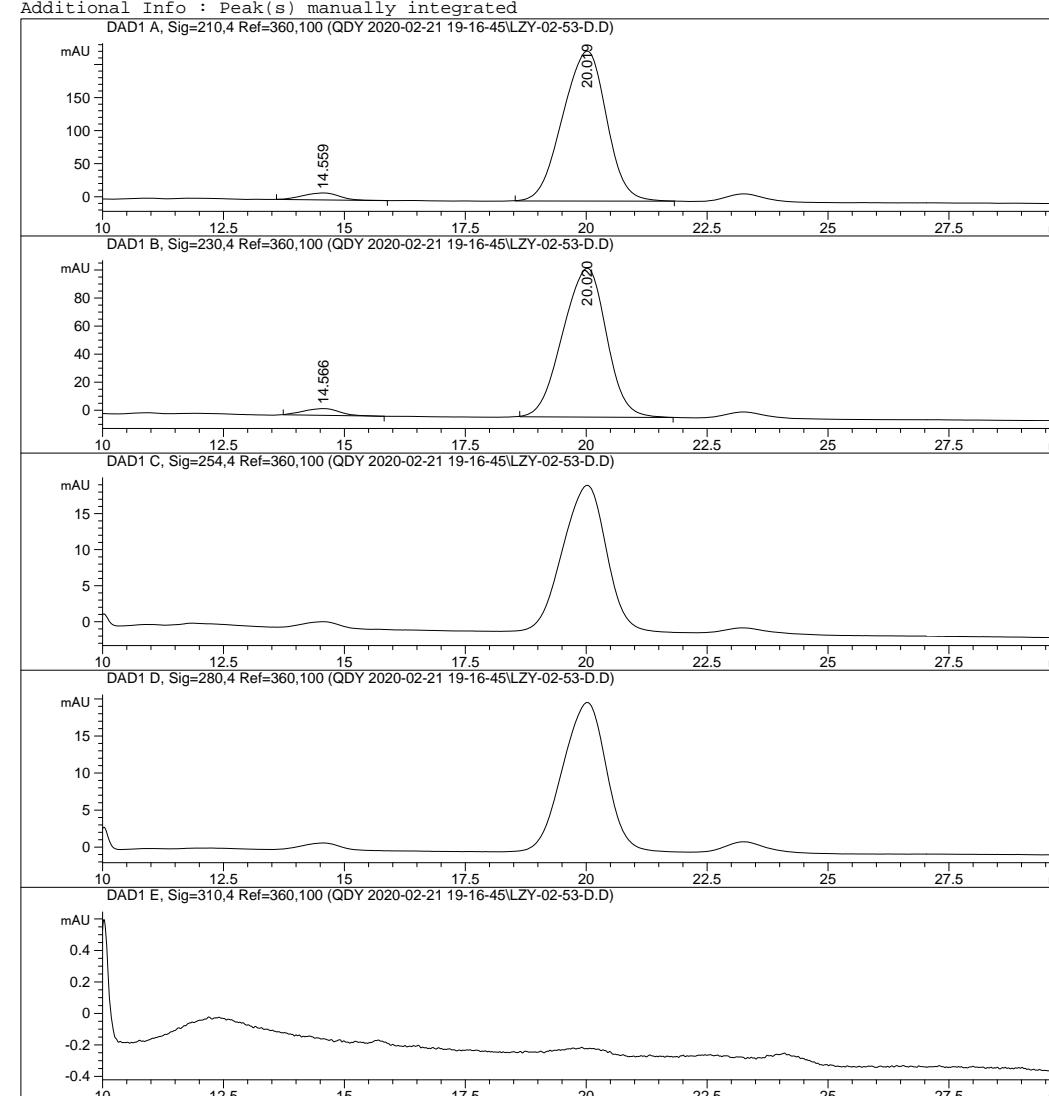
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100

=====
*** End of Report ***
=====



=====
 Acq. Operator : Seq. Line : 2
 Acq. Instrument : Instrument 1 Location : Vial 89
 Injection Date : 2/21/2020 7:30:02 PM Inj : 1
 Inj Volume : 5.000 μ l
 Different Inj Volume from Sequence ! Actual Inj Volume : 6.000 μ l
 Acq. Method : C:\CHEM32\1\DATA\QDY 2020-02-21 19-16-45\IC-03-30.M
 Last changed : 2/21/2020 7:29:10 PM (modified after loading)
 Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
 Last changed : 11/24/2020 2:52:38 PM (modified after loading)
 Additional Info : Peak(s) manually integrated



===== Area Percent Report =====

Sorted By : Signal
 Multiplier : 1.0000
 Dilution : 1.0000
 Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	14.559	BB	0.6502	544.63507	10.51568	3.6713
2	20.019	BB	1.0048	1.42902e4	226.88597	96.3287

Totals : 1.48348e4 237.40165

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	14.566	BB	0.7005	243.73860	4.81535	3.5336
2	20.020	BB	0.9974	6653.98584	106.14471	96.4664

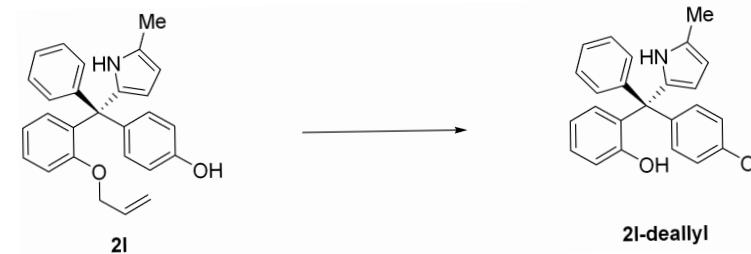
Totals : 6897.72444 110.96005

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

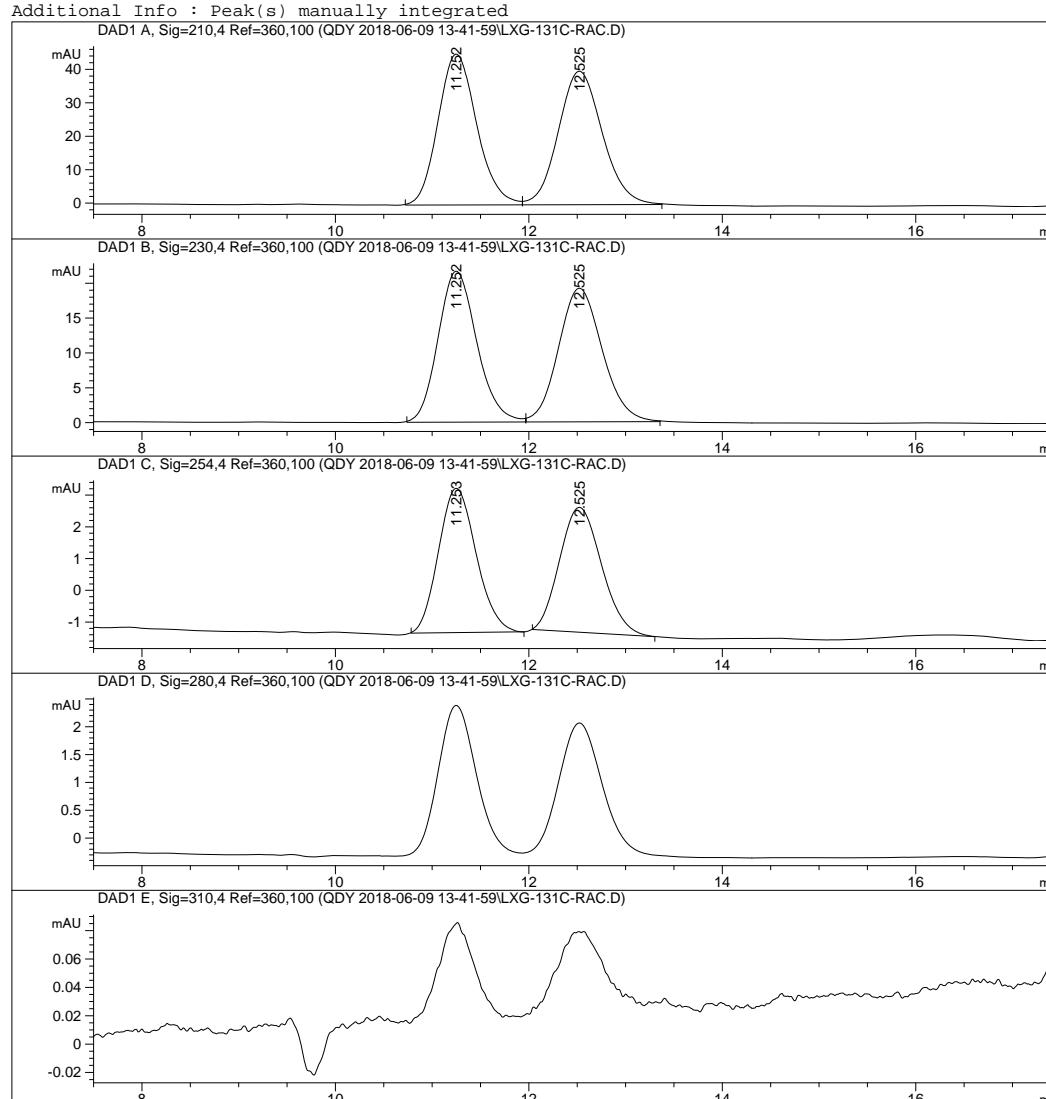
Signal 5: DAD1 E, Sig=310,4 Ref=360,100

=====
 *** End of Report ***
 =====



Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 2
Acq. Instrument : Instrument 1             Location : Vial 62
Injection Date : 6/9/2018 1:55:33 PM        Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 2.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2018-06-09 13-41-59\IC-30-30.M
Last changed : 6/9/2018 1:54:41 PM          (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:15:03 PM         (modified after loading)
Additional Info : Peak(s) manually integrated
```



Data File C:\CHEM32\1\DATA\QDY 2018-06-09 13-41-59\LXG-131C-RAC.D
Sample Name:

```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.252	BV	0.4298	1252.26025	45.12457	49.8303
2	12.525	VB	0.4924	1260.78906	39.93031	50.1697

Totals : 2513.04932 85.05489

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.252	BB	0.4300	601.50604	21.65733	49.9837
2	12.525	BB	0.4905	601.89801	19.15991	50.0163

Totals : 1203.40405 40.81724

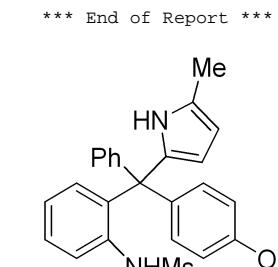
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.253	BB	0.4190	122.72398	4.54591	50.8898
2	12.525	BB	0.4671	118.43242	3.93561	49.1102

Totals : 241.15640 8.48152

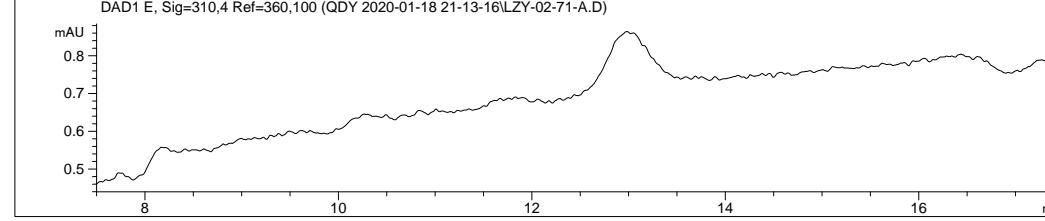
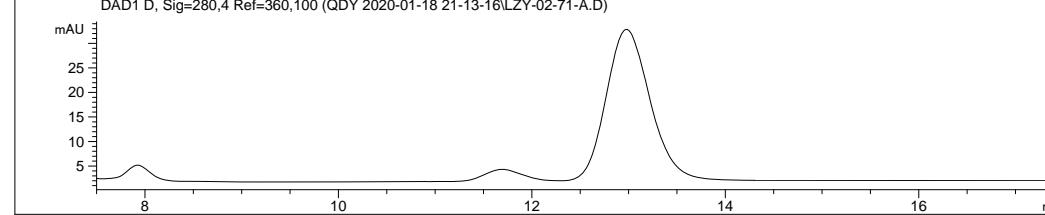
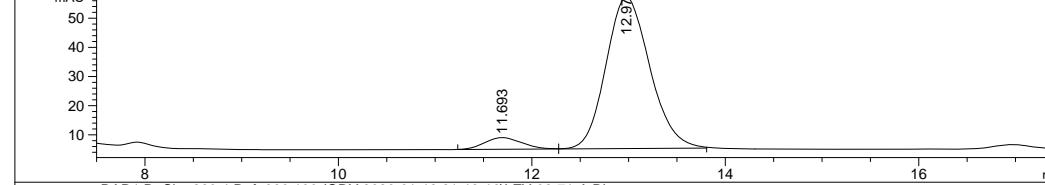
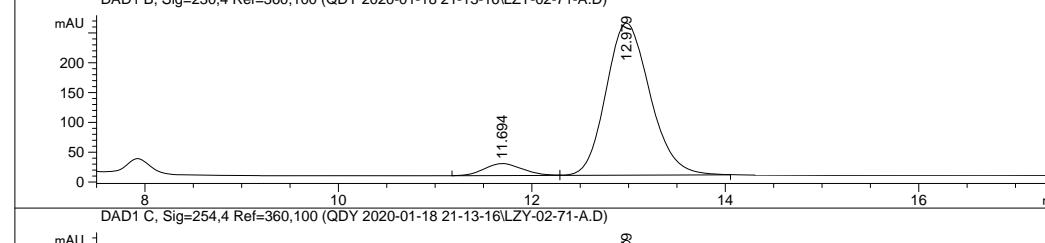
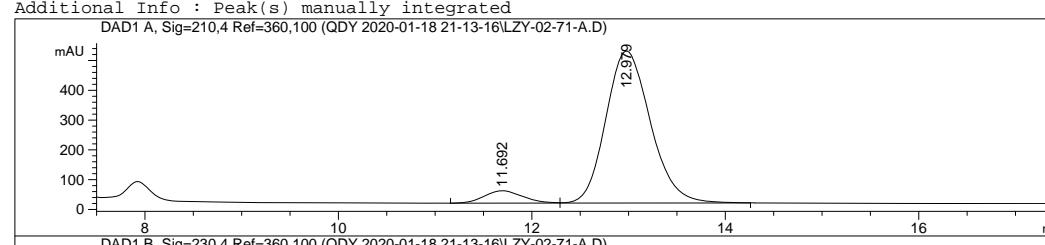
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 3
Acq. Instrument : Instrument 1             Location : Vial 94
Injection Date : 1/18/2020 9:57:32 PM       Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-01-18 21-13-16\IC-30-30.M
Last changed : 1/18/2020 9:56:41 PM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:15:03 PM        (modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.692	BV	0.4255	1137.97705	41.55007	6.5206
2	12.979	VB	0.4973	1.63141e4	509.80856	93.4794
Totals :					1.74521e4	551.35863

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

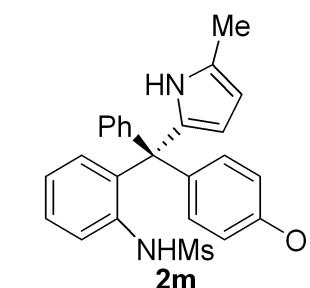
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.694	BV	0.4218	547.47168	20.09611	6.3694
2	12.979	VB	0.4884	8047.87207	254.83559	93.6306
Totals :					8595.34375	274.93169

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

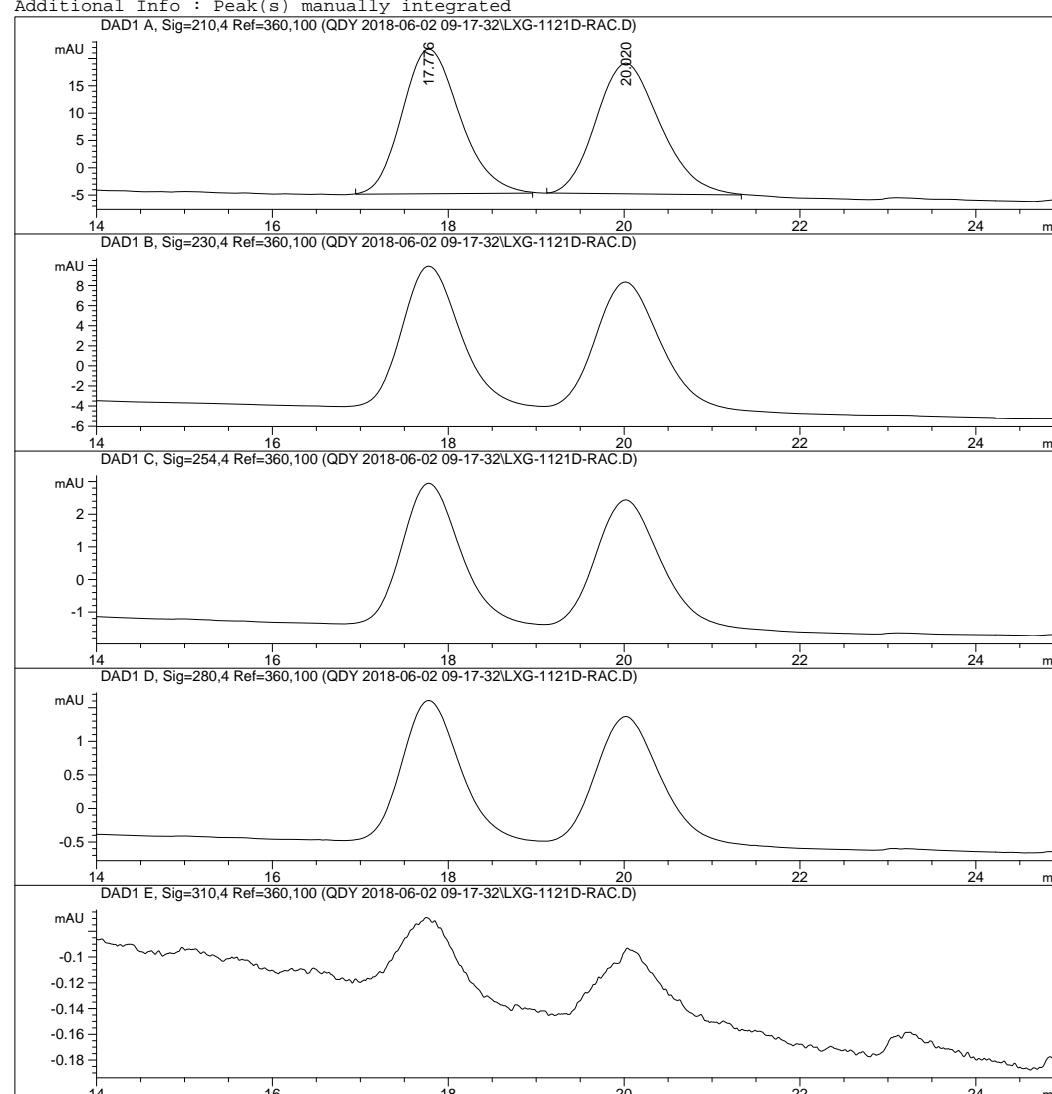
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.693	BV	0.4161	105.81114	4.00634	6.0913
2	12.979	VB	0.4868	1631.28723	51.89074	93.9087
Totals :					1737.09837	55.89708

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



```
=====
Acq. Operator :                               Seq. Line : 6
Acq. Instrument : Instrument 1             Location : Vial 51
Injection Date : 6/2/2018 10:41:01 AM        Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !       Actual Inj Volume : 2.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2018-06-02 09-17-32\IC-15-30.M
Last changed : 6/2/2018 10:40:10 AM          (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:25:15 PM          (modified after loading)
Additional Info : Peak(s) manually integrated
```



```
=====
Area Percent Report
=====
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100
```

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	17.776	BB	0.7033	1210.11121	26.47231	50.0003
2	20.020	BB	0.7725	1210.09485	23.90942	49.9997

Totals : 2420.20605 50.38173

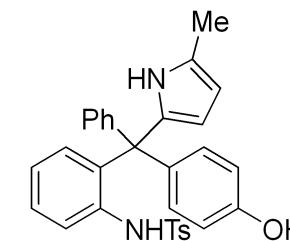
Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

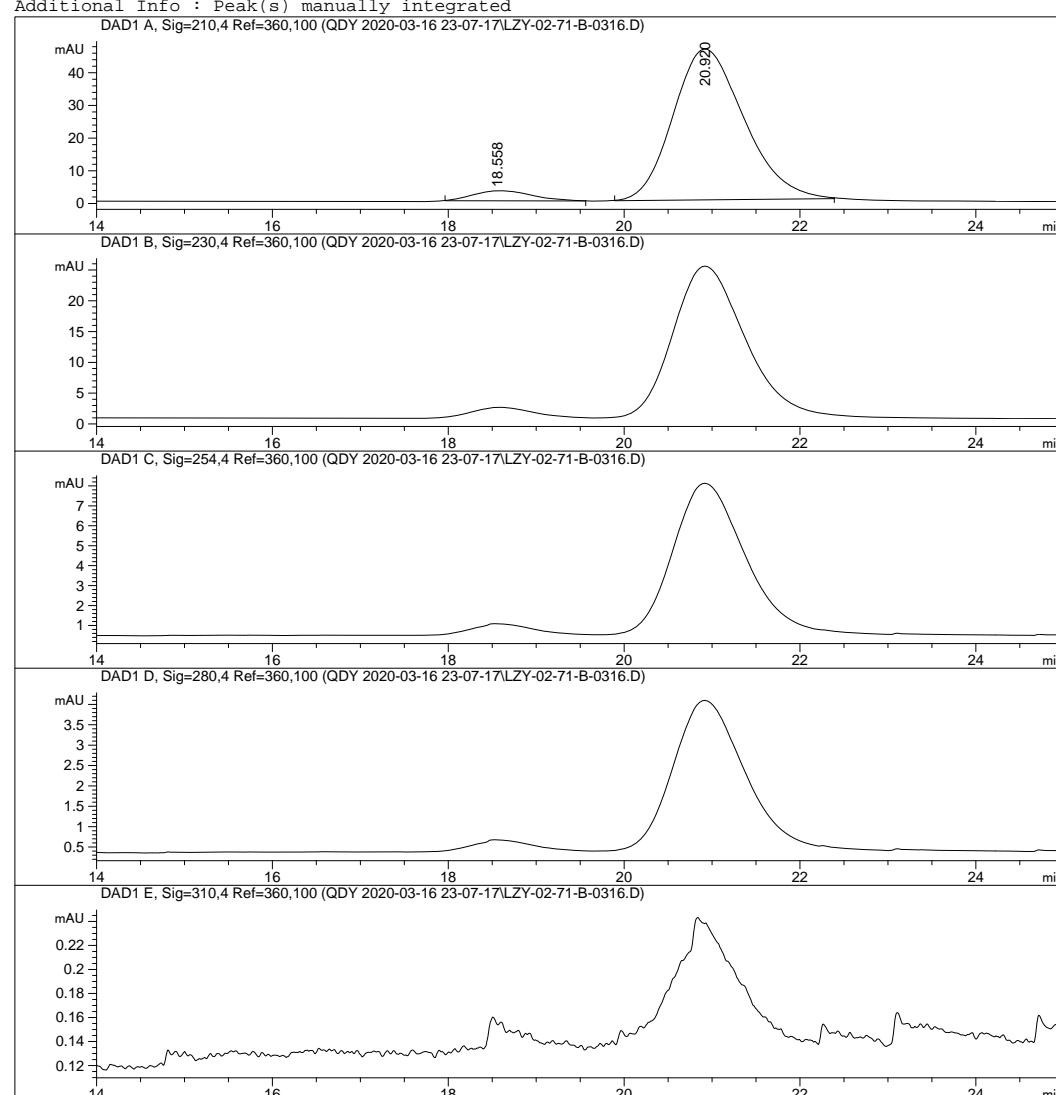
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100

===== *** End of Report *** =====

**2n-Rac**

```
=====
Acq. Operator :                               Seq. Line : 9
Acq. Instrument : Instrument 1             Location : Vial 93
Injection Date : 3/17/2020 1:08:41 AM        Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 8.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-03-16 23-07-17\IC-15-30.M
Last changed : 3/17/2020 1:07:47 AM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:25:15 PM        (modified after loading)
Additional Info : Peak(s) manually integrated
```

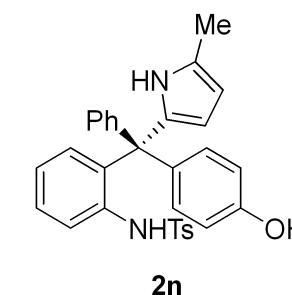


```
=====
Area Percent Report
=====
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100
Peak RetTime Type Width Area Height Area
# [min] [min] [mAU*s] [mAU] %
-----|-----|-----|-----|-----|
1 18.558 BB 0.5665 143.21919 3.06842 5.2412
2 20.920 BB 0.8648 2589.35400 46.13736 94.7588

Totals : 2732.57320 49.20577

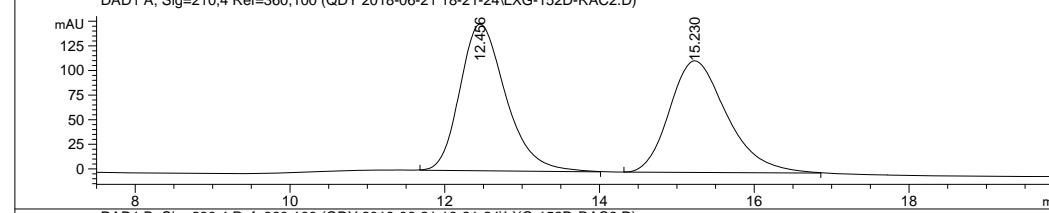
Signal 2: DAD1 B, Sig=230,4 Ref=360,100
Signal 3: DAD1 C, Sig=254,4 Ref=360,100
Signal 4: DAD1 D, Sig=280,4 Ref=360,100
Signal 5: DAD1 E, Sig=310,4 Ref=360,100
=====
```

