

Synthesis and glycosidase inhibition of *N*-substituted derivatives of 1,4-dideoxy-1,4-imino-D-mannitol (DIM)

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Supplementary Information

Experimental data for
compound 13b-i, 14b-i, 16b-g, 17b-g, 19b-h, 20b-h, 28b-g and 29b-g
and
NMR spectra and Infrared spectra for
compound 2, 7-11, 13a-i, 14a-i, 16a-g and 17a-g

1. Experimental data for compounds

***N*-Ethyl-1,4-dideoxy-2,3,5,6-di-*O*-isopropylidene-1,4-imino-D-mannitol (13b)**

According to the general reductive amination procedures, product **13b** (99.5 mg, 89% yield) was obtained from **10** (0.10 g, 0.41 mmol) and 40% aq. acetaldehyde (142.6 μ L, 1.23 mmol) as a light yellow syrup. $[\alpha]_D^{27}$ -41.0 (c 1.80 in CH_2Cl_2); $\nu_{\text{max}}/\text{cm}^{-1}$ 2985 s, 2933 s, 2800 w, 1704 w, 1455 w, 1380 s, 1267 m, 1208 vs, 1157 s, 1103 m, 1058 s, 991 w, 931 w, 860 m, 844 m, 514 w; δ_{H} (500 MHz; CDCl_3) 4.60-4.56 (2H, m), 4.46 (1H, td, $J = 7.5$ Hz, 2.0 Hz), 4.19 (1H, t, $J = 7.8$ Hz), 3.99 (1H, t, $J = 7.5$ Hz), 3.26 (1H, d, $J = 11.1$ Hz), 3.11 (1H, dq, $J = 12.4$ Hz, 7.5 Hz), 2.62 (1H, d, $J = 2.3$ Hz), 2.15-2.08 (2H, m), 1.45 (3H, s), 1.43 (3H, s), 1.33 (3H, s), 1.28 (3H, s), 1.04 (3H, t, $J = 7.5$ Hz); δ_{C} (125 MHz; CDCl_3) 110.9, 107.1, 81.2, 77.6, 75.2, 67.6, 65.8, 58.5, 47.8, 26.2, 25.8, 24.7, 24.0, 12.5; HRMS(ESI) calcd for $\text{C}_{14}\text{H}_{26}\text{NO}_4^+$ $[\text{M}+\text{H}]^+$ 272.18563, found 272.18515.

***N*-Ethyl-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (14b)**

According to the general deprotection procedures, product **14b** (16.7 mg, 100% yield) was obtained from **13b** (20.0 mg, 0.07 mmol) as a white solid. M.p. 142-143 $^{\circ}\text{C}$; $[\alpha]_D^{27}$ -36.8 (c 0.35 in CH_2Cl_2); $\nu_{\text{max}}/\text{cm}^{-1}$ 3420 vs, 2926 vw, 1653 w, 1132 w, 1098 w, 1044 m, 625 vw; δ_{H} (500 MHz; D_2O) 4.57-4.51 (2H, m), 4.22 (1H, dd, $J = 11.0$ Hz, 5.0 Hz), 3.83 (2H, m), 3.65 (1H, t, $J = 4.1$ Hz), 3.59 (3H, m), 3.19-3.12 (1H, m), 1.35 (3H, t, $J = 7.3$ Hz); δ_{C} (125 MHz; D_2O) 70.8, 68.5, 68.1, 67.3, 62.5, 54.6, 49.3, 9.3; HRMS(ESI) calcd for $\text{C}_8\text{H}_{18}\text{NO}_4^+$ $[\text{M}+\text{H}]^+$ 192.12303, found 192.12280.

***N*-Propyl-1,4-dideoxy-2,3,5,6-di-*O*-isopropylidene-1,4-imino-D-mannitol (13c)**

According to the general reductive amination procedures, product **13c** (45.2 mg, 77% yield) was obtained from **10** (50.0 mg, 0.21 mmol) and propanal (45.7 μ L, 0.63 mmol) as a light yellow syrup. $[\alpha]_D^{28}$ -57.1 (c 1.90 in CH_2Cl_2); $\nu_{\text{max}}/\text{cm}^{-1}$ 2985 s, 2934 s, 2874 m, 2783 w, 1709 w, 1456 m, 1380 s, 1267 m, 1208 vs, 1157 s, 1098 m, 1059 s, 990 w, 935 w, 860 s, 842 m, 514 w; δ_{H} (400 MHz; CDCl_3) 4.59-4.54 (2H, m), 4.46 (1H, td, $J = 7.2$ Hz, 1.6 Hz), 4.21 (1H, t, $J = 8.0$ Hz), 3.98 (1H, t, $J = 7.6$ Hz), 3.25 (1H, d, $J = 10.8$ Hz), 3.05-2.98 (1H, m), 2.61 (1H, d, $J = 2.8$ Hz), 2.07-1.99 (2H, m), 1.51-1.42 (8H, m), 1.32 (3H, s), 1.27 (3H, s), 0.89 (3H, t, $J = 7.4$ Hz); δ_{C} (100 MHz; CDCl_3) 110.0, 107.1, 81.1, 77.7, 75.3, 67.9, 65.7, 59.1, 56.0, 26.2, 25.9, 24.9, 24.0, 20.9, 11.8; HRMS(ESI) calcd for $\text{C}_{15}\text{H}_{28}\text{NO}_4^+$ $[\text{M}+\text{H}]^+$ 286.20168, found 286.20103.

***N*-Propyl-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (14c)**

According to the general deprotection procedures, product **14c** (25.3 mg, 100% yield) was obtained from **13c** (30.0 mg, 0.10 mmol) as a light yellow syrup. $[\alpha]_D^{27}$ -36.1 (c 0.85 in CH_2Cl_2); $\nu_{\text{max}}/\text{cm}^{-1}$ 3392 vs, 2927 m, 2856 w, 1653 w, 1405 w, 1108 w, 1044 m, 560 w; δ_{H} (400 MHz; D_2O) 4.55-4.51 (2H, m), 4.23-4.22 (1H, m), 3.88-3.80 (2H, m), 3.67 (1H, s), 3.56 (2H, d, $J = 7.3$ Hz), 3.40-3.33 (1H, m), 3.10-3.03 (1H, m), 1.77-1.75 (2H, m), 0.98 (3H, t, $J = 7.2$ Hz); δ_{C} (100 MHz; D_2O) 70.7, 68.8, 68.6, 67.3, 62.6, 55.0, 55.4, 18.1, 10.1; HRMS(ESI) calcd for $\text{C}_9\text{H}_{20}\text{NO}_4^+$ $[\text{M}+\text{H}]^+$ 203.13868, found 206.13845.

***N*-Butyl-1,4-dideoxy-2,3,5,6-di-*O*-isopropylidene-1,4-imino-D-mannitol (13d)**

According to the general reductive amination procedures, product **13d** (51.1 mg, 87% yield) was obtained from **10** (48.7 mg, 0.20 mmol) and n-butanal (54.1 μ L, 0.60 mmol) as a colorless syrup. $[\alpha]_D^{24}$ -46.8 (c 1.12 in CH_2Cl_2); $\nu_{\text{max}}/\text{cm}^{-1}$ 3032 w, 2985 m, 2959 m, 2933 s, 2873 w, 2783 w, 1707 w, 1457 w, 1370 s, 1266 m, 1208 vs, 1156 s, 1089 m, 1058 m, 860 m, 844 m, 514 w; δ_{H} (400 MHz; CDCl_3) 4.59-4.54 (2H, m), 4.46 (1H, t, $J = 7.2$ Hz), 4.20 (1H, t, $J = 7.6$ Hz), 3.98 (1H, t, $J = 7.6$ Hz), 3.25 (1H, d, $J = 11.0$ Hz), 3.11-3.04 (1H, m), 2.59 (1H, d, $J = 2.0$ Hz), 2.06-1.98 (2H, m), 1.43-1.27 (16H, m), 0.90 (3H, t, $J = 7.2$ Hz); δ_{C} (100 MHz; CDCl_3) 111.0, 107.1, 81.1, 77.7, 75.3, 68.1, 65.7, 59.1, 54.0, 30.0, 26.3, 25.9, 24.9, 24.0, 20.6, 14.1; HRMS(ESI) calcd for $\text{C}_{16}\text{H}_{30}\text{NO}_4^+$ $[\text{M}+\text{H}]^+$ 300.21693, found 300.21697.

***N*-Butyl-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (14d)**

According to the general deprotection procedures, product **14d** (13.6 mg, 99% yield) was obtained from **13d** (16.0 mg, 0.05 mmol) as a light yellow syrup. $[\alpha]_D^{26}$ -41.4 (c 0.35 in CH_3OH); $\nu_{\text{max}}/\text{cm}^{-1}$ 3347 s, 2961 m, 2937 w, 2873 w, 1111 m, 1034 m; δ_{H} (400 MHz; CDCl_3) 4.58-4.50 (2H, m), 4.22 (1H, q, $J = 5.1$ Hz), 3.88-3.80 (2H, m), 3.67 (1H, t, $J = 4.1$ Hz), 3.56 (2H, d, $J = 7.4$ Hz), 3.44-3.37 (1H, m), 1.81-1.66 (2H, m), 1.46-1.33 (2H, m), 0.94 (3H, t, $J = 7.3$ Hz); δ_{C} (100 MHz; CDCl_3) 70.7, 68.8, 68.6, 67.3, 62.6, 55.4, 54.4, 26.4, 19.2, 12.7; ^1H HRMS(ESI) calcd for $\text{C}_{10}\text{H}_{22}\text{NO}_4^+$ $[\text{M}+\text{H}]^+$ 220.15433, found 220.15401.

***N*-Pentyl-1,4-dideoxy-2,3,5,6-di-*O*-isopropylidene-1,4-imino-D-mannitol (13e)**

According to the general reductive amination procedures, product **13e** (108.2 mg, 84% yield) was obtained from **10** (100.0 mg, 0.41 mmol) and n-pentanal (137.8 μ L, 1.23 mmol) as a colorless syrup. $[\alpha]_D^{24}$ -34.0 (c 1.97 in CH_2Cl_2); $\nu_{\text{max}}/\text{cm}^{-1}$ 2990 m, 2956 m, 2927 vs, 2858 m, 2790 w, 1684 m, 1458 m, 1379 s, 1208 s, 1155 m, 1094 m, 1044 m, 861 m, 419 w; δ_{H} (400 MHz; CDCl_3) 4.59-4.54 (2H, m), 4.46 (1H, td, $J = 7.2$ Hz, 1.6 Hz), 4.19 (1H, t, $J = 7.6$ Hz), 3.98 (1H, t, $J = 7.6$ Hz), 3.25 (1H, d, $J = 11.0$ Hz), 3.09-3.01 (1H, m), 2.59 (1H, d, $J = 2.7$ Hz), 2.06-1.97 (2H, m), 1.46-1.21 (18H, m), 0.88 (3H, t, $J = 7.2$ Hz); δ_{C} (100 MHz; CDCl_3) 111.0, 107.1, 81.1, 77.7, 75.3, 68.0, 65.7, 59.1, 54.2, 29.6, 27.4, 26.3, 25.9, 24.9, 24.0, 22.6, 14.1; HRMS(ESI) calcd for $\text{C}_{17}\text{H}_{32}\text{NO}_4^+$ $[\text{M}+\text{H}]^+$ 314.23258, found 314.23238.

***N*-Pentyl-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (14e)**

According to the general deprotection procedures, product **14e** (13.6 mg, 99% yield) was obtained from **13e** (16.0 mg, 0.05 mmol) as a light yellow syrup. $[\alpha]_D^{26}$ -41.9 (c 0.53 in CH₃OH); $\nu_{\max}/\text{cm}^{-1}$ 3362 s, 2958 m, 2937 m, 2873 w, 1653 m, 1458 w, 1113 m, 1050 m, 517 w; δ_{H} (400 MHz; D₂O) 4.53-4.47 (2H, m), 4.18 (1H, q, $J = 5.2$ Hz), 3.86-3.78 (2H, m), 3.53 (1H, s), 3.45-3.44 (2H, d, $J = 6.8$ Hz), 3.32-3.24 (1H, m), 3.02-2.95 (1H, m), 1.72-1.67 (2H, m), 1.36-1.31 (4H, m), 0.89 (3H, t, $J = 7.2$ Hz); δ_{C} (100 MHz; D₂O) 71.0, 68.8, 68.4, 67.8, 62.7, 55.4, 54.8, 28.0, 24.4, 21.5, 13.0; HRMS(ESI) calcd for C₁₁H₂₄NO₄⁺ [M+H]⁺ 234.16998, found 234.16964.

N-Hexyl-1,4-dideoxy-2,3:5,6-di-O-isopropylidene-1,4-imino-D-mannitol (13f)

According to the general reductive amination procedures, product **13f** (66.7 mg, 84% yield) was obtained from **10** (59.0 mg, 0.24 mmol) and n-hexanal (88.4 μ L, 0.72 mmol) as a colorless syrup. $[\alpha]_D^{24}$ -54.6 (c 1.86 in CH₂Cl₂); $\nu_{\max}/\text{cm}^{-1}$ 2984 m, 2956 m, 2928 s, 2859 m, 2782 w, 1456 w, 1369 s, 1265 m, 1206 vs, 1154 s, 1092 m, 1047s, 989 w, 860 m, 843 m, 419 w; δ_{H} (400 MHz; CDCl₃) 4.60-4.55 (2H, m), 4.46 (1H, td, $J = 7.3$ Hz, 1.6 Hz), 4.20 (1H, t, $J = 7.6$ Hz), 3.98 (1H, t, $J = 7.6$ Hz), 3.26 (1H, d, $J = 11.0$ Hz), 3.09-3.02 (1H, m), 2.59 (1H, d, $J = 2.6$ Hz), 2.07-1.98 (2H, m), 1.44-1.27 (20H, m), 0.88 (3H, t, $J = 6.4$ Hz); δ_{C} (100 MHz; CDCl₃) 111.0, 107.1, 81.1, 77.7, 75.3, 68.0, 65.8, 59.1, 54.2, 31.8, 27.7, 27.0, 26.3, 25.9, 24.9, 24.0, 22.6, 14.0; HRMS(ESI) calcd for C₁₈H₃₄NO₄⁺ [M+H]⁺ 328.24824, found 328.24786.

N-Hexyl-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (14f)

According to the general deprotection procedures, product **14f** (23.4 mg, 100% yield) was obtained from **13f** (27.0 mg, 0.08 mmol) as a light yellow syrup. $[\alpha]_D^{26}$ -46.9 (c 1.05 in CH₃OH); $\nu_{\max}/\text{cm}^{-1}$ 3348 vs, 2956 m, 2930 m, 2859 w, 1457 w, 1115 m, 1034 w; δ_{H} (500 MHz; D₂O) 4.57-4.51 (2H, m), 4.22 (1H, q, $J = 4.5$ Hz), 3.84 (2H, qd, $J = 11.9$ Hz, 6.2 Hz), 3.66 (1H, s), 3.55 (2H, d, $J = 7.4$ Hz), 3.42-3.35 (1H, m), 3.12-3.07 (1H, m), 1.78-1.69 (2H, m), 1.39-1.32 (6H, m), 0.87 (3H, t, $J = 5.5$ Hz); δ_{C} (125 MHz; D₂O) 70.7, 68.7, 68.6, 67.2, 62.5, 55.3, 54.5, 30.3, 25.3, 24.2, 21.6, 13.1; HRMS(ESI) calcd for C₁₂H₂₆NO₄⁺ [M+H]⁺ 248.18563, found 248.18526.

N-Heptyl-1,4-dideoxy-2,3:5,6-di-O-isopropylidene-1,4-imino-D-mannitol (13g)

According to the general reductive amination procedures, product **13g** (106.7 mg, 76% yield) was obtained from **10** (100 mg, 0.41 mmol) and n-heptanal (165.2 μ L, 1.23 mmol) as a colorless syrup. $[\alpha]_D^{27}$ -45.6 (c 1.49 in CH₂Cl₂); $\nu_{\max}/\text{cm}^{-1}$ 2985 m, 2956 m, 2928 vs, 2858 m, 2783 w, 1457 w, 1379 s, 1267 m, 1208 vs, 1156 s, 1094 m, 1049 s, 990 w, 931 s, 860 m, 844 m, 514 w; δ_{H} (500 MHz; CDCl₃) 4.58-4.54 (2H, m), 4.46 (1H, t, $J = 7.5$ Hz), 4.18 (1H, t, $J = 7.5$ Hz), 3.98 (1H, t, $J = 7.5$ Hz, H1), 3.25 (1H, d, $J = 11.0$ Hz), 3.08-3.02 (1H, m), 2.59 (1H, d, $J = 3.5$ Hz), 2.06-1.98 (2H, m), 1.43-1.27 (22H, m), 0.87 (3H, t, $J = 6.3$ Hz); δ_{C} (125 MHz; CDCl₃) 111.0, 107.1, 81.1, 77.7, 75.3, 68.0, 65.8, 59.1, 54.3, 31.9, 29.3, 27.7, 27.4, 26.3, 25.9, 24.9, 24.0, 22.7, 14.1; HRMS(ESI) calcd for C₁₉H₃₆NO₄⁺ [M+H]⁺ 342.26389, found 342.26332.

N-Heptyl-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (14g)

According to the general deprotection procedures, product **14g** (30.5 mg, 100% yield) was obtained from **13g** (35.0 mg, 0.10 mmol) as a light yellow syrup. $[\alpha]_D^{26}$ -28.7 (c 0.28 in CH₃OH); $\nu_{\max}/\text{cm}^{-1}$ 3349 s, 2956 m, 2929 s, 2859 w, 1684 w, 1458 w, 1117 w, 1041 m; δ_{H} (400 MHz; D₂O) 4.59-4.51 (2H, m), 4.23 (1H, q, $J = 5.1$ Hz), 3.90-3.81 (2H, m), 3.67 (1H, t, $J = 4.0$ Hz), 3.56 (2H, d, $J = 7.6$ Hz), 3.44-3.37 (1H, m), 3.15-3.07 (1H, m), 1.76-1.71 (2H, m), 1.37-1.30 (8H, m), 0.88 (3H, t, $J = 7.2$ Hz); δ_{C} (100 MHz; D₂O) 70.7, 68.8, 68.6, 67.3, 62.6, 55.4, 54.6, 30.7, 27.8, 24.3, 21.8, 13.3; HRMS(ESI) calcd for C₁₃H₂₈NO₄⁺ [M+H]⁺ 262.20128, found 262.20098.

N-Octyl-1,4-dideoxy-2,3:5,6-di-O-isopropylidene-1,4-imino-D-mannitol (13h)

According to the general reductive amination procedures, product **13h** (57.0 mg, 78% yield) was obtained from **10** (50.0 mg, 0.21 mmol) and n-octanal (98.4 μ L, 0.63 mmol) as a light yellow syrup. $[\alpha]_D^{26}$ -38.1 (c 0.53 in CH₂Cl₂); $\nu_{\max}/\text{cm}^{-1}$ 2986 m, 2956 m, 2927 vs, 2856 m, 2790 w, 1684 m, 1460 m, 1380 s, 1267 m, 1208 s, 1157 m, 1095 m, 1052 m, 860 m, 844 w, 511 w; δ_{H} (500 MHz; CDCl₃) 4.59-4.55 (2H, m), 4.46 (1H, td, $J = 7.5$ Hz, 1.5 Hz), 4.19 (1H, t, $J = 7.5$ Hz), 3.98 (1H, t, $J = 7.5$ Hz), 3.25 (1H, d, $J = 11.0$ Hz), 3.08-3.02 (1H, m), 2.59 (1H, d, $J = 3.0$ Hz), 2.06-1.98 (2H, m), 1.47-1.27 (24H, m), 0.87 (3H, t, $J = 6.5$ Hz); δ_{C} (125 MHz; CDCl₃) 111.0, 107.1, 81.1, 77.71, 75.3, 68.0, 65.8, 59.1, 54.3, 31.9, 29.6, 29.3, 27.8, 27.4, 26.3, 25.87, 24.9, 24.1, 22.7, 14.1; HRMS(ESI) calcd for C₂₀H₃₈NO₄⁺ [M+H]⁺ 356.27954, found 356.27921.

N-Octyl-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (14h)

According to the general deprotection procedures, product **14h** (21.9 mg, 100% yield) was obtained from **13h** (25.0 mg, 0.07 mmol) as a light yellow syrup. $[\alpha]_D^{26}$ -39.2 (c 0.39 in CH₃OH); $\nu_{\max}/\text{cm}^{-1}$ 3377 vs, 2956 m, 2926 s, 2856 m, 1647 w, 1456 w, 1119 m, 1041 w, 517 w; δ_{H} (400 MHz; D₂O) 4.58-4.50 (2H, m), 4.22 (1H, q, $J = 5.1$ Hz), 3.88-3.80 (2H, m), 3.66 (1H, t, $J = 4.2$ Hz), 3.56 (2H, d, $J = 7.4$ Hz), 3.43-3.36 (1H, m), 3.13-3.06 (1H, m), 1.79-1.67 (2H, m), 1.36-1.28 (10H, m), 0.86 (3H, t, $J = 6.4$ Hz); δ_{C} (100 MHz; D₂O) 70.7, 68.7, 68.6, 67.3, 62.5, 55.3, 54.6, 30.9, 28.11, 28.08, 25.7, 24.3, 21.9, 13.4; HRMS(ESI) calcd for C₁₄H₃₀NO₄⁺ [M+H]⁺ 276.21693, found 276.21684.

N-Nonyl-1,4-dideoxy-2,3:5,6-di-O-isopropylidene-1,4-imino-D-mannitol (13i)

According to the general reductive amination procedures, product **13i** (64.7 mg, 71% yield) was obtained from **10** (60.0 mg, 0.25 mmol) and n-nonanal (128.9 μ L, 0.75 mmol) as a colorless syrup. $[\alpha]_D^{20}$ -45.4 (c 0.55 in CH₂Cl₂); $\nu_{\max}/\text{cm}^{-1}$ 2984 w, 2925 vs, 2855 m, 2783 w, 1457 w, 1369 m, 1266 w, 1027 s, 1155 m, 1050 m, 1105 vs, 860 w, 843 w; δ_{H} (500 MHz; CDCl₃)

4.59-4.55 (2H, m), 4.46 (1H, t, $J = 7.0$ Hz), 4.19 (1H, t, $J = 7.5$ Hz), 3.98 (1H, t, $J = 7.5$ Hz), 3.25 (1H, d, $J = 11.5$ Hz), 3.08-3.02 (1H, m), 2.59 (1H, d, $J = 3.0$ Hz), 2.06-1.98 (2H, m), 1.43 (8H, s), 1.33-1.25 (18H, m), 0.87 (3H, t, $J = 6.5$ Hz); δ_c (125 MHz; CDCl₃) 111.0, 107.1, 81.1, 77.7, 75.3, 68.0, 65.7, 59.1, 54.3, 31.9, 29.6, 29.3, 27.7, 27.4, 26.3, 25.9, 24.9, 24.1, 22.7, 14.1; HRMS(ESI) calcd for C₂₁H₄₀NO₄⁺ [M+H]⁺ 370.29519, found 370.29464.

***N*-Nonyl-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (14i)**

According to the general deprotection procedures, product **14i** (26.4 mg, 100% yield) was obtained from **13i** (30.1 mg, 0.08 mmol) as a light yellow syrup. $[\alpha]_D^{26} -32.1$ (c 0.86 in CH₃OH); $\nu_{\max}/\text{cm}^{-1}$ 3308 s, 2925 vs, 2855 m, 1116 m, 1044 w, 419 m; δ_H (400 MHz; D₂O) 4.53-4.50 (2H, m), 4.20 (1H, m), 3.82 (2H, s), 3.64 (1H, s), 3.53 (2H, d, $J = 6.7$ Hz), 3.38-3.34 (1H, m), 3.11-3.07 (1H, m), 1.73 (2H, s), 1.32-1.26 (12H, m), 0.84 (3H, s); δ_c (100 MHz; D₂O) 70.7, 68.7, 68.5, 67.2, 62.5, 55.3, 54.5, 31.0, 28.3, 28.2, 28.1, 25.6, 24.2, 21.9, 13.3; HRMS(ESI) calcd for C₁₅H₃₂NO₄⁺ [M+H]⁺ 290.23258, found 290.23267.

***N*-(But-3-en-1-yl)-1,4-dideoxy-2,3:5,6-di-O-isopropylidene-1,4-imino-D-mannitol (16b)**

According to the general nucleophilic substitution procedures, product **16b** (63.1 mg, 86% yield) was obtained from **10** (60.0 mg, 0.25 mmol) and 1-bromo-4-butene (38.3 μ L, 0.38 mmol) as a colorless syrup. $[\alpha]_D^{26} -53.1$ (c 0.36 in CH₂Cl₂); $\nu_{\max}/\text{cm}^{-1}$ 2984 m, 2934 m, 2797 w, 1370 s, 1267 w, 1207 vs, 1155 m, 1098 m, 1047 s, 990 w, 912 w, 860 m, 844 w; δ_H (400 MHz; CDCl₃) 5.81 (1H, ddt, $J = 17.0$ Hz, 10.3 Hz, 6.5 Hz), 5.06-4.95 (2H, m), 4.60-4.55 (2H, m), 4.46 (1H, td, $J = 6.4$ Hz, 1.6 Hz), 4.21 (1H, t, $J = 8.0$ Hz), 3.98 (1H, t, $J = 7.6$ Hz), 3.27 (1H, d, $J = 10.8$ Hz), 3.20-3.13 (1H, m), 2.67 (1H, dd, $J = 4.3$ Hz, 1.1 Hz), 2.24-2.11 (4H, m), 1.43 (6H, s), 1.32 (3H, s), 1.27 (3H, s); δ_c (100 MHz; CDCl₃) 136.7, 115.3, 111.1, 107.1, 81.0, 77.7, 75.3, 67.6, 65.7, 59.1, 53.3, 31.9, 26.2, 25.8, 24.8, 24.0; HRMS(ESI) calcd for C₁₆H₂₈NO₄⁺ [M+H]⁺ 298.20128, found 298.20101.

***N*-(But-3-en-1-yl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (17b)**

According to the general deprotection procedures, product **17b** (34.1 mg, 100% yield) was obtained from **16b** (40.0 mg, 0.13 mmol) as a light yellow syrup. $[\alpha]_D^{26} -30.6$ (c 0.60 in CH₃OH); $\nu_{\max}/\text{cm}^{-1}$ 3356 vs, 2926 w, 2379 w, 2320 w, 1732 w, 1510 w, 1458 w, 1338 w, 1127 m, 1033 w, 1002 w, 933 w; δ_H (400 MHz; D₂O) 5.83 (1H, ddt, $J = 17.1$ Hz, 10.3 Hz, 6.7 Hz), 5.30-5.21 (2H, m), 4.59-4.52 (2H, m), 4.24 (1H, q, $J = 5.0$ Hz), 3.90 (1H, dd, $J = 11.9$ Hz, 5.9 Hz), 3.83 (1H, dd, $J = 11.8$ Hz, 4.8 Hz), 3.72 (1H, t, $J = 4.3$ Hz), 3.59-3.51 (3H, m), 3.25-3.18 (1H, m), 2.61-2.49 (2H, m); δ_c (100 MHz; D₂O) 132.4, 118.8, 70.6, 69.3, 68.6, 67.1, 62.6, 55.4, 53.6, 28.9; HRMS(ESI) calcd for C₁₀H₂₀NO₄⁺ [M+H]⁺ 218.13868, found 218.13830.

***N*-(Pent-4-en-1-yl)-1,4-dideoxy-2,3:5,6-di-O-isopropylidene-1,4-imino-D-mannitol (16c)**

According to the general nucleophilic substitution procedures, product **16c** (908.7 mg, 71% yield) was obtained from **10** (1.0 g, 4.11 mmol) and 1-bromo-5-pentene (0.73 mL, 6.17 mmol) as a colorless syrup. $[\alpha]_D^{24} -26.1$ (c 1.72 in CH₂Cl₂); $\nu_{\max}/\text{cm}^{-1}$ 3392 w, 3081 w, 2986 s, 2935 s, 2801 w, 1698 s, 1642 m, 1458 w, 1381 s, 1268 m, 1209 vs, 1157 m, 1097 m, 1056 s, 991 w, 916 s, 860 m, 517 w; δ_H (400 MHz; CDCl₃) 5.88-5.78 (1H, m), 5.03-5.02 (1H, m), 4.95-4.92 (1H, m), 4.60-4.55 (2H, m), 4.46 (1H, td, $J = 7.8$ Hz, 1.7 Hz), 3.98 (1H, t, $J = 7.5$ Hz), 3.98 (1H, t, $J = 7.6$ Hz), 3.26 (1H, d, $J = 11.0$ Hz), 3.08 (1H, dt, $J = 12.2$ Hz, 8.5 Hz), 2.60 (1H, d, $J = 2.7$ Hz), 2.17-1.99 (4H, m), 1.60-1.51 (2H, m), 1.43 (6H, s), 1.32 (3H, s), 1.27 (3H, s); δ_c (100 MHz; CDCl₃) 138.8, 114.4, 111.0, 107.1, 81.1, 77.7, 75.3, 68.1, 65.7, 59.1, 53.7, 31.5, 27.1, 26.3, 25.9, 24.8, 24.0; HRMS(ESI) calcd for C₁₇H₃₀NO₄⁺ [M+H]⁺ 312.21693, found 312.21667.

***N*-(Pent-4-en-1-yl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (17c)**

According to the general deprotection procedures, product **17c** (85.0 mg, 99% yield) was obtained from **16c** (100.0 mg, 0.32 mmol) as a light yellow syrup. $[\alpha]_D^{26} -16.8$ (c 0.75 in CH₃OH); $\nu_{\max}/\text{cm}^{-1}$ 3420 vs, 2973 w, 1684 w, 1559 w, 1402 m, 1127 w, 1038 w, 594 w; δ_H (400 MHz; D₂O) 5.89-5.82 (1H, m), 5.14-5.06 (2H, m), 4.53-4.50 (2H, m), 4.20 (1H, s), 3.86-3.82 (2H, m), 3.68-3.62 (1H, m), 3.56-3.54 (2H, m), 3.40-3.36 (1H, m), 3.13-3.09 (1H, m), 2.16-2.15 (2H, m), 1.95-1.78 (2H, m); δ_c (100 MHz; D₂O) 137.0, 116.0, 70.7, 68.8, 68.6, 67.2, 62.5, 55.4, 53.9, 29.8, 23.5; HRMS(ESI) calcd for C₁₁H₂₂NO₄⁺ [M+H]⁺ 232.15433, found 232.15405.

***N*-(Hex-5-en-1-yl)-1,4-dideoxy-2,3:5,6-di-O-isopropylidene-1,4-imino-D-mannitol (16d)**

According to the general nucleophilic substitution procedures, product **16d** (40.0 mg, 56% yield) was obtained from **10** (54.0 mg, 0.22 mmol) and 1-bromo-6-hexene (44.5 μ L, 0.33 mmol) as a colorless syrup. $[\alpha]_D^{26} -51.1$ (c 1.17 in CH₂Cl₂); $\nu_{\max}/\text{cm}^{-1}$ 2984 m, 2930 s, 2859 w, 2783 w, 1456 w, 1370 m, 1266 w, 1207 vs, 1155 m, 1051 s, 990 w, 910 w, 860 m, 844 w; δ_H (400 MHz; CDCl₃) 5.85-5.75 (1H, m), 5.02-4.92 (2H, m), 4.59-4.55 (2H, m), 4.47 (1H, t, $J = 7.2$ Hz), 3.98 (1H, t, $J = 7.6$ Hz), 3.25 (1H, d, $J = 11.0$ Hz), 3.11-3.04 (1H, dt, $J = 12.0$ Hz, 8.0 Hz), 2.60 (1H, d, $J = 2.7$ Hz), 2.07-2.00 (4H, m), 1.47-1.35 (10H, m), 1.33 (3H, s), 1.27 (3H, s); δ_c (100 MHz; CDCl₃) 139.0, 114.3, 111.0, 107.1, 81.1, 77.7, 75.3, 68.1, 65.7, 59.1, 54.0, 33.7, 27.2, 26.7, 26.3, 25.9, 24.8, 23.9; HRMS(ESI) calcd for C₁₈H₃₂NO₄⁺ [M+H]⁺ 326.23258, found 326.23225.

***N*-(Hex-5-en-1-yl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (17d)**

According to the general deprotection procedures, product **17d** (25.9 mg, 100% yield) was obtained from **16d** (30.0 mg, 0.09 mmol) as a light yellow syrup. $[\alpha]_D^{26} -40.9$ (c 0.44 in CH₃OH); $\nu_{\max}/\text{cm}^{-1}$ 3362 vs, 2943 m, 2836 w, 1640 m, 1447 m, 1126 m, 1024 s, 918 w, 556 m; δ_H (400 MHz; D₂O) 5.88 (1H, ddt, $J = 17.1$ Hz, 10.3 Hz, 6.7 Hz), 5.11-5.03 (2H, m), 4.57-4.50 (2H, m), 4.22 (1H, q, $J = 5.0$ Hz), 3.88-3.79 (2H, m), 3.66 (1H, t, $J = 4.1$ Hz), 3.56-3.54 (2H, m), 3.45-3.35 (1H, m), 3.15-3.07 (1H,

m), 2.12 (2H, q, $J = 6.9$ Hz), 1.78-1.75 (2H, m), 1.51-1.47 (2H, m); δ_c (100 MHz; D₂O) 138.6, 115.0, 70.6, 68.7, 68.5, 67.2, 62.5, 55.3, 54.17, 32.2, 24.8, 23.6; HRMS(ESI) calcd for C₁₂H₂₄NO₄⁺ [M+H]⁺ 246.16998, found 246.16957.

***N*-(Hept-6-en-1-yl)-1,4-dideoxy-2,3:5,6-di-*O*-isopropylidene-1,4-imino-D-mannitol (16e)**

According to the general nucleophilic substitution procedures, product **16e** (398 mg, 87% yield) was obtained from **10** (328.0 mg, 1.35 mmol) and 1-bromo-7-heptene (358.6 mg, 2.03 mmol) as a colorless syrup. $[\alpha]_D^{26}$ -51.5 (c 1.13 in CH₂Cl₂); $\nu_{\max}/\text{cm}^{-1}$ 2985 m, 2930 s, 2859 w, 2783 w, 1643 vw, 1458 w, 1380 s, 1267 m, 1208 vs, 1156 s, 1102 m, 1051 s, 991 m, 910 w, 860 m, 844 m, 514 vw; δ_H (400 MHz; CDCl₃) 5.85-5.75 (1H, m), 5.00-4.90 (2H, m), 4.59-4.54 (2H, m), 4.46(1H, t, $J = 7.3$ Hz), 4.18 (1H, t, $J = 7.8$ Hz), 3.98 (1H, t, $J = 7.6$ Hz), 3.25 (1H, d, $J = 11.0$ Hz), 3.20-3.13 (1H, dt, $J = 12.0$ Hz, 8.1 Hz), 2.59 (1H, d, $J = 2.4$ Hz), 2.06-1.98 (4H, m), 1.51-1.27 (18H, m); δ_c (100 MHz; CDCl₃) 139.0, 114.2, 111.0, 107.1, 81.1, 77.7, 75.3, 68.0, 65.7, 59.1, 54.1, 33.8, 28.8, 27.6, 26.87, 26.28, 25.86, 24.9, 24.0; HRMS(ESI) calcd for C₁₉H₃₄NO₄⁺ [M+H]⁺ 340.24824, found 340.24800.

***N*-(Hept-6-en-1-yl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (17e)**

According to the general deprotection procedures, product **17e** (20.0 mg, 100% yield) was obtained from **16e** (22.0 mg, 0.06 mmol) as a light yellow syrup. $[\alpha]_D^{26}$ -40.5 (c 0.77 in CH₃OH); $\nu_{\max}/\text{cm}^{-1}$ 3349 vs, 2935 m, 2859 w, 1640 w, 1417 m, 1126 m, 1043 m, 998 w, 911 w, 554 w; δ_H (400 MHz; D₂O) 5.93 (1H, ddt, $J = 17.1$ Hz, 10.3 Hz, 6.7 Hz), 5.11-5.01 (2H, m), 4.60-4.53 (2H, m), 4.24 (1H, q, $J = 5.0$ Hz), 3.91-3.82 (2H, m), 3.68 (1H, t, $J = 4.1$ Hz), 3.58 (2H, d, $J = 7.4$ Hz), 3.46-3.39 (1H, m), 3.16-3.09 (1H, m), 2.11 (2H, q, $J = 6.5$ Hz), 1.80-1.74 (2H, m), 1.53-1.39 (4H, m); δ_c (100 MHz; D₂O) 139.6, 114.5, 70.7, 68.8, 68.6, 67.3, 62.6, 55.4, 54.5, 32.7, 24.4, 25.2, 24.2; HRMS(ESI) calcd for C₁₃H₂₆NO₄⁺ [M+H]⁺ 260.18563, found 260.18520.

***N*-(Oct-7-en-1-yl)-1,4-dideoxy-2,3:5,6-di-*O*-isopropylidene-1,4-imino-D-mannitol (16f)**

According to the general nucleophilic substitution procedures, product **16f** (134.8 mg, 64% yield) was obtained from **10** (145.0 mg, 0.60 mmol) and 1-bromo-8-octene (172.0 mg, 0.90 mmol) as a colorless syrup. $[\alpha]_D^{24}$ -36.2 (c 0.31 in CH₂Cl₂); $\nu_{\max}/\text{cm}^{-1}$ 2984 m, 2928 vs, 2657 m, 2783 w, 1456 w, 1369 m, 1267 m, 1207 s, 1156 m, 1103 w, 1050 m, 990 w, 910 w, 860 m, 843 w; δ_H (400 MHz; CDCl₃) 5.79 (1H, ddt, $J = 17.0$ Hz, 10.2 Hz, 6.6 Hz), 5.00-4.90 (2H, m), 4.59-4.54 (2H, m), 4.45 (1H, td, $J = 7.3$ Hz, 1.6 Hz), 4.19 (1H, t, $J = 7.7$ Hz), 3.97 (1H, t, $J = 7.5$ Hz), 3.24 (1H, d, $J = 11.0$ Hz), 3.20-3.13 (1H, dt, $J = 11.8$ Hz, 8.1 Hz), 2.59 (1H, d, $J = 2.8$ Hz), 2.06-1.98 (4H, m), 1.48-1.24 (20H, m); δ_c (100 MHz; CDCl₃) 139.2, 114.1, 111.0, 107.1, 81.1, 77.7, 75.3, 68.0, 65.7, 59.1, 54.2, 33.8, 29.1, 28.9, 27.7, 27.2, 26.3, 25.9, 24.9, 24.0; HRMS(ESI) calcd for C₂₀H₃₆NO₄⁺ [M+H]⁺ 354.26389, found 354.26343.

***N*-(Oct-7-en-1-yl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (17f)**

According to the general deprotection procedures, product **17f** (43.3 mg, 99% yield) was obtained from **16f** (50.0 mg, 0.14 mmol) as a light yellow syrup. $[\alpha]_D^{26}$ -36.9 (c 0.96 in CH₃OH); $\nu_{\max}/\text{cm}^{-1}$ 3368 vs, 2928 s, 2859 m, 1640 w, 1458 w, 1126 m, 1044 w, 913 w; δ_H (400 MHz; D₂O) 5.93 (1H, ddt, $J = 17.2$ Hz, 10.3 Hz, 6.7 Hz), 5.11-5.00 (2H, m), 4.59-4.52 (2H, m), 4.23 (1H, q, $J = 5.1$ Hz), 3.90-3.81 (2H, m), 3.66 (1H, t, $J = 4.2$ Hz), 3.56-3.54 (2H, m), 3.43-3.36 (1H, m), 3.14-3.06 (1H, m), 2.11 (2H, dd, $J = 12.4$ Hz, 7.8 Hz), 1.82-1.71 (2H, m), 1.47-1.40 (6H, m); δ_c (100 MHz; D₂O) 140.1, 114.2, 70.8, 68.7, 68.6, 67.4, 62.6, 55.4, 54.6, 32.9, 27.7, 27.6, 25.5, 24.3; HRMS(ESI) calcd for C₁₄H₂₈NO₄⁺ [M+H]⁺ 274.20128, found 274.20118.

***N*-(Non-8-en-1-yl)-1,4-dideoxy-2,3:5,6-di-*O*-isopropylidene-1,4-imino-D-mannitol (16g)**

According to the general nucleophilic substitution procedures, product **16g** (66.7 mg, 69% yield) was obtained from **10** (64.0 mg, 0.26 mmol) and 1-bromo-9-nonene (61.6 mg, 0.30 mmol) as a colorless syrup. $[\alpha]_D^{24}$ -23.4 (c 0.57 in CH₂Cl₂); $\nu_{\max}/\text{cm}^{-1}$ 2985 m, 2928 vs, 2856 m, 2784 vw, 1704 w, 1456 w, 1380 m, 1266 w, 1209 s, 1158 m, 1093 w, 1054 m, 910 w, 860 m; δ_H (400 MHz; CDCl₃) 5.79 (1H, ddt, $J = 17.0$ Hz, 10.2 Hz, 6.7 Hz), 5.00-4.90 (2H, m), 4.58-4.54 (2H, m), 4.45 (1H, td, $J = 7.3$ Hz, 1.6 Hz), 4.18 (1H, t, $J = 7.8$ Hz), 3.97 (1H, t, $J = 7.5$ Hz), 3.24 (1H, d, $J = 11.0$ Hz), 3.05 (1H, dt, $J = 12.1$ Hz, 8.0 Hz), 2.58 (1H, d, $J = 2.4$ Hz), 2.06-1.97 (4H, m), 1.43-1.26 (22H, m); δ_c (100 MHz; CDCl₃) 139.2, 114.1, 111.0, 107.1, 81.1, 77.7, 75.3, 68.1, 65.8, 59.1, 54.2, 33.8, 29.4, 29.1, 28.9, 27.7, 27.3, 26.3, 25.9, 24.8, 24.0; HRMS(ESI) calcd for C₂₁H₃₈NO₄⁺ [M+H]⁺ 368.27954, found 368.27932.

***N*-(Non-8-en-1-yl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (17g)**

According to the general deprotection procedures, product **17g** (22.6 mg, 99% yield) was obtained from **16g** (26.0 mg, 0.07 mmol) as a light yellow syrup. $[\alpha]_D^{26}$ -29.4 (c 0.98 in CH₃OH); $\nu_{\max}/\text{cm}^{-1}$ 3367 vs, 2928 s, 2854 m, 1653 w, 1458 m, 1125 m, 1041 w, 910 w, 726 w, 628 w; δ_H (400 MHz; D₂O) 5.95 (1H, ddt, $J = 17.1$ Hz, 10.3 Hz, 6.7 Hz), 5.11-5.00 (2H, m), 4.61-4.54 (2H, m), 4.25 (1H, q, $J = 5.1$ Hz), 3.92-3.83 (2H, m), 3.69 (1H, t, $J = 4.2$ Hz), 3.59-3.58 (2H, m), 3.46-3.39 (1H, m), 3.17-3.10 (1H, m), 2.10 (2H, q, $J = 6.5$ Hz), 1.83-1.73 (2H, m), 1.54-1.29 (8H, m); δ_c (100 MHz; D₂O) 140.4, 114.1, 70.8, 68.8, 68.6, 67.3, 62.6, 55.4, 54.6, 33.0, 28.0, 25.6, 24.3; HRMS(ESI) calcd for C₁₅H₃₀NO₄⁺ [M+H]⁺ 288.21693, found 288.21672.

***N*-(Phenethyl)-1,4-dideoxy-2,3:5,6-di-*O*-isopropylidene-1,4-imino-D-mannitol (19b)**

According to the general nucleophilic substitution procedures, product **19b** (116.1 mg, 74% yield) was obtained from **10** (110.0 mg, 0.45 mmol) and 2-phenyl-1-bromoethane (125.8 mg, 0.68 mmol) as a white solid. M.p. 86-87 °C; $[\alpha]_D^{24}$ -38.3 (c 0.42 in CH₂Cl₂); $\nu_{\max}/\text{cm}^{-1}$ 2985 m, 2922 vs, 2853 m, 2809 w, 2373 w, 1685 w, 1455 w, 1369 m, 1272 w, 1207 s, 1154 m, 1099 m, 1044 s, 991 w, 930 w, 859 w, 755 w, 700 m, 522 m; δ_H (400 MHz; CDCl₃) 7.30-7.26 (2H, m), 7.23-7.17 (3H, m),

4.64-4.57 (2H, m), 4.47 (1H, td, $J = 7.3$ Hz, 1.4 Hz), 4.06 (1H, t, $J = 7.8$ Hz), 3.93 (1H, t, $J = 7.5$ Hz), 3.41-3.34 (2H, m), 2.80-2.77 (2H, m), 2.73 (1H, d, $J = 3.8$ Hz), 2.40-2.34 (1H, m), 2.23 (1H, dd, $J = 10.9$ Hz, 4.6 Hz), 1.42 (3H, s), 1.41 (3H, s), 1.33 (3H, s), 1.29 (3H, s); δ_c (100 MHz; CDCl_3) 140.6, 128.8, 128.2, 125.9, 111.1, 107.1, 81.0, 77.7, 75.2, 67.4, 65.6, 59.2, 55.5, 34.1, 26.3, 25.9, 24.8, 23.9; HRMS(ESI) calcd for $\text{C}_{20}\text{H}_{30}\text{NO}_4^+$ [$\text{M}+\text{H}$] $^+$ 348.21693, found 348.21667.

***N*-(Phenethyl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (20b)**

According to the general deprotection procedures, product **20b** (17.5 mg, 100% yield) was obtained from **19b** (20.0 mg, 0.06 mmol) as a light yellow syrup. $[\alpha]_D^{26} -34.1$ (c 0.38 in CH_3OH); $\nu_{\text{max}}/\text{cm}^{-1}$ 3361 vs, 2923 w, 1651 w, 1460 w, 1419 w, 1127 m, 1033 w, 753 w, 703w, 503 w; δ_H (500 MHz; D_2O) 7.44-7.41 (2H, m), 7.37-7.36 (3H, m), 4.59-4.55 (1H, m), 4.53-4.52 (1H, m), 4.22-4.21 (1H, m), 3.86 (1H, dd, $J = 11.8$ Hz, 6.8 Hz), 3.79 (1H, dd, $J = 11.8$ Hz, 4.6 Hz), 3.72-3.58 (4H, m), 3.41-3.55 (1H, m), 3.18-3.07 (2H, m); δ_c (125 MHz; D_2O) 136.0, 129.1, 128.8, 127.4, 70.5, 69.3, 68.6, 67.0, 62.5, 55.4, 55.3, 30.7; HRMS(ESI) calcd for $\text{C}_{14}\text{H}_{22}\text{NO}_4^+$ [$\text{M}+\text{H}$] $^+$ 268.15433, found 268.15397.

***N*-(Phenylpropyl)-1,4-dideoxy-2,3:5,6-di-O-isopropylidene-1,4-imino-D-mannitol (19c)**

According to the general nucleophilic substitution procedures, product **19c** (117.7 mg, 72% yield) was obtained from **10** (110.0 mg, 0.45 mmol) and 3-phenyl-1-bromopropane (135.4 mg, 0.68 mmol) as a colorless syrup. $[\alpha]_D^{26} -42.8$ (c 2.1 in CH_2Cl_2); $\nu_{\text{max}}/\text{cm}^{-1}$ 3062 w, 3026 w, 2985 s, 2931 vs, 2855 m, 2785 m, 1706m, 1603 w, 1497 w, 1454 m, 1340 s, 1268 m, 1208 vs, 1157 m, 1100 m, 1049 s, 990 w, 931 w, 860 m, 842 m, 747 m, 699 m, 514 w; δ_H (400 MHz; CDCl_3) 7.28-7.25 (2H, m), 7.19-7.14 (3H, m), 4.60-4.56 (2H, m), 4.45 (2H, td, $J = 7.3$ Hz, 1.9 Hz), 4.16 (1H, t, $J = 7.8$ Hz), 3.95 (1H, t, $J = 7.5$ Hz), 3.29 (1H, d, $J = 10.9$ Hz), 3.11 (1H, dt, $J = 12.2$ Hz, 8.4 Hz), 2.75-2.68 (1H, m), 2.62-2.54 (2H, m), 2.13-2.06 (2H, m), 1.85-1.73 (2H, m), 1.45 (3H, s), 1.36 (3H, s), 1.32 (3H, s), 1.28 (3H, s); δ_c (100 MHz; CDCl_3) 142.4, 128.4, 128.3, 125.7, 111.0, 107.1, 81.1, 77.7, 75.3, 68.1, 65.8, 59.1, 53.8, 33.6, 29.5, 26.2, 25.9, 24.9, 24.0; HRMS(ESI) calcd for $\text{C}_{21}\text{H}_{32}\text{NO}_4^+$ [$\text{M}+\text{H}$] $^+$ 362.223258, found 362.23229.

***N*-(Phenylpropyl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (20c)**

According to the general deprotection procedures, product **20c** (26.3 mg, 100% yield) was obtained from **19c** (30.0 mg, 0.08 mmol) as a light yellow syrup. $[\alpha]_D^{26} -36.1$ (c 0.19 in CH_3OH); $\nu_{\text{max}}/\text{cm}^{-1}$ 3361 vs, 2922 m, 2854 w, 1653 w, 1455 m, 1419 w, 1130 m, 755 m, 701 m; δ_H (400 MHz; D_2O) 7.43-7.39 (2H, m), 7.34-7.32 (3H, m), 4.52-4.49 (2H, m), 4.17-4.14 (1H, m), 3.84-3.76 (2H, m), 3.65-3.58 (1H, m), 3.57-3.48 (2H, m), 3.45-3.38 (1H, m), 3.13-3.05 (1H, m), 2.82-2.69 (2H, m), 2.14-2.06 (2H, m); δ_c (100 MHz; D_2O) 140.5, 128.9, 128.5, 126.6, 70.7, 68.8, 68.6, 67.2, 62.5, 55.4, 53.7, 31.7, 26.0; HRMS(ESI) calcd for $\text{C}_{15}\text{H}_{24}\text{NO}_4^+$ [$\text{M}+\text{H}$] $^+$ 282.16998, found 282.16991.

***N*-(Phenylbutyl)-1,4-dideoxy-2,3:5,6-di-O-isopropylidene-1,4-imino-D-mannitol (19d)**

According to the general nucleophilic substitution procedures, product **19d** (225.8 mg, 77% yield) was obtained from **10** (190.0 mg, 0.78 mmol) and 4-phenyl-1-bromobutane (249.3 mg, 1.17 mmol) as a colorless syrup. $[\alpha]_D^{26} -43.5$ (c 0.96 in CH_2Cl_2); $\nu_{\text{max}}/\text{cm}^{-1}$ 2985 m, 2934 m, 2784 w, 1705 w, 1497 vw, 1454 w, 1370 m, 1267 w, 1207 vs, 1155 m, 1050 s, 860 m, 746 w, 699 m, 514 vw; δ_H (400 MHz; CDCl_3) 7.29-7.25 (2H, m), 7.19-7.15 (3H, m), 4.59-4.55 (2H, m), 4.48 (1H, td, $J = 7.2$ Hz, 1.3 Hz), 4.22 (1H, t, $J = 7.7$ Hz), 3.99 (1H, t, $J = 7.6$ Hz), 3.25 (1H, d, $J = 11.0$ Hz), 3.12 (1H, dt, $J = 12.2$ Hz, 8.4 Hz), 2.69-2.58 (3H, m), 2.08-2.02 (2H, m), 1.77-1.47 (4H, m), 1.45-1.44 (6H, m), 1.33 (3H, s), 1.28 (3H, s); δ_c (100 MHz; CDCl_3) 142.7, 128.4, 128.3, 125.6, 111.0, 107.1, 81.1, 77.7, 75.3, 68.1, 65.7, 59.1, 54.0, 35.8, 29.1, 27.4, 26.3, 25.9, 24.8, 23.9; HRMS(ESI) calcd for $\text{C}_{22}\text{H}_{34}\text{NO}_4^+$ [$\text{M}+\text{H}$] $^+$ 376.24824, found 376.24770.

***N*-(Phenylbutyl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (20d)**

According to the general deprotection procedures, product **20d** (44.1 mg, 99% yield) was obtained from **19d** (50.0 mg, 0.13 mmol) as a light yellow syrup. $[\alpha]_D^{26} -36.0$ (c 0.53 in CH_3OH); $\nu_{\text{max}}/\text{cm}^{-1}$ 3355 vs, 2940 m, 1647 w, 1496 w, 1455 m, 1341 w, 1130 m, 1056 w, 1044 w, 748 w, 700 m, 502 w; δ_H (400 MHz; D_2O) 7.40-7.36 (2H, m), 7.31-7.26 (3H, m), 4.48-4.40 (2H, m), 4.12-4.11 (1H, m), 3.78 (2H, d, $J = 5.2$ Hz), 3.46-3.39 (1H, m), 3.38-3.27 (2H, m), 3.22-3.20 (1H, m), 2.99-2.92 (1H, m), 2.73-7.63 (2H, m), 1.75-1.61 (4H, m); δ_c (100 MHz; D_2O) 142.3, 128.7, 128.6, 126.1, 71.1, 68.9, 68.3, 68.1, 62.7, 55.4, 54.7, 34.4, 27.7, 24.4; HRMS(ESI) calcd for $\text{C}_{16}\text{H}_{26}\text{NO}_4^+$ [$\text{M}+\text{H}$] $^+$ 296.18563, found 296.18530.

***N*-(Phenylpentyl)-1,4-dideoxy-2,3:5,6-di-O-isopropylidene-1,4-imino-D-mannitol (19e)**

According to the general nucleophilic substitution procedures, product **19e** (169.5 mg, 73% yield) was obtained from **10** (145.0 mg, 0.60 mmol) and 5-phenyl-1-bromopentane (204.4 mg, 0.90 mmol) as a colorless syrup. $[\alpha]_D^{26} -42.5$ (c 1.88 in CH_2Cl_2); $\nu_{\text{max}}/\text{cm}^{-1}$ 3026 vw, 2985 m, 2932 vs, 2858 m, 2783 w, 1496 w, 1454 w, 1379 s, 1367 m, 1207 vs, 1156 m, 1102 m, 1053 s, 860 m, 843 w, 699 m, 514 vw; δ_H (400 MHz; CDCl_3) 7.28-7.24 (2H, m), 7.17-7.14 (3H, m), 4.59-4.54 (2H, m), 4.46 (1H, td, $J = 7.3$ Hz, 1.6 Hz), 4.19 (1H, t, $J = 7.8$ Hz), 3.98 (1H, t, $J = 7.5$ Hz), 3.25 (1H, d, $J = 11.0$ Hz), 3.07 (1H, dt, $J = 11.7$ Hz, 8.1 Hz), 2.62-2.58 (3H, m), 2.06-1.99 (2H, m), 1.68-1.58 (2H, m), 1.54-1.27 (16H, m); δ_c (100 MHz; CDCl_3) 142.8, 128.4, 128.2, 125.6, 111.0, 107.1, 81.1, 77.7, 75.2, 68.0, 65.7, 59.1, 54.1, 36.0, 31.5, 27.6, 27.1, 26.3, 25.9, 24.8, 24.0; HRMS(ESI) calcd for $\text{C}_{23}\text{H}_{36}\text{NO}_4^+$ [$\text{M}+\text{H}$] $^+$ 390.26389, found 390.26359.

***N*-(Phenylpentyl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (20e)**

According to the general deprotection procedures, product **20e** (35.5 mg, 100% yield) was obtained from **19e** (40 mg, 0.10 mmol) as a light yellow syrup. $[\alpha]_D^{26}$ -34.9 (c 0.57 in CH₃OH); $\nu_{\max}/\text{cm}^{-1}$ 3347 vs, 3025 w, 2934 s, 2857 w, 1653 w, 1452 m, 1128 m, 1042 m, 748 w, 700 m, 576 w; δ_{H} (500 MHz; D₂O) 7.39-7.36 (2H, m), 7.31-7.26 (3H, m), 4.54-4.49 (2H, m), 4.19 (1H, q, $J = 5.0$ Hz), 3.83 (2H, dq, $J = 21.0$ Hz, 6.1 Hz), 3.62 (1H, t, $J = 4.1$ Hz), 3.55-3.47 (2H, m), 3.39-3.33 (1H, m), 3.08-3.02 (1H, m), 2.67 (2H, t, $J = 7.5$ Hz), 1.81-1.72 (2H, m), 1.71-1.65 (2H, m), 1.45-1.33 (2H, m); δ_{C} (125 MHz; D₂O) 142.9, 128.6, 125.9, 70.7, 68.7, 68.6, 67.2, 62.5, 55.3, 54.4, 34.7, 30.0, 25.2, 24.2; HRMS(ESI) calcd for C₁₇H₂₈NO₄⁺ [M+H]⁺ 310.20128, found 310.20107.

N-(Phenylhexyl)-1,4-dideoxy-2,3:5,6-di-O-isopropylidene-1,4-imino-D-mannitol (19f)

According to the general nucleophilic substitution procedures, product **19f** (147.8 mg, 81% yield) was obtained from **10** (110.0 mg, 0.45 mmol) and 6-phenyl-1-bromohexane (162.8 mg, 0.68 mmol) as a colorless syrup. $[\alpha]_D^{26}$ -36.5 (c 1.59 in CH₂Cl₂); $\nu_{\max}/\text{cm}^{-1}$ 2986 m, 2932 s, 2859 w, 2784 vw, 1705 vw, 1603 vw, 1497 vw, 1454 w, 1380 s, 1267 m, 1207 vs, 1157 m, 1102 m, 1053 s, 860 m, 747 w, 699 m, 517 w; δ_{H} (500 MHz; CDCl₃) 7.28-7.25 (2H, m), 7.18-7.15 (3H, m), 4.59-4.55 (2H, m), 4.46 (1H, td, $J = 7.3$ Hz, 1.3 Hz), 4.19 (1H, t, $J = 7.8$ Hz), 3.98 (1H, t, $J = 7.5$ Hz), 3.25 (1H, d, $J = 11.0$ Hz), 3.07 (1H, dt, $J = 11.9$ Hz, 8.2 Hz), 2.61-2.58 (3H, m), 2.06-1.99 (2H, m), 1.64-1.56 (2H, m), 1.48-1.27 (18H, m); δ_{C} (125 MHz; CDCl₃) 142.9, 128.4, 128.2, 125.5, 112.0, 107.1, 81.1, 77.7, 75.3, 68.0, 65.7, 59.1, 54.2, 35.94, 31.5, 29.3, 27.7, 27.3, 26.3, 25.9, 24.8, 24.0; HRMS(ESI) calcd for C₂₄H₃₈NO₄⁺ [M+H]⁺ 404.27954, found 404.27899.

N-(Phenylhexyl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (20f)

According to the general deprotection procedures, product **20f** (44.5 mg, 100% yield) was obtained from **19f** (50.0 mg, 0.12 mmol) as a light yellow syrup. $[\alpha]_D^{26}$ -32.9 (c 1.01 in CH₃OH); $\nu_{\max}/\text{cm}^{-1}$ 3332 vs, 3025 w, 2931 s, 2857 w, 1724 w, 1495 w, 1453 m, 1339 w, 1288 w, 1127 m, 1089 w, 1041 w, 748 w, 699 m, 492 w; δ_{H} (500 MHz; D₂O) 7.39-7.36 (2H, m), 7.31-7.25 (3H, m), 4.55-4.50 (2H, m), 4.20 (1H, q, $J = 5.0$ Hz), 3.87-3.80 (2H, m), 3.66 (1H, t, $J = 4.2$ Hz), 3.56-3.49 (2H, m), 3.39-3.34 (1H, m), 3.08-3.03 (1H, m), 2.67 (2H, t, $J = 7.5$ Hz), 1.76-1.71 (2H, m), 1.67-1.61 (2H, m), 1.45-1.33 (4H, m); δ_{C} (125 MHz; D₂O) 143.3, 128.6, 128.5, 125.8, 70.7, 68.7, 68.6, 67.2, 62.5, 55.3, 54.4, 34.8, 30.3, 27.6, 25.5, 24.2; HRMS(ESI) calcd for C₁₈H₃₀NO₄⁺ [M+H]⁺ 324.21693, found 324.21586.

N-(Phenylheptyl)-1,4-dideoxy-2,3:5,6-di-O-isopropylidene-1,4-imino-D-mannitol (19g)

According to the general nucleophilic substitution procedures, product **19g** (177.6 mg, 69% yield) was obtained from **10** (150.0 mg, 0.62 mmol) and 7-phenyl-1-bromoheptane (237.3 mg, 0.93 mmol) as a colorless syrup. $[\alpha]_D^{26}$ -34.3 (c 1.90 in CH₂Cl₂); $\nu_{\max}/\text{cm}^{-1}$ 2986 m, 2930 vs, 2856 m, 2787 w, 1710 m, 1458 w, 1379 m, 1267 w, 1208 s, 1157 m, 1103 m, 1058 m, 860 m, 699 m, 511 w; δ_{H} (400 MHz; CDCl₃) 7.29-7.25 (2H, m), 7.18-7.15 (3H, m), 4.59-4.55 (2H, m), 4.46 (1H, td, $J = 7.2$ Hz, 1.2 Hz), 4.20 (1H, t, $J = 7.7$ Hz), 3.98 (1H, t, $J = 7.5$ Hz), 3.25 (1H, d, $J = 11.0$ Hz), 3.06 (1H, dt, $J = 11.9$ Hz, 8.0 Hz), 2.61-2.57 (3H, m), 2.06-1.97 (2H, m), 1.66-1.57 (2H, m), 1.44-1.27 (20H, m); δ_{C} (100 MHz; CDCl₃) 142.9, 128.4, 128.2, 125.5, 111.0, 107.1, 81.1, 77.7, 75.3, 68.0, 65.7, 59.1, 54.2, 36.0, 31.5, 29.5, 29.3, 27.7, 27.3, 26.3, 25.9, 24.9, 24.0; HRMS(ESI) calcd for C₂₅H₄₀NO₄⁺ [M+H]⁺ 418.29519, found 418.29492.

N-(Phenylheptyl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (20g)

According to the general deprotection procedures, product **20g** (41.1 mg, 100% yield) was obtained from **19g** (46.0 mg, 0.11 mmol) as a light yellow syrup. $[\alpha]_D^{26}$ -30.4 (c 1.45 in CH₃OH); $\nu_{\max}/\text{cm}^{-1}$ 3420 vs, 2926 m, 2854 w, 1655 w, 1559 w, 1460 w, 1127 w, 1047 w, 747 w, 700 w; δ_{H} (500 MHz; D₂O) 7.37-7.34 (2H, m), 7.30-7.23 (3H, m), 4.53-4.50 (2H, m), 4.20 (1H, d, $J = 4.9$ Hz), 3.87-3.80 (2H, m), 3.62-3.61 (1H, m), 3.54-3.46 (2H, m), 3.38-3.32 (1H, m), 3.07-3.01 (1H, m), 2.63 (2H, t, $J = 7.4$ Hz), 1.76-1.66 (2H, m), 1.65-1.59 (2H, m), 1.35-1.33 (6H, m); δ_{C} (125 MHz; D₂O) 143.4, 128.6, 128.5, 125.8, 70.7, 68.7, 68.6, 67.3, 62.6, 55.3, 54.5, 34.9, 30.5, 27.9, 25.6, 24.3; HRMS(ESI) calcd for C₁₉H₃₂NO₄⁺ [M+H]⁺ 338.23258, found 338.23227.

N-(Phenylloctyl)-1,4-dideoxy-2,3:5,6-di-O-isopropylidene-1,4-imino-D-mannitol (19h)

According to the general nucleophilic substitution procedures, product **19h** (140.7 mg, 61% yield) was obtained from **10** (130.0 mg, 0.53 mmol) and 8-phenyl-1-bromooctane (201.9 mg, 0.75 mmol) as a colorless syrup. $[\alpha]_D^{26}$ -32.7 (c 1.35 in CH₂Cl₂); $\nu_{\max}/\text{cm}^{-1}$ 3026 vw, 2985 m, 2928 vs, 2783 vw, 1705 w, 1454 w, 1379 m, 1267 w, 1208 s, 1157 m, 1047 m, 860 w, 699 w, 514 vw; δ_{H} (500 MHz; CDCl₃) 7.28-7.25 (2H, m), 7.18-7.15 (3H, m), 4.59-4.55 (2H, m), 4.46 (1H, td, $J = 7.3$ Hz, 1.6 Hz), 4.19 (1H, t, $J = 7.8$ Hz), 3.98 (1H, t, $J = 7.5$ Hz), 3.25 (1H, d, $J = 11.0$ Hz), 3.05 (1H, dt, $J = 12.3$ Hz, 8.4 Hz), 2.61-2.58 (3H, m), 2.06-1.98 (2H, m), 1.62-1.56 (2H, m), 1.45-1.27 (22H, m); δ_{C} (125 MHz; CDCl₃) 142.9, 128.4, 128.2, 125.5, 111.0, 107.1, 81.1, 77.7, 75.3, 68.0, 65.7, 59.1, 54.2, 36.0, 31.5, 29.5, 29.5, 29.3, 27.7, 27.4, 26.3, 25.9, 24.9, 24.0; HRMS(ESI) calcd for C₂₆H₄₂NO₄⁺ [M+H]⁺ 432.31084, found 432.30954.

N-(Phenylloctyl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (20h)

According to the general deprotection procedures, product **20h** (54.7 mg, 100% yield) was obtained from **19h** (61.0 mg, 0.14 mmol) as a light yellow syrup. $[\alpha]_D^{26}$ -34.5 (c 0.44 in CH₃OH); $\nu_{\max}/\text{cm}^{-1}$ 3342 vs, 3025 w, 2931 s, 2856 m, 1653 w, 1453 m, 1129 m, 1044 w, 749 w, 699 m, 576 w; δ_{H} (500 MHz; D₂O) 7.22-7.18 (2H, m), 7.13-7.08 (3H, m), 4.53-4.51 (2H, m), 4.18 (1H, q, $J = 4.8$ Hz), 3.89-3.81 (2H, m), 3.59 (1H, s), 3.52-3.49 (1H, m), 3.41-3.37 (1H, m), 3.32-3.37 (1H, m), 2.98-2.93 (1H, m), 2.51 (2H, t, $J = 7.5$ Hz), 1.64-1.63 (2H, m), 1.52 (2H, s), 1.23 (8H, m); δ_{C} (125 MHz; D₂O) 143.0, 128.4, 128.3, 125.6, 70.6,

69.0, 68.6, 67.3, 62.6, 55.3, 54.4, 35.4, 31.1, 28.8, 28.6, 28.5, 26.0, 24.5; HRMS(ESI) calcd for C₂₀H₃₄NO₄⁺ [M+H]⁺ 352.24824, found 352.24780.

***N*-3-((*tert*-Butyldimethylsilyloxy)propyl)-1,4-dideoxy-2,3:5,6-di-*O*-isopropylidene-1,4-imino-D-mannitol (28b)**

According to the general nucleophilic substitution procedures, product **28b** (204.5 mg, 57% yield) was obtained from **10** (200.0 mg, 0.82 mmol) and 3-((*tert*-butyldimethylsilyloxy)-1-bromopropane (311.5 mg, 1.23 mmol) as a light yellow syrup. [α]_D²⁶ -29.1 (c 0.93 in CH₂Cl₂); ν_{max} /cm⁻¹ 3392 m, 2986 s, 2935 s, 2805 w, 1688 m, 1456 w, 1373 s, 1381 m, 1211 vs, 1156 s, 1059 vs, 990 w, 858 m, 842 m, 777 w, 514 w; δ_{H} (400 MHz; CDCl₃) 4.59-4.55 (2H, m), 4.43 (1H, td, *J* = 7.3 Hz, 1.9 Hz), 4.13 (1H, t, *J* = 7.8 Hz), 3.98 (1H, t, *J* = 7.4 Hz), 3.73-3.60 (2H, m), 3.24 (1H, d, *J* = 11.0 Hz), 3.13 (1H, dt, *J* = 12.3 Hz, 8.2 Hz), 2.59-2.58 (1H, m), 2.11-2.05 (2H, m), 1.71-1.65 (2H, m), 1.42 (6H, s), 1.32 (3H, s), 1.27 (3H, s), 0.88 (9H, s), 0.04 (6H, s); δ_{C} (100 MHz; CDCl₃) 110.0, 107.1, 81.2, 77.6, 75.3, 68.1, 65.9, 61.6, 59.3, 50.8, 31.0, 26.4, 26.0, 25.87, 24.9, 24.1, 18.4, -5.3; HRMS(ESI) calcd for C₂₁H₄₂NO₅Si⁺ [M+H]⁺ 416.28268, found 416.28230.

***N*-3-Hydroxypropyl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (29b)**

According to the general deprotection procedures, product **29b** (37.1 mg, 99% yield) was obtained from **28b** (60.0 mg, 0.14 mmol) as a light yellow solid. M. p. 78-80 °C; [α]_D²⁶ -26.1 (c 1.00 in CH₃OH); ν_{max} /cm⁻¹ 3397 vs, 2964 w, 1647 w, 1405 w, 1126 w, 1050 w, 561 w; δ_{H} (400 MHz; D₂O) 4.63-4.55 (2H, m), 4.26 (1H, q, *J* = 5.0 Hz), 3.93-3.84 (2H, m), 3.82-3.71 (3H, m), 3.63-3.54 (3H, m), 3.30-3.23 (1H, m), 2.12-1.96 (2H, m); δ_{C} (100 MHz; D₂O) 70.8, 69.1, 68.65, 67.3, 62.6, 56.2, 55.6, 52.6, 26.9; HRMS(ESI) calcd for C₉H₂₀NO₅⁺ [M+H]⁺ 222.13360, found 222.13360.

***N*-4-((*tert*-Butyldimethylsilyloxy)butyl)-1,4-dideoxy-2,3:5,6-di-*O*-isopropylidene-1,4-imino-D-mannitol (28c)**

According to the general nucleophilic substitution procedures, product **28c** (108.8 mg, 77% yield) was obtained from **10** (80.0 mg, 0.33 mmol) and 4-((*tert*-butyldimethylsilyloxy)-1-bromobutane (132.3 mg, 0.50 mmol) as a light yellow syrup. [α]_D²⁸ -32.3 (c 1.50 in CH₂Cl₂); ν_{max} /cm⁻¹ 2986 w, 2934 s, 2858 vw, 1705 w, 1463 w, 1380 m, 1256 m, 1208 s, 1157 m, 1100 vs, 1058 m, 837 s, 776 m, 515 vw; δ_{H} (400 MHz; CDCl₃) 4.59-4.54 (2H, m), 4.45 (1H, td, *J* = 7.2 Hz, 1.2 Hz), 4.18 (1H, t, *J* = 7.8 Hz), 3.98 (1H, t, *J* = 7.6 Hz), 3.61 (2H, t, *J* = 5.2 Hz), 3.25 (1H, d, *J* = 11.0 Hz), 3.06 (1H, dt, *J* = 12.1 Hz, 7.4 Hz), 2.60 (1H, d, *J* = 2.4 Hz), 2.08-2.02 (2H, m), 1.61-1.46 (4H, m), 1.43 (6H, s), 1.32 (3H, s), 1.27 (3H, s), 0.88 (9H, s), 0.04 (6H, s); δ_{C} (100 MHz; CDCl₃) 110.0, 107.1, 81.1, 77.7, 75.3, 68.0, 65.7, 63.2, 59.1, 54.1, 30.7, 26.3, 26.0, 25.9, 24.9, 24.1, 24.0, 18.4, -5.2, -5.3; HRMS(ESI) calcd for C₂₂H₄₄NO₅Si⁺ [M+H]⁺ 430.29833, found 430.29768.

***N*-4-Hydroxybutyl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (29c)**

According to the general deprotection procedures, product **29c** (11.4 mg, 100% yield) was obtained from **35c** (18.0 mg, 0.04 mmol) as a light yellow syrup. [α]_D²⁶ -29.3 (c 0.28 in CH₃OH); ν_{max} /cm⁻¹ 3359 w, 2921 m, 2854 w, 2373 w, 2326 w, 1647 w, 1458 w, 1130 w, 1022 w, 417 w; δ_{H} (400 MHz; D₂O) 4.57-4.50 (2H, m), 4.22 (1H, q, *J* = 5.1 Hz), 3.89-3.79 (2H, m), 3.68-3.60 (3H, m), 3.56 (2H, d, *J* = 7.4 Hz), 3.47-3.40 (1H, m), 3.17-3.10 (1H, m), 1.88-1.74 (2H, m), 1.66-1.59 (2H, m); δ_{C} (100 MHz; D₂O) 70.7, 68.8, 68.6, 67.2, 62.5, 60.8, 55.3, 54.2, 28.4, 21.4; HRMS(ESI) calcd for C₁₀H₂₂NO₅⁺ [M+H]⁺ 236.14925, found 236.14906.

***N*-5-((*tert*-Butyldimethylsilyloxy)pentyl)-1,4-dideoxy-2,3:5,6-di-*O*-isopropylidene-1,4-imino-D-mannitol (28d)**

According to the general nucleophilic substitution procedures, product **35d** (240.6 mg, 66% yield) was obtained from **10** (200.0 mg, 0.82 mmol) and 5-((*tert*-butyldimethylsilyloxy)-1-bromopentane (346.0 mg, 1.23 mmol) as a light yellow syrup. [α]_D²⁰ -16.6 (c 0.56 in CH₂Cl₂); ν_{max} /cm⁻¹ 3447 w, 2931 vs, 2858 m, 2801 w, 1729 m, 1684 m, 1458 w, 1380 m, 1257 m, 1208 s, 1156 m, 1098 s, 1073 m, 990 w, 932 w, 860 m, 836 s, 776 m, 512 w; δ_{H} (400 MHz; CDCl₃) 4.59-4.54 (2H, m), 4.45 (1H, td, *J* = 7.3 Hz, 1.4 Hz), 4.18 (1H, t, *J* = 7.7 Hz), 3.98 (1H, t, *J* = 7.5 Hz), 3.59 (2H, t, *J* = 6.5 Hz), 3.25 (1H, d, *J* = 11.0 Hz), 3.06 (1H, dt, *J* = 11.8 Hz, 8.1 Hz), 2.59 (1H, d, *J* = 2.7 Hz), 2.07-1.99 (2H, m), 1.56-1.27 (18H, m), 0.88 (9H, s), 0.03 (6H, s); δ_{C} (100 MHz; CDCl₃) 111.1, 107.1, 81.1, 77.7, 75.3, 68.0, 65.8, 63.2, 54.3, 32.9, 27.6, 26.3, 26.0, 25.9, 24.9, 24.0, 23.8, 18.3, -5.3; HRMS(ESI) calcd for C₂₃H₄₆NO₅Si⁺ [M+H]⁺ 444.3140, found 444.3137.

***N*-5-Hydroxypentyl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (29d)**

According to the general deprotection procedures, product **29d** (38.6 mg, 100% yield) was obtained from **28d** (60.0 mg, 0.14 mmol) as a light yellow syrup. [α]_D²⁰ -41.4 (c 0.76 in CH₂Cl₂); ν_{max} /cm⁻¹ 3365 vs, 2942 m, 1647 w, 1418 m, 1130 m, 1046 m, 560 w; δ_{H} (500 MHz; D₂O) 4.56-4.52 (1H, m), 4.50 (1H, t, *J* = 4.0 Hz), 4.21 (1H, q, *J* = 5.0 Hz), 3.87-3.79 (2H, m), 3.66 (1H, t, *J* = 4.0 Hz), 3.60 (2H, t, *J* = 6.5 Hz), 3.55 (2H, d, *J* = 7.5 Hz), 3.44-3.38 (1H, m), 3.13-3.08 (1H, m), 1.84-1.71 (2H, m), 1.61-1.56 (2H, m), 1.48-1.35 (2H, m); δ_{C} (125 MHz; D₂O) 70.7, 68.8, 68.6, 67.2, 62.5, 61.2, 55.3, 54.3, 30.6, 24.1, 22.2; HRMS(ESI) calcd for C₁₁H₂₄NO₅⁺ [M+H]⁺ 250.1649, found 250.1649.

***N*-6-((*tert*-Butyldimethylsilyloxy)hexyl)-1,4-dideoxy-2,3:5,6-di-*O*-isopropylidene-1,4-imino-D-mannitol (28e)**

According to the general nucleophilic substitution procedures, product **28e** (259.5 mg, 69% yield) was obtained from **10** (200.0 mg, 0.82 mmol) and 6-((*tert*-butyldimethylsilyloxy)-1-bromohexane (363.3 mg, 1.23 mmol) as a light yellow syrup. [α]_D²⁵ -40.8 (c 1.82 in CH₂Cl₂); ν_{max} /cm⁻¹ 2986 m, 2931 vs, 2858 s, 2783 w, 1463 w, 1380 m, 1256 m, 1208 s, 1156 m, 1101 s, 1048 m, 860 m, 837 s, 775 m, 514 w; δ_{H} (500 MHz; CDCl₃) 4.59-4.54 (2H, m), 4.45 (1H, td, *J* = 7.3 Hz, 1.4 Hz), 4.18 (1H, t, *J* = 7.8 Hz), 3.98 (1H, t, *J* = 7.5 Hz), 3.58 (2H, t, *J* = 6.7 Hz), 3.25 (1H, d, *J* = 11.0 Hz), 3.05 (1H, dt, *J* = 12.0 Hz, 8.1 Hz), 2.59 (1H, d,

$J = 2.8$ Hz), 2.06-1.99 (2H, m), 1.53-1.48 (2H, m), 1.43-1.27 (18H, m), 0.89 (9H, s), 0.04 (6H, s); δ_c (125 MHz; CDCl_3) 111.0, 107.1, 81.1, 77.7, 75.3, 68.0, 65.7, 63.3, 59.1, 54.2, 32.9, 27.7, 26.3, 26.3, 26.0, 25.9, 25.8, 24.8, 24.0, 18.4, -5.3; HRMS(ESI) calcd for $\text{C}_{24}\text{H}_{48}\text{NO}_5\text{Si}^+$ [M+H]⁺ 458.32963, found 458.32918.

***N*-(6-Hydroxyhexyl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (29e)**

According to the general deprotection procedures, product **29e** (39.2 mg, quantitative yield) was obtained from **28e** (60 mg, 0.13 mmol) as a light yellow syrup. $[\alpha]_D^{26}$ -40.2 (c 0.60 in CH_2Cl_2); $\nu_{\text{max}}/\text{cm}^{-1}$ 3365 vs, 2937 m, 2865 w, 1647 w, 1458 m, 1130 m, 1046 m, 550 w; δ_H (500 MHz; D_2O) 4.57-4.50 (2H, m), 4.22 (1H, q, $J = 5.0$ Hz), 3.88-3.80 (2H, m), 3.67 (1H, t, $J = 4.0$ Hz), 3.60 (2H, t, $J = 6.5$ Hz), 3.56 (2H, d, $J = 7.5$ Hz), 3.44-3.38 (1H, m), 3.14-3.08 (1H, m), 1.84-1.71 (2H, m), 1.57-1.55 (2H, m), 1.41-1.40 (4H, m); δ_c (125 MHz; D_2O) 70.7, 68.7, 68.6, 67.2, 62.5, 61.5, 55.3, 54.4, 30.9, 24.4, 24.5, 24.2; HRMS(ESI) calcd for $\text{C}_{12}\text{H}_{26}\text{NO}_5^+$ [M+H]⁺ 264.180555, found 264.18035.

***N*-(7-((*tert*-Butyldimethylsilyloxy)heptyl)-1,4-dideoxy-2,3:5,6-di-*O*-isopropylidene-1,4-imino-D-mannitol (28f)**

According to the general nucleophilic substitution procedures, product **28f** (143.6 mg, 57% yield) was obtained from **10** (130.0 mg, 0.53 mmol) and 7-((*tert*-butyldimethylsilyloxy)-1-bromoheptane (245.9 mg, 0.80 mmol) as a pale yellow syrup. $[\alpha]_D^{26}$ -33.1 (c 1.57 in CH_2Cl_2); $\nu_{\text{max}}/\text{cm}^{-1}$ 3514 w, 2986 m, 2931 vs, 2857 m, 2784 w, 1707 w, 1463 w, 1380 m, 1256 m, 1208 s, 1156 m, 1099 s, 1050 m, 990 w, 932 w, 860 w, 837 m, 775 m, 514 w; δ_H (400 MHz; CDCl_3) 4.58-4.53 (2H, m), 4.45 (1H, td, $J = 7.2$ Hz, 1.5 Hz), 4.18 (1H, t, $J = 7.8$ Hz), 3.98 (1H, t, $J = 7.5$ Hz), 3.58 (2H, t, $J = 6.6$ Hz), 3.24 (1H, d, $J = 11.0$ Hz), 3.04 (1H, td, $J = 12.1$ Hz, 8.1 Hz), 2.58 (1H, d, $J = 2.6$ Hz), 2.05-1.98 (2H, m), 1.50-1.42 (10H, m), 1.34-1.23 (12H, m), 0.88 (9H, s), 0.04 (6H, s); δ_c (100 MHz; CDCl_3) 111.0, 107.1, 81.1, 77.7, 75.3, 68.0, 65.7, 63.3, 59.1, 54.2, 32.9, 29.4, 27.7, 27.4, 26.3, 26.0, 25.9, 25.8, 24.9, 24.0, 18.4, -5.3; HRMS(ESI) calcd for $\text{C}_{25}\text{H}_{50}\text{NO}_5\text{Si}^+$ [M+H]⁺ 472.34528, found 472.34512.

***N*-(7-Hydroxyheptyl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (29f)**

According to the general deprotection procedures, product **29f** (46.5 mg, 100% yield) was obtained from **28f** (70.0 mg, 0.15 mmol) as a pale yellow syrup. $[\alpha]_D^{28}$ -26.7 (c 1.64 in CH_2Cl_2); $\nu_{\text{max}}/\text{cm}^{-1}$ 3419 vs, 2935 m, 2931 vs, 2862 w, 1647 m, 1458 w, 1133 w, 1026 w, 597 w; δ_H (400 MHz; CDCl_3) 4.57-4.50 (2H, m), 4.21 (1H, q, $J = 5.0$ Hz), 3.87-3.79 (2H, m), 3.66 (1H, t, $J = 3.8$ Hz), 3.60-3.54 (4H, m), 3.43-3.35 (1H, m), 3.13-3.06 (1H, m), 1.80-1.68 (2H, m), 1.55-1.52 (2H, m), 1.42-1.30 (6H, m); δ_c (100 MHz; CDCl_3) 70.7, 68.8, 68.6, 67.3, 62.6, 61.7, 55.4, 54.5, 31.1, 27.9, 25.6, 24.7, 24.3; HRMS(ESI) calcd for $\text{C}_{13}\text{H}_{28}\text{NO}_5^+$ [M+H]⁺ 278.19620, found 278.19604.

***N*-(8-((*tert*-Butyldimethylsilyloxy)octyl)-1,4-dideoxy-2,3:5,6-di-*O*-isopropylidene-1,4-imino-D-mannitol (28g)**

According to the general nucleophilic substitution procedures, product **28g** (133.7 mg, 67% yield) was obtained from **10** (100.0 mg, 0.41 mmol) and 8-((*tert*-butyldimethylsilyloxy)-1-bromooctane (198.9 mg, 0.62 mmol) as a light yellow syrup. $[\alpha]_D^{28}$ -33.1 (c 1.57 in CH_2Cl_2); $\nu_{\text{max}}/\text{cm}^{-1}$ 2986 m, 2933 s, 2859 w, 2784 w, 1705 w, 1455 w, 1380 s, 1267 m, 1208 vs, 1157 m, 1053 s, 860 m, 843 w, 699 m, 517 w; δ_H (500 MHz; CDCl_3) 4.59-4.54 (2H, m), 4.45 (1H, td, $J = 7.3$ Hz, 1.5 Hz), 4.18 (1H, t, $J = 7.8$ Hz), 3.98 (1H, t, $J = 7.5$ Hz), 3.58 (2H, t, $J = 6.7$ Hz), 3.24 (1H, d, $J = 11.1$ Hz), 3.05 (1H, dt, $J = 12.2$ Hz, 8.3 Hz), 2.59 (1H, d, $J = 2.8$ Hz), 2.06-1.98 (2H, m), 1.51-1.43 (10H, m), 1.32-1.27 (14H, m), 0.88 (9H, s), 0.04 (6H, s); δ_c (125 MHz; CDCl_3) 111.0, 107.1, 81.1, 77.7, 75.3, 68.0, 65.8, 63.3, 59.1, 54.3, 32.9, 29.6, 29.5, 27.7, 27.4, 26.3, 26.0, 25.9, 25.8, 24.9, 24.0, 18.4, -5.2; HRMS(ESI) calcd for $\text{C}_{26}\text{H}_{52}\text{NO}_5\text{Si}^+$ [M+H]⁺ 486.36093, found 486.36082.

***N*-(8-Hydroxyoctyl)-1,4-dideoxy-1,4-imino-D-mannitol hydrochloride (29g)**

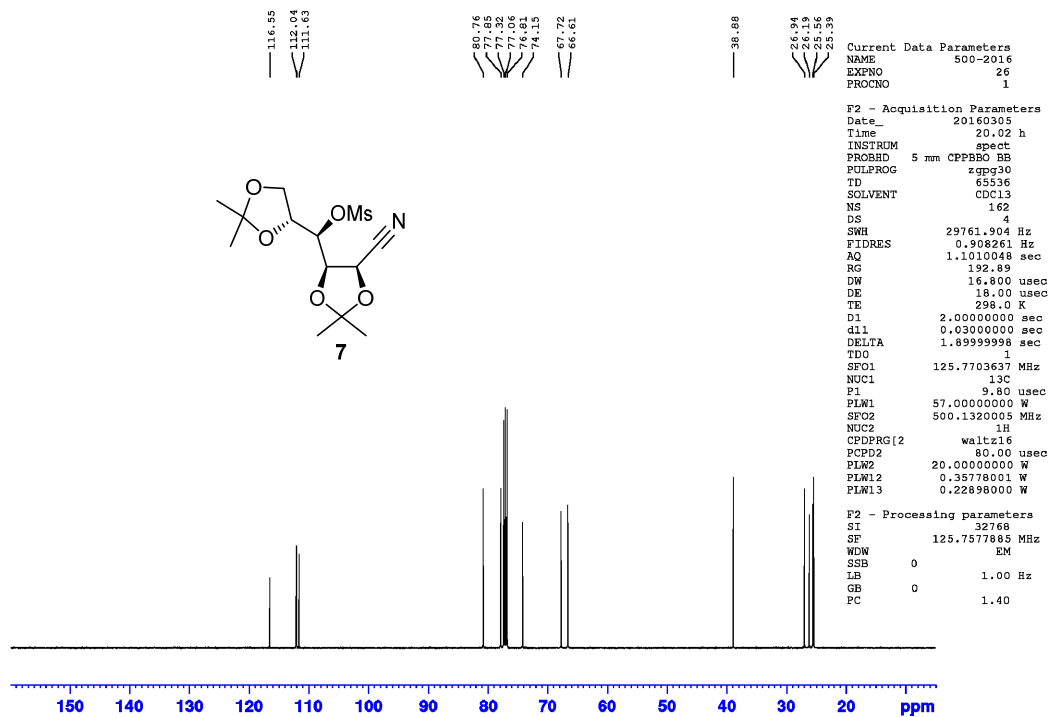
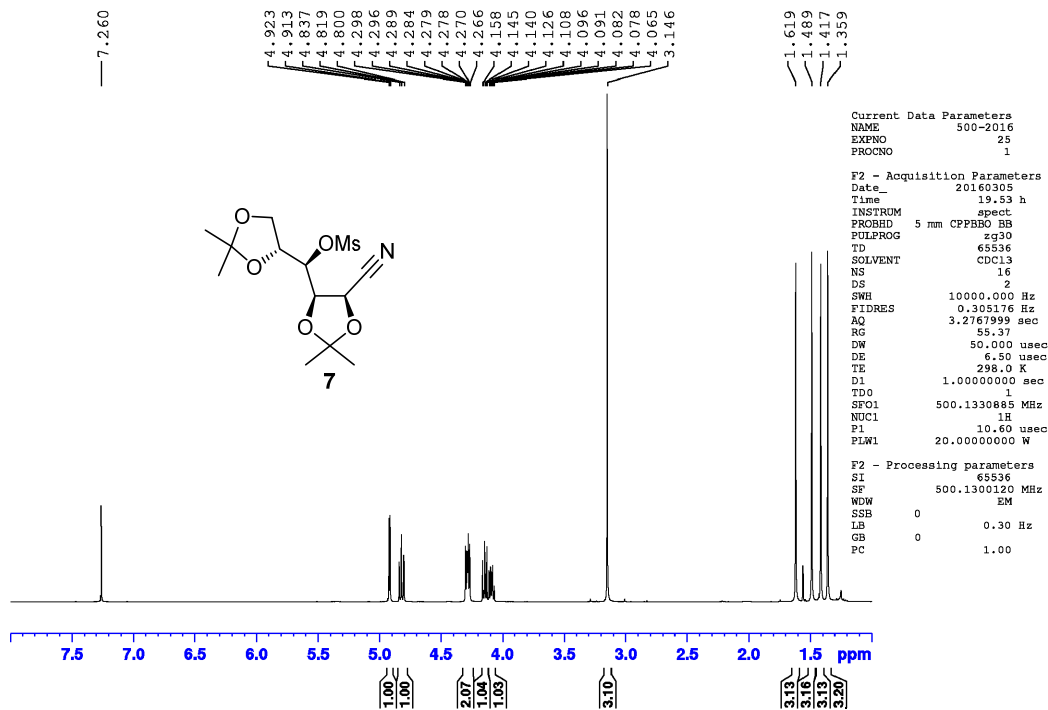
According to the general deprotection procedures, product **29g** (33.7 mg, 100% yield) was obtained from **28g** (50.0 mg, 0.10 mmol) as a pale yellow syrup. $[\alpha]_D^{26}$ -35.2 (c 0.77 in CH_2Cl_2); $\nu_{\text{max}}/\text{cm}^{-1}$ 3336 vs, 2933 s, 2856 m, 1716 w, 1404 s, 1277 w, 1128 m, 1046 m, 714 w, 517 w; δ_H (500 MHz; CDCl_3) 4.57-4.50 (2H, m), 4.22 (1H, q, $J = 5.0$ Hz), 3.88-3.80 (2H, m), 3.66 (1H, t, $J = 4.2$ Hz), 3.59 (2H, t, $J = 6.7$ Hz), 3.55 (2H, d, $J = 7.6$ Hz), 3.43-3.35 (1H, m), 3.13-3.07 (1H, m), 1.79-1.69 (2H, m), 1.55-1.53 (2H, m), 1.36-1.33 (8H, m); δ_c (125 MHz; CDCl_3) 70.7, 68.7, 68.6, 67.3, 62.5, 61.8, 55.3, 54.5, 31.1, 28.1, 28.0, 25.6, 24.8, 24.5; HRMS(ESI) calcd for $\text{C}_{14}\text{H}_{30}\text{NO}_5^+$ [M+H]⁺ 292.21185, found 292.21140.

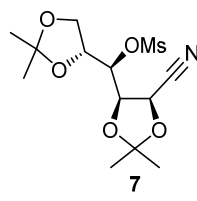
References

1 G. W. J. Fleet, B. Winchester, N. M. Carpenter, US4996329A, 1991.

2. NMR spectra and Infrared spectra for compounds

Compound 7:





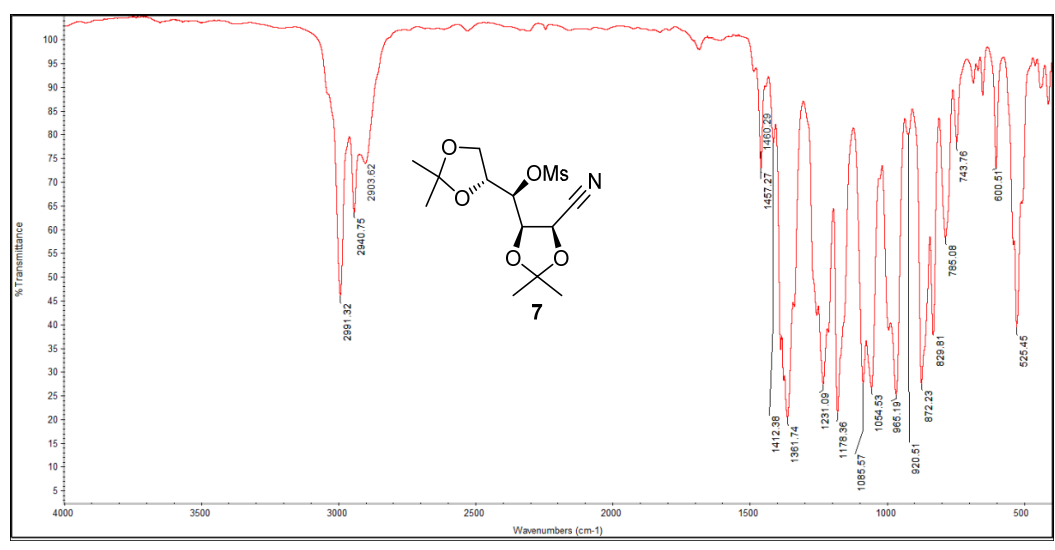
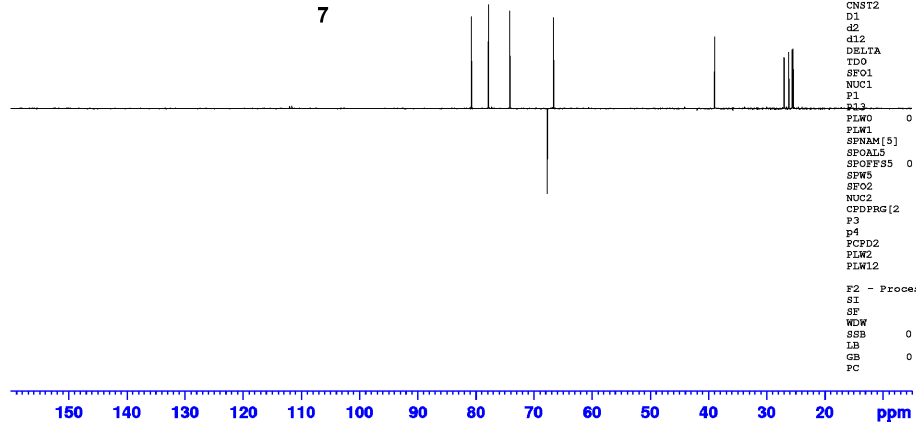
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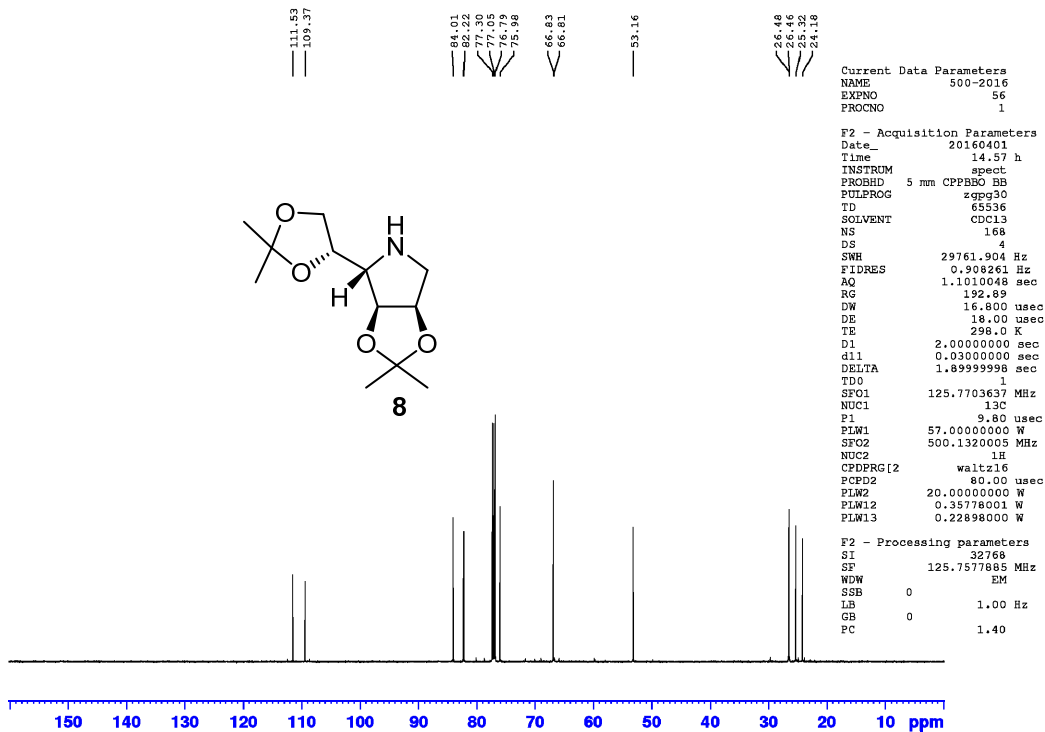
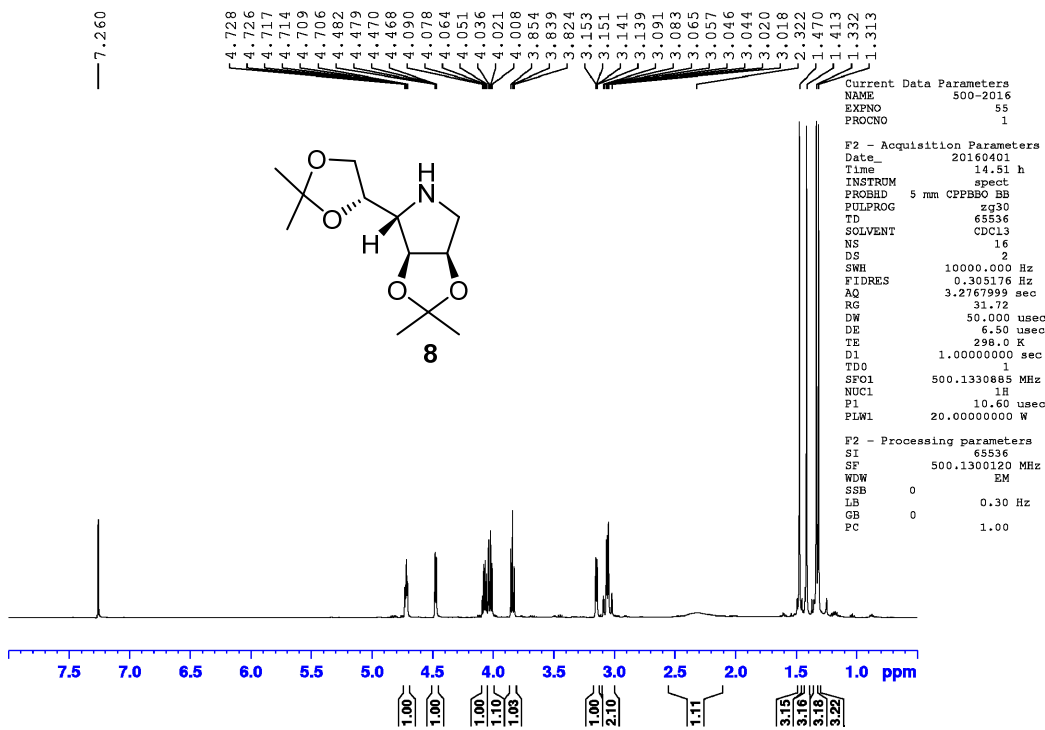
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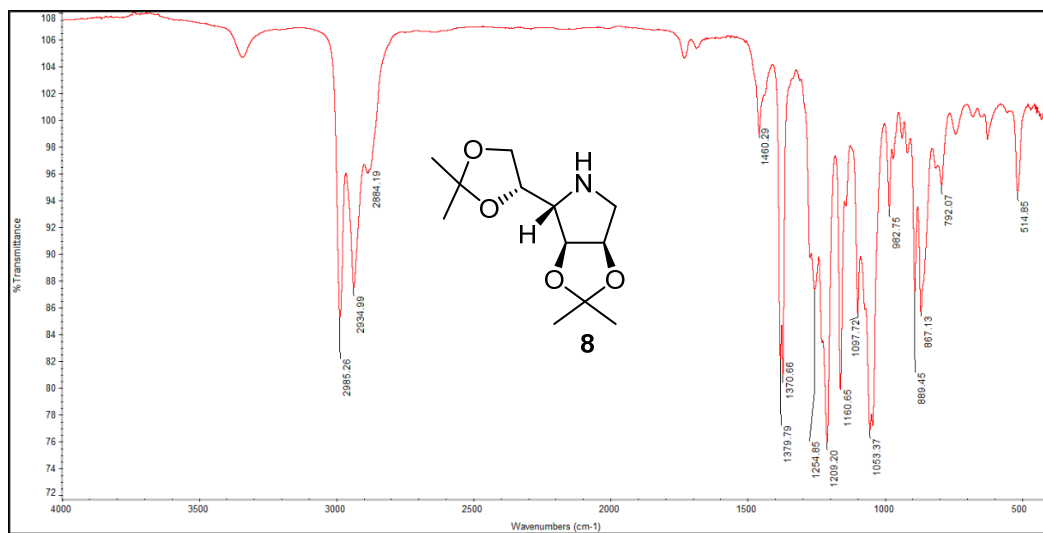
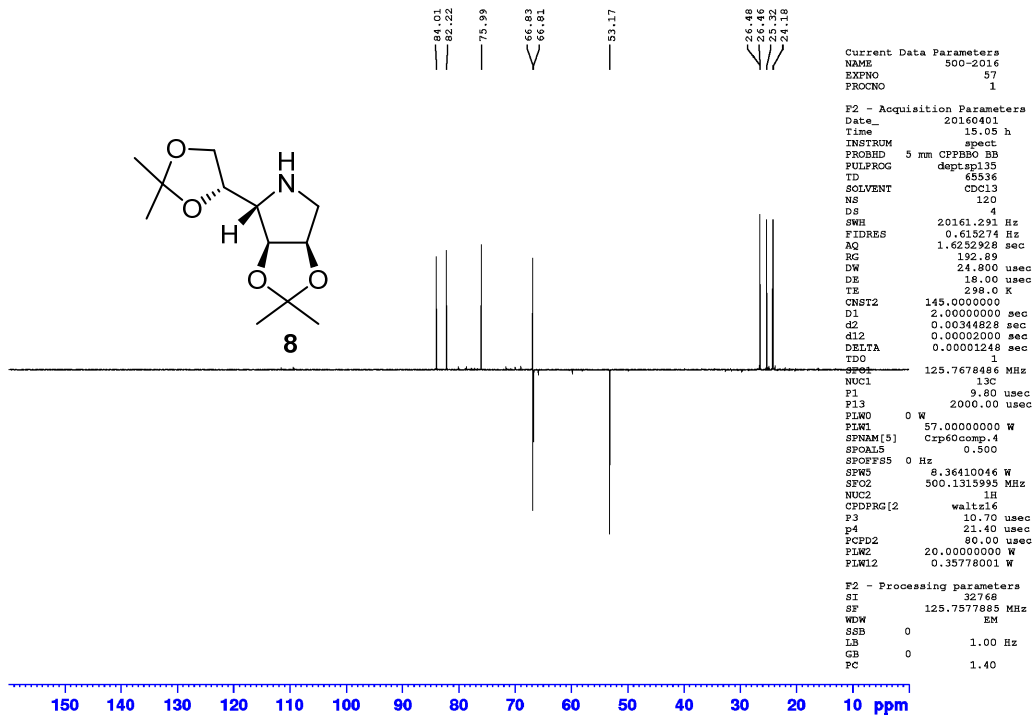
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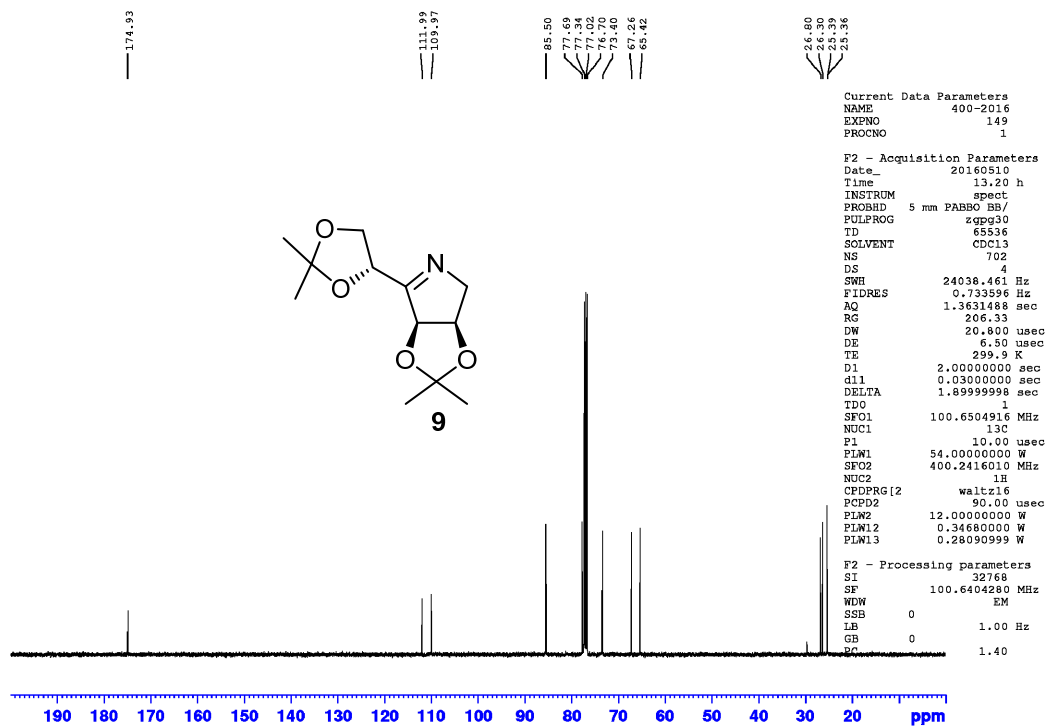
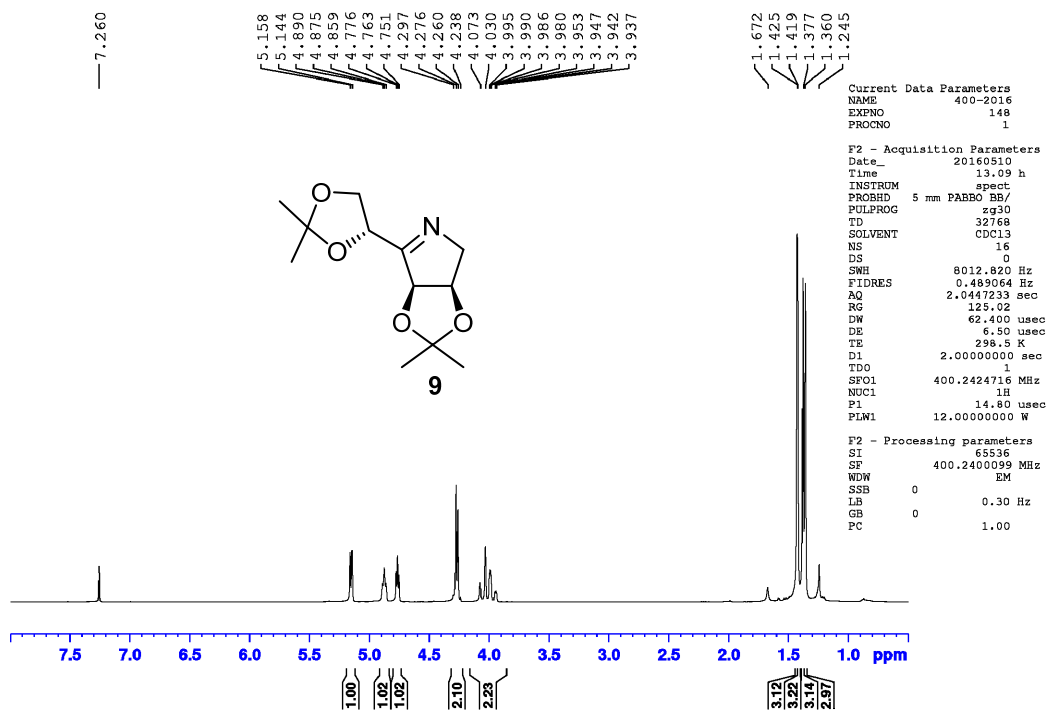


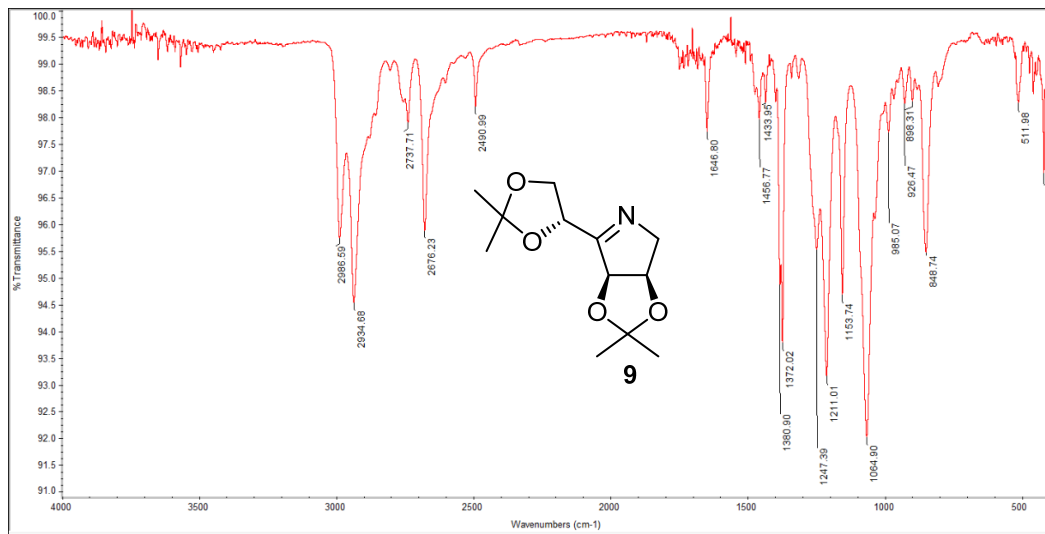
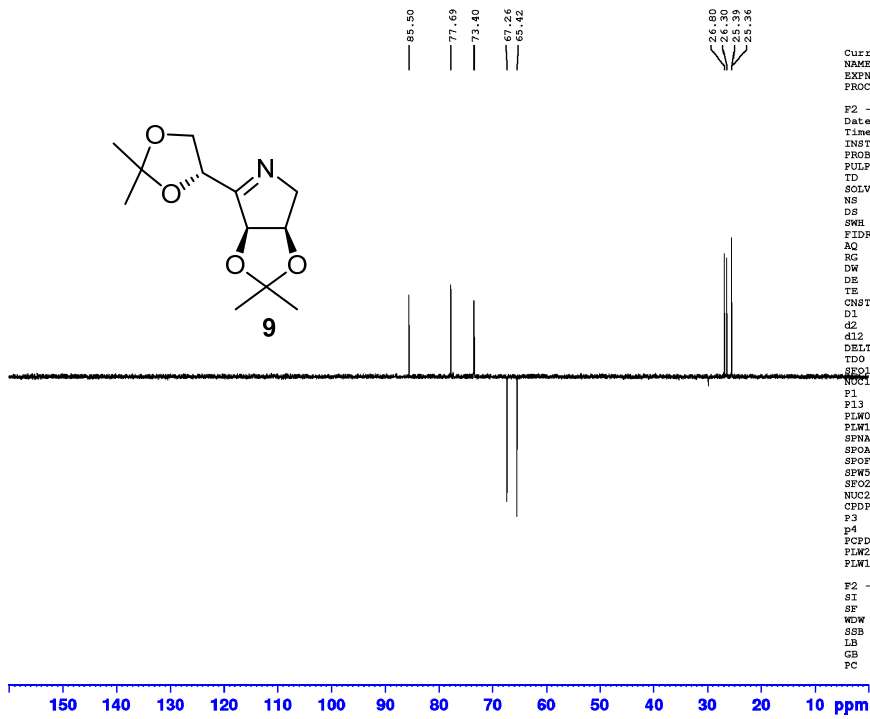
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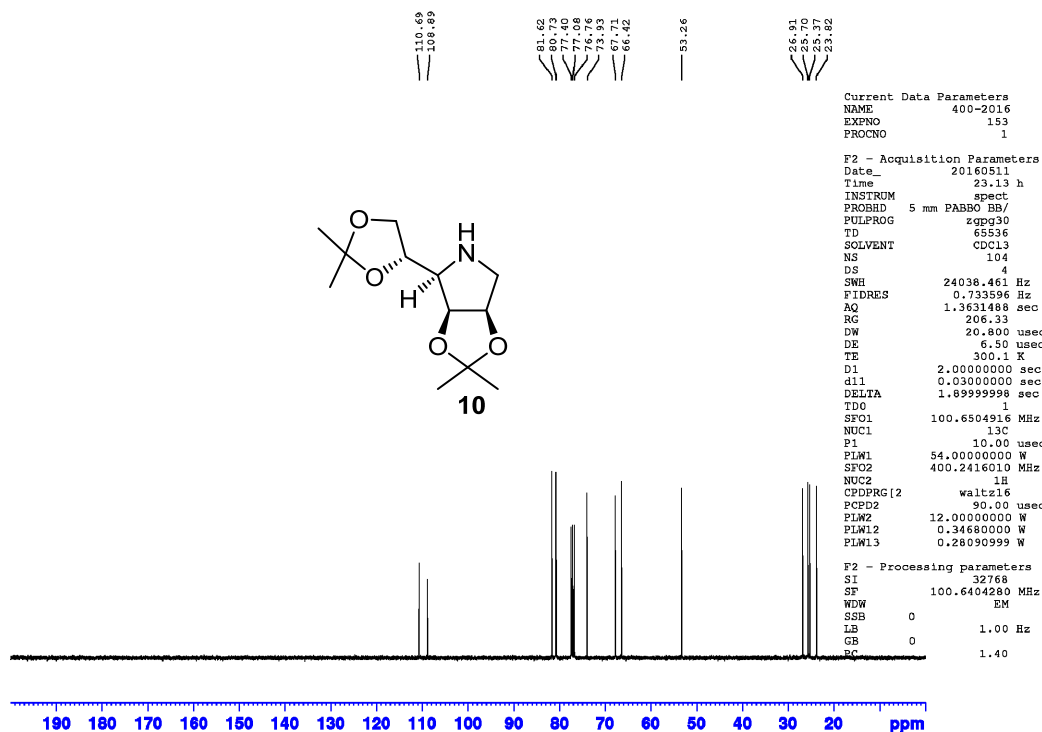
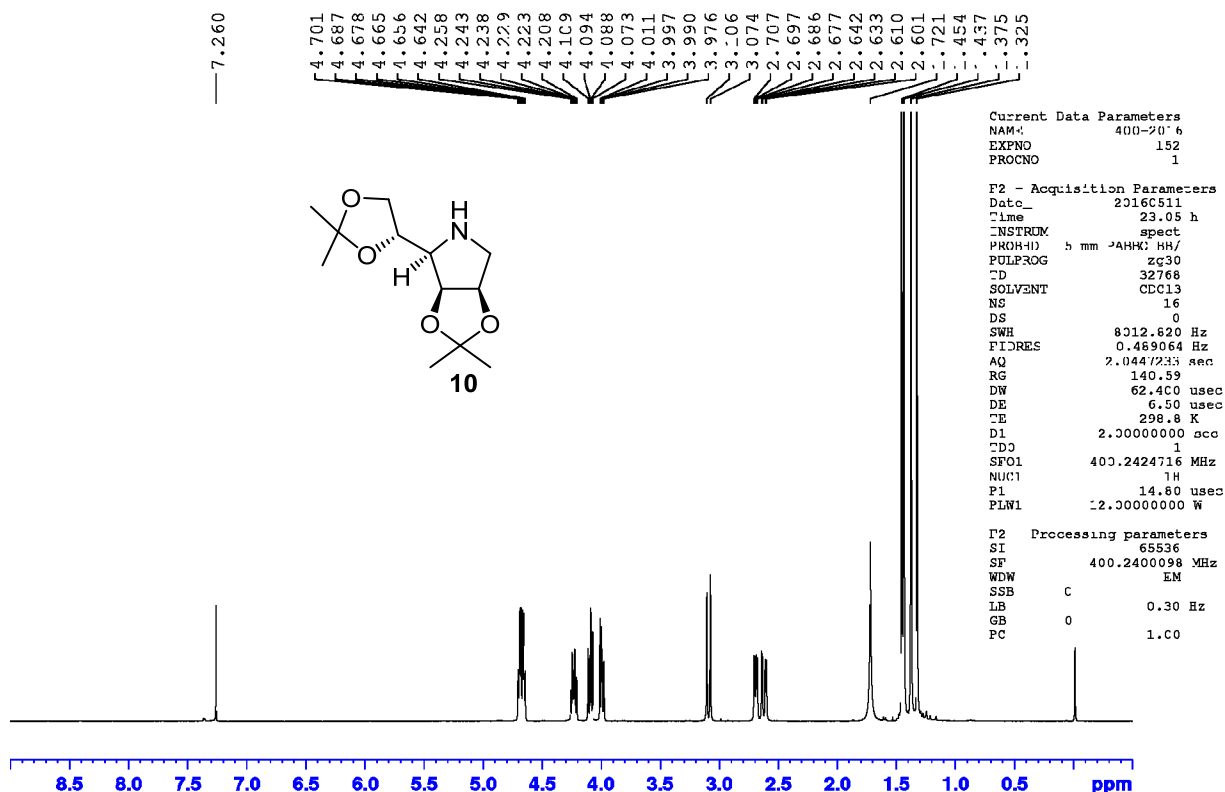


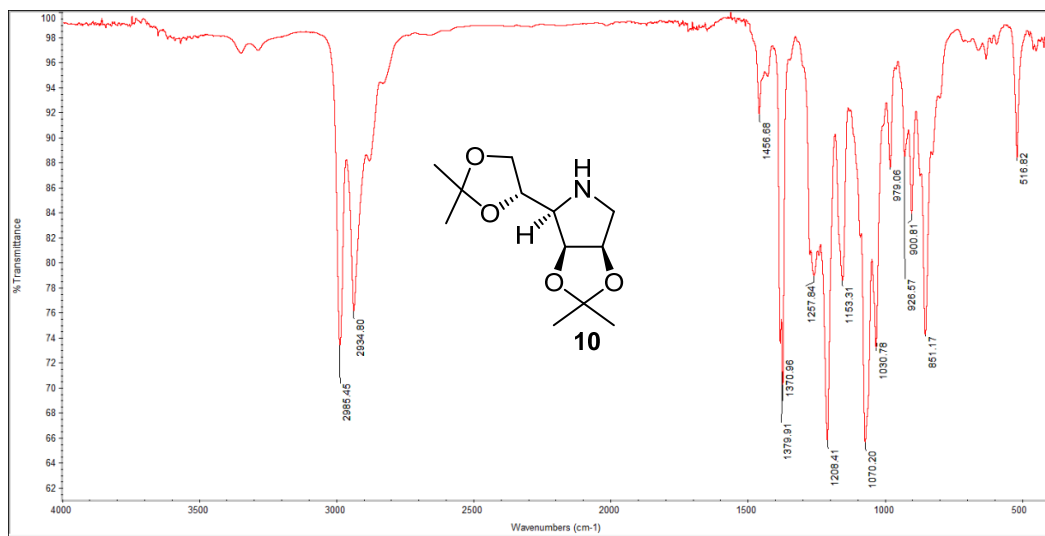
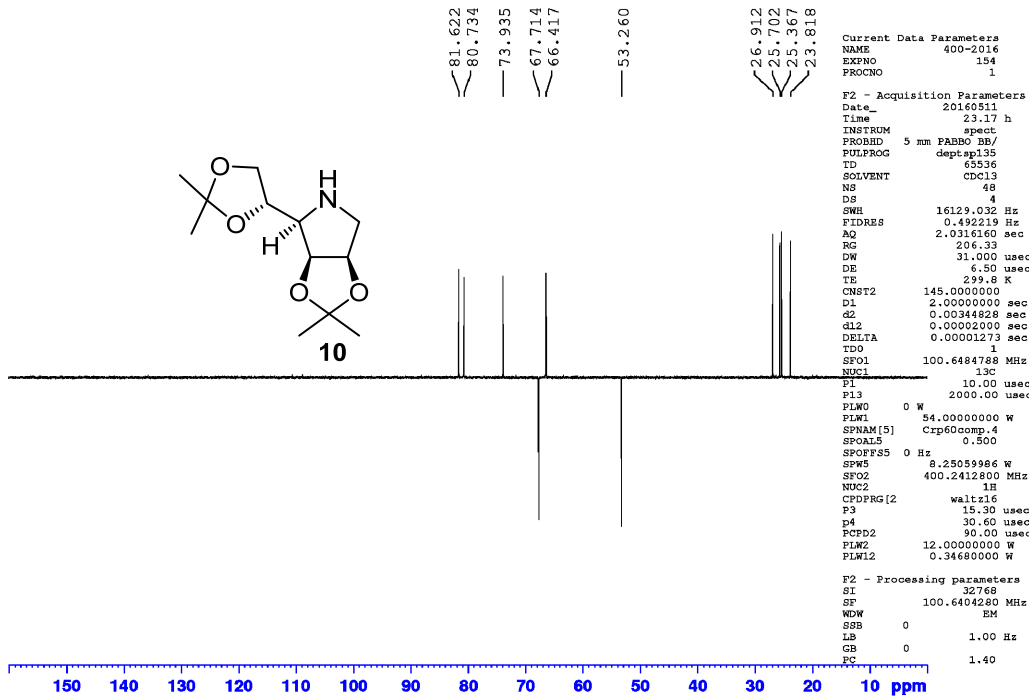
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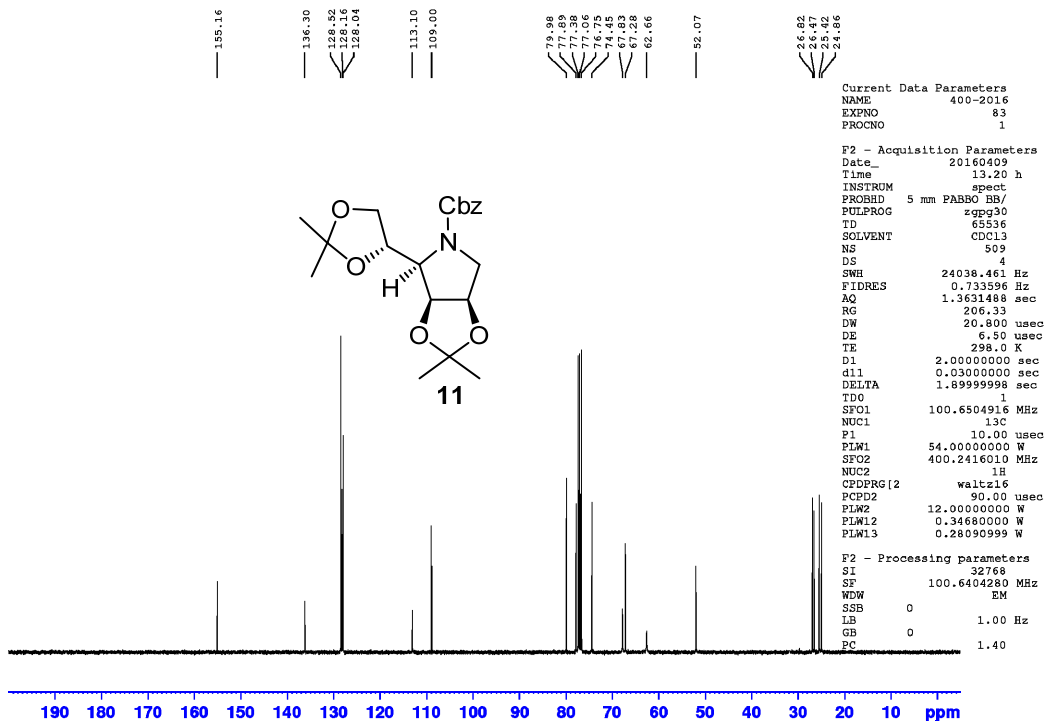
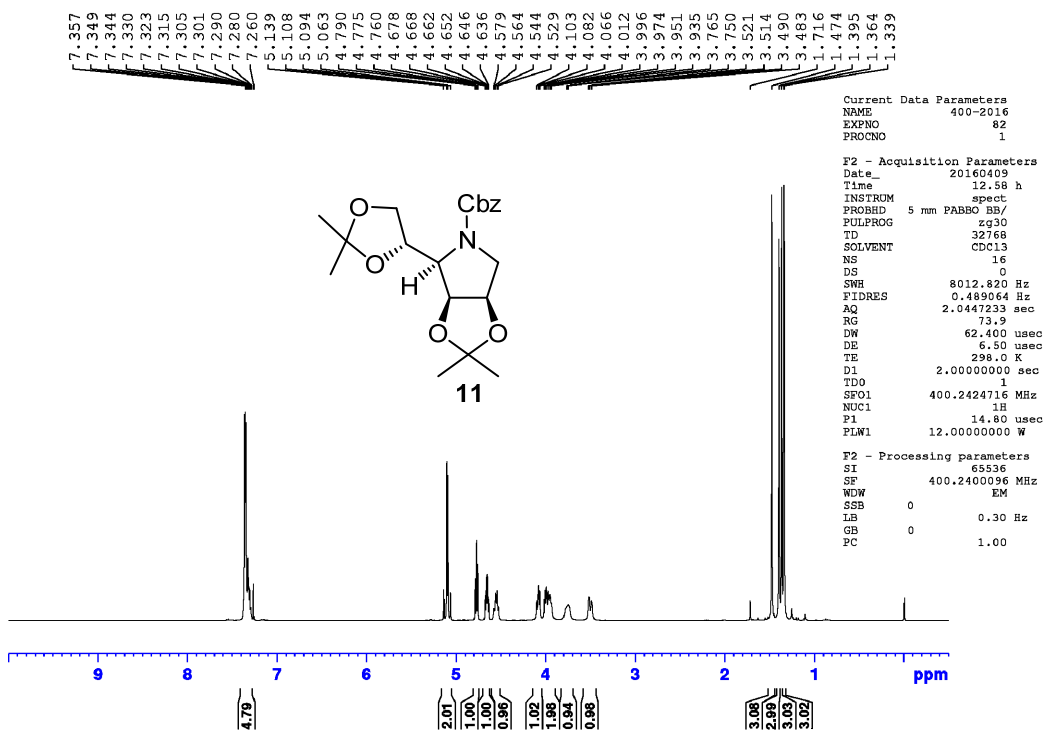


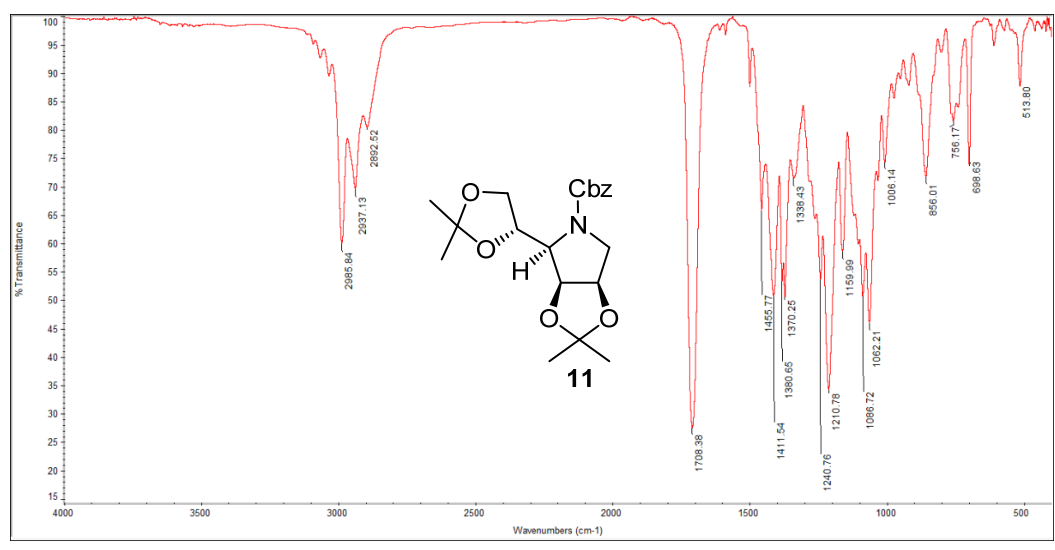
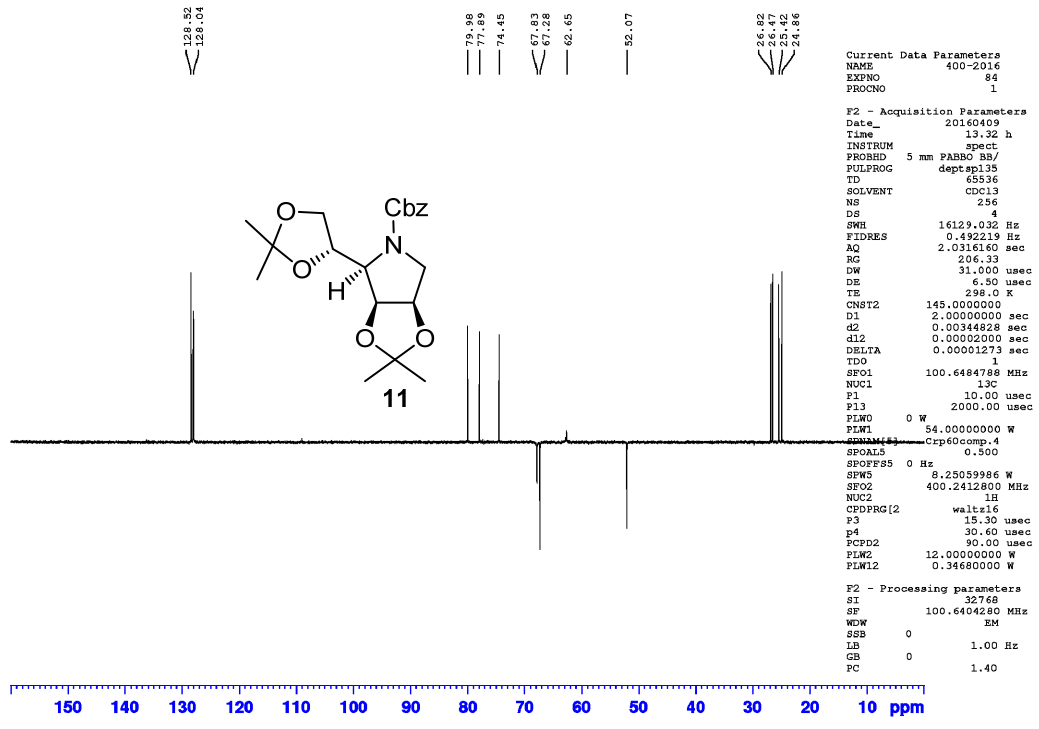
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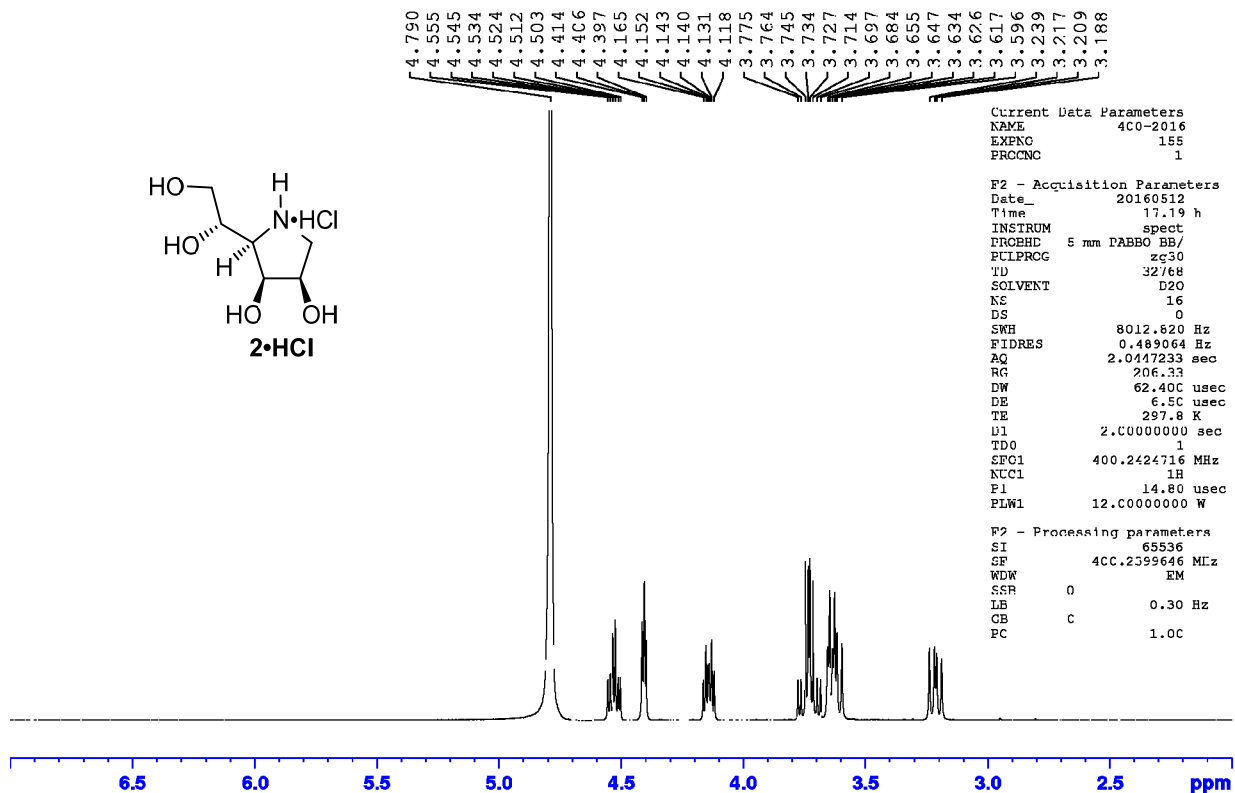
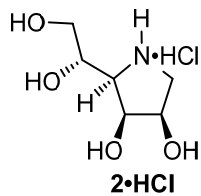


Compound 11:





Compound 2 • HCl:

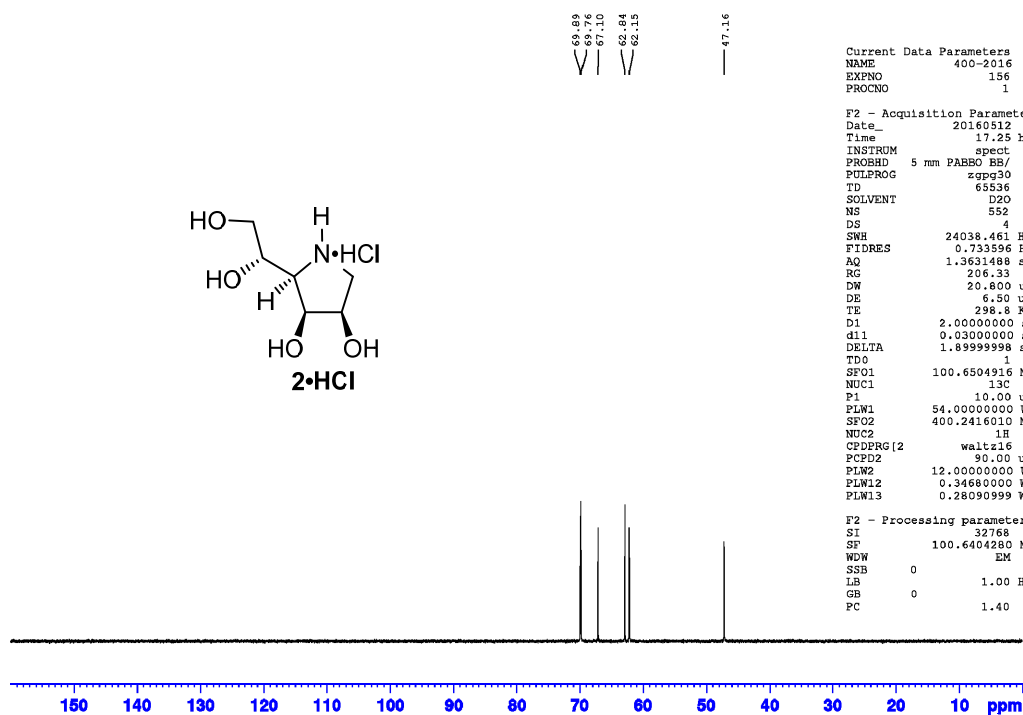
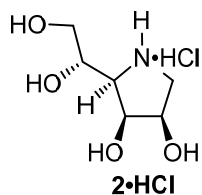


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PC            1.0C
    
```

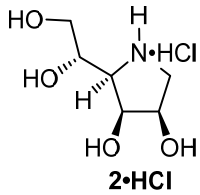


```

Current Data Parameters
NAME          400-2016
EXPNO         156
PROCNO        1

F2 - Acquisition Parameters
Date_         20160512
Time          17.25 h
INSTRUM       spect
PROBHD        5 mm PABBO BB/
PULPROG       zgpg30
TD            65536
SOLVENT       D2O
NS            552
DS            4
SWH           24038.461 Hz
FIDRES        0.733596 Hz
AQ            1.3631488 sec
RG            206.33
DE            20.800 usec
TE            298.8 K
D1            2.0000000 sec
d11           0.0300000 sec
DELTA         1.8999999 sec
TD0           1
SFO1          100.6504916 MHz
NUC1          13C
P1            10.00 usec
PLW1          54.0000000 W
SFO2          400.2416010 MHz
NUC2          1H
CPDPRG2       waltz16
PCPD2         90.00 usec
PLW2          12.0000000 W
PLW12         0.3468000 W
PLW13         0.28090999 W

F2 - Processing parameters
SI            32768
SF            100.6404280 MHz
WDW           EM
SSB           0
LB            1.00 Hz
GB            0
PC            1.40
    
```



69.87
 69.74
 67.07
 62.82
 62.12
 47.13

Current Data Parameters

NAME 400-2016
 EXPNO 157
 PROCNO 1

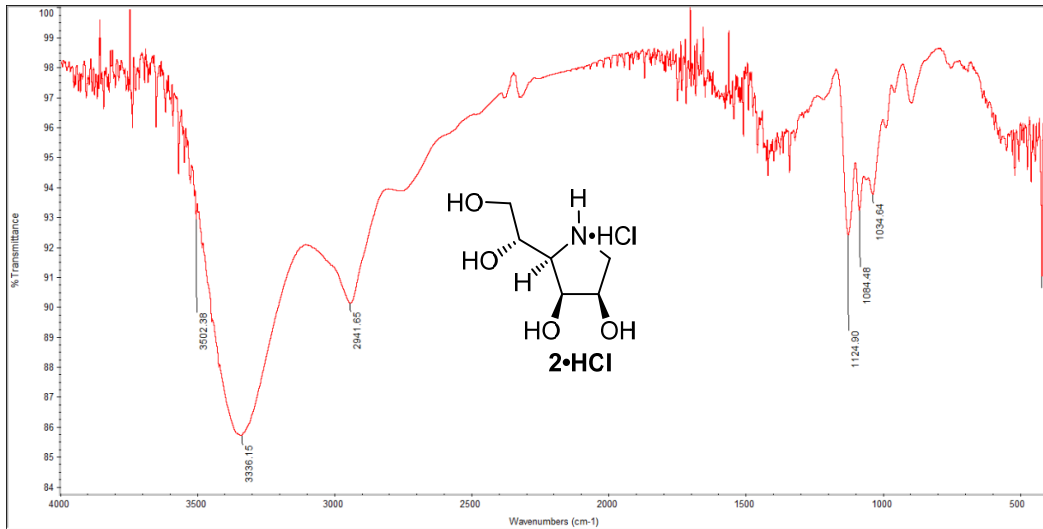
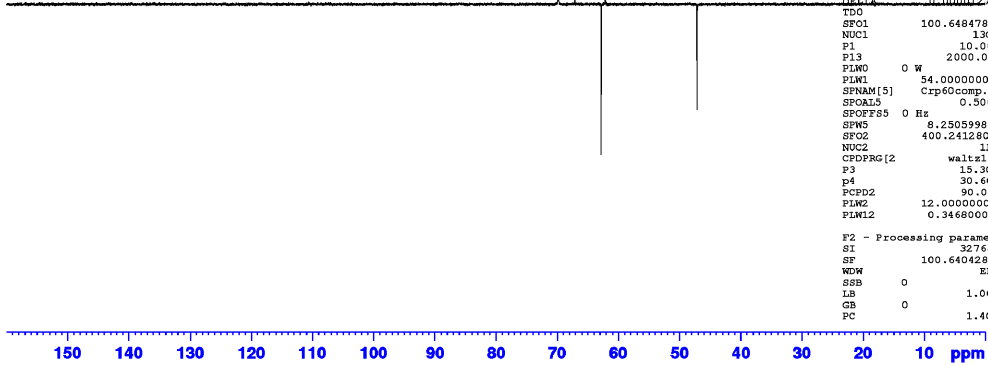
F2 - Acquisition Parameters

Date_ 20160512
 Time 17.56 h
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG deptspl33
 TD 65536
 SOLVENT D2O
 NS 96
 DS 4
 SWH 16129.032 Hz
 FIDRES 0.492219 Hz
 AQ 2.0316160 sec
 RG 206.33
 DW 31.000 usec
 DE 6.50 usec
 TE 298.7 K
 CNST2 145.0000000
 d1 2.00000000 sec
 d2 0.00344828 sec
 d12 0.00002000 sec
 DELTA 0.00001273 sec

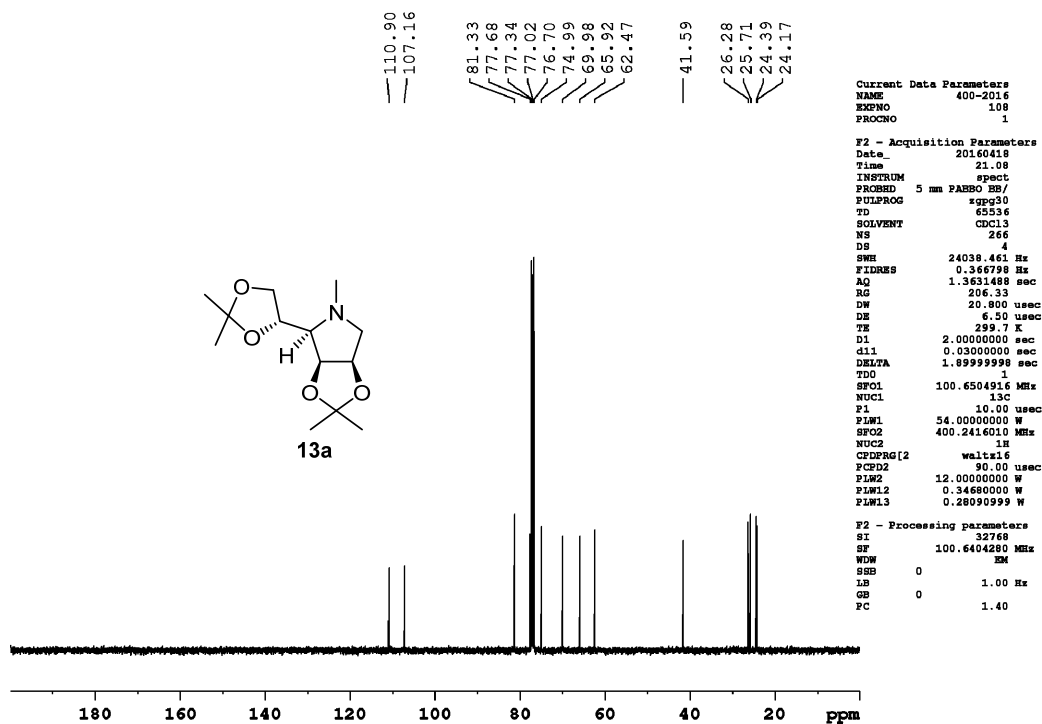
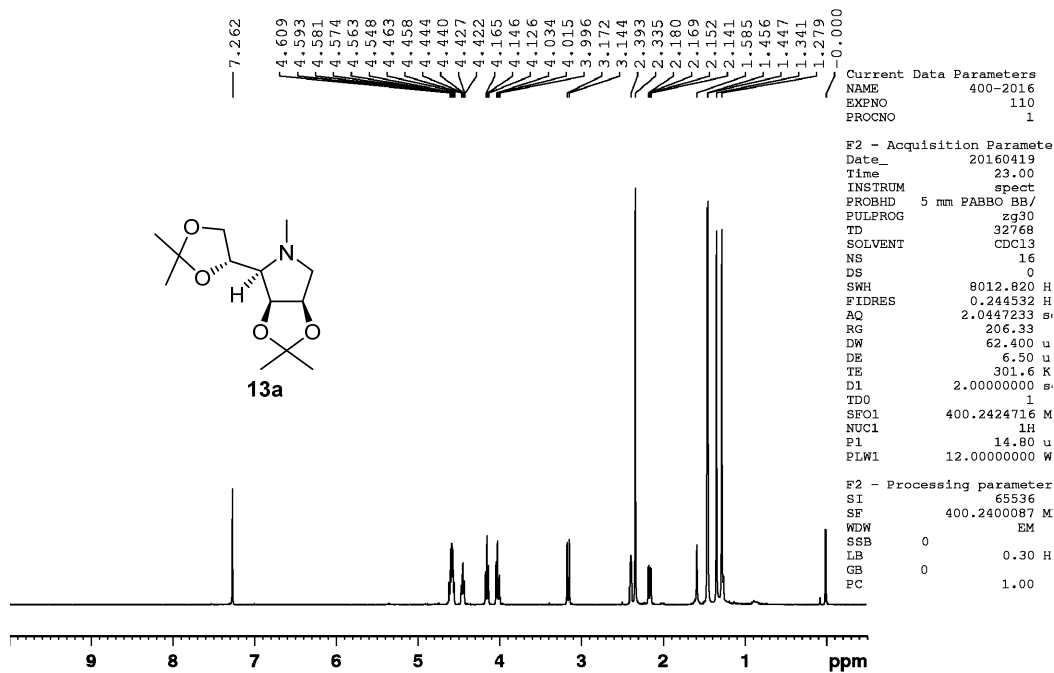
TDO 1
 SFO1 100.6484788 MHz
 NUC1 13C
 P1 10.00 usec
 P13 2000.00 usec
 PLW0 0 W
 PLW1 54.00000000 W
 SPNAM[5] Crp60comp.4
 SFOAL5 0.500
 SFOF53 0 Hz
 SPW5 8.25059986 W
 SFO2 400.2412800 MHz
 NUC2 1H
 CPDPRG2 waltz16
 P3 15.30 usec
 p4 30.60 usec
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.34680000 W

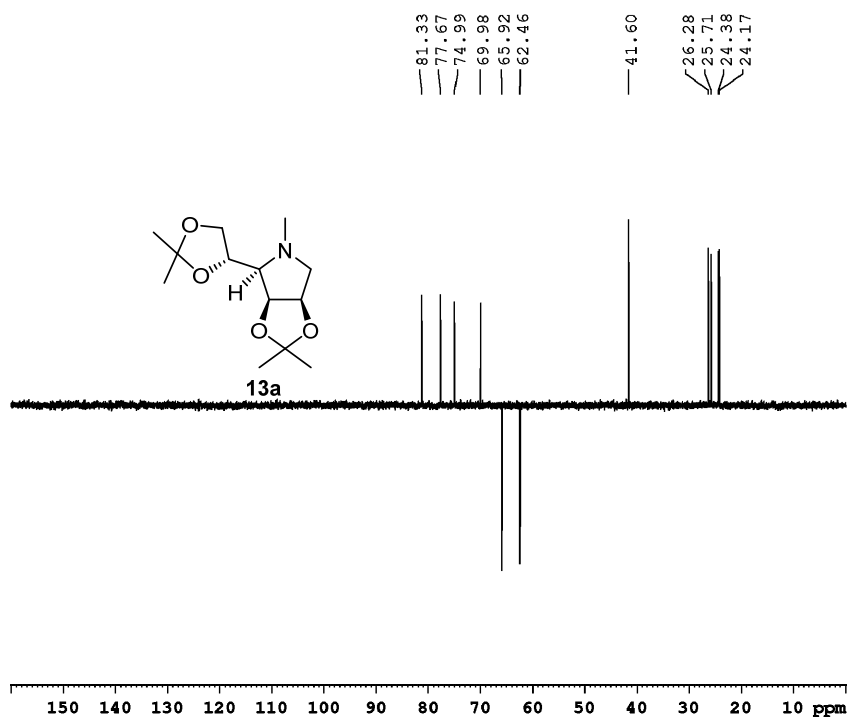
F2 - Processing parameters

SI 32768
 SF 100.6404280 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40



Compound 13a:





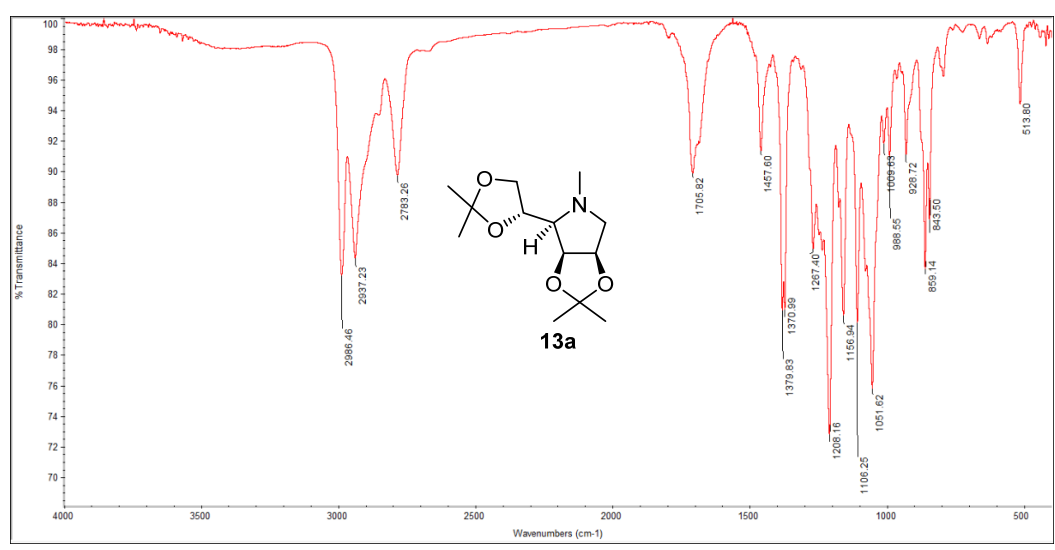
```

Current Data Parameters
NAME      400-2016
EXPNO    109
PROCNO    1

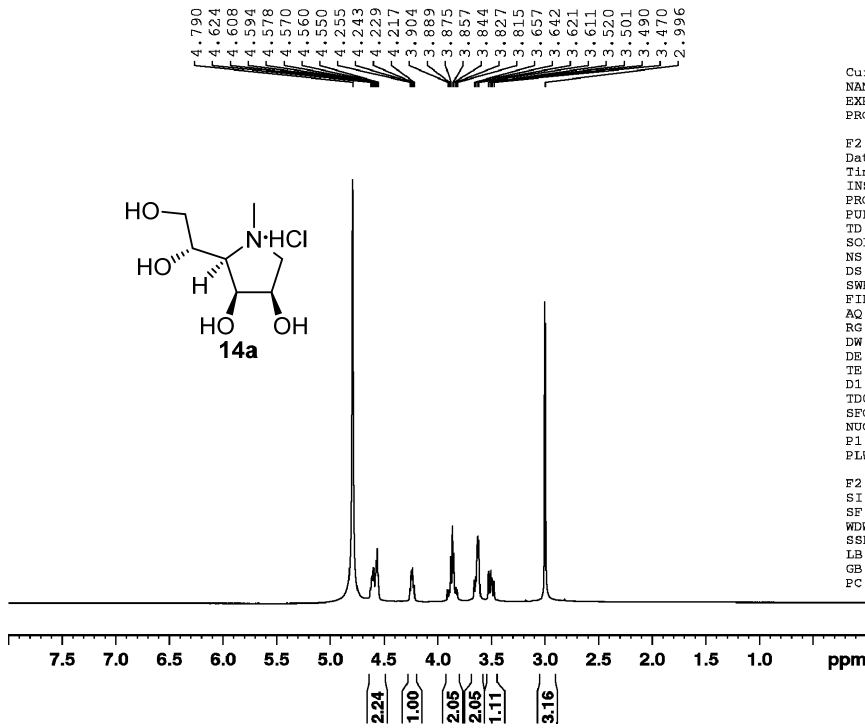
F2 - Acquisition Parameters
Date_     20160418
Time      21.19
INSTRUM   spect
PROBHD    5 mm PABBO BB/
PULPROG   zgpg30
TD         65536
SOLVENT   CDCl3
NS         81
DS         4
SWH        16129.032 Hz
FIDRES     0.246110 Hz
AQ         2.0216180 sec
RG         206.39
DW         31.000 usec
DE         6.50 usec
TE         299.2 K
CQF2       145.000000
SI         2.0000000 sec
d2         0.00344828 sec
d12        0.00002000 sec
DELTA     0.00001273 sec
VFO        1
VFO1       100.6464788 MHz
NUC1        13C
P1         10.00 usec
P13        2000.00 usec
PLMG       0 W
PLM1       54.0000000 W
SFOFF(5)   Csp50comp.4
SFOFF5     0 Hz
SFOFFAS    0 Hz
SFW5       8.25099986 W
SFO2       400.2412800 MHz
NUC2        1H
CHPRG2     waltz16
V3         15.30 usec
p4         30.60 usec
PCPD2      90.00 usec
PLM2       12.0000000 W
PLML2      0.34690000 W

F2 - Processing parameters
SI         32768
SF         100.6404280 MHz
WDW        EM
SSB         0
LB         1.00 Hz
GB         0
PC         1.40

```



Compound 14a:

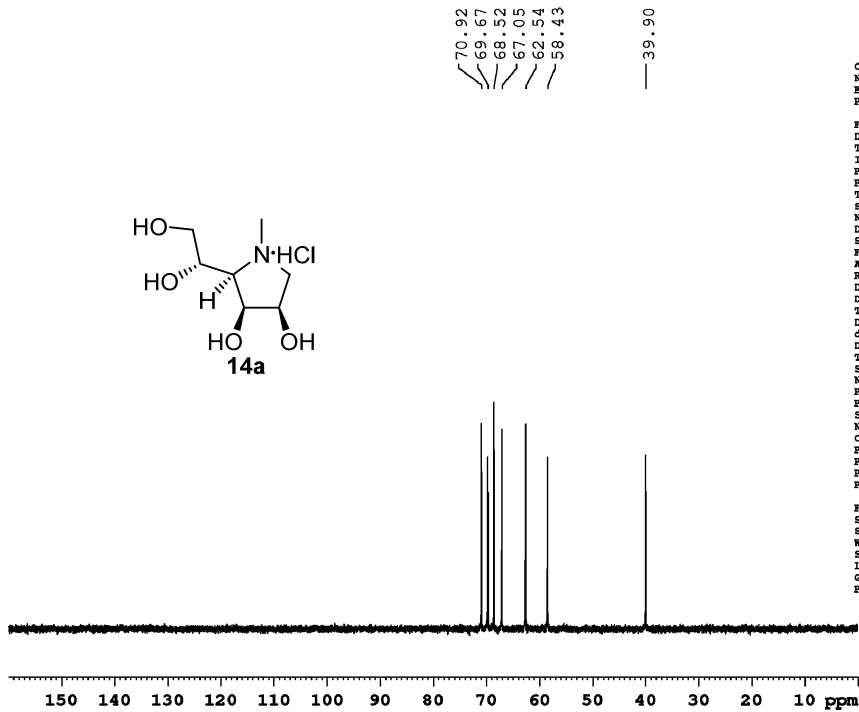


```

Current Data Parameters
NAME      400-2016
EXPNO    114
PROCNO   1

F2 - Acquisition Paramete
Date_    20160420
Time     16.19
INSTRUM spect
PROBHD   5 mm PABBO BB/
PULPROG  zg30
TD        32768
SOLVENT  D2O
NS        16
DS        0
SWH       8012.820 H
FIDRES    0.244532 H
AQ        2.0447233 s
RG        102.73
DW        62.400 u
DE        6.50 u
TE        301.2 K
D1        2.00000000 s
TDO       1
SFO1     400.2424716 M
NUC1     1H
P1       14.80 u
PLW1     12.00000000 W

F2 - Processing parameter
SI       65536
SF       400.2399658 M
WDW      EM
SSB      0
LB       0.30 H
GB       0
PC       1.00
    
```

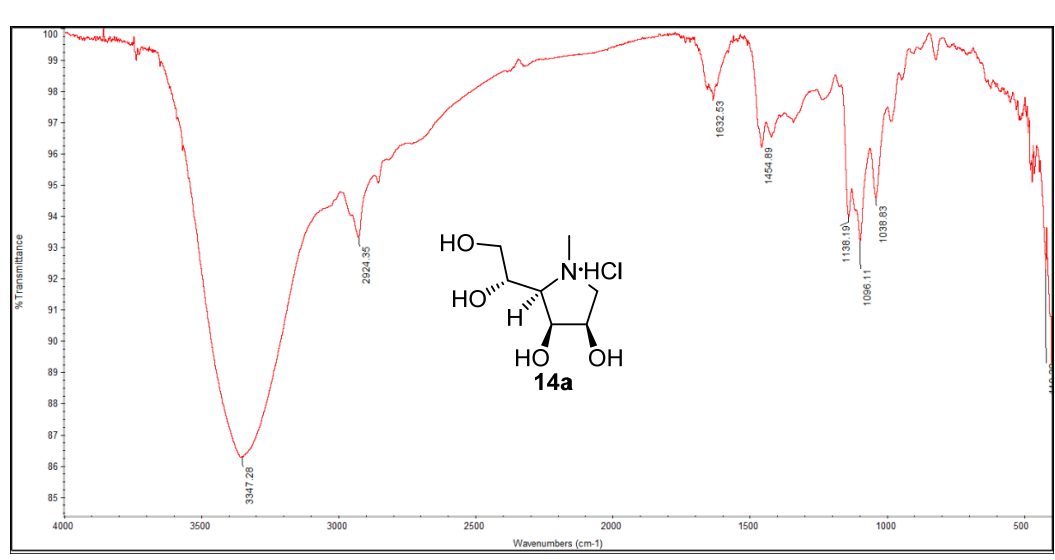
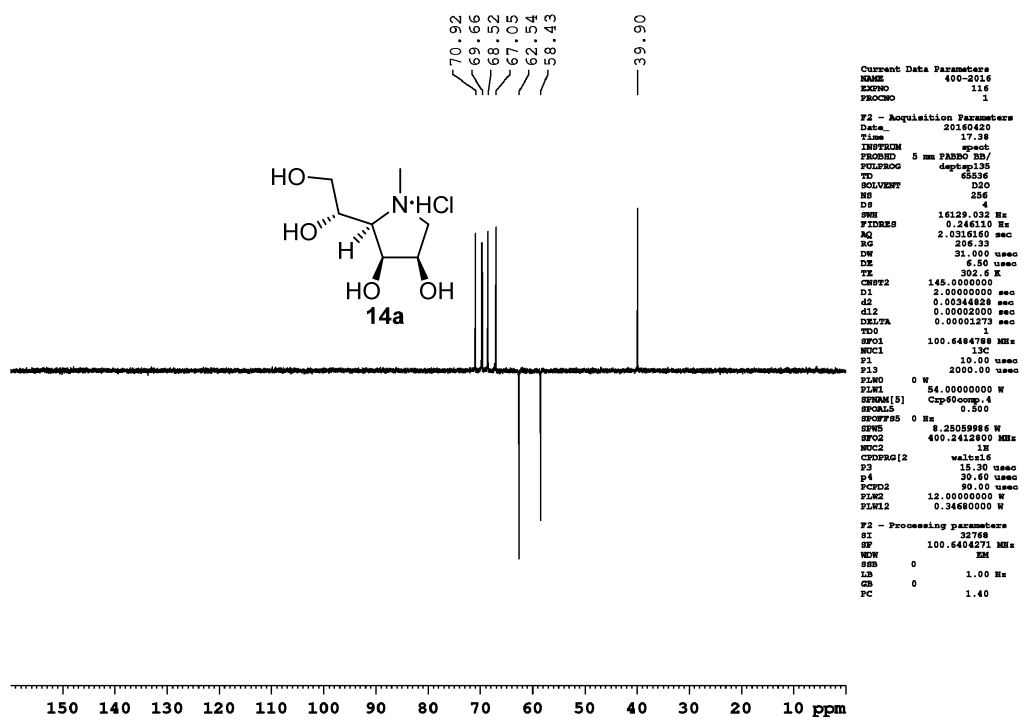


```

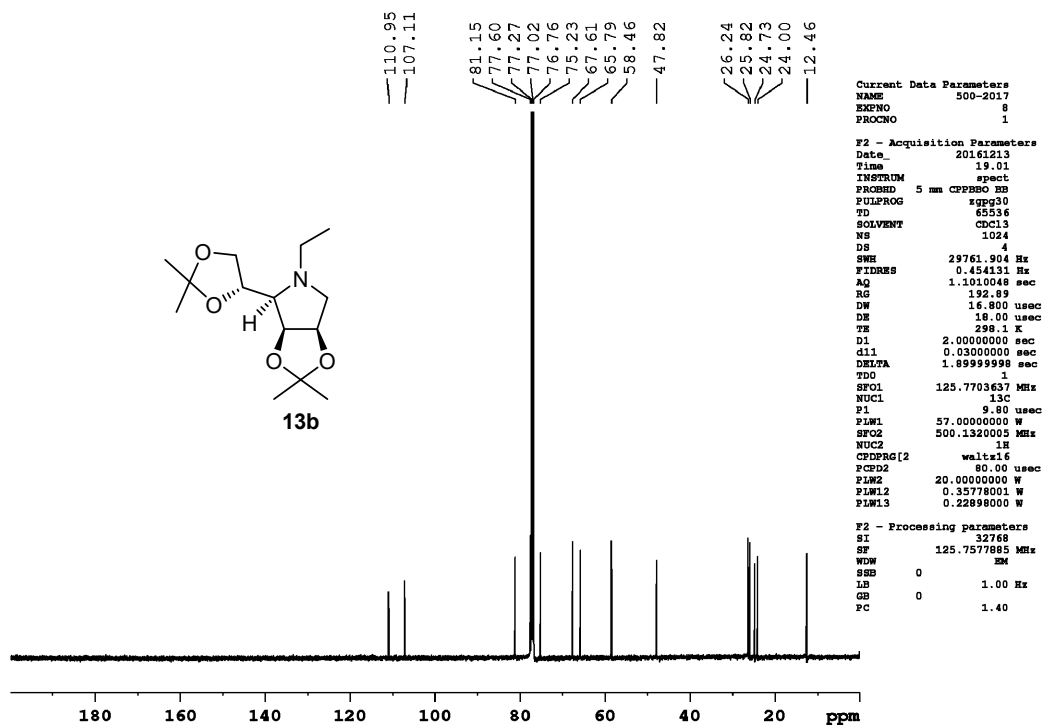
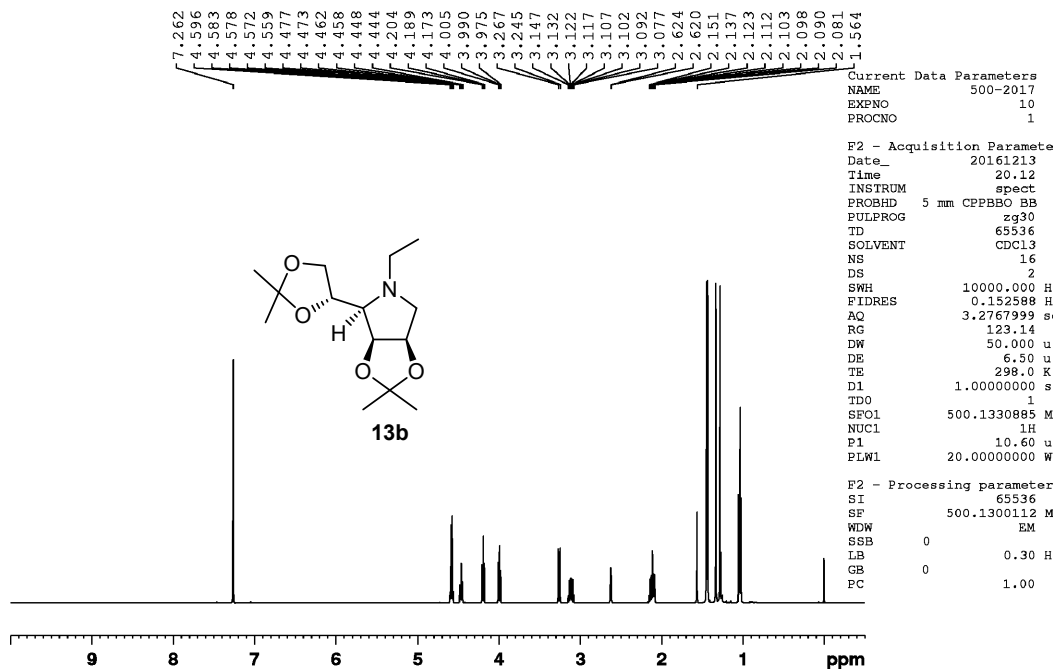
Current Data Parameters
NAME      400-2016
EXPNO    115
PROCNO   1

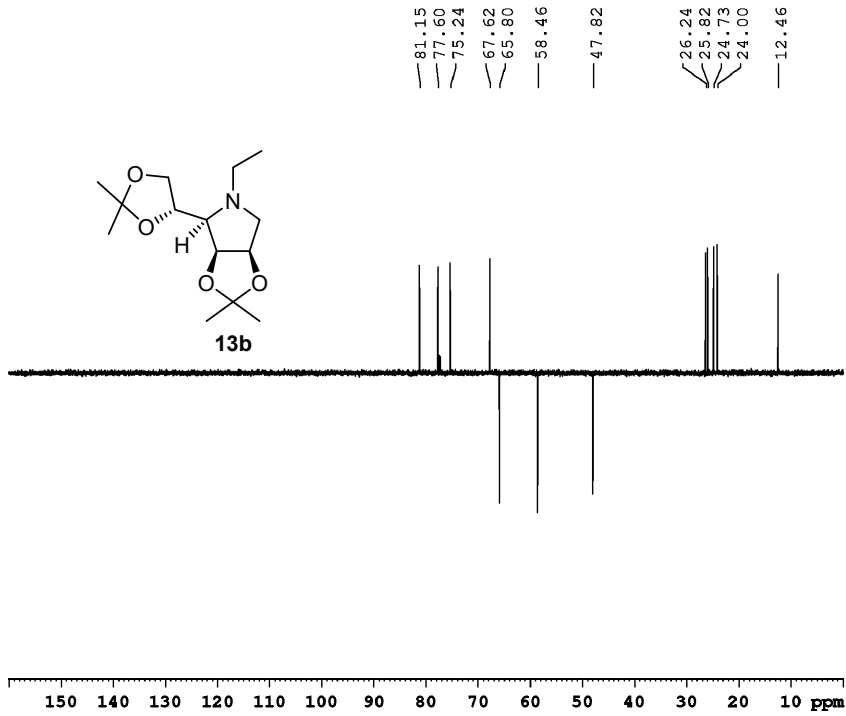
F2 - Acquisition Parameters
Date_    20160420
Time     17.19
INSTRUM spect
PROBHD   5 mm PABBO BB/
PULPROG  zgpg30
TD        55536
SOLVENT  D2O
NS        1024
DS        4
SWH       24038.461 Hx
FIDRES    0.366798 Hx
AQ        1.3631488 sec
RG        206.33
DW        20.800 usec
DE        6.50 usec
TE        303.0 K
D1        2.00000000 sec
d11       0.03000000 sec
DELTA     1.89999998 sec
TDO       1
SFO1     100.6504916 MHz
NUC1     13C
P1       10.00 usec
PLW1     54.00000000 W
SFO2     400.2416010 MHz
NUC2     1H
CPDPRG2  waltz16
PCPD2    90.00 usec
PLW2     12.00000000 W
PLW12    0.34680000 W
PLW13    0.28090999 W

F2 - Processing parameters
SI       32768
SF       100.6404280 MHz
WDW      EM
SSB      0
LB       1.00 Hx
GB       0
PC       1.40
    
```

Compound 13b:



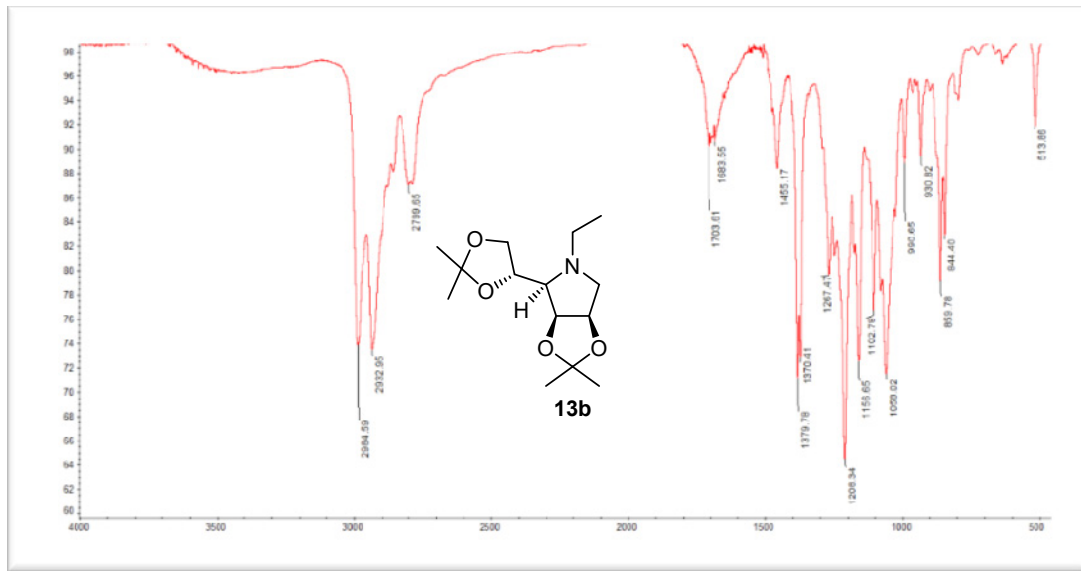


```

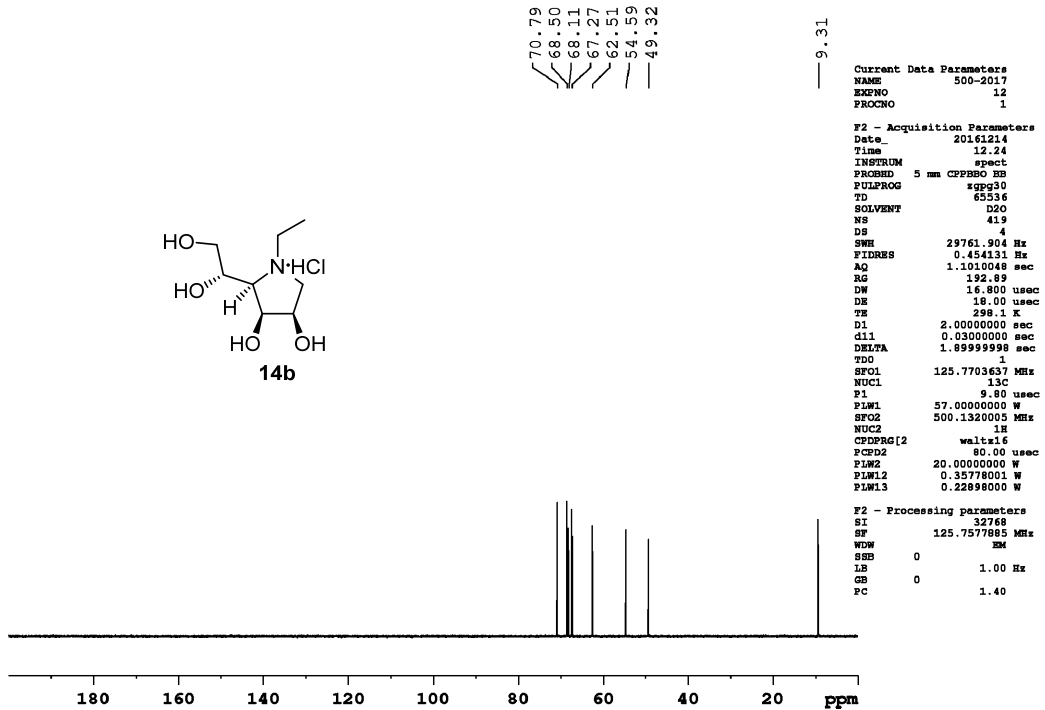
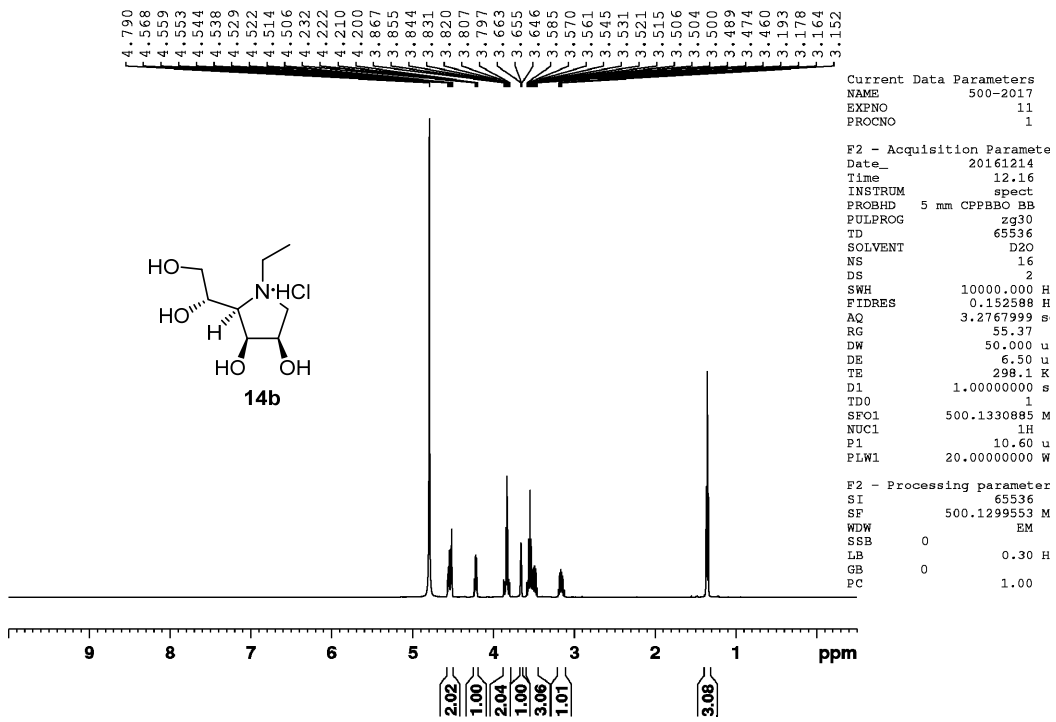
Current Data Parameters
NAME      800-2017
EXPNO     9
PROCNO    1

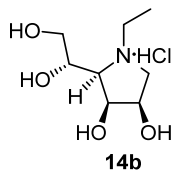
F2 - Acquisition Parameters
Date_     20161213
Time     20.10
INSTRUM   spect
PROBHD    5 mm CPYPRB0 BB
PULPROG   zgpg30
TD         65536
SOLVENT   CDCl3
NS         256
DS         4
SHE       20161.291 Hz
FIDRES    0.307637 Hz
AQ         1.6252928 sec
RG         192.89
DW         24.800 usec
DE         18.00 usec
TE         298.0 K
CHFT2     145.0000000
D1         2.0000000 sec
d2         0.00344828 sec
d12        0.0002000 sec
DELTA     0.0001248 sec
TDO        1
SFO1      125.767846 MHz
NUC1       13C
P1         9.80 usec
P2         2000.00 usec
PLM0      0 W
PLM1      37.0000000 W
SFO2      8.36410046 MHz
SFO3      500.1315995 MHz
SFO4      0.800
SFO5      0 Hz
SFO6      8.36410046 MHz
SFO7      500.1315995 MHz
NUC2       1H
CPDPRG2   waltz16
S3         10.70 usec
P4         21.40 usec
PCPD2     80.00 usec
PLM2      20.0000000 W
PLM3      0.35778001 W

F2 - Processing parameters
SI         32768
SF         125.767846 MHz
WDW        EM
SSB        0
LB         1.00 Hz
GB         0
PC         1.40
  
```



Compound **14b**:





70.79
68.50
68.11
67.27
62.51
54.59
49.32

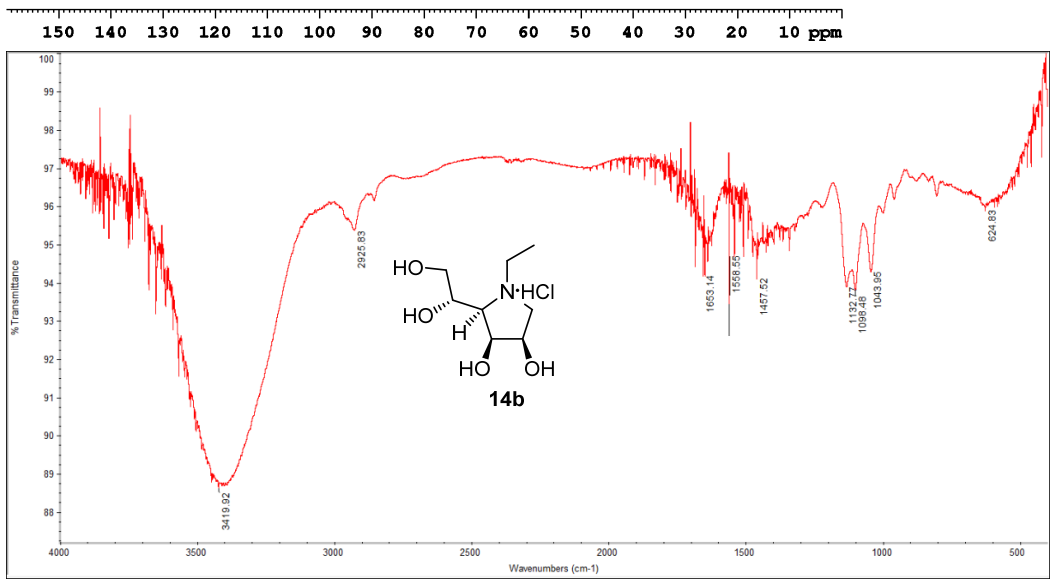
9.31

```

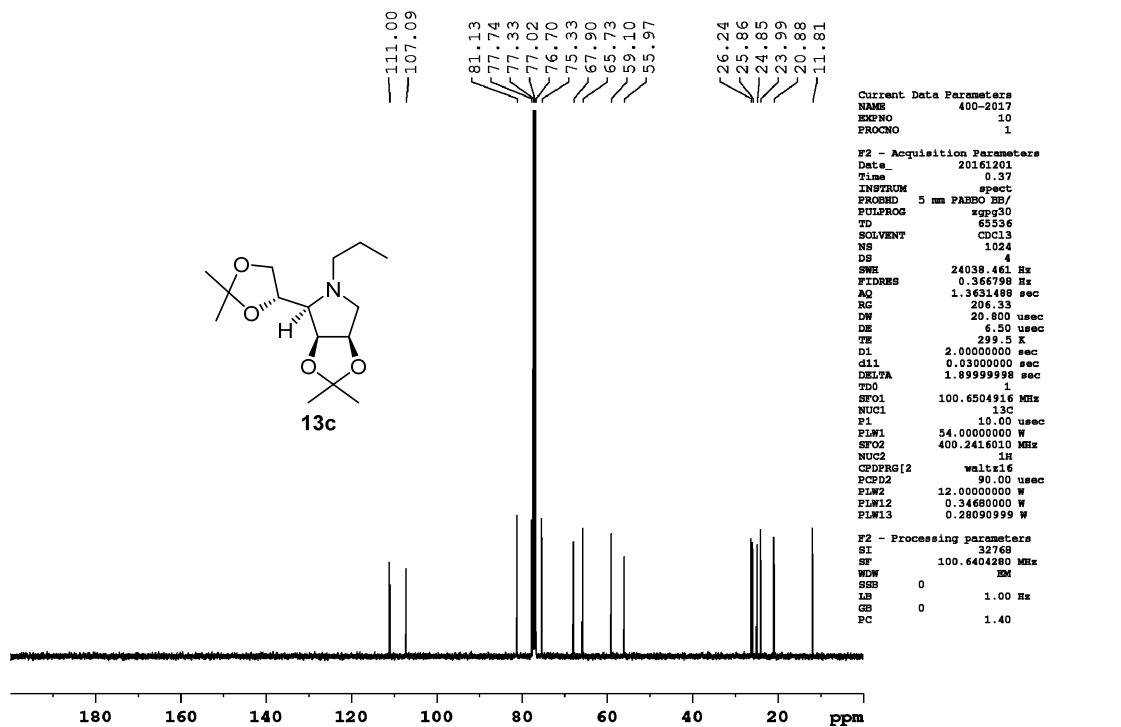
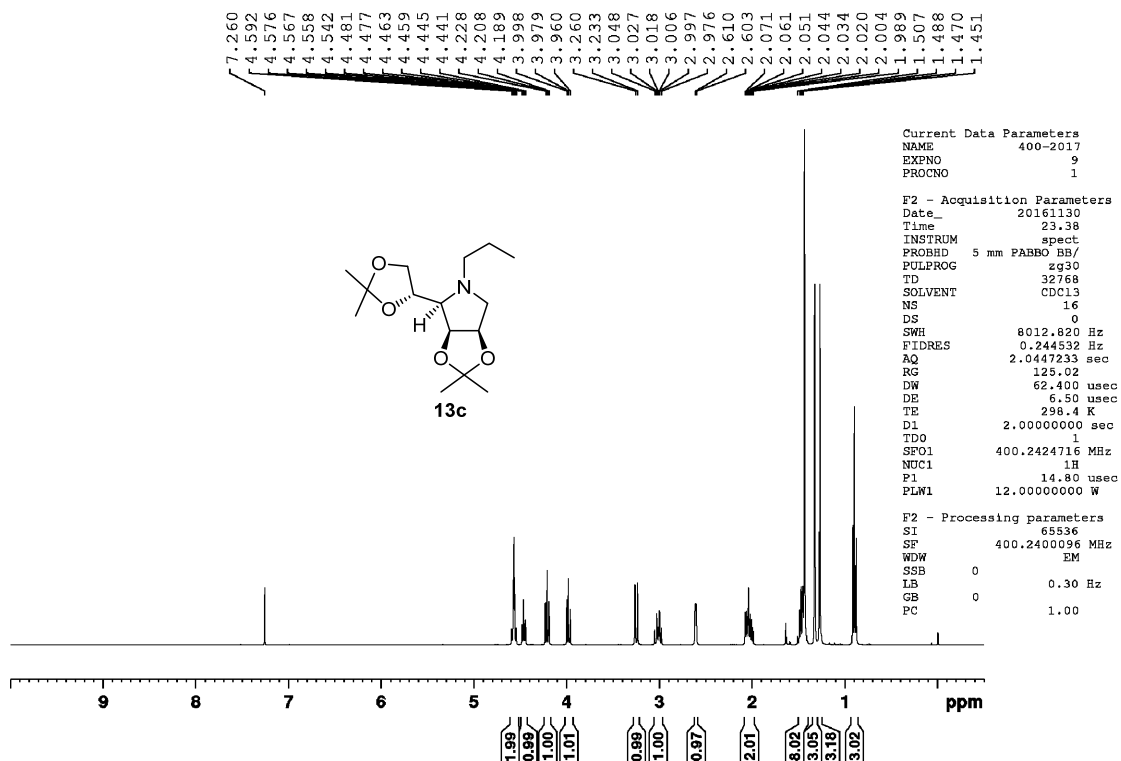
Current Data Parameters
NAME      500-2017
EXPNO    13
PROCNO   1

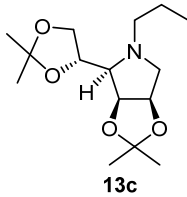
F2 - Acquisition Parameters
Date_    20161214
Time     12.44
INSTRUM  spect
PROBHD   5 mm CPYBBO BB
PULPROG  zgpg30
TD        65536
SOLVENT  D2O
NS        98
DS        4
SWH       20161.391 Hz
FIDRES   0.307637 Hz
AQ        1.6252928 sec
RG         192.89
AQ        24.800 usec
DE         18.00 usec
VE         298.0 K
CHFT2    145.000000
D1        2.00000000 sec
d2        0.00344828 sec
d12       0.00002000 sec
DELTA    0.0001248 sec
TDO       1
SFO1     125.7678486 MHz
NUC1      13C
P1         9.80 usec
P13       2000.00 usec
PLMO      0 W
PLM1     57.00000000 W
SFO2      Crp60comp_4
SFO2S     0.500
SFO2S5    0 Hz
SFO5      8.36410046 W
SFO5C     500.1218996 MHz
NUC2      1H
CHOPRO2[2] waltz16
P3         10.70 usec
p4         21.40 usec
PCPD2     80.00 usec
PLM2     20.00000000 W
PLM12    0.35778001 W

F2 - Processing parameters
SI         32768
SF         125.7577885 MHz
WDW        EM
SSB        0
LB         1.00 Hz
GB         0
PC         1.40
  
```



Compound 13c:





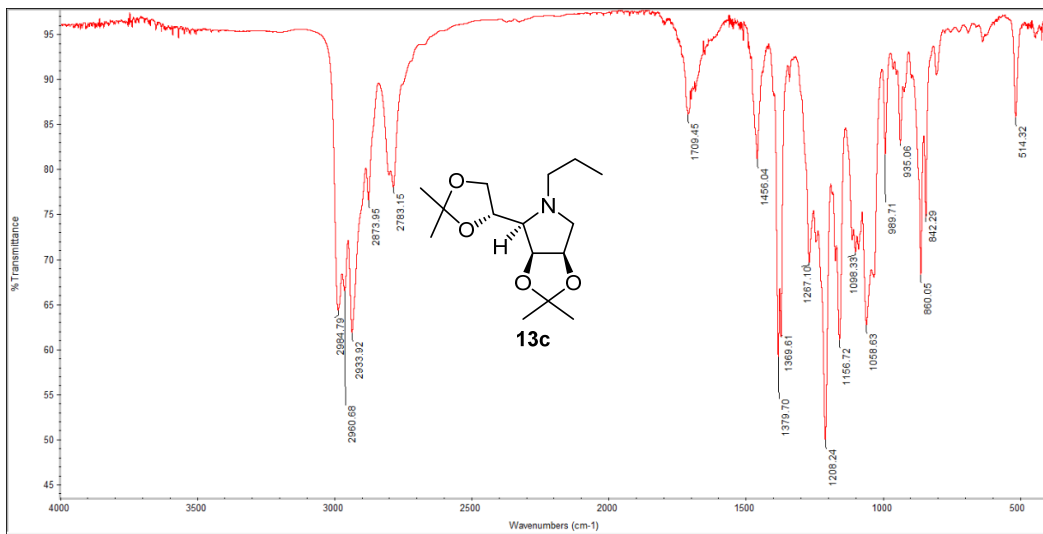
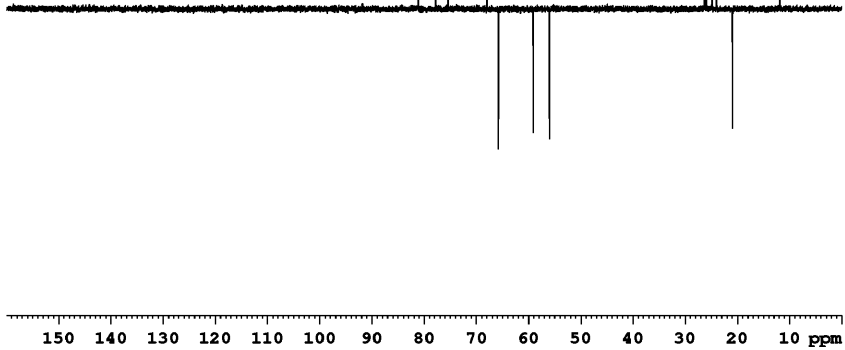
81.13
77.74
75.33
67.89
65.73
59.10
55.97
26.24
25.86
24.85
23.99
20.88
11.81

```

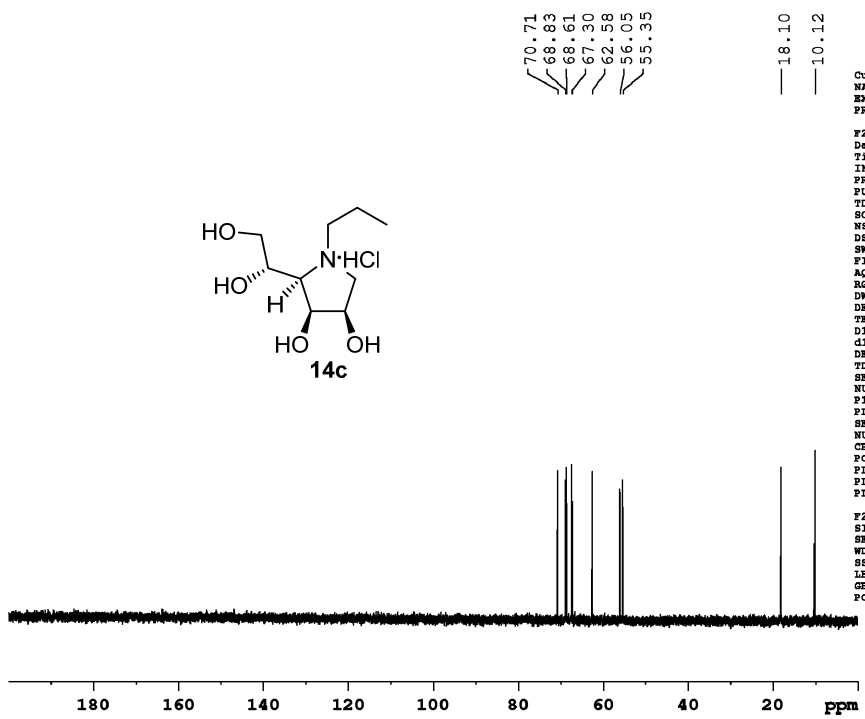
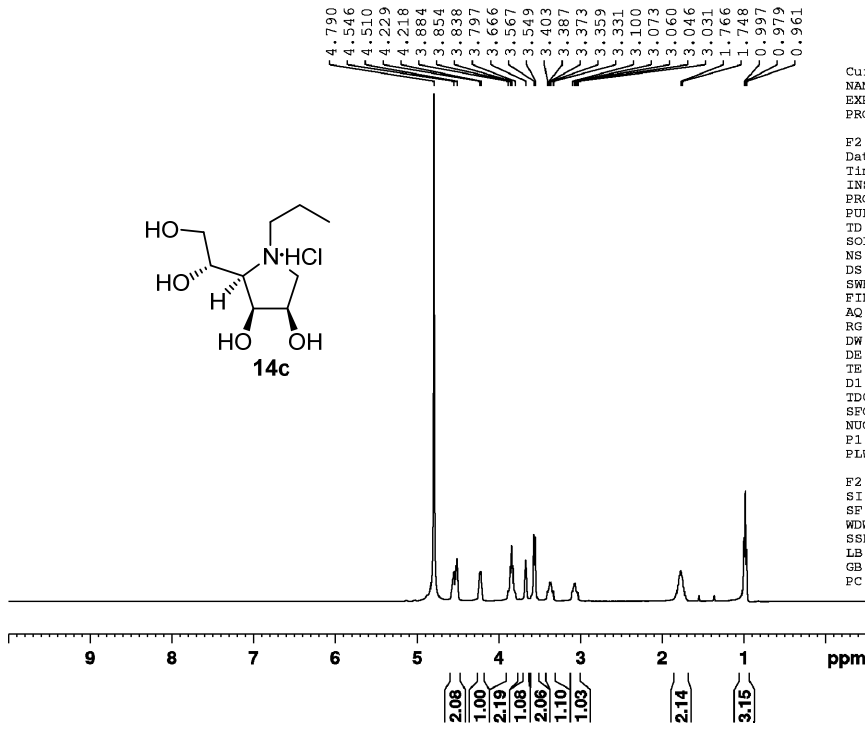
Current Data Parameters
NAME      400-2017
EXPNO    11
PROCNO   1

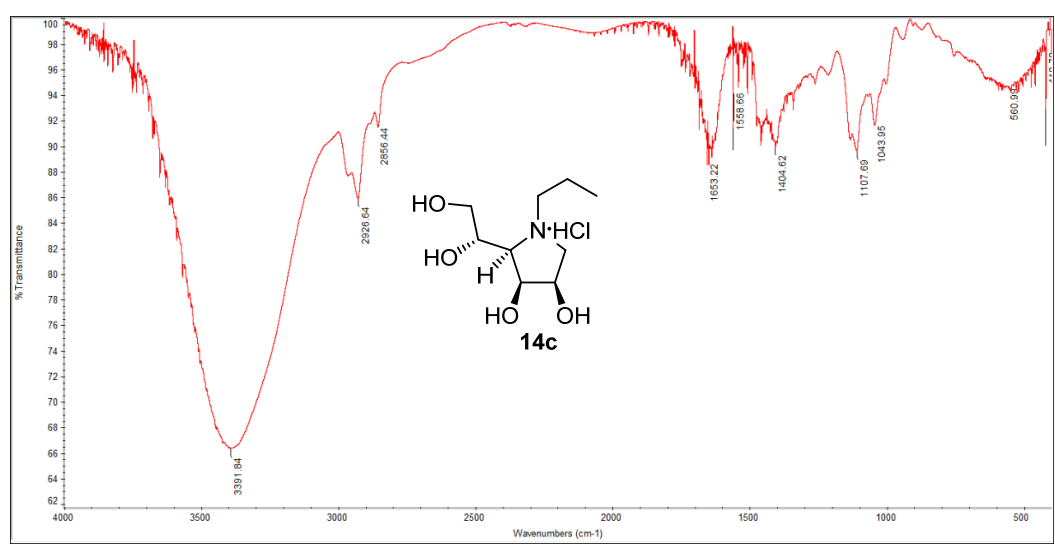
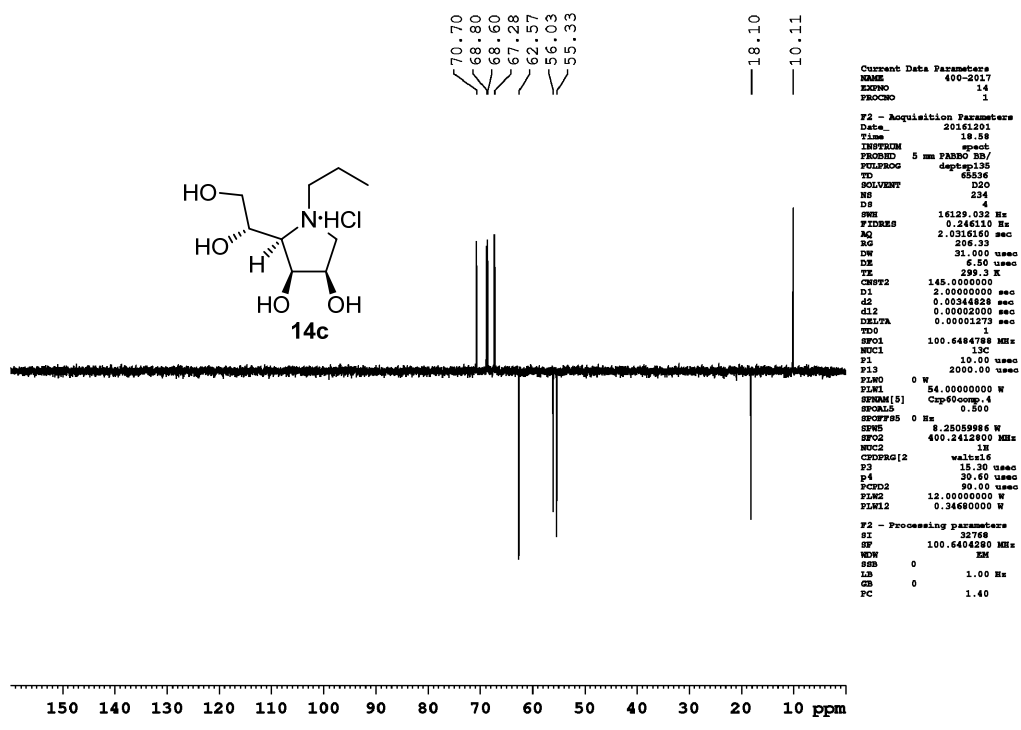
F2 - Acquisition Parameters
Date_    20161201
Time     0.55
INSTRUM  spect
PROBHD   5 mm PABBO BB/
PULPROG  zgpg30
TD        65536
SOLVENT  CDCl3
NS        256
DS        4
SWH       16129.932 Hz
FIDRES   0.246110 Hz
AQ        2.0216180 sec
RG         206.39
DW        31.000 usec
DE        6.50 usec
TE        299.0 K
CQF2     145.000000
D1        2.0000000 sec
d2        0.00344628 sec
d12       0.00002000 sec
DELTA    0.00001273 sec
VFO1     100.6464788 MHz
NUC1      13C
P1        10.00 usec
E13       2000.00 usec
PLM0      0 W
PLM1     54.0000000 W
SFOFF(5) Csp60comp.4
SFOFF(5) 0.500
SFOFF(5) 0 Hz
SFOFF(5) 8.25059988 W
SFOFF(5) 400.2412800 MHz
NUC2      1H
CQFPG[2] waltz16
P2        15.30 usec
P4        30.60 usec
PCPD2    90.00 usec
PLM2     12.0000000 W
PLM12    0.34680000 W

F2 - Processing parameters
SI        32768
SF        100.6404280 MHz
WDW       EM
SSB       0
LB        1.00 Hz
GB        0
PC        1.40
  
```

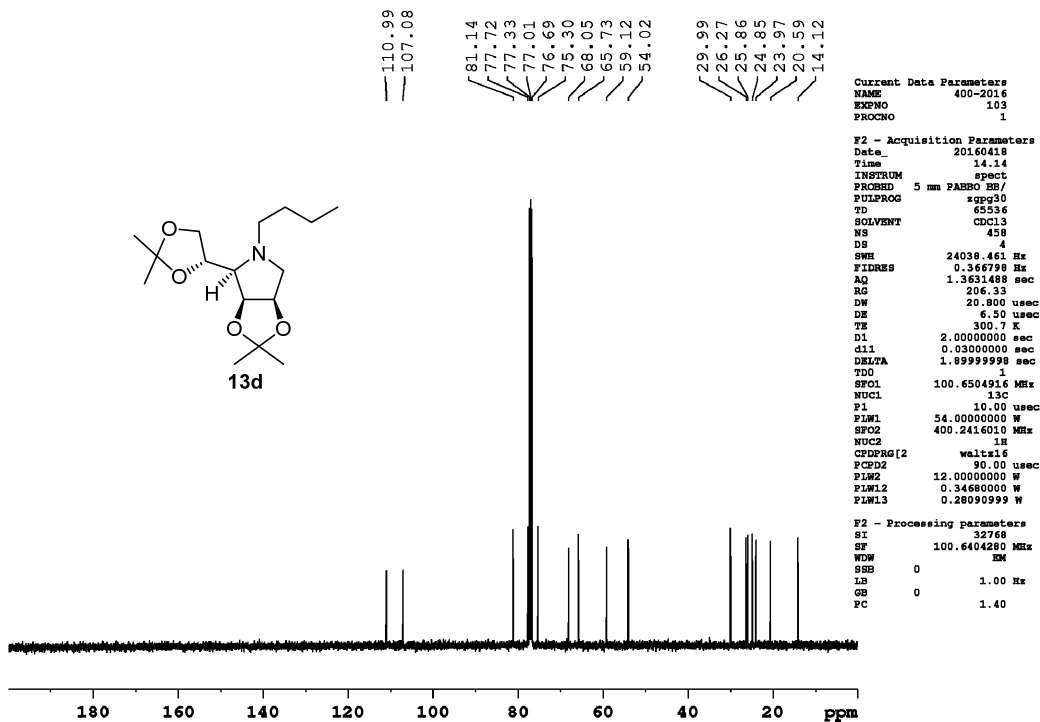
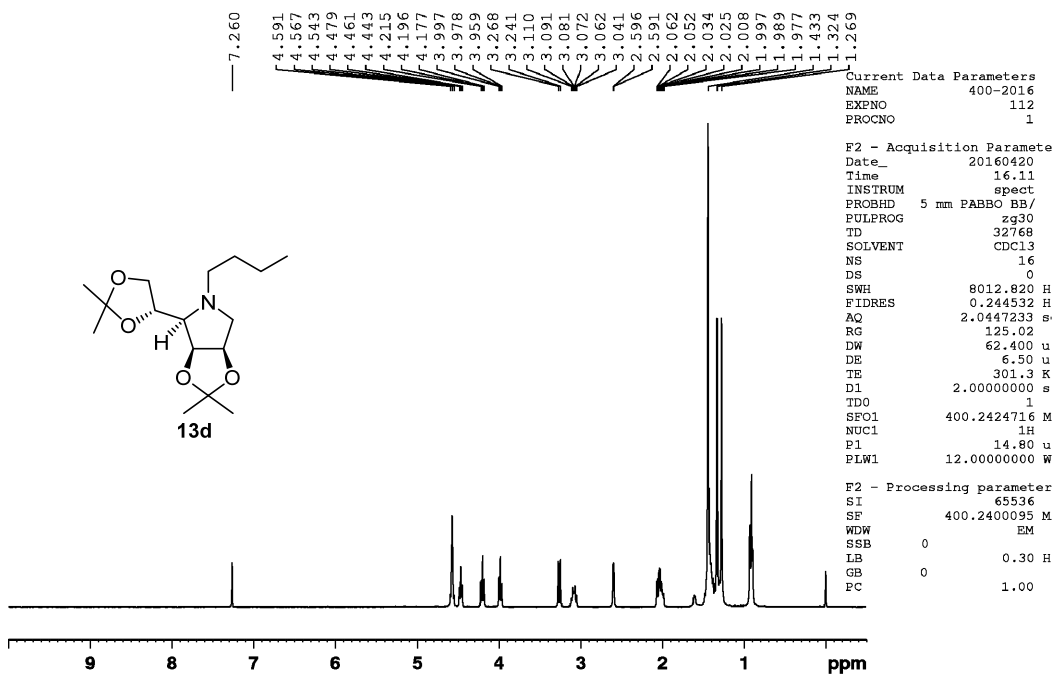


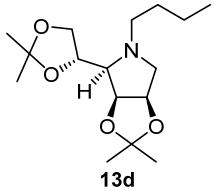
Compound 14c:



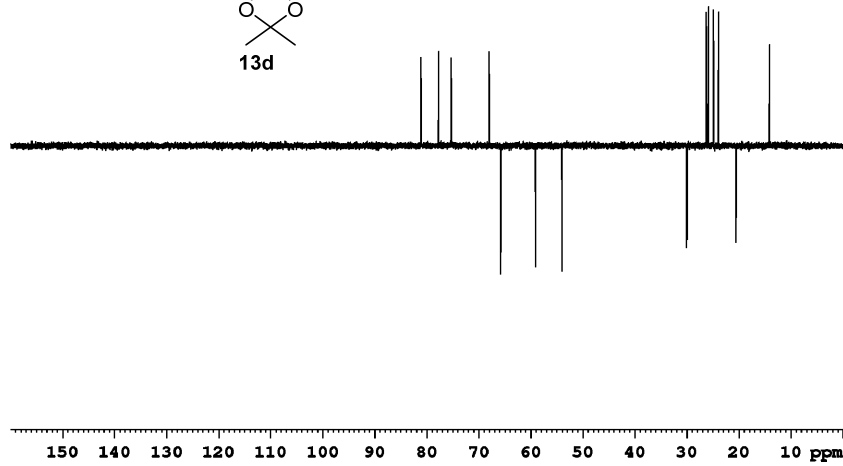


Compound 13d:





- 81.14
- 77.71
- 75.30
- 68.05
- 65.73
- 59.11
- 54.03
- 29.99
- 26.27
- 25.86
- 24.85
- 23.96
- 20.59
- 14.12

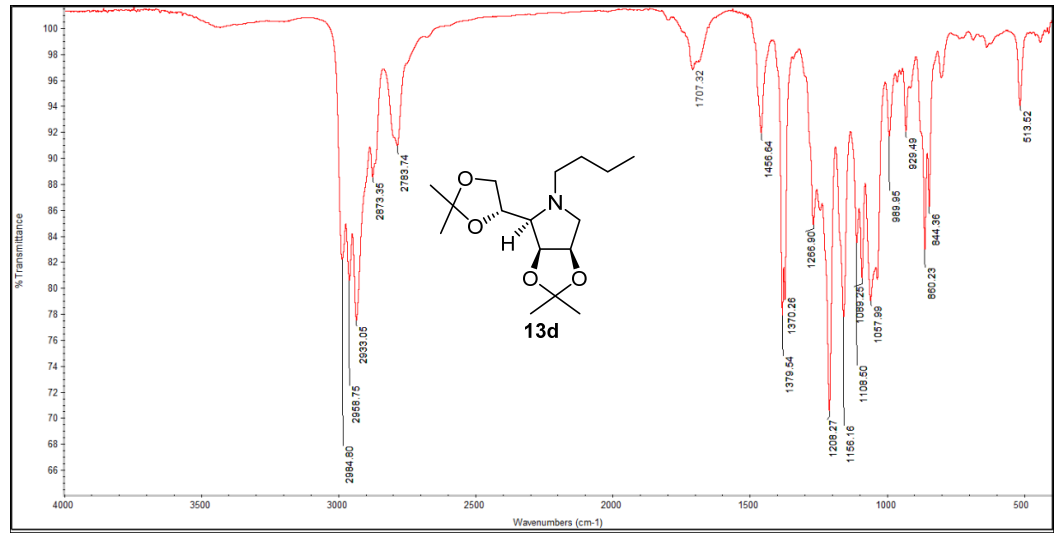


```

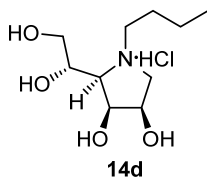
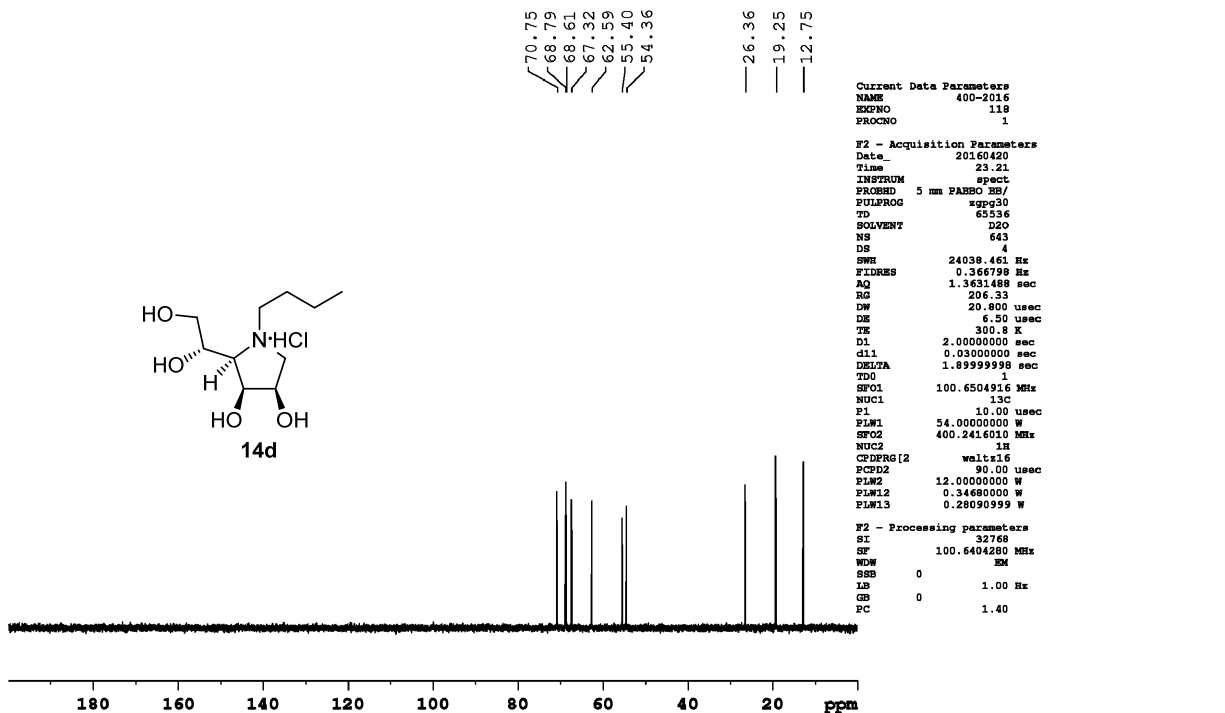
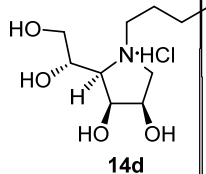
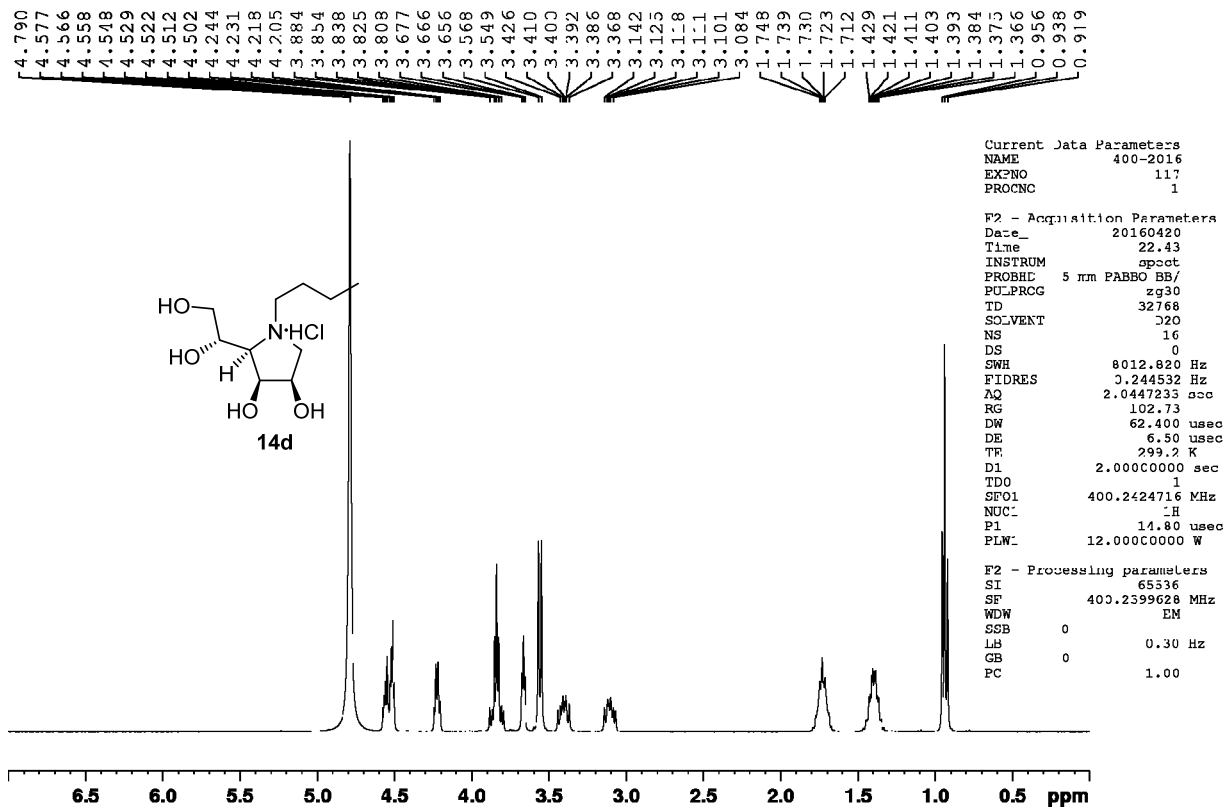
Current Data Parameters
NAME      400-2016
XDRNO    106
PROCNO    1

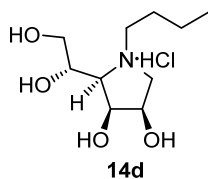
F2 - Acquisition Parameters
Date_     20160418
Time      16.24
INSTRUM   spect
PROBHD    5 mm PABBO BB/
PULPROG   zgpg30
TD         65536
SOLVENT   CDCl3
NS         162
DS         4
SWH        16129.022 Hz
FIDRES    0.246110 Hz
AQ         2.0316180 sec
RG         206.33
DW         31.000 usec
DE         6.50 usec
TE         298.6 K
CHWT2     148.000000
D1         2.0000000 sec
d2         0.00344828 sec
d12        0.00000000 sec
DELTA     0.00001273 sec
SFO1       100.6484788 MHz
NUC1       13C
P1         10.00 usec
P13        2000.00 usec
O W        0
ELW0       54.0000000 W
SFM1       Csp60comp.4
SFOFFAS    0 Hz
SWS        8.25059986 W
SFO2       400.2412800 MHz
NUC2       1H
CHPRG[2]   waltz16
Y3         15.30 usec
p4         30.60 usec
NCH2       90.00 usec
ELW2       12.0000000 W
ELWL2     0.34680000 W

F2 - Processing parameters
SI         32768
SF         100.6404280 MHz
WDW        EM
SSB        0
LB         1.00 Hz
GB         0
PC         1.40
  
```



Compound 14d:





70.73
68.77
68.60
67.30
62.58
55.37
54.34

— 26.35
— 19.24
— 12.74

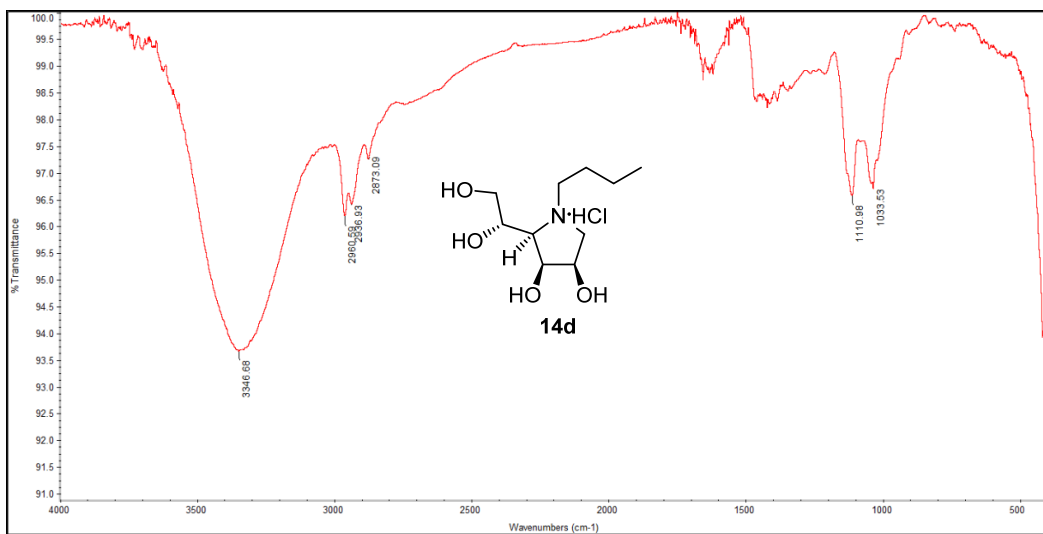
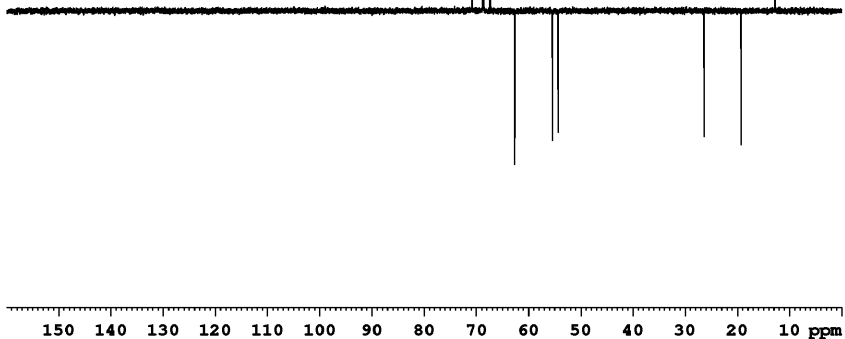
```

Current Data Parameters
NAME      400-2014
EXPNO    119
PROCNO   1

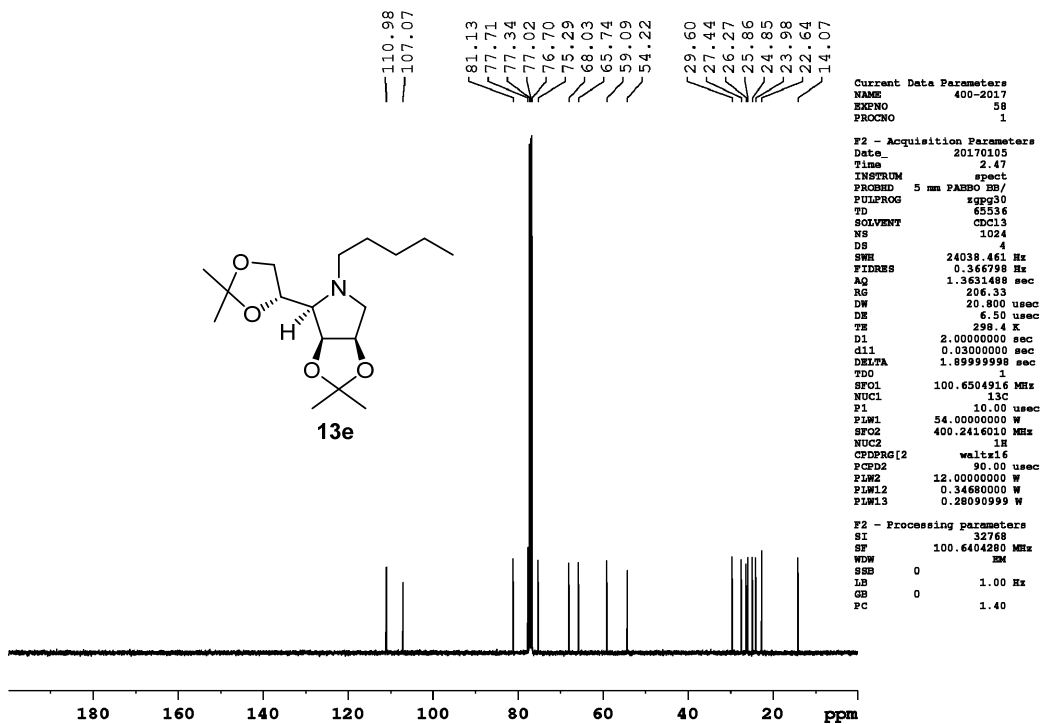
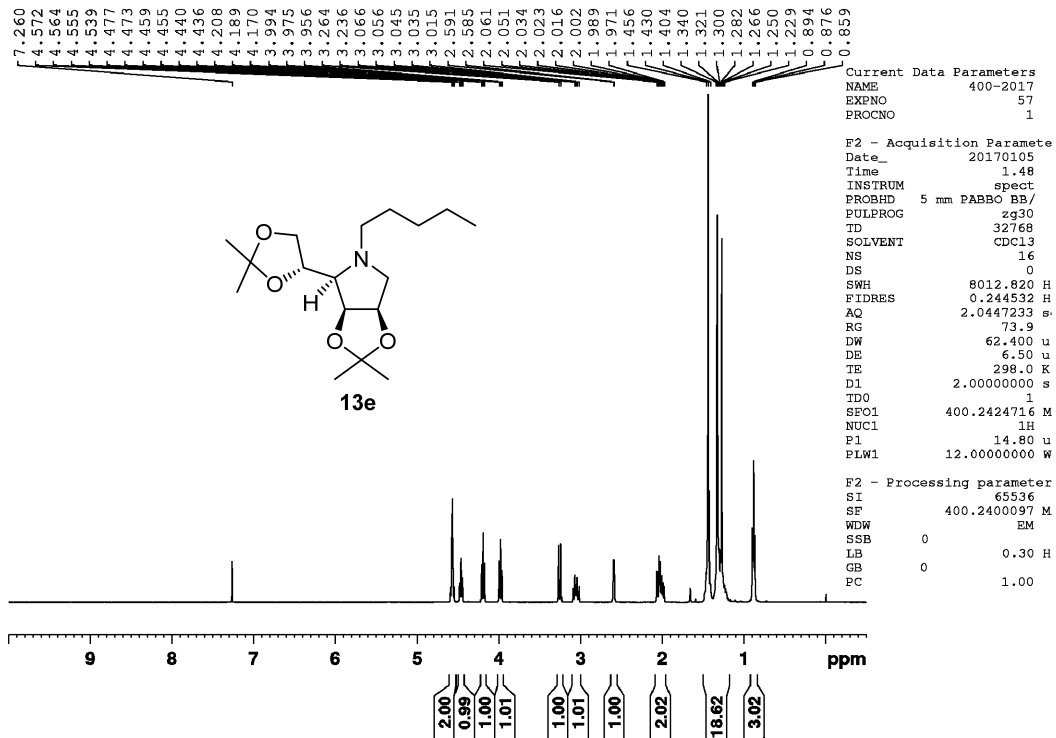
F2 - Acquisition Parameters
Date_    20160420
Time     23.24
INSTRUM  spect
PROBHD   5 mm F400 DD/
PULPROG  zgpg30
TD        65536
SOLVENT  D2O
NS        147
DS        4
SWH       16129.022 Hz
FIDRES   0.246110 Hz
AQ        2.0316180 sec
RG         206.33
DW         31.000 usec
DE         6.50 usec
TE        300.3 K
CQV2     145.000000
D1        2.0000000 sec
d2        0.00346828 sec
d12       0.00002000 sec
DELTA    0.00001273 sec
VDO       1
SFO1     100.6464788 MHz
NUC1      13C
P1        10.00 usec
P13       2000.00 usec
PLM0      0 W
PLM1     54.0000000 W
SFO1M[5] Csp60comp.4
SFO2M[5] 0 Hz
SFO2M[5] 0.500
SFO5     8.25059986 W
SFO6     400.2412800 MHz
NUC2      1H
CQV2G[2] waltz16
P2        15.30 usec
P4        30.60 usec
PCPD2    90.00 usec
PLM2     12.0000000 W
PLM12    0.34680000 W

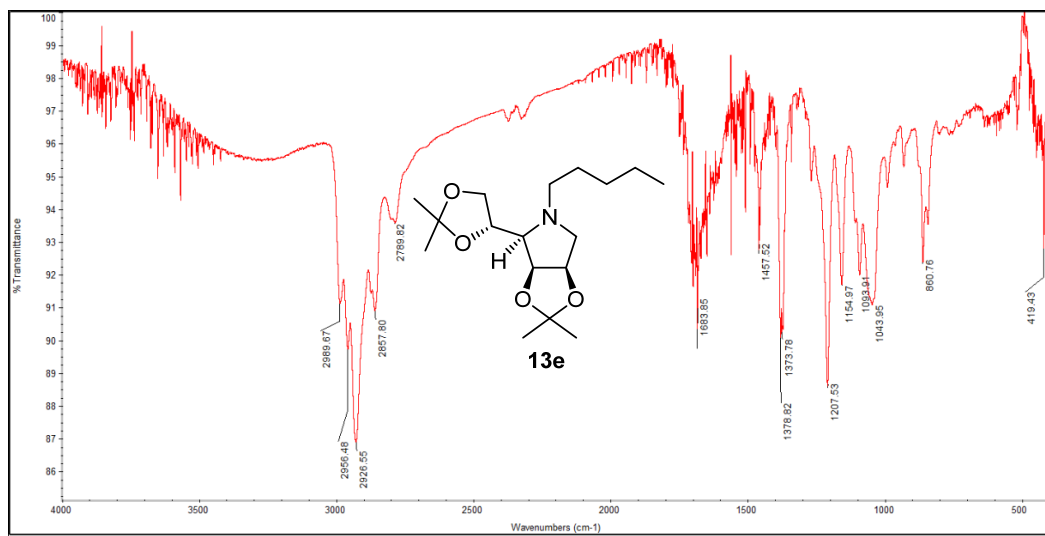
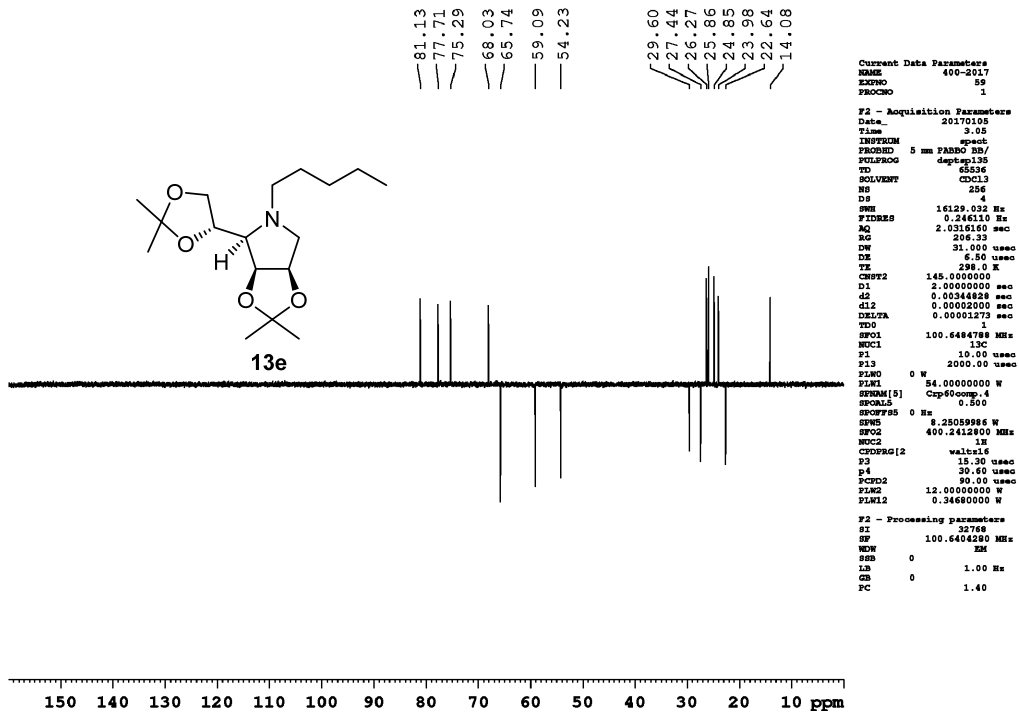
F2 - Processing parameters
SI        32768
SF        100.6404280 MHz
WDW       EM
SSB       0
LB        1.00 Hz
GB        0
PC        1.40

```

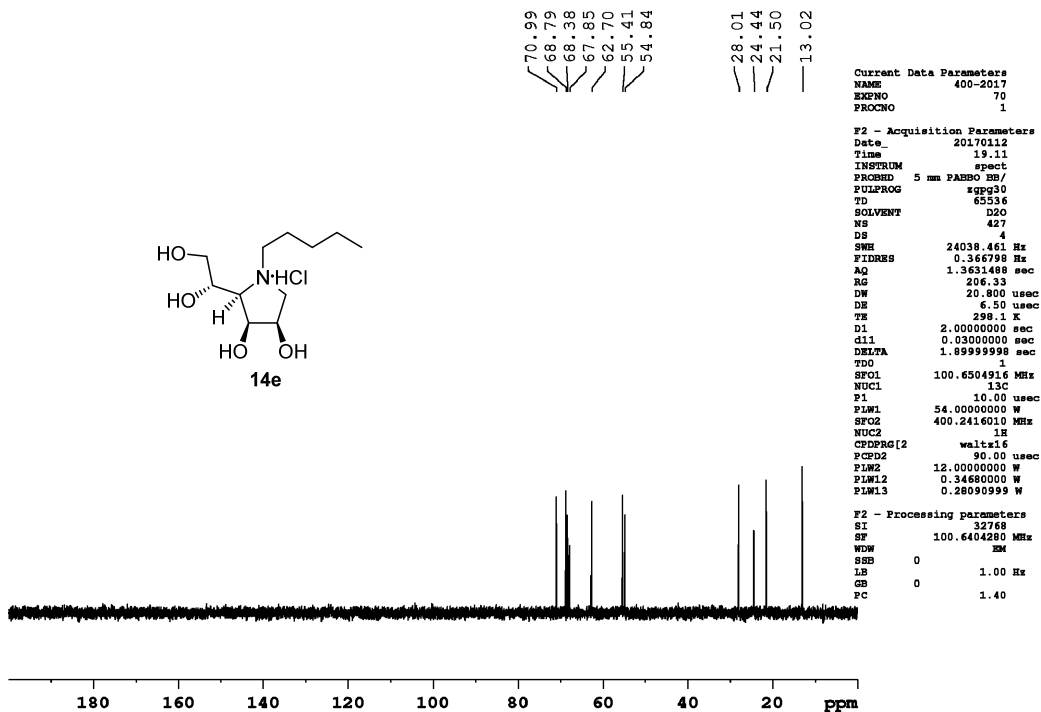
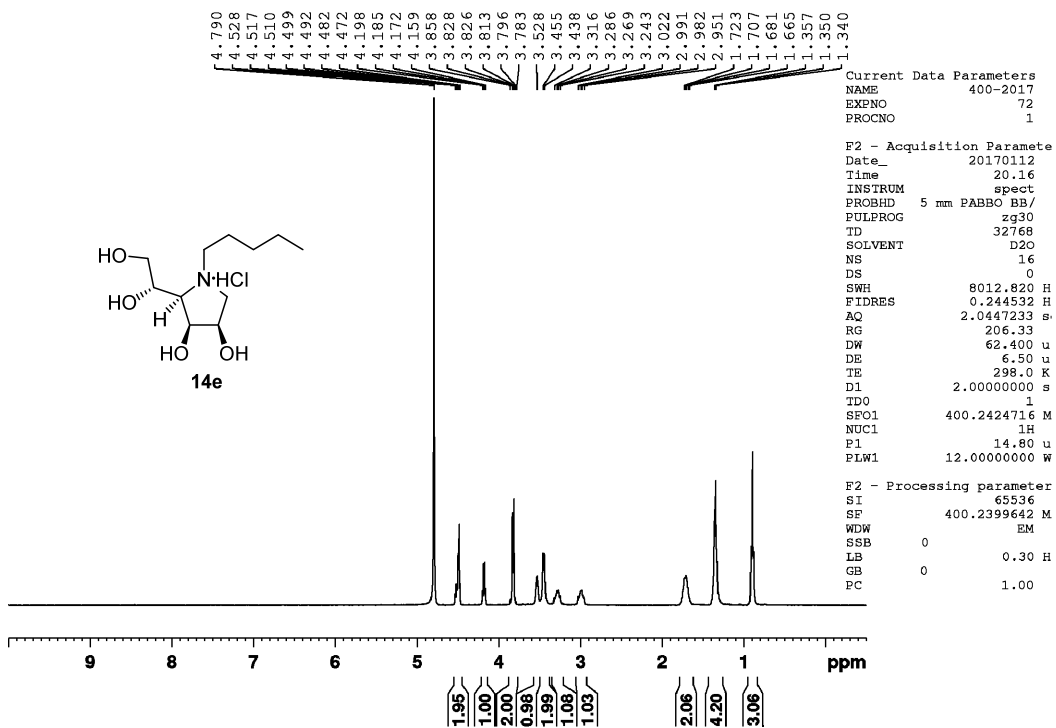


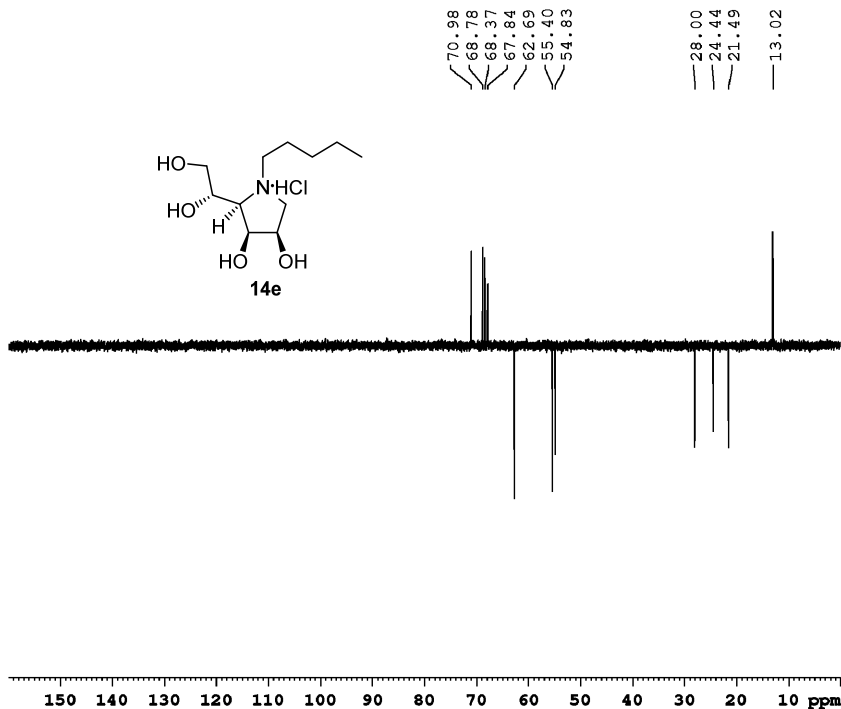
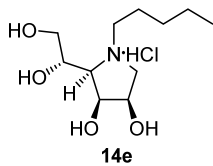
Compound 13e:





Compound 14e:



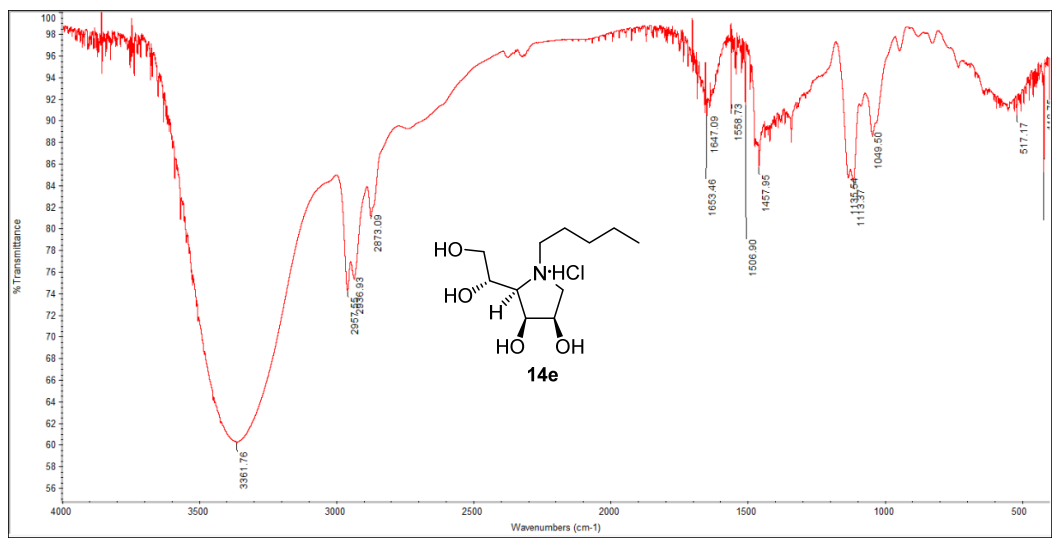


```

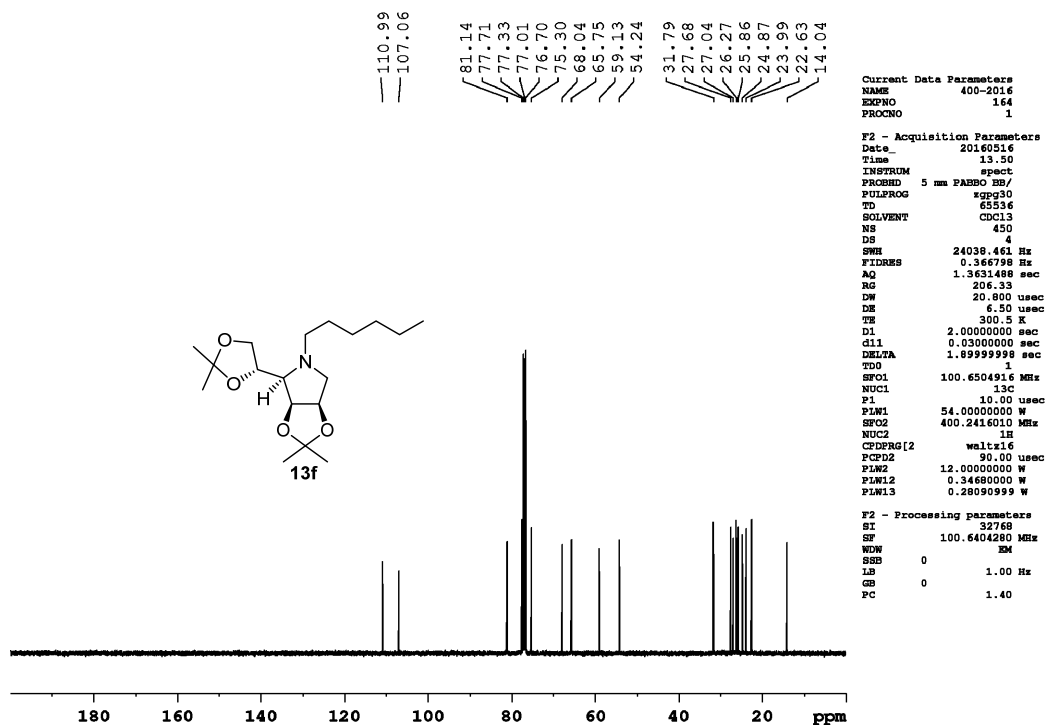
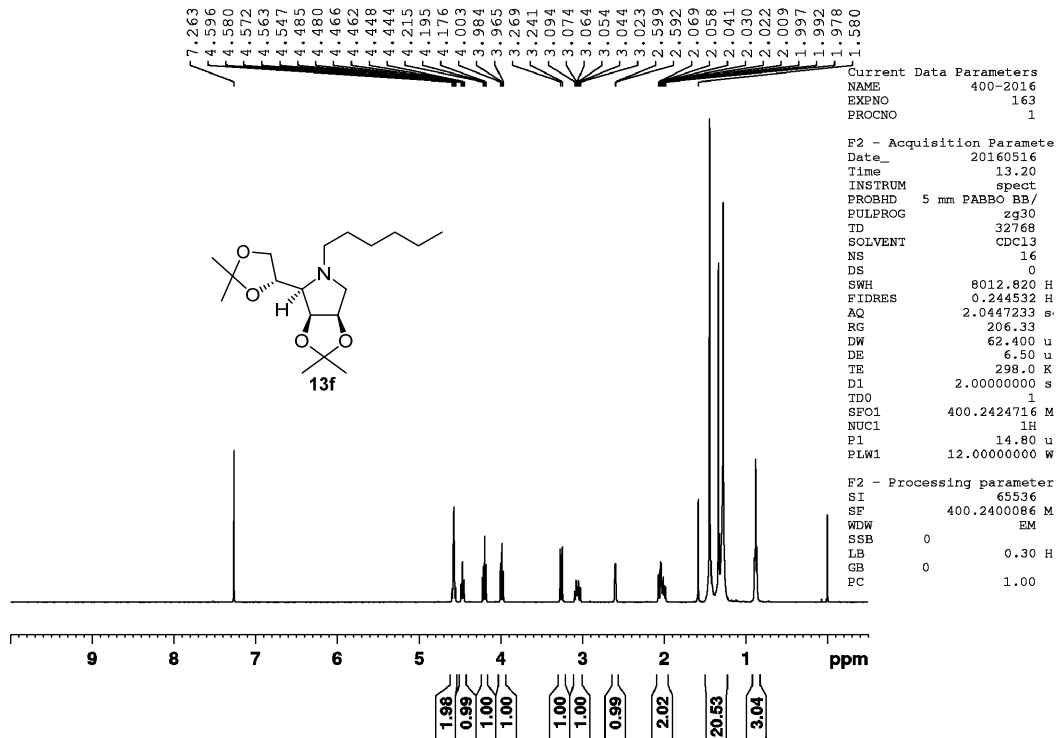
Current Data Parameters
NAME      400-2017
EXPNO    71
PROCNO   1

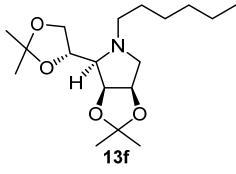
F2 - Acquisition Parameters
Date_    20170112
Time     19.26
INSTRUM  spect
PROBHD   5 mm PABBO BB/
PULPROG  zgpg30
TD       65536
SOLVENT  D2O
NS       157
DS       4
SWH      16129.932 Hz
FIDRES   0.246110 Hz
AQ       2.0316180 sec
RG       206.33
DN       31.000 usec
DE       6.50 usec
TE       298.0 K
CQF2     145.0000000
SI       2.00000000 sec
d2       0.00344628 sec
d12      0.00000000 sec
DELTA    0.0001273 sec
TDO      2
SFO1     100.6484788 MHz
NUC1      13C
SI       10.00 usec
P13      2000.00 usec
PLMO     0 W
PLM1     54.0000000 W
SFO1M[5] Crp60comp_4
SFO1M[5] 0.500
SFOFF[5] 0 Hz
SWS      8.25059986 W
SFO2     400.2412180 MHz
NUC2      1H
CHPRG[2] waltz16
P3       15.30 usec
p4       30.60 usec
PCPD2    90.00 usec
PLM2     12.0000000 W
PLM12    0.3468000 W

F2 - Processing parameters
SI       32768
SF       100.6404280 MHz
NS       65536
SFR      0
LB       1.00 Hz
GB       0
PC       1.40
  
```



Compound 13f:





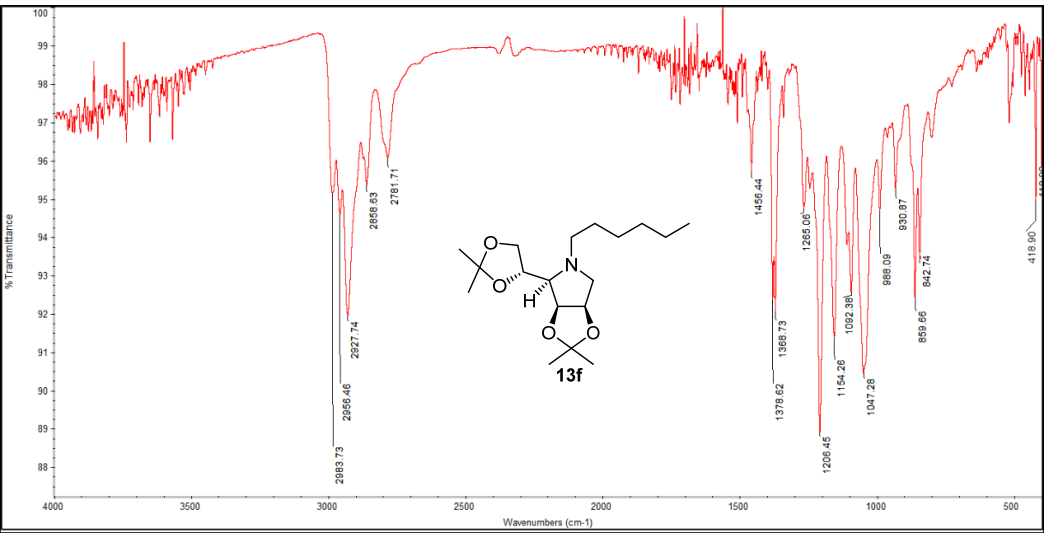
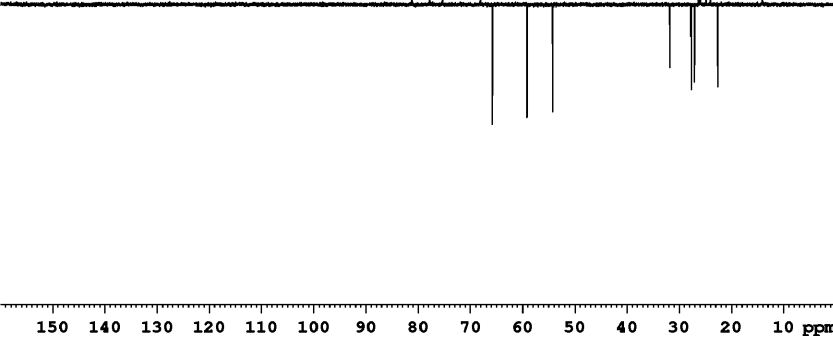
81.14
77.71
75.30
68.04
65.75
59.13
54.24
31.79
27.68
27.05
26.27
25.86
24.87
23.99
22.63
14.04

```

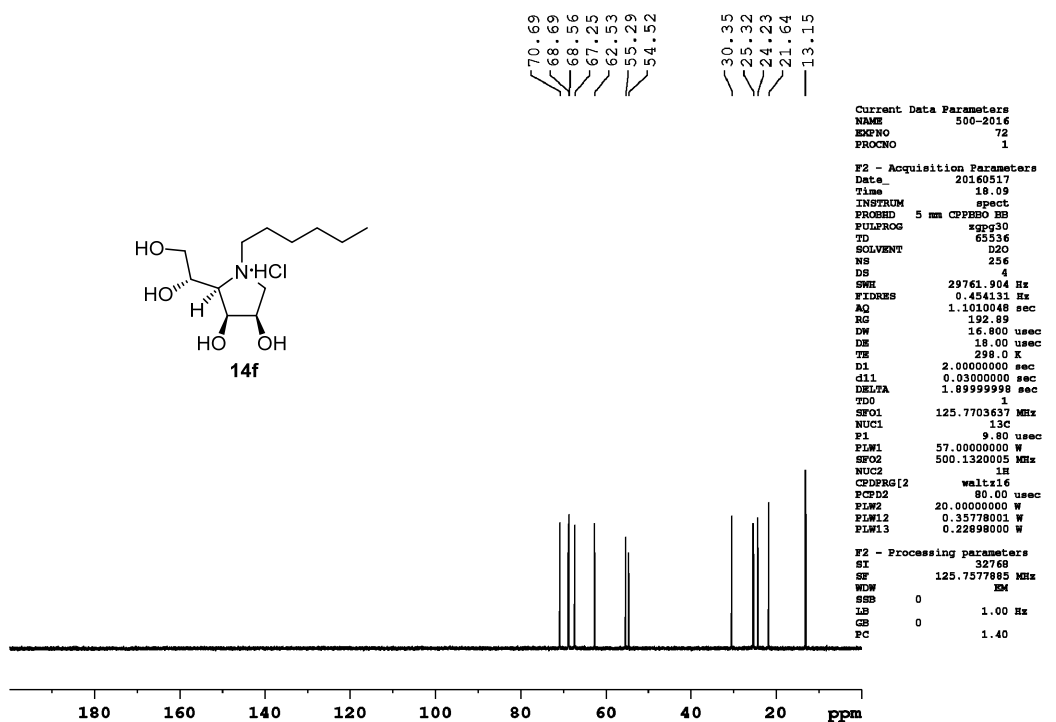
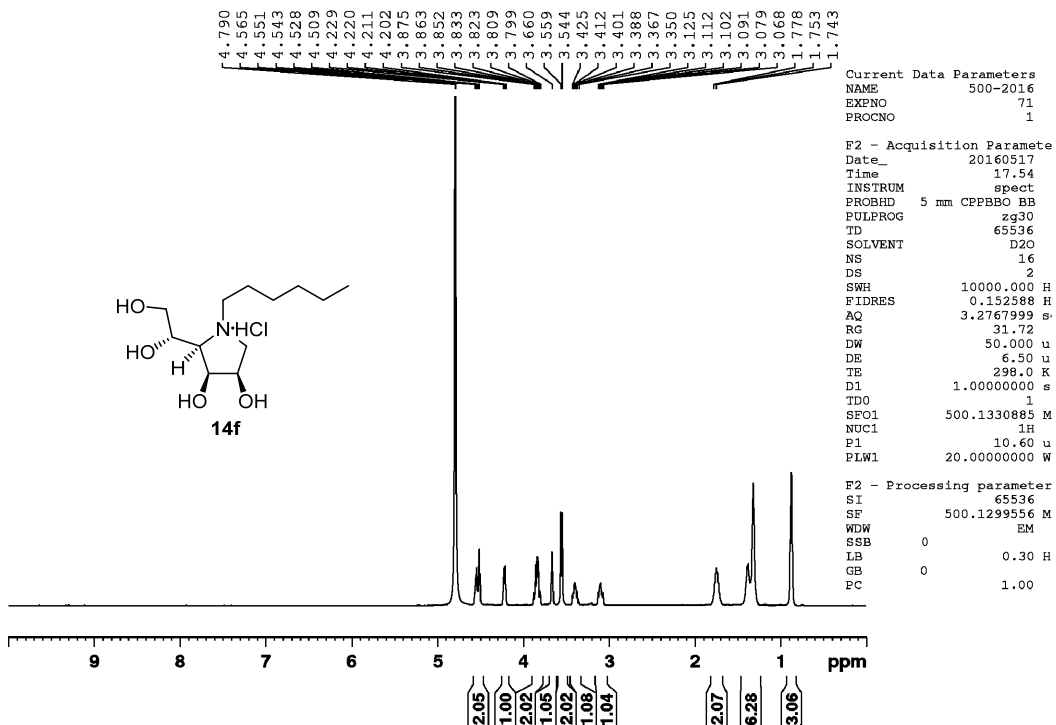
Current Data Parameters
NAME      400-2016
EXPNO     165
PROCNO    1

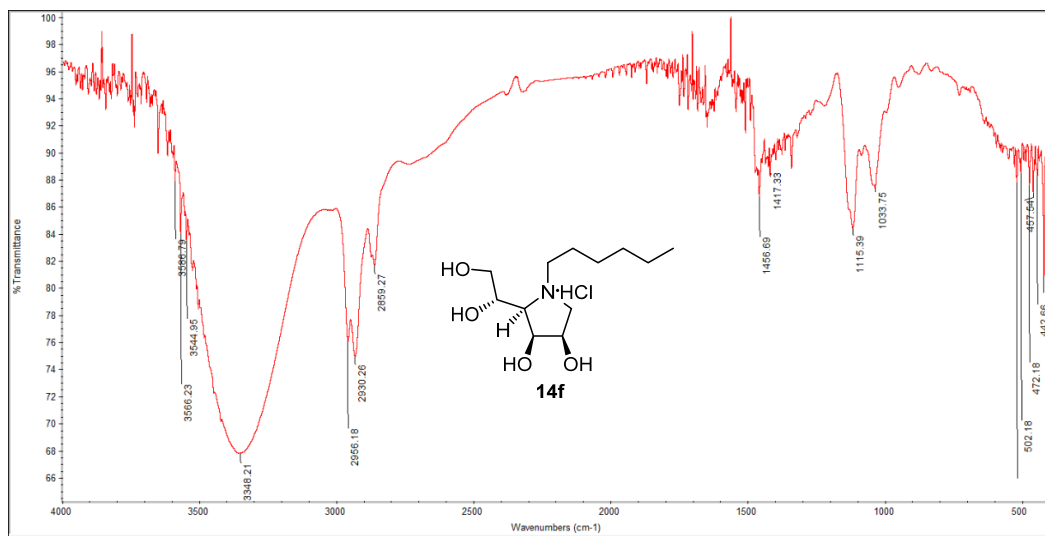
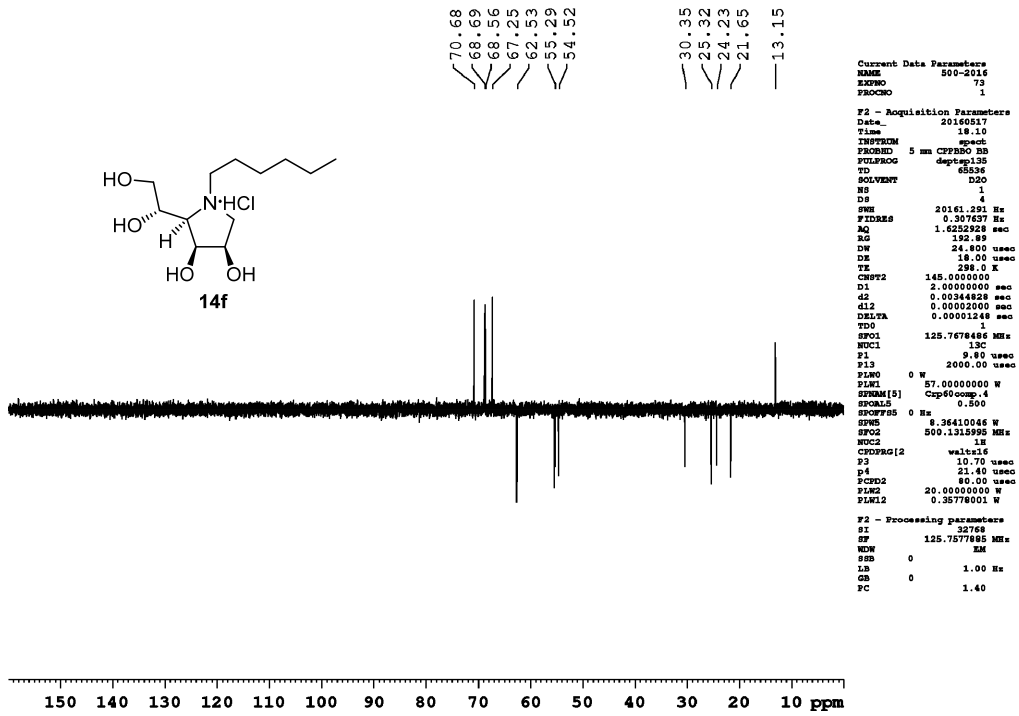
F2 - Acquisition Parameters
Date_     20160316
Time      14.01
INSTRUM   spect
PROBHD    5 mm PABBO BB/
PULPROG   zgpg30
VD        6536
SOLVENT   CDCl3
NS        150
DS        4
SWH       16129.932 Hz
FIDRES    0.246110 Hz
AQ        2.0216180 sec
RG        206.33
DW        31.000 usec
DE        6.50 usec
TE        300.3 K
CQF2      145.000000
D1        2.0000000 sec
d2        0.00344828 sec
d12       0.00000000 sec
DELTA    0.00001273 sec
YD0       1
SFO1      100.646478 MHz
NUC1      13C
P1        13.00 usec
F13       2000.00 usec
PLMO      0 W
PLM1      54.0000000 W
SFO2      Csp60comp.4
SFOALS    0.500
SFOFFAS   0 Hz
SWS       8.25099986 W
SFC2      400.2412800 MHz
MUC2      1H
CHOPROG2  waltz16
V3        15.30 usec
p4        30.60 usec
NCPD2     90.00 usec
PLM2      12.0000000 W
PLML2     0.34680000 W

F2 - Processing Parameters
SI        32768
SF        100.6404280 MHz
WDW       EM
SSB       0
LB        1.00 Hz
GB        0
PC        1.40
  
```

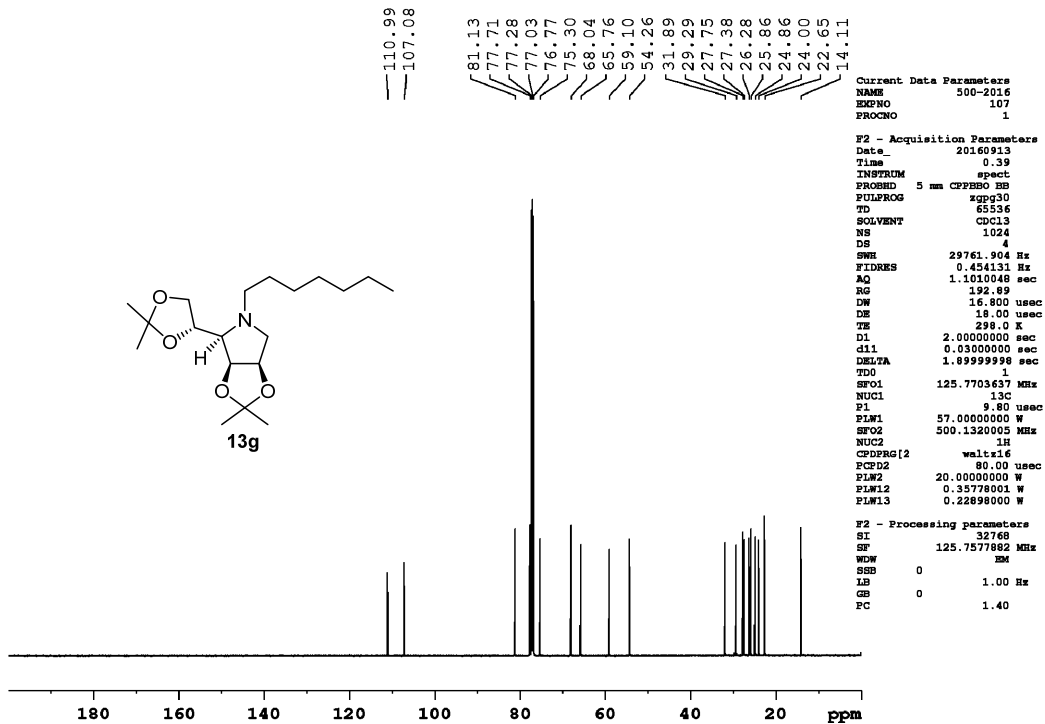
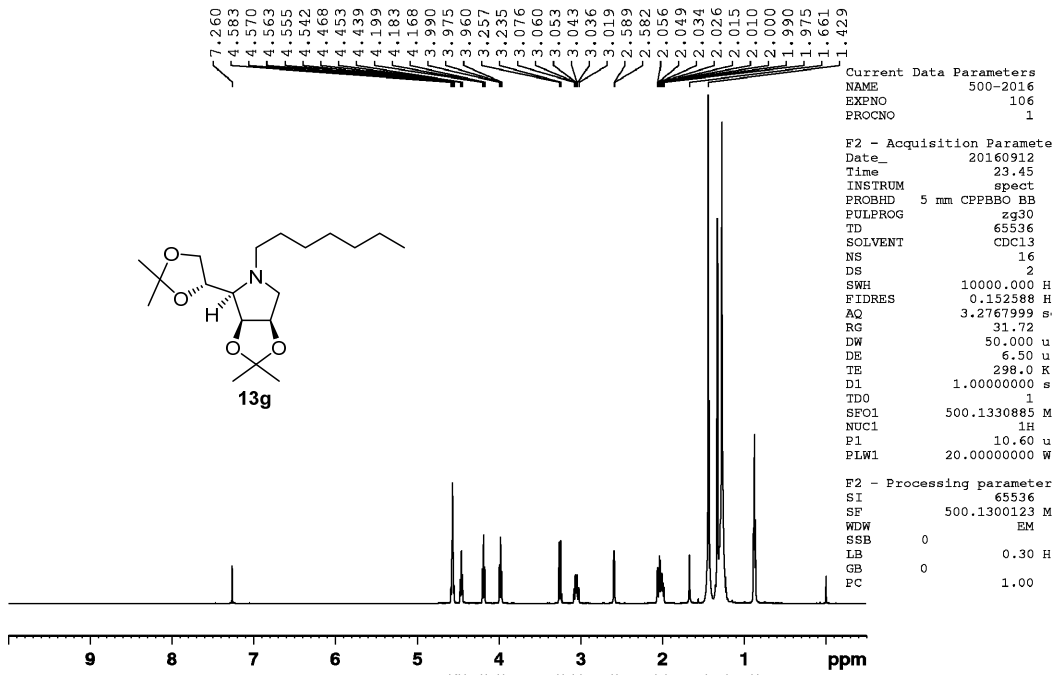


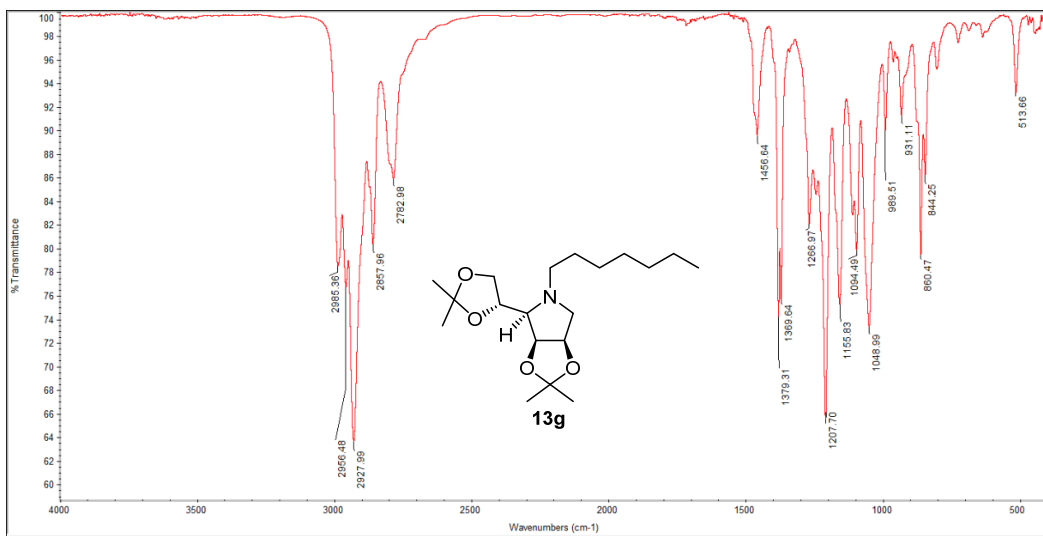
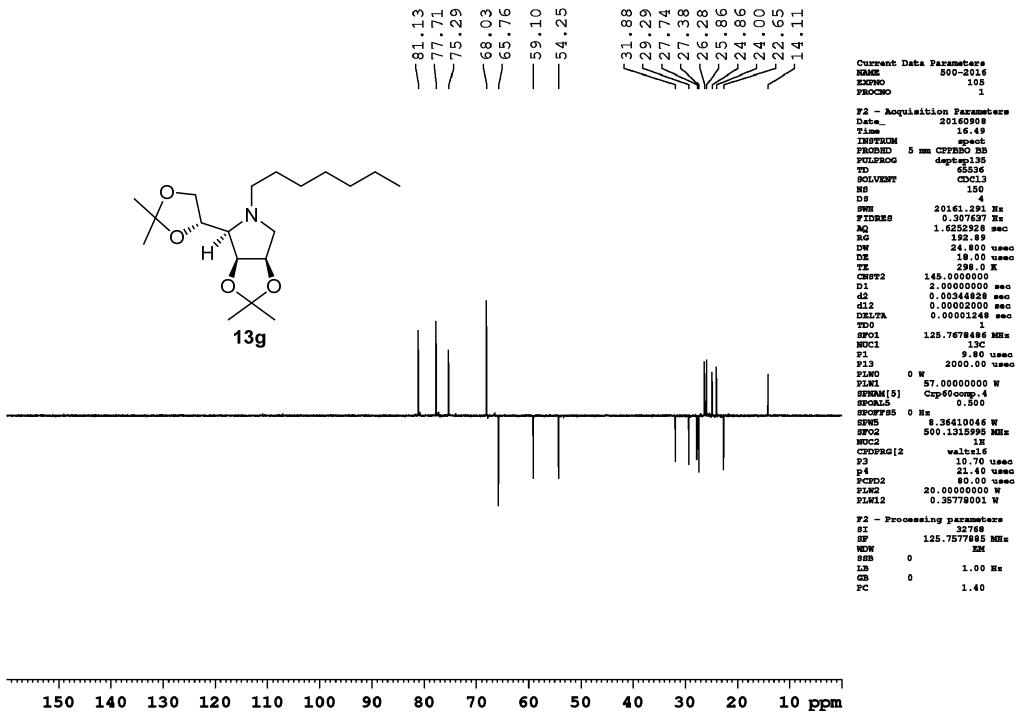
Compound 14f:



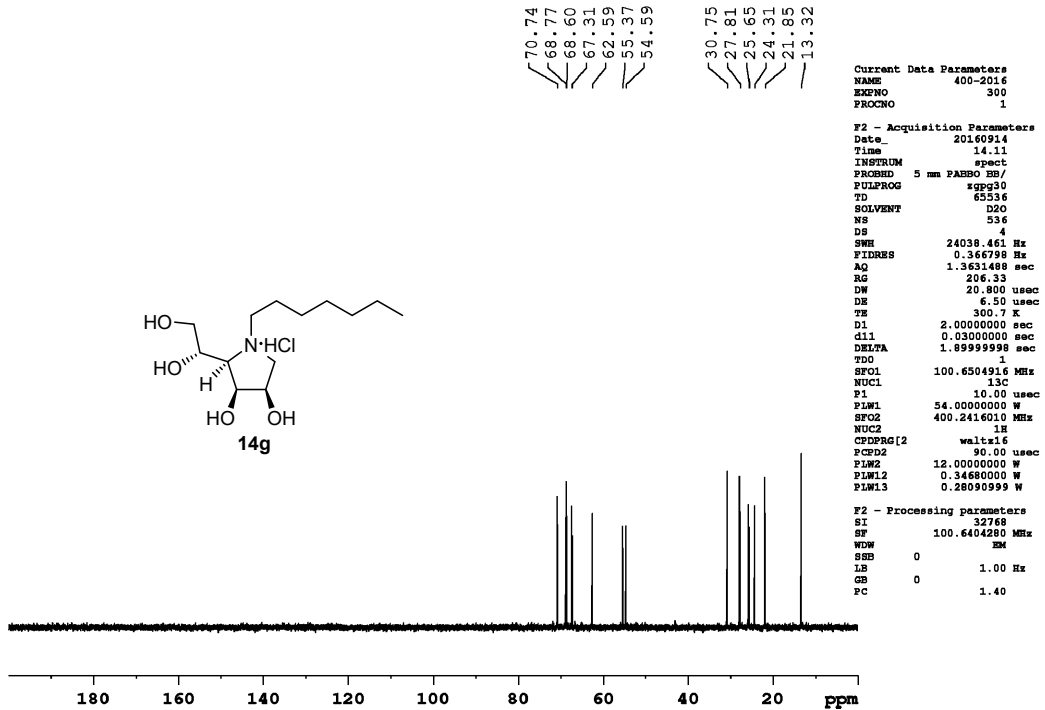
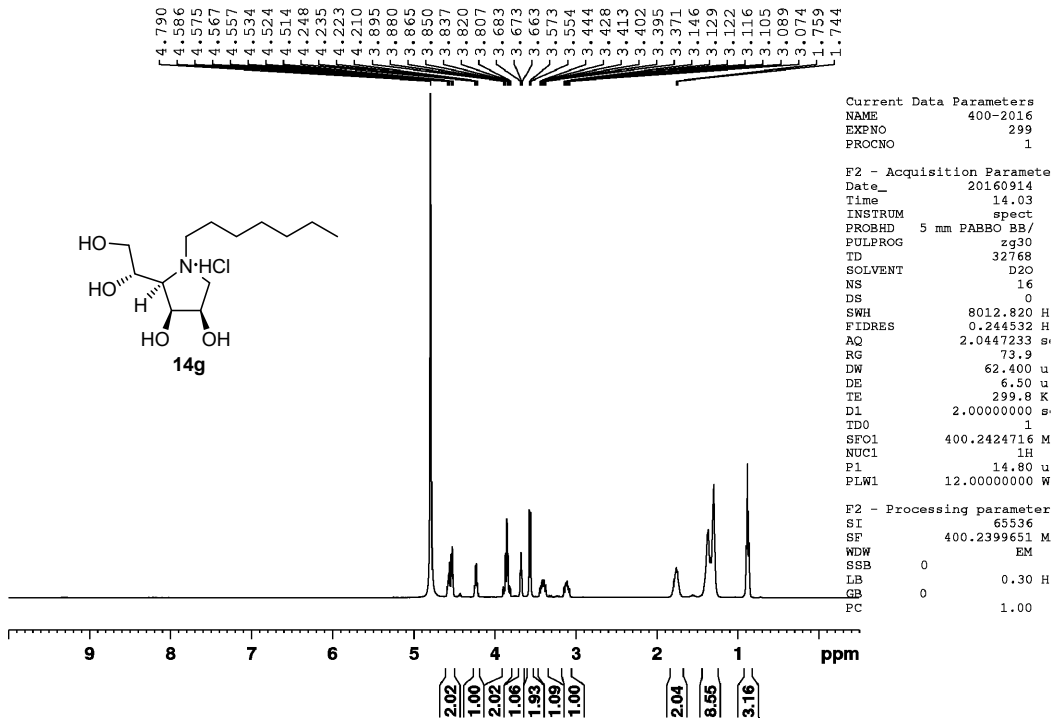


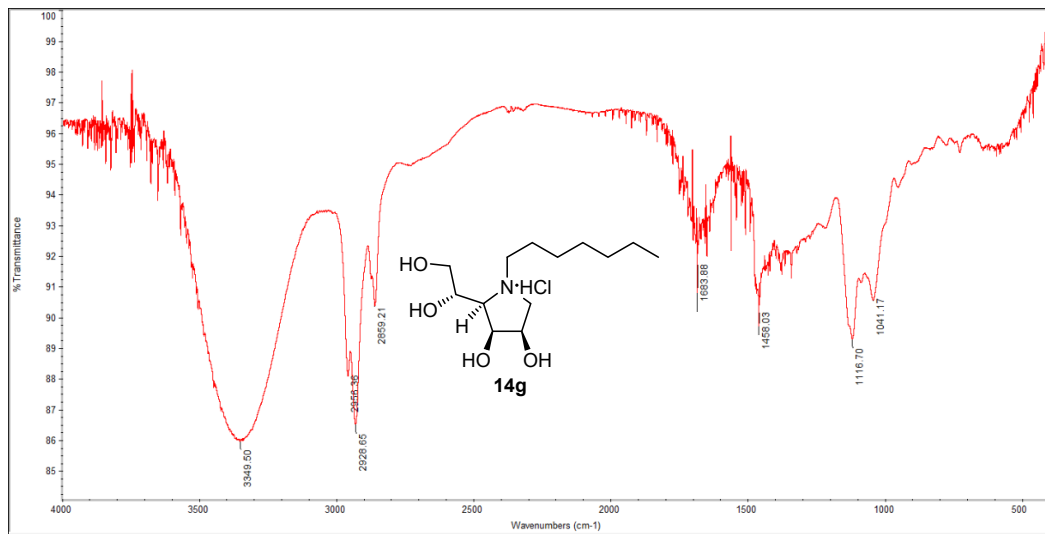
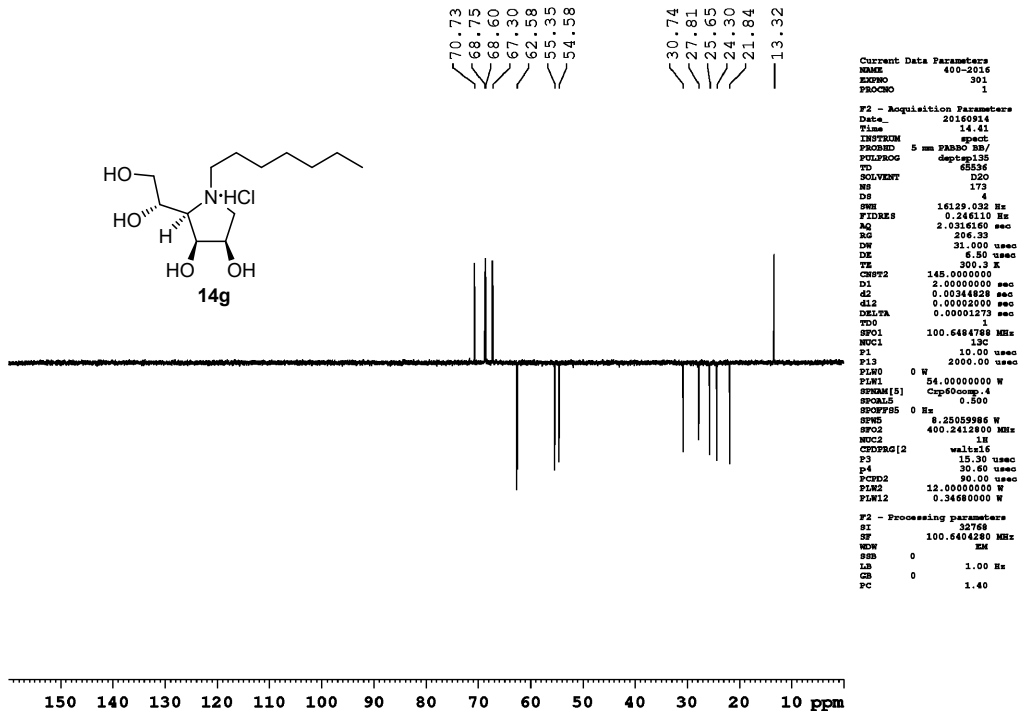
Compound 13g:



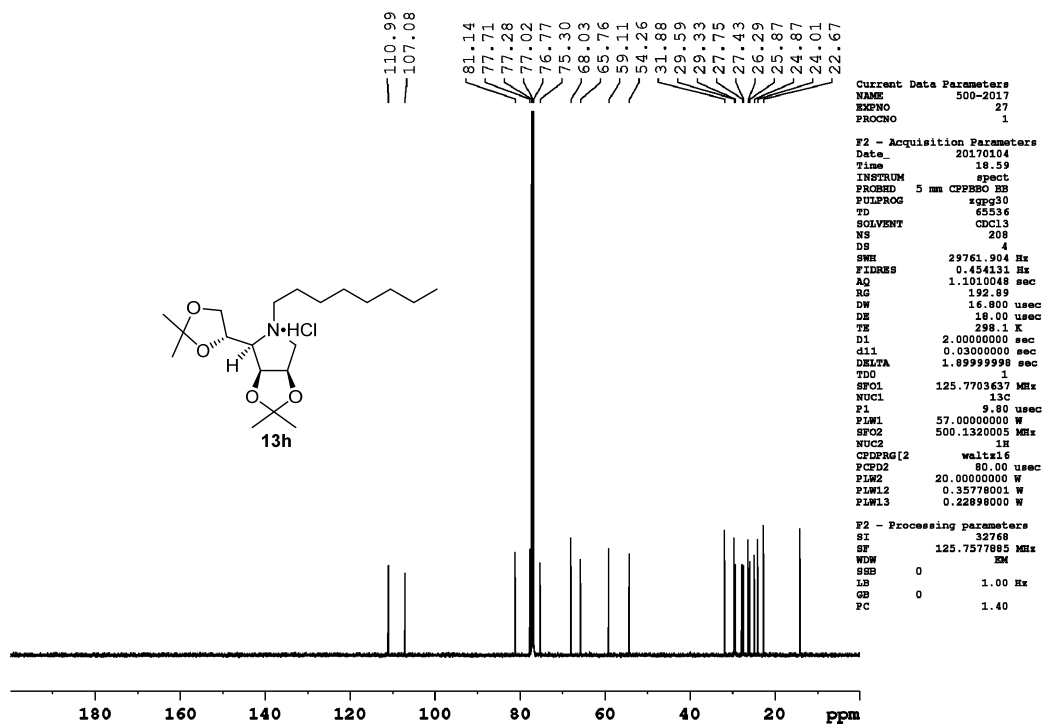
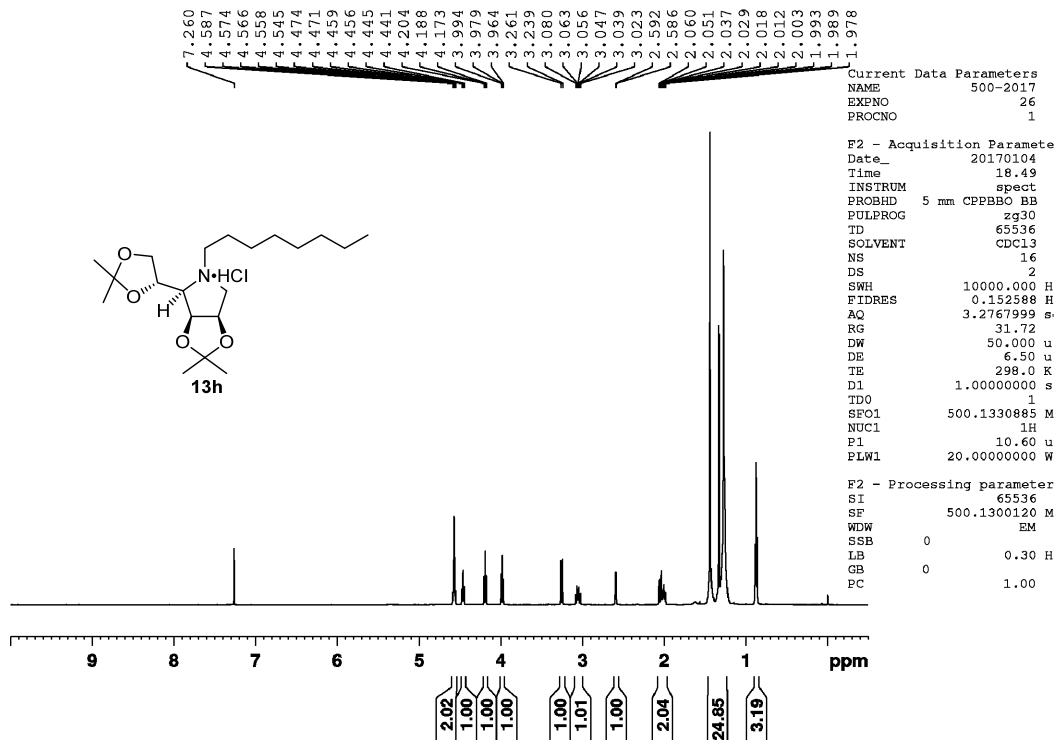


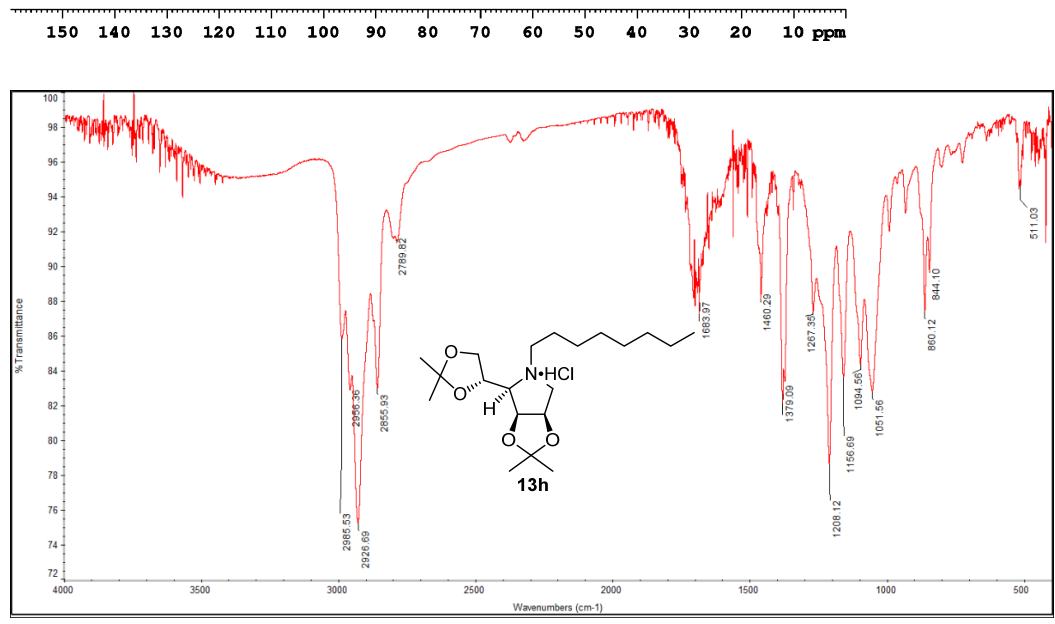
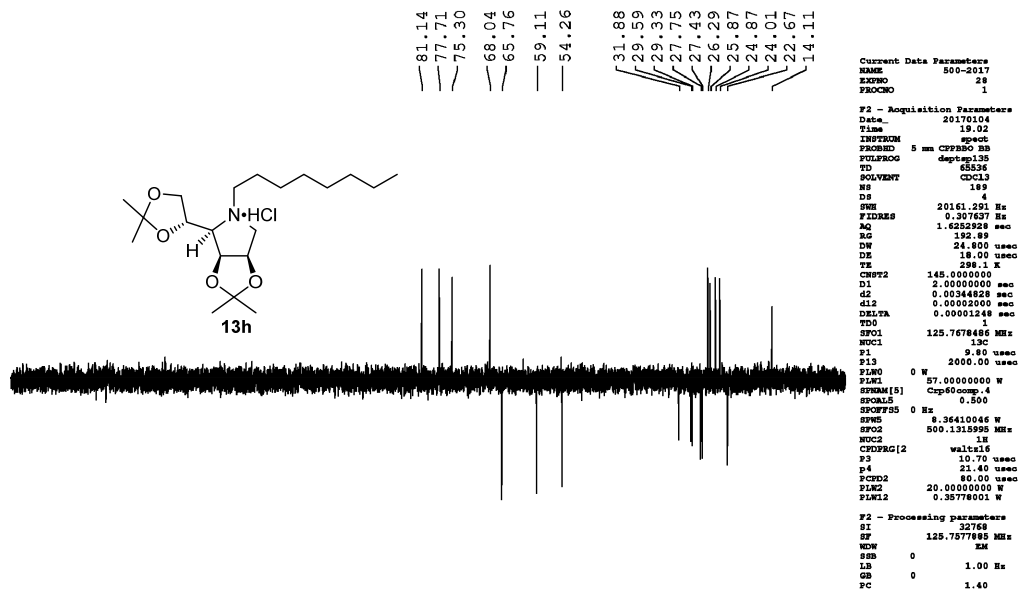
Compound 14g:



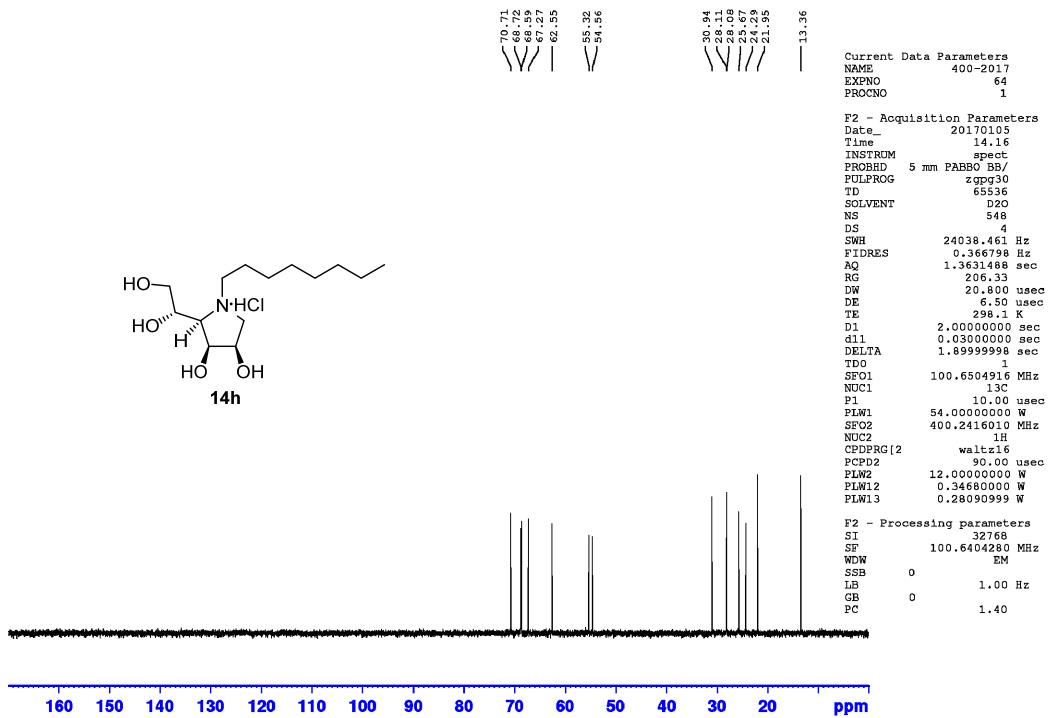
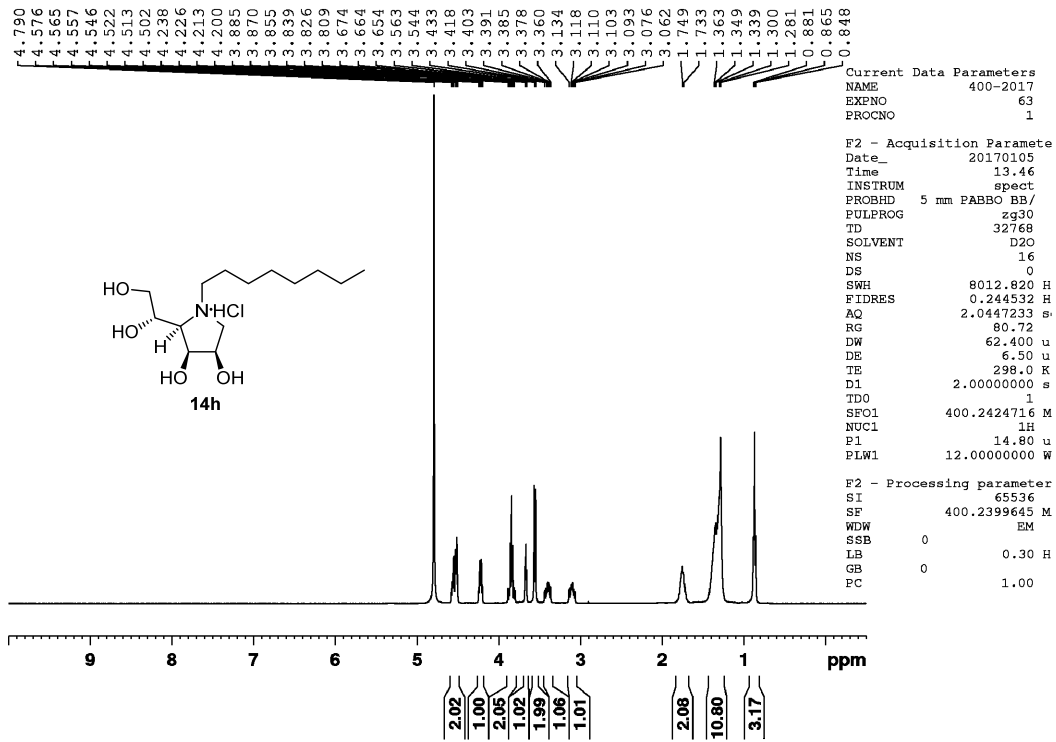


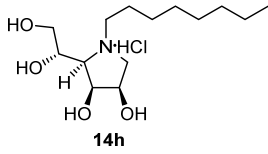
Compound 13h:





Compound 14h:





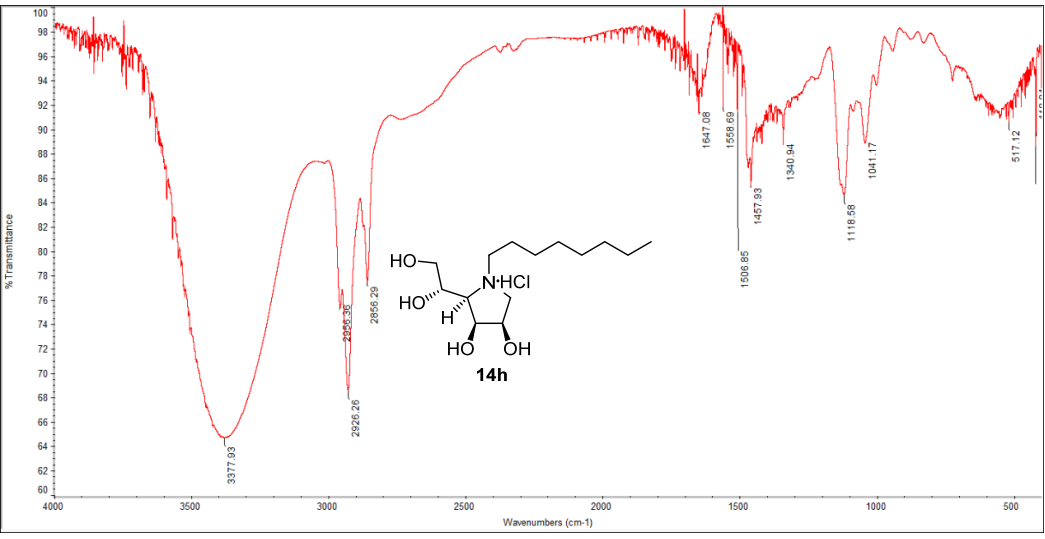
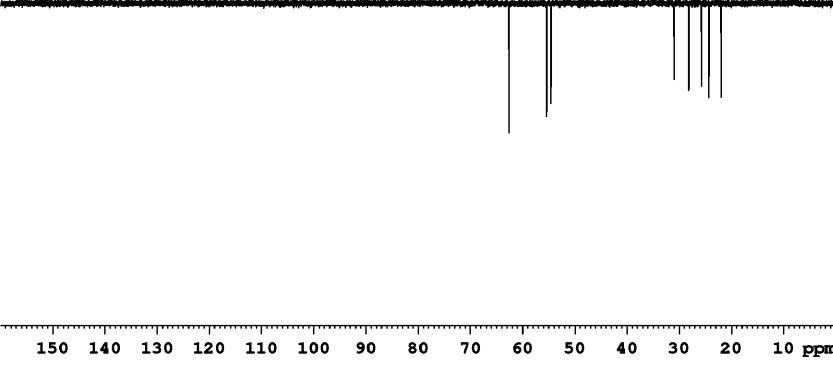
70.71
68.71
68.58
67.27
62.54
55.31
54.55
30.93
28.11
28.08
25.67
24.28
21.94
13.35

```

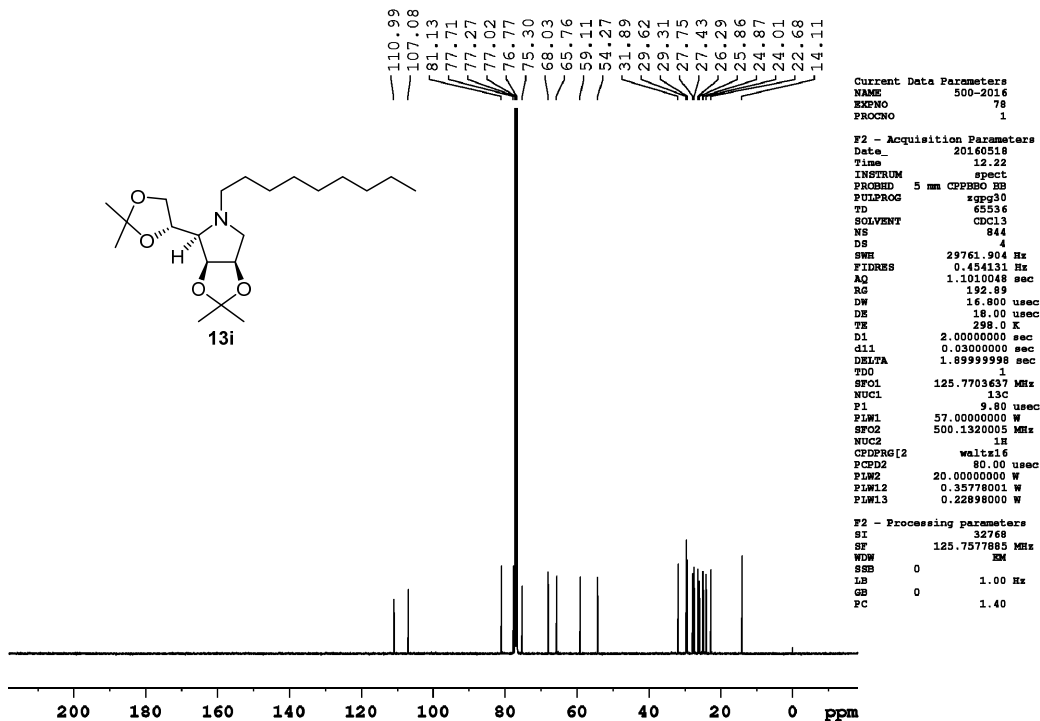
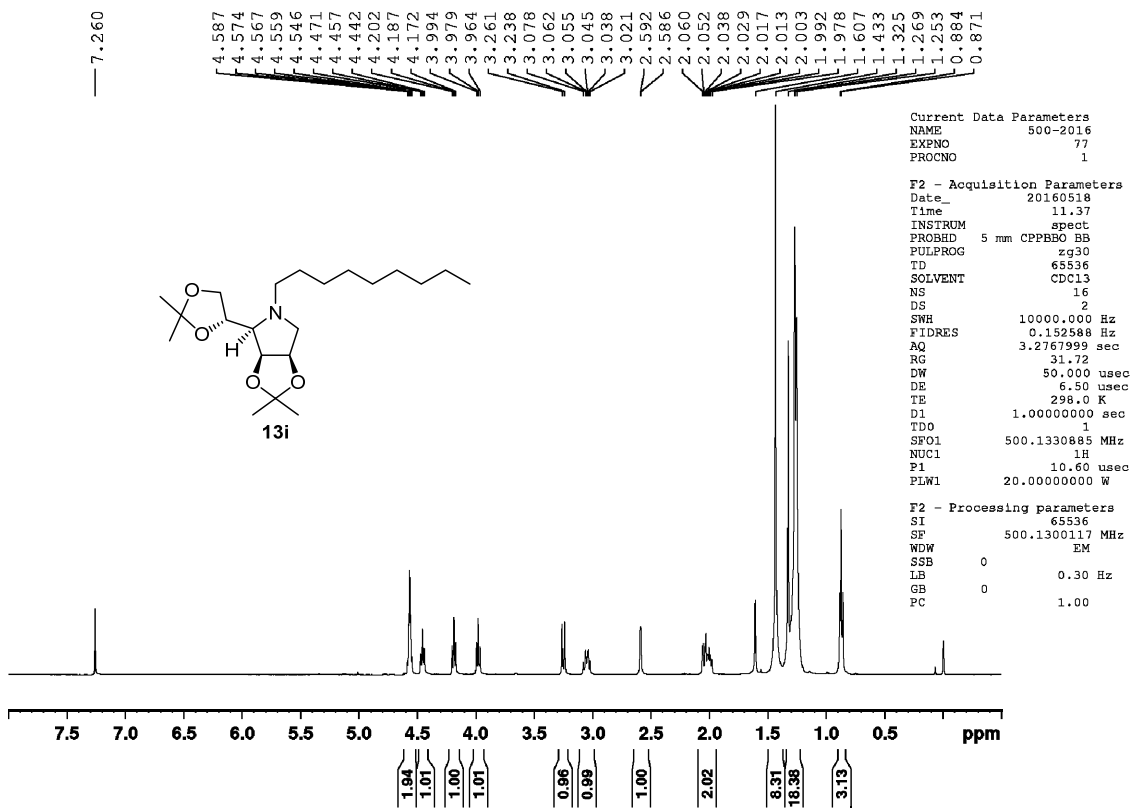
Current Data Parameters
NAME      400-2017
EXPNO     65
PROCNO    1

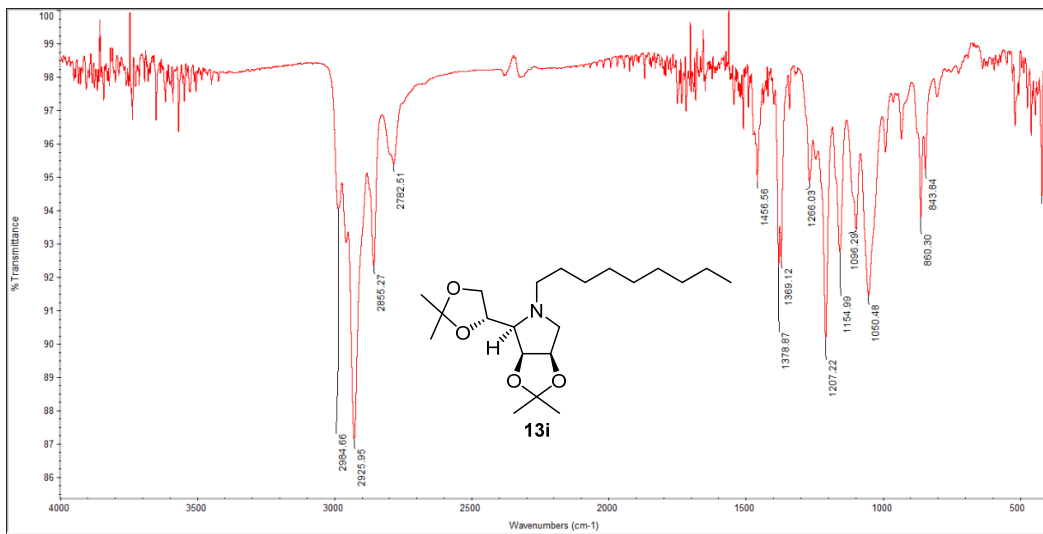
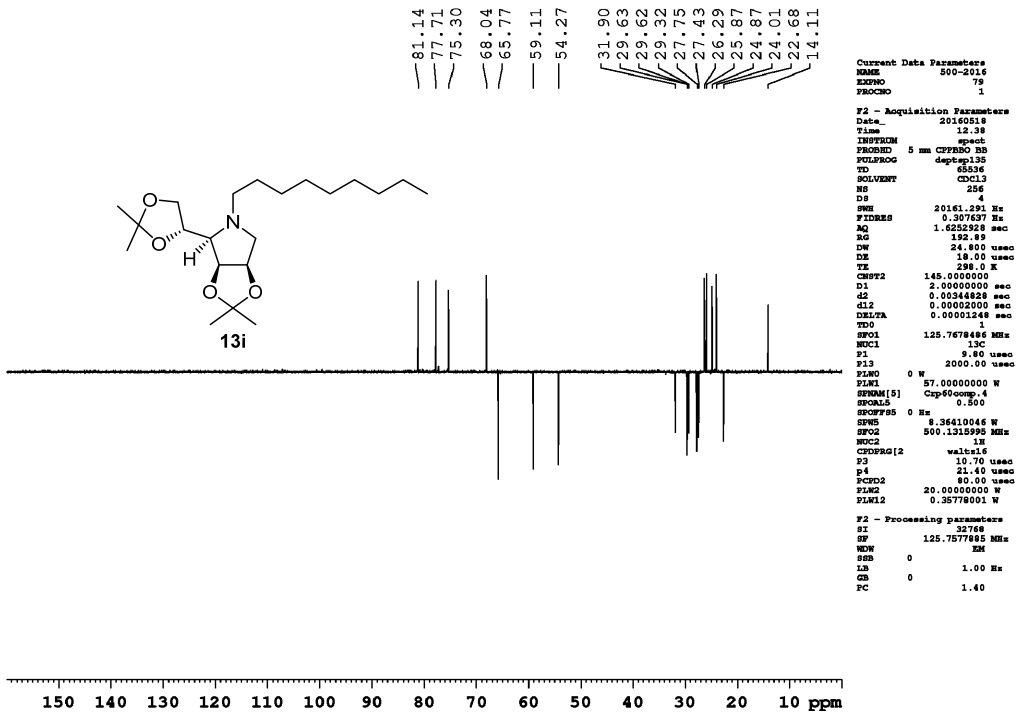
F2 - Acquisition Parameters
Date_     20170105
Time      14.23
INSTRUM   spect
PROBHD    5 mm PABBO BB/
PULPROG   zgpg30
TD         65536
SOLVENT   D2O
NS         148
DS         4
SWH        16129.932 Hz
FIDRES    0.246110 Hz
AQ         2.0216180 sec
RG         206.39
DW         31.000 usec
DE         6.50 usec
TE         298.0 K
CQF2      145.000000
D1         2.0000000 sec
d2         0.00344828 sec
d12        0.00002000 sec
DELTA     0.00001273 sec
VFO        1
NUC1       100.6464788 MHz
NUC2       13C
P1         10.00 usec
P13        2000.00 usec
PL1        0 W
PL12       54.0000000 W
SFOFF(5)  Csp60comp.4
SFOFF(5)  0.500
SFOFF(5)  0 Hz
SFRS       8.25059986 W
SFOFF      400.2412800 MHz
NUC22      1H
CHOPROG[2] waltz16
V3         15.30 usec
p4         30.60 usec
SFOFF2     90.00 usec
PLM2       12.0000000 W
PLML2      0.34680000 W

F2 - Processing parameters
SI         32768
SF         100.6404280 MHz
WDW        EM
SSB         0
LB         1.00 Hz
GB         0
PC         1.40
  
```

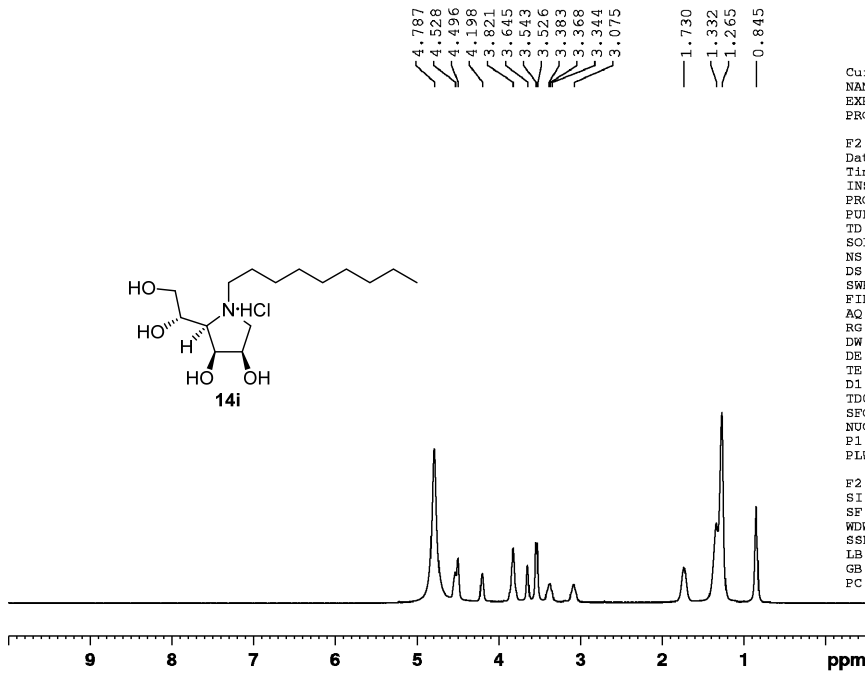


Compound 13i:





Compound 14i:

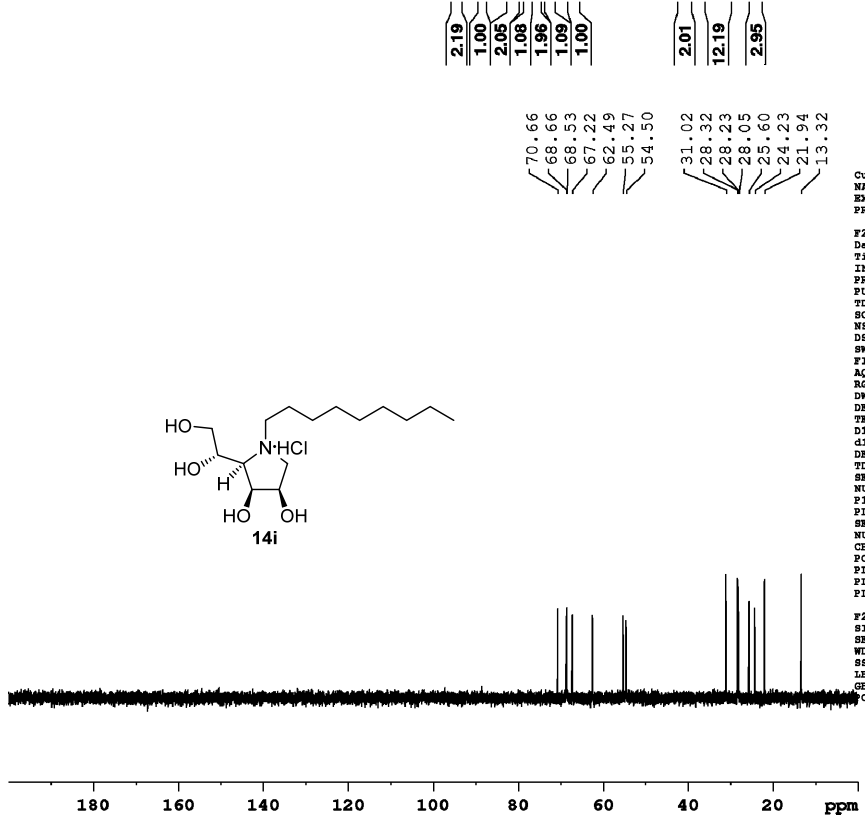


```

Current Data Parameters
NAME      400-2016
EXPNO    173
PROCNO   1

F2 - Acquisition Parameters
Date_    20160519
Time     13.21
INSTRUM spect
PROBHD   5 mm PABBO BB/
PULPROG zg30
TD       32768
SOLVENT  D2O
NS       16
DS       0
SWH      8012.820 Hz
FIDRES   0.244532 Hz
AQ       2.0447233 sec
RG       162.77
DW       62.400 usec
DE       6.50 usec
TE       298.1 K
D1       2.0000000 sec
TDO      1
SFO1    400.2424716 MHz
NUC1     1H
P1       14.80 usec
PLW1    12.00000000 W

F2 - Processing parameters
SI       65536
SF       400.2399501 MHz
WDW      EM
SSB      0
LB       0.30 Hz
GB       0
PC       1.00
    
```

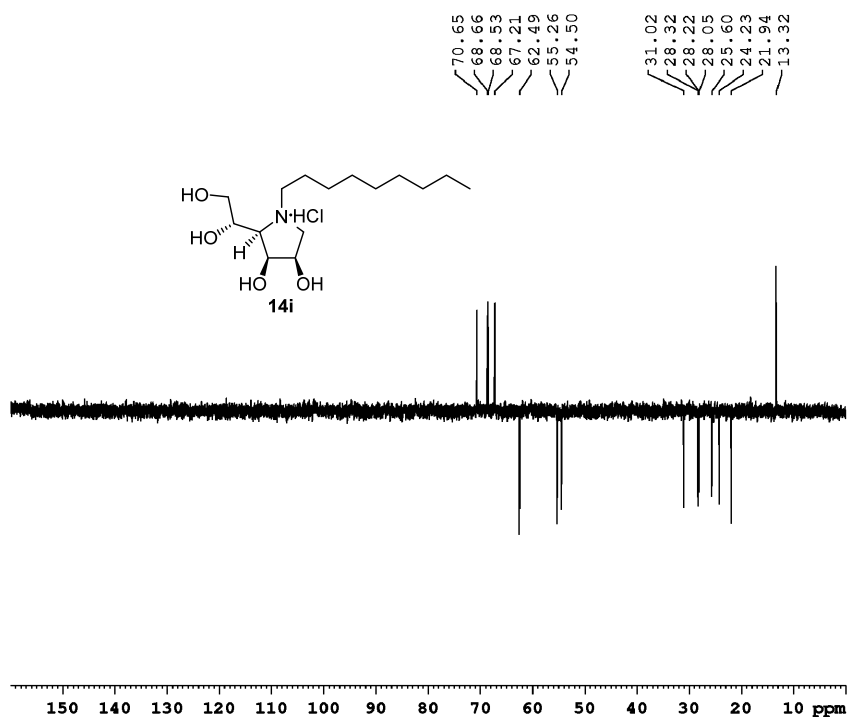


```

Current Data Parameters
NAME      400-2016
EXPNO    174
PROCNO   1

F2 - Acquisition Parameters
Date_    20160519
Time     13.29
INSTRUM spect
PROBHD   5 mm PABBO BB/
PULPROG zgpg30
TD       65536
SOLVENT  D2O
NS       600
DS       4
SWH      24038.481 Hz
FIDRES   0.366798 Hz
AQ       1.3631488 sec
RG       206.33
DW       20.800 usec
DE       6.50 usec
TE       298.2 K
D1       2.0000000 sec
d11      0.03000000 sec
DELTA    1.89999998 sec
TDO      1
SFO1    100.6504916 MHz
NUC1     13C
P1       10.00 usec
PLW1    54.00000000 W
SFO2    400.2416010 MHz
NUC2     1H
CPDPRG2 waltz16
PCPD2   90.00 usec
PLW2    12.00000000 W
PLW12   0.34680000 W
PLW13   0.28090999 W

F2 - Processing parameters
SI       32768
SF       100.6404280 MHz
WDW      EM
SSB      0
LB       1.00 Hz
GB       0
PC       1.40
    
```

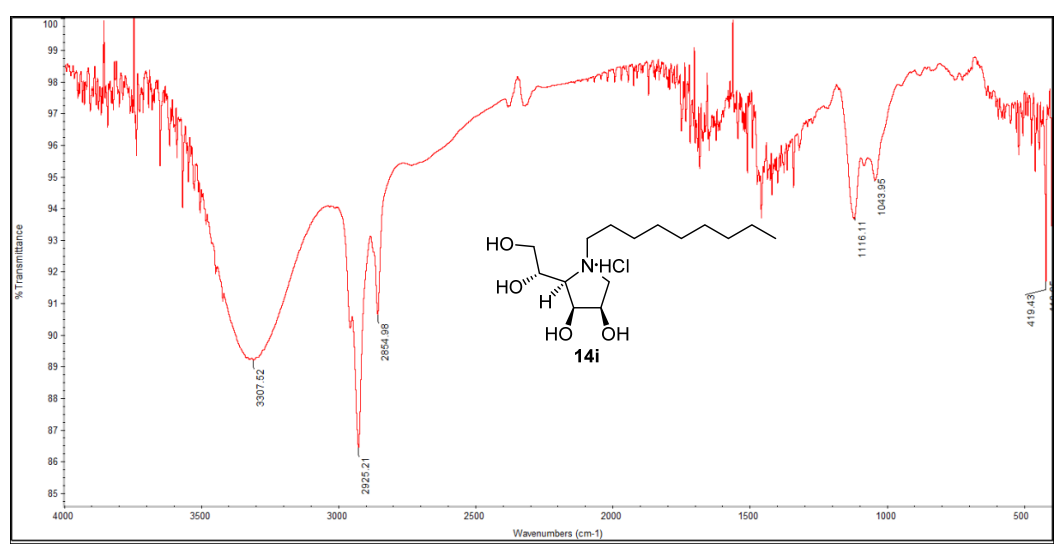



```

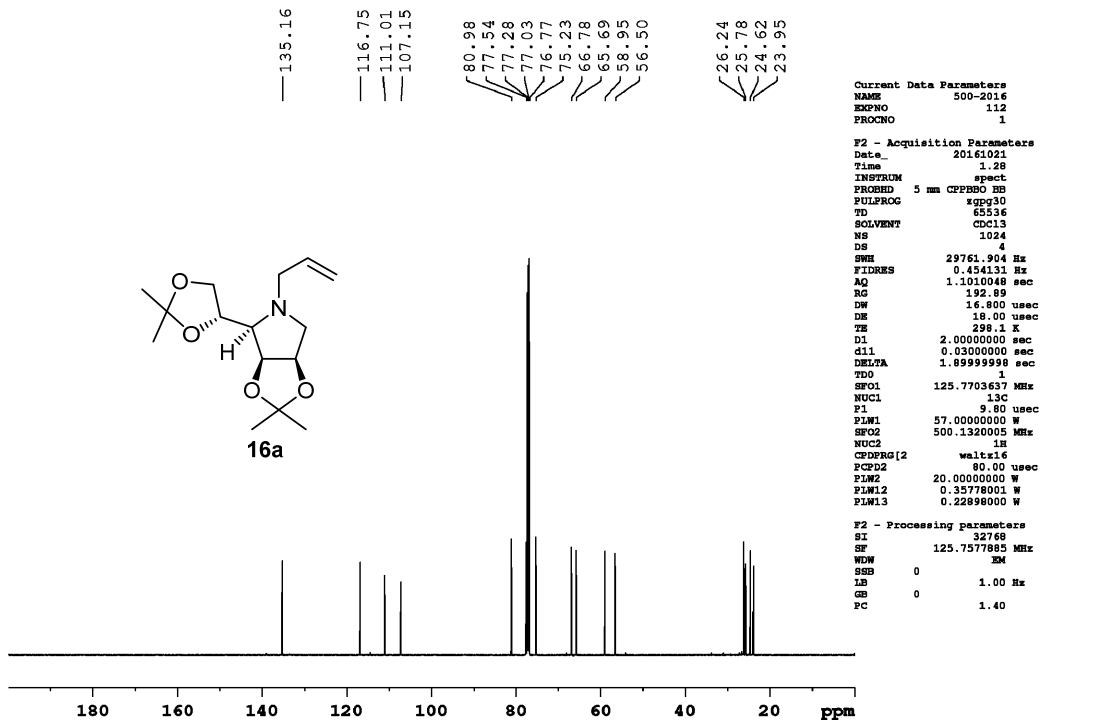
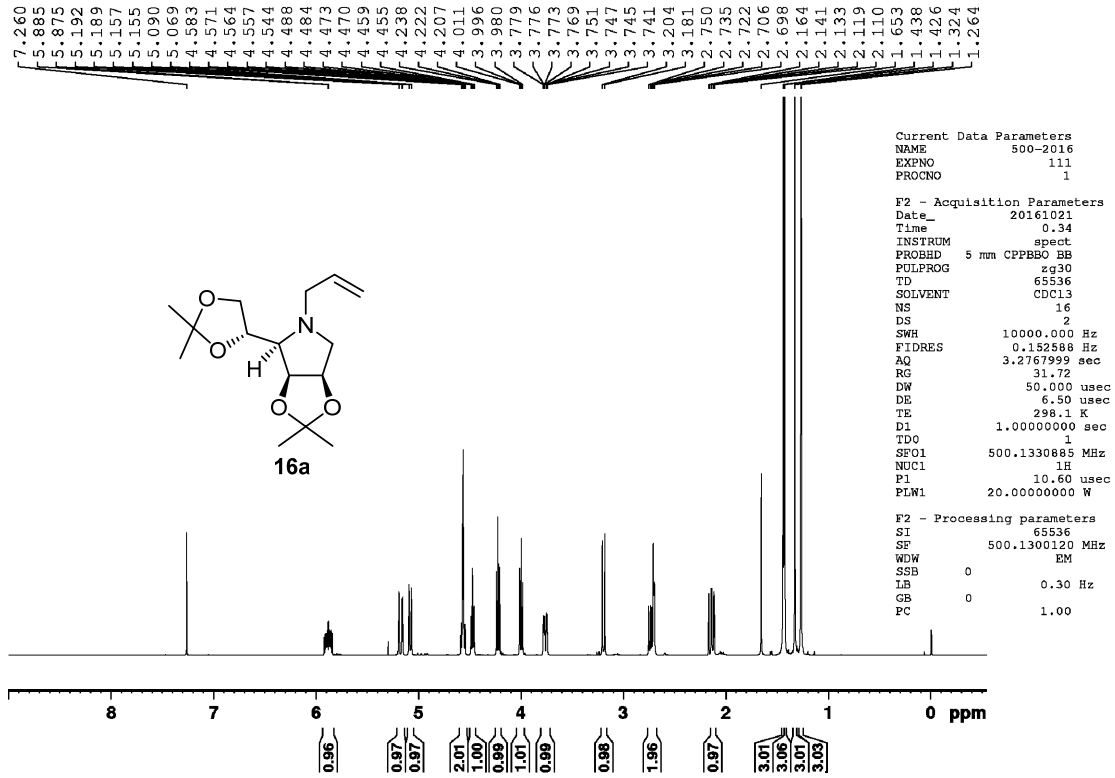
Current Data Parameters
NAME      400-2014
EXPNO     175
PROCNO    1

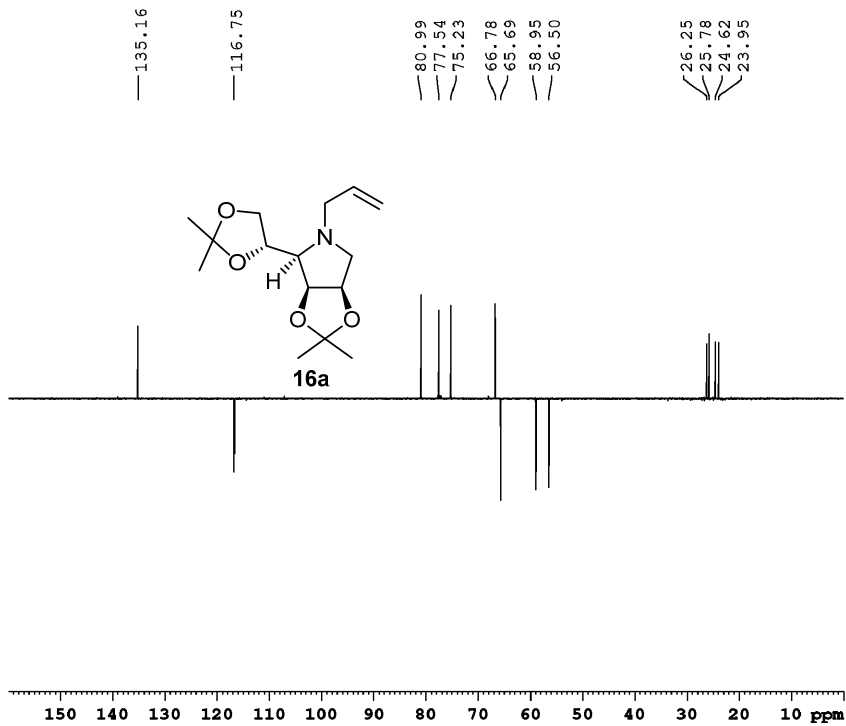
F2 - Acquisition Parameters
Date_     20160819
Time      14.08
INSTRUM   spect
PROBHD    5 mm PABBO BB/
PULPROG   zgpg30
TD         65536
SOLVENT   D2O
NS         150
DS         4
SWH        16129.832 Hz
FIDRES     0.246110 Hz
AQ         2.0216180 sec
RG         206.39
DW         31.000 usec
DE         6.50 usec
TE         298.3 K
CQF2       145.000000
SI         2.0000000 sec
d2         0.00344828 sec
d12        0.00000000 sec
DELTA      0.0001273 sec
WDW        2
SFO1       100.6484788 MHz
NUC1        13C
P1         10.00 usec
E13         2000.00 usec
PLMO       0 W
PLM1       54.0000000 W
SFOALS      Csp60comp.4
SFOALS      0.500
SFOFFS0     0 Hz
SWS        8.25059988 W
SFO2       400.2612800 MHz
NUC2        1H
CQFPG2[2]  waltz16
P2         15.30 usec
P4         30.60 usec
SFO22      50.00 usec
PLM2       12.0000000 W
PLML2      0.34680000 W

F2 - Processing parameters
SI         32768
SF         100.6404280 MHz
WDW        EM
SSB         0
LB         1.00 Hz
GB         0
PC         1.40
  
```



Compound 16a:



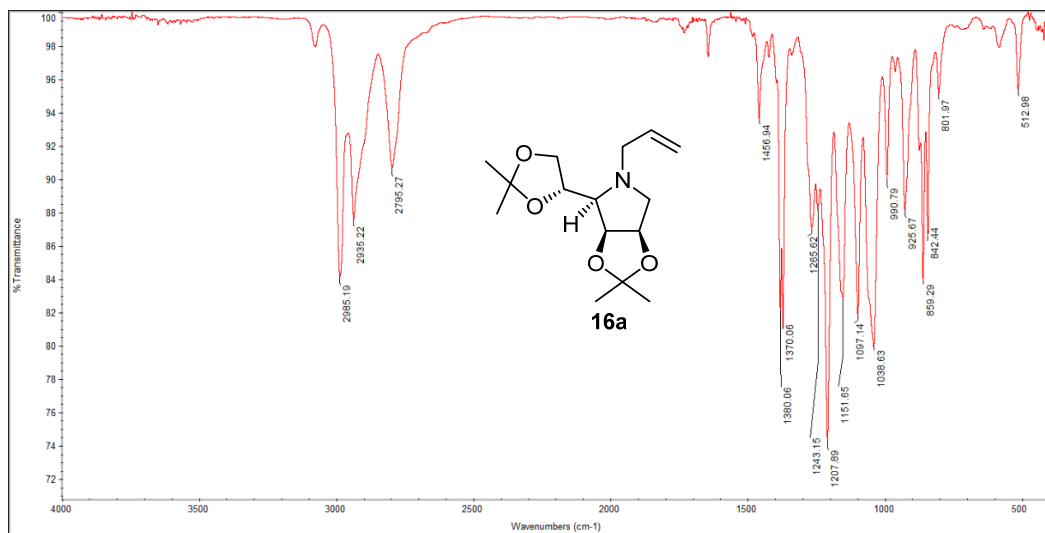


```

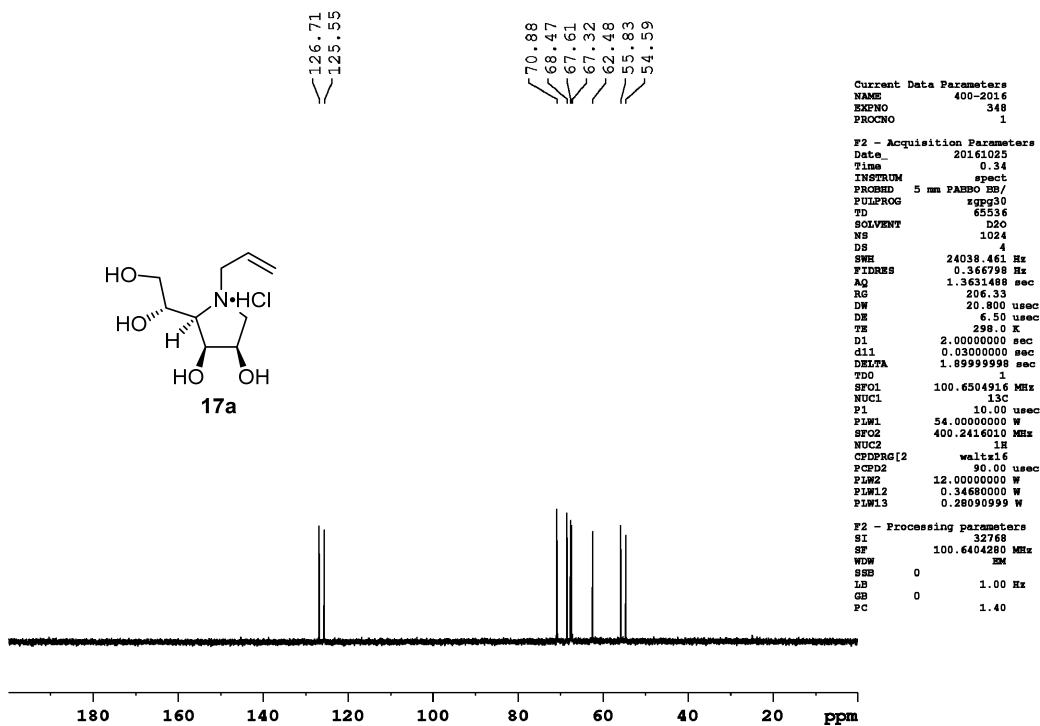
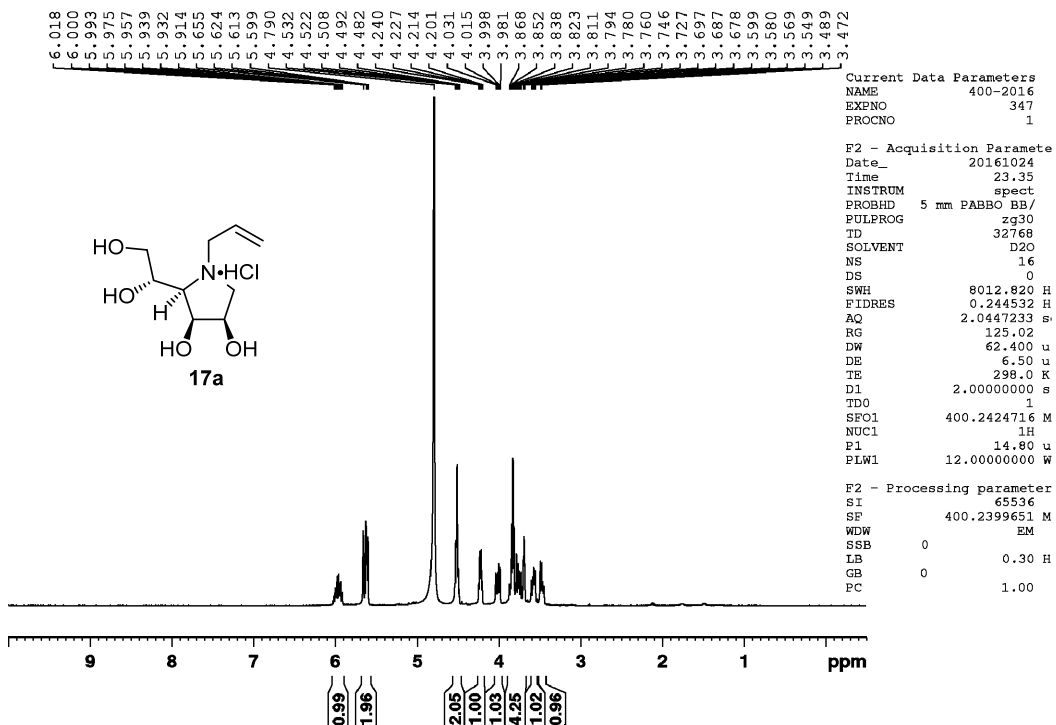
Current Data Parameters
NAME      500-2016
EXPNO    113
PROCNO   1

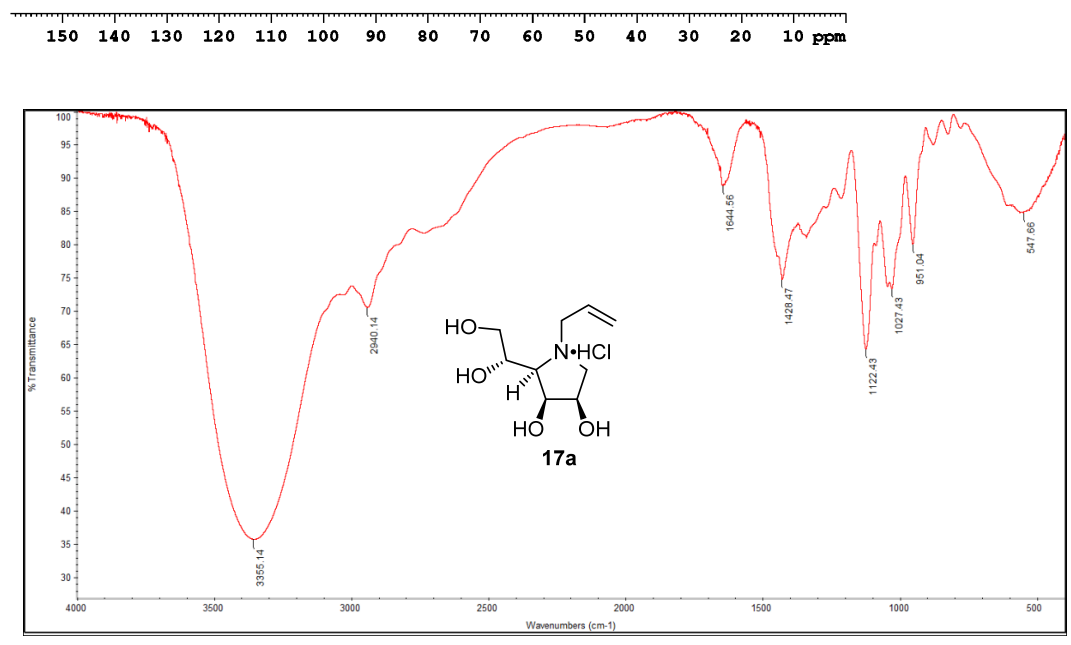
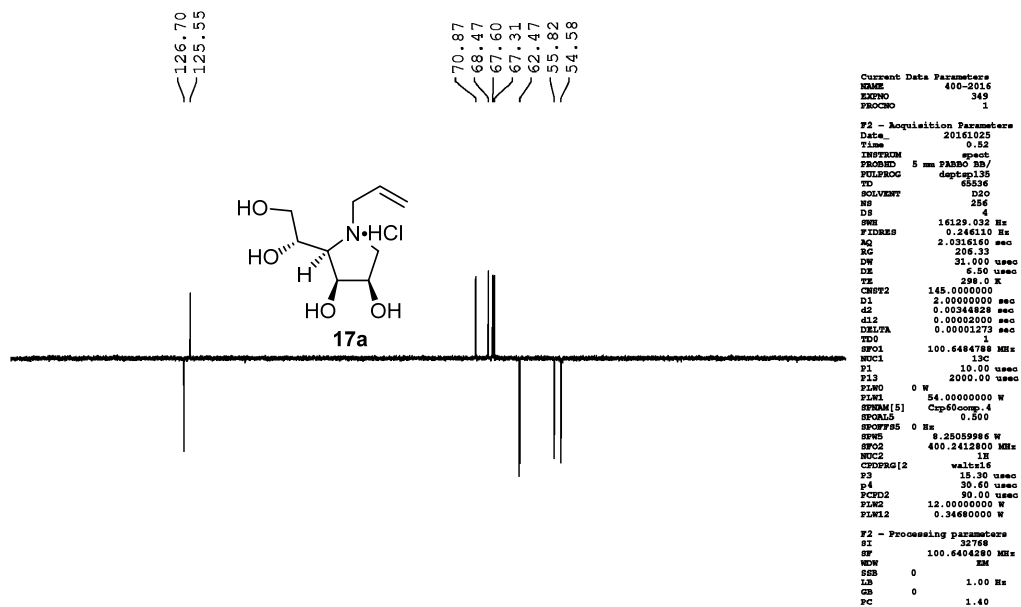
F2 - Acquisition Parameters
Date_    20161021
Time     1.44
INSTRUM  spect
PROBHD   5 mm CPBPR0 3B
PULPROG  deptgpr135
TD        65536
SOLVENT  CDCl3
NS        256
DS        4
SWH       20161.291 Hz
FIDRES   0.307637 Hz
AQ        1.6282928 sec
RG        192.88
SQ        24.400 usec
DE        18.00 usec
TE        298.0 K
CPCP2    145.000000
D1        2.0000000 sec
d2        0.00044928 sec
d12       0.0002000 sec
DELTA    0.00001248 sec
TD0       1
SFO1     125.7678486 MHz
SFC1     13C
P1        9.80 usec
P12      2000.00 usec
P120     0 W
PLM1     57.0000000 W
SFO2     500.1315995 MHz
SFC2     1H
CPCP2    waltz16
P3        10.70 usec
P4        21.40 usec
PCPD2    80.00 usec
PLM2     20.0000000 W
PLM12    0.35778001 W

F2 - Processing parameters
SI        32768
SF        125.7577885 MHz
WDW       EM
SSB       0
LB        1.00 Hz
GB        0
PC        1.40
  
```

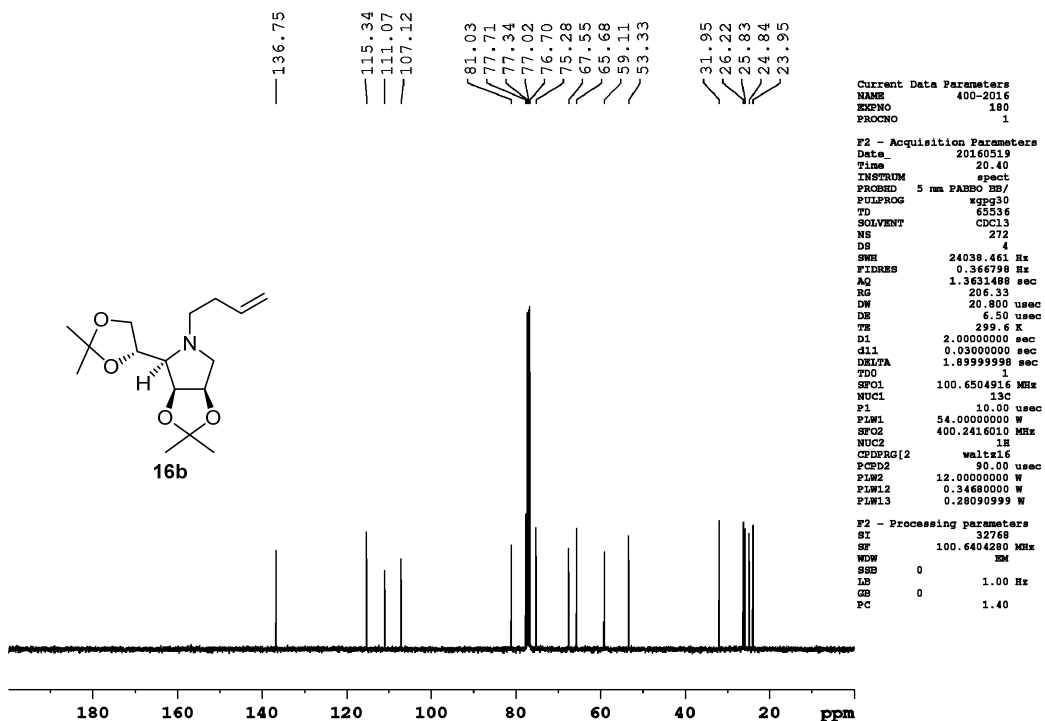
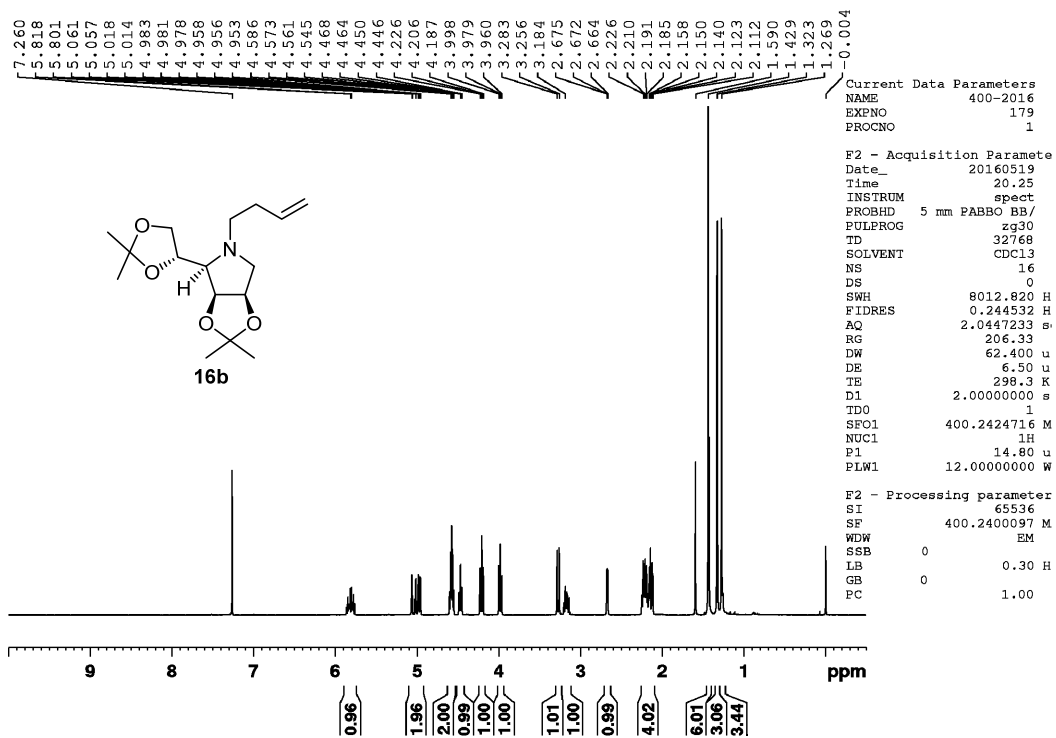


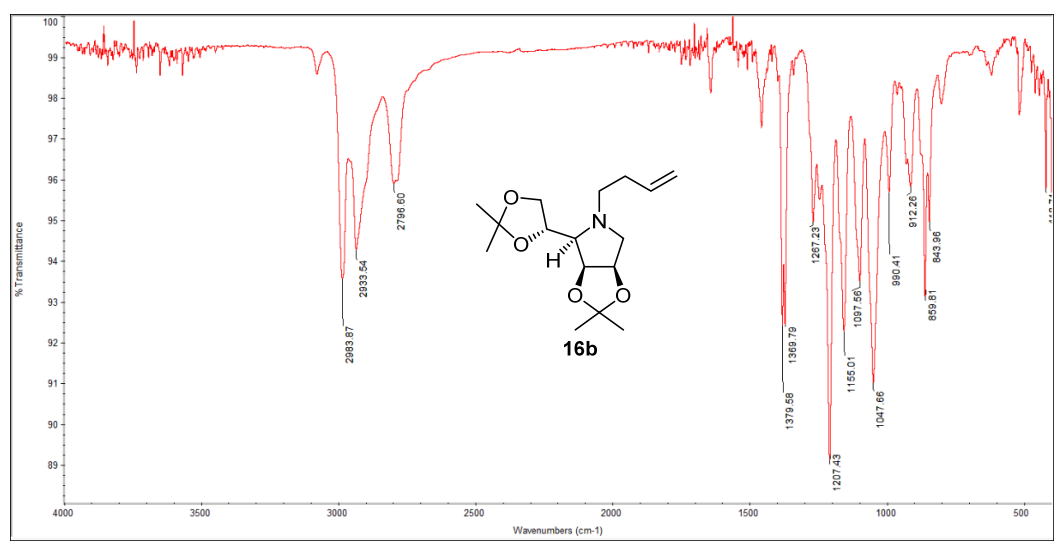
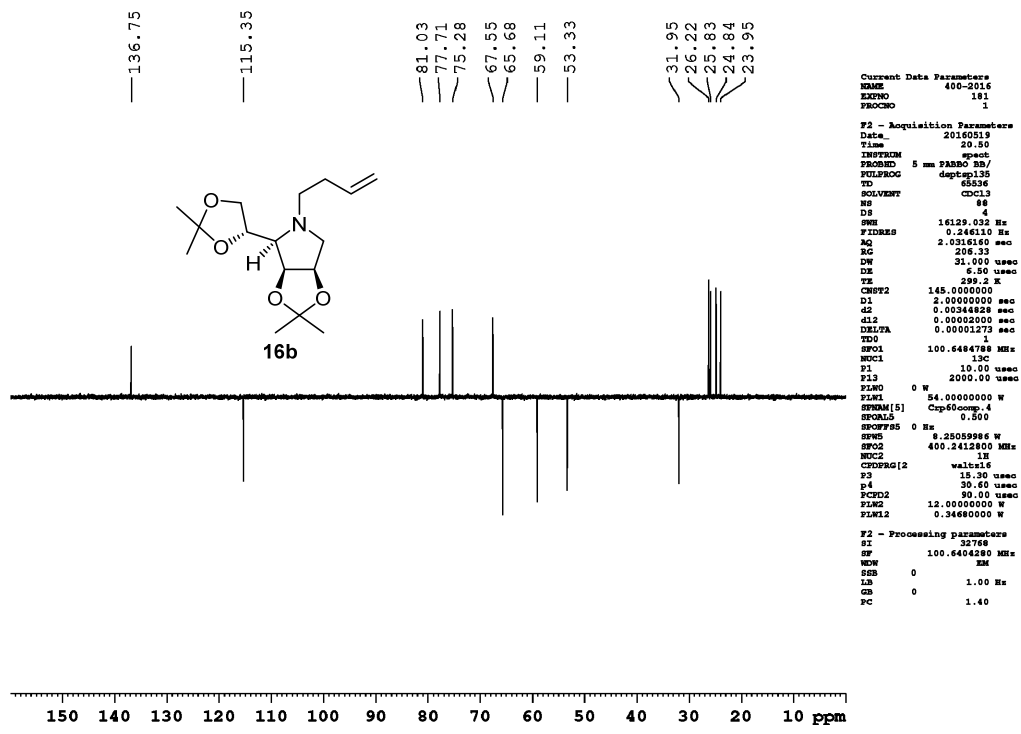
Compound 17a:



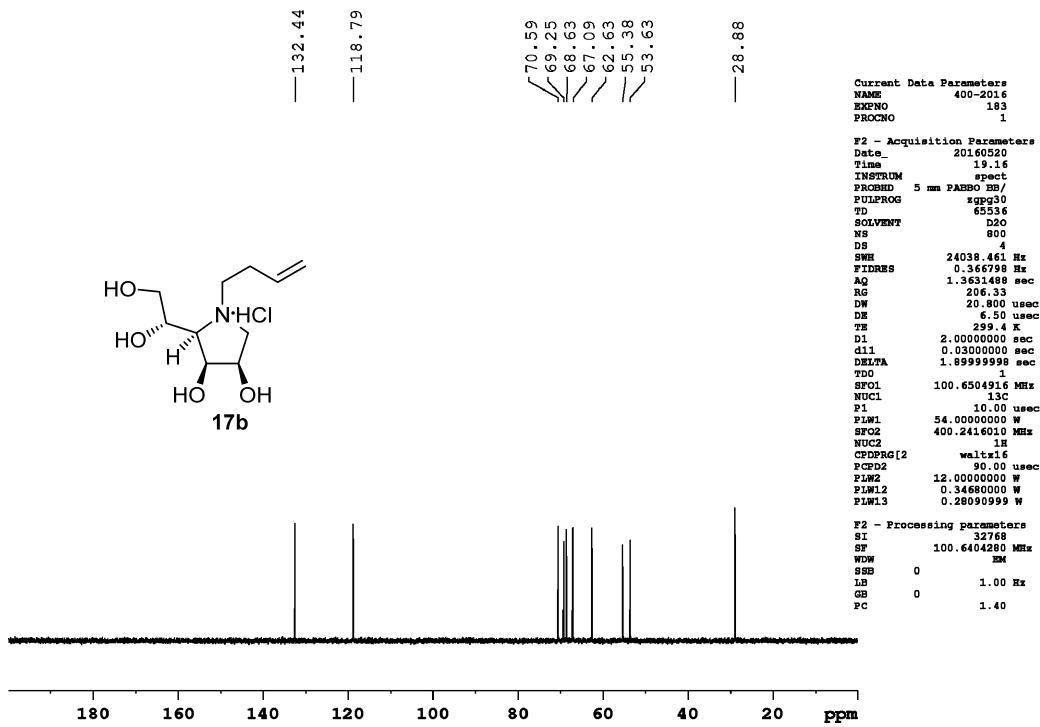
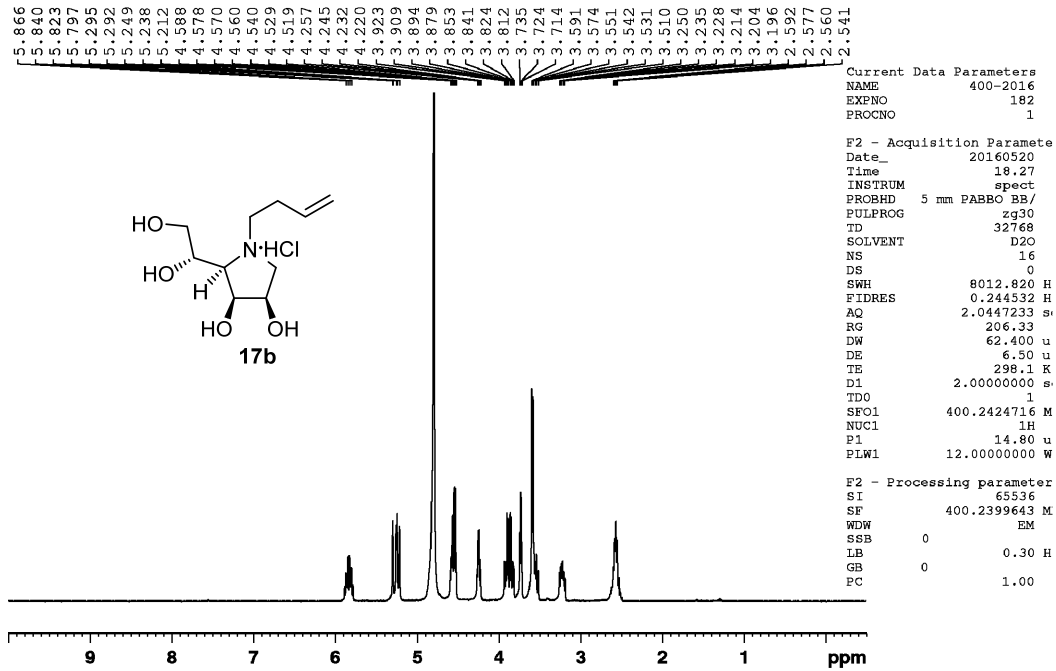


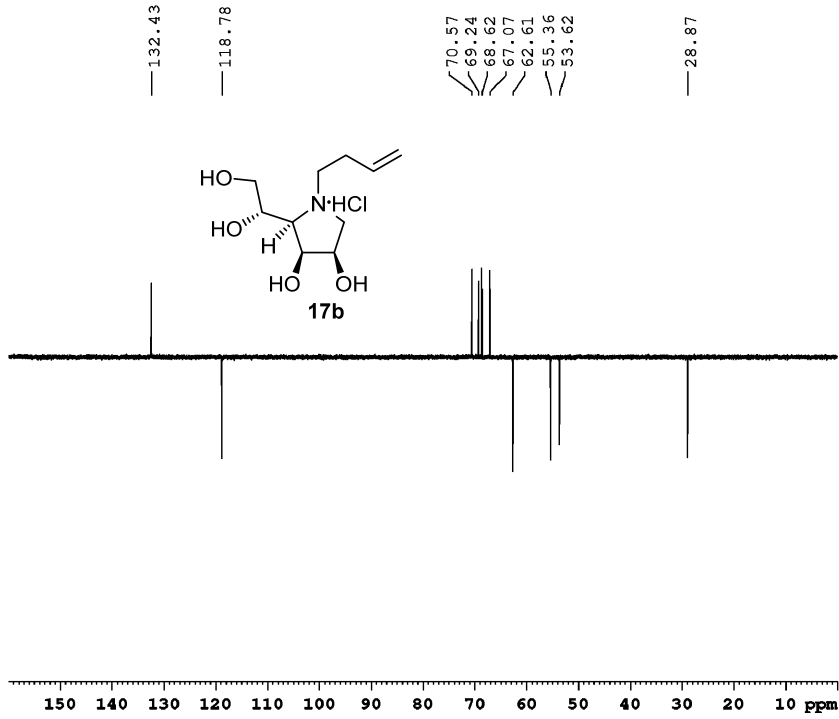
Compound 16b:





Compound 17b:



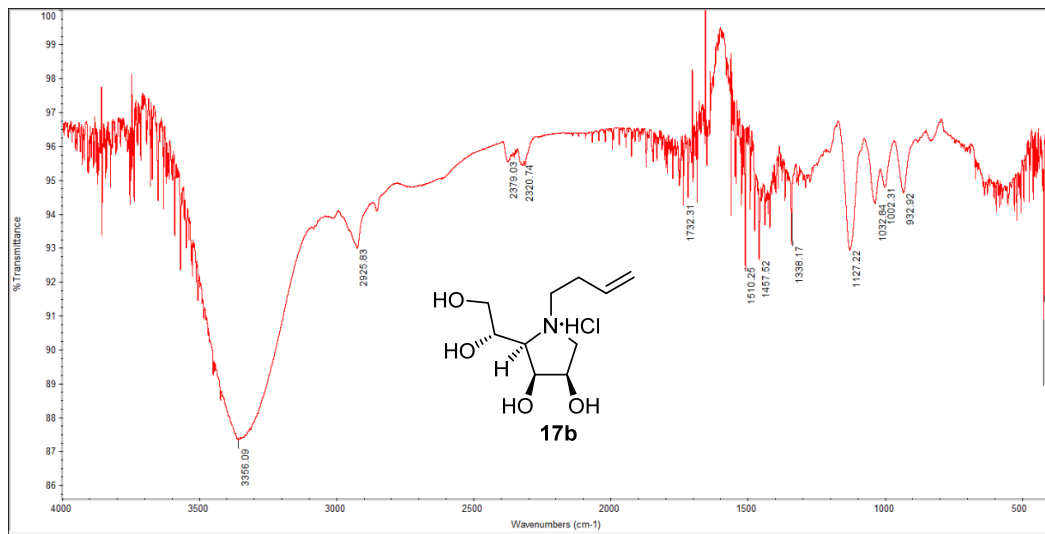


```

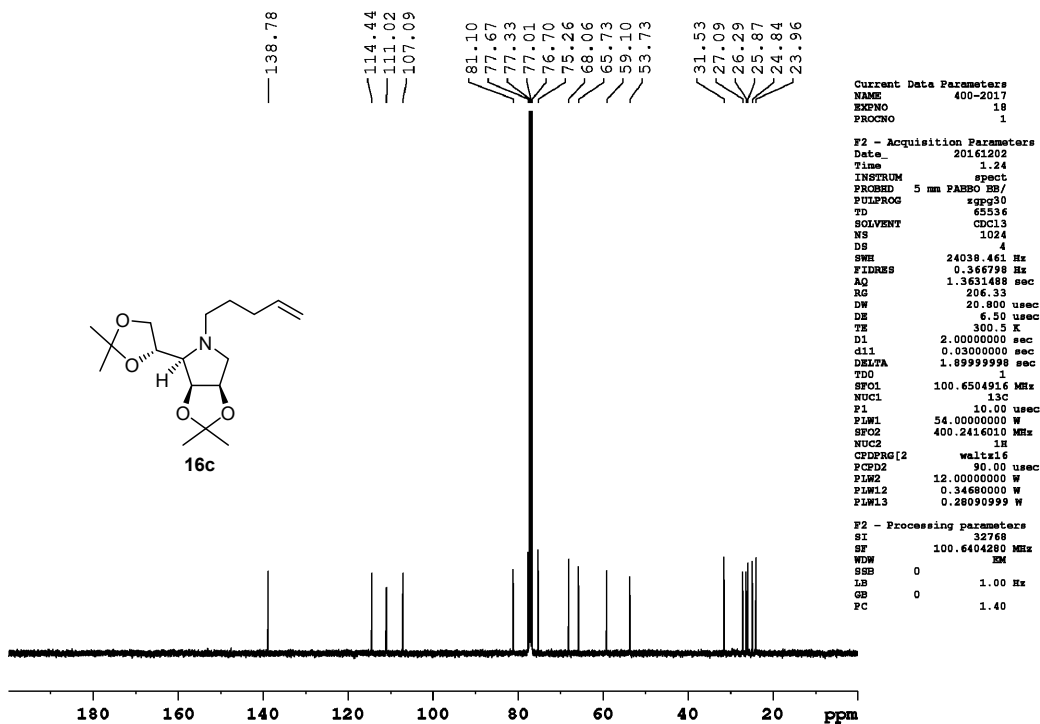
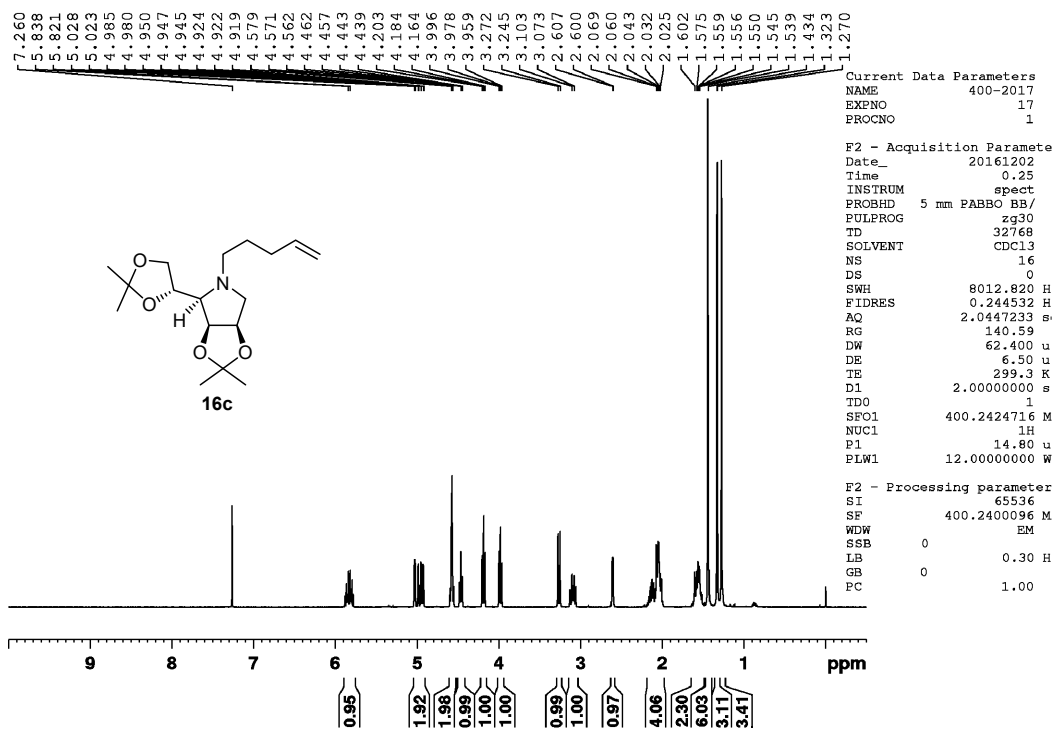
Current Data Parameters
NAME      400-2016
EXPNO     184
PROCNO    1

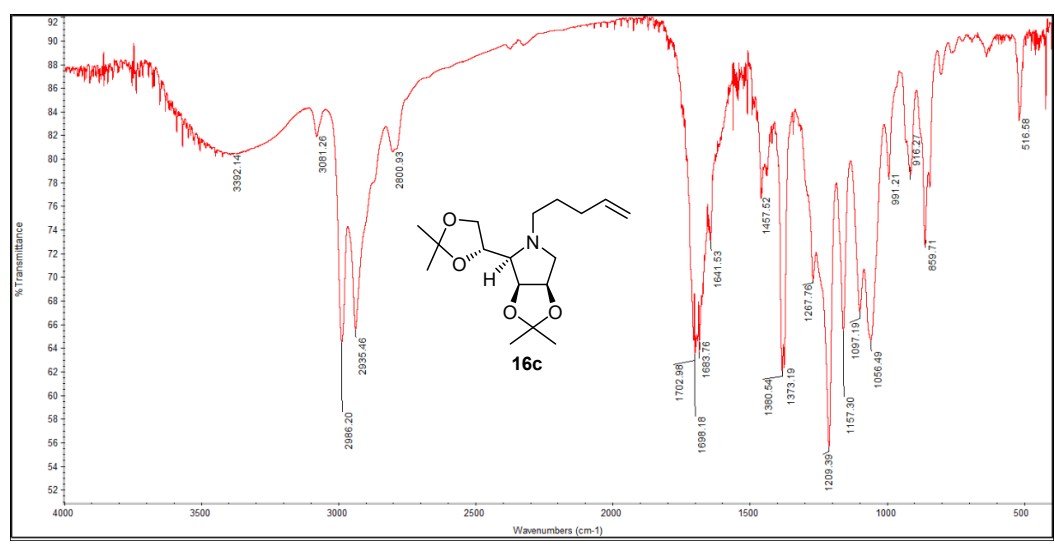
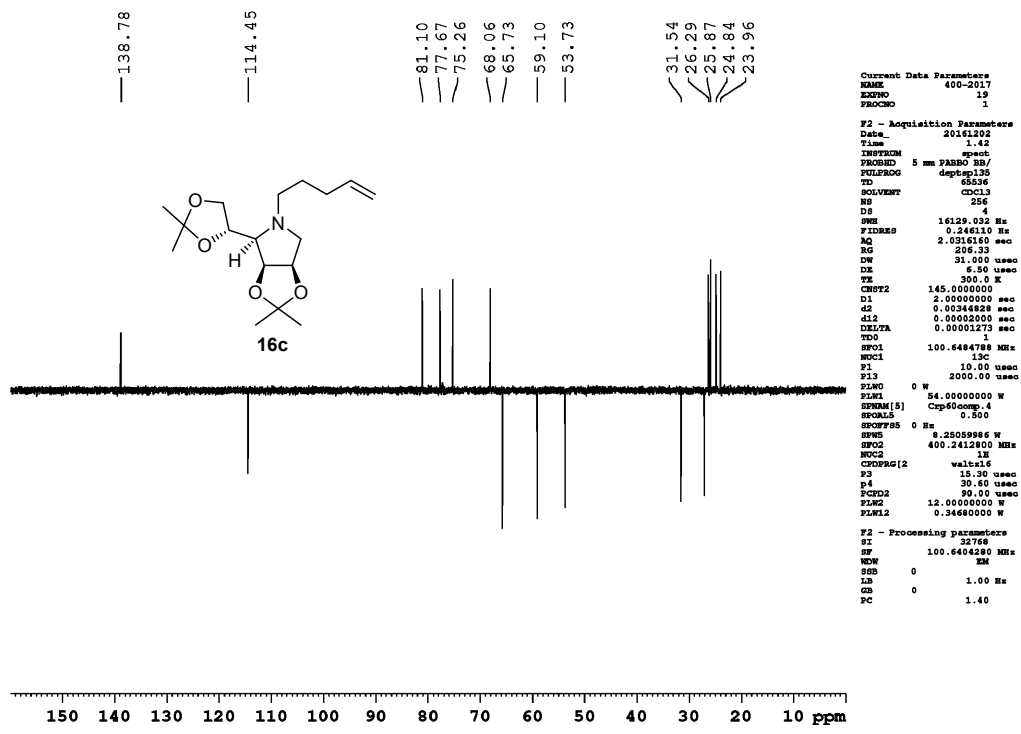
F2 - Acquisition Parameters
Date_     20160520
Time      19.29
INSTRUM   spect
PROBHD    5 mm F3BBO BB7
PULPROG   zgpg30
DEPTPRG   zgpg30
TD        65536
SOLVENT   D2O
NS         180
DS         4
SWH        16129.032 Hz
FIDRES    0.246130 Hz
AQ        2.0316160 sec
RG         206.33
DW        31.000 usec
DE        6.80 usec
TE        298.2 K
CMT2      145.0000000
D1         2.0000000 sec
d2         0.00344828 sec
d12        0.00002000 sec
DELTA     0.00001273 sec
TDO        1
SFO1      100.6484788 MHz
NUC1       13C
P1         10.00 usec
P2         2000.00 usec
PL1        0 W
PL2        54.0000000 W
SFO2      400.2412800 MHz
NUC2       1H
PCPDPRG2  waltz16
P3         15.30 usec
P4         30.60 usec
PCPD2     90.00 usec
PL3       12.0000000 W
PL4       0.3468000 W

F2 - Processing parameters
SI         32768
SF         100.6404280 MHz
WDW        EM
SSB        0
GB         0
PC         1.40
  
```

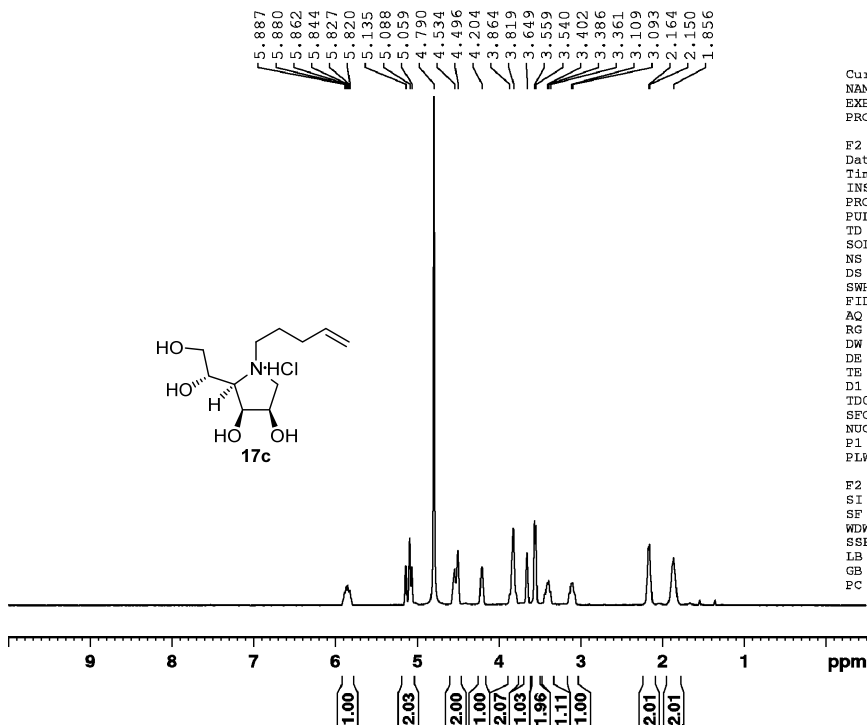


Compound 16c:





Compound 17c:

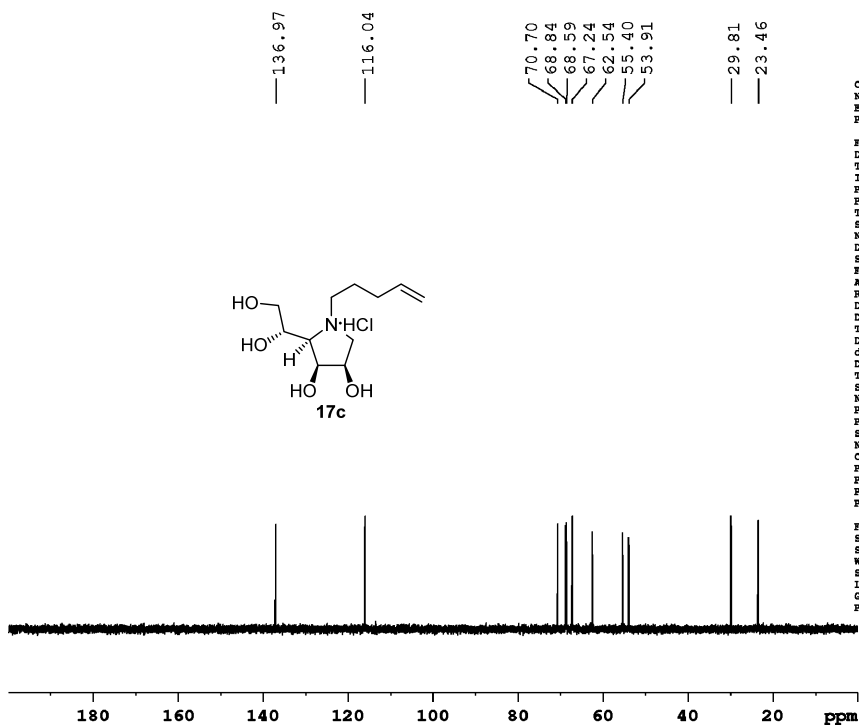


```

Current Data Parameters
NAME      400-2017
EXPNO    20
PROCNO   1

F2 - Acquisition Paramete
Date_    20161204
Time     21.26
INSTRUM spect
PROBHD   5 mm PABBO BB/
PULPROG zg30
TD       32768
SOLVENT  D2O
NS       16
DS       0
SWH      8012.820 H
FIDRES   0.244532 H
AQ       2.0447233 s
RG       206.33
DW       62.400 u
DE       6.50 u
TE       298.1 K
D1       2.0000000 s
TDO      1
SFO1     400.2424716 M
NUC1     1H
P1       14.80 u
PLW1     12.0000000 W

F2 - Processing parameter
SI       65536
SF       400.2399643 M
WDW      EM
SSB      0
LB       0.30 H
GB       0
PC       1.00
    
```

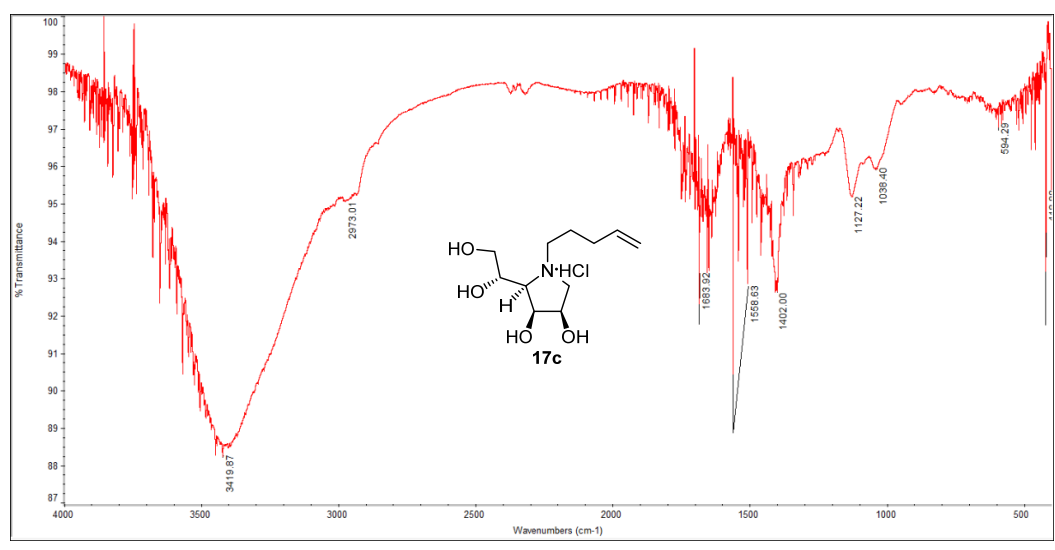
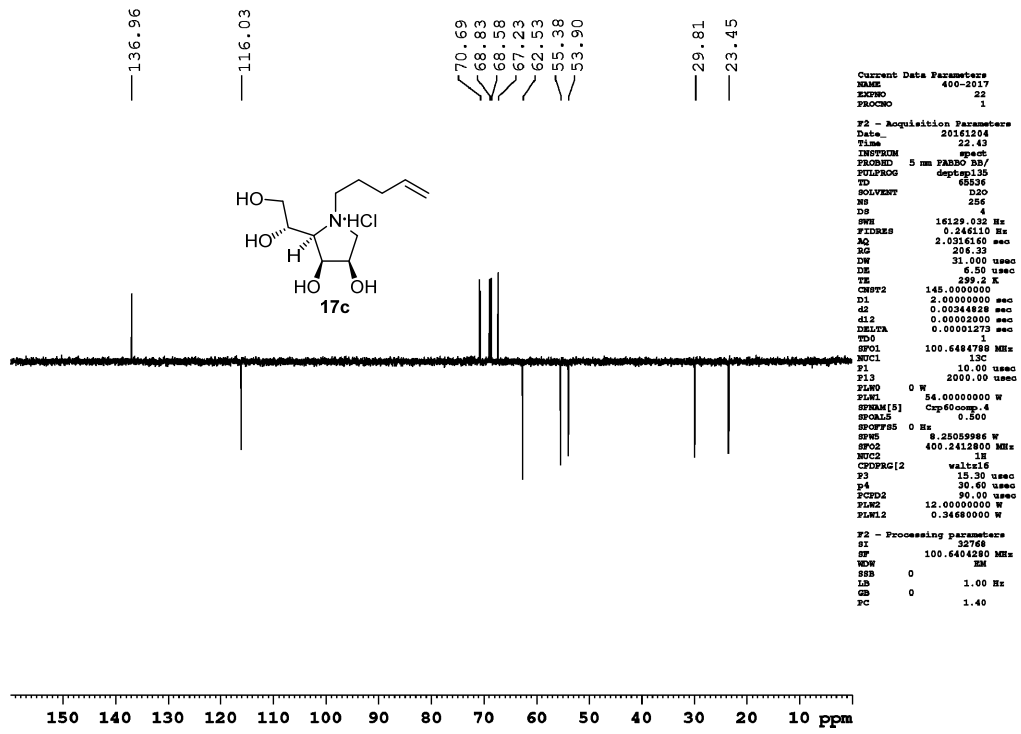


```

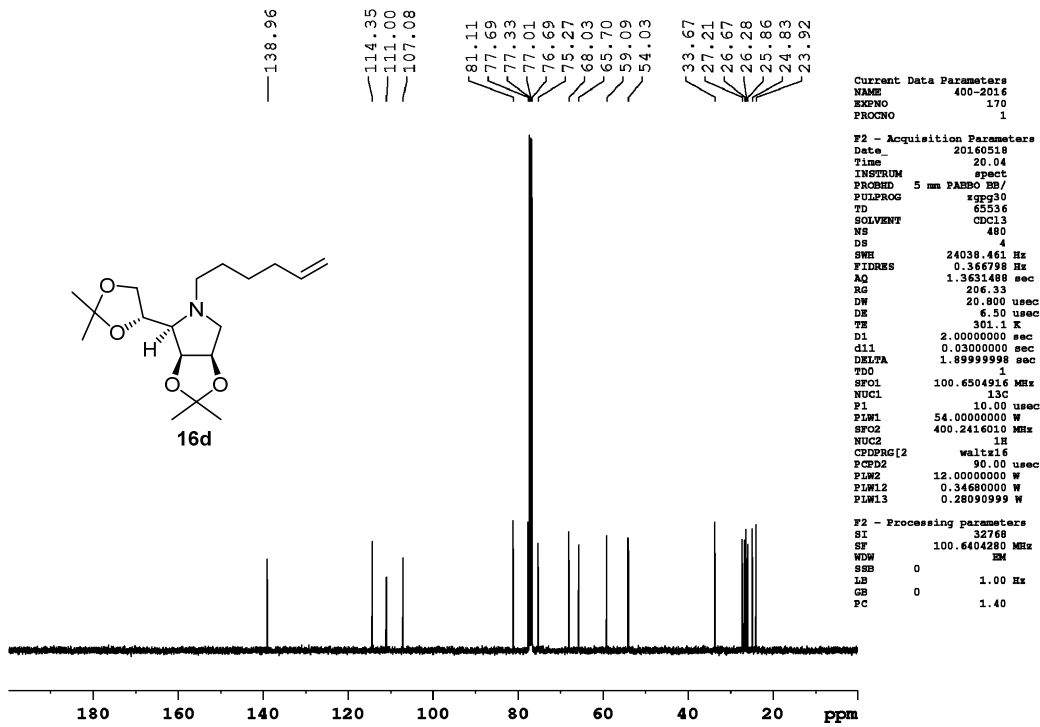
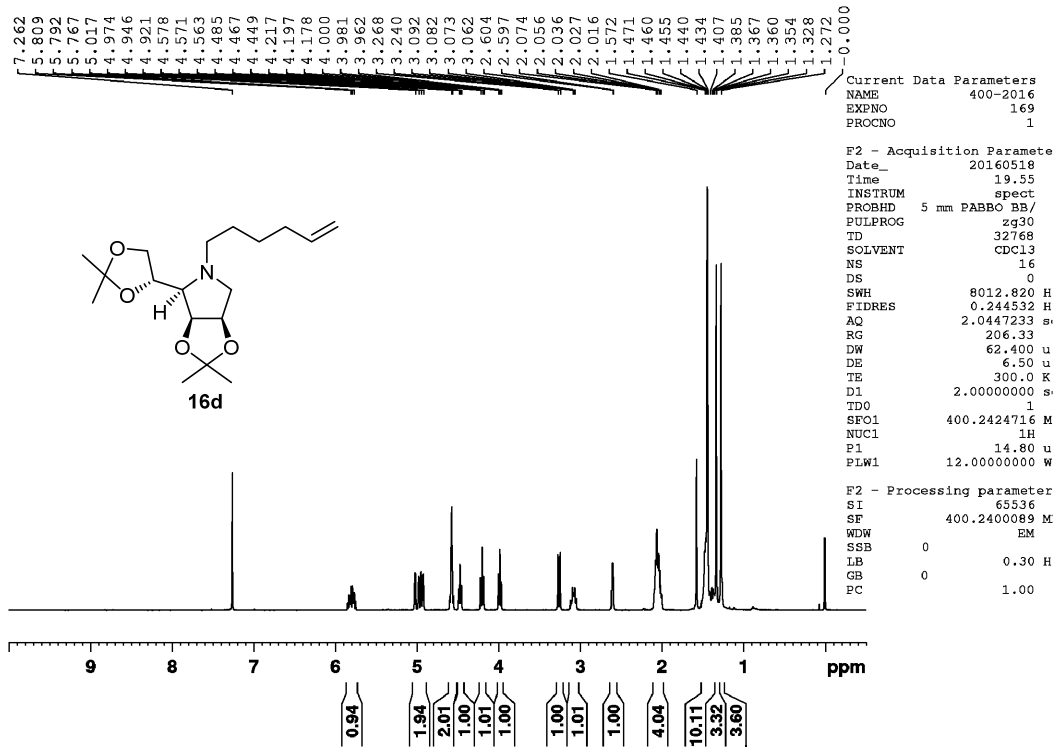
Current Data Parameters
NAME      400-2017
EXPNO    21
PROCNO   1

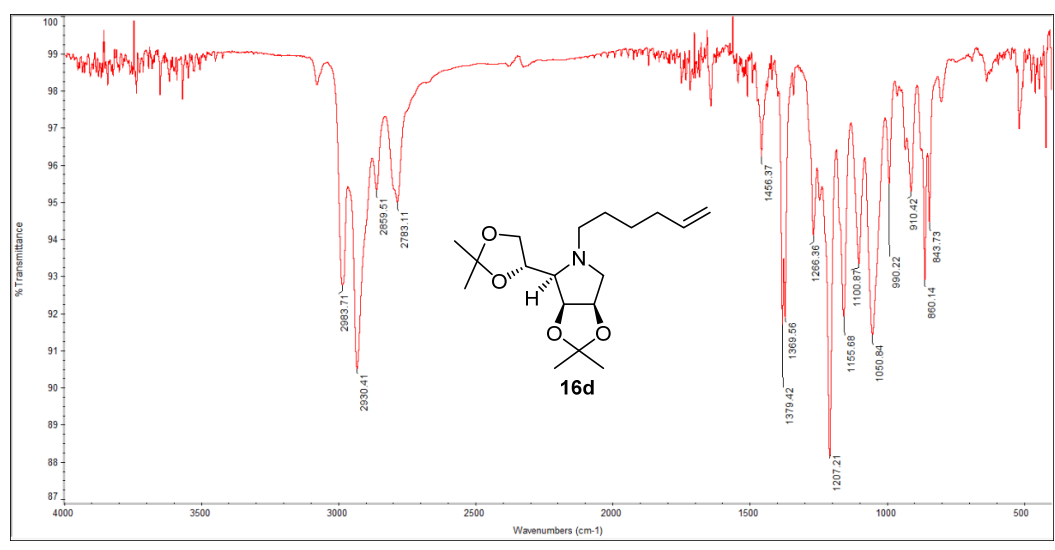
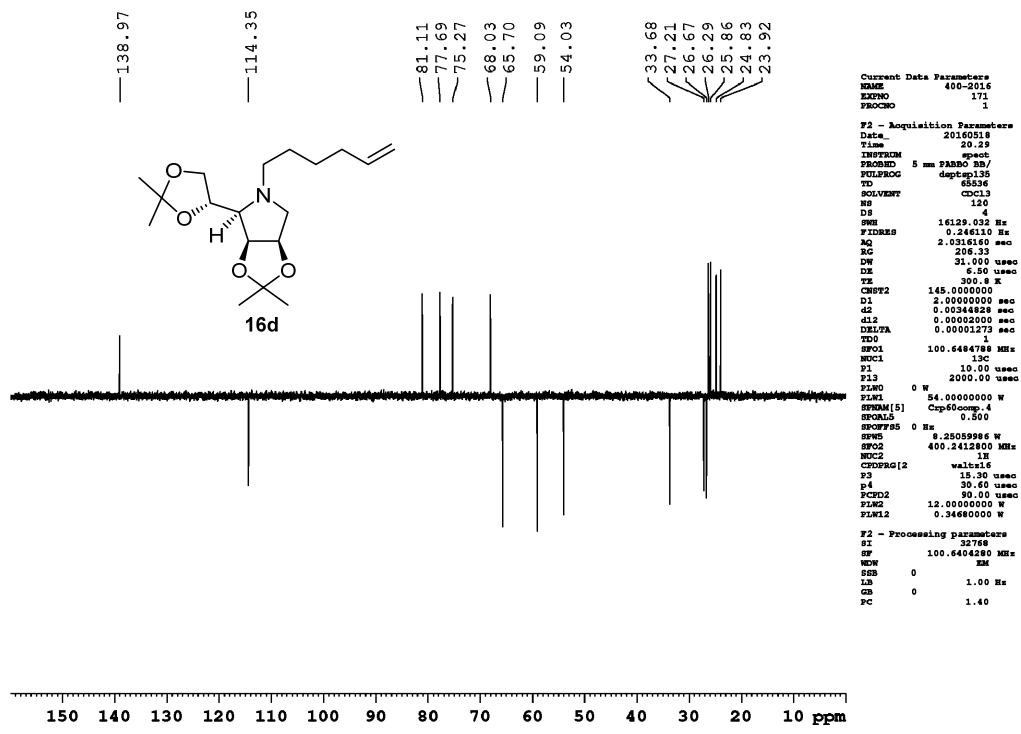
F2 - Acquisition Parameters
Date_    20161204
Time     22.25
INSTRUM spect
PROBHD   5 mm PABBO BB/
PULPROG zgpg30
TD       65536
SOLVENT  D2O
NS       1024
DS       0
SWH      24038.461 Hz
FIDRES   0.366798 Hz
AQ       1.3631488 sec
RG       206.33
DW       20.800 usec
DE       6.50 usec
TE       299.7 K
D1       2.0000000 sec
d11      0.0300000 sec
DELTA    1.89999998 sec
TDO      1
SFO1     100.6504916 MHz
NUC1     13C
P1       10.00 usec
PLW1     54.0000000 W
SFO2     400.2416010 MHz
NUC2     1H
CPDPRG2 waltz16
PCPD2    90.00 usec
PLW2     12.0000000 W
PLW12    0.3468000 W
PLW13    0.28090999 W

F2 - Processing parameters
SI       32768
SF       100.6404280 MHz
WDW      EM
SSB      0
LB       1.00 Hz
GB       0
PC       1.40
    
```

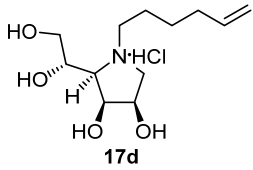
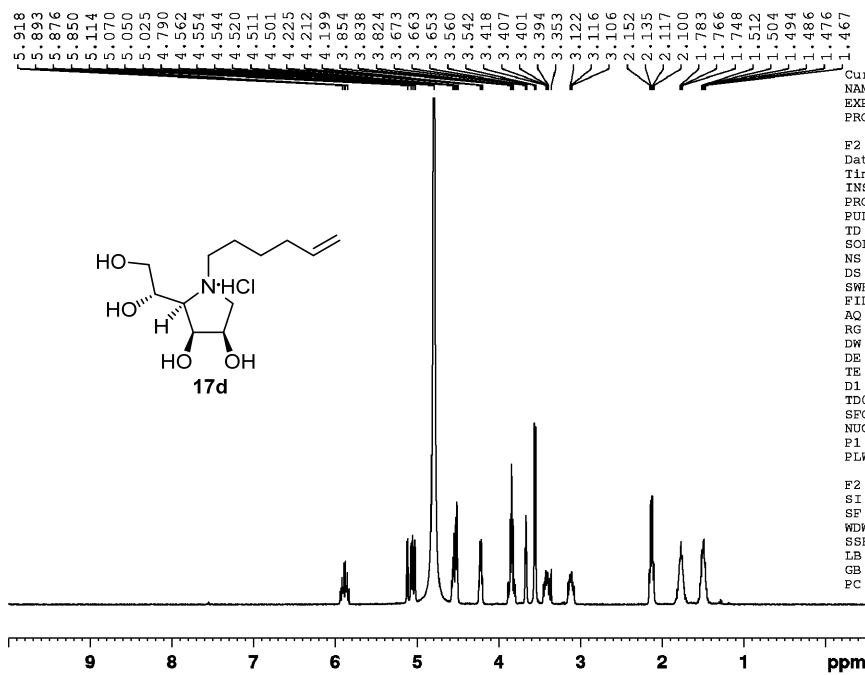


Compound 16d:





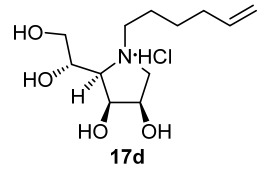
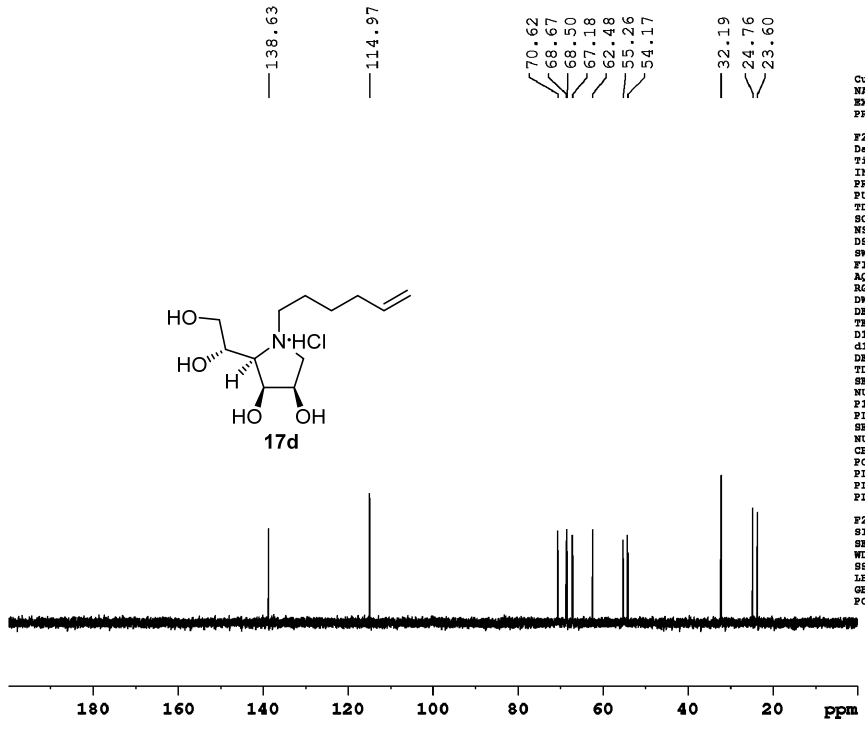
Compound 17d:



Current Data Parameters
 NAME 400-2016
 EXPNO 172
 PROCNO 1

F2 - Acquisition Parameters
 Date_ 20160519
 Time 13.17
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zg30
 TD 32768
 SOLVENT D2O
 NS 16
 DS 0
 SWH 8012.820 H
 FIDRES 0.244532 H
 AQ 2.0447233 s
 RG 206.33
 DW 62.400 u
 DE 6.50 u
 TE 298.1 K
 D1 2.00000000 s
 TD0 1
 SFO1 400.2424716 M
 NUC1 1H
 P1 14.80 u
 PLW1 12.00000000 W

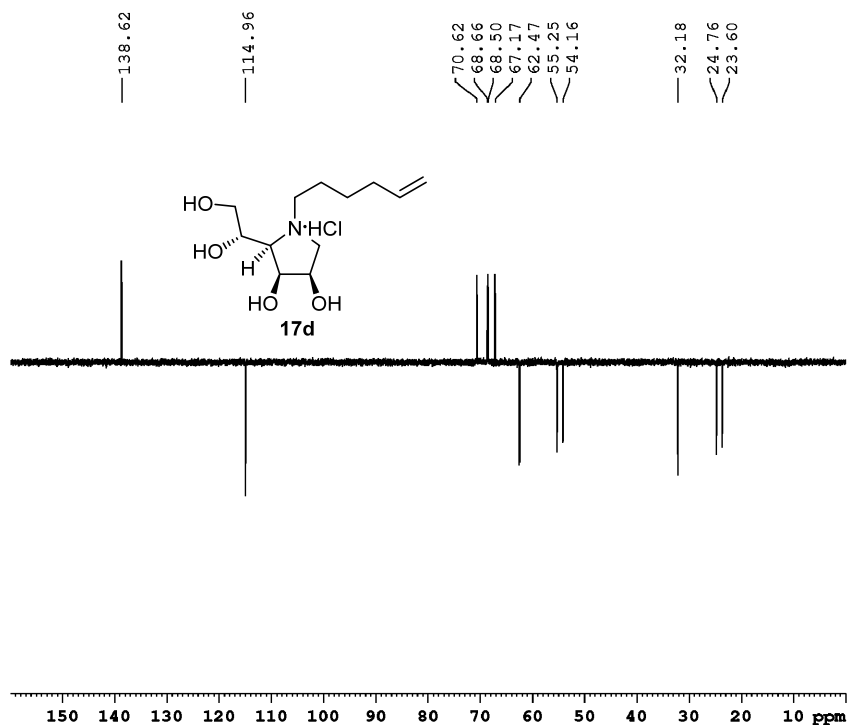
F2 - Processing parameter
 SI 65536
 SF 400.2399591 M
 WDW EM
 SSB 0
 LB 0.30 H
 GB 0
 PC 1.00



Current Data Parameters
 NAME 400-2016
 EXPNO 176
 PROCNO 1

F2 - Acquisition Parameters
 Date_ 20160519
 Time 14.40
 INSTRUM spect
 PROBHD 5 mm PABBO BB/
 PULPROG zgpg30
 TD 65536
 SOLVENT D2O
 NS 700
 DS 4
 SWH 24038.461 Hz
 FIDRES 0.366798 Hz
 AQ 1.3631488 sec
 RG 206.33
 DW 20.800 usec
 DE 6.50 usec
 TE 299.1 K
 D1 2.00000000 sec
 d11 0.03000000 sec
 DELTA 1.89999998 sec
 TD0 1
 SFO1 100.6504916 MHz
 NUC1 13C
 P1 10.00 usec
 PLW1 54.00000000 W
 SFO2 400.2416010 MHz
 NUC2 1H
 CPDPRG[2] waltz16
 PCPD2 90.00 usec
 PLW2 12.00000000 W
 PLW12 0.34680000 W
 PLW13 0.28090999 W

F2 - Processing parameters
 SI 32768
 SF 100.6404280 MHz
 WDW EM
 SSB 0
 LB 1.00 Hz
 GB 0
 PC 1.40

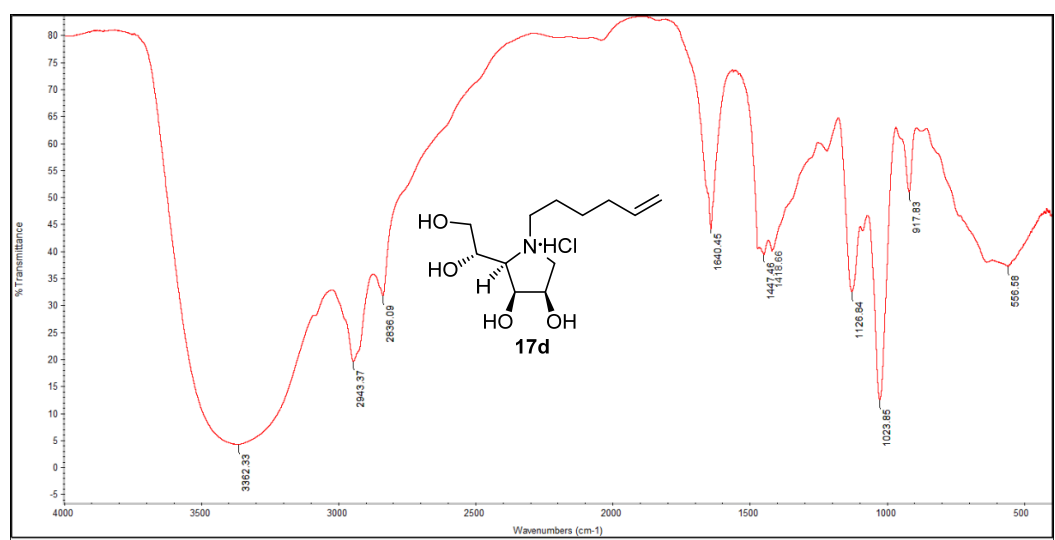


```

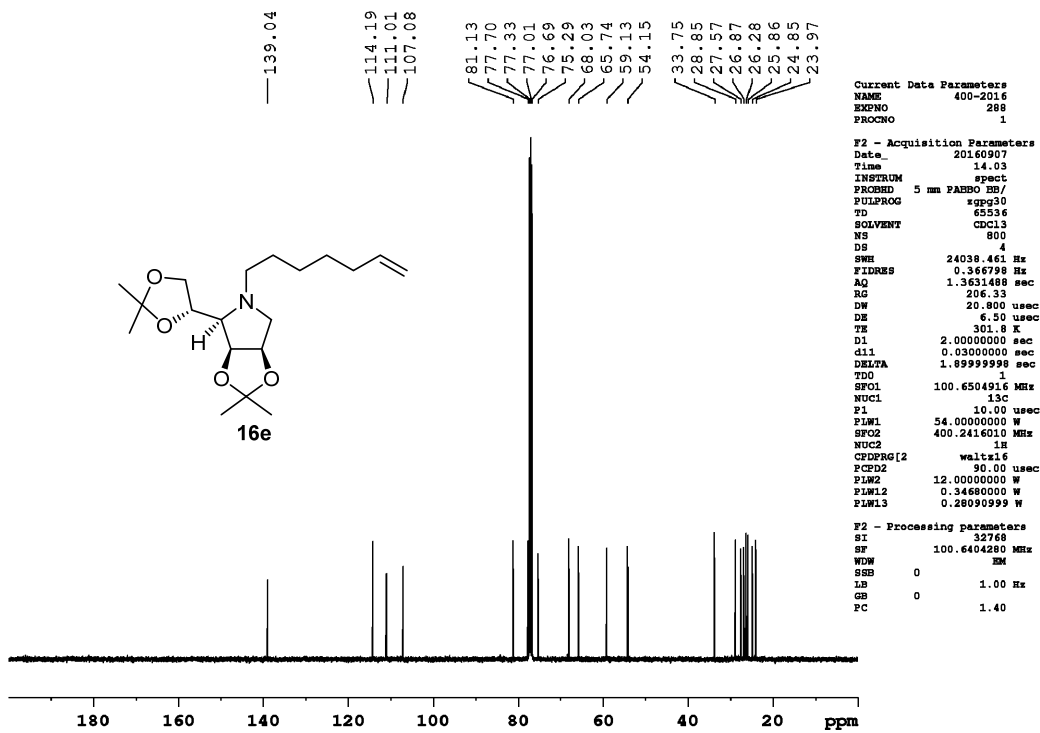
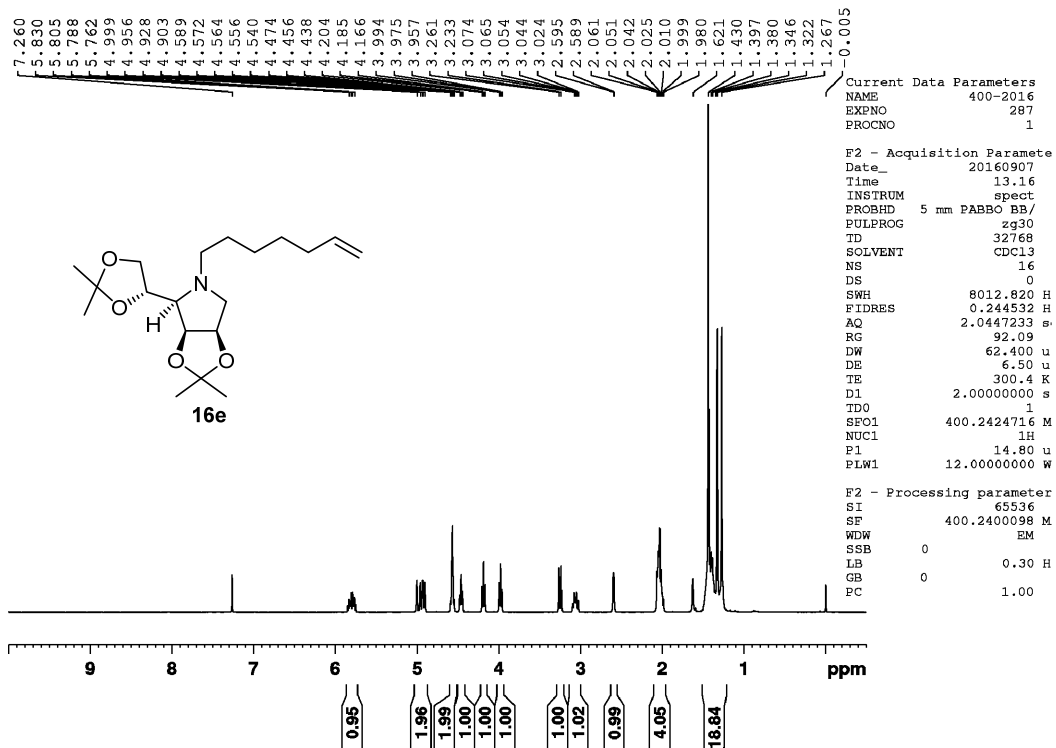
Current Data Parameters
NAME      400-2016
EXPNO    177
PROCNO   1

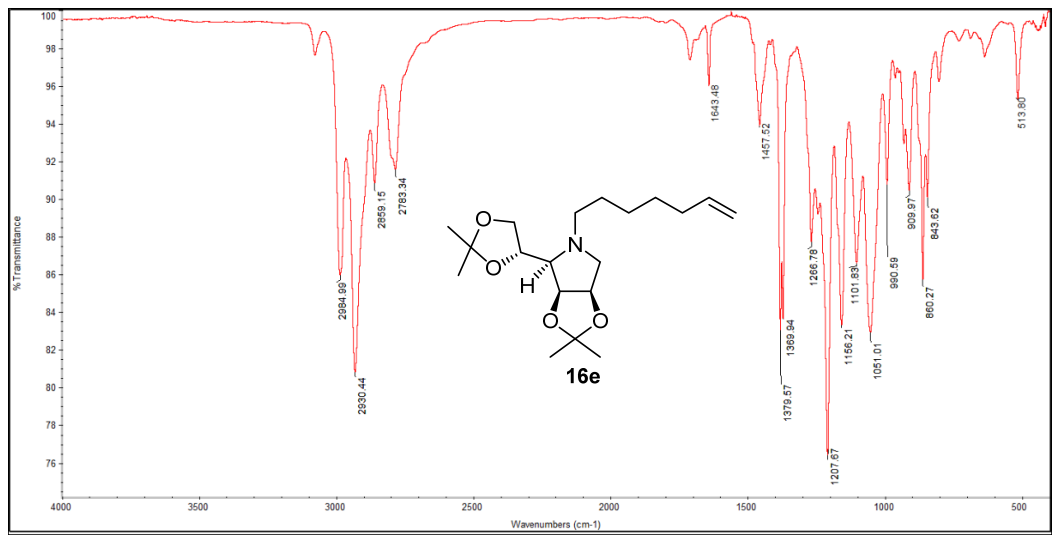
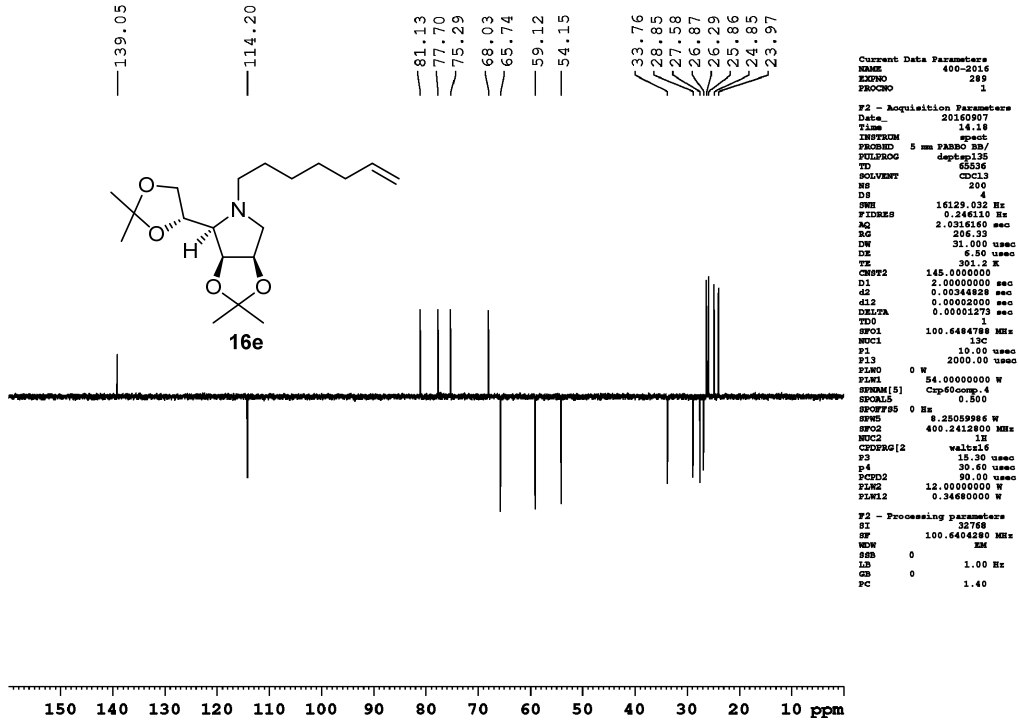
F2 - Acquisition Parameters
Date_    20160519
Time     15.06
INSTRUM  spect
PROBHD   5 mm PABBO BB7
PULPROG  zgpg30
TD        65536
SOLVENT  D2O
NS        200
DS        4
SWH       16129.032 Hz
FIDRES   0.246110 Hz
AQ        2.0316160 sec
RG         206.33
AQ        31.300 usec
DE         6.80 usec
TE        300.2 K
CQPC2    145.0000000
D1        2.0000000 sec
d2        0.00344828 sec
d12       0.0002000 sec
DELTA    0.0001273 sec
TD0       1
SFO1     100.6484788 MHz
NUC1      13C
P1        10.00 usec
P13       1.00 usec
P130     0 W
PL1      54.0000000 W
SFO1M15] Exp60comp-4
SFOAL5   0.500
SFOFFAS  0 Hz
SWS      8.25059986 W
SFO2     400.2412800 MHz
NUC2      1H
CQPC2[2] waltz16
P3        15.30 usec
P4         30.60 usec
PCPD2    80.00 usec
PLAS     12.0000000 W
PLM12    0.34680000 W

F2 - Processing parameters
SI        32768
SF        100.6404280 MHz
WDW       EM
SSB       0
GB        0
PC        1.40
  
```

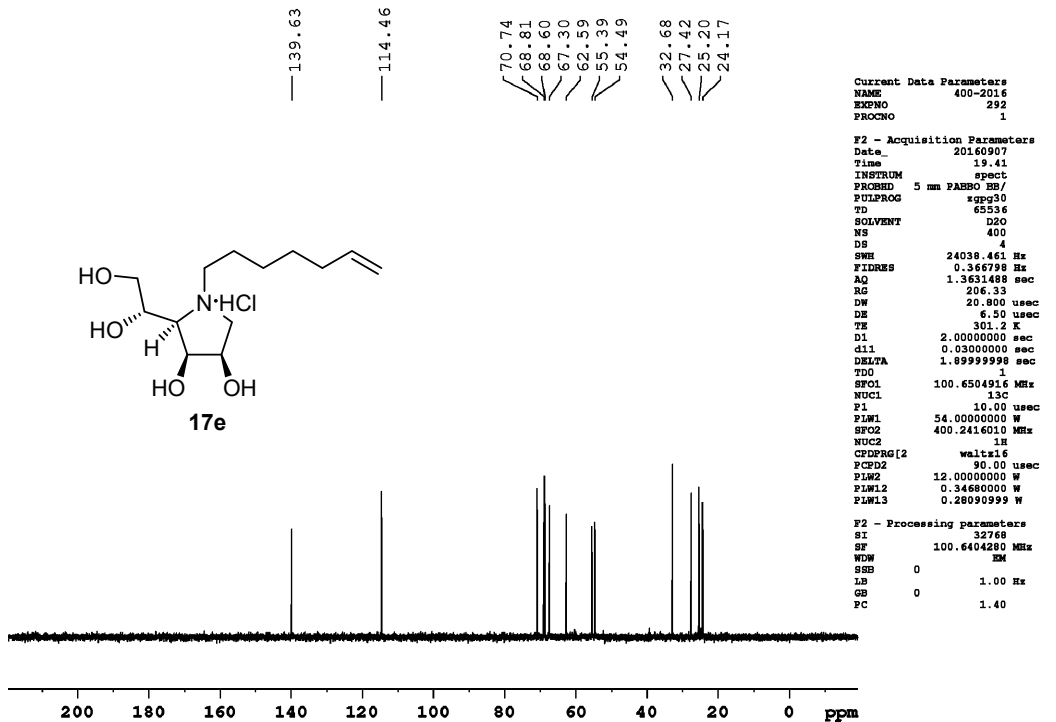
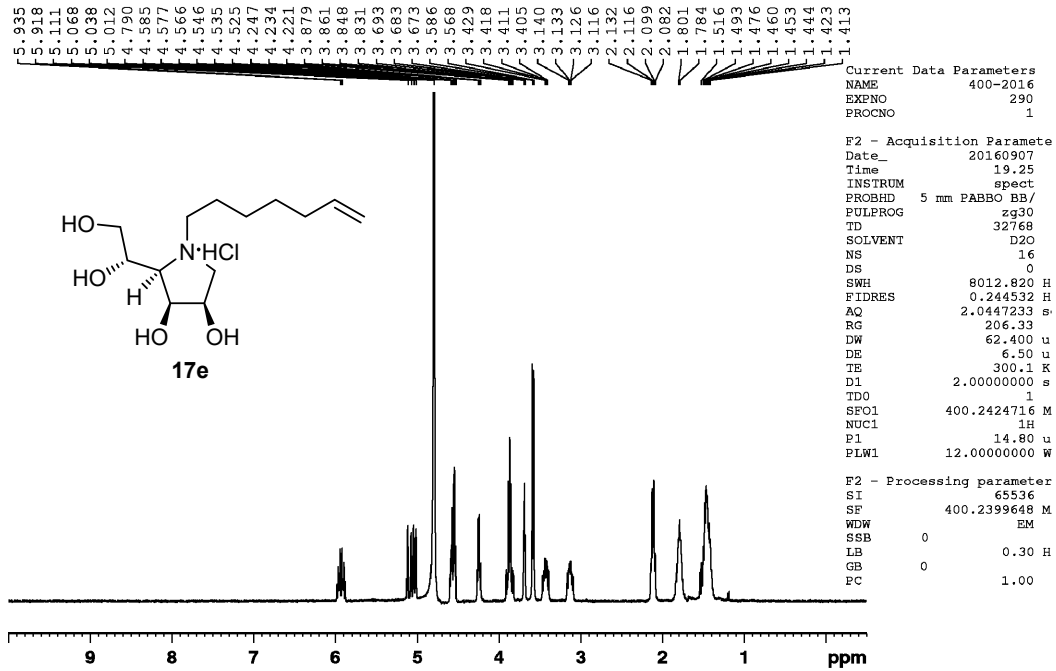


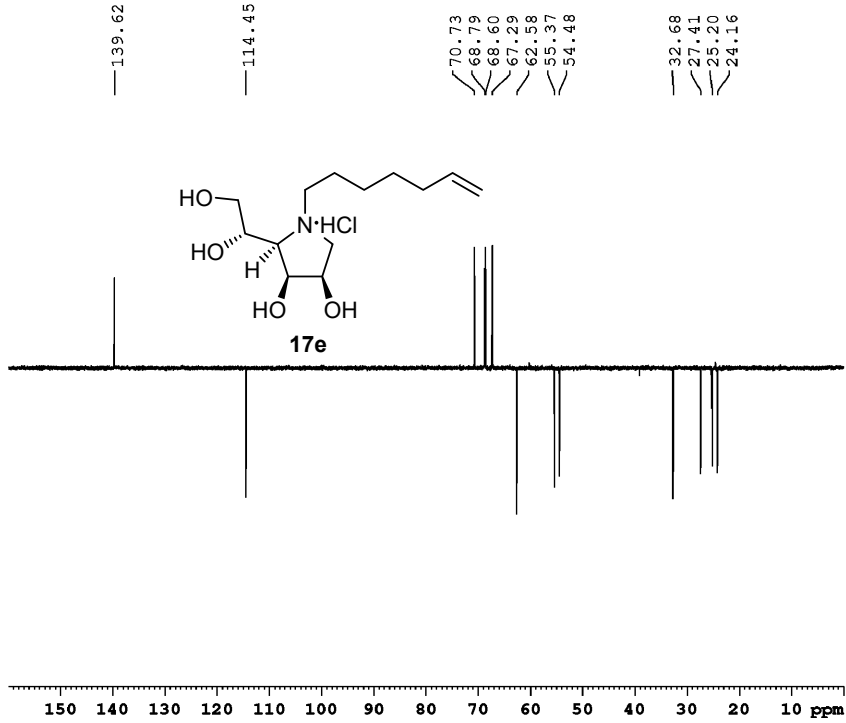
Compound 16e:





Compound 17e:



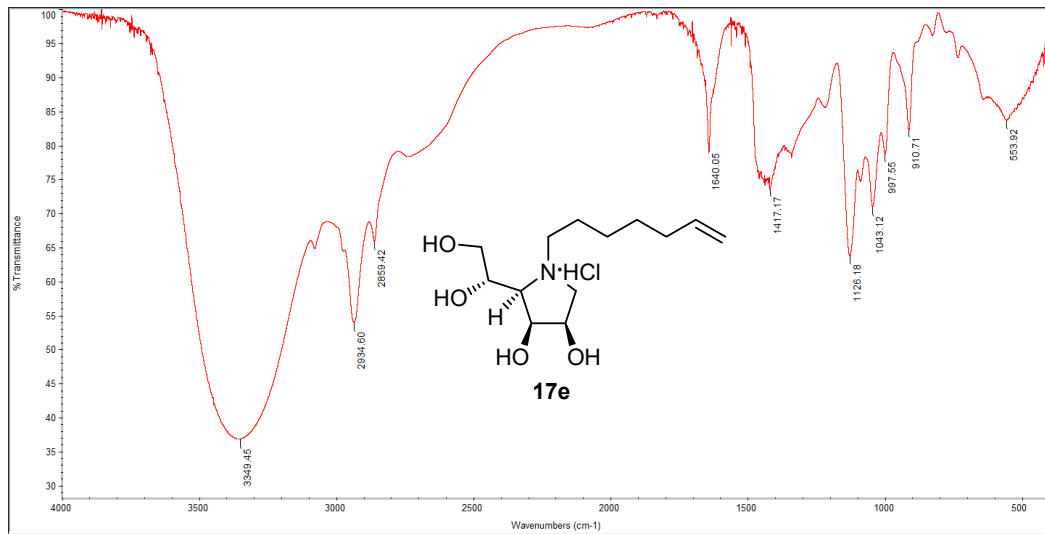


```

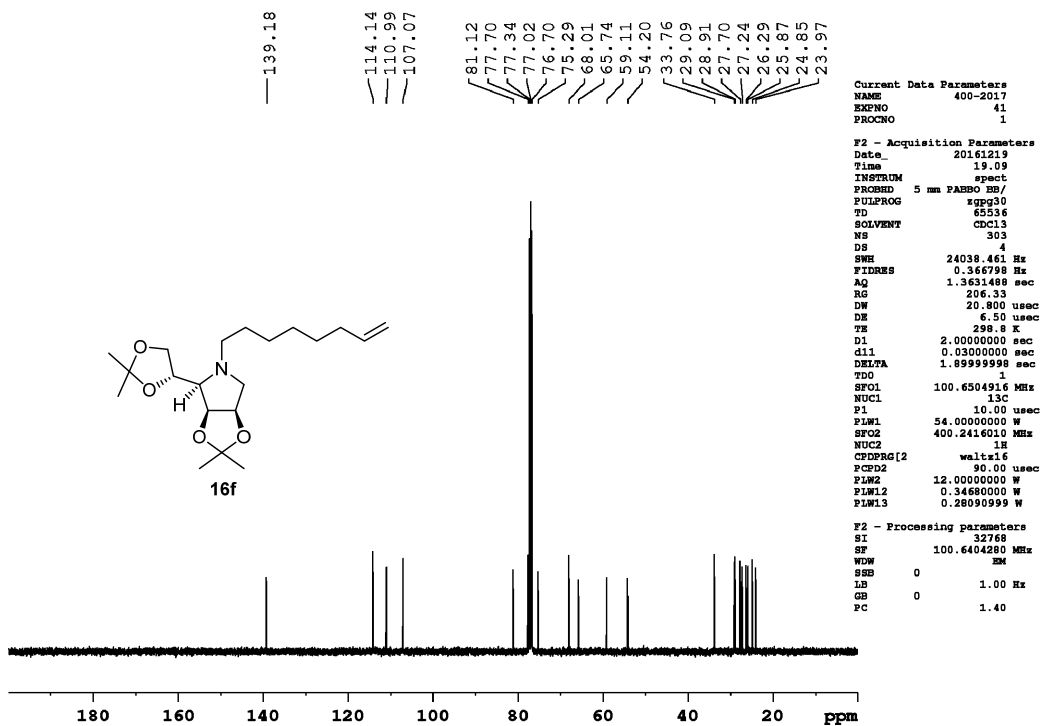
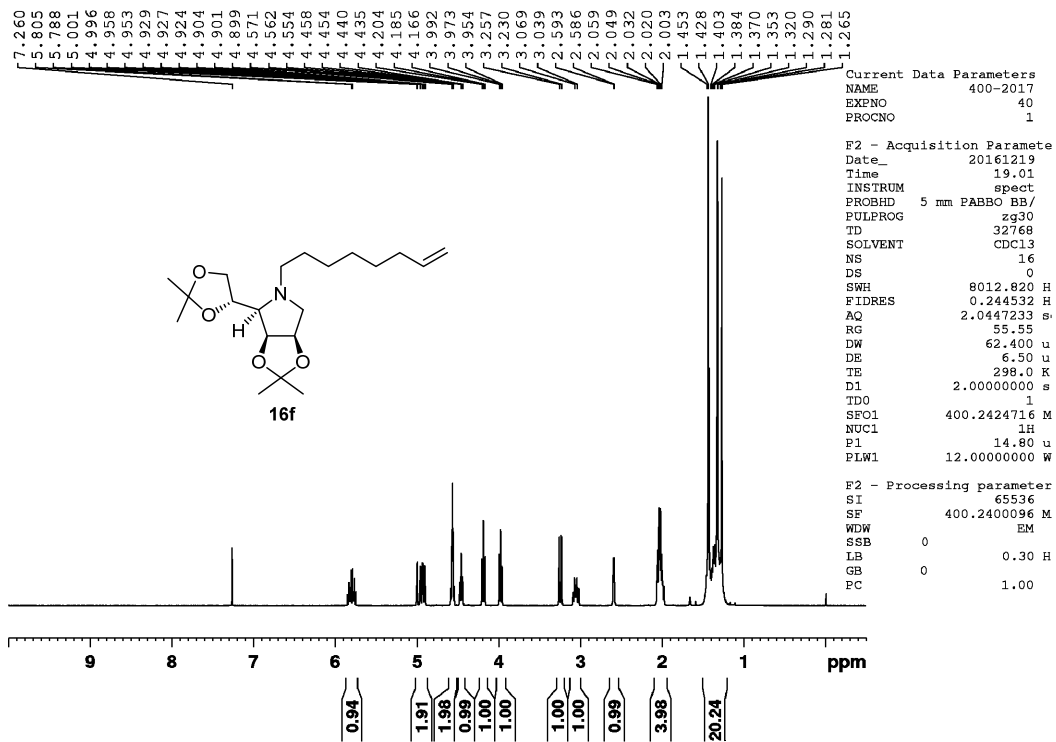
Current Data Parameters
NAME      400-2016
EXPNO    293
PROCNO    1

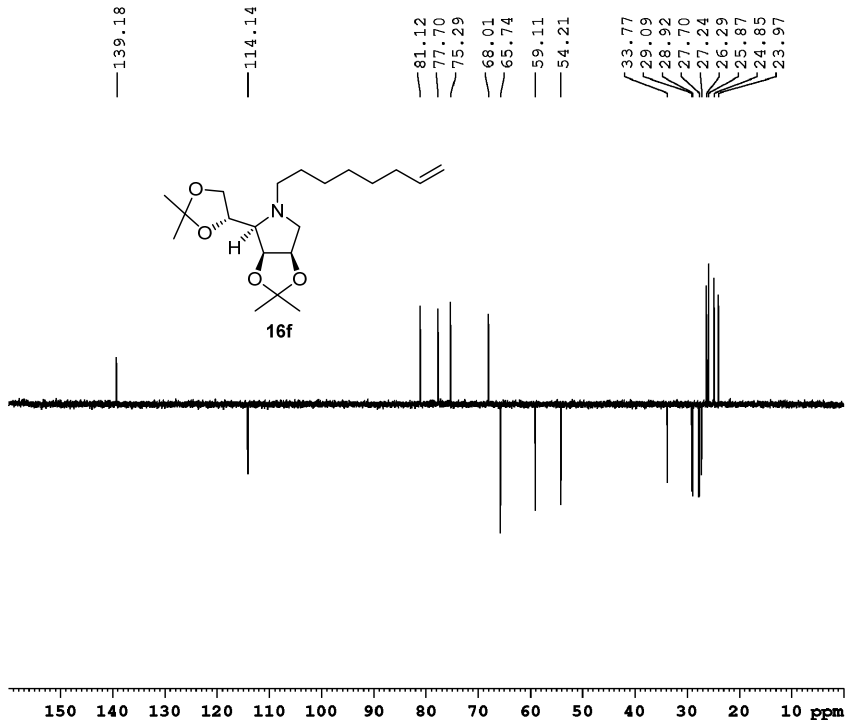
F2 - Acquisition Parameters
Date_     20160907
Time      19.54
INSTRUM   spect
PROBHD    5 mm PABBO 5B/
PULPROG   deptap136
TD         65536
SOLVENT   D2O
NS         256
DS         4
SWH        16129.032 Hz
FIDRES    0.246110 Hz
AQ         2.0316160 sec
RG         206.33
SQ         31.300 usec
DE         6.80 usec
TE         300.2 K
CQF2      145.000000
D1         2.0000000 sec
d2         0.00344828 sec
d12        0.0002000 sec
DELTA     0.0001273 sec
TD0        1
SFO1      100.6484788 MHz
NUC1       13C
P1         10.00 usec
P13        2000.00 usec
P180       0 W
PLM1      54.0000000 W
SFO(M) [5] Csp50omp-4
SFOALD [5] 0.500
SFOFFAS [0] 8.25059986 W
SFO2      400.2412800 MHz
NUC2       1H
CQDFPR [2] waltz16
P3         15.30 usec
P4         30.60 usec
PCPD2     80.00 usec
PLM2      12.0000000 W
PLM12     0.34680000 W

F2 - Processing parameters
SI         32768
SF         100.6404280 MHz
WDW        EM
SSB        0
GB         0
PC         1.40
  
```



Compound 16f



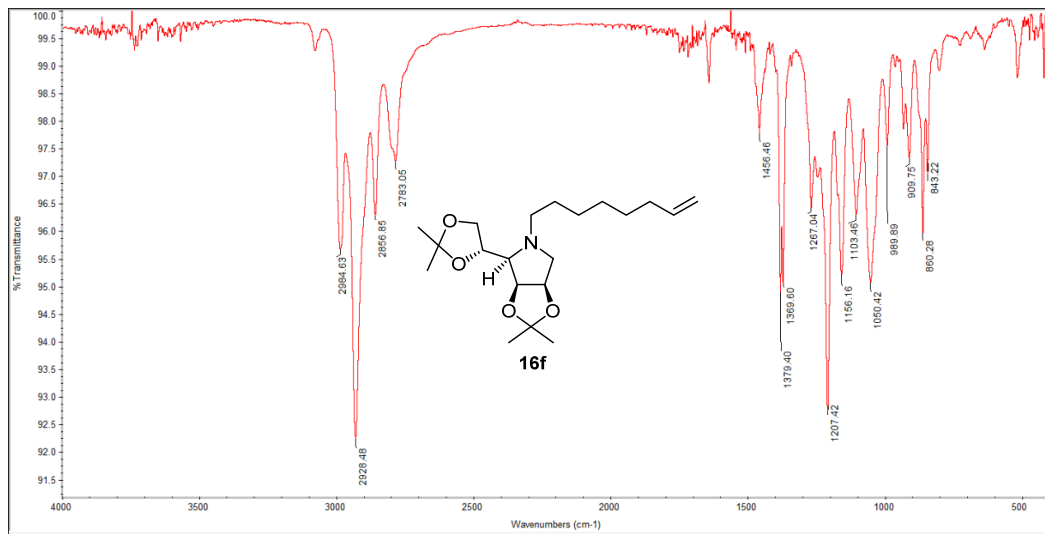


```

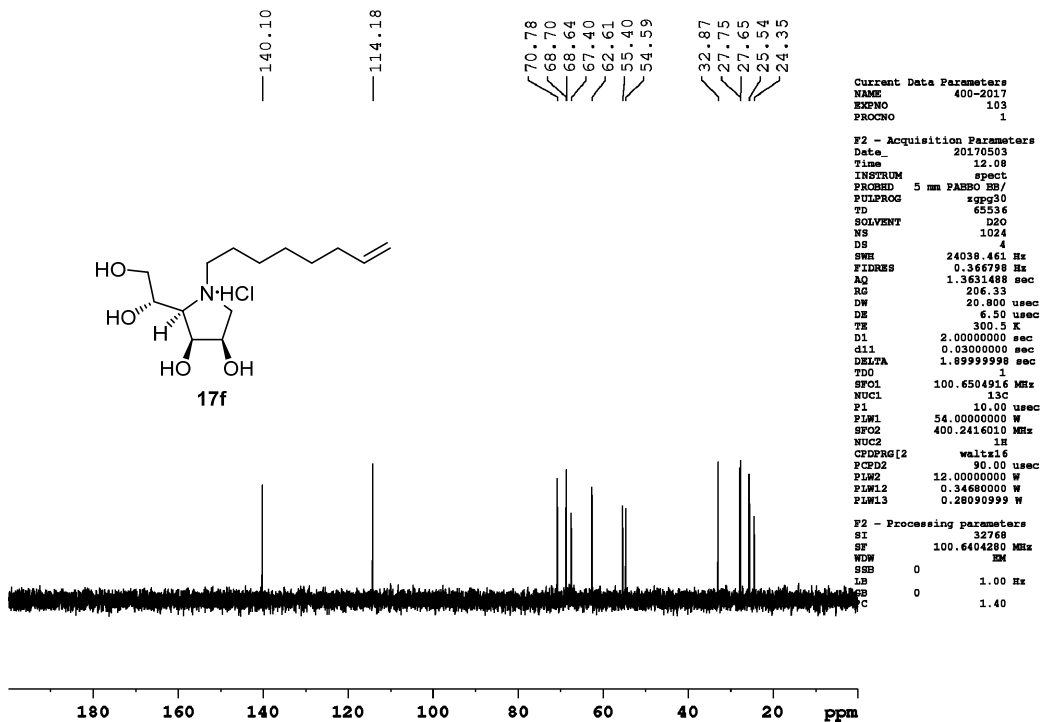
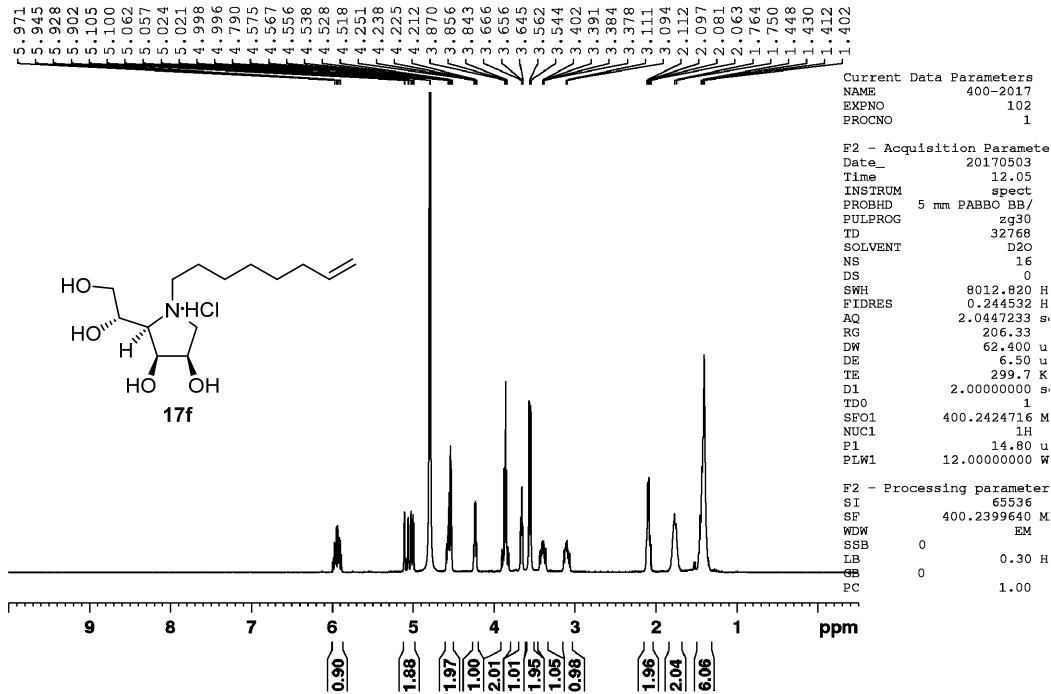
Current Data Parameters
NAME      400-2017
EXPNO    42
PROCNO    1

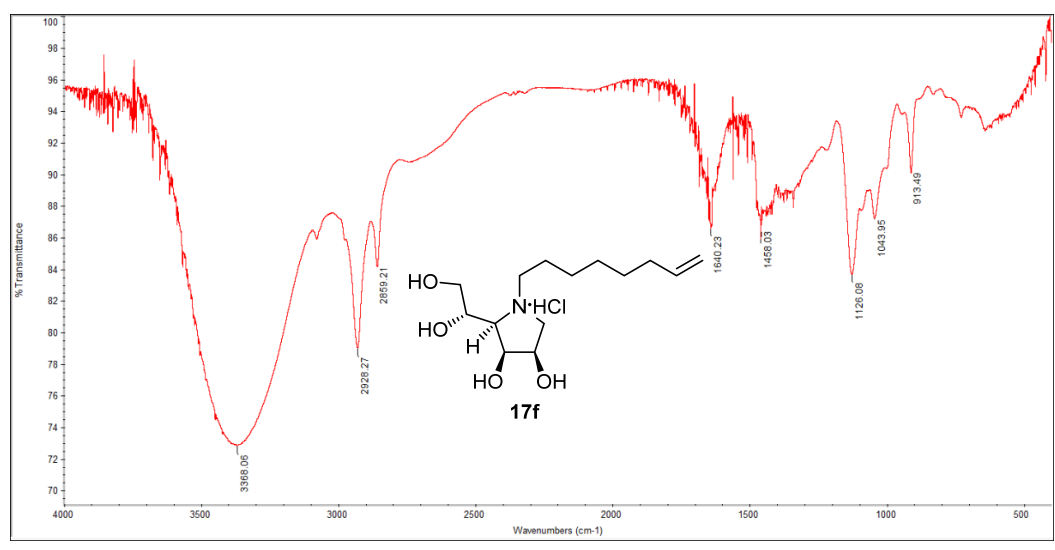
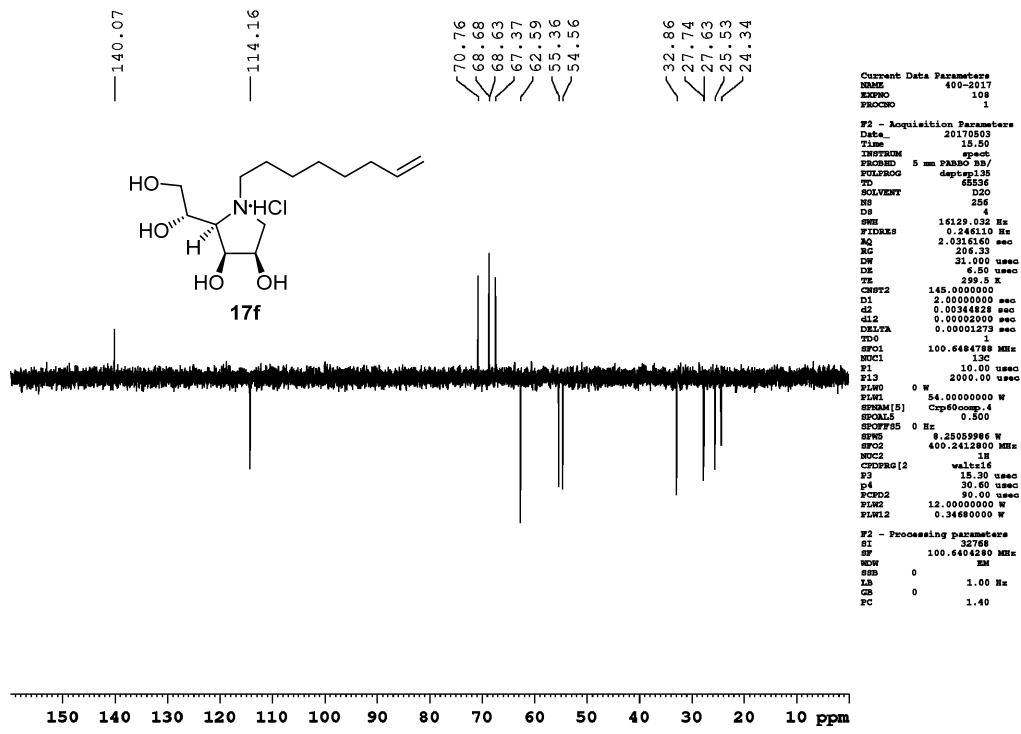
F2 - Acquisition Parameters
Date_     20161219
Time      19.22
INSTRUM   spect
PROBHD    5 mm F300 SBL
PULPROG   zgpg30
TD         65536
SOLVENT   CDCl3
NS         116
DS         4
SWH        16129.032 Hz
FIDRES    0.246110 Hz
AQ         2.0316160 sec
RG         206.33
SQ         31.300 usec
DE         6.80 usec
TE         298.3 K
CQF2       145.000000
D1         2.0000000 sec
d2         0.00344828 sec
d12        0.0002000 sec
DELTA     0.0001273 sec
TD0        1
SFO1       100.6484788 MHz
NUC1       13C
P1         10.00 usec
P12        2000.00 usec
SFO        0 W
PLM1       54.0000000 W
SFO5(1)    Exp50comp 4
SFOAL5     0.500
SFOFF5     0 Hz
SWS        8.25059986 W
SFO2       400.2412800 MHz
NUC2       1H
CQFPR2(2)  waltz16
P3         15.30 usec
P4         30.60 usec
PCPD2      90.00 usec
PLM2       12.0000000 W
PLM12      0.34680000 W

F2 - Processing parameters
SI         32768
SF         100.6404280 MHz
WDW        EM
SSB        0
GB         1.00 Hz
PC         1.40
  
```

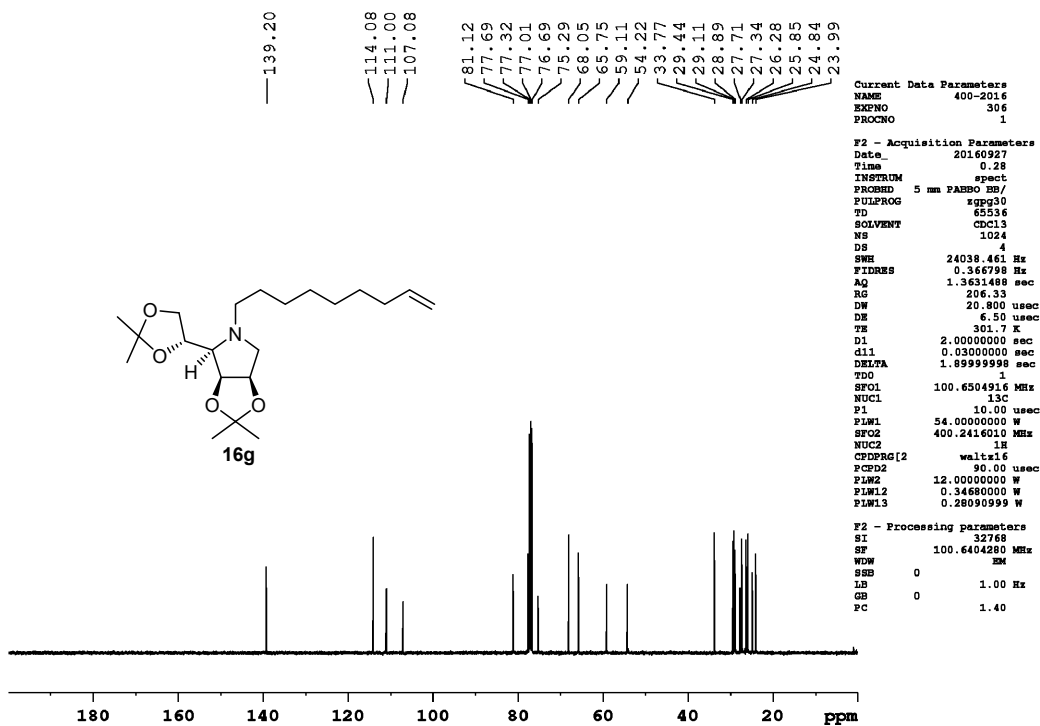
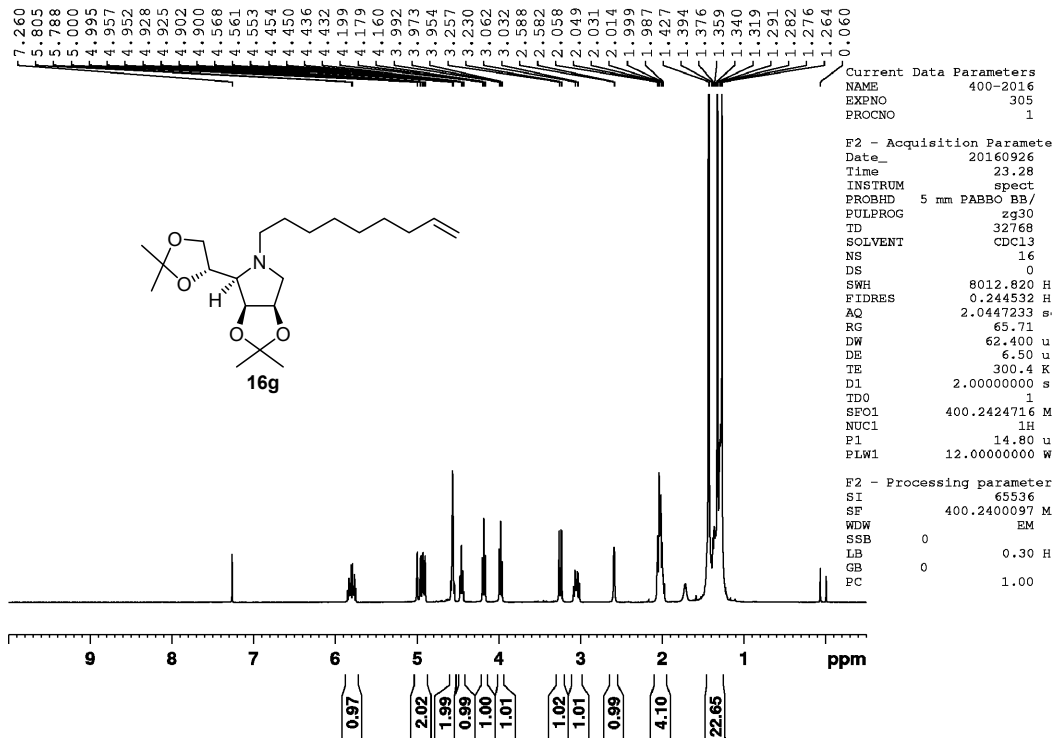


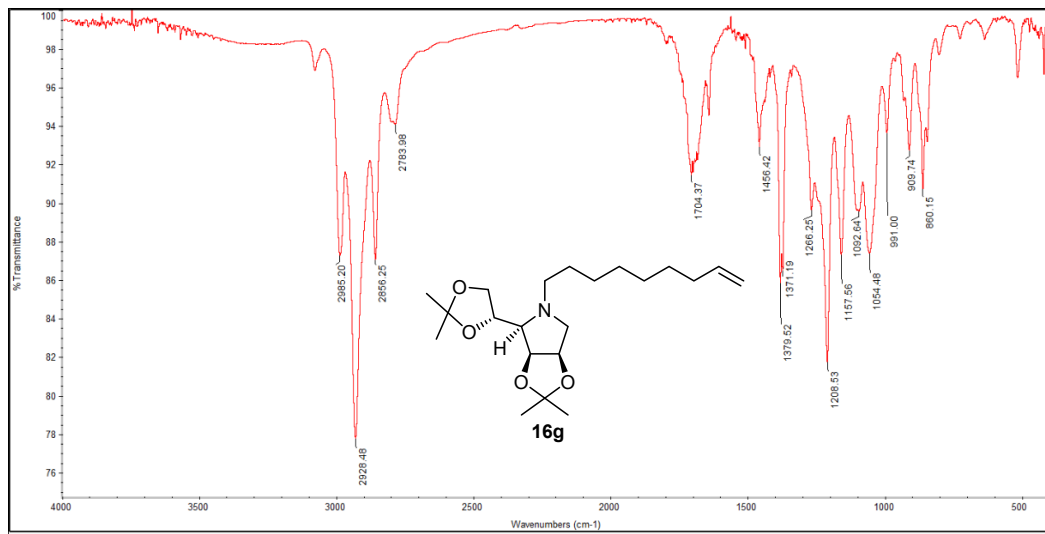
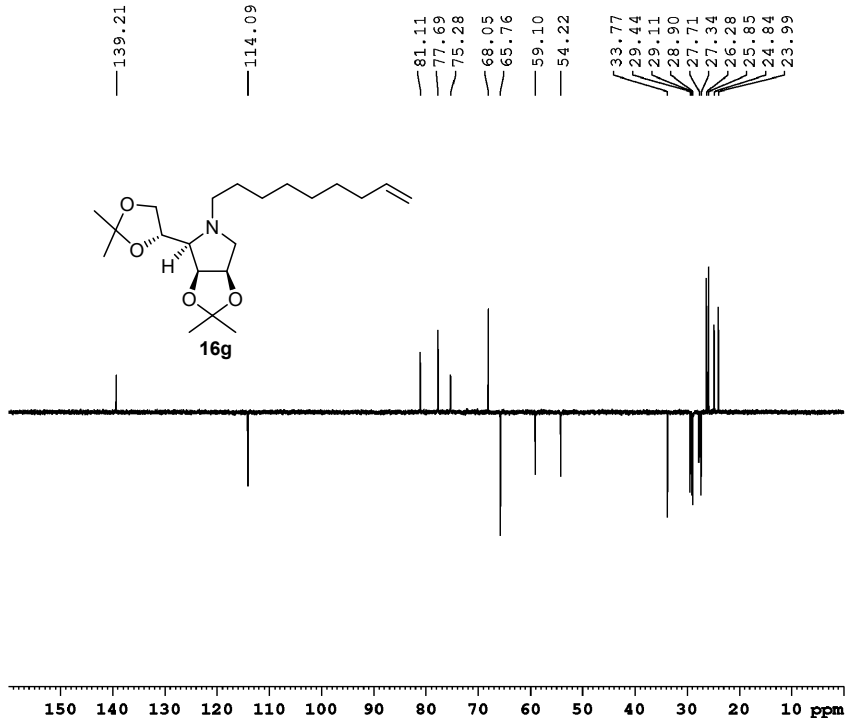
Compound 17f:





Compound 16g:





Compound 17g:

