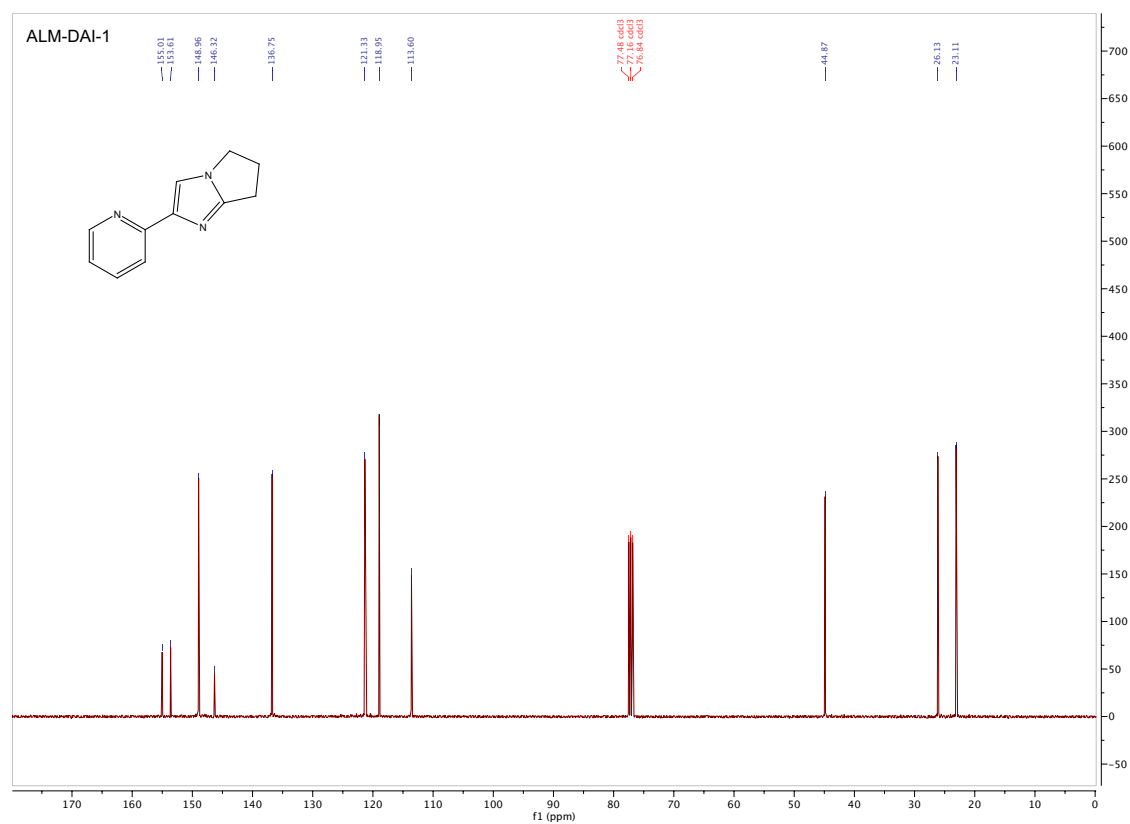
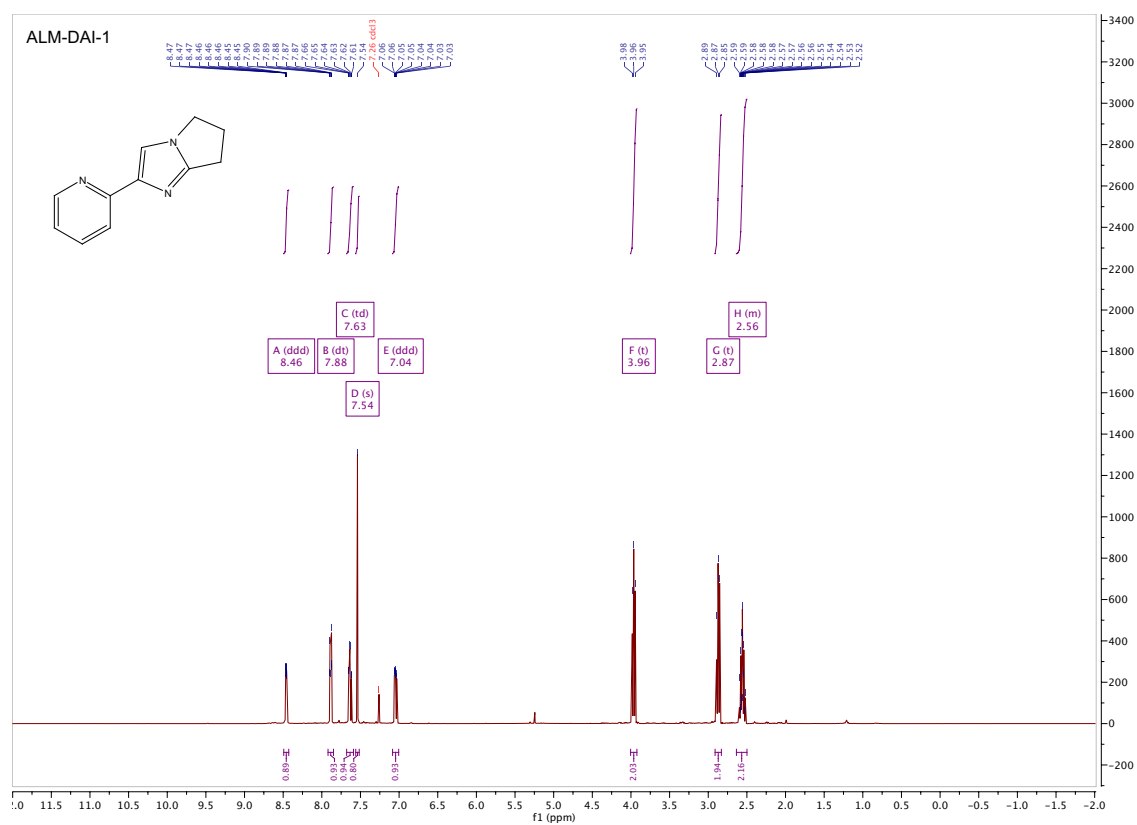
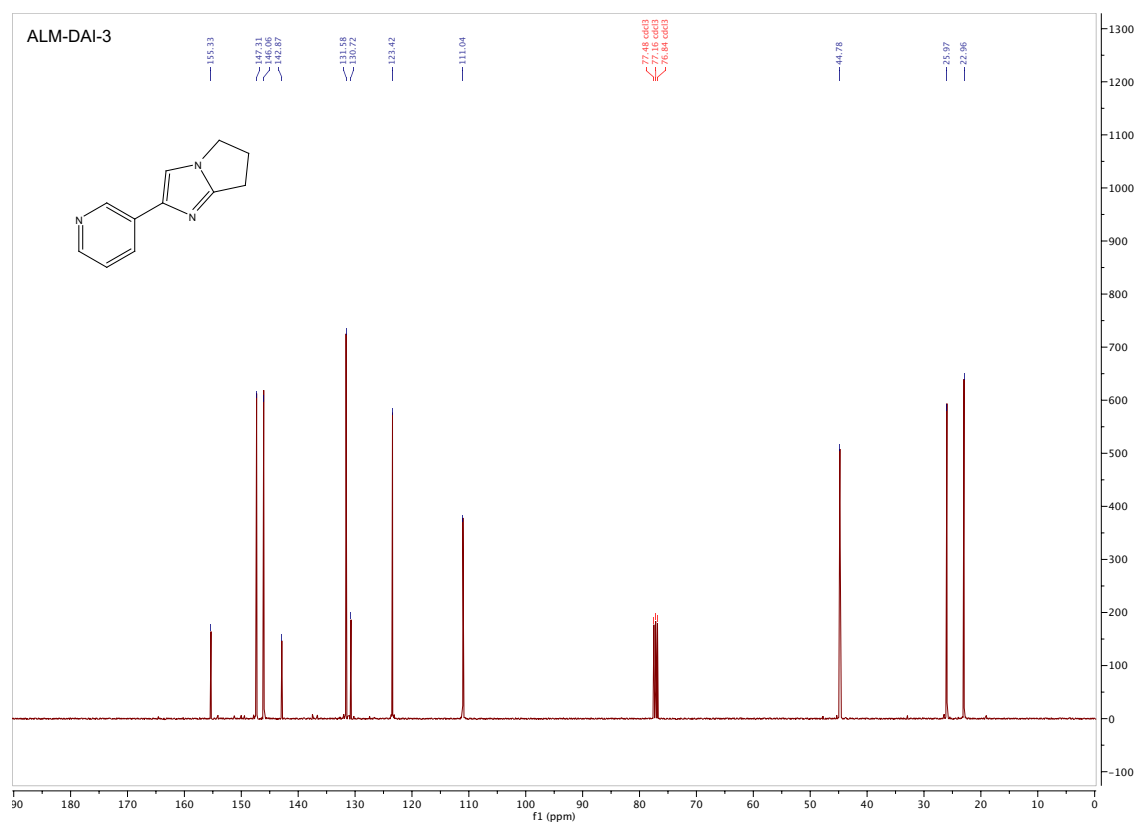
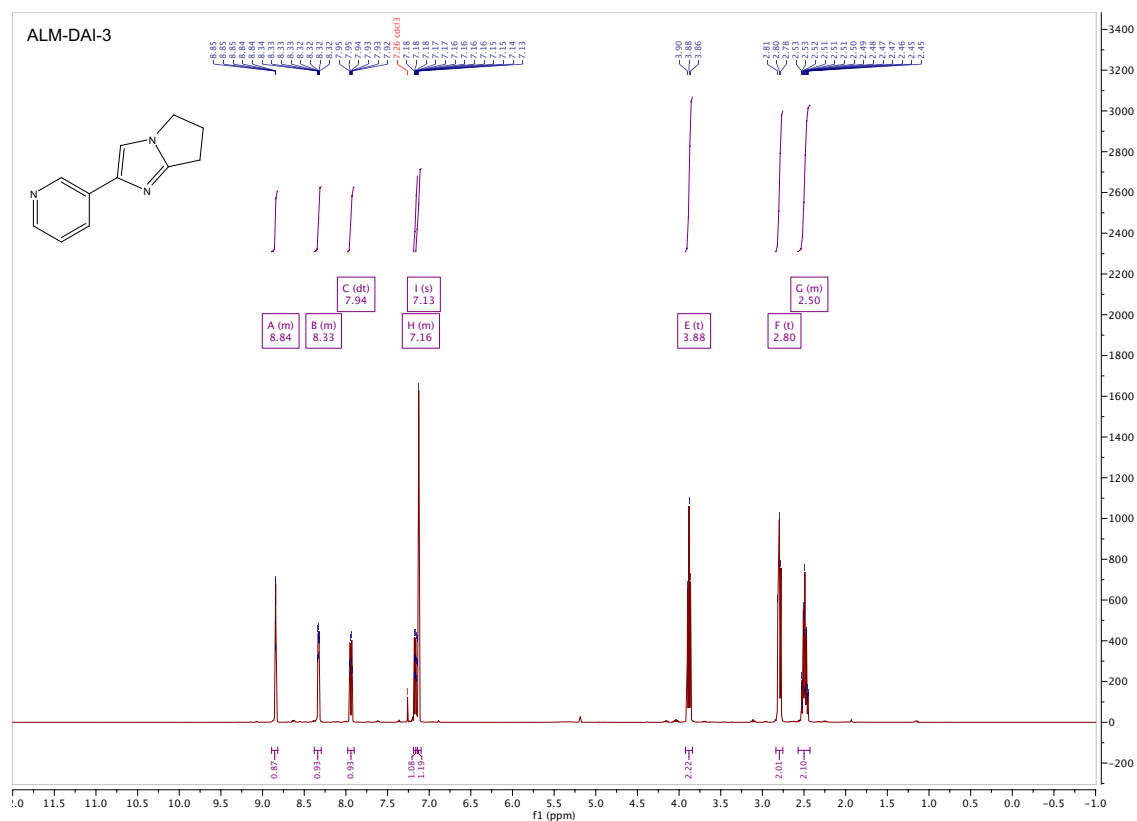


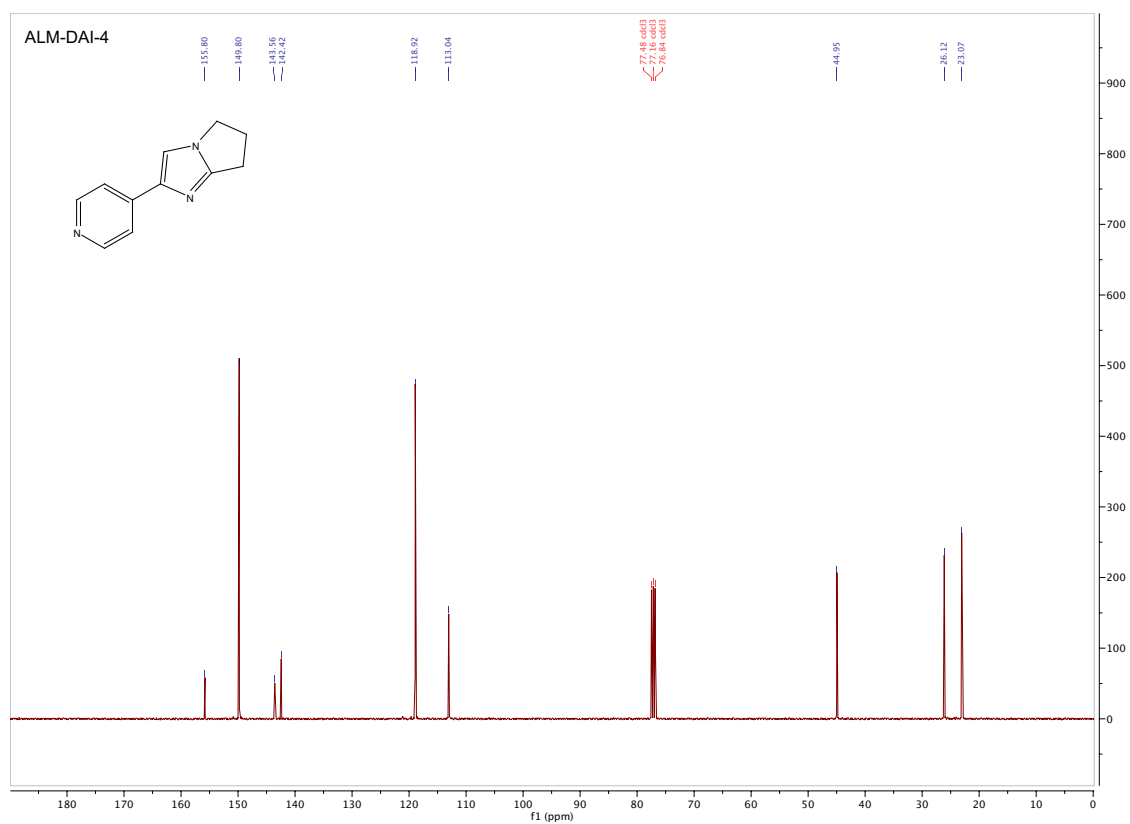
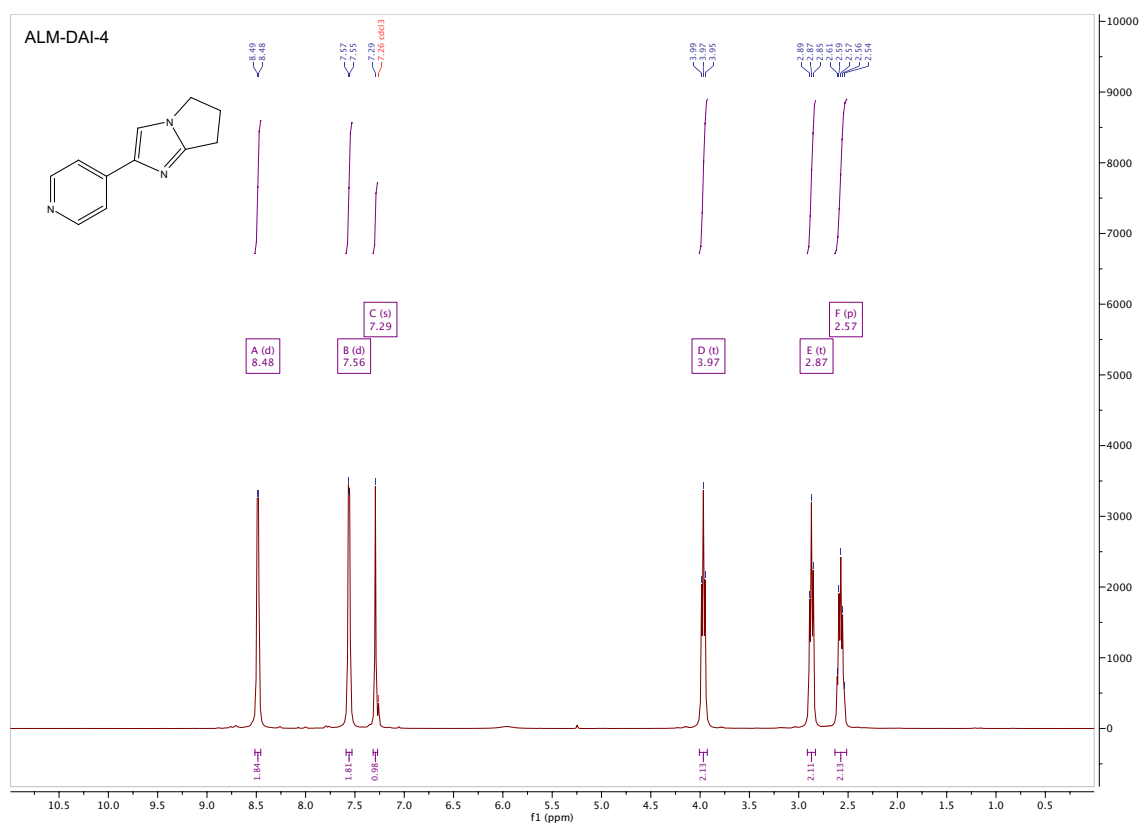
2-(pyridin-2-yl)-6,7-dihydro-5H-pyrrolo[1,2-a]imidazole (ALM-DAI-1).



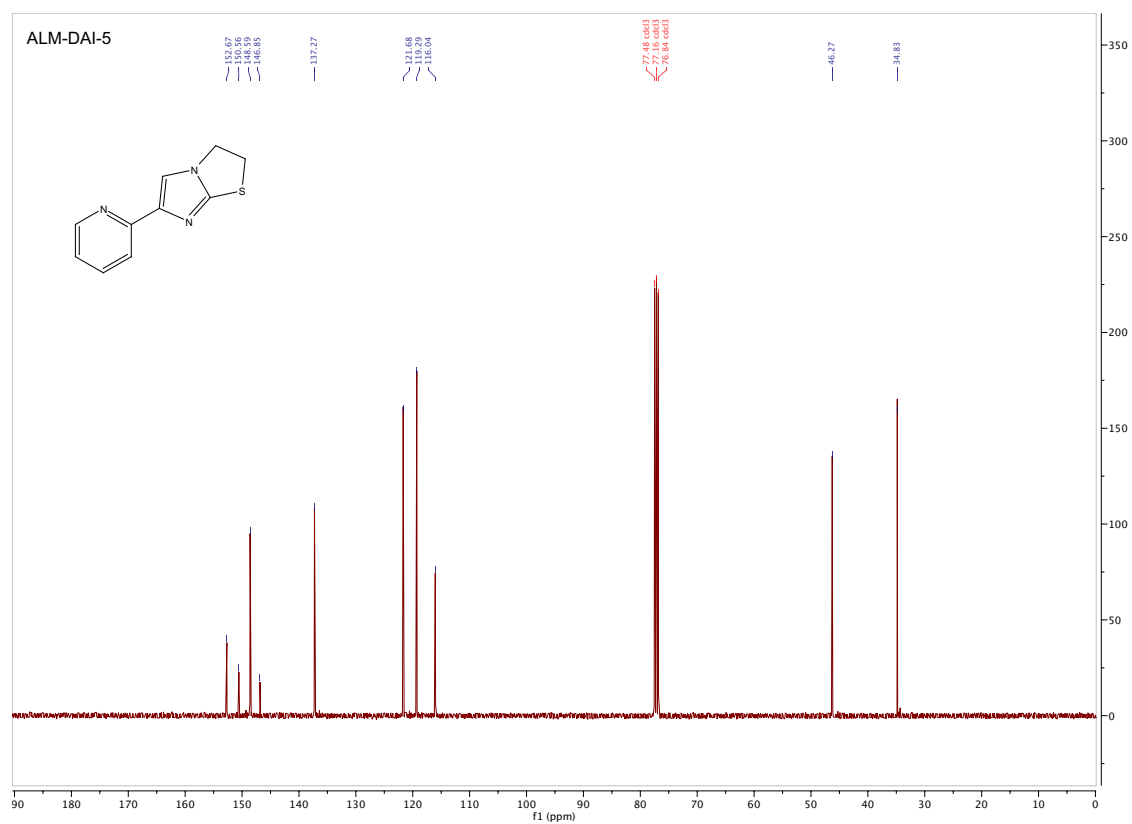
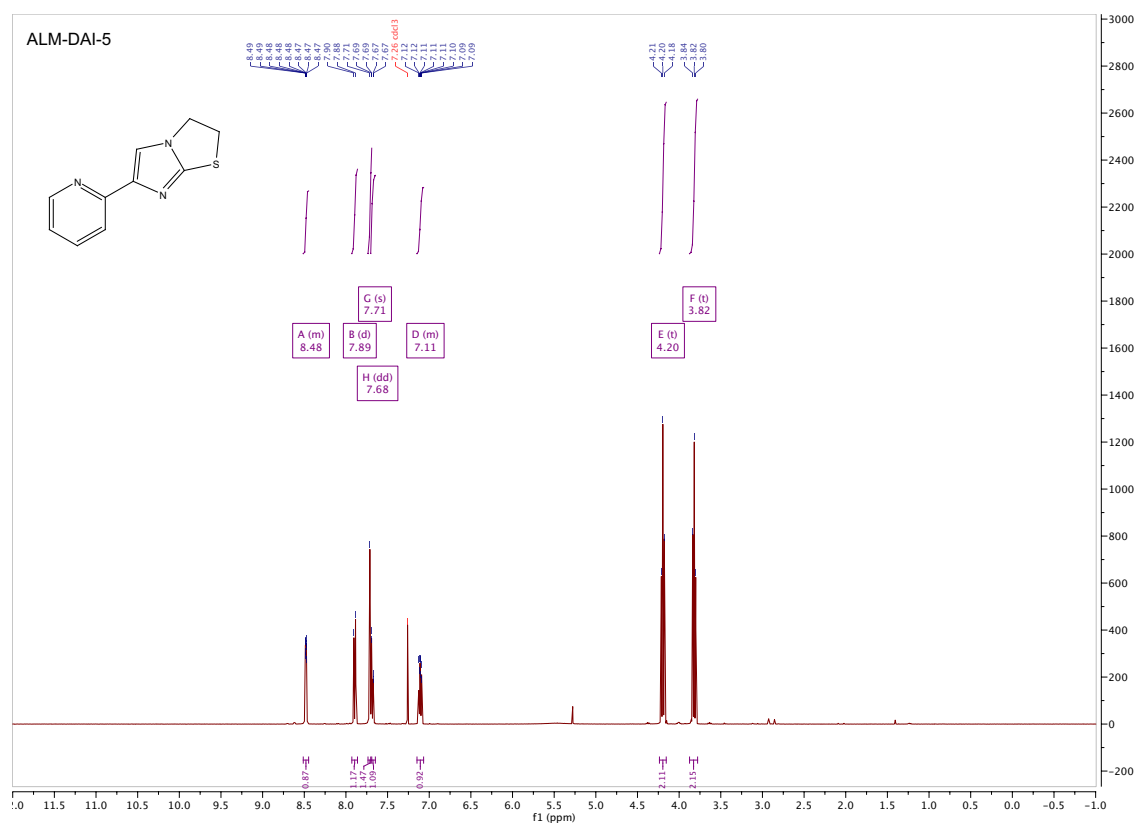
2-(pyridin-3-yl)-6,7-dihydro-5H-pyrrolo[1,2-a]imidazole (ALM-DAI-3).



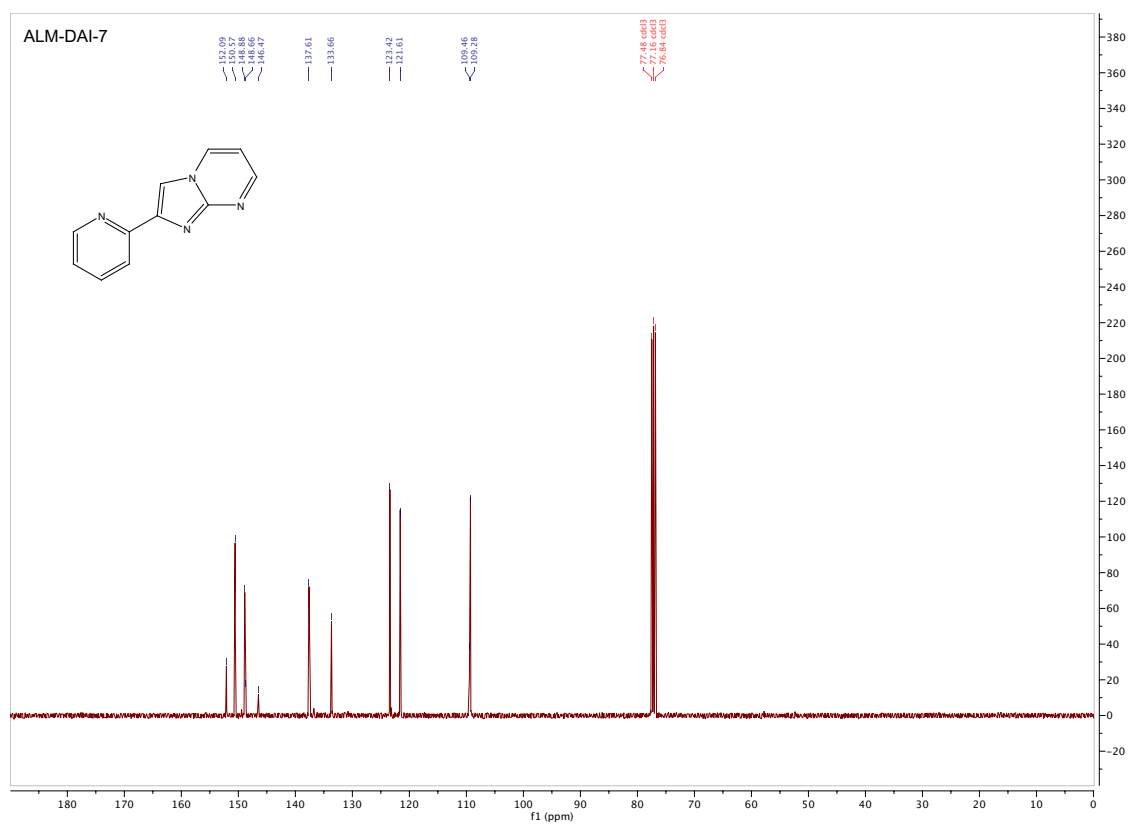
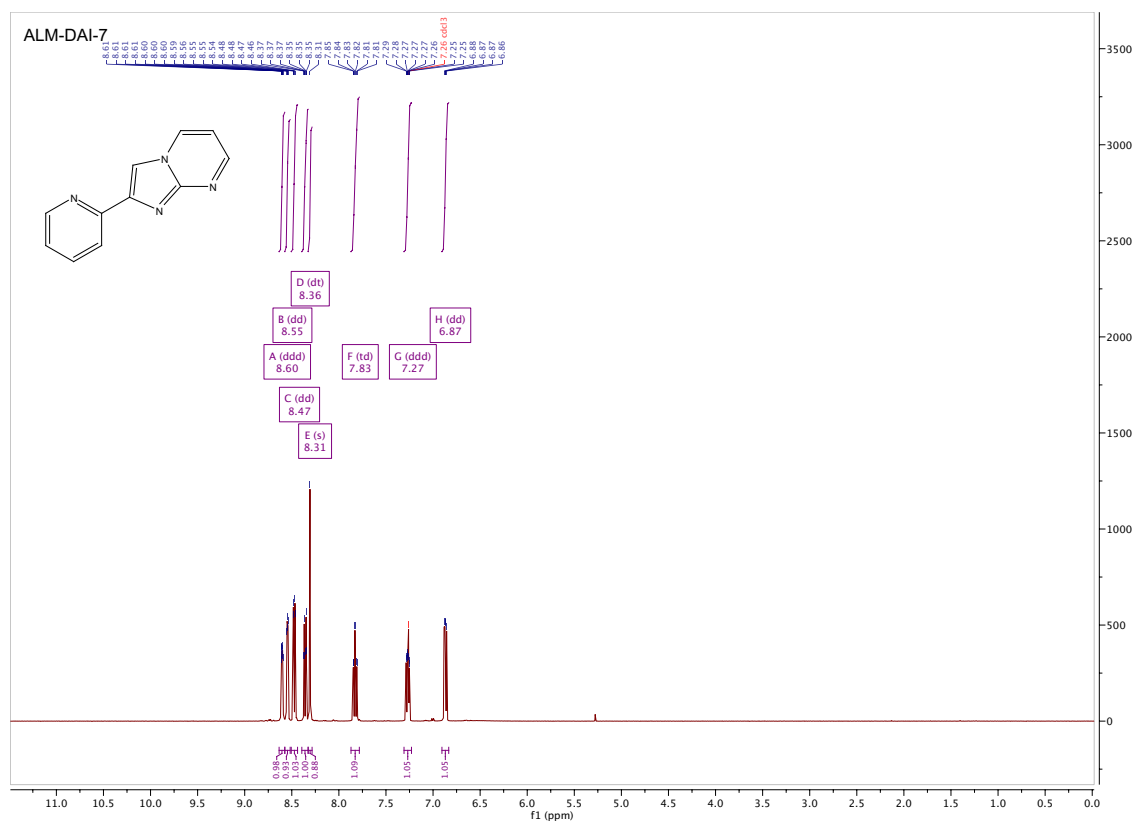
2-(pyridin-4-yl)-6,7-dihydro-5H-pyrrolo[1,2-a]imidazole (ALM-DAI-4).



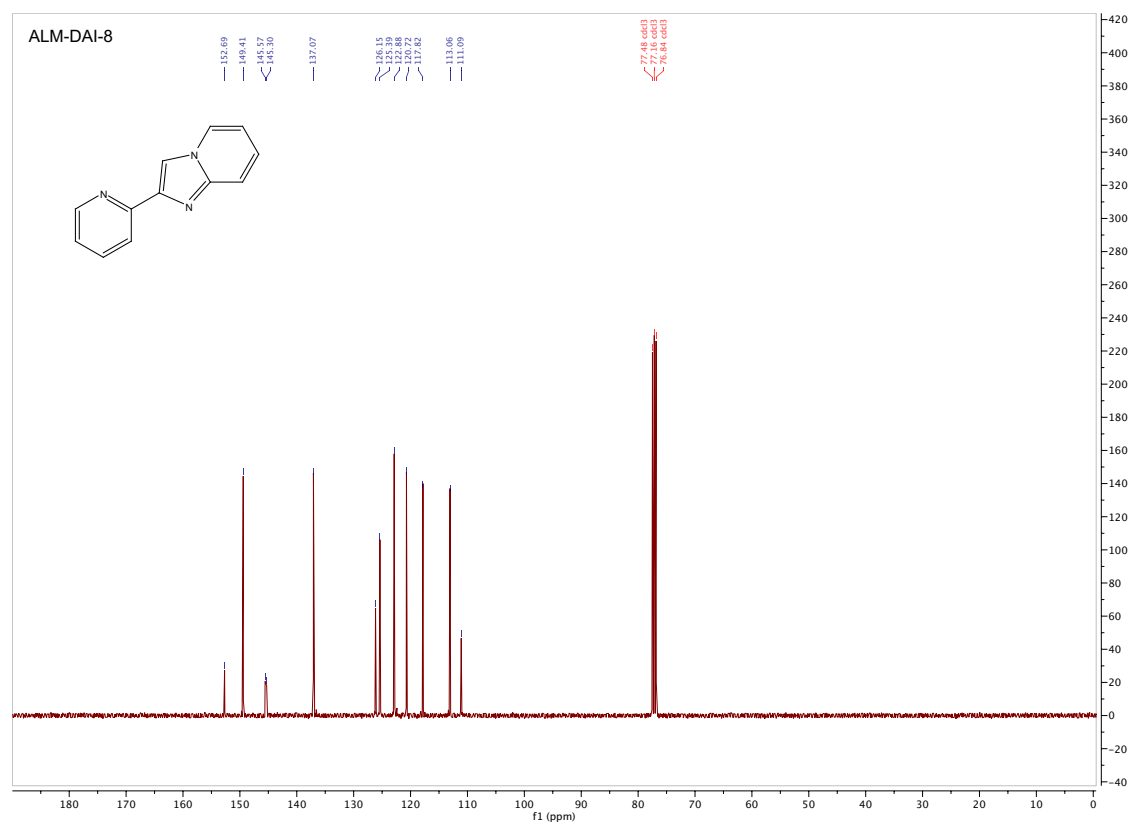
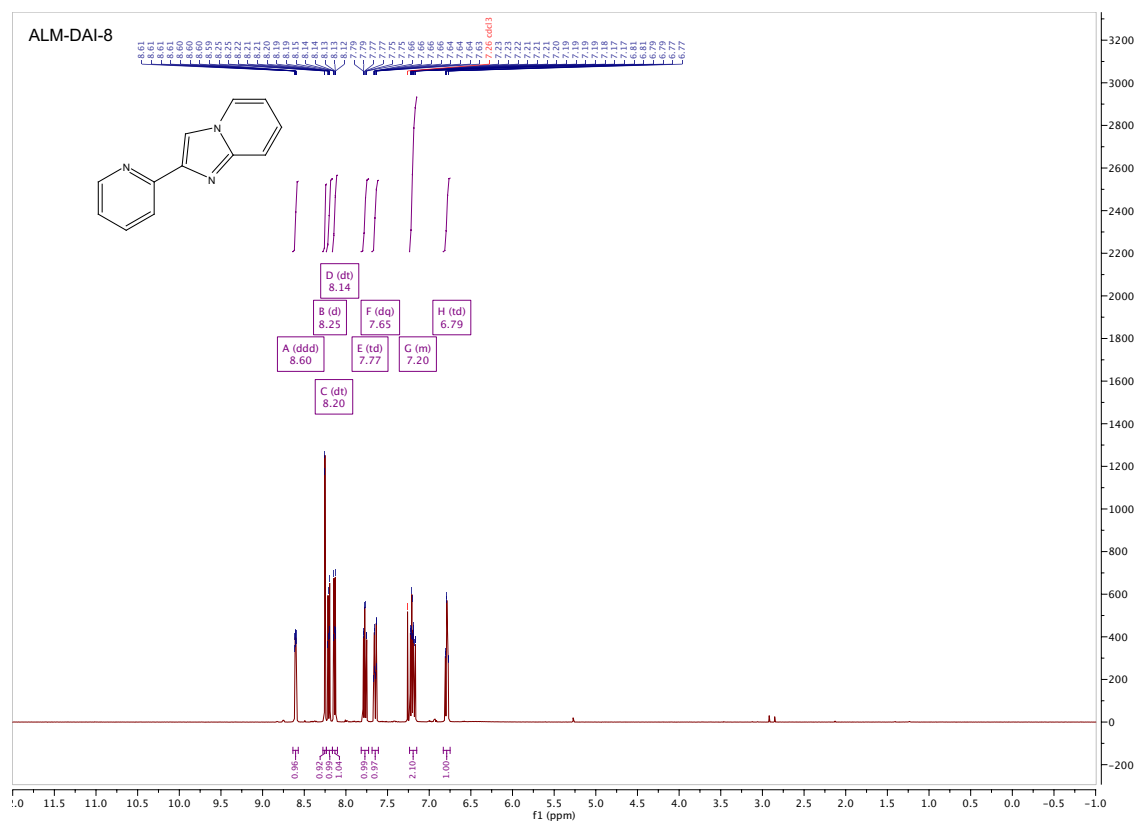
6-(pyridin-2-yl)-2,3-dihydroimidazo[2,1-*b*]thiazole (ALM-DAI-5).



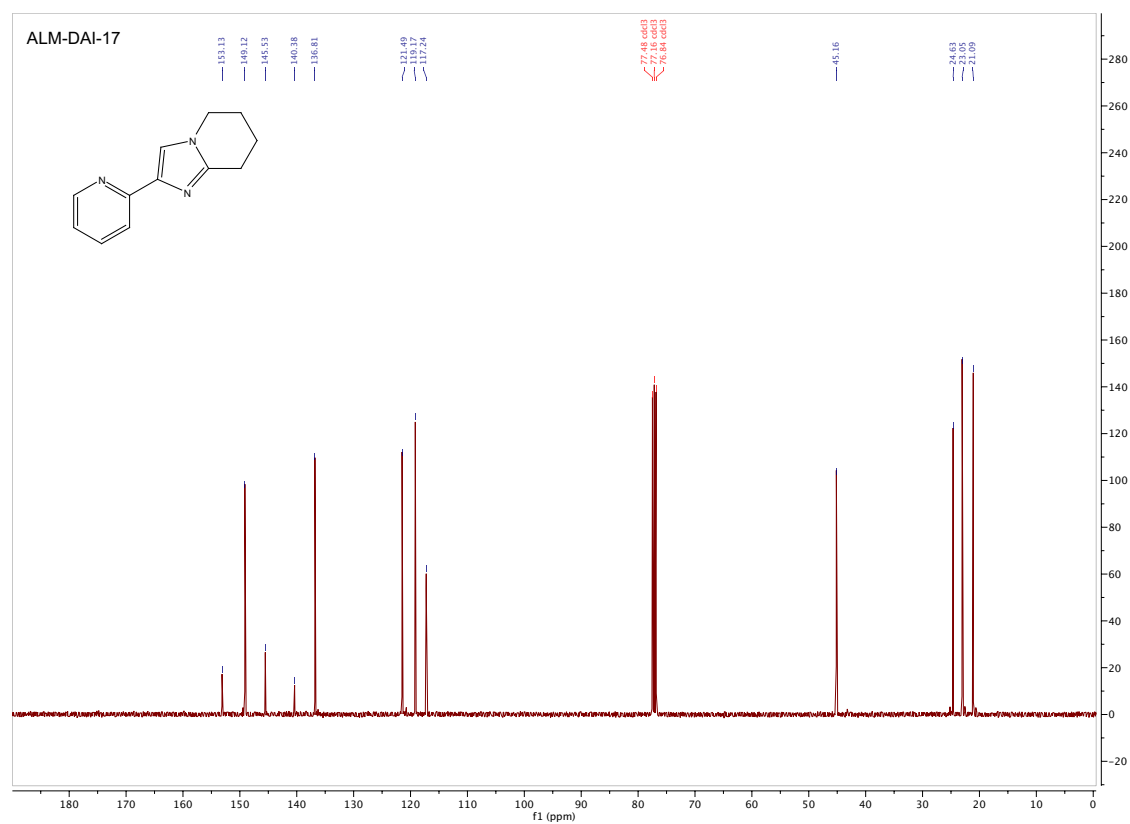
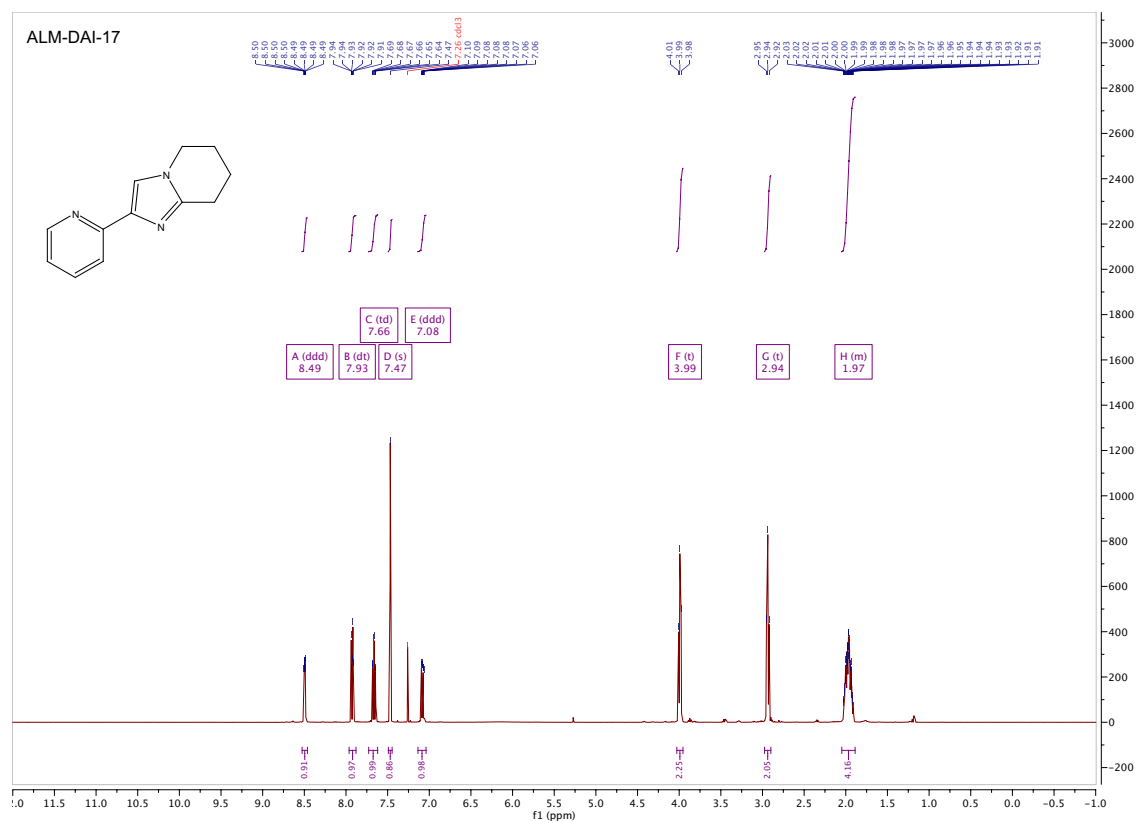
2-(pyridin-2-yl)imidazo[1,2-a]pyrimidine (ALM-DAI-7).



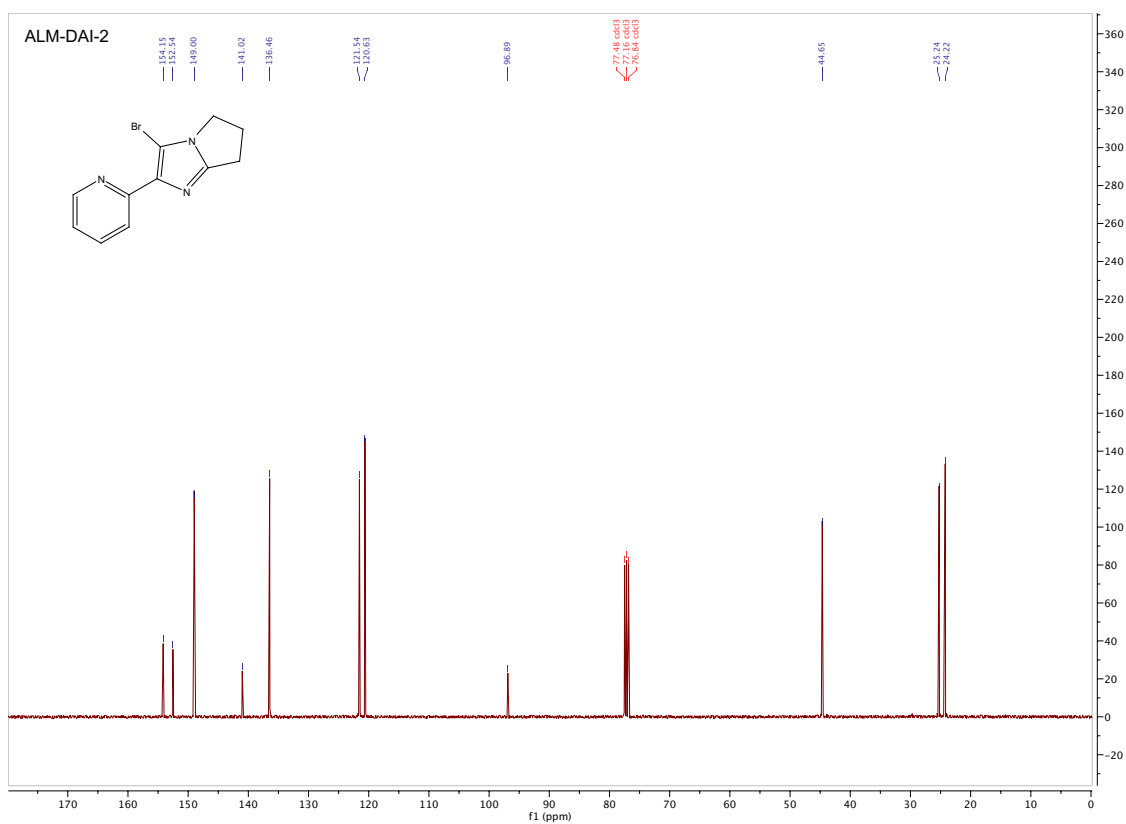
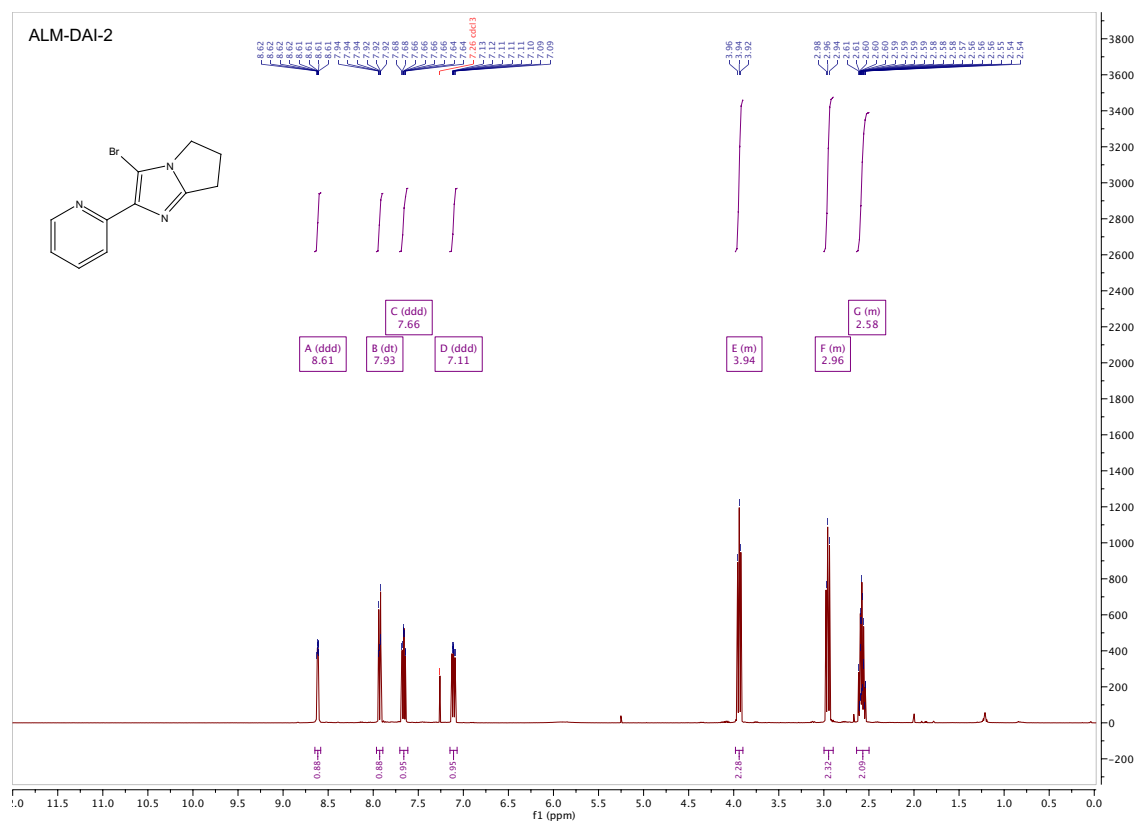
2-(pyridin-2-yl)imidazo[1,2-a]pyridine (ALM-DAI-8).



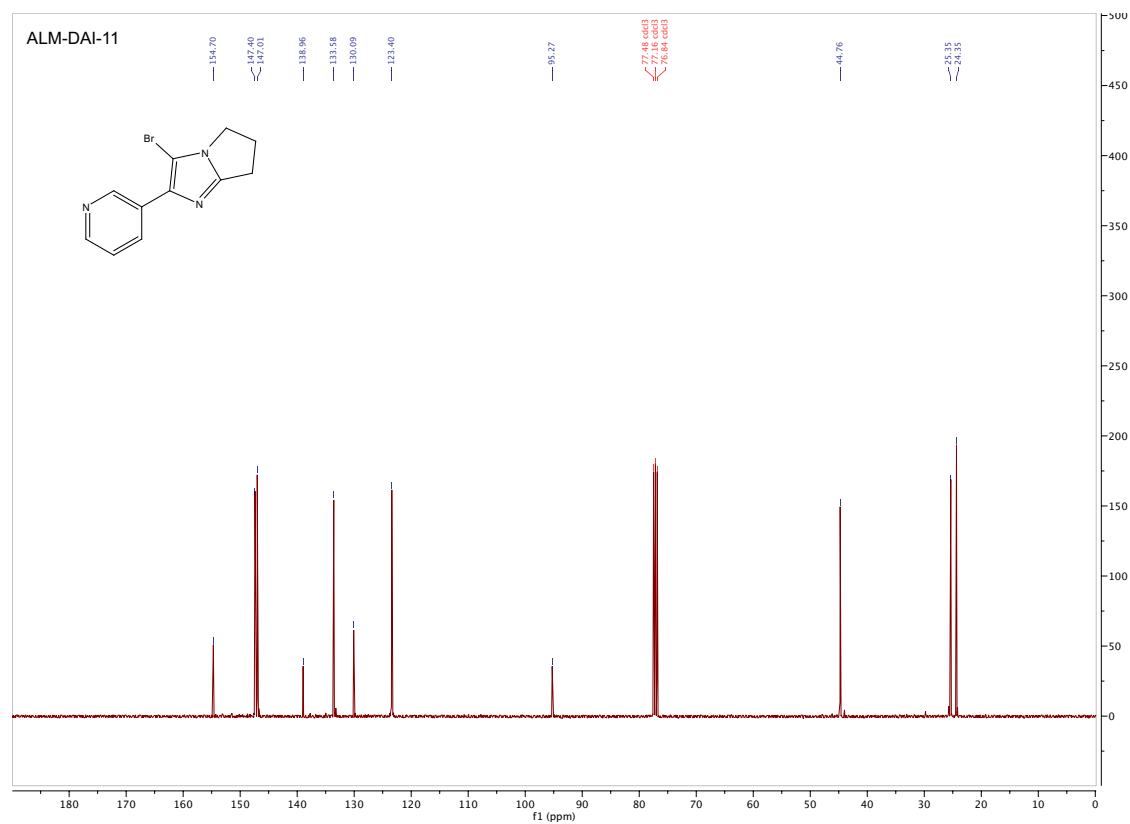
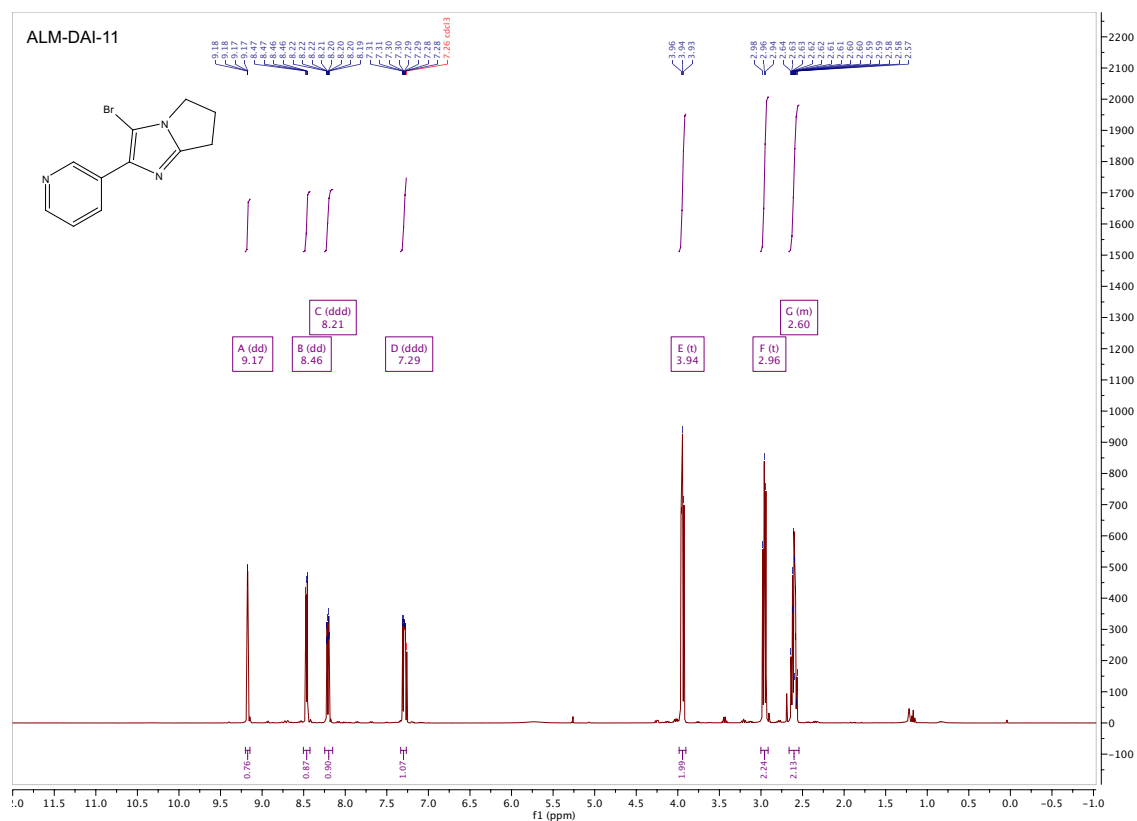
2-(pyridin-2-yl)-5,6,7,8-tetrahydroimidazo[1,2-*a*]pyridine (ALM-DAI-17).



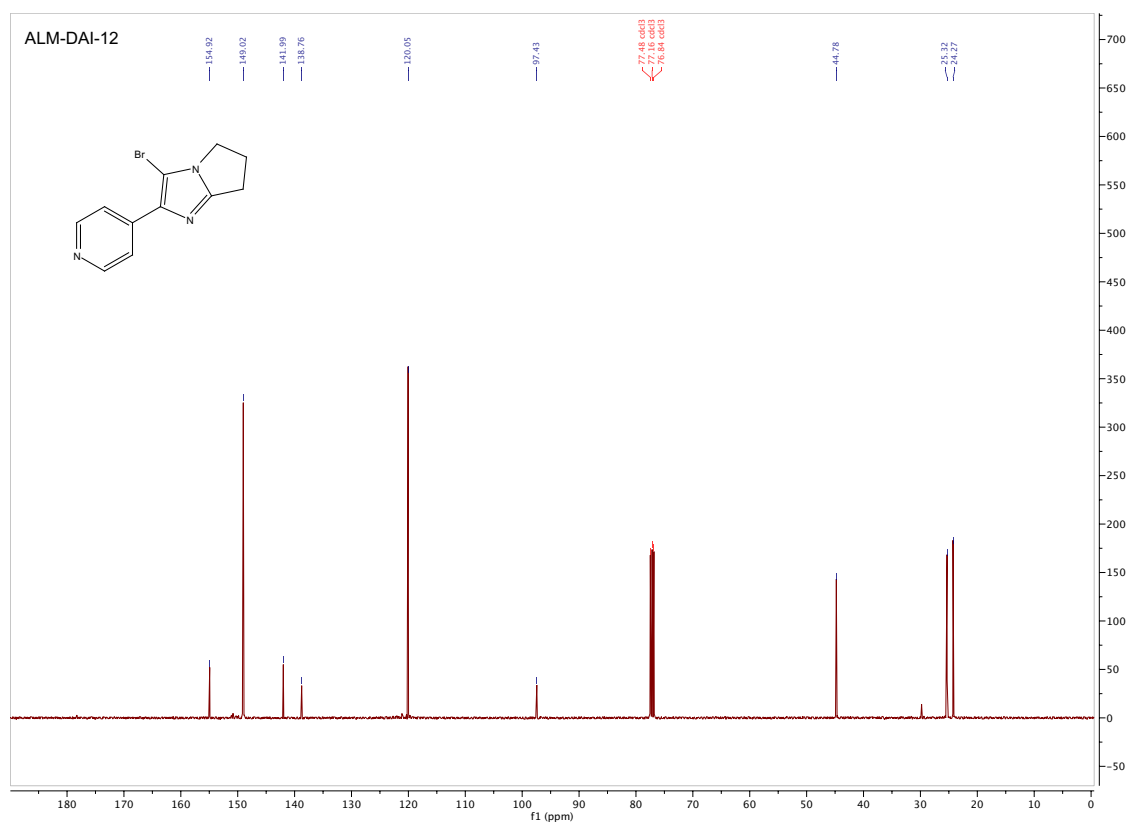
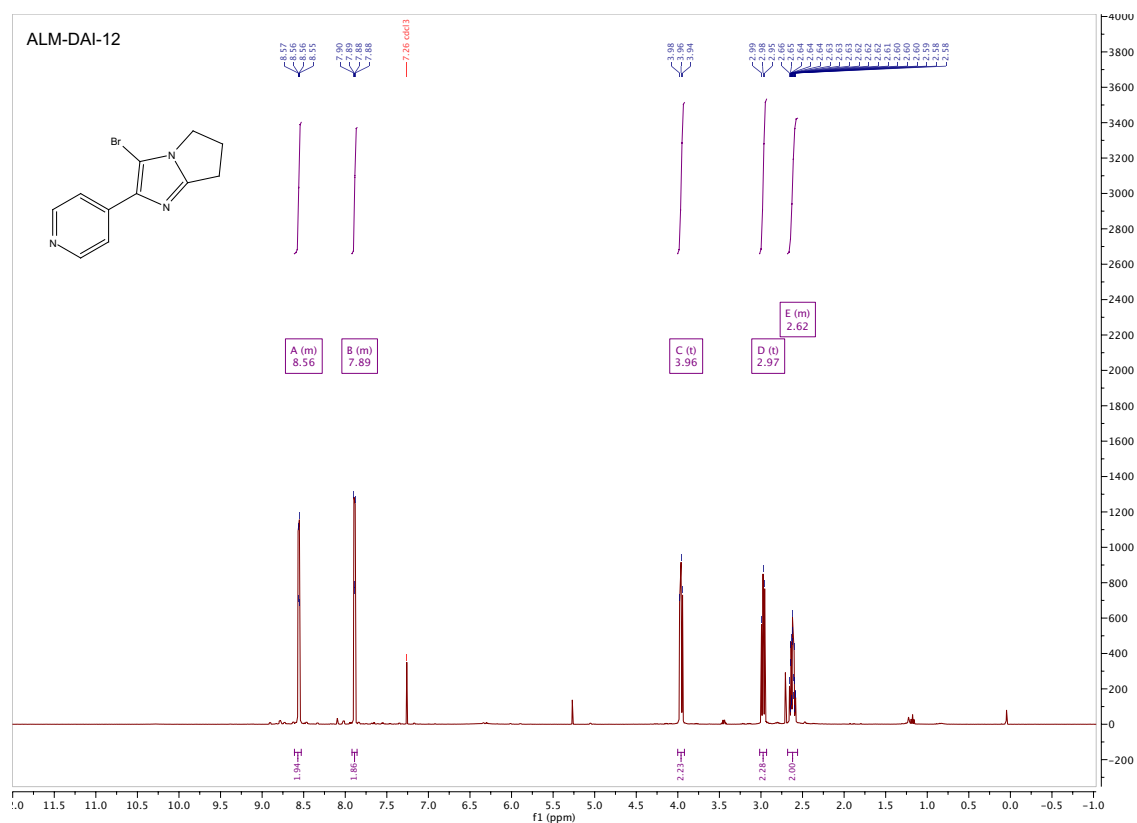
3-bromo-2-(pyridin-2-yl)-6,7-dihydro-5H-pyrrolo[1,2-a]imidazole (ALM-DAI-2).



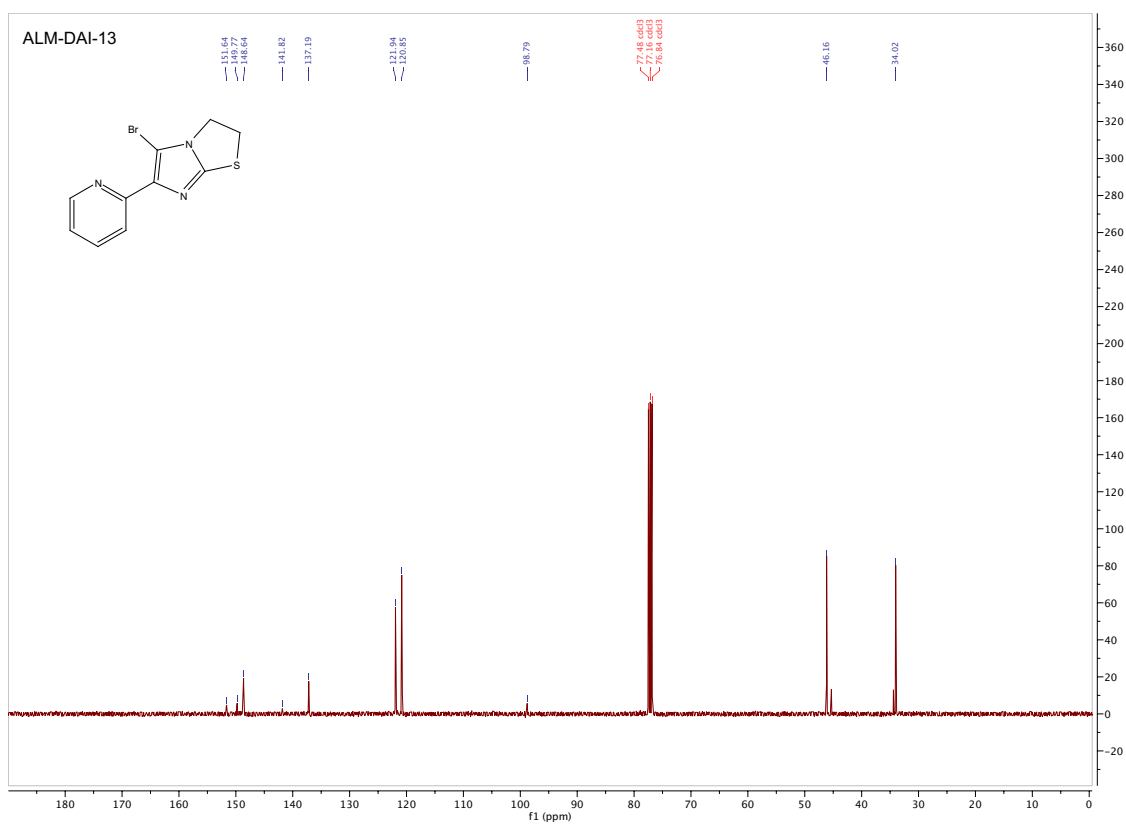
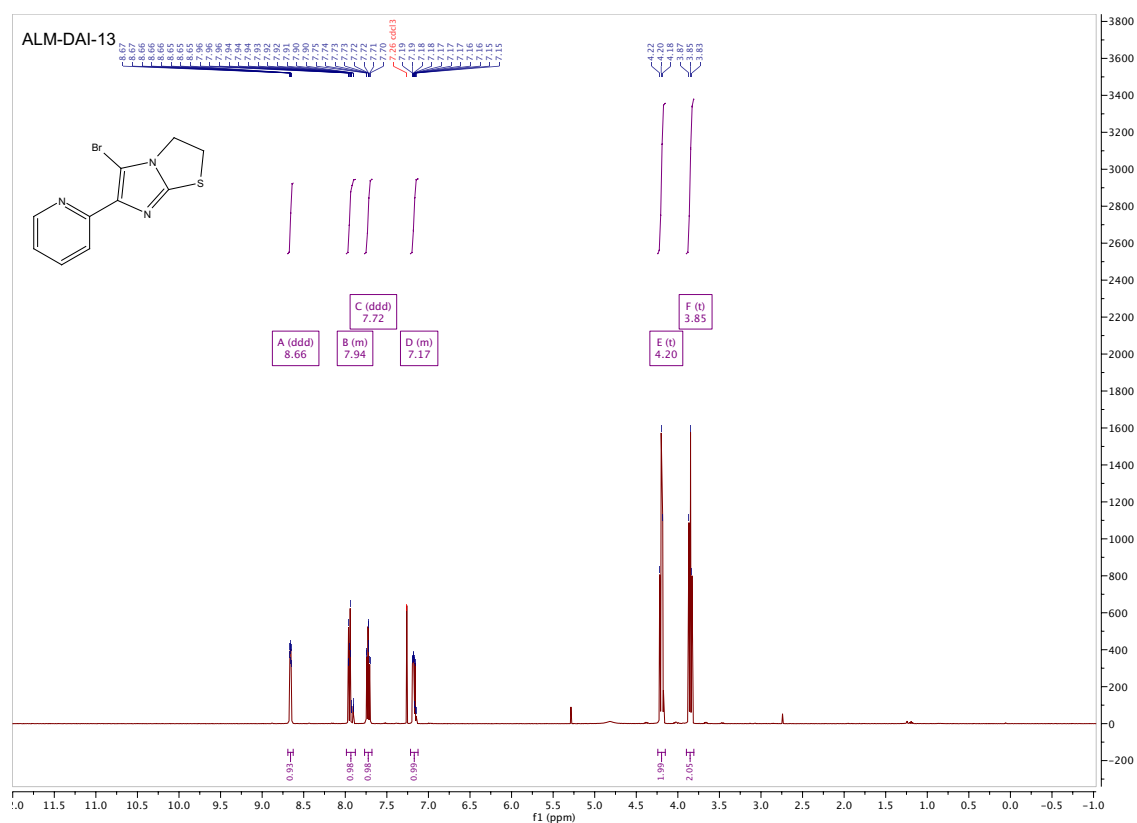
3-bromo-2-(pyridin-3-yl)-6,7-dihydro-5H-pyrrolo[1,2-a]imidazole (ALM-DAI-11).



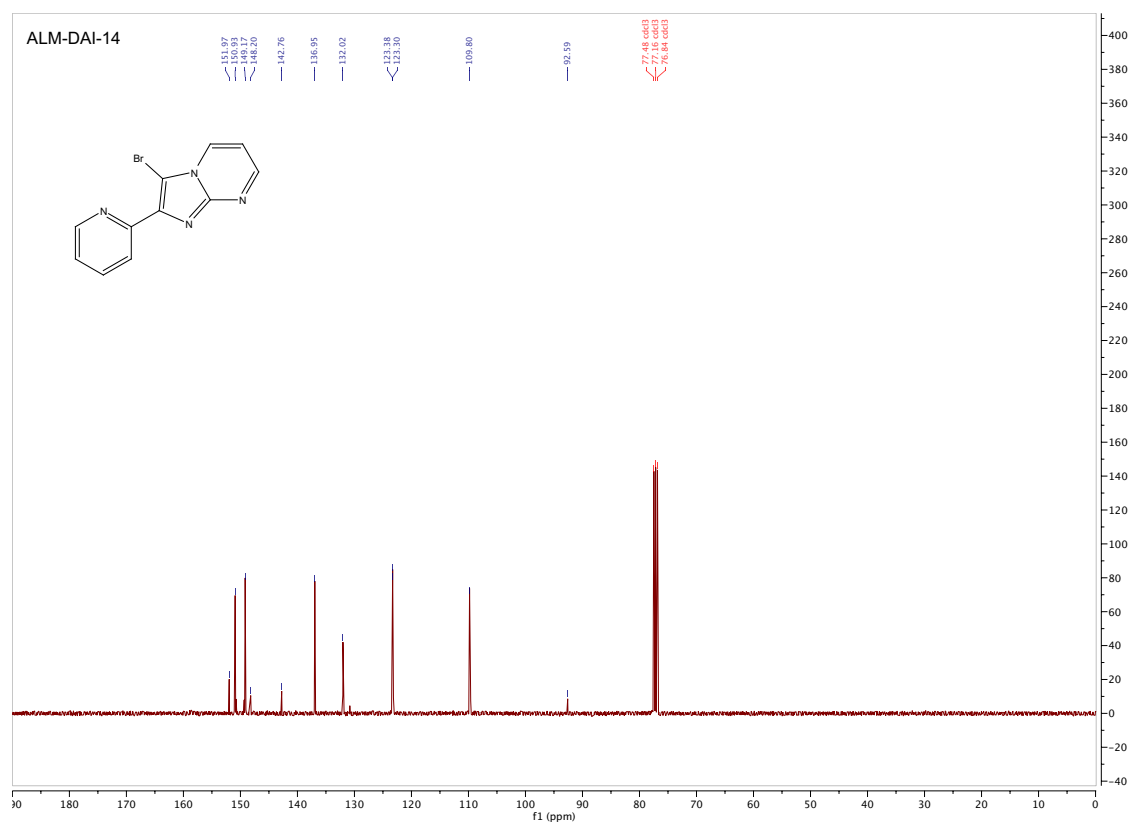
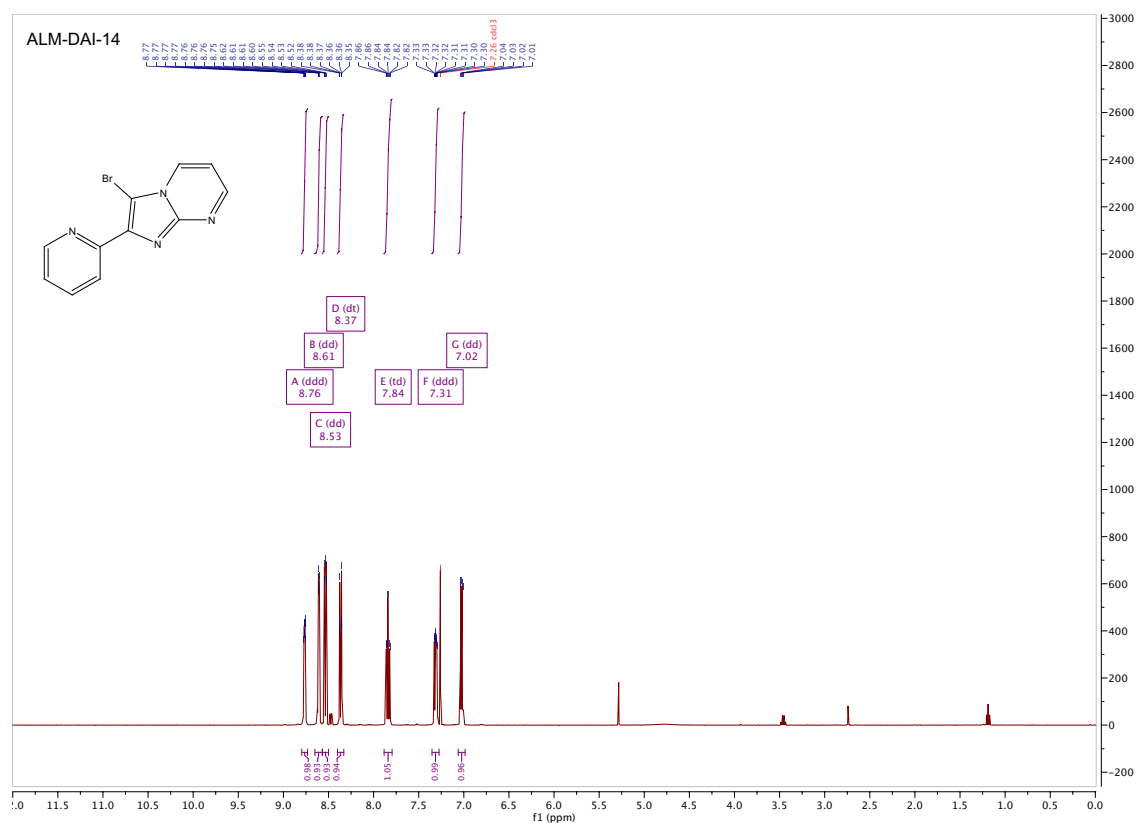
3-bromo-2-(pyridin-4-yl)-6,7-dihydro-5H-pyrrolo[1,2-a]imidazole (ALM-DAI-12).



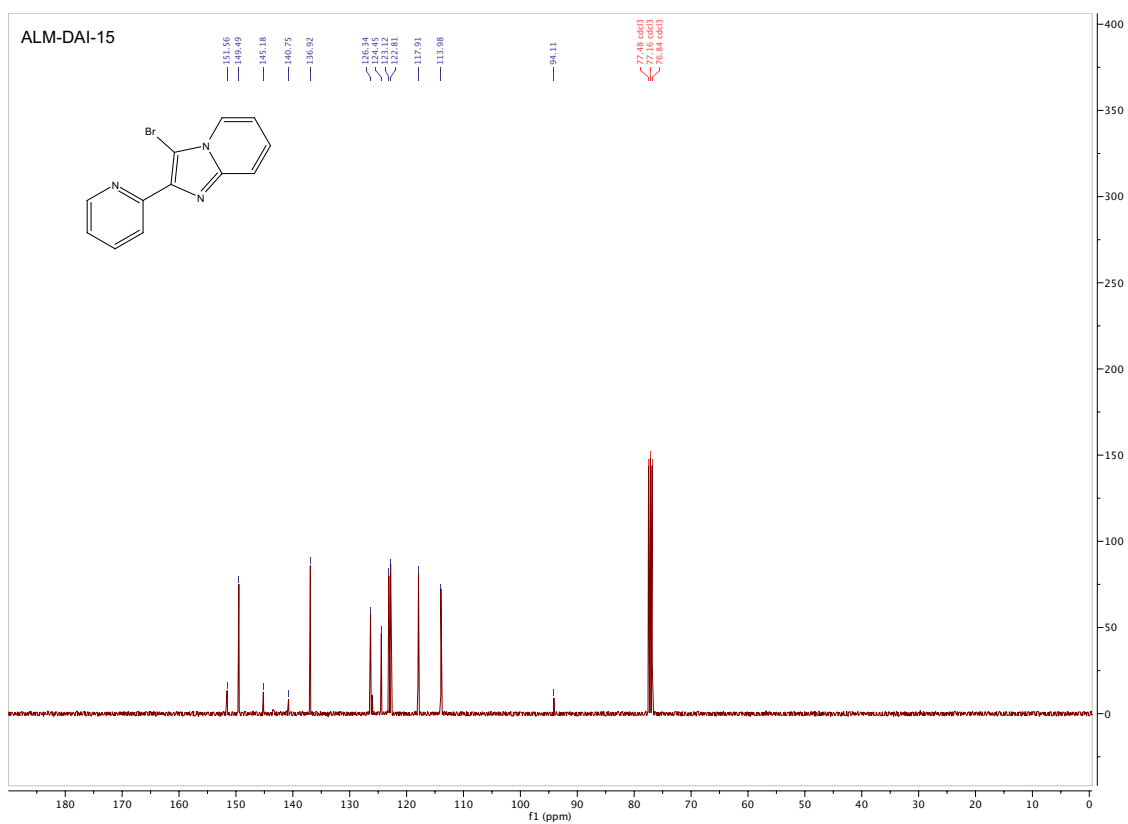
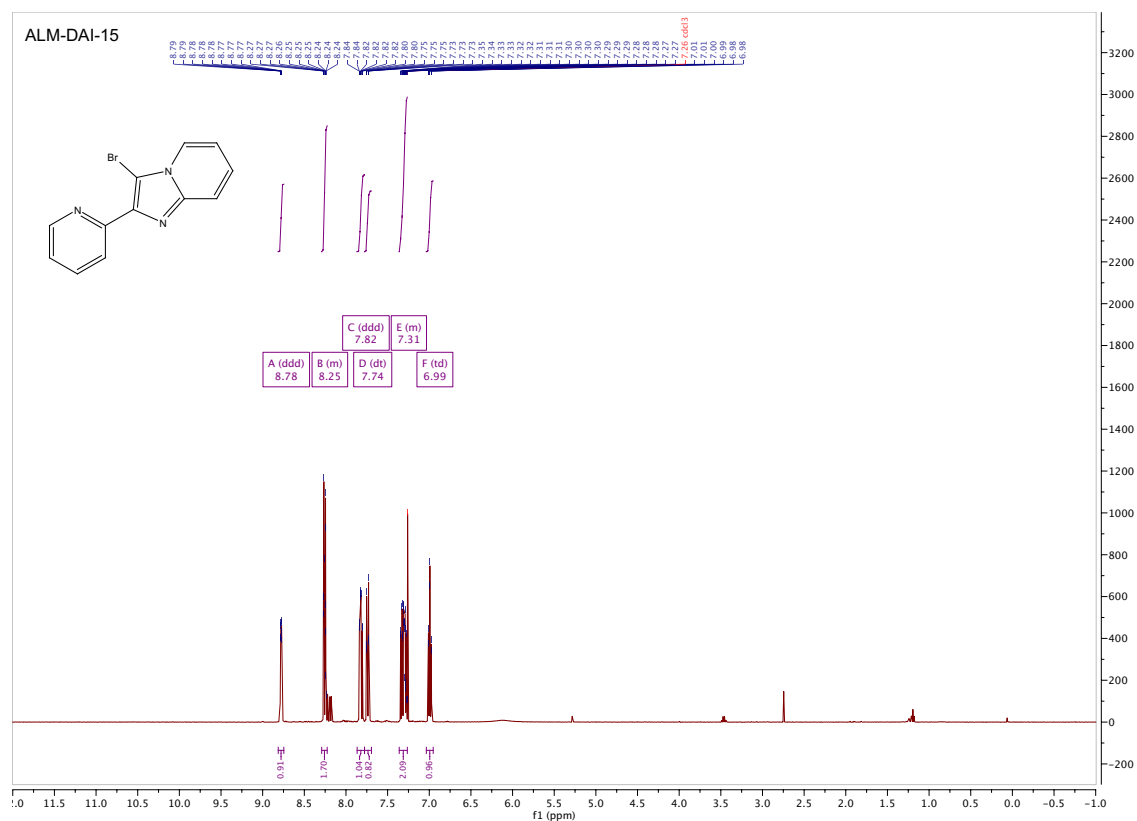
5-bromo-6-(pyridin-2-yl)-2,3-dihydroimidazo[2,1-*b*]thiazole (ALM-DAI-13).



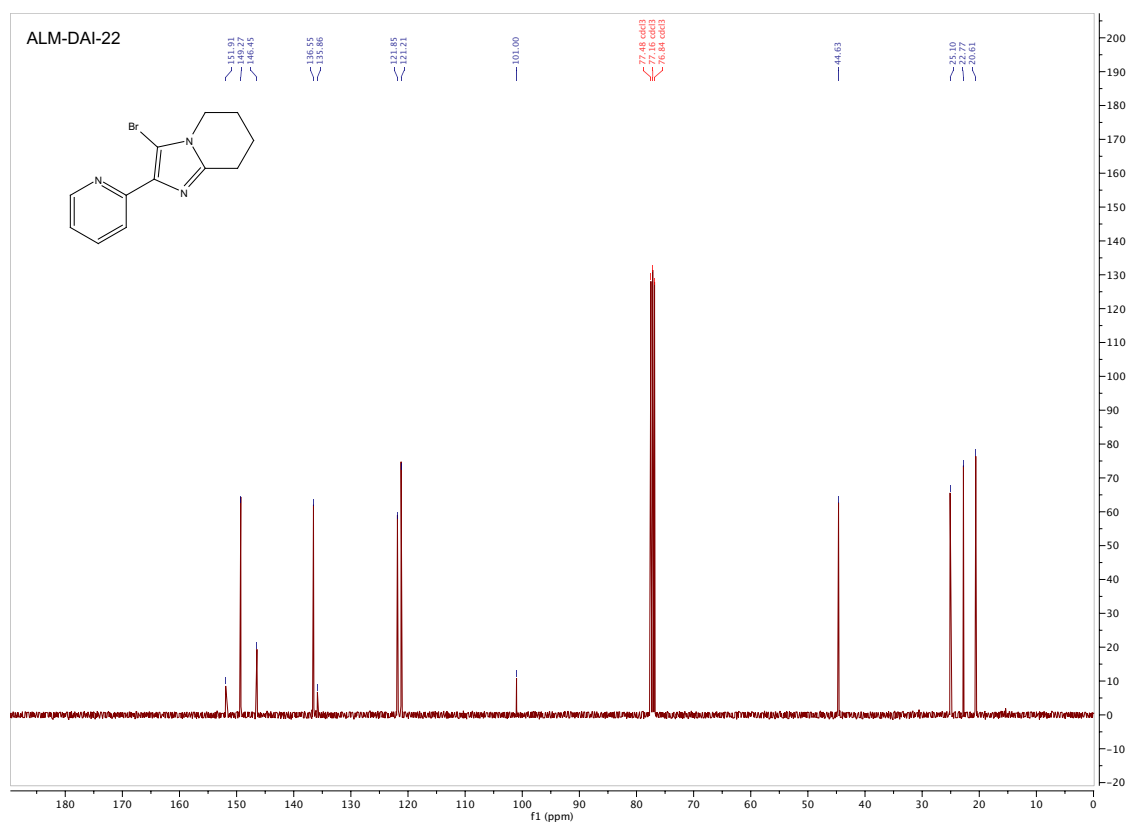
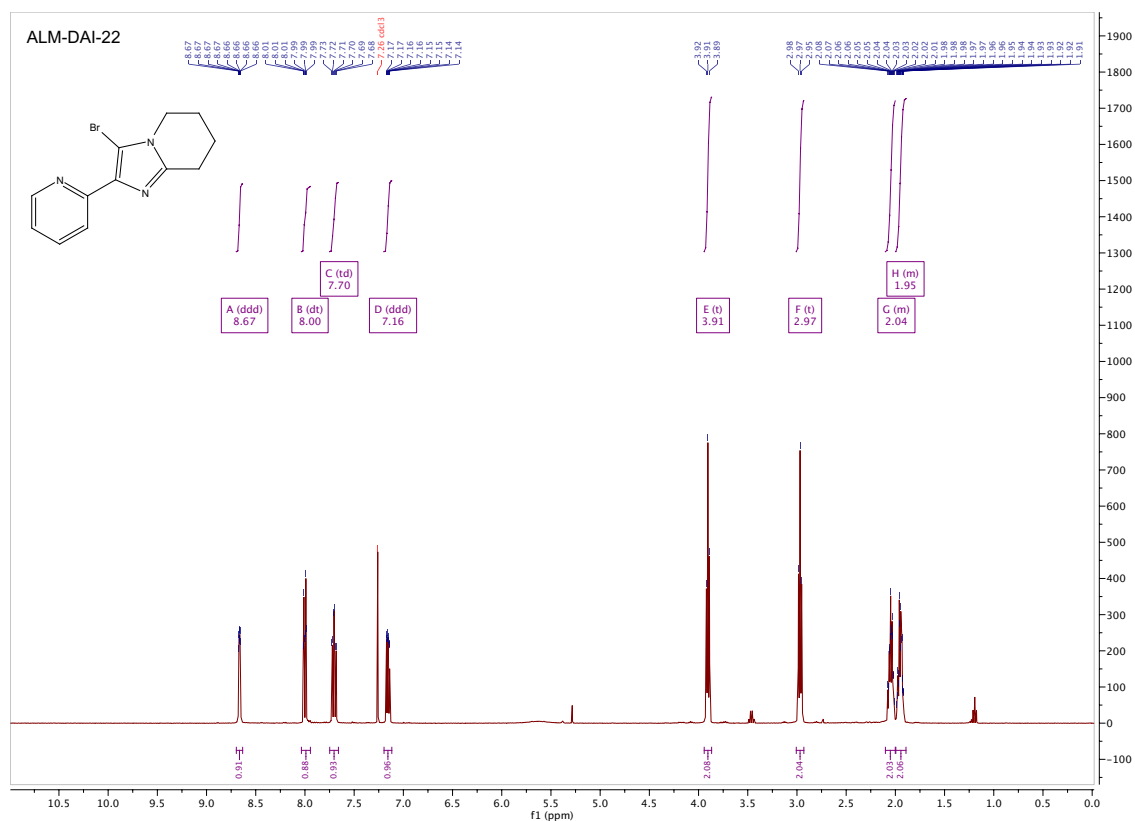
3-bromo-2-(pyridin-2-yl)imidazo[1,2-*a*]pyrimidine (ALM-DAI-14).



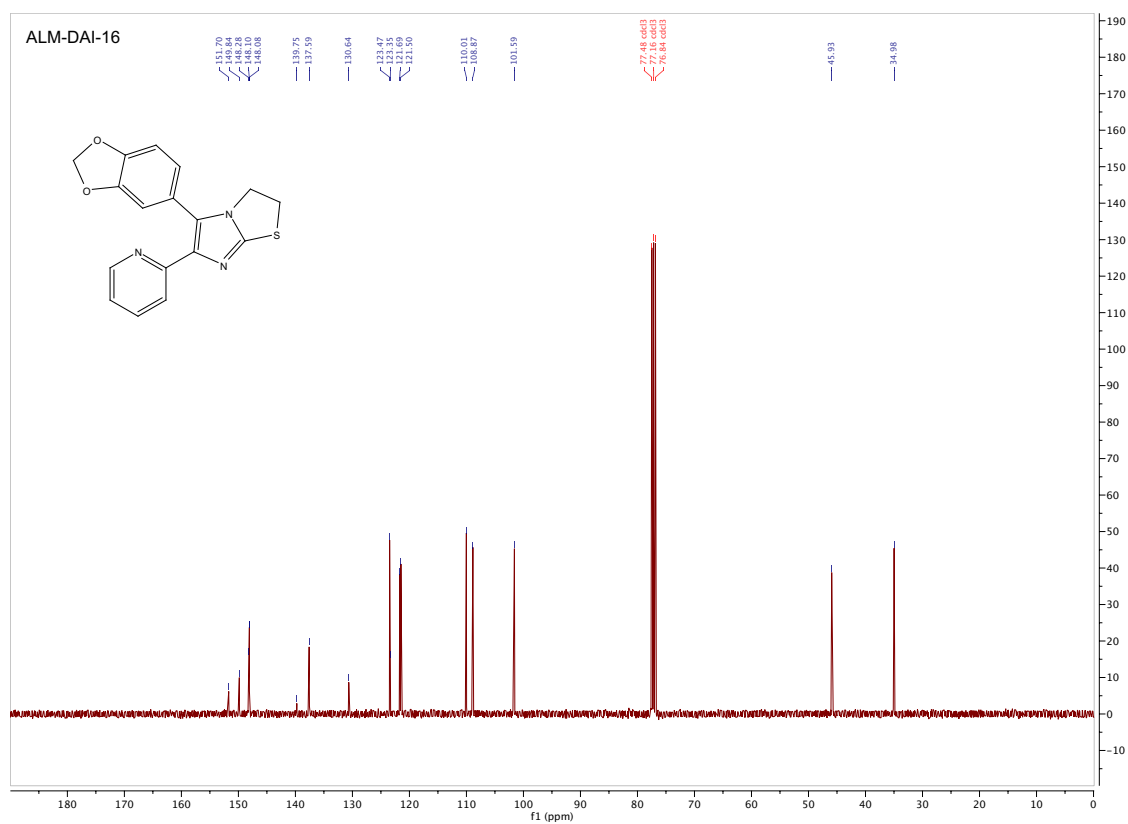
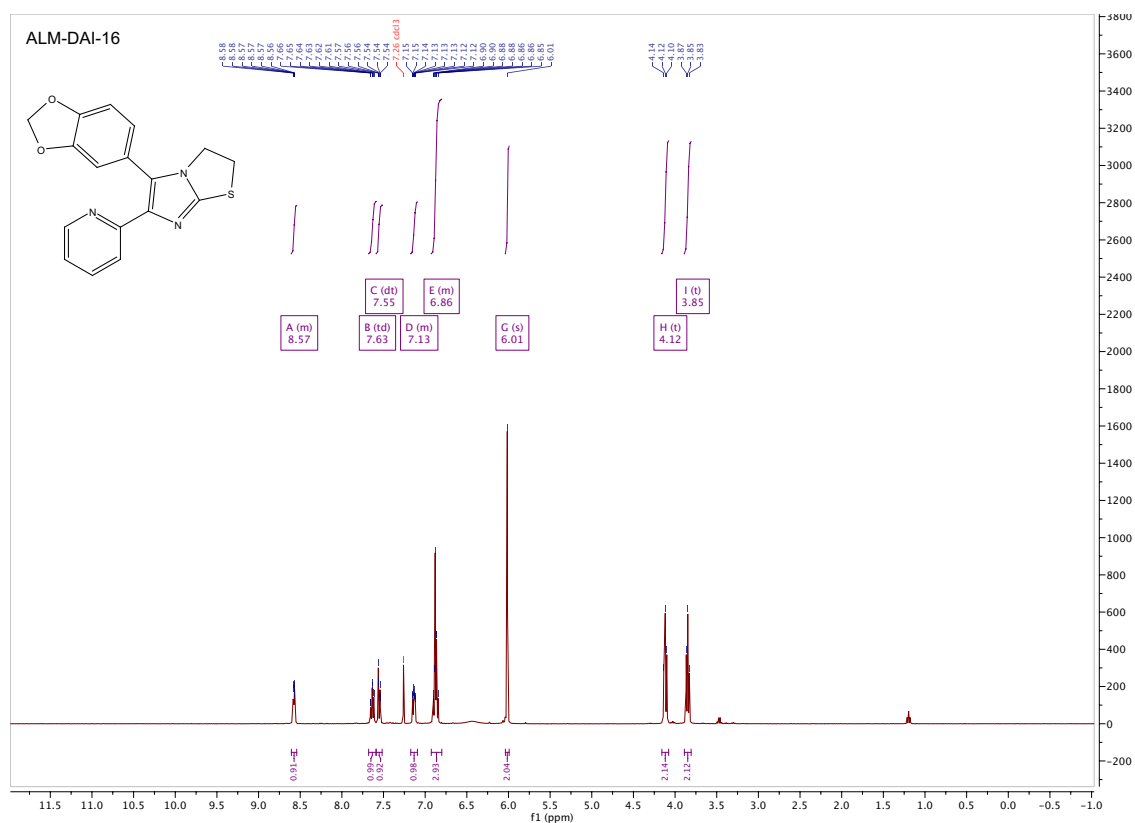
3-bromo-2-(pyridin-2-yl)imidazo[1,2-*a*]pyridine (ALM-DAI-15).



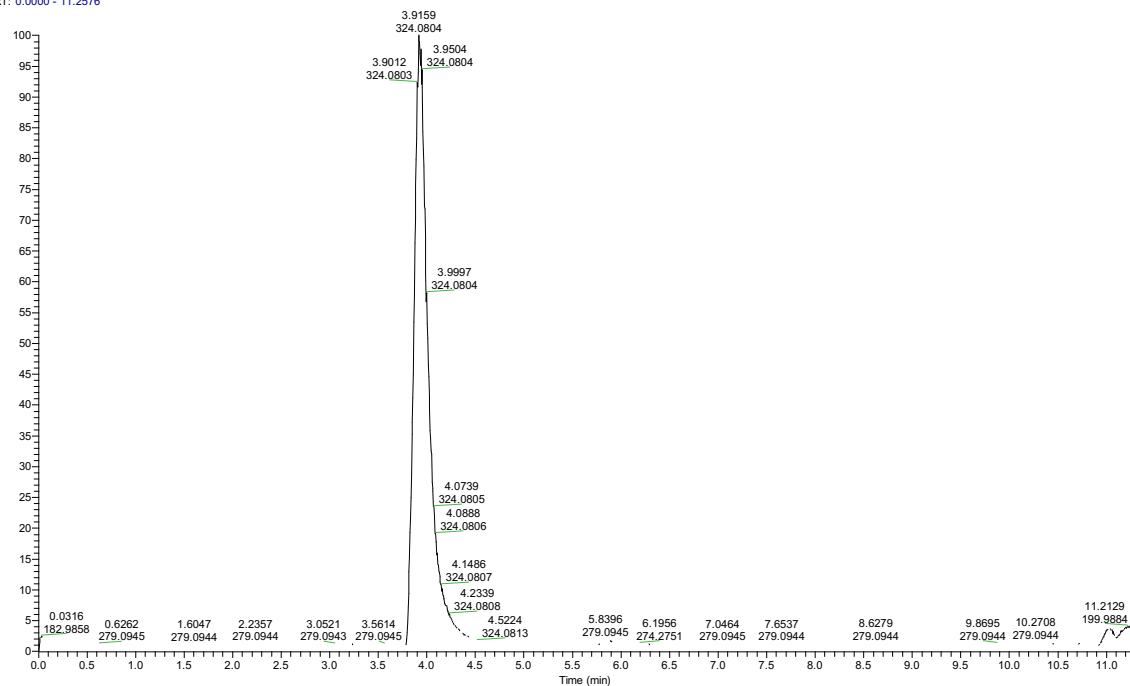
3-bromo-2-(pyridin-2-yl)-5,6,7,8-tetrahydroimidazo[1,2-a]pyridine (ALM-DAI-22).



5-(benzo[d][1,3]dioxol-5-yl)-6-(pyridin-2-yl)-2,3-dihydroimidazo[2,1-b]thiazole (ALM-DAI-16).

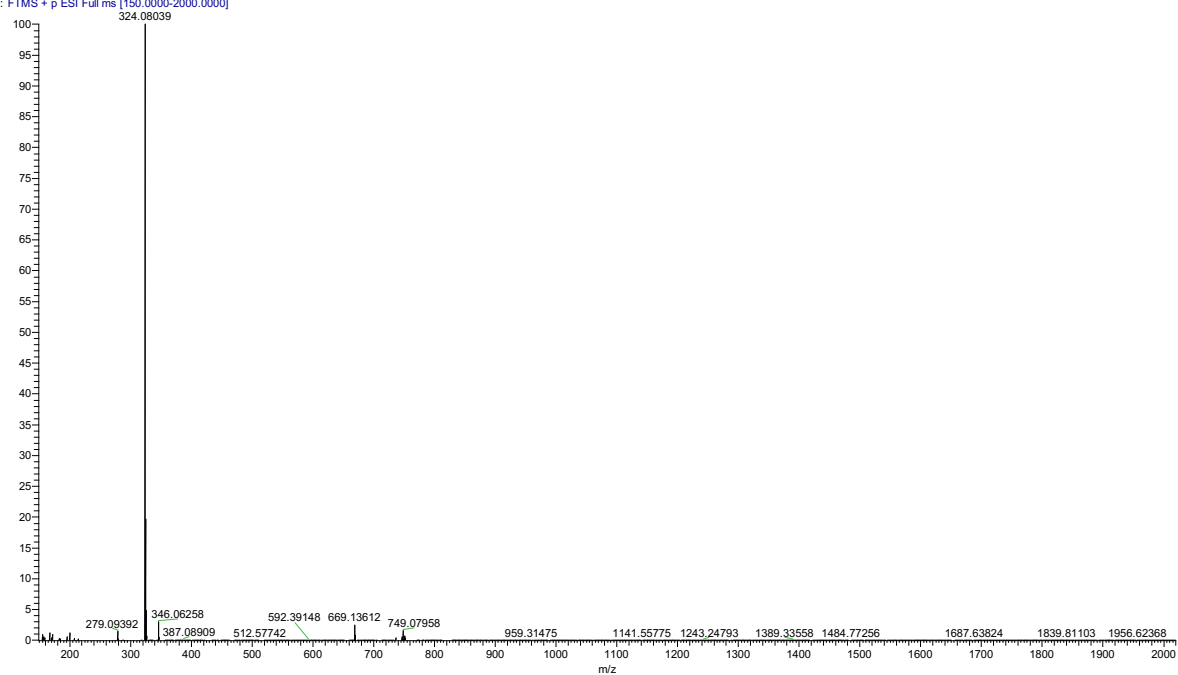


RT: 0.0000 - 11.2576

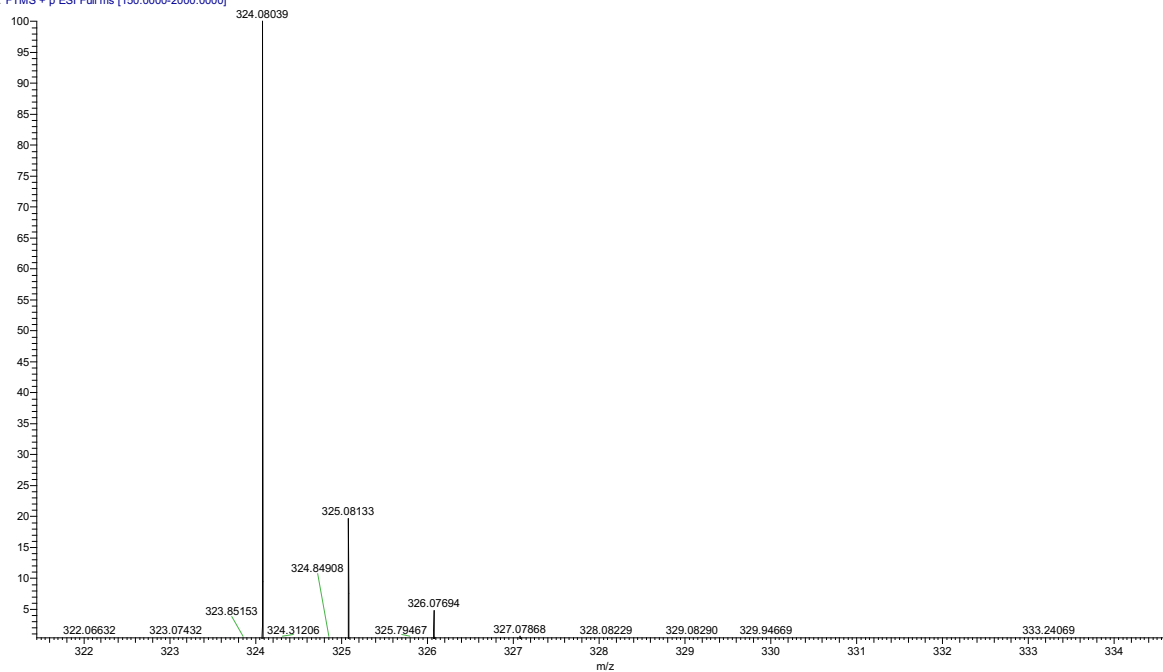


NL:
4.30E9
Base Peak
MS
ALM_DAI_1
6

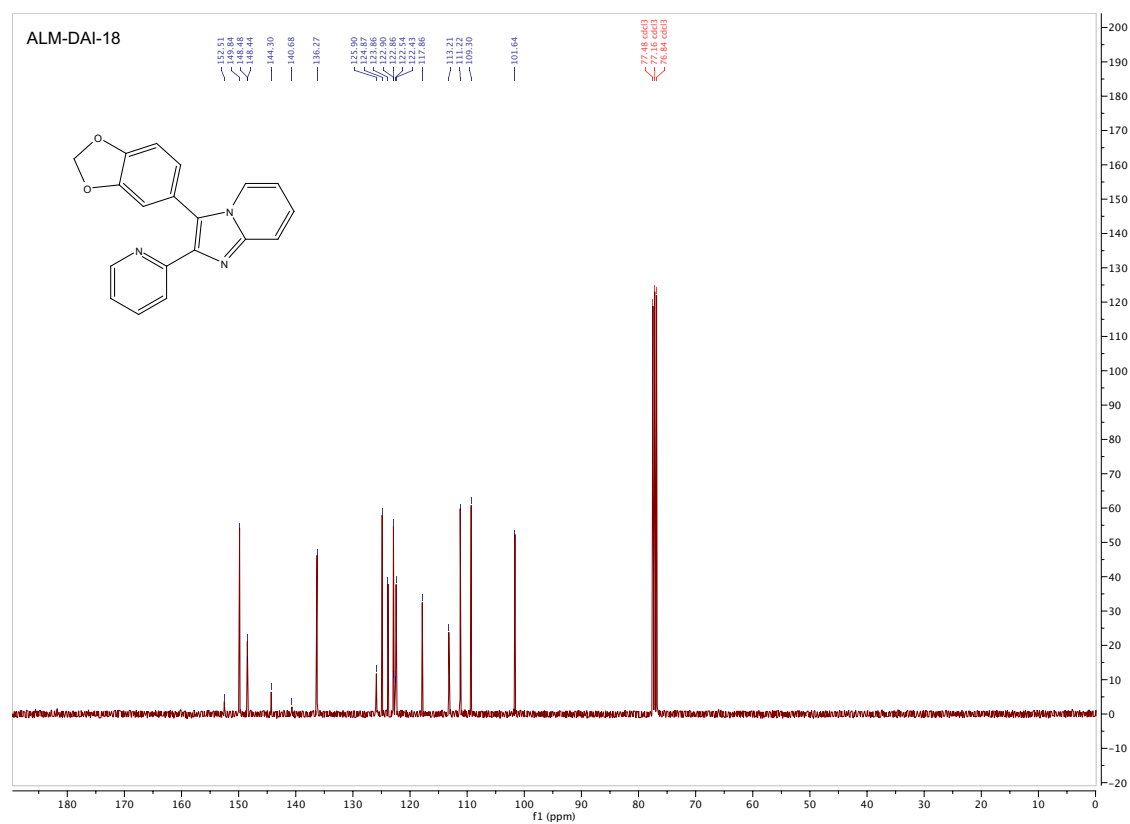
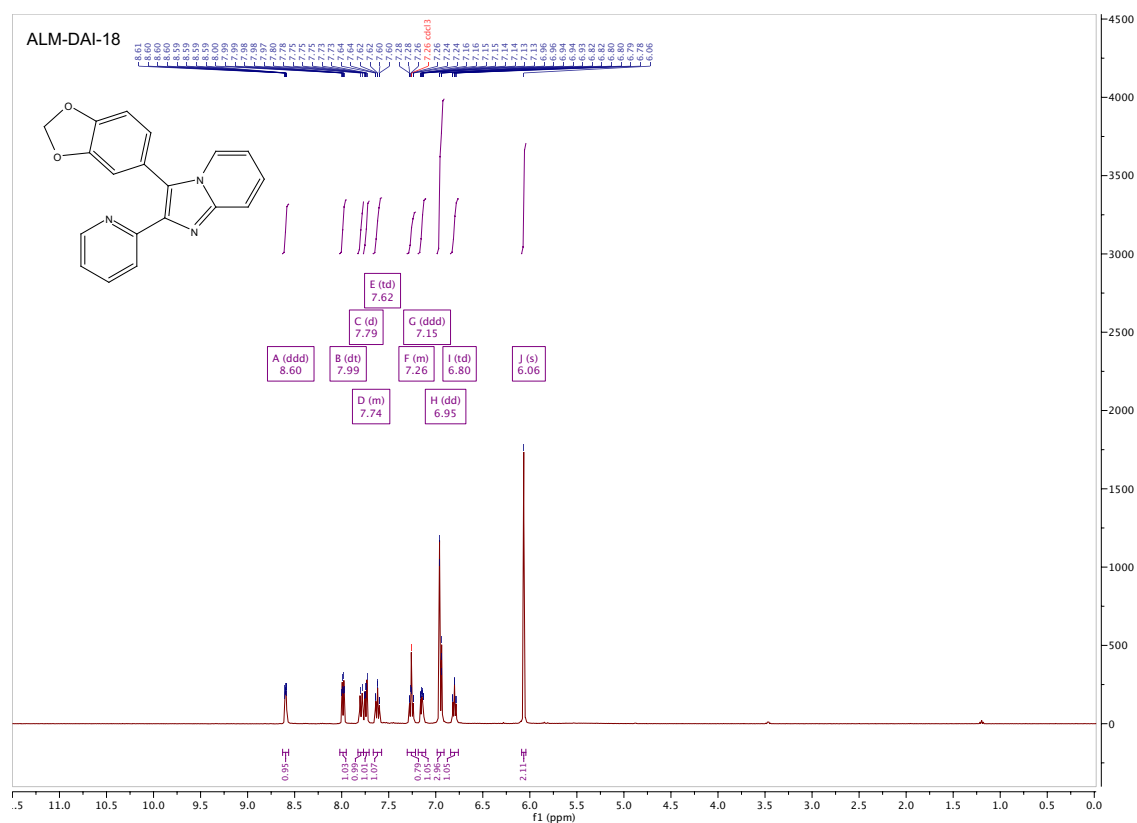
ALM_DAI_16 #749-789 RT: 3.85-4.05 AV: 41 NL: 2.73E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



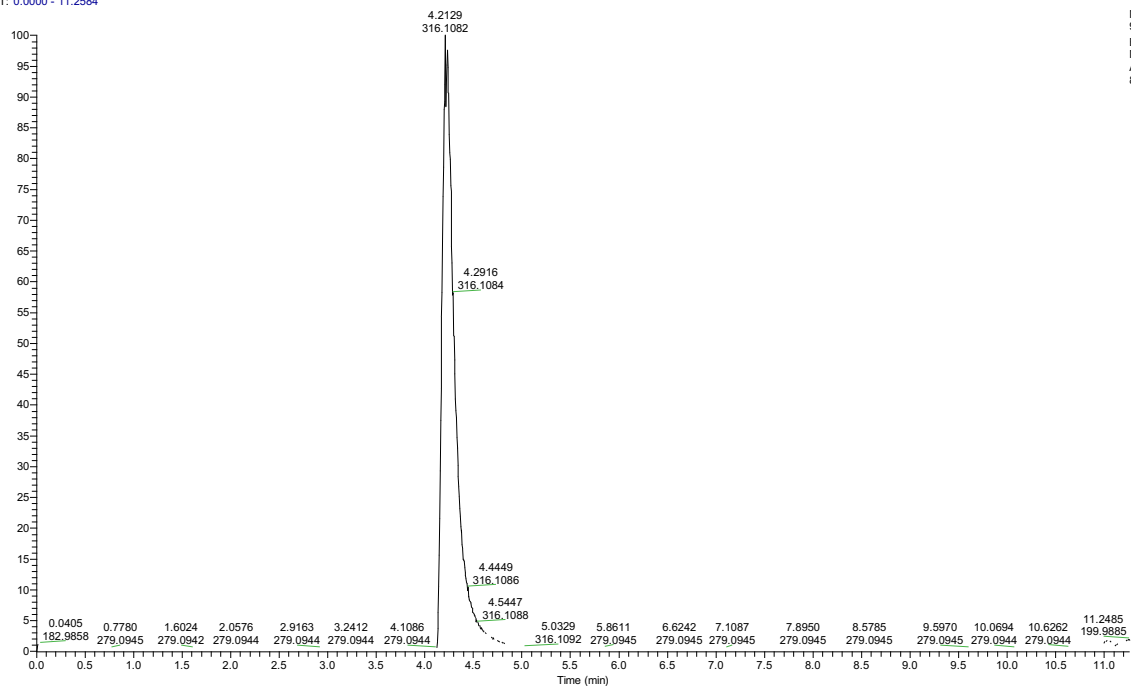
ALM_DAI_16 #749-789 RT: 3.85-4.05 AV: 41 NL: 2.73E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



3-(benzo[d][1,3]dioxol-5-yl)-2-(pyridin-2-yl)imidazo[1,2-*a*]pyridine (ALM-DAI-18).

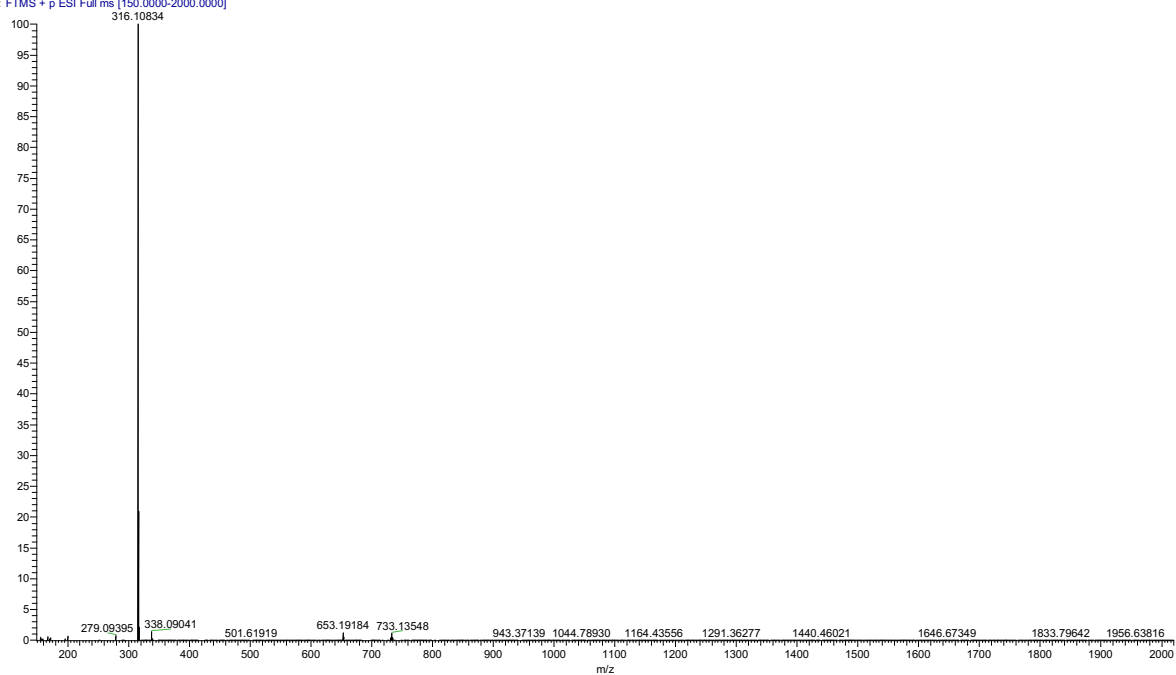


RT: 0.0000 - 11.2584

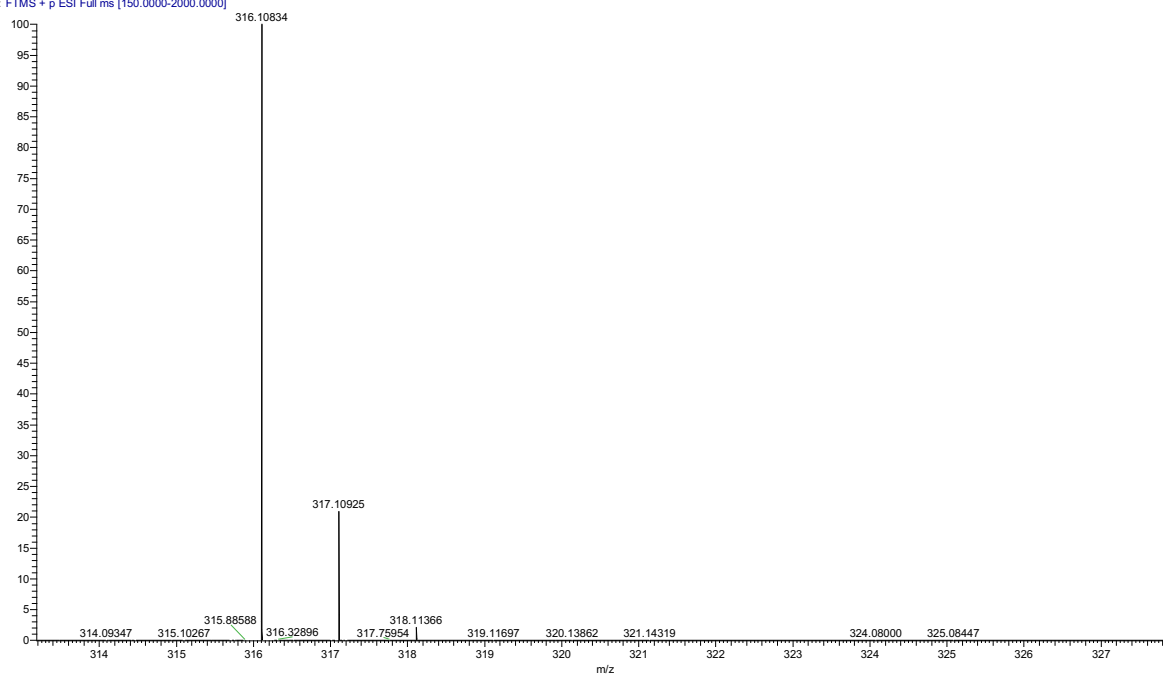


NL:
9.14E9
Base Peak
MS
ALM_DAI_1
8

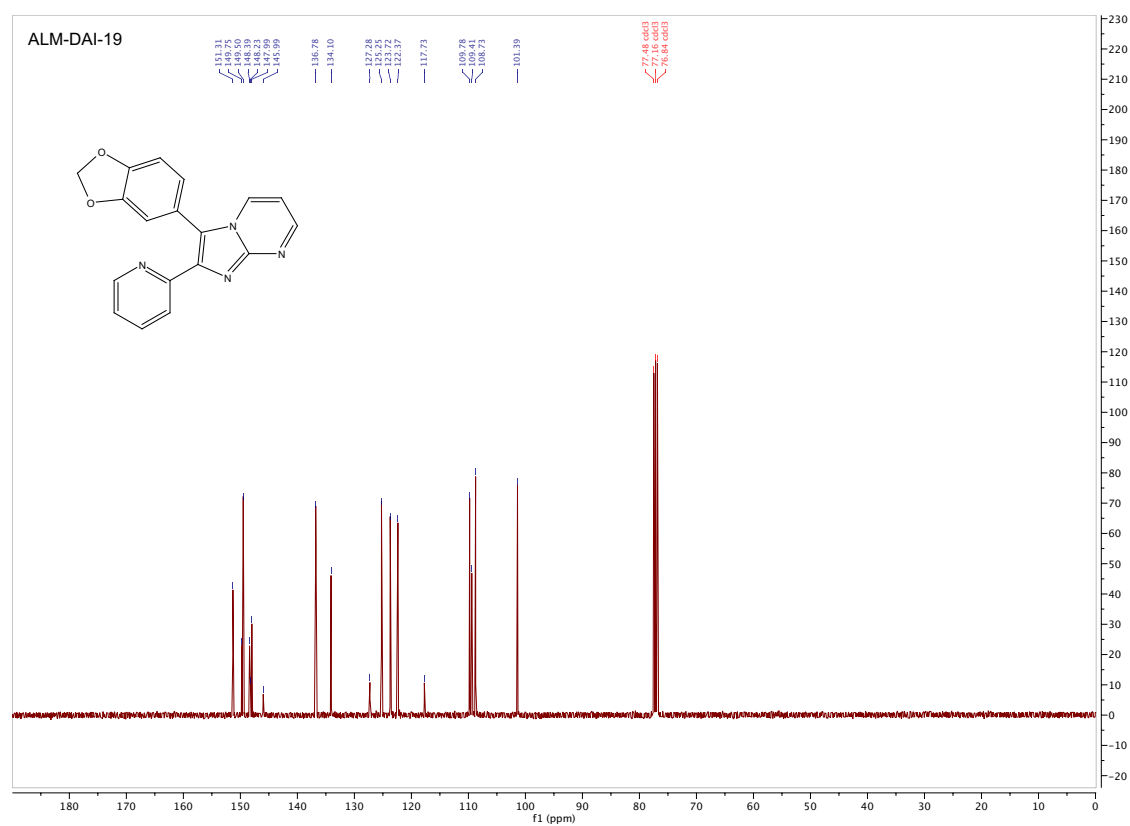
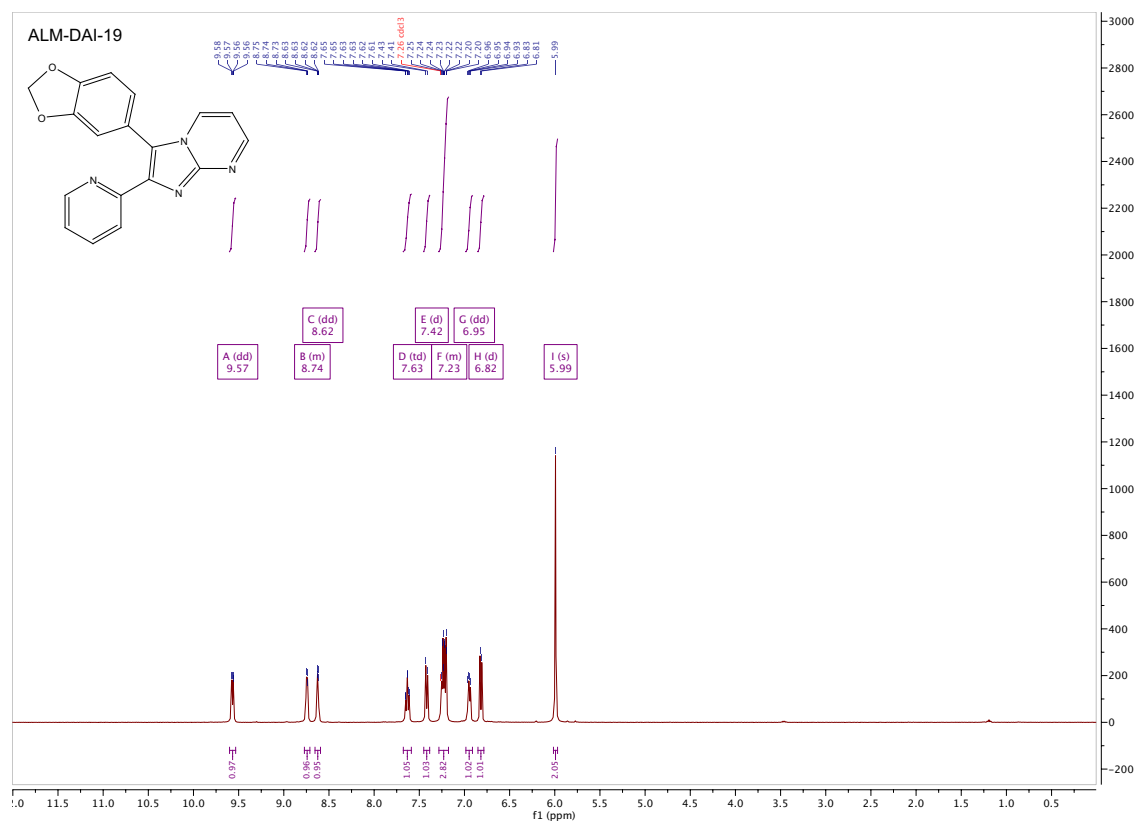
ALM_DAI_18 #810-852 RT: 4.16-4.37 AV: 43 NL: 5.24E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



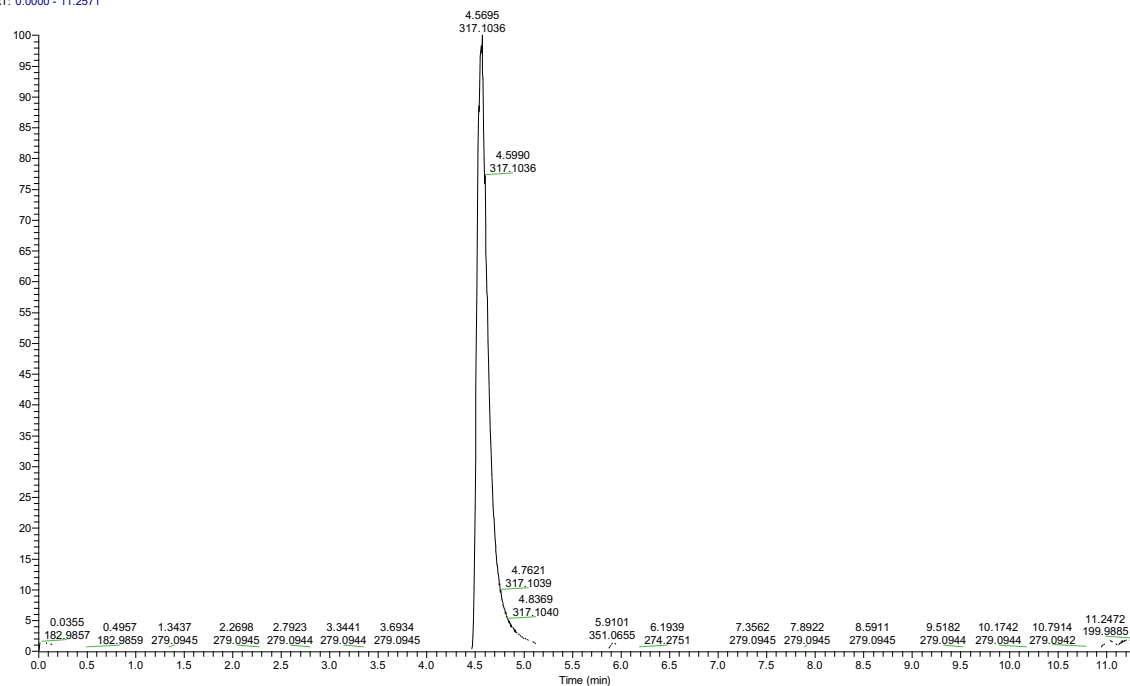
ALM_DAI_18 #810-852 RT: 4.16-4.37 AV: 43 NL: 5.24E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



3-(benzo[d][1,3]dioxol-5-yl)-2-(pyridin-2-yl)imidazo[1,2-*a*]pyrimidine (ALM-DAI-19).

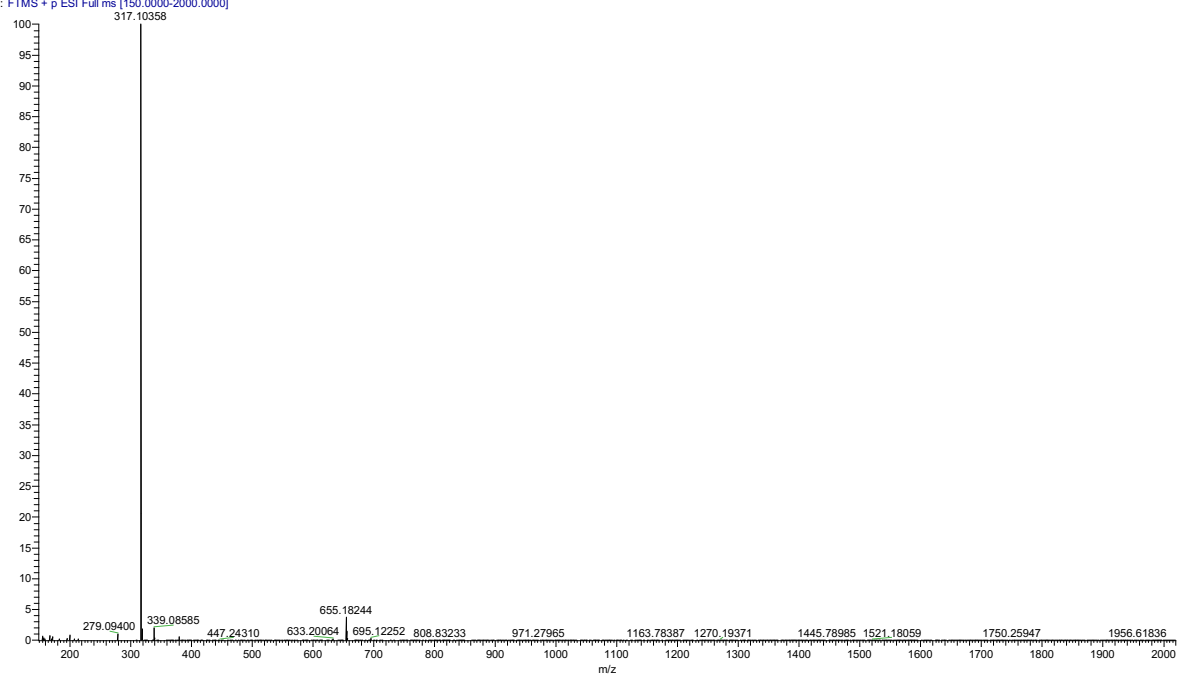


RT: 0.0000 - 11.2571

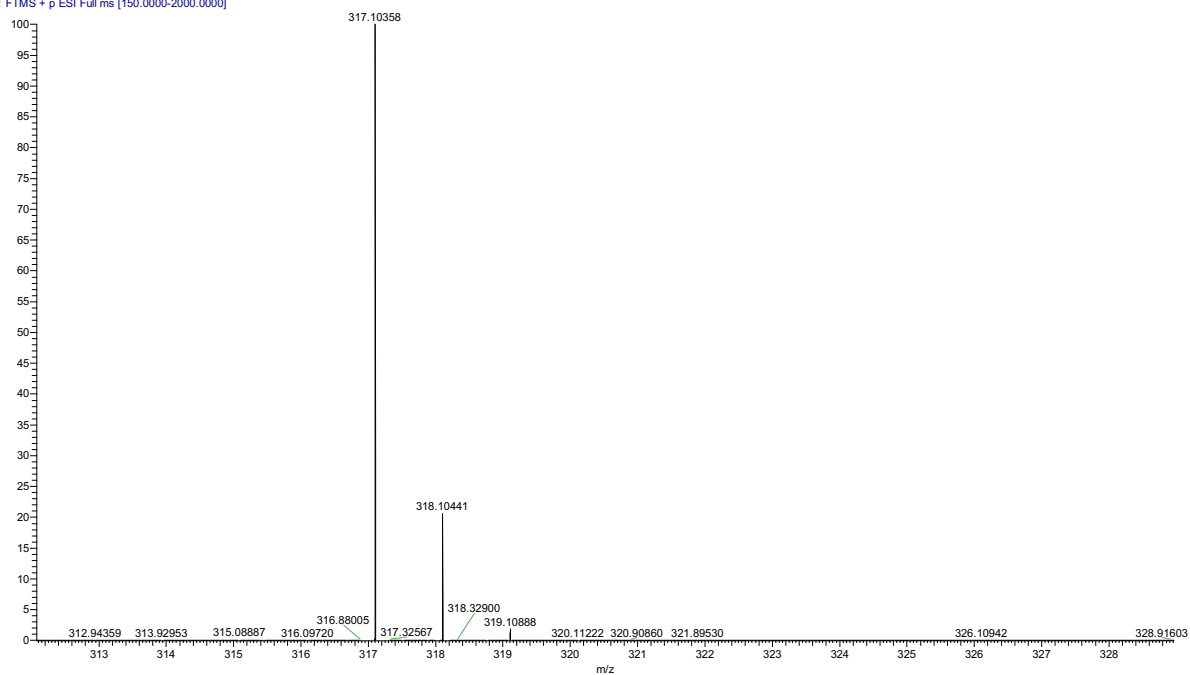


NL:
8.93E9
Base Peak
MS
ALM_DAI_1
9

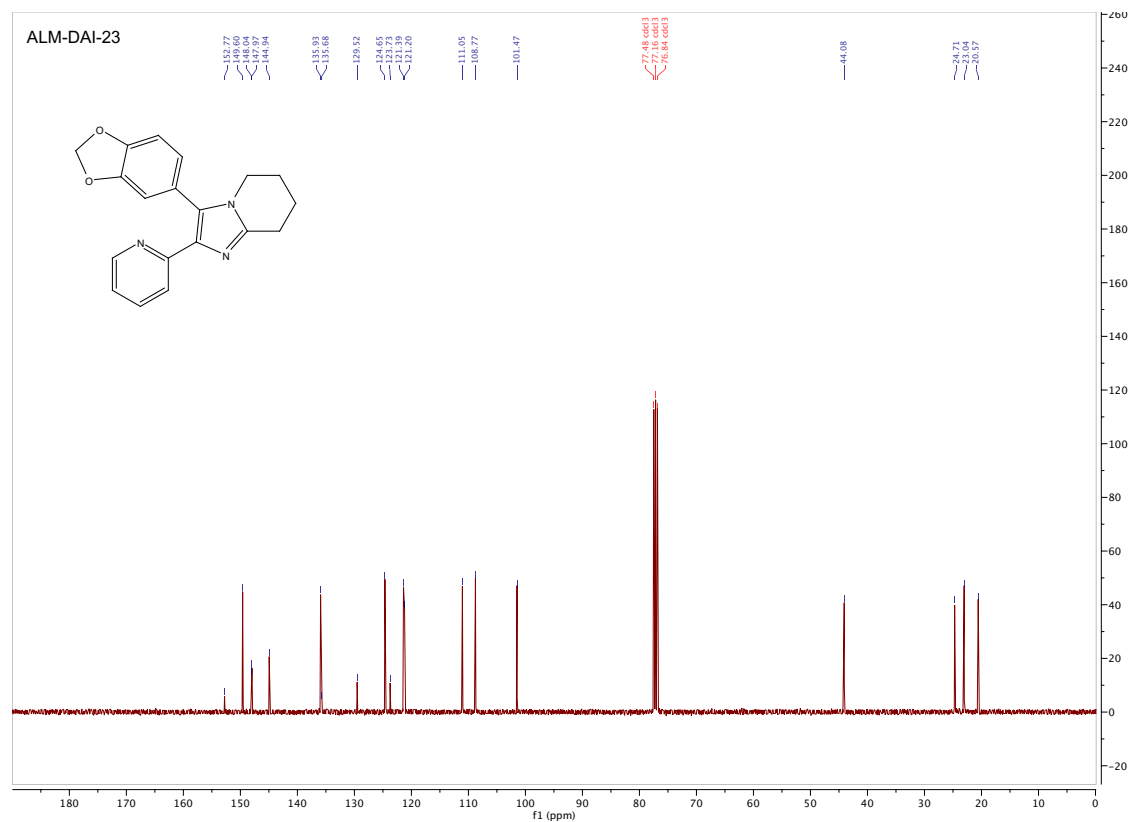
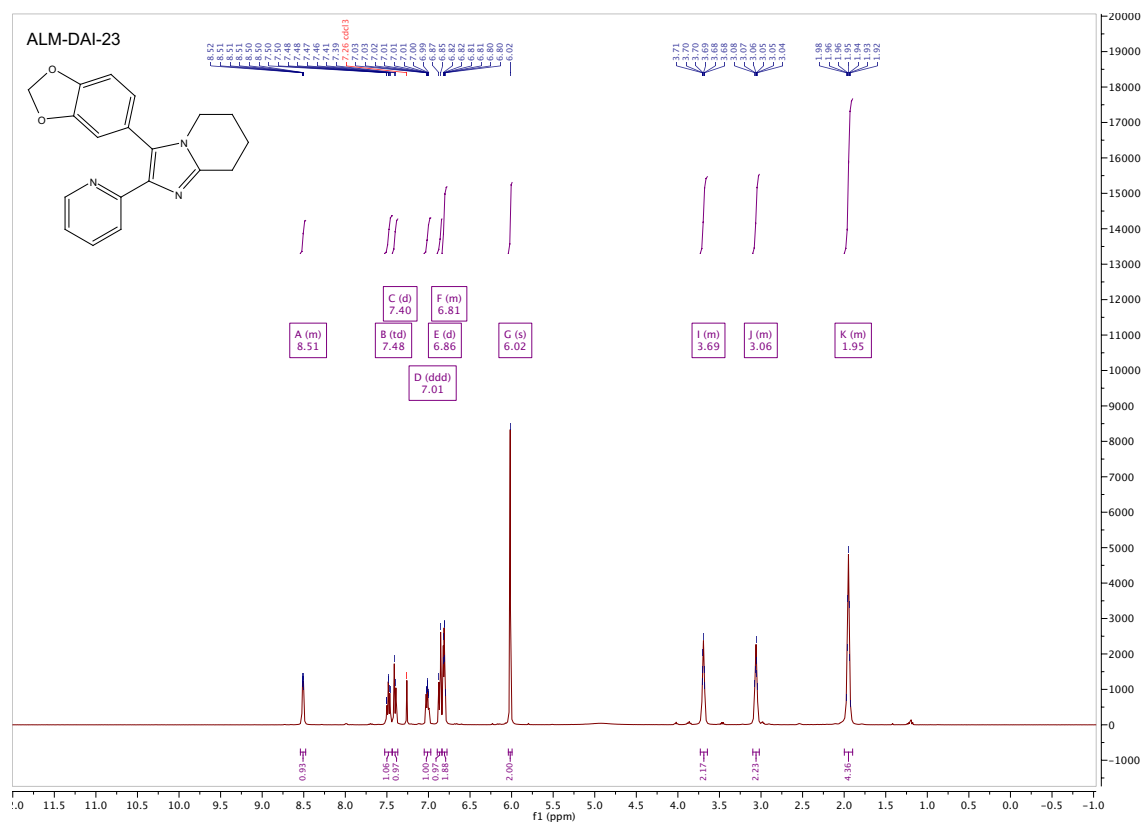
ALM_DAI_19 #874-925 RT: 4.49-4.74 AV: 52 NL: 4.19E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



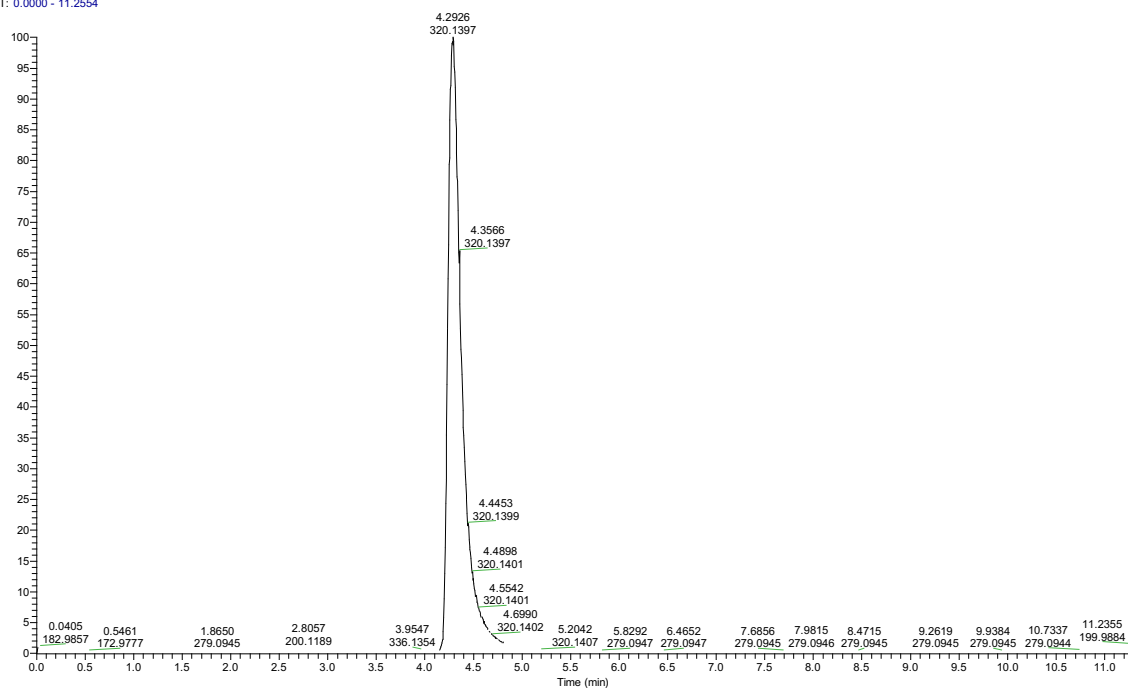
ALM_DAI_19 #874-925 RT: 4.49-4.74 AV: 52 NL: 4.19E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



3-(benzo[d][1,3]dioxol-5-yl)-2-(pyridin-2-yl)-5,6,7,8-tetrahydroimidazo[1,2-a]pyridine (ALM-DAI-23).

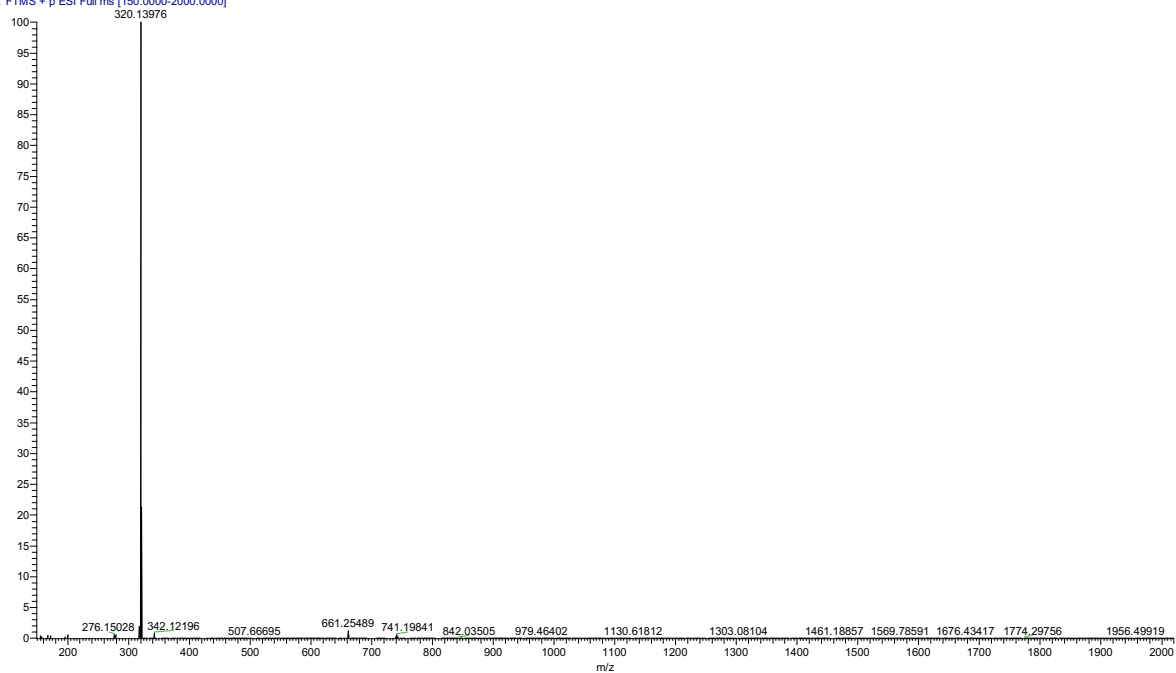


RT: 0.0000 - 11.2554

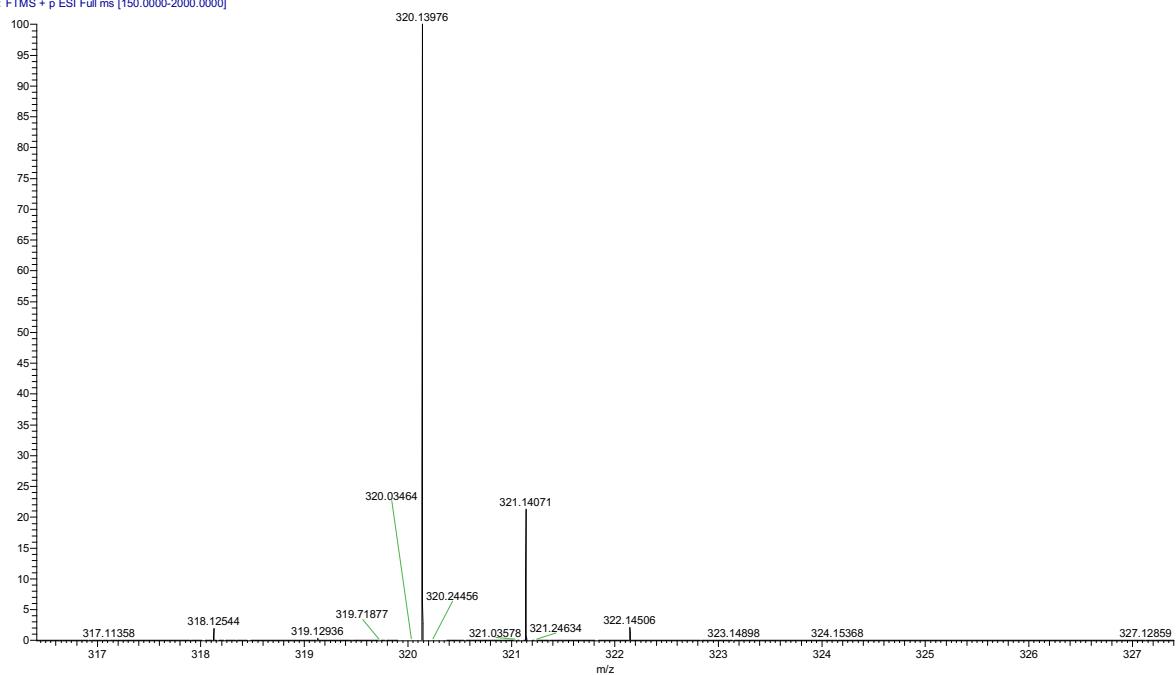


NL:
1.21E10
Base Peak
MS
ALM_DAI_2
3

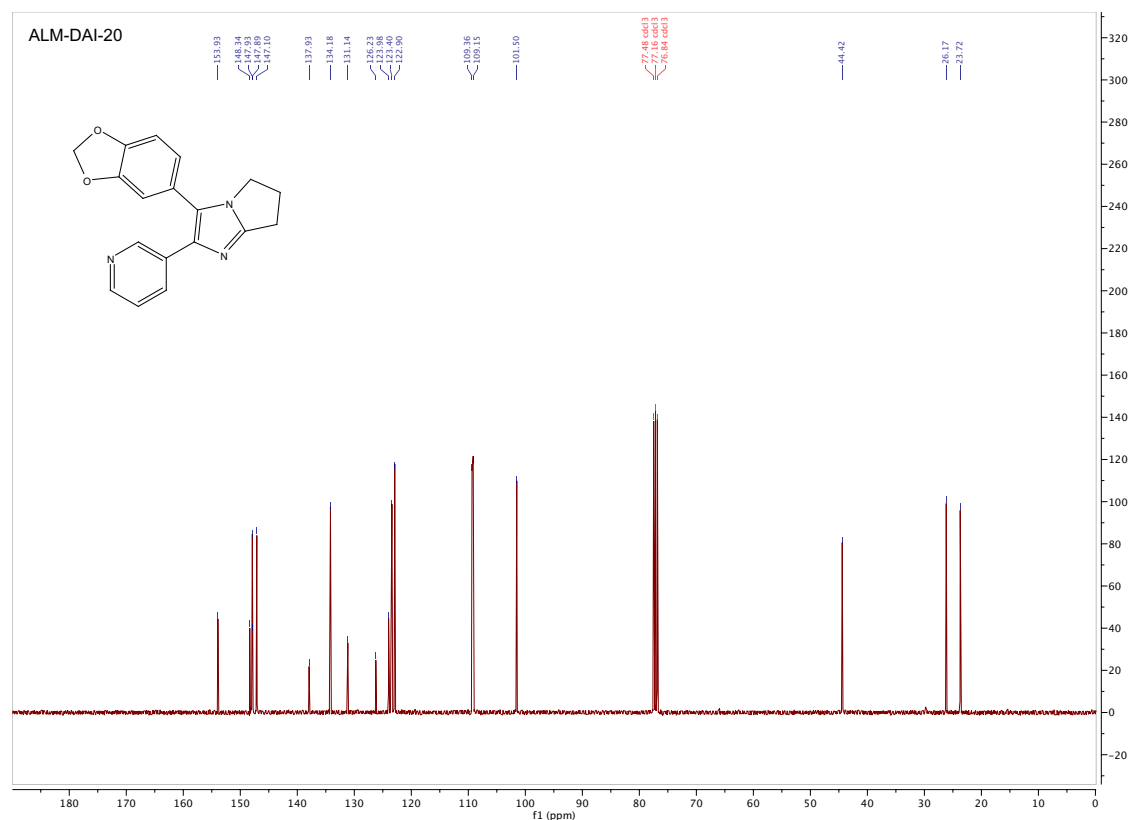