**2n**

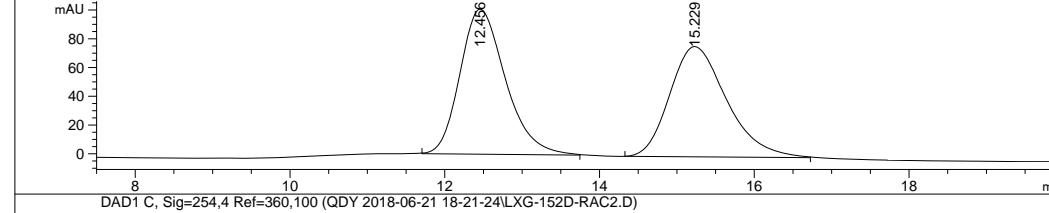
Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 2
Acq. Instrument : Instrument 1             Location : Vial 67
Injection Date : 6/21/2018 6:34:26 PM       Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2018-06-21 18-21-24\IC-30-30.M
Last changed : 6/21/2018 6:33:33 PM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:33:55 PM        (modified after loading)
Additional Info : Peak(s) manually integrated
```

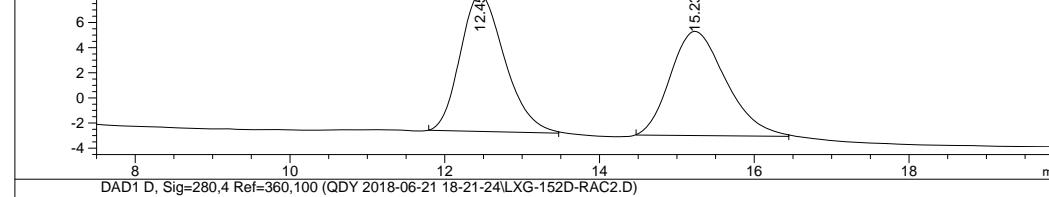
DAD1 A, Sig=210,4 Ref=360,100 (QDY 2018-06-21 18-21-24\LXG-152D-RAC2.D)



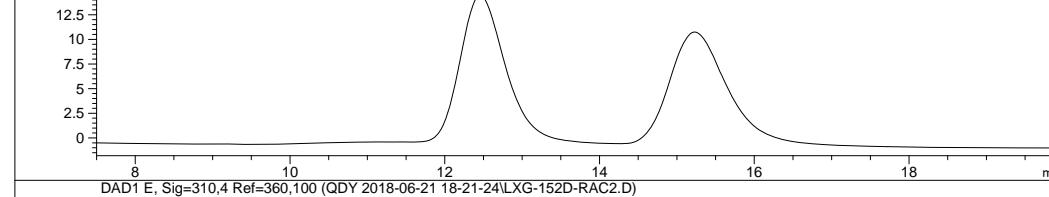
DAD1 B, Sig=230,4 Ref=360,100 (QDY 2018-06-21 18-21-24\LXG-152D-RAC2.D)



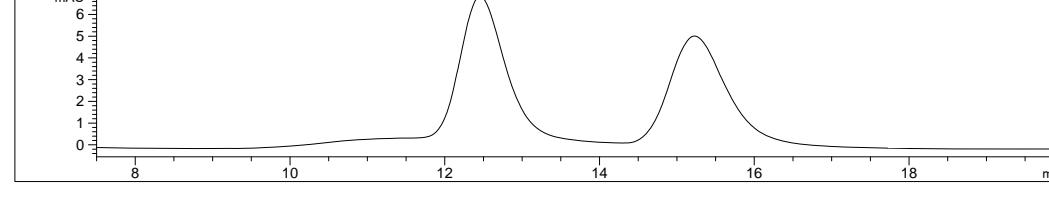
DAD1 C, Sig=254,4 Ref=360,100 (QDY 2018-06-21 18-21-24\LXG-152D-RAC2.D)



DAD1 D, Sig=280,4 Ref=360,100 (QDY 2018-06-21 18-21-24\LXG-152D-RAC2.D)



DAD1 E, Sig=310,4 Ref=360,100 (QDY 2018-06-21 18-21-24\LXG-152D-RAC2.D)



Data File C:\CHEM32\1\DATA\QDY 2018-06-21 18-21-24\LXG-152D-RAC2.D
Sample Name:

===== Area Percent Report =====

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.456	BB	0.6317	6126.23535	149.00308	51.1842
2	15.230	BB	0.7913	5842.75342	113.36085	48.8158

Totals : 1.19690e4 262.36393

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.456	BB	0.6277	4113.04004	100.86504	51.1741
2	15.229	BB	0.7851	3924.30298	76.67237	48.8259

Totals : 8037.34302 177.53741

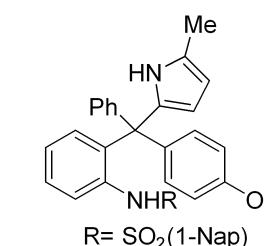
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.455	BB	0.6158	433.83267	10.91284	51.1764
2	15.231	BB	0.7643	413.88727	8.29309	48.8236

Totals : 847.71994 19.20593

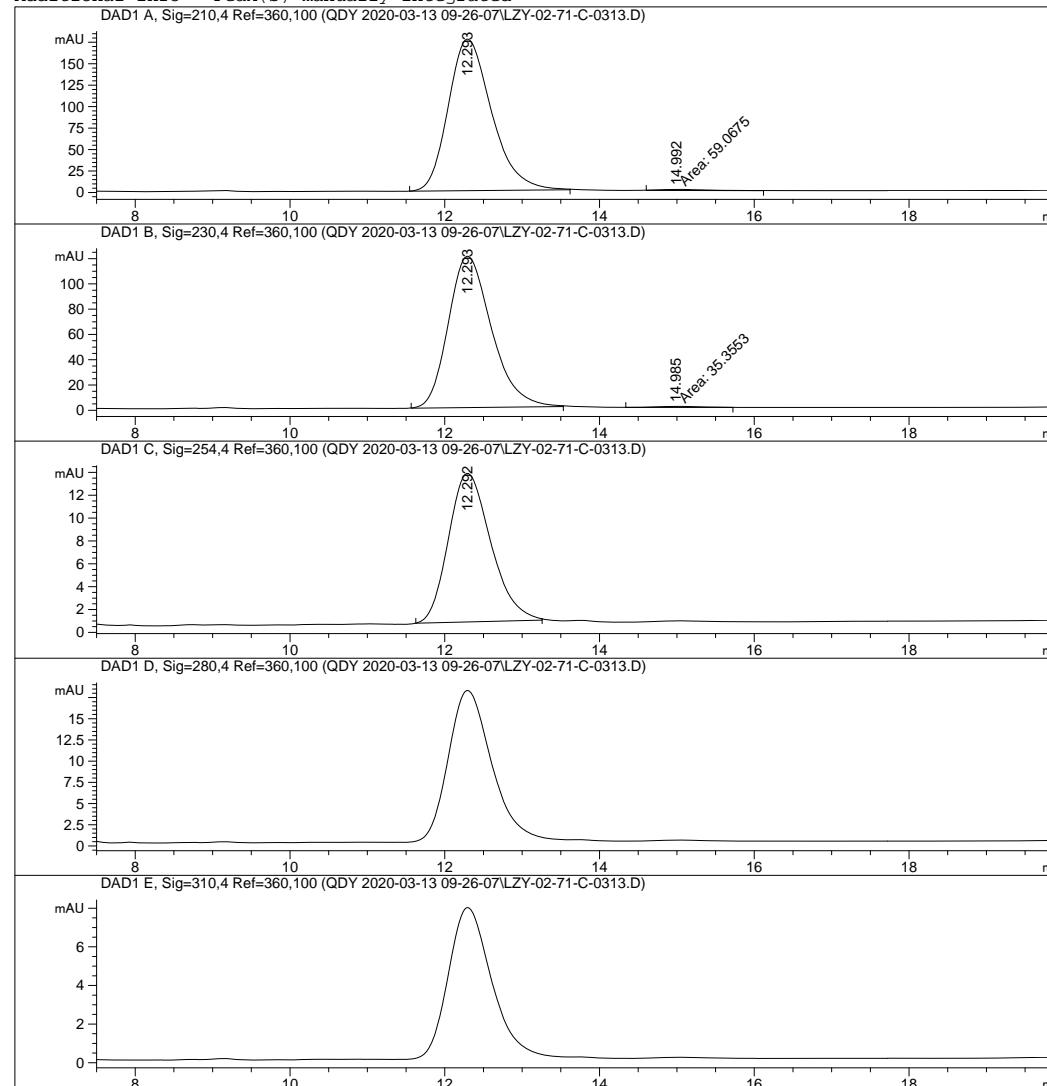
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

=====
 Acq. Operator : Seq. Line : 23
 Acq. Instrument : Instrument 1 Location : Vial 84
 Injection Date : 3/13/2020 4:34:57 PM Inj : 1
 Inj Volume : 5.000 μ l
 Different Inj Volume from Sequence ! Actual Inj Volume : 1.000 μ l
 Acq. Method : C:\CHEM32\1\DATA\QDY 2020-03-13 09-26-07\IC-30-20.M
 Last changed : 3/13/2020 9:34:34 AM
 Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
 Last changed : 11/24/2020 8:39:23 PM
 (modified after loading)
 Additional Info : Peak(s) manually integrated



Data File C:\CHEM32\1\DATA\QDY 2020-03-13 09-26-07\LZY-02-71-C-0313.D
Sample Name:

=====
 Area Percent Report
 =====

Sorted By : Signal
 Multiplier : 1.0000
 Dilution : 1.0000
 Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.293	BB	0.5978	6870.16016	176.63481	99.1476
2	14.992	MM	0.7524	59.06746	1.30838	0.8524

Totals : 6929.22762 177.94319

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.293	BB	0.5963	4645.24121	119.82249	99.2446
2	14.985	MM	0.7220	35.35526	8.16122e-1	0.7554

Totals : 4680.59647 120.63861

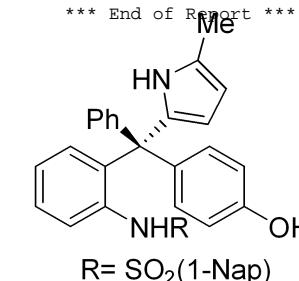
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.292	BB	0.5851	493.85947	13.00298	100.0000

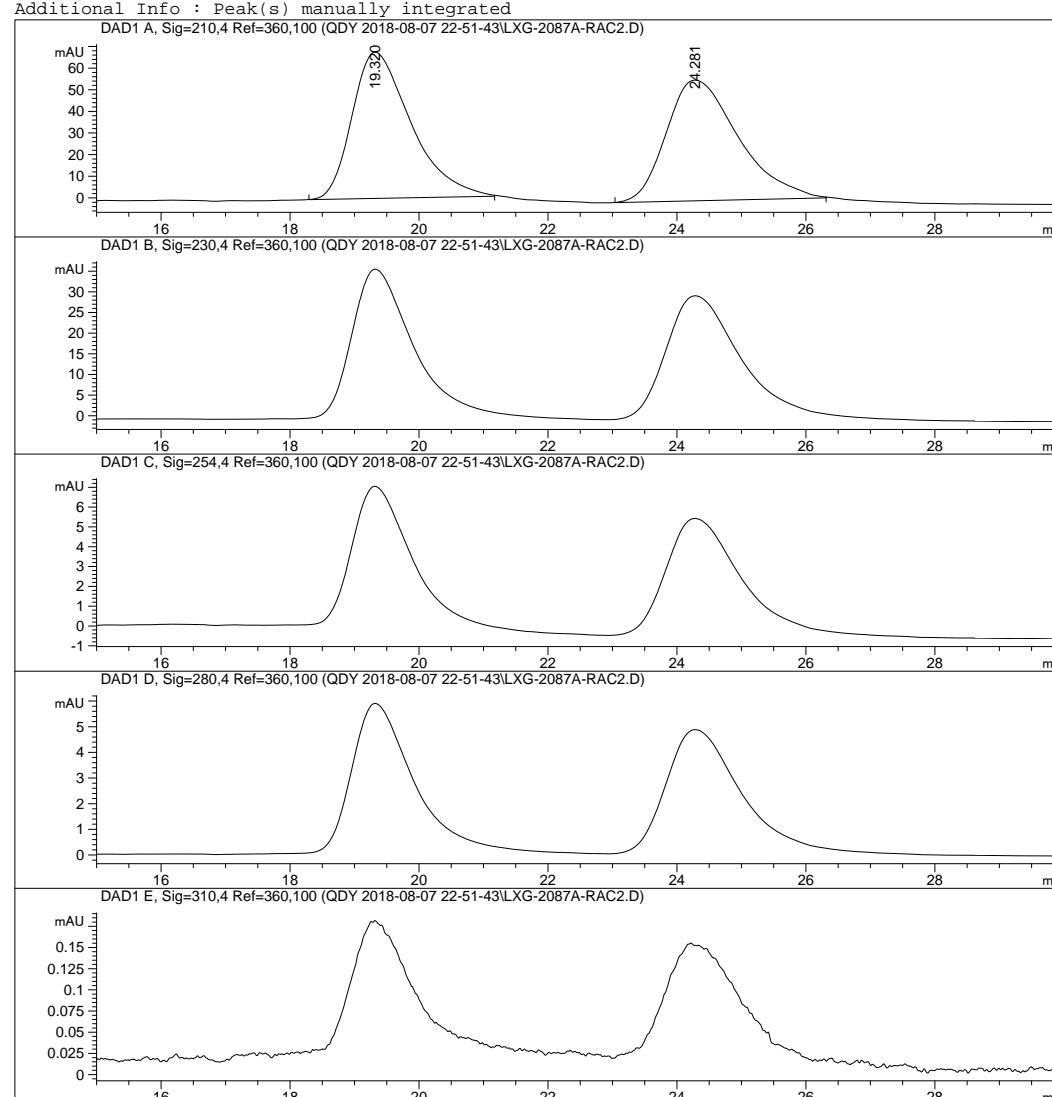
Totals : 493.85947 13.00298

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



```
=====
Acq. Operator :                               Seq. Line : 6
Acq. Instrument : Instrument 1             Location : Vial 68
Injection Date : 8/8/2018 12:12:24 AM       Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2018-08-07 22-51-43\OD-05-30.M
Last changed : 8/8/2018 12:11:31 AM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:37:01 PM        (modified after loading)
Additional Info : Peak(s) manually integrated
```



```
=====
Area Percent Report
=====
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100
```

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	19.320	BB	0.9428	4283.52246	67.63930	49.7465
2	24.281	BB	1.1572	4327.17969	55.77165	50.2535

Totals : 8610.70215 123.41095

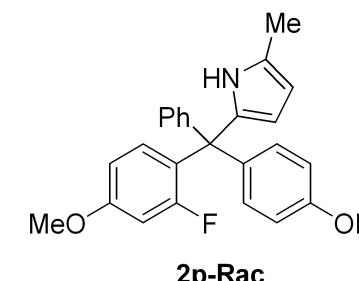
Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

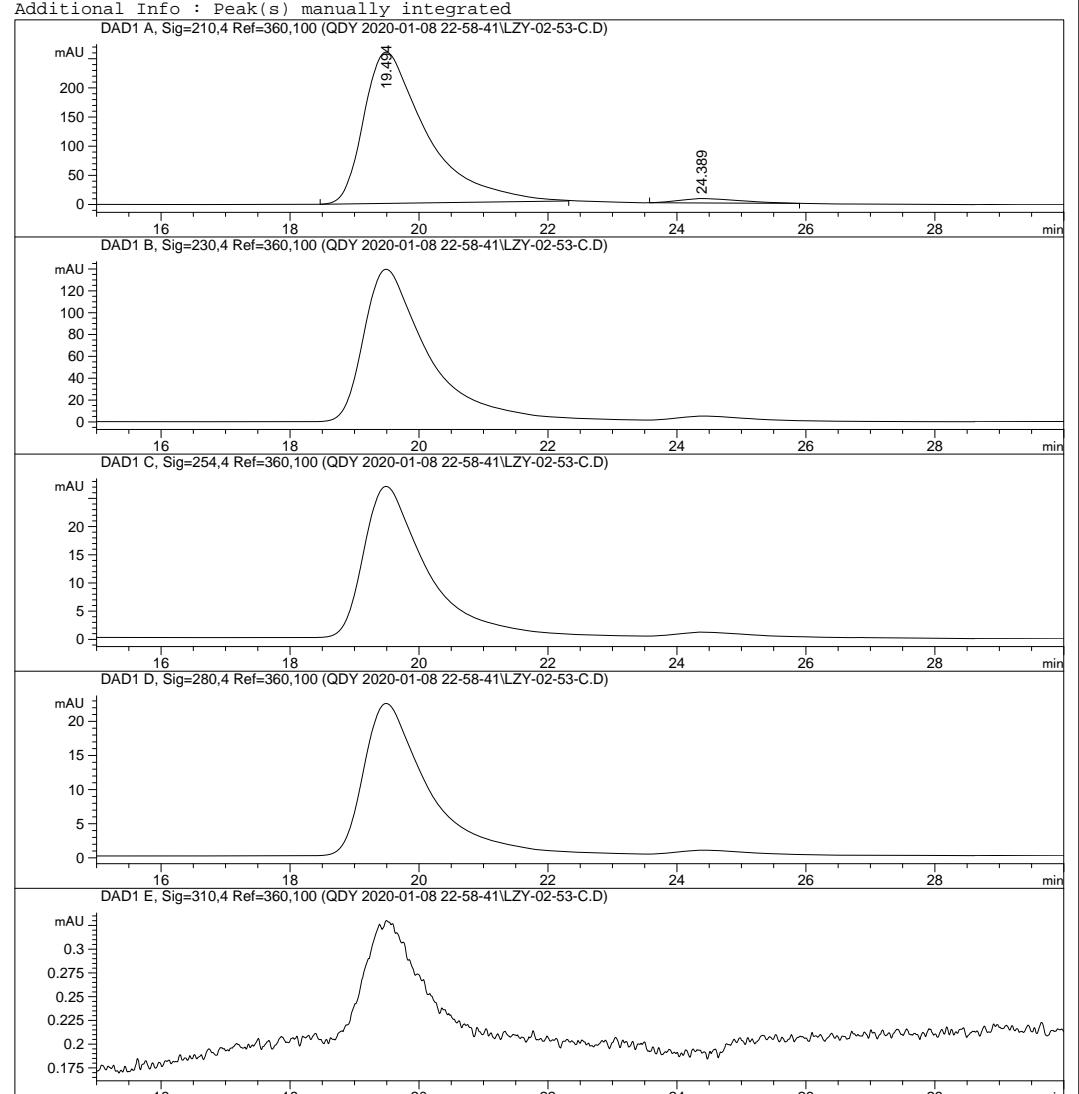
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100

=====
*** End of Report ***



```
=====
Acq. Operator :                               Seq. Line : 15
Acq. Instrument : Instrument 1             Location : Vial 96
Injection Date : 1/9/2020 2:36:31 AM        Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 4.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-01-08 22-58-41\OD-05-30.M
Last changed : 1/9/2020 2:35:39 AM          (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:37:01 PM          (modified after loading)
Additional Info : Peak(s) manually integrated
```



```
=====
Area Percent Report
=====
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100
Peak RetTime Type Width Area Height Area
# [min] [min] [mAU*s] [mAU] %
-----|-----|-----|-----|-----|-----|
1 19.494 BB 0.9786 1.75318e4 259.38416 97.2576
2 24.389 BB 0.7989 494.34113 7.57937 2.7424

Totals : 1.80262e4 266.96353
```

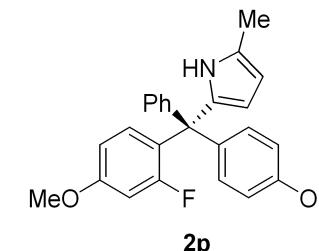
Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

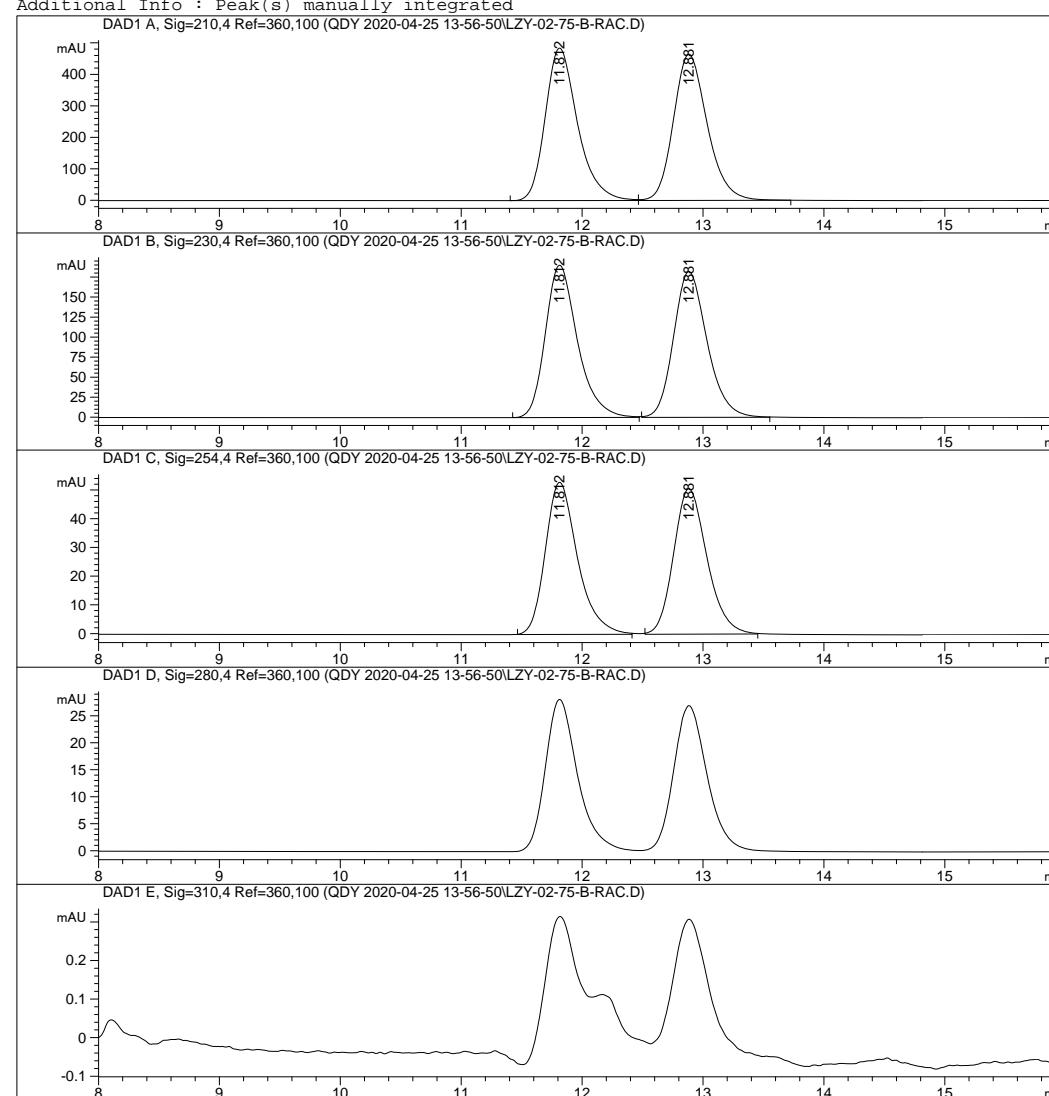
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100

=====
*** End of Report ***



```
=====
Acq. Operator : Seq. Line : 25
Acq. Instrument : Instrument 1 Location : Vial 81
Injection Date : 4/25/2020 9:18:37 PM Inj : 1
Inj Volume : 5.000 µl
Different Inj Volume from Sequence ! Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-04-25 13-56-50\AD-05-20.M
Last changed : 4/25/2020 9:17:45 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 2:55:56 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.812	BV	0.2826	8987.99902	484.12137	49.9256
2	12.881	VB	0.2985	9014.78516	464.31113	50.0744

Totals : 1.80028e4 948.43250

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.812	BB	0.2794	3512.41675	190.20287	50.0723
2	12.881	BB	0.2943	3502.27344	182.13647	49.9277

Totals : 7014.69019 372.33934

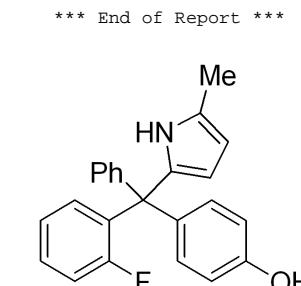
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.812	BB	0.2818	979.65533	52.96997	50.2141
2	12.881	BB	0.2935	971.30011	50.69146	49.7859

Totals : 1950.95544 103.66143

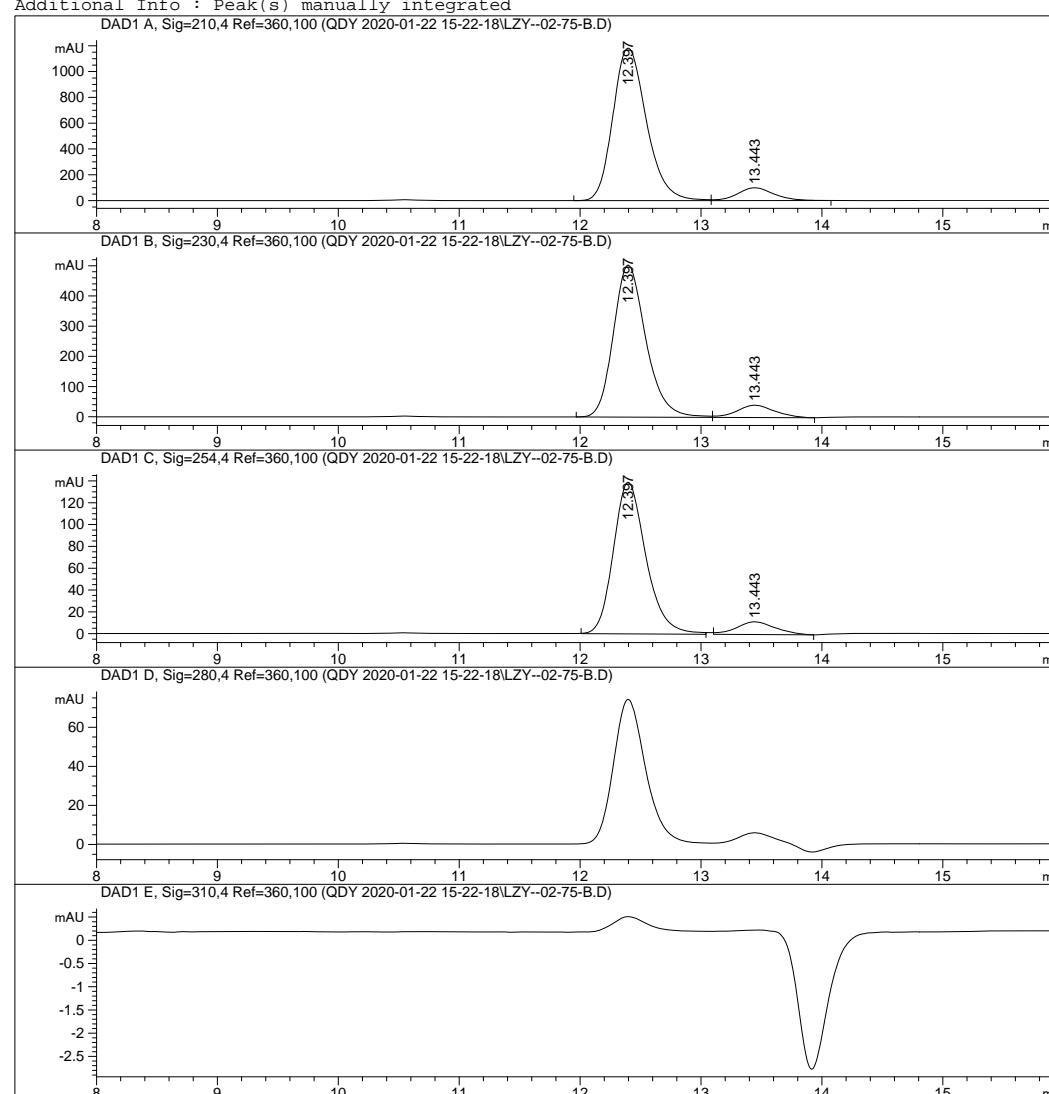
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 3
Acq. Instrument : Instrument 1             Location : Vial 88
Injection Date : 1/22/2020 3:56:25 PM       Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 2.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-01-22 15-22-18\AD-05-20.M
Last changed : 1/22/2020 3:34:31 PM
                                                (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 2:55:56 PM
                                                (modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.397	BV	0.2966	2.27745e4	1182.88330	91.6953
2	13.443	VB	0.3183	2062.66333	99.32850	8.3047

Totals : 2.48372e4 1282.21180

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.397	BB	0.2860	9405.17676	503.36429	91.3841
2	13.443	BV	0.3262	886.74402	41.02750	8.6159

Totals : 1.02919e4 544.39179

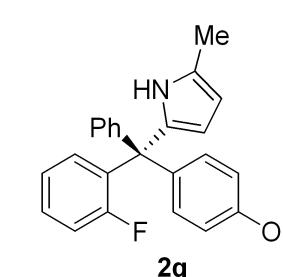
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.397	BB	0.2861	2597.93774	138.97246	90.8687
2	13.443	BV	0.3361	261.06342	11.71148	9.1313

Totals : 2859.00116 150.68394

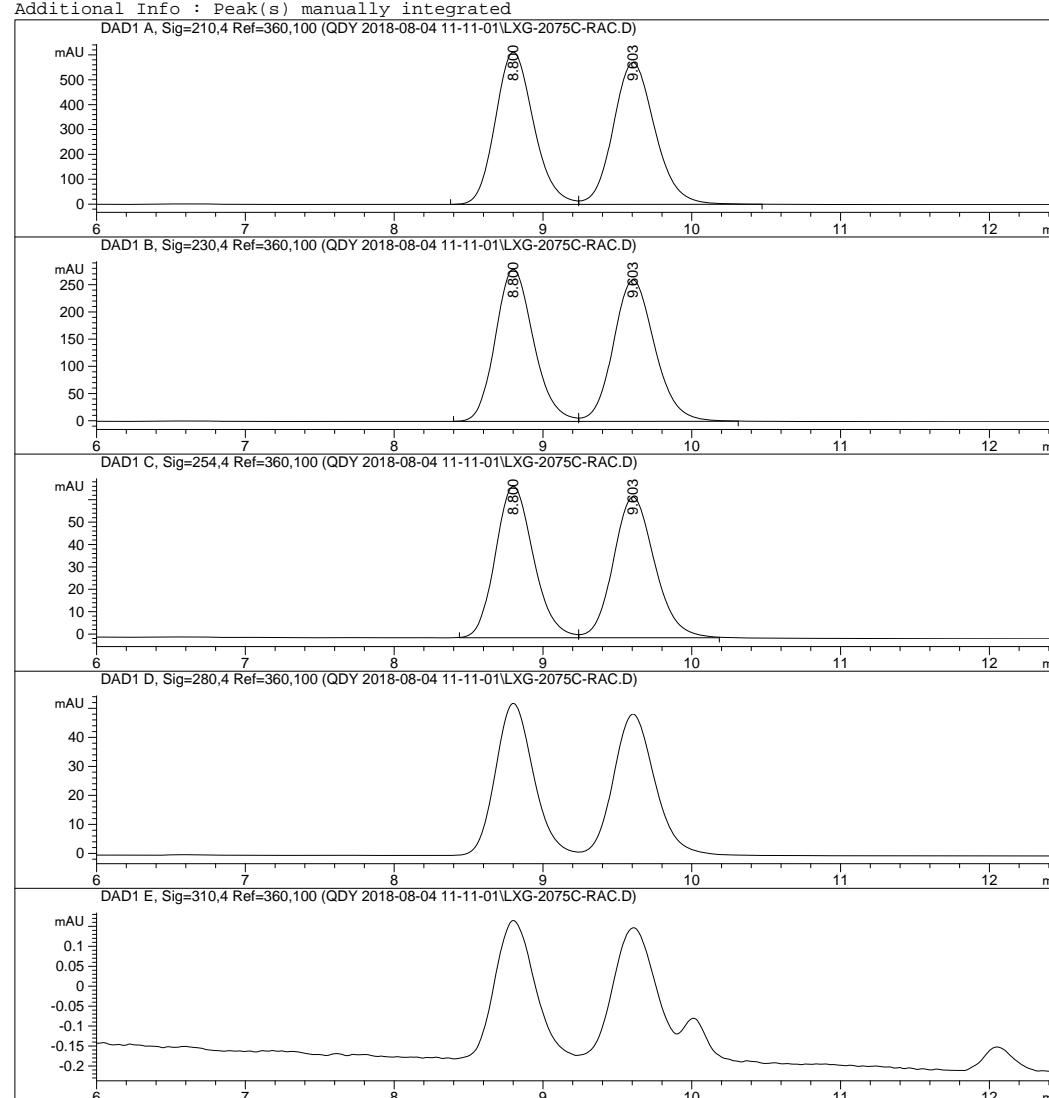
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

```
=====
Acq. Operator : Seq. Line : 2
Acq. Instrument : Instrument 1 Location : Vial 66
Injection Date : 8/4/2018 11:24:41 AM Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence ! Actual Inj Volume : 2.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2018-08-04 11-11-01\IC-03-20.M
Last changed : 8/4/2018 11:23:49 AM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:40:50 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



Data File C:\CHEM32\1\DATA\QDY 2018-08-04 11-11-01\LXG-2075C-RAC.D
Sample Name:

```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.800	BV	0.2782	1.09314e4	612.50708	49.6216
2	9.603	VB	0.2988	1.10981e4	570.79449	50.3784

Totals : 2.20296e4 1183.30157

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.800	BV	0.2749	4960.99951	279.71667	49.7215
2	9.603	VB	0.2969	5016.57324	260.20297	50.2785

Totals : 9977.57275 539.91965

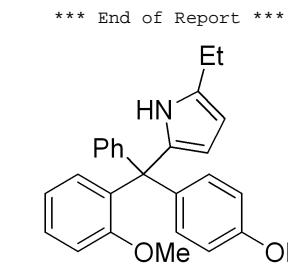
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.800	BV	0.2747	1199.28052	67.67670	49.7427
2	9.603	VB	0.2965	1211.68494	62.96595	50.2573

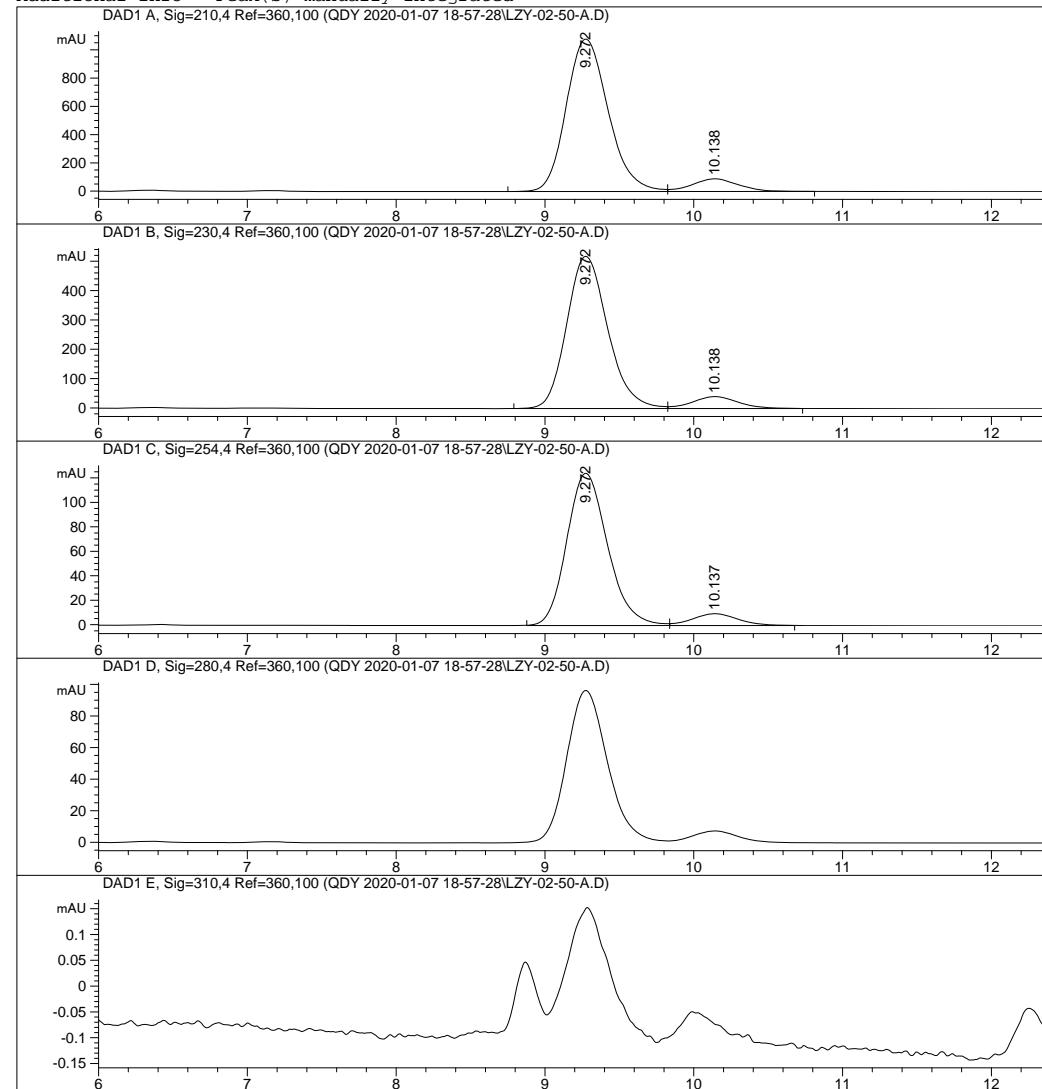
Totals : 2410.96545 130.64266

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



```
=====
Acq. Operator : Seq. Line : 8
Acq. Instrument : Instrument 1 Location : Vial 91
Injection Date : 1/7/2020 10:09:07 PM Inj : 1
Inj Volume : 5.000 µl
Different Inj Volume from Sequence ! Actual Inj Volume : 2.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-01-07 18-57-28\IC-03-20.M
Last changed : 9/18/2012 8:18:47 PM
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:40:50 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.272	BV	0.3058	2.1064e4	1078.78027	91.6011
2	10.138	VB	0.3294	1931.35620	89.62775	8.3989

Totals : 2.29954e4 1168.40803

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.272	BV	0.2968	9913.86035	519.22614	91.9132
2	10.138	VB	0.3283	872.24756	40.65038	8.0868

Totals : 1.07861e4 559.87651

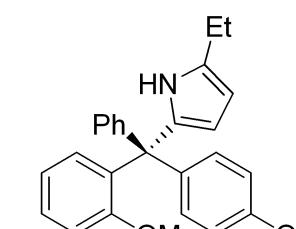
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.272	BB	0.2941	2369.02222	124.46080	91.9652
2	10.137	BB	0.3247	206.97551	9.70764	8.0348

Totals : 2575.99773 134.16844

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

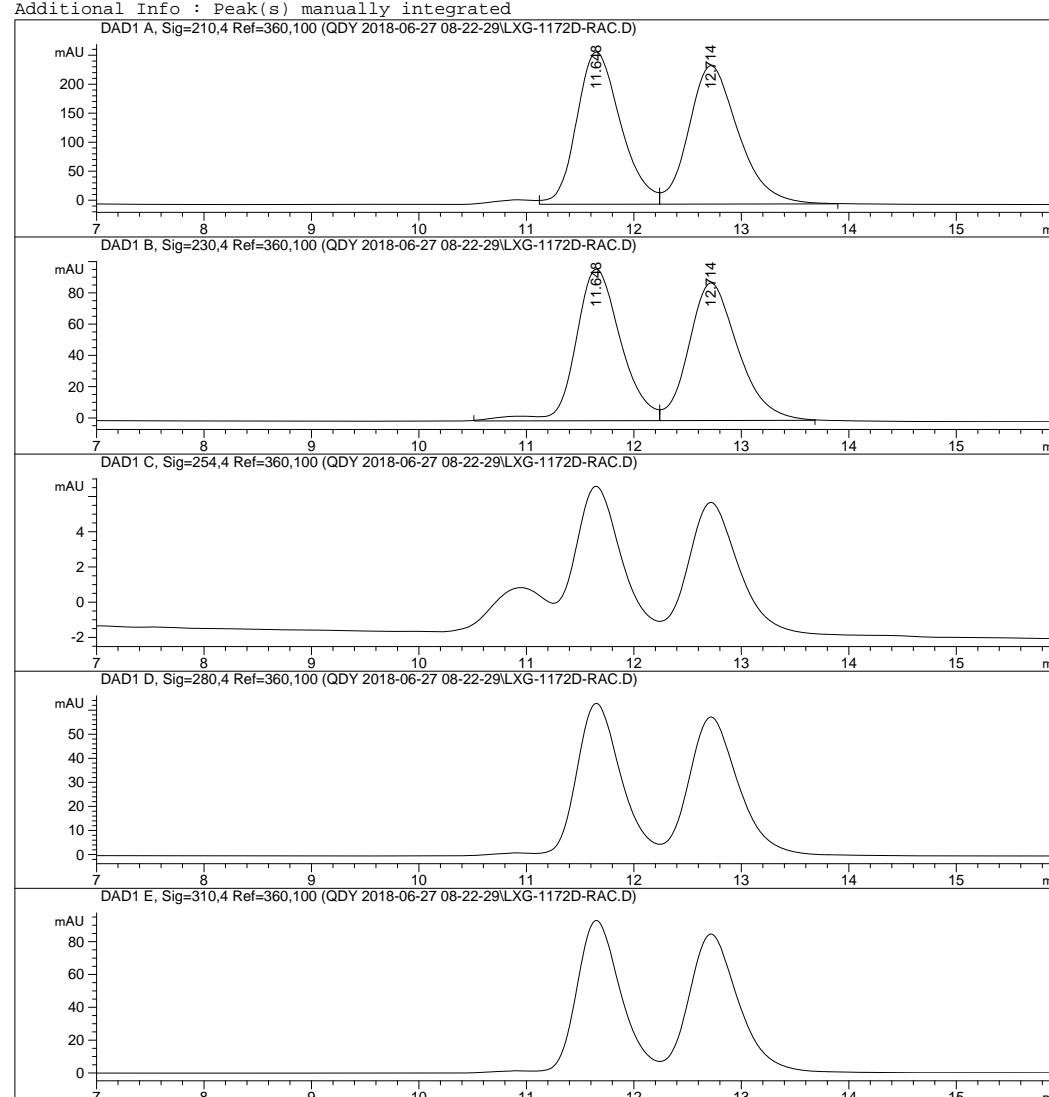
Signal 5: DAD1 E, Sig=310,4 Ref=360,100



2r

Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 2
Acq. Instrument : Instrument 1             Location : Vial 63
Injection Date : 6/27/2018 8:36:04 AM        Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 2.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2018-06-27 08-22-29\OD-05-20.M
Last changed : 6/27/2018 8:35:12 AM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:43:47 PM        (modified after loading)
Additional Info : Peak(s) manually integrated
```



Data File C:\CHEM32\1\DATA\QDY 2018-06-27 08-22-29\LXG-1172D-RAC.D
Sample Name:

```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.648	VV	0.4399	7463.65869	262.36359	49.4228
2	12.714	VB	0.4895	7637.99951	238.55977	50.5772

Totals : 1.51017e4 500.92336

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	11.648	BV	0.4462	2826.79395	96.93911	50.3677
2	12.714	VB	0.4854	2785.52222	87.98379	49.6323

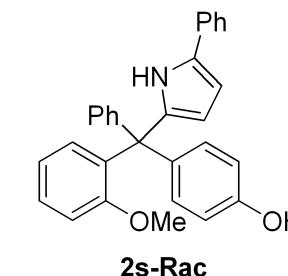
Totals : 5612.31616 184.92290

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

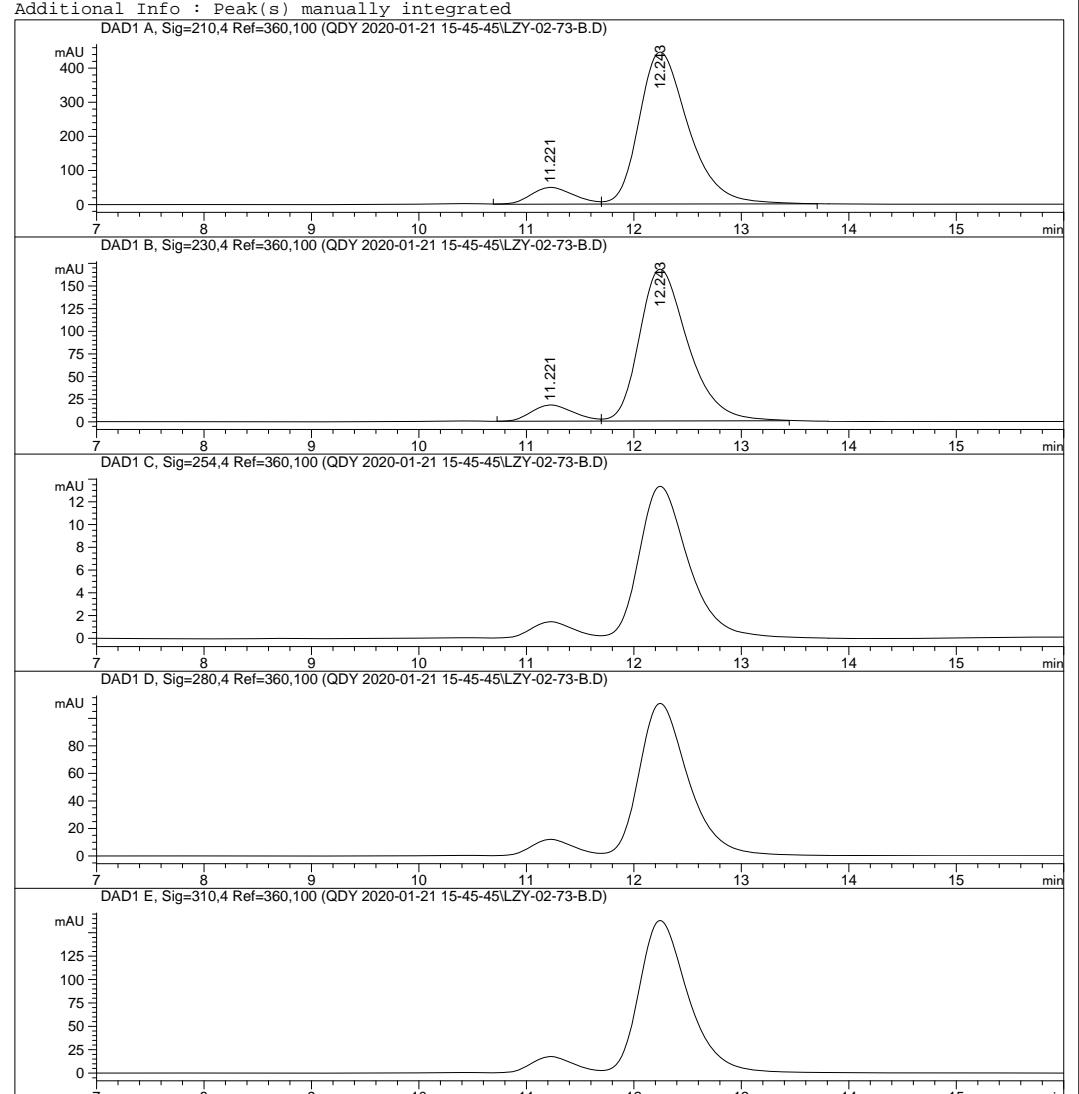
Signal 5: DAD1 E, Sig=310,4 Ref=360,100

===== *** End of Report ***



Sample Name:

```
=====
Acq. Operator : Seq. Line : 17
Acq. Instrument : Instrument 1 Location : Vial 85
Injection Date : 1/21/2020 9:47:23 PM Inj : 1
Inj Volume : 5.000 µl
Different Inj Volume from Sequence ! Actual Inj Volume : 2.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-01-21 15-45-45\OD-05-20.M
Last changed : 1/21/2020 9:46:30 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:43:47 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

```
=====
Area Percent Report
=====
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs

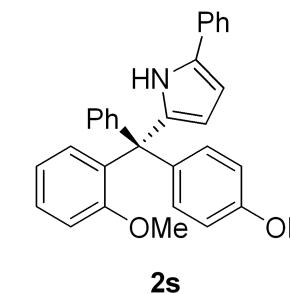
Signal 1: DAD1 A, Sig=210,4 Ref=360,100
Peak RetTime Type Width Area Height Area
# [min] [min] [mAU*s] [mAU] %
-----|-----|-----|-----|-----|-----|
1 11.221 VV 0.4256 1362.21912 49.41045 8.7555
2 12.243 VB 0.4876 1.41963e4 445.65451 91.2445

Totals : 1.55585e4 495.06496

Signal 2: DAD1 B, Sig=230,4 Ref=360,100
Peak RetTime Type Width Area Height Area
# [min] [min] [mAU*s] [mAU] %
-----|-----|-----|-----|-----|
1 11.221 BV 0.4203 484.59872 17.99050 8.4364
2 12.243 VB 0.4800 5259.56787 167.64790 91.5636

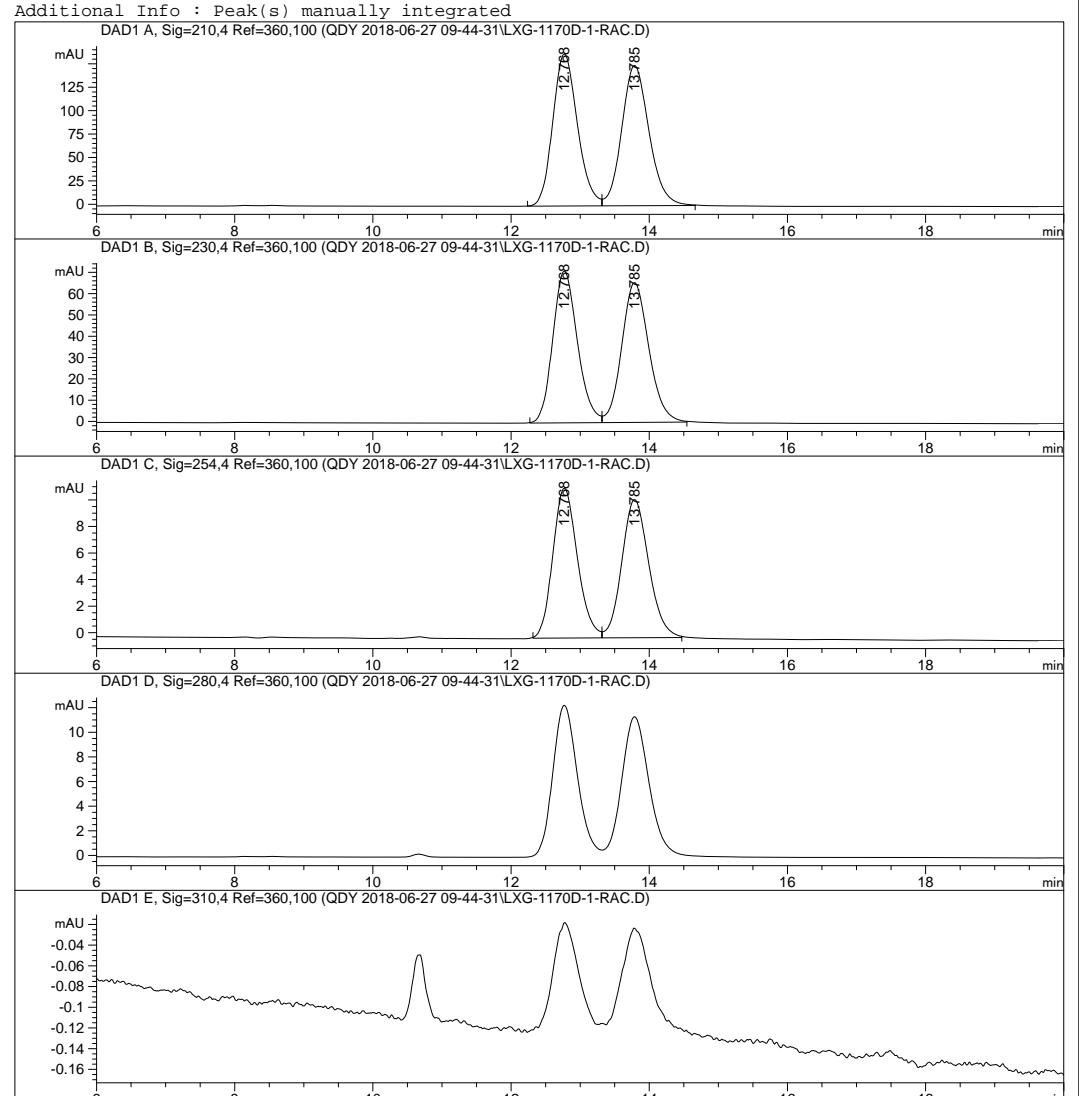
Totals : 5744.16660 185.63840

Signal 3: DAD1 C, Sig=254,4 Ref=360,100
Signal 4: DAD1 D, Sig=280,4 Ref=360,100
Signal 5: DAD1 E, Sig=310,4 Ref=360,100
=====
```



Sample Name:

```
=====
Acq. Operator : Seq. Line : 45
Acq. Instrument : Instrument 1 Location : Vial 56
Injection Date : 6/27/2018 11:02:49 PM Inj : 1
Inj Volume : 5.000 µl
Different Inj Volume from Sequence ! Actual Inj Volume : 2.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2018-06-27 09-44-31\IC-02-30.M
Last changed : 6/27/2018 11:01:56 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 9:08:48 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



Data File C:\CHEM32\1\DATA\QDY 2018-06-27 09-44-31\LXG-1170D-1-RAC.D
Sample Name:

```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.768	BV	0.3839	4035.83545	162.42874	49.5455
2	13.785	VB	0.4220	4109.88281	149.84027	50.4545

Totals : 8145.71826 312.26901

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.768	BV	0.3832	1765.27112	71.21311	49.7066
2	13.785	VB	0.4196	1786.11365	65.61869	50.2934

Totals : 3551.38477 136.83180

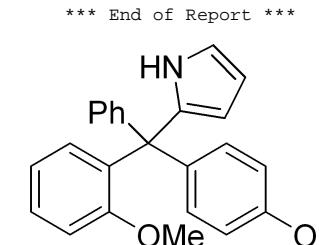
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	12.768	BV	0.3801	278.57065	11.27951	49.7152
2	13.785	VB	0.4180	281.76282	10.40222	50.2848

Totals : 560.33347 21.68172

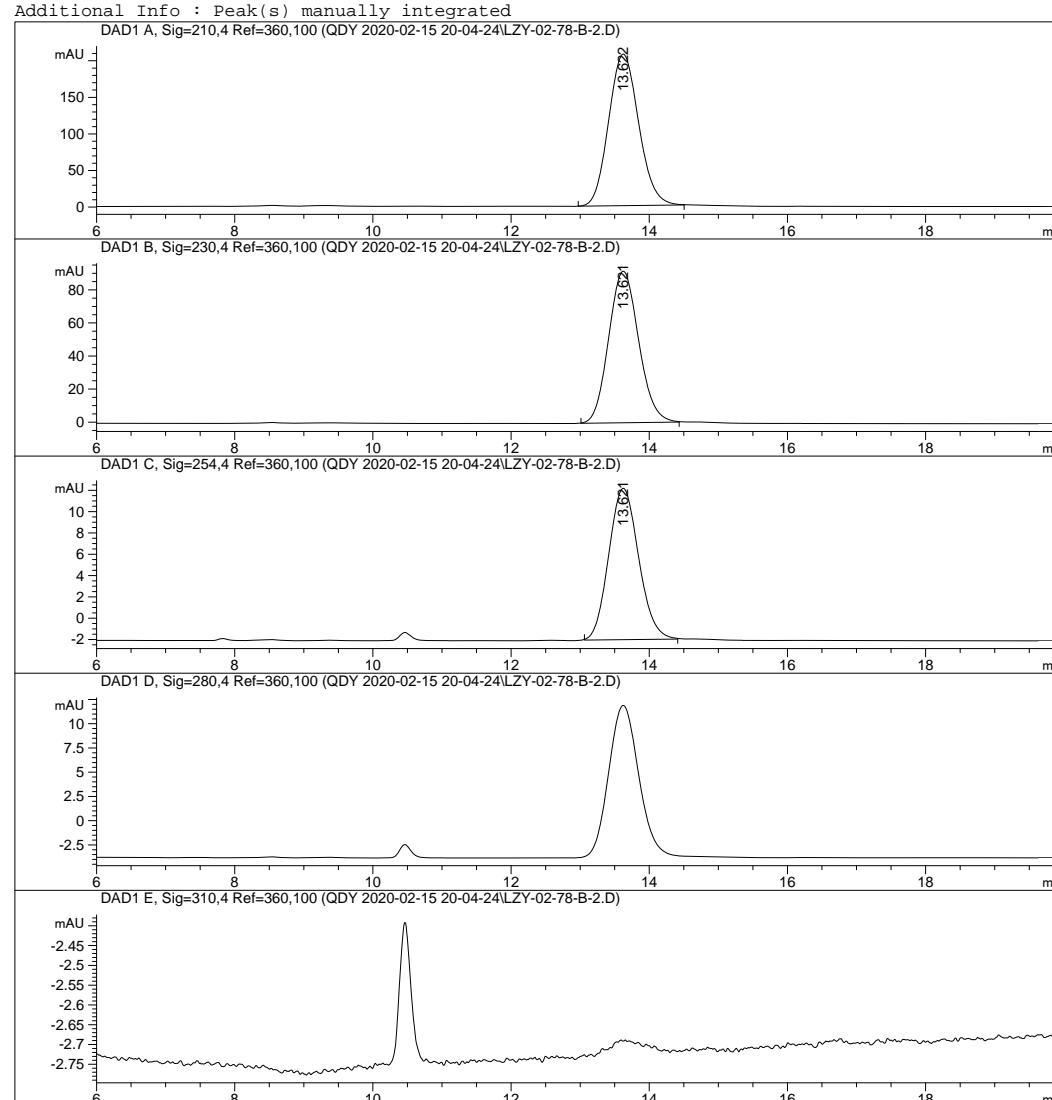
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 1
Acq. Instrument : Instrument 1             Location : Vial 94
Injection Date : 2/15/2020 8:06:16 PM        Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !       Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-02-15 20-04-24\IC-02-20.M
Last changed : 2/15/2020 8:04:36 PM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:49:19 PM        (modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

===== Area Percent Report =====

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	13.622	BB	0.4725	6260.75342	207.19859	100.0000

Totals : 6260.75342 207.19859

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	13.621	BB	0.4694	2761.62769	91.70020	100.0000

Totals : 2761.62769 91.70020

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

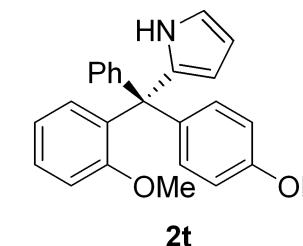
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	13.621	BB	0.4684	426.49472	14.19864	100.0000

Totals : 426.49472 14.19864

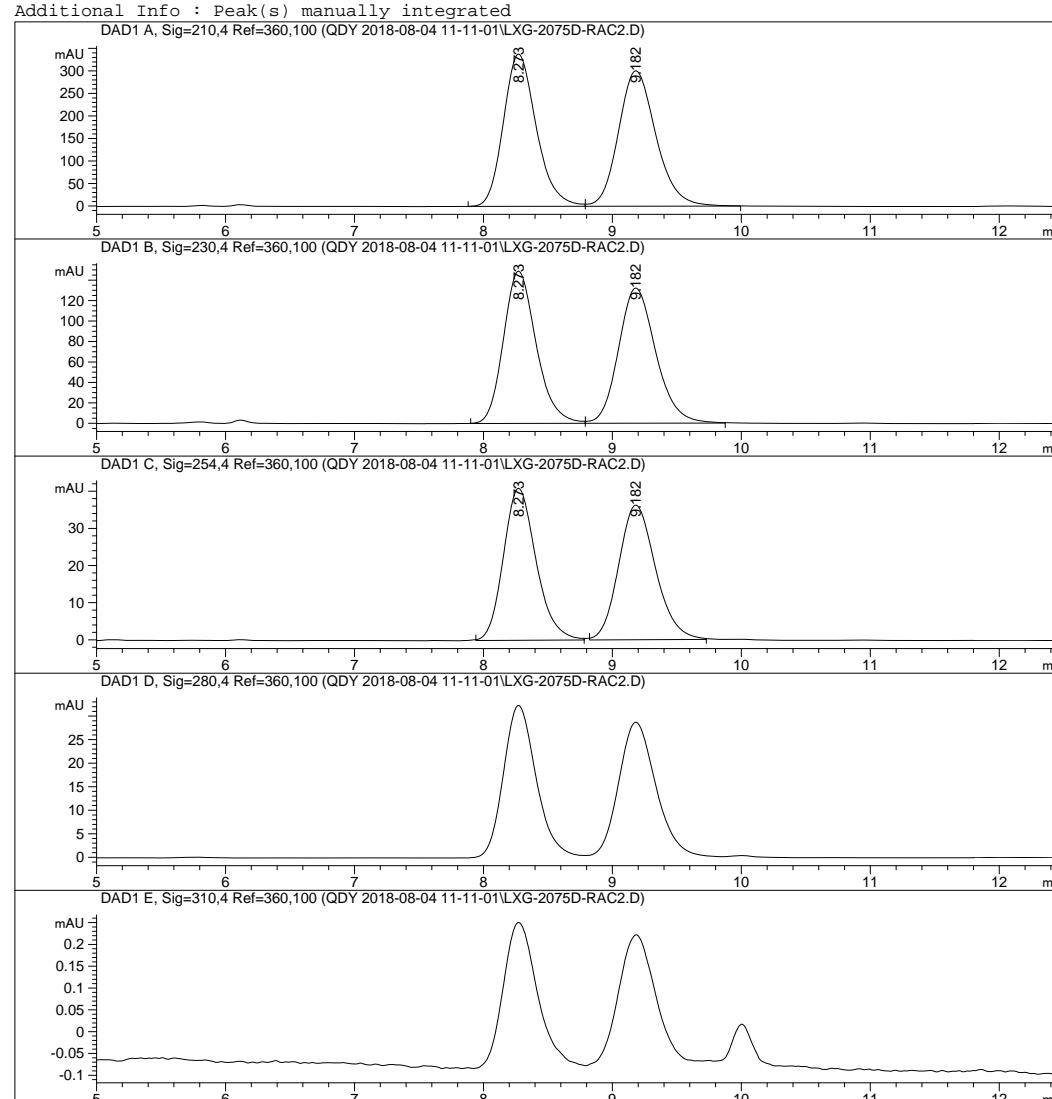
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100

=====
*** End of Report ***
=====



```
=====
Acq. Operator : Seq. Line : 7
Acq. Instrument : Instrument 1 Location : Vial 67
Injection Date : 8/4/2018 12:40:42 PM Inj : 1
Inj Volume : 5.000 µl
Different Inj Volume from Sequence ! Actual Inj Volume : 2.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2018-08-04 11-11-01\IC-03-30.M
Last changed : 8/4/2018 12:39:51 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:53:05 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.273	BV	0.2723	5981.02783	338.28845	49.6196
2	9.182	VB	0.3096	6072.72754	300.65823	50.3804

Totals : 1.20538e4 638.94669

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.273	BV	0.2696	2624.16455	148.91055	49.7284
2	9.182	VB	0.3082	2652.83350	132.15411	50.2716

Totals : 5276.99805 281.06467

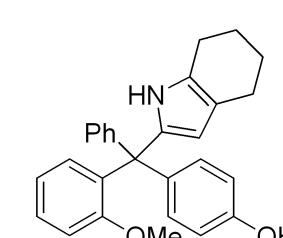
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.273	BB	0.2690	719.12695	40.92221	49.9726
2	9.182	BB	0.3057	719.91656	36.24767	50.0274

Totals : 1439.04352 77.16987

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

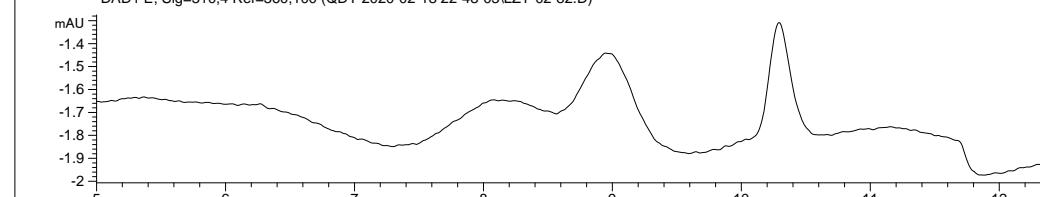
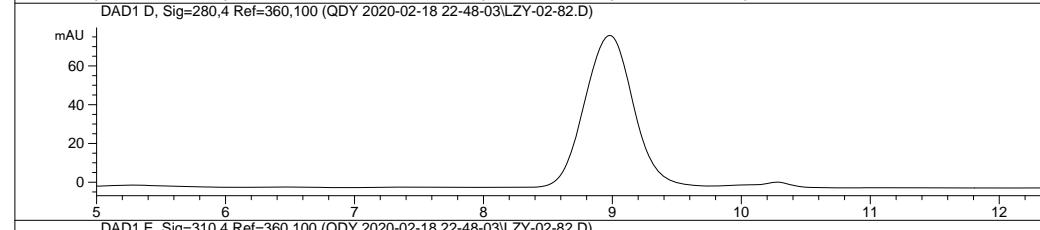
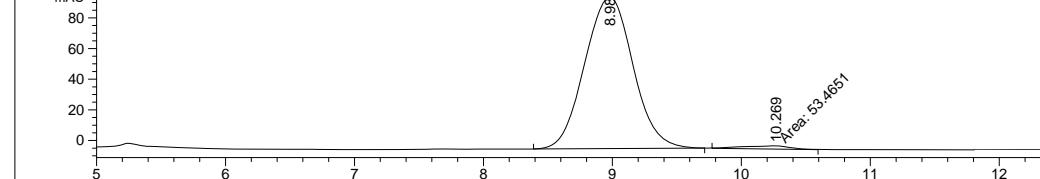
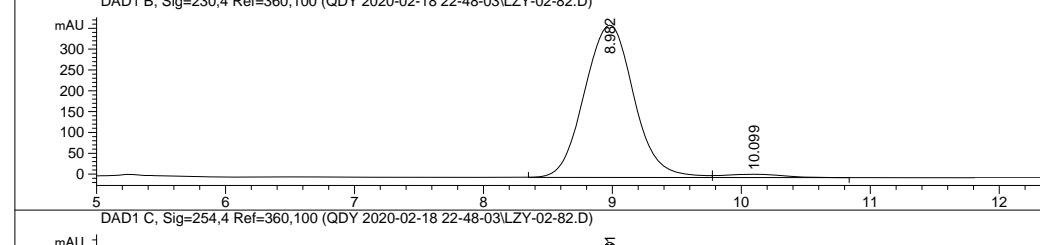
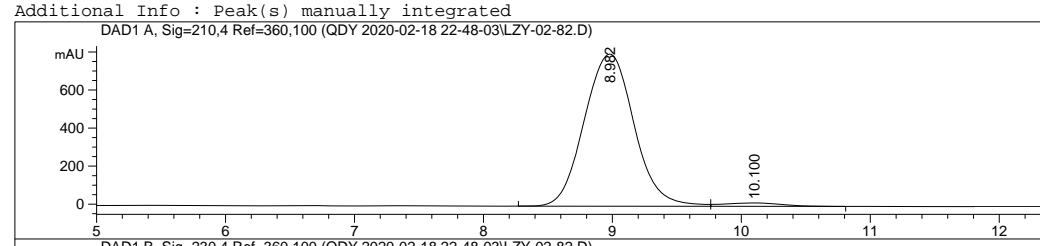
Signal 5: DAD1 E, Sig=310,4 Ref=360,100



2u-Rac

Sample Name:

```
=====
Acq. Operator : Seq. Line : 2
Acq. Instrument : Instrument 1 Location : Vial 81
Injection Date : 2/18/2020 11:01:16 PM Inj : 1
Inj Volume : 5.000 µl
Different Inj Volume from Sequence ! Actual Inj Volume : 4.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-02-18 22-48-03\IC-03-20.M
Last changed : 2/18/2020 11:00:23 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:53:05 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.982	VV	0.4233	2.14795e4	799.94556	97.4914
2	10.100	VB	0.4687	552.70178	17.77789	2.5086

Totals : 2.20322e4 817.72345

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.982	BB	0.4199	9706.77539	365.40839	97.5331
2	10.099	BB	0.4697	245.51053	7.87633	2.4669

Totals : 9952.28592 373.28471

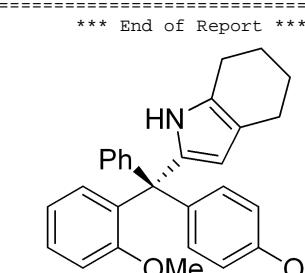
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.981	BB	0.4133	2565.37451	98.69510	97.9584
2	10.269	MM	0.4166	53.46511	2.13881	2.0416

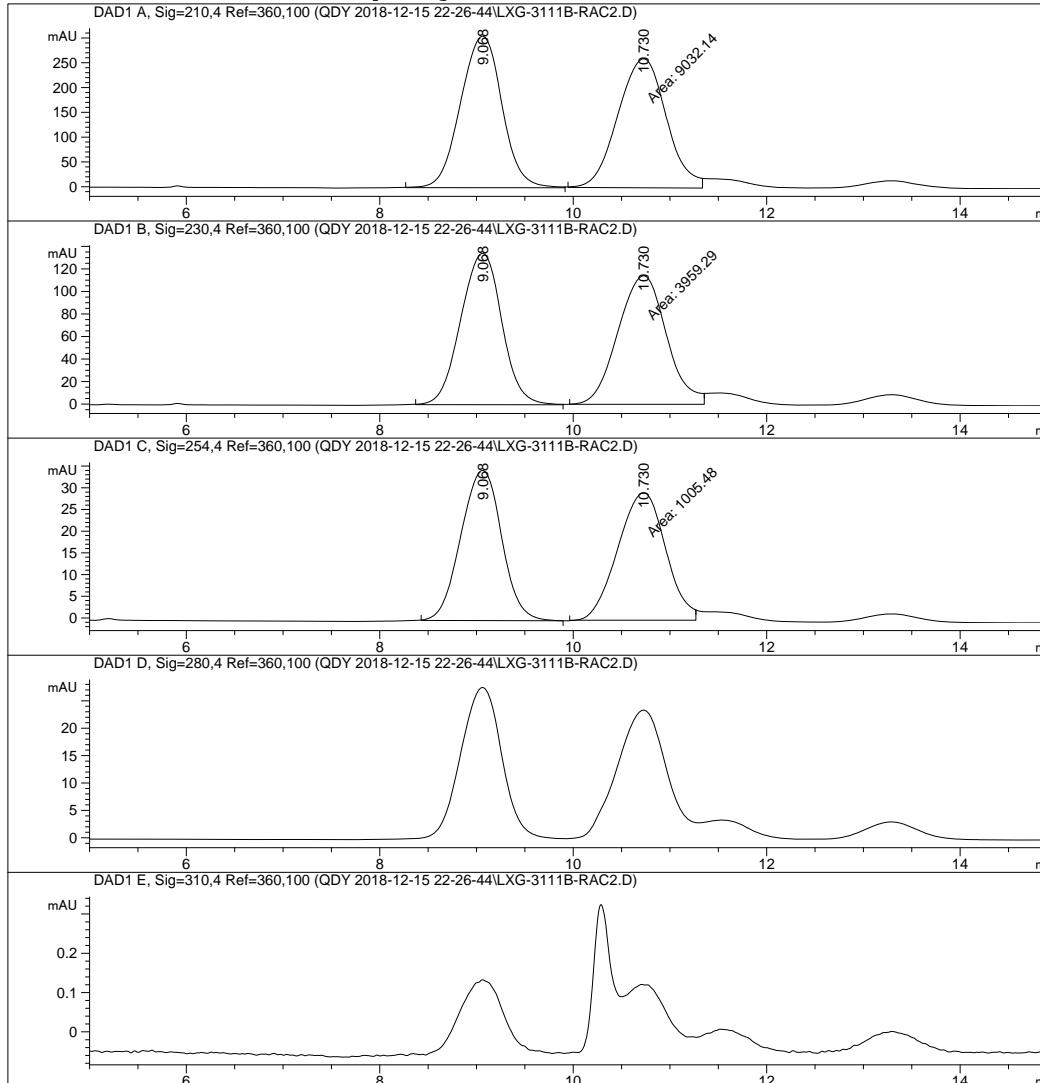
Totals : 2618.83963 100.83391

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



```
=====
Acq. Operator :                               Seq. Line : 2
Acq. Instrument : Instrument 1             Location : Vial 70
Injection Date : 12/15/2018 10:40:33 PM      Inj : 1
                                                Inj Volume : 5.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2018-12-15 22-26-44\IC-03-20.M
Last changed : 12/15/2018 10:39:38 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:54:52 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.068	BB	0.4701	9071.71387	305.83694	50.1093
2	10.730	MM	0.5784	9032.13965	260.24973	49.8907
Totals :					1.81039e4	566.08667

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

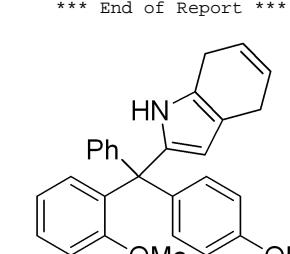
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.068	BB	0.4673	3956.71411	134.47079	49.9838
2	10.730	MM	0.5783	3959.28564	114.11115	50.0162
Totals :					7915.99976	248.58194

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.068	BB	0.4649	1011.59222	34.62283	50.1515
2	10.730	MM	0.5703	1005.48236	29.38600	49.8485
Totals :					2017.07458	64.00883

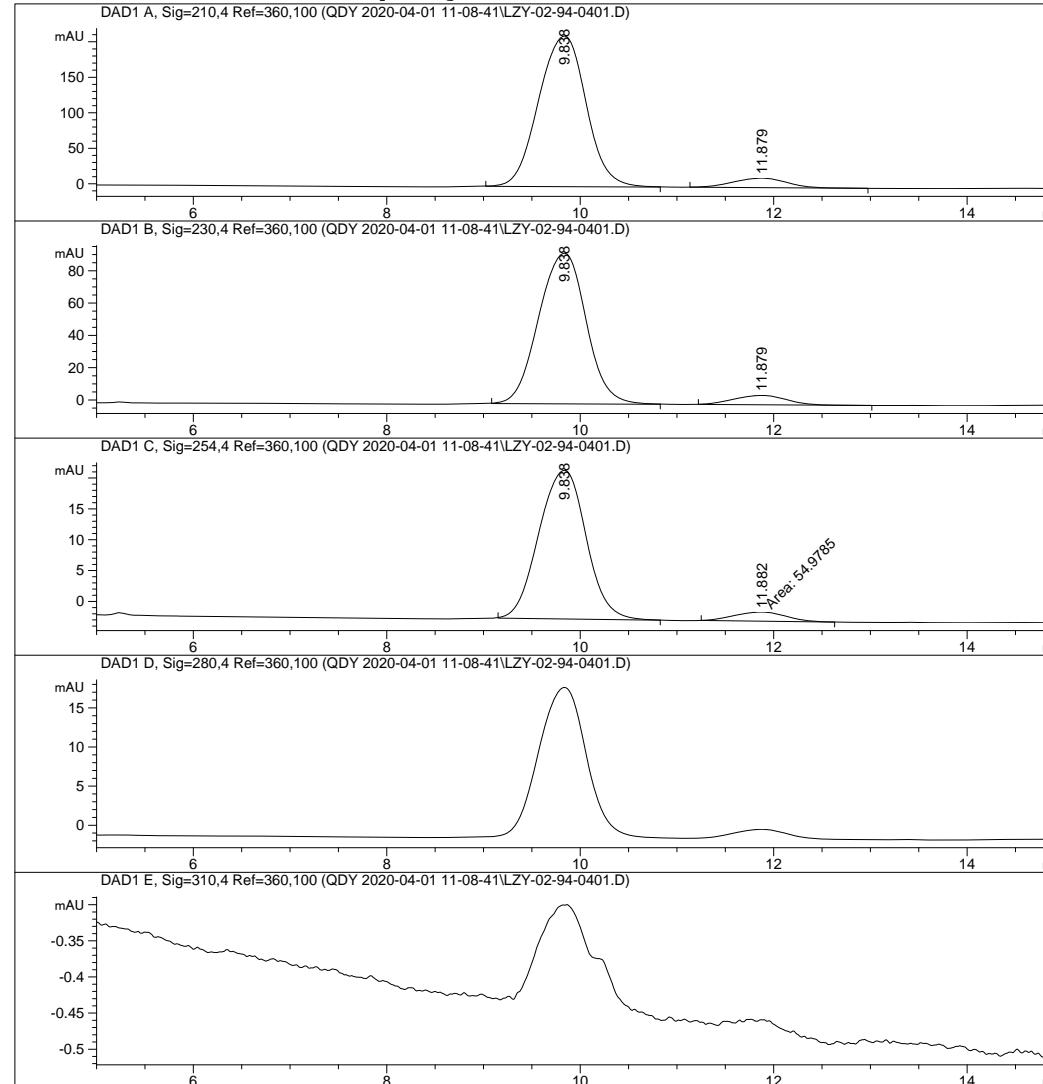
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



2v-Rac

=====
 Acq. Operator : Seq. Line : 2
 Acq. Instrument : Instrument 1 Location : Vial 85
 Injection Date : 4/1/2020 11:22:08 AM Inj : 1
 Inj Volume : 5.000 μ l
 Acq. Method : C:\CHEM32\1\DATA\QDY 2020-04-01 11-08-41\IC-03-20.M
 Last changed : 4/1/2020 11:21:15 AM
 (modified after loading)
 Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
 Last changed : 11/24/2020 6:54:52 PM
 (modified after loading)
 Additional Info : Peak(s) manually integrated



===== Area Percent Report =====

Sorted By : Signal
 Multiplier : 1.0000
 Dilution : 1.0000
 Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.838	BB	0.5457	7230.84375	212.11845	93.4383
2	11.879	BB	0.6092	507.78485	13.30731	6.5617
Totals :					7738.62860	225.42577

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

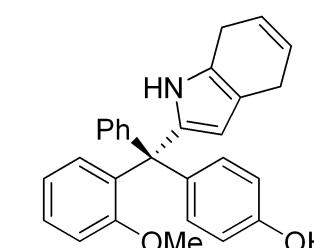
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.838	BB	0.5450	3174.27417	93.29274	93.5923
2	11.879	BB	0.6091	217.32181	5.77410	6.4077
Totals :					3391.59598	99.06684

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.838	BB	0.5440	817.92877	24.09974	93.7017
2	11.882	MM	0.6185	54.97850	1.48138	6.2983
Totals :					872.90727	25.58113

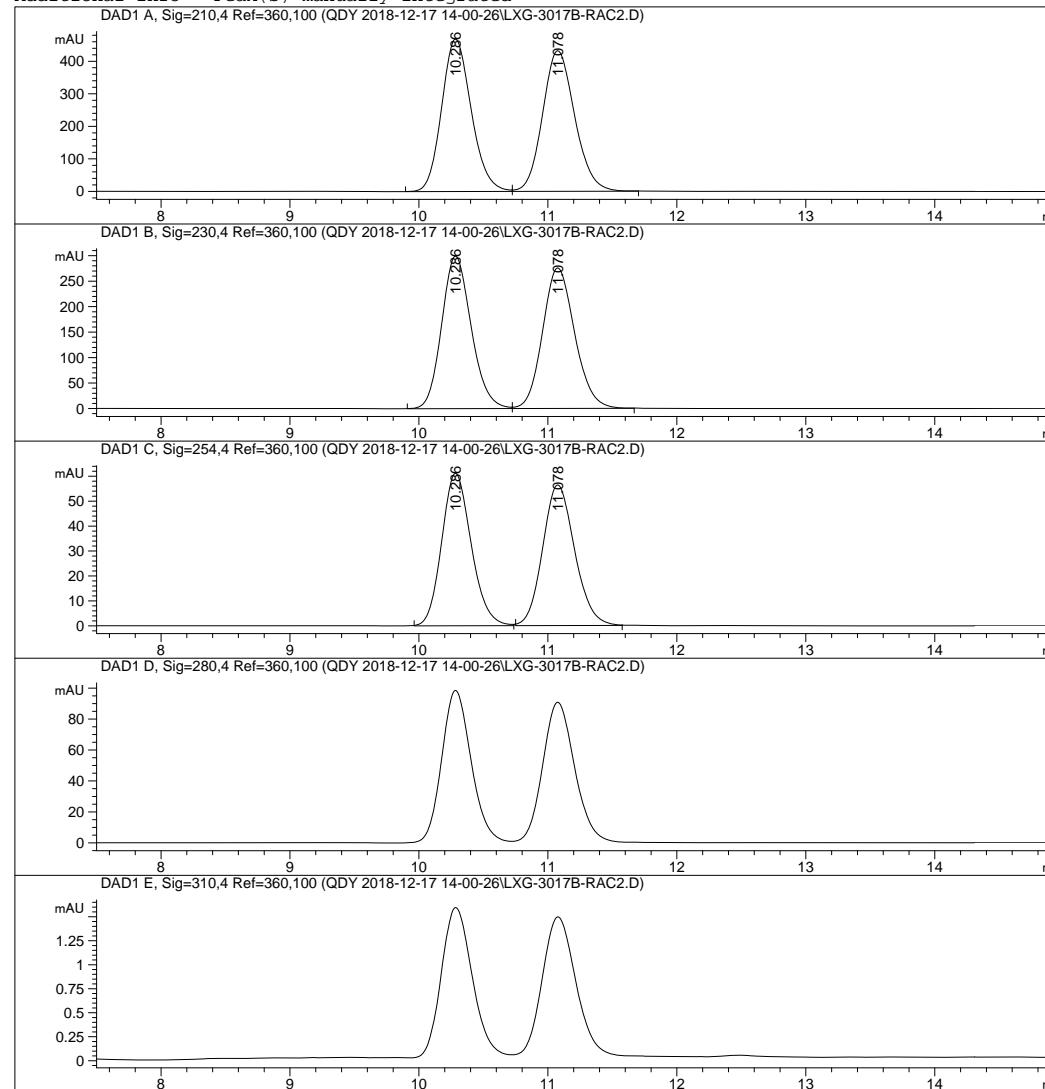
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100

**2v**

Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 5
Acq. Instrument : Instrument 1             Location : Vial 61
Injection Date : 12/17/2018 2:57:29 PM       Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 4.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2018-12-17 14-00-26\AD-03-20.M
Last changed : 7/2/2016 3:26:02 PM
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:58:51 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



Data File C:\CHEM32\1\DATA\QDY 2018-12-17 14-00-26\LXG-3017B-RAC2.D
Sample Name:

```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.286	BV	0.2457	7460.69238	468.72238	49.9435
2	11.078	VB	0.2678	7477.58594	432.23337	50.0565

Totals : 1.49383e4 900.95575

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.286	BV	0.2455	4760.53760	299.33954	49.9784
2	11.078	VB	0.2675	4764.64502	275.89572	50.0216

Totals : 9525.18262 575.23526

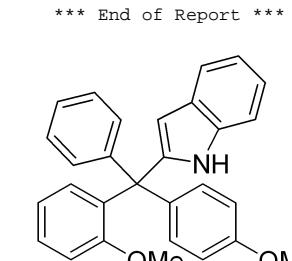
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.286	BB	0.2452	970.58643	61.14531	50.0038
2	11.078	BB	0.2669	970.43921	56.36258	49.9962

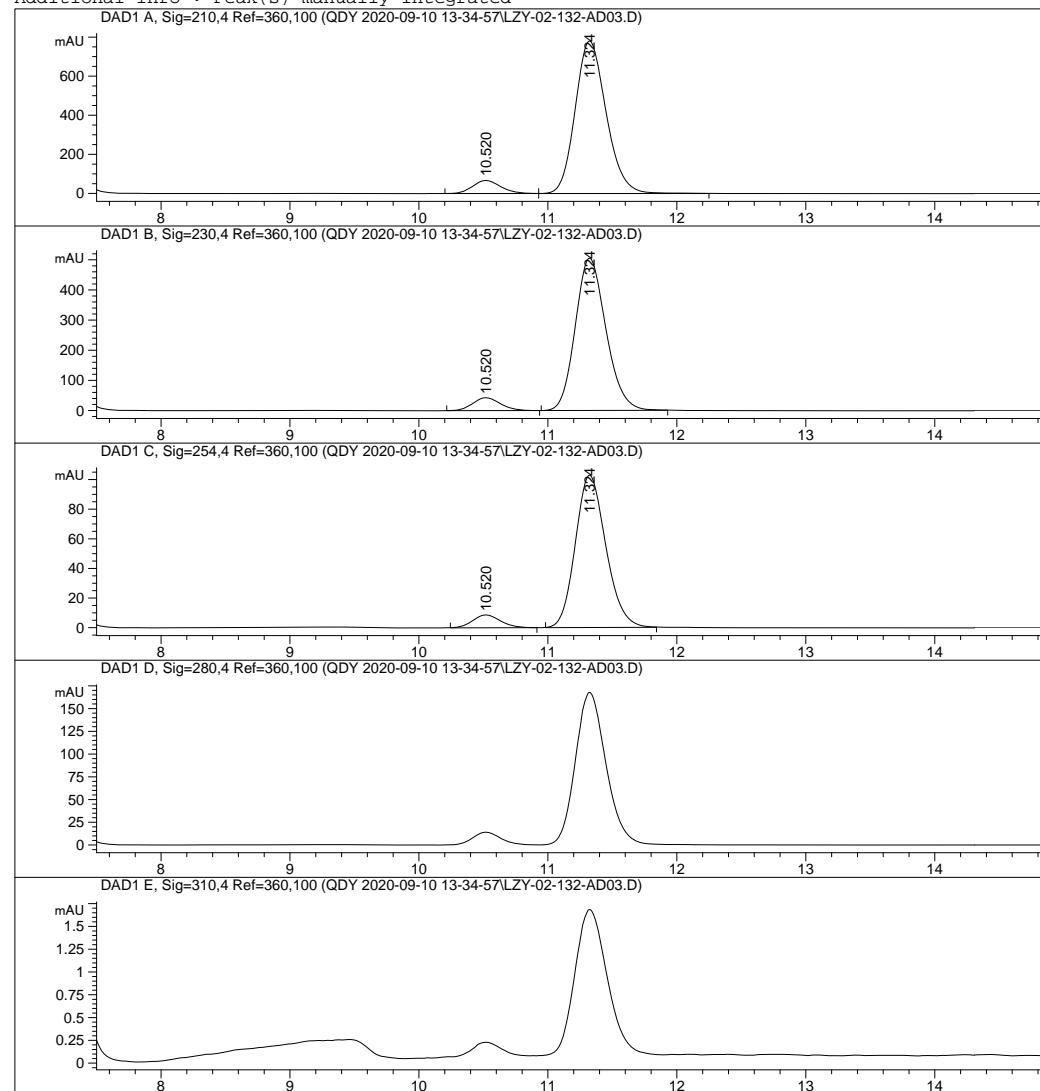
Totals : 1941.02563 117.50788

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100

**2w-Rac**

```
=====
Acq. Operator : Seq. Line : 33
Acq. Instrument : Instrument 1 Location : Vial 93
Injection Date : 9/10/2020 11:30:53 PM Inj : 1
Inj Volume : 5.000 µl
Different Inj Volume from Sequence ! Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-09-10 13-34-57\AD-03-20.M
Last changed : 7/2/2016 3:26:02 PM
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 6:58:51 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.520	BV	0.2322	1011.62958	66.99722	7.2128
2	11.324	VB	0.2586	1.30139e4	780.21997	92.7872
Totals :				1.40256e4	847.21719	

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

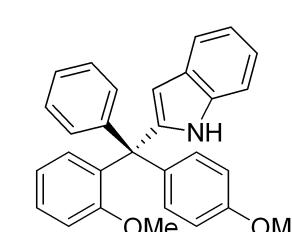
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.520	BB	0.2305	640.46625	42.81888	7.1278
2	11.324	BB	0.2546	8345.02441	505.40024	92.8722
Totals :				8985.49066	548.21912	

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

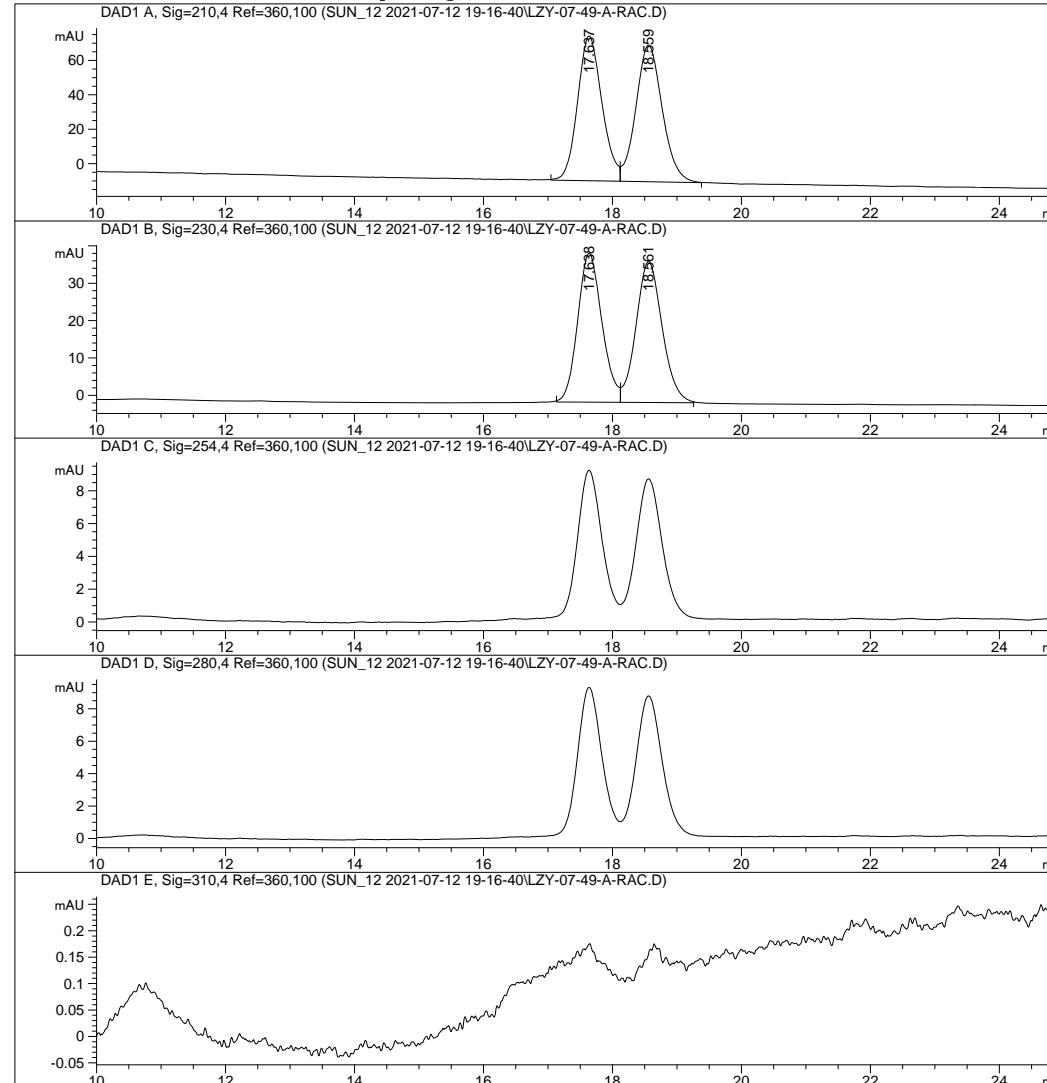
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.520	BB	0.2304	129.39140	8.65633	7.1097
2	11.324	BB	0.2538	1690.52771	102.83469	92.8903
Totals :				1819.91911	111.49102	

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



```
=====
Acq. Operator :                               Seq. Line : 2
Acq. Instrument : Instrument 1             Location : Vial 81
Injection Date : 7/12/2021 7:29:58 PM        Inj : 1
                                                Inj Volume : 5.000 µl
Acq. Method : C:\CHEM32\1\DATA\SUN_12 2021-07-12 19-16-40\AD-02-30.M
Last changed : 7/12/2021 7:29:05 PM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\AS-18-45-0.3.M
Last changed : 7/19/2021 6:21:35 PM        (modified after loading)
Additional Info : Peak(s) manually integrated
```



```
=====
Area Percent Report
=====

Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak RetTime Type Width Area Height Area
# [min] [min] [mAU*s] [mAU] %
-----|-----|-----|-----|-----|
1 17.637 BV 0.4044 2207.37720 84.05779 49.4282
2 18.559 VB 0.4398 2258.44849 79.41764 50.5718

Totals : 4465.82568 163.47543
```

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

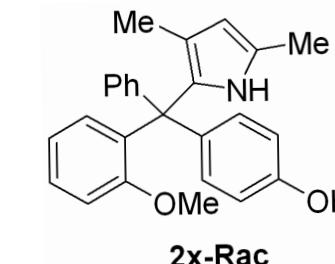
```
Peak RetTime Type Width Area Height Area
# [min] [min] [mAU*s] [mAU] %
-----|-----|-----|-----|-----|
1 17.638 BV 0.4022 1039.56750 39.86308 49.4553
2 18.561 VB 0.4308 1062.46765 37.70211 50.5447

Totals : 2102.03516 77.56519
```

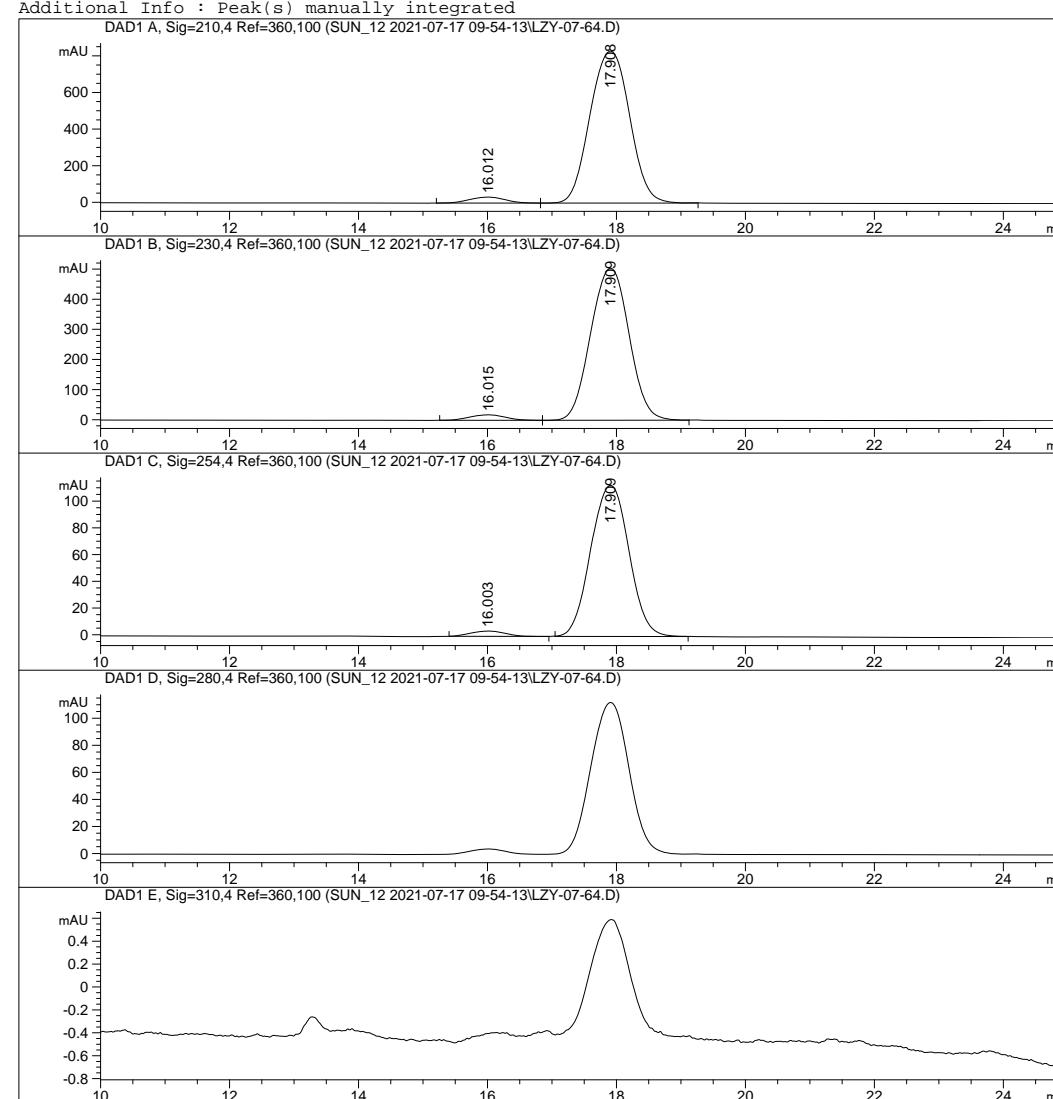
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



```
=====
Acq. Operator : Seq. Line : 33
Acq. Instrument : Instrument 1 Location : Vial 82
Injection Date : 7/17/2021 7:37:34 PM Inj : 1
Inj Volume : 5.000 µl
Different Inj Volume from Sequence ! Actual Inj Volume : 4.000 µl
Acq. Method : C:\CHEM32\1\DATA\SUN_12 2021-07-17 09-54-13\AD-02-30.M
Last changed : 7/17/2021 7:36:41 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\AS-18-45-0.3.M
Last changed : 7/19/2021 6:21:35 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	16.012	BV	0.5067	1255.66016	32.50308	3.4268
2	17.908	VB	0.6832	3.53865e4	830.29260	96.5732

Totals : 3.66422e4 862.79568

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	16.015	BV	0.6106	695.39563	18.08652	3.2334
2	17.909	VB	0.6631	2.08113e4	504.65451	96.7666

Totals : 2.15067e4 522.74103

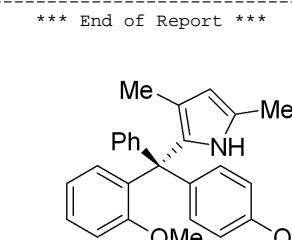
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	16.003	BB	0.4995	147.08163	3.92190	3.0667
2	17.909	BB	0.6596	4648.98389	113.09949	96.9333

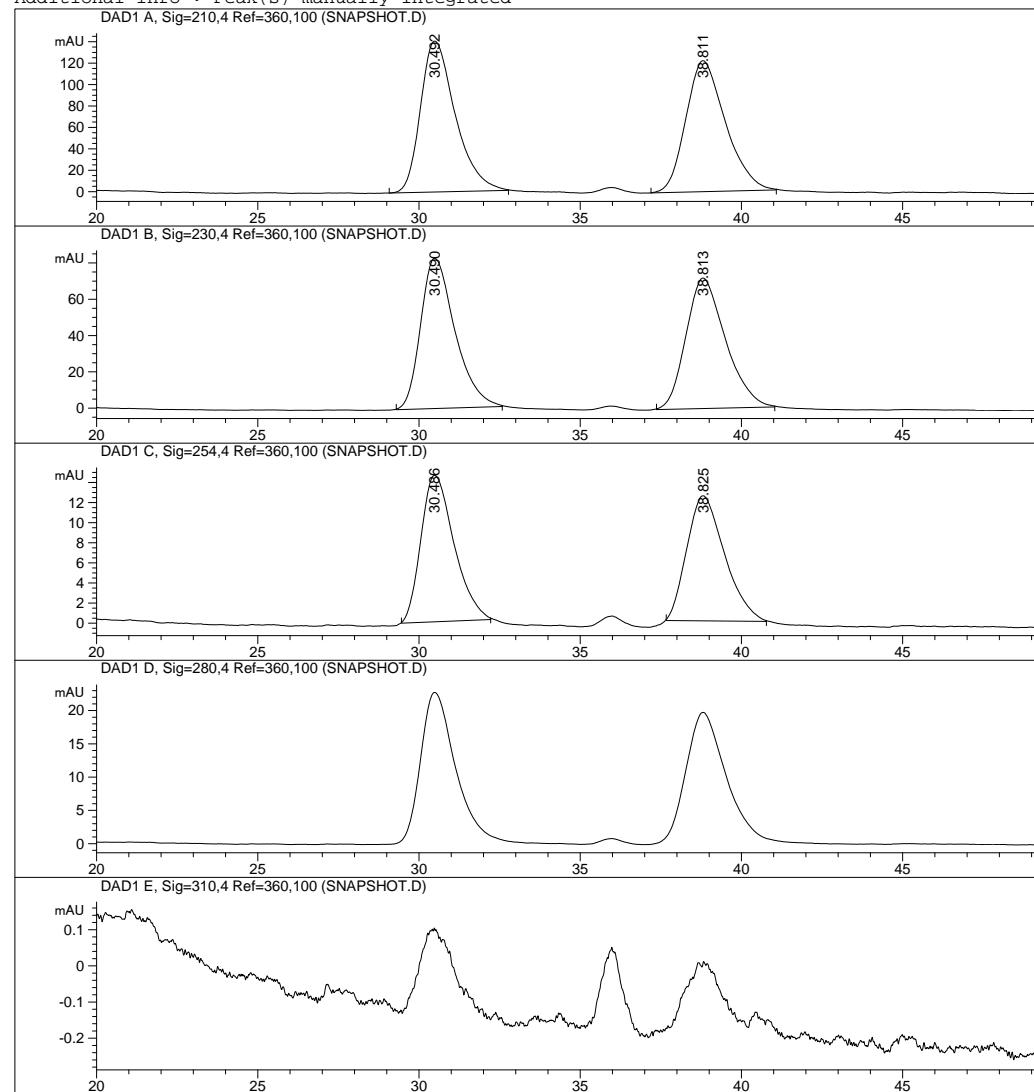
Totals : 4796.06552 117.02139

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



=====
 Acq. Operator : Seq. Line : 11
 Location : Vial 83
 Injection Date : 7/19/2021 10:05:41 PM Inj : 1
 Acq. Method : IC-03-60.M
 Analysis Method : C:\CHEM32\1\METHODS\IC-07-10.M
 Last changed : 7/19/2021 10:50:43 PM
 (modified after loading)
 Additional Info : Peak(s) manually integrated



===== Area Percent Report =====

Sorted By : Signal
 Multiplier : 1.0000
 Dilution : 1.0000
 Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	30.492	BB	0.9897	1.05487e4	140.85474	50.2662
2	38.811	BB	1.0483	1.04370e4	122.21844	49.7338

Totals : 2.09856e4 263.07318

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	30.490	BB	1.0987	6123.68701	82.81453	50.0182
2	38.813	BB	1.2227	6119.22803	71.91388	49.9818

Totals : 1.22429e4 154.72841

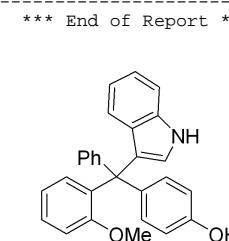
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	30.486	BB	0.9396	1038.62061	14.57118	50.6285
2	38.825	BB	0.9698	1012.83453	12.44062	49.3715

Totals : 2051.45514 27.01181

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

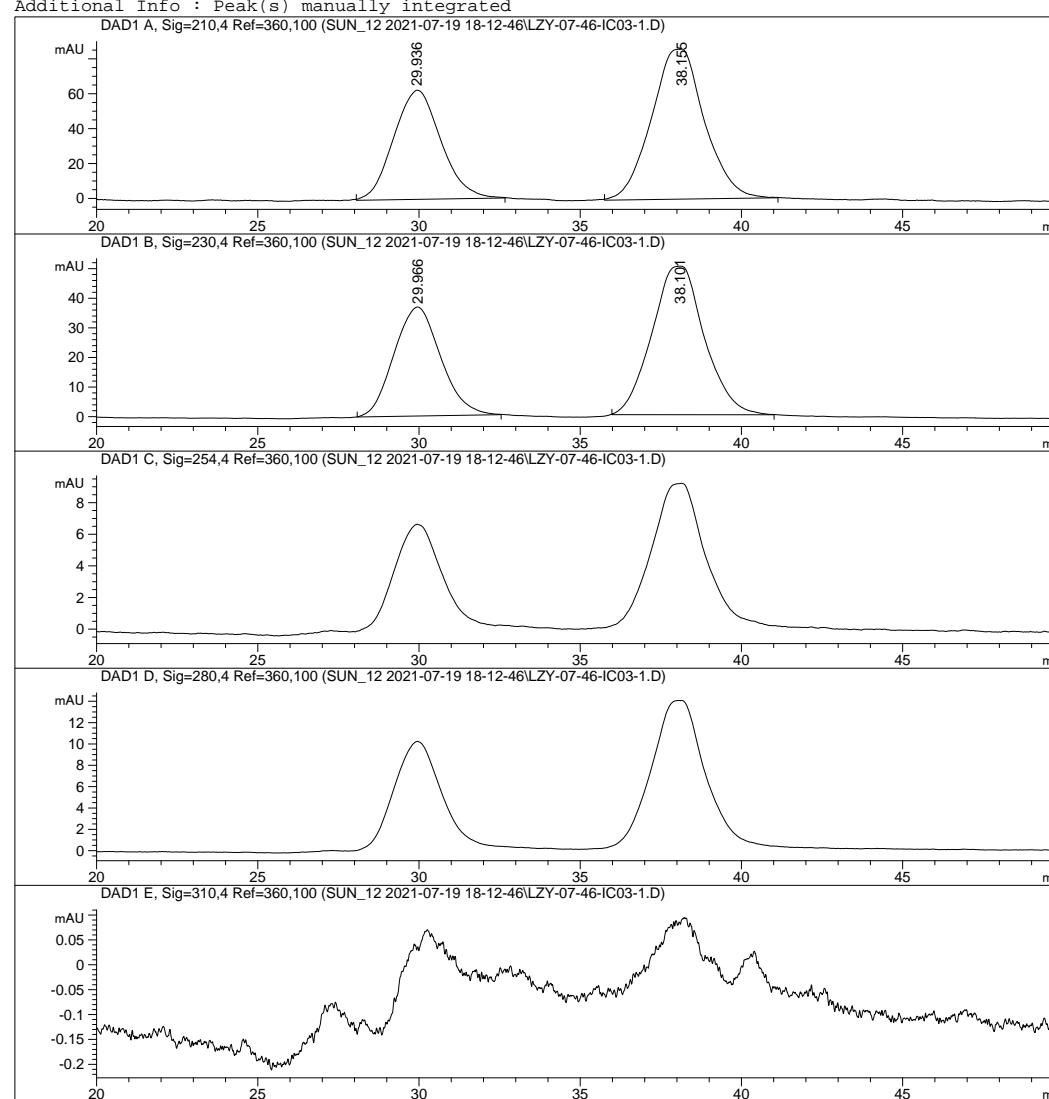
Signal 5: DAD1 E, Sig=310,4 Ref=360,100



2y-Rac

Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 10
Acq. Instrument : Instrument 1             Location : Vial 83
Injection Date : 7/19/2021 9:04:34 PM       Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 4.000 µl
Acq. Method : C:\CHEM32\1\DATA\SUN_12 2021-07-19 18-12-46\IC-03-60.M
Last changed : 5/7/2019 10:27:30 AM
Analysis Method : C:\CHEM32\1\METHODS\IC-07-10.M
Last changed : 7/19/2021 10:50:43 PM
                                         (modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

```
=====
Area Percent Report
=====
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs

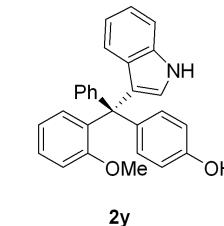
Signal 1: DAD1 A, Sig=210,4 Ref=360,100
Peak RetTime Type Width Area Height Area
# [min]   [min] [mAU*s] [mAU] %
-----|-----|-----|-----|-----|
1 29.936 BB 1.2006 6361.79541 62.54972 39.3273
2 38.155 BB 1.3440 9814.74414 85.96726 60.6727

Totals :           1.61765e4 148.51698

Signal 2: DAD1 B, Sig=230,4 Ref=360,100
Peak RetTime Type Width Area Height Area
# [min]   [min] [mAU*s] [mAU] %
-----|-----|-----|-----|-----|
1 29.966 BB 1.2606 3699.13989 36.80454 39.4746
2 38.101 BB 1.3249 5671.80713 50.17292 60.5254

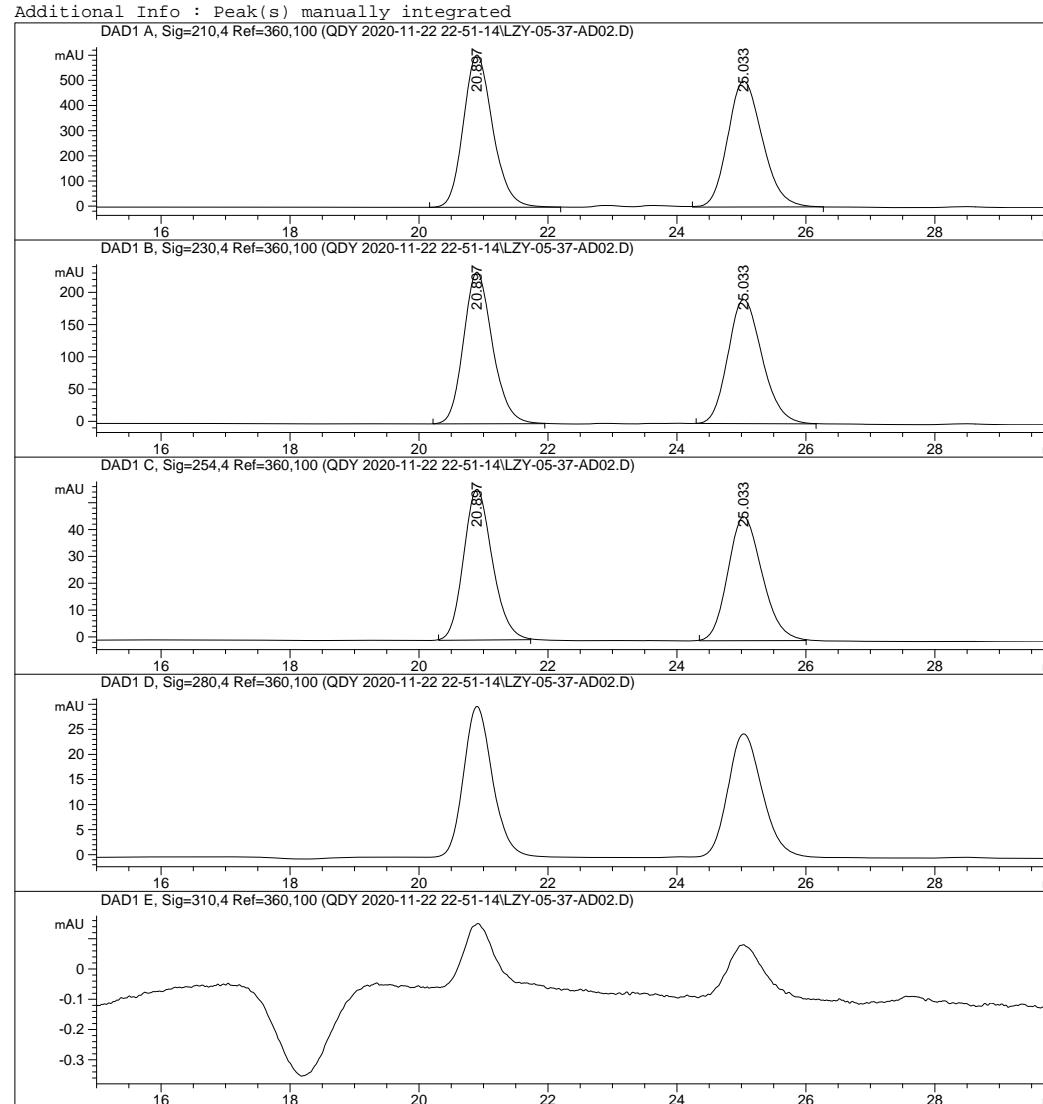
Totals :           9370.94702 86.97746

Signal 3: DAD1 C, Sig=254,4 Ref=360,100
Signal 4: DAD1 D, Sig=280,4 Ref=360,100
Signal 5: DAD1 E, Sig=310,4 Ref=360,100
=====
```



Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 5
Acq. Instrument : Instrument 1             Location : Vial 83
Injection Date : 11/23/2020 12:02:15 AM      Inj : 1
                                                Inj Volume : 5.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-11-22 22-51-14\AD-02-60.M
Last changed : 9/23/2020 10:59:45 PM
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 7:37:31 PM
                                                (modified after loading)
Additional Info : Peak(s) manually integrated
```



Data File C:\CHEM32\1\DATA\QDY 2020-11-22 22-51-14\LZY-05-37-AD02.D
Sample Name:

```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	20.897	BB	0.4891	1.91413e4	604.95667	50.6223
2	25.033	VB	0.5794	1.86707e4	500.23126	49.3777

Totals : 3.78119e4 1105.18793

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	20.897	BB	0.4797	7327.97266	235.03383	50.7969
2	25.033	BB	0.5702	7098.05029	192.51373	49.2031

Totals : 1.44260e4 427.54756

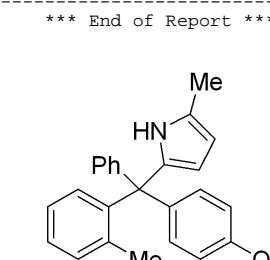
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	20.897	BB	0.4770	1737.50293	56.15725	50.5272
2	25.033	BB	0.5701	1701.24512	46.15036	49.4728

Totals : 3438.74805 102.30761

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

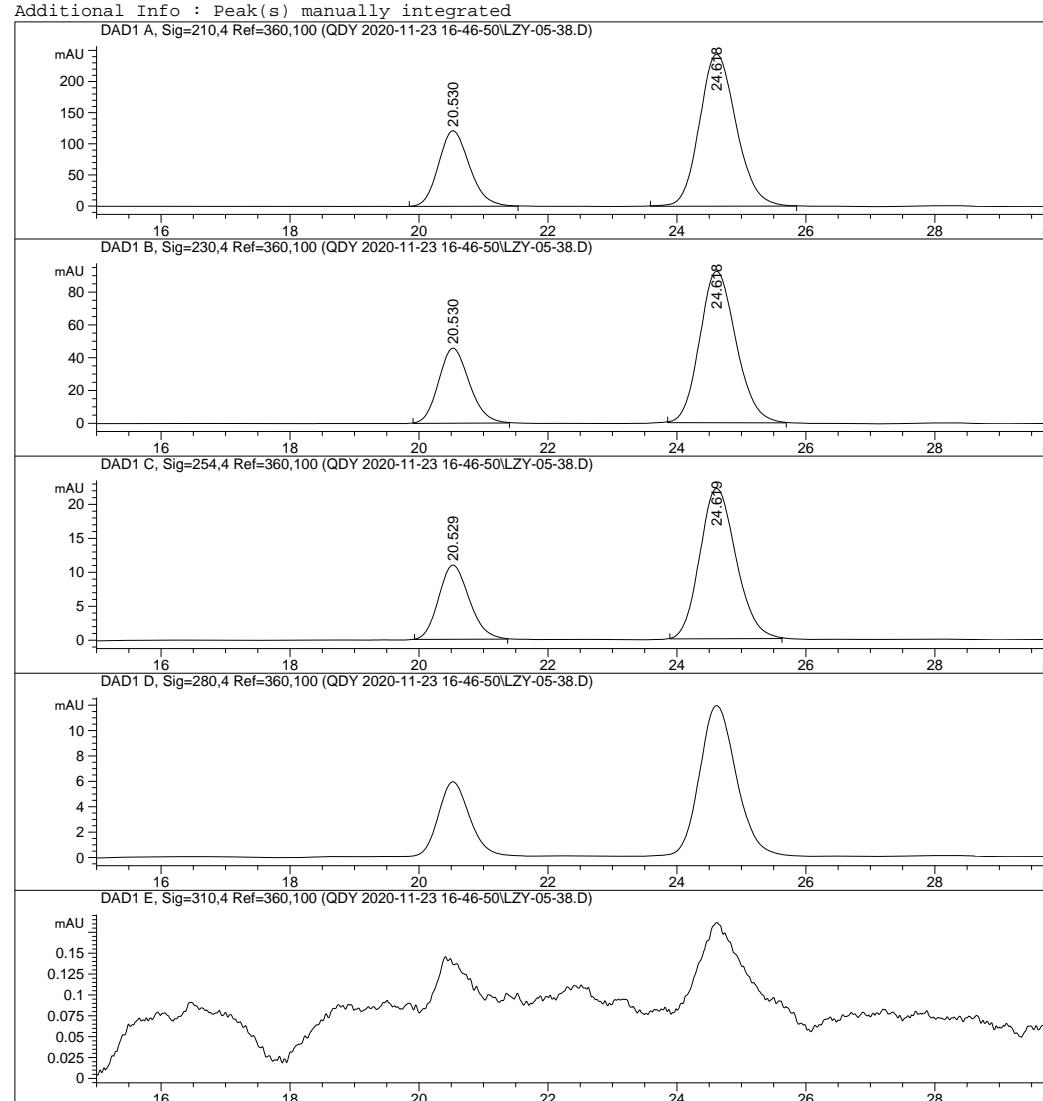
Signal 5: DAD1 E, Sig=310,4 Ref=360,100



2aa-Rac

Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 19
Acq. Instrument : Instrument 1             Location : Vial 84
Injection Date : 11/23/2020 10:47:28 PM      Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !       Actual Inj Volume : 2.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-11-23 16-46-50\AD-02-60.M
Last changed : 9/23/2020 10:59:45 PM
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 7:37:31 PM
                                         (modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	20.530	BB	0.5201	4059.53271	121.40965	29.5948
2	24.618	BB	0.6154	9657.50586	244.19589	70.4052

Totals : 1.37170e4 365.60554

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	20.530	BB	0.5127	1517.53760	45.78773	29.6057
2	24.618	BB	0.6058	3608.29224	92.36298	70.3943

Totals : 5125.82983 138.15071

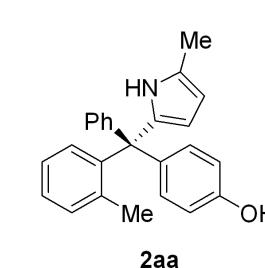
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	20.529	BB	0.5065	360.24033	10.92999	29.4489
2	24.619	BB	0.6046	863.03198	22.14850	70.5511

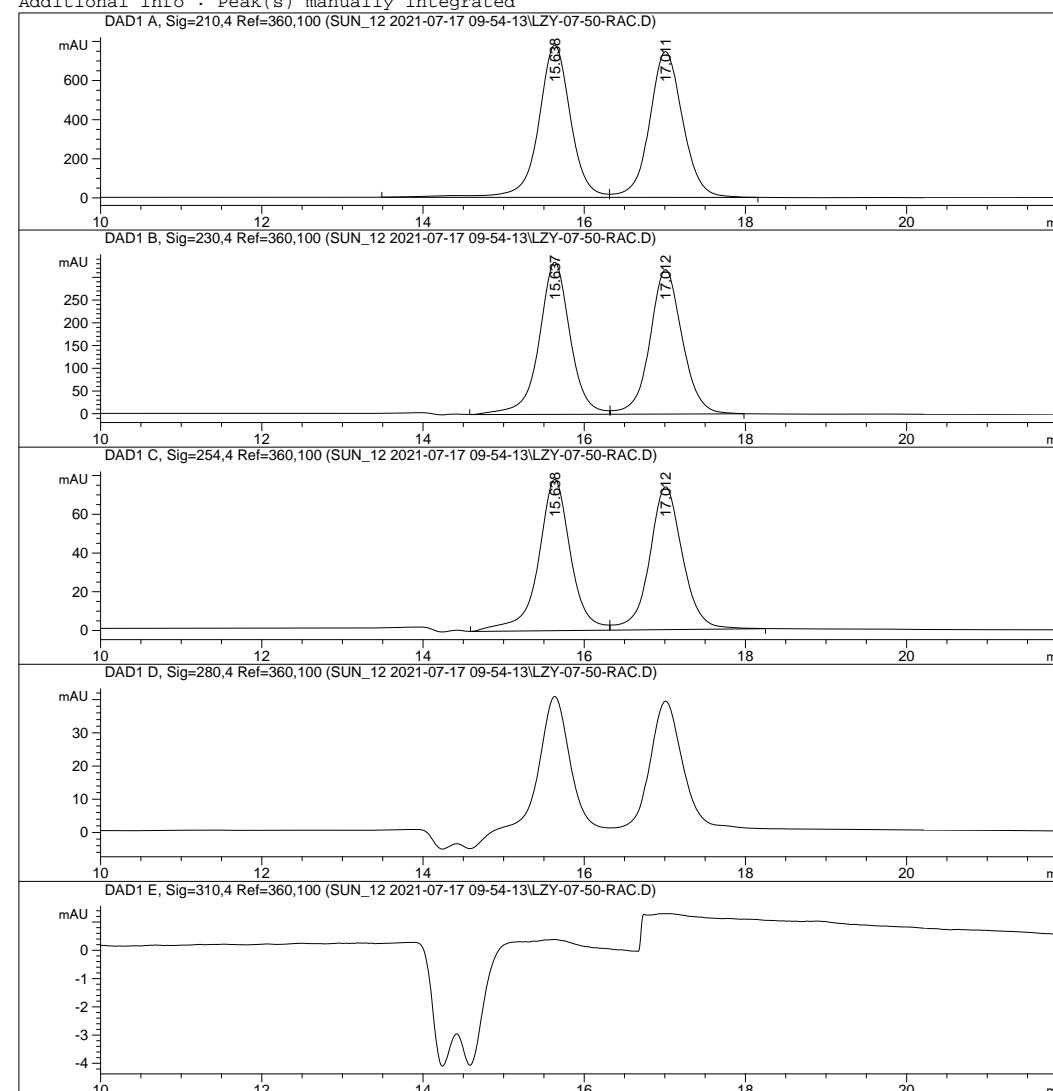
Totals : 1223.27231 33.07850

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



```
=====
Acq. Operator :                               Seq. Line : 34
                                                Location : Vial 84
Injection Date : 7/17/2021 8:08:41 PM      Inj : 1
Acq. Method : AD-03-30.M
Analysis Method : C:\CHEM32\1\METHODS\IC-07-10.M
Last changed : 7/19/2021 10:26:00 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	15.638	BV	0.4196	2.15957e4	778.59137	50.3315
2	17.011	VB	0.4356	2.13113e4	745.19684	49.6685

Totals : 4.29070e4 1523.78821

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	15.637	VV	0.4015	8933.09766	334.55411	50.2286
2	17.012	VB	0.4231	8851.79395	317.68359	49.7714

Totals : 1.77849e4 652.23770

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	15.638	VV	0.4128	2161.18896	78.11736	50.7231
2	17.012	VB	0.4282	2099.57007	73.72577	49.2769

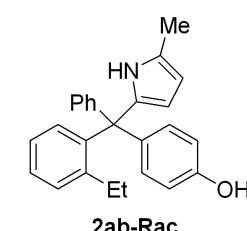
Totals : 4260.75903 151.84312

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100

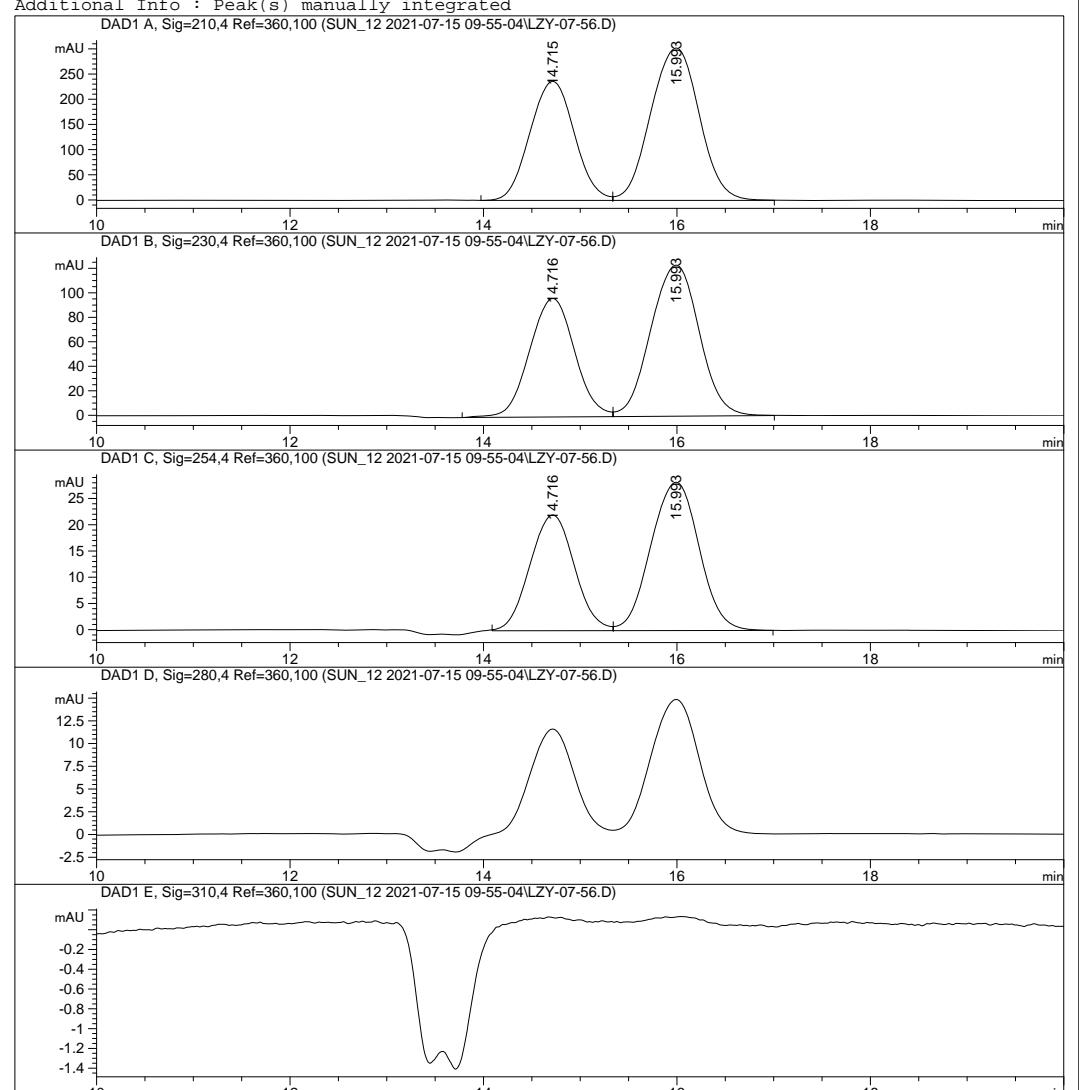
=====

*** End of Report ***



Sample Name:

```
=====
Acq. Operator : Seq. Line : 25
Acq. Instrument : Instrument 1 Location : Vial 81
Injection Date : 7/15/2021 4:55:09 PM Inj : 1
Inj Volume : 5.000 µl
Different Inj Volume from Sequence ! Actual Inj Volume : 4.000 µl
Acq. Method : C:\CHEM32\1\DATA\SUN_12 2021-07-15 09-55-04\AD-03-30.M
Last changed : 7/15/2021 4:54:16 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\AS-15-15.M
Last changed : 7/17/2021 9:58:58 AM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	14.715	BV	0.5034	7521.92773	236.28804	41.6410
2	15.993	VB	0.5588	1.05418e4	302.36267	58.3590

Totals : 1.80637e4 538.65071

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	14.716	BV	0.5115	3131.12671	96.77523	42.1442
2	15.993	VB	0.5587	4298.43506	123.29113	57.8558

Totals : 7429.56177 220.06636

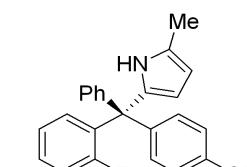
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	14.716	BV	0.5022	703.77405	22.05937	41.7689
2	15.993	VB	0.5550	981.14893	28.26350	58.2311

Totals : 1684.92297 50.32287

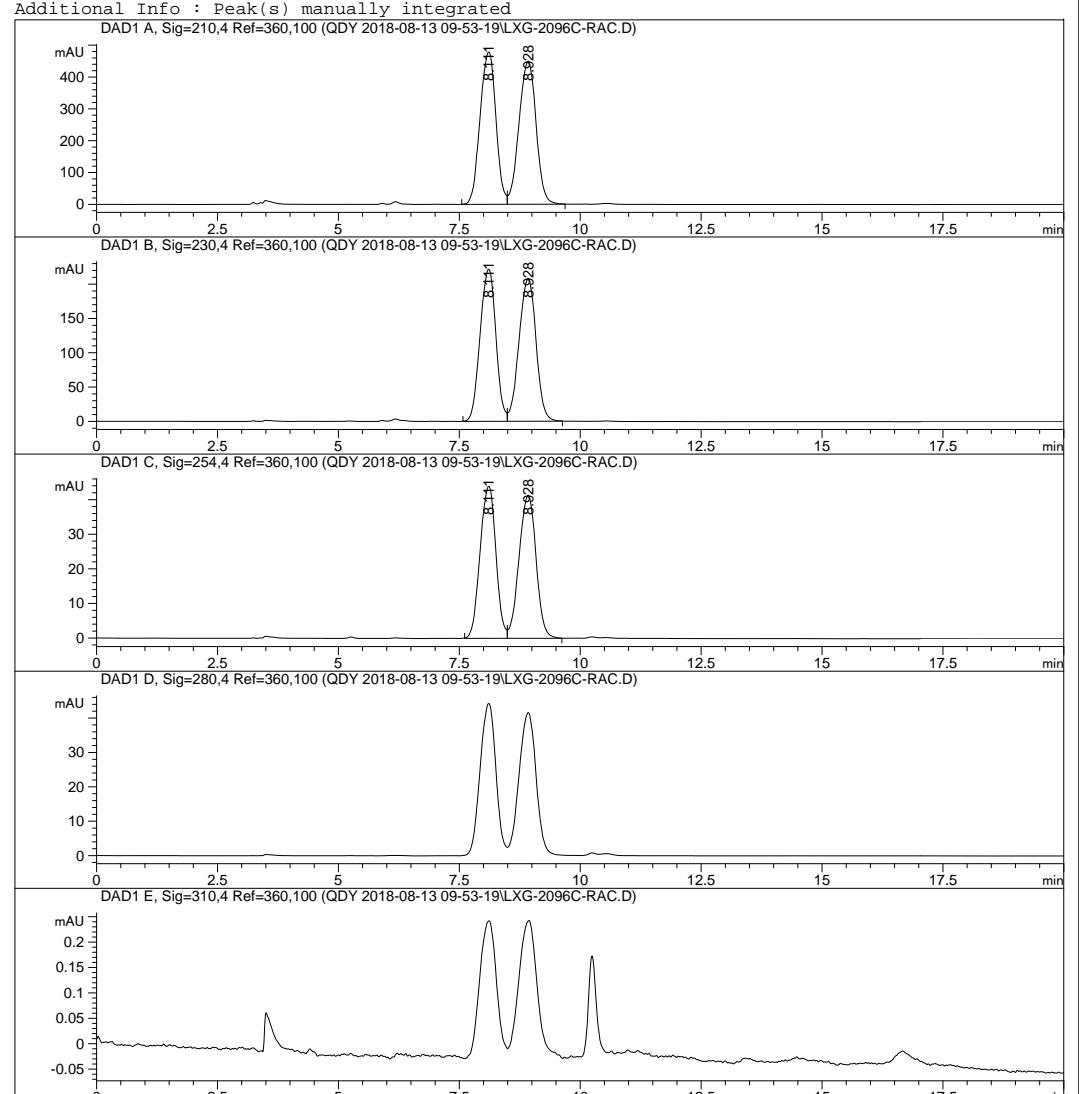
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 12
Acq. Instrument : Instrument 1             Location : Vial 56
Injection Date : 8/13/2018 2:39:53 PM       Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2018-08-13 09-53-19\IC-02-20.M
Last changed : 8/13/2018 2:39:01 PM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\AD-10-10.M
Last changed : 1/16/2020 9:52:19 AM        (modified after loading)
Additional Info : Peak(s) manually integrated
```



Data File C:\CHEM32\1\DATA\QDY 2018-08-13 09-53-19\LXG-2096C-RAC.D
Sample Name:

```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.111	BV	0.3655	1.08839e4	478.20779	49.6404
2	8.928	VB	0.3945	1.10416e4	449.64343	50.3596
Totals :					2.19255e4	927.85123

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

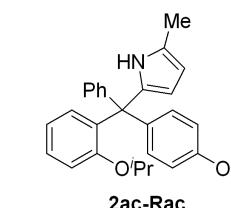
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.111	BV	0.3645	5026.81592	221.75375	49.6977
2	8.928	VB	0.3910	5087.97168	208.26950	50.3023
Totals :					1.01148e4	430.02325

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.111	BV	0.3646	997.15350	43.95827	49.6376
2	8.928	VB	0.3893	1011.71417	41.37379	50.3624
Totals :					2008.86768	85.33206

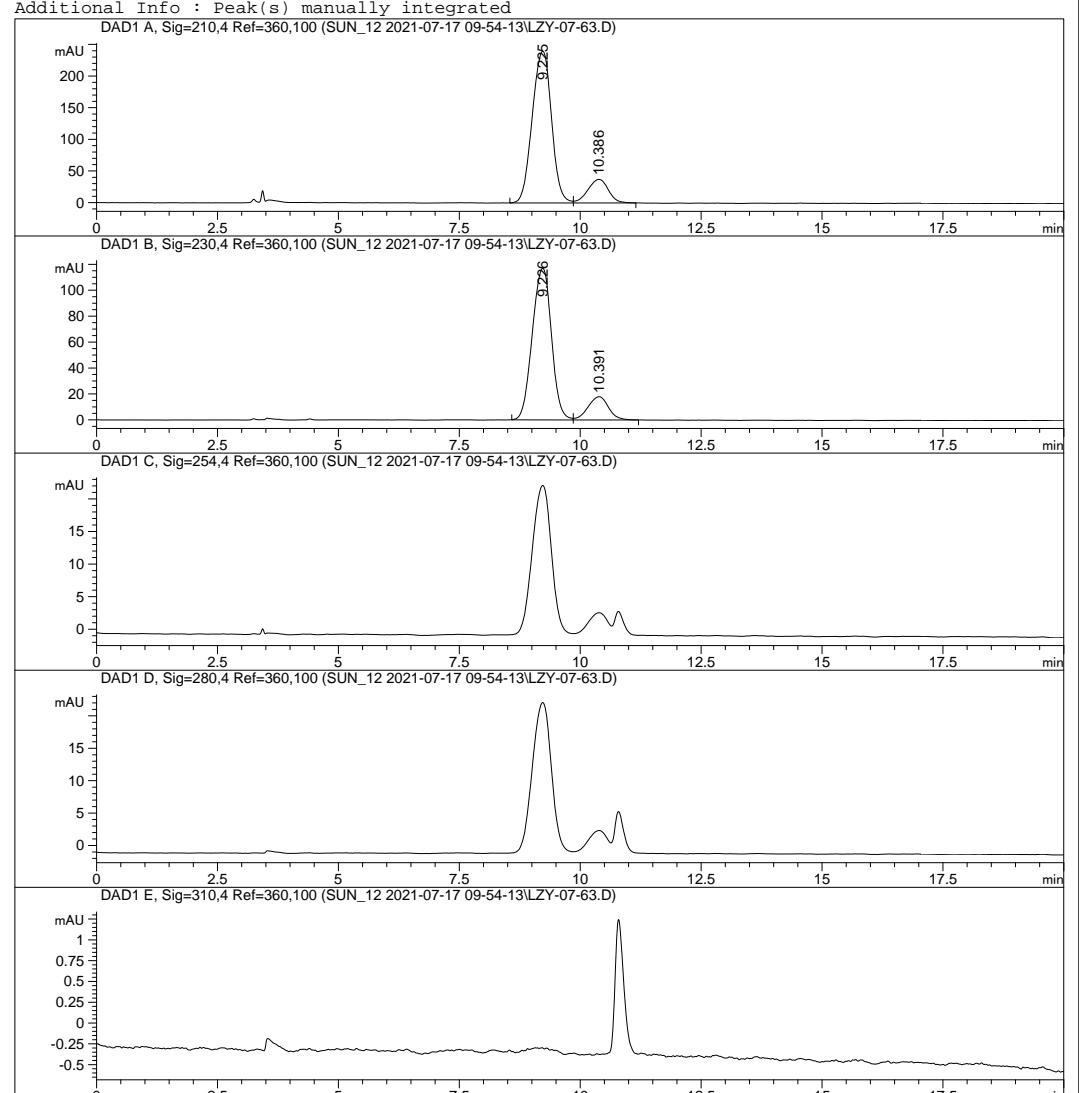
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 29
Acq. Instrument : Instrument 1             Location : Vial 81
Injection Date : 7/17/2021 6:37:47 PM        Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\SUN_12 2021-07-17 09-54-13\IC-02-20.M
Last changed : 7/17/2021 6:36:54 PM
                                                (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\AS-18-45-0.3.M
Last changed : 7/19/2021 6:28:40 PM
                                                (modified after loading)
Additional Info : Peak(s) manually integrated
```



Data File C:\CHEM32\1\DATA\SUN_12 2021-07-17 09-54-13\LZY-07-63.D
Sample Name:

===== Area Percent Report =====

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.225	BV	0.4423	6631.34863	240.00844	85.8247
2	10.386	VB	0.4203	1095.27039	37.33301	14.1753

Totals : 7726.61902 277.34145

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

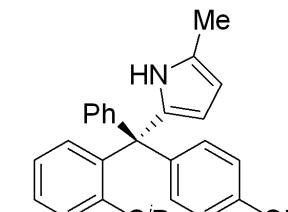
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.226	BV	0.4385	3216.06519	117.06415	85.9157
2	10.391	VB	0.4650	527.21429	17.92997	14.0843

Totals : 3743.27948 134.99412

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

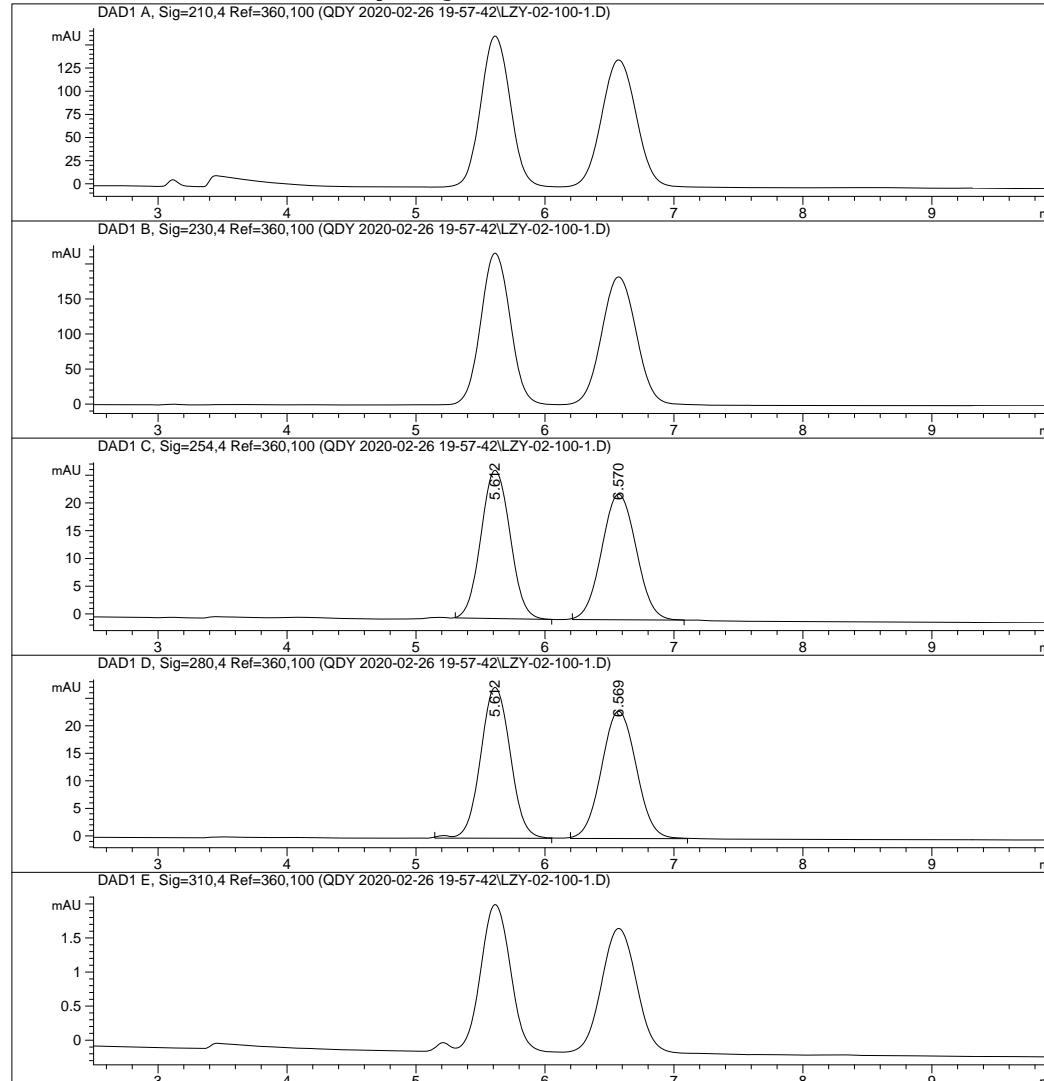
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 2
Acq. Instrument : Instrument 1             Location : Vial 81
Injection Date : 2/26/2020 8:11:05 PM        Inj : 1
                                                Inj Volume : 5.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-02-26 19-57-42\AD-01-30.M
Last changed : 2/26/2020 8:10:10 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 2:57:57 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



Data File C:\CHEM32\1\DATA\QDY 2020-02-26 19-57-42\LZY-02-100-1.D
Sample Name:

```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	5.612	BB	0.2492	419.41470	26.70506	49.9197
2	6.570	BB	0.2930	420.76358	22.61930	50.0803

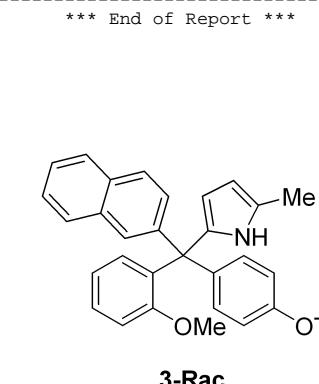
Totals : 840.17828 49.32436

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	5.612	BB	0.2525	437.83301	27.39108	50.2826
2	6.569	BB	0.2984	432.91190	23.12917	49.7174

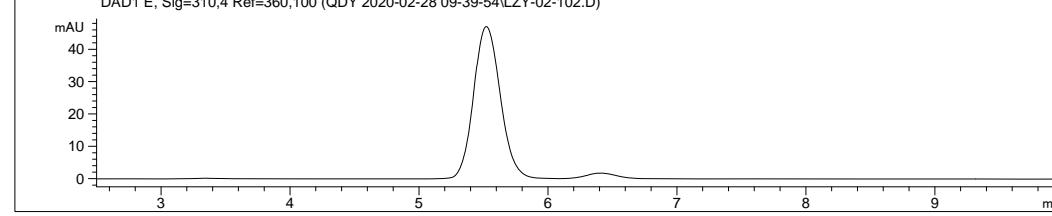
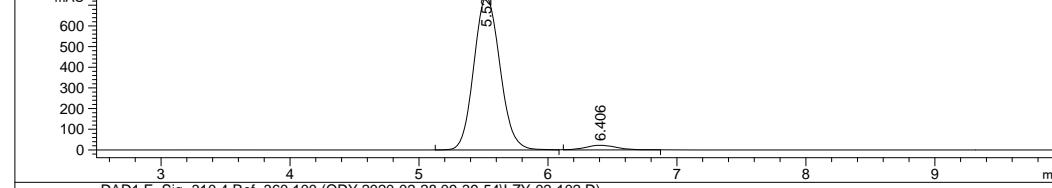
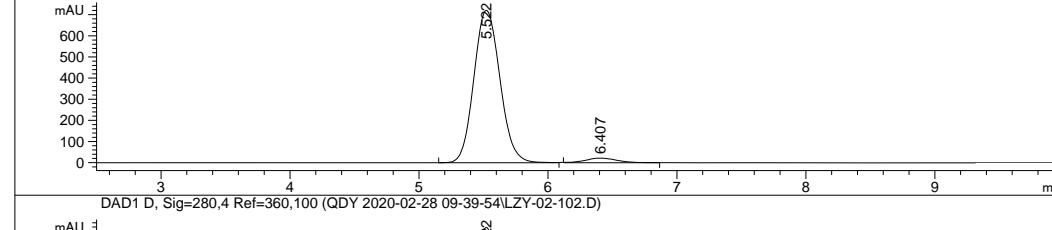
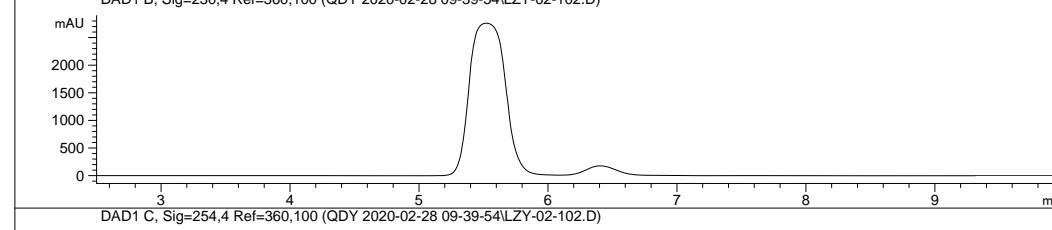
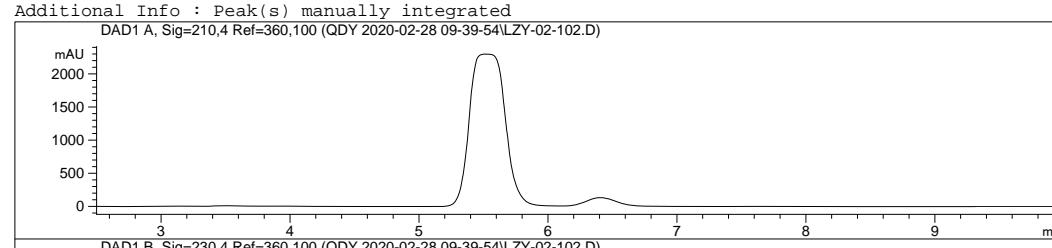
Totals : 870.74490 50.52024

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

```
=====
Acq. Operator :                               Seq. Line :   6
Acq. Instrument : Instrument 1             Location : Vial 81
Injection Date : 2/28/2020 10:57:41 AM       Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 4.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-02-28 09-39-54\AD-01-20.M
Last changed : 2/28/2020 10:56:49 AM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 2:57:57 PM        (modified after loading)
Additional Info : Peak(s) manually integrated
```



Data File C:\CHEM32\1\DATA\QDY 2020-02-28 09-39-54\LZY-02-102.D

Sample Name:

```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	5.522	BB	0.2219	1.01315e4	721.28223	96.4846
2	6.407	BB	0.2647	369.14151	21.89908	3.5154

Totals : 1.05007e4 743.18131

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

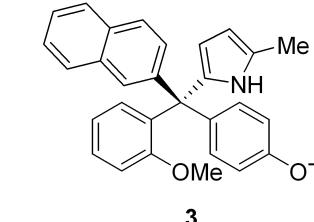
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	5.522	BB	0.2220	1.03741e4	738.08148	96.5137
2	6.406	BB	0.2639	374.73767	22.31883	3.4863

Totals : 1.07488e4 760.40031

Signal 5: DAD1 E, Sig=310,4 Ref=360,100

=====

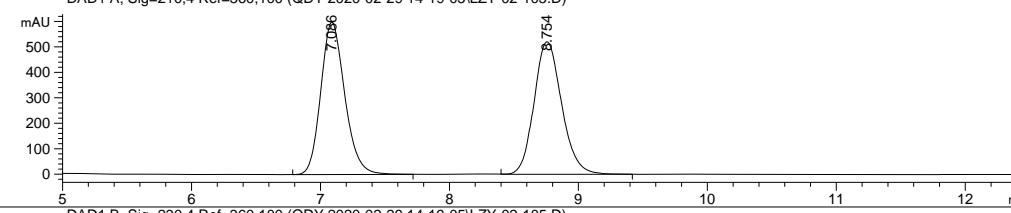
*** End of Report ***



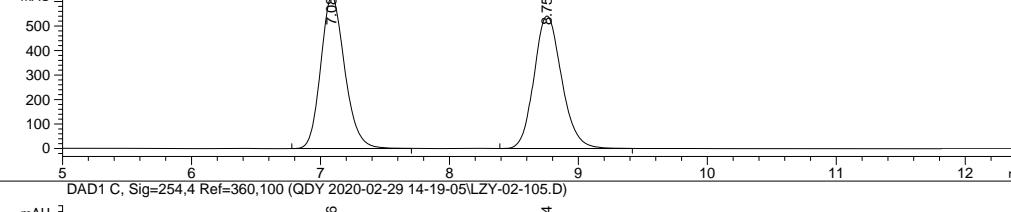
Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 17
Acq. Instrument : Instrument 1             Location : Vial 91
Injection Date : 2/29/2020 7:35:51 PM        Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-02-29 14-19-05\AD-01-20.M
Last changed : 2/29/2020 7:34:58 PM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 8:50:44 PM        (modified after loading)
Additional Info : Peak(s) manually integrated
```

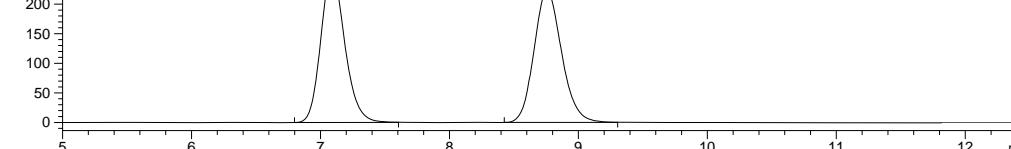
DAD1 A, Sig=210,4 Ref=360,100 (QDY 2020-02-29 14-19-05\LZY-02-105.D)



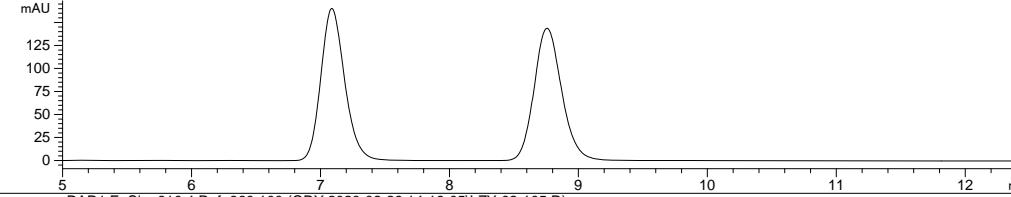
DAD1 B, Sig=230,4 Ref=360,100 (QDY 2020-02-29 14-19-05\LZY-02-105.D)



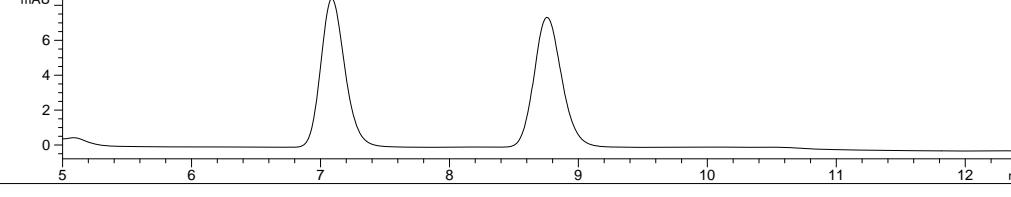
DAD1 C, Sig=254,4 Ref=360,100 (QDY 2020-02-29 14-19-05\LZY-02-105.D)



DAD1 D, Sig=280,4 Ref=360,100 (QDY 2020-02-29 14-19-05\LZY-02-105.D)



DAD1 E, Sig=310,4 Ref=360,100 (QDY 2020-02-29 14-19-05\LZY-02-105.D)



Data File C:\CHEM32\1\DATA\QDY 2020-02-29 14-19-05\LZY-02-105.D
Sample Name:

===== Area Percent Report =====

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.086	BB	0.1983	7712.21045	598.72394	49.9275
2	8.754	VB	0.2291	7734.60010	521.38654	50.0725

Totals : 1.54468e4 1120.11047

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.086	BB	0.1982	8002.52881	621.77722	49.9326
2	8.754	VB	0.2289	8024.14697	541.38800	50.0674

Totals : 1.60267e4 1163.16522

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

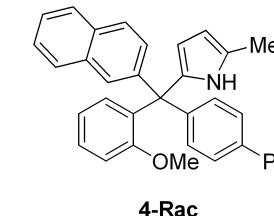
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.086	BB	0.1975	3297.33911	257.24408	49.9841
2	8.754	BB	0.2282	3299.43335	223.58458	50.0159

Totals : 6596.77246 480.82866

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

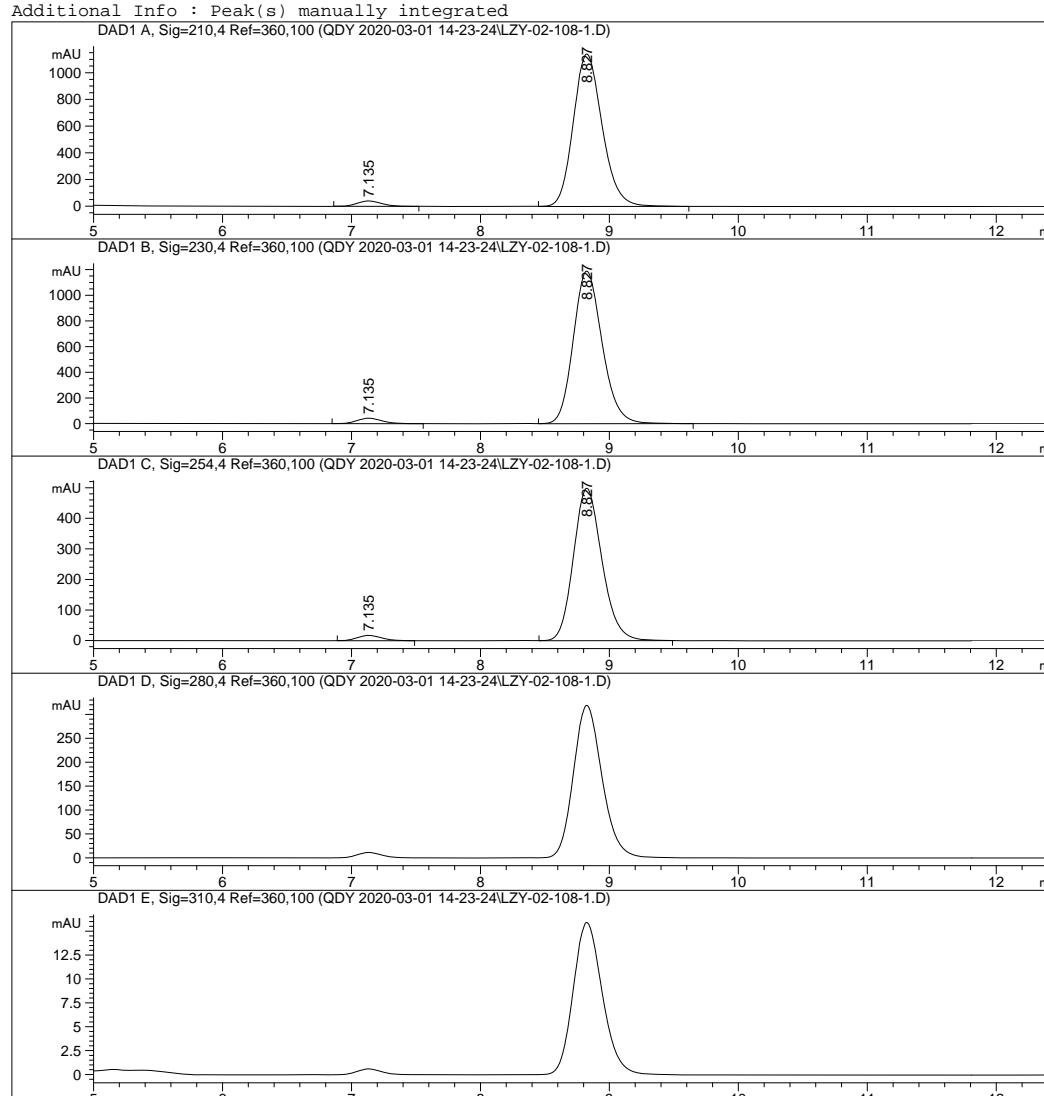
Signal 5: DAD1 E, Sig=310,4 Ref=360,100

=====
*** End of Report ***
=====



Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 2
Acq. Instrument : Instrument 1             Location : Vial 82
Injection Date : 3/1/2020 2:36:27 PM        Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 3.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-03-01 14-23-24\AD-01-20.M
Last changed : 3/1/2020 2:35:34 PM          (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 3:01:11 PM         (modified after loading)
Additional Info : Peak(s) manually integrated
```



Data File C:\CHEM32\1\DATA\QDY 2020-03-01 14-23-24\LZY-02-108-1.D
Sample Name:

```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.135	BB	0.2037	535.52502	40.65454	2.9252
2	8.827	VB	0.2415	1.77716e4	1142.56104	97.0748

Totals : 1.83072e4 1183.21558

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.135	BB	0.2060	573.54810	42.88038	3.0125
2	8.827	VB	0.2412	1.84655e4	1188.60522	96.9875

Totals : 1.90391e4 1231.48561

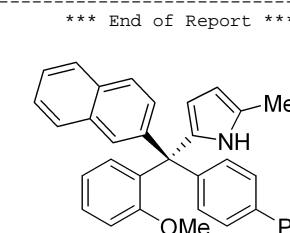
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.135	BB	0.2040	230.95700	17.49571	2.9220
2	8.827	VB	0.2354	7673.16943	498.90271	97.0780

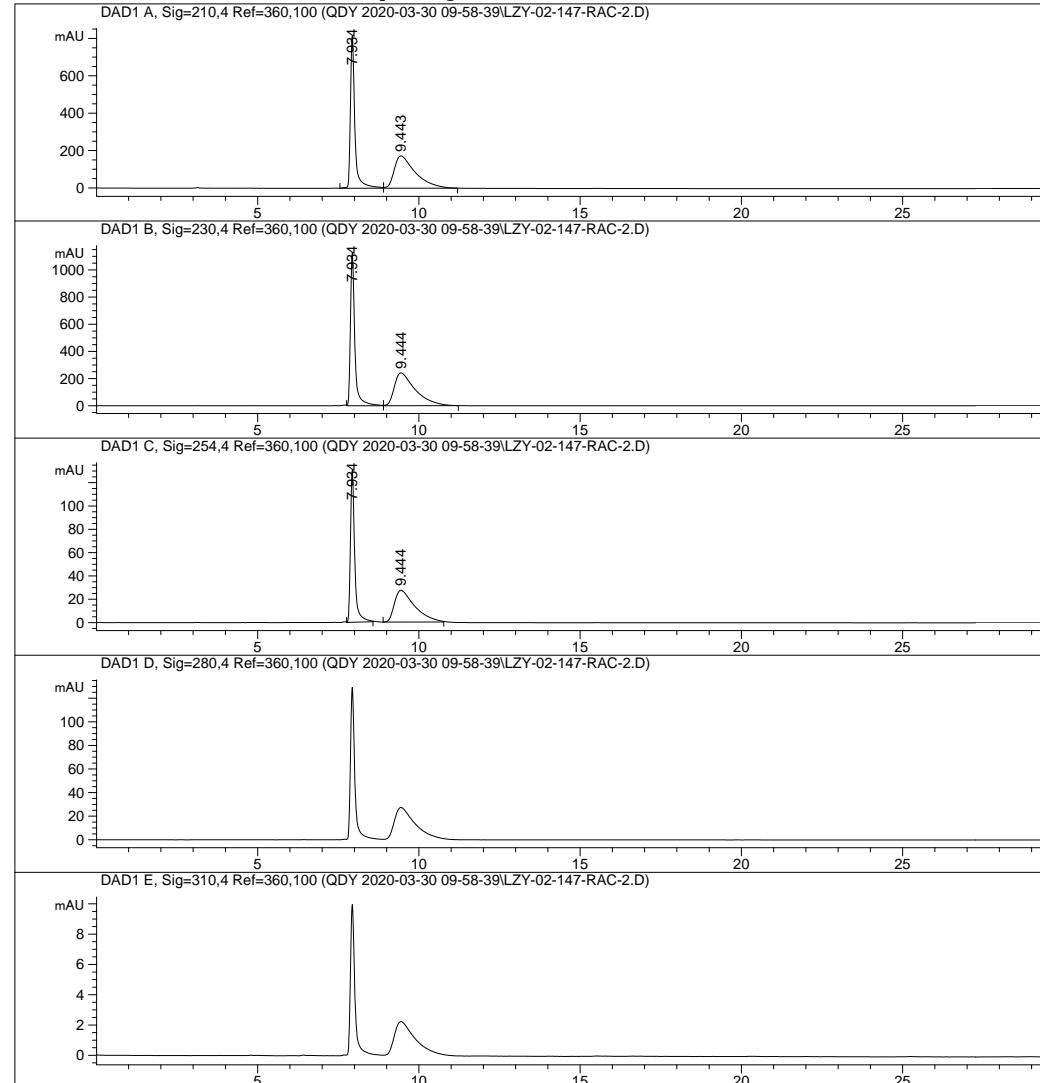
Totals : 7904.12643 516.39842

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



=====
 Acq. Operator : Seq. Line : 19
 Acq. Instrument : Instrument 1 Location : Vial 91
 Injection Date : 3/30/2020 4:17:45 PM Inj : 1
 Inj Volume : 5.000 μ l
 Different Inj Volume from Sequence ! Actual Inj Volume : 2.000 μ l
 Acq. Method : C:\CHEM32\1\DATA\QDY 2020-03-30 09-58-39\AD-007-30.M
 Last changed : 3/30/2020 4:16:53 PM
 (modified after loading)
 Analysis Method : C:\CHEM32\1\METHODS\IC-10-10.M
 Last changed : 6/25/2018 2:48:18 PM
 Additional Info : Peak(s) manually integrated



===== Area Percent Report =====

Sorted By : Signal
 Multiplier : 1.0000
 Dilution : 1.0000
 Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.934	VV	0.1418	7815.05322	815.45728	49.8456
2	9.443	VB	0.6602	7863.45947	173.02028	50.1544

Totals : 1.56785e4 988.47755

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.934	VV	0.1420	1.06001e4	1124.64417	49.3228
2	9.444	VB	0.6535	1.08912e4	241.77637	50.6772

Totals : 2.14912e4 1366.42053

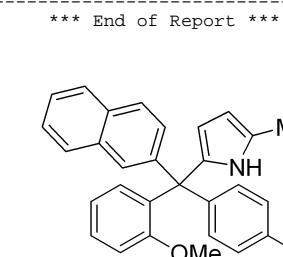
Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	7.934	VB	0.1365	1190.39087	130.45436	49.9835
2	9.444	BB	0.6411	1191.17749	27.29478	50.0165

Totals : 2381.56836 157.74914

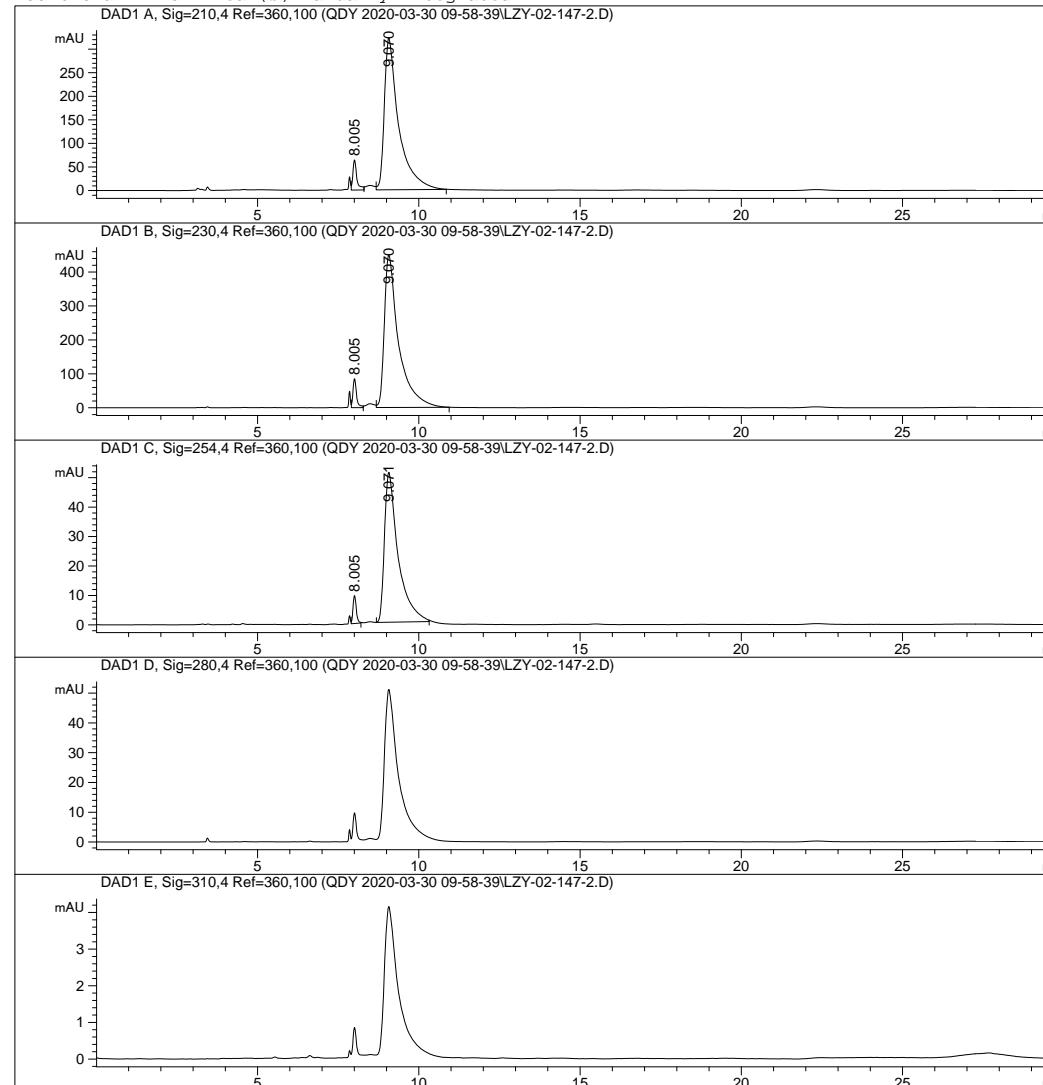
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 20
Acq. Instrument : Instrument 1             Location : Vial 92
Injection Date : 3/30/2020 4:49:03 PM        Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 2.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-03-30 09-58-39\AD-007-30.M
Last changed : 3/30/2020 4:16:53 PM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-10-10.M
Last changed : 6/25/2018 2:48:18 PM
Additional Info : Peak(s) manually integrated
```



Data File C:\CHEM32\1\DATA\QDY 2020-03-30 09-58-39\LZY-02-147-2.D
Sample Name:

```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.005	VB	0.1259	546.06274	63.76646	5.2021
2	9.070	VB	0.4406	9950.96387	321.95505	94.7979
Totals :					1.04970e4	385.72150

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

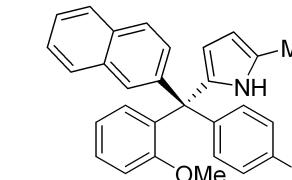
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.005	VV	0.1155	655.39935	85.40343	4.5344
2	9.070	VB	0.4366	1.37987e4	448.97861	95.4656
Totals :					1.44541e4	534.38204

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

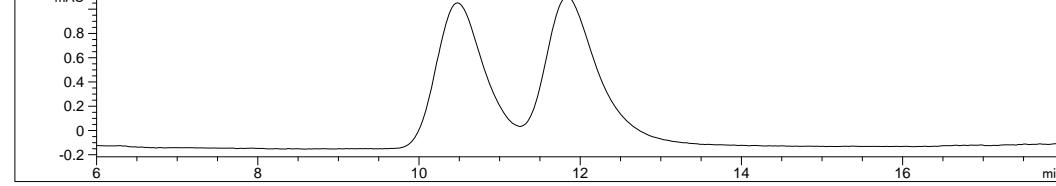
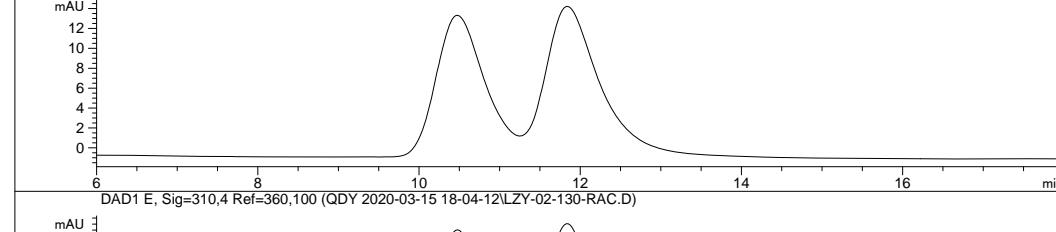
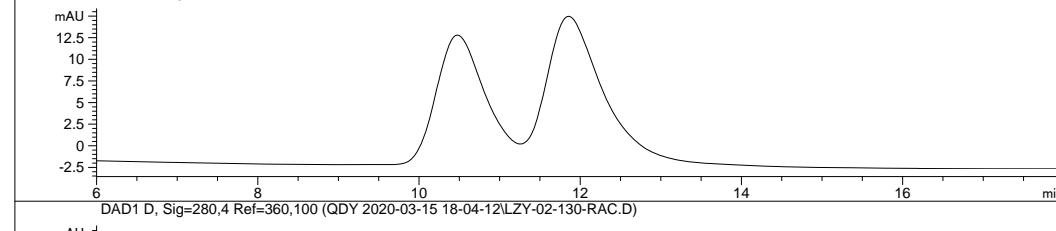
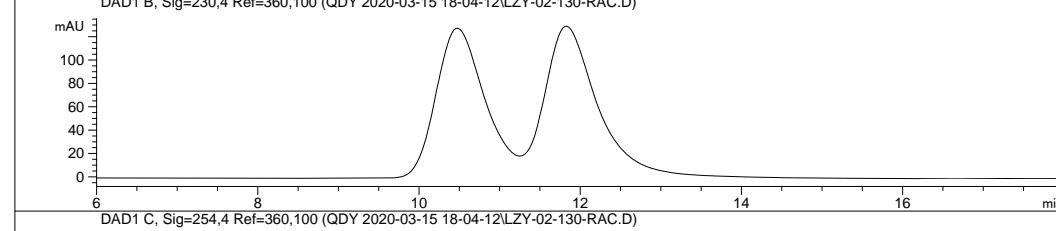
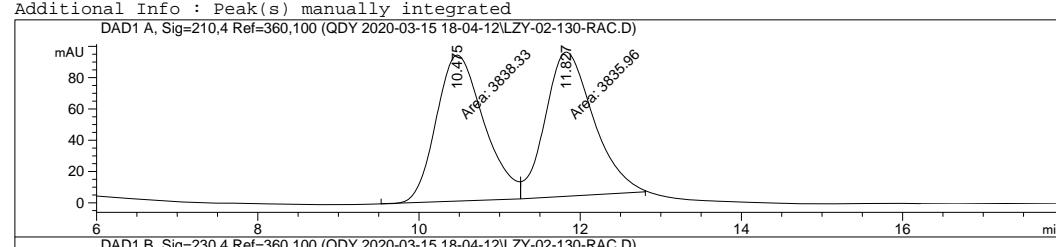
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	8.005	VB	0.1078	66.51704	9.48703	4.2567
2	9.071	BB	0.4228	1496.14343	50.92101	95.7433
Totals :					1562.66047	60.40804

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100



```
=====
Acq. Operator :                               Seq. Line : 2
Acq. Instrument : Instrument 1             Location : Vial 81
Injection Date : 3/15/2020 6:17:25 PM       Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !      Actual Inj Volume : 2.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-03-15 18-04-12\OD-007-30.M
Last changed : 3/15/2020 6:16:32 PM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 7:28:44 PM        (modified after loading)
Additional Info : Peak(s) manually integrated
```



```
=====
                                         Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	10.475	MM	0.6854	3838.33325	93.34061	50.0154
2	11.827	MM	0.6933	3835.96191	92.21267	49.9846

Totals : 7674.29517 185.55328

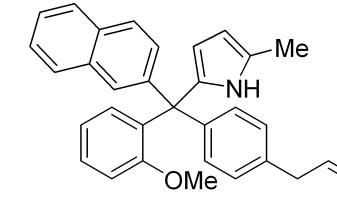
Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

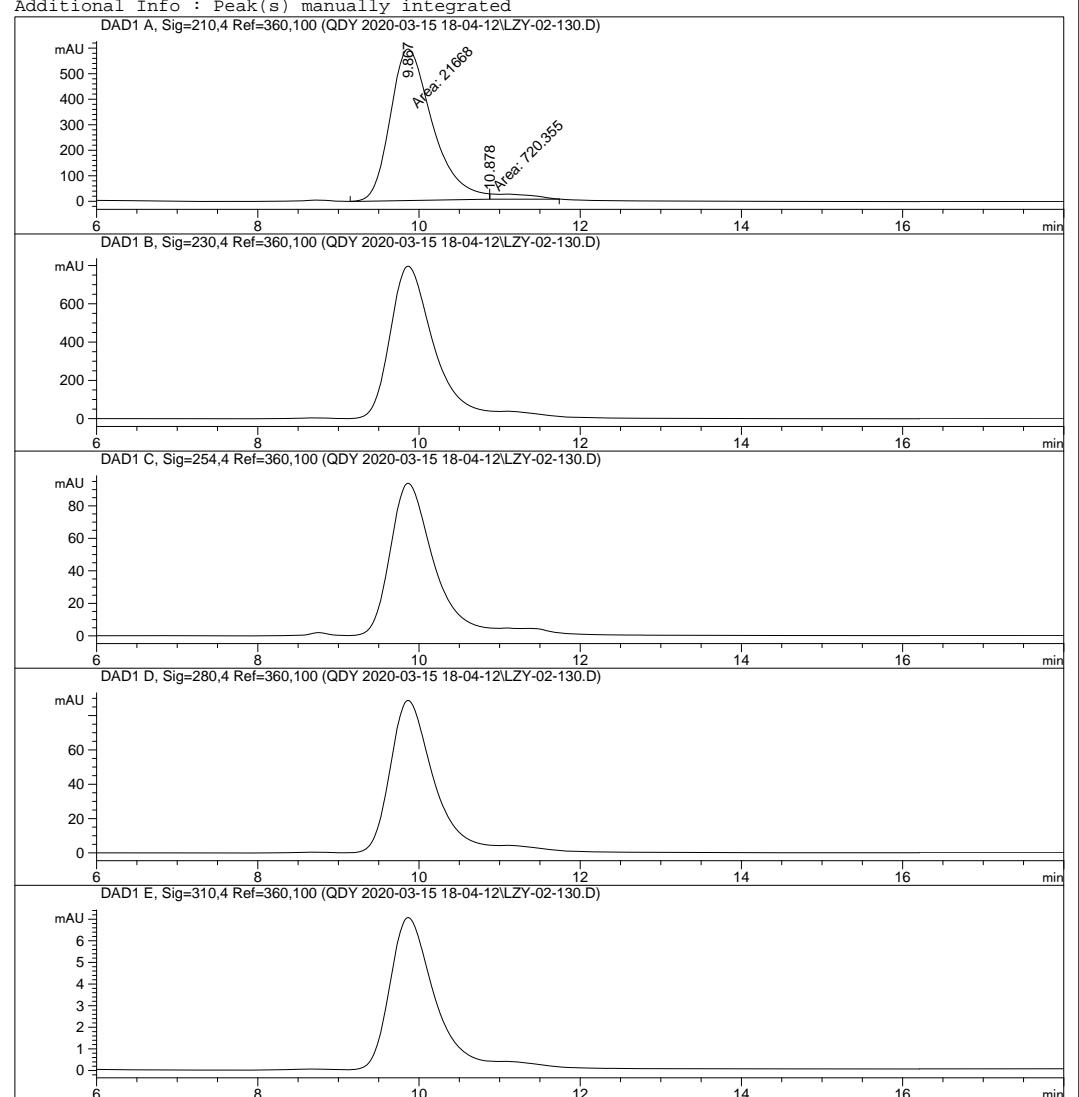
Signal 5: DAD1 E, Sig=310,4 Ref=360,100

===== *** End of Report *** =====



Sample Name:

```
=====
Acq. Operator :                               Seq. Line : 3
Acq. Instrument : Instrument 1             Location : Vial 82
Injection Date : 3/15/2020 6:48:27 PM        Inj : 1
                                                Inj Volume : 5.000 µl
Different Inj Volume from Sequence !       Actual Inj Volume : 2.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-03-15 18-04-12\OD-007-30.M
Last changed : 3/15/2020 6:16:32 PM        (modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 7:28:44 PM        (modified after loading)
Additional Info : Peak(s) manually integrated
```



Data File C:\CHEM32\1\DATA\QDY 2020-03-15 18-04-12\LZY-02-130.D
Sample Name:

```
=====
Area Percent Report
=====
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: DAD1 A, Sig=210,4 Ref=360,100
Peak RetTime Type Width Area Height Area
# [min] [min] [mAU*s] [mAU] %
-----|-----|-----|-----|-----|-----|
1 9.867 MM 0.6083 2.16680e4 593.68268 96.7825
2 10.878 MM 0.4773 720.35461 20.78572 3.2175

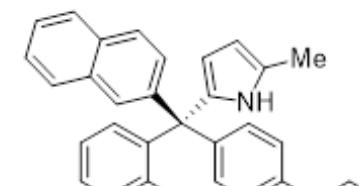
Totals : 2.23884e4 614.46840
```

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

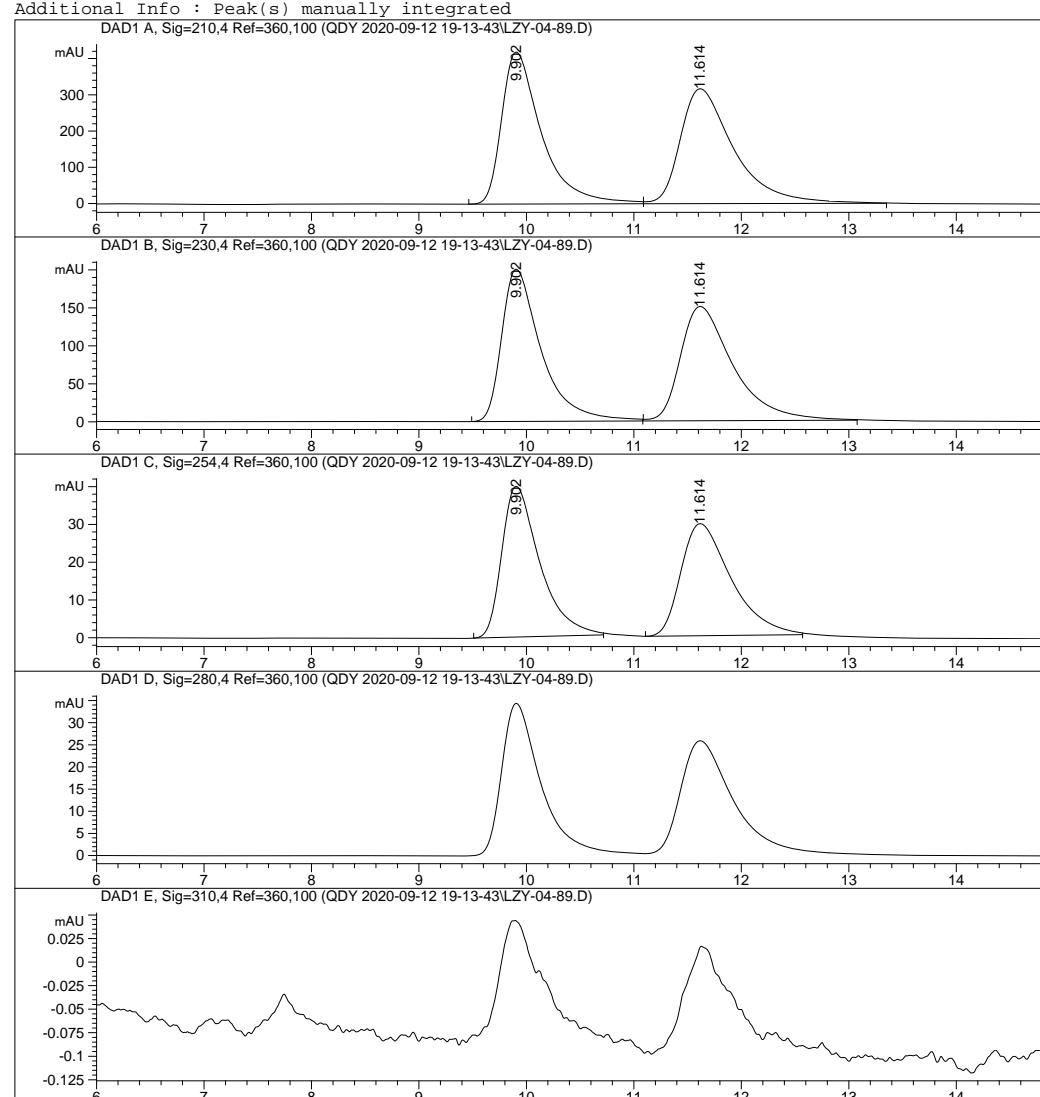
Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100

=====
*** End of Report ***

Sample Name:

```
=====
Acq. Operator : Seq. Line : 6
Acq. Instrument : Instrument 1 Location : Vial 92
Injection Date : 9/12/2020 9:11:20 PM Inj : 1
Inj Volume : 5.000 µl
Different Inj Volume from Sequence ! Actual Inj Volume : 2.000 µl
Acq. Method : C:\CHEM32\1\DATA\QDY 2020-09-12 19-13-43\OD-007-15.M
Last changed : 9/12/2020 9:10:28 PM
(modified after loading)
Analysis Method : C:\CHEM32\1\METHODS\IC-007-90.M
Last changed : 11/24/2020 9:09:57 PM
(modified after loading)
Additional Info : Peak(s) manually integrated
```



Sample Name:

```
=====
Area Percent Report
=====
```

```
Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs
```

Signal 1: DAD1 A, Sig=210,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.902	BV	0.3826	1.08027e4	419.20486	49.5885
2	11.614	VB	0.5184	1.09819e4	316.84763	50.4115
Totals :				2.17846e4	736.05249	

Signal 2: DAD1 B, Sig=230,4 Ref=360,100

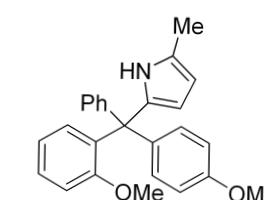
Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.902	BB	0.3798	5116.91504	200.43622	49.9390
2	11.614	BB	0.5097	5129.41602	150.46278	50.0610
Totals :				1.02463e4	350.89900	

Signal 3: DAD1 C, Sig=254,4 Ref=360,100

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	9.902	BB	0.3681	978.92847	39.92512	50.4088
2	11.614	BB	0.4906	963.04956	29.67351	49.5912
Totals :				1941.97803	69.59863	

Signal 4: DAD1 D, Sig=280,4 Ref=360,100

Signal 5: DAD1 E, Sig=310,4 Ref=360,100

**2a-Me**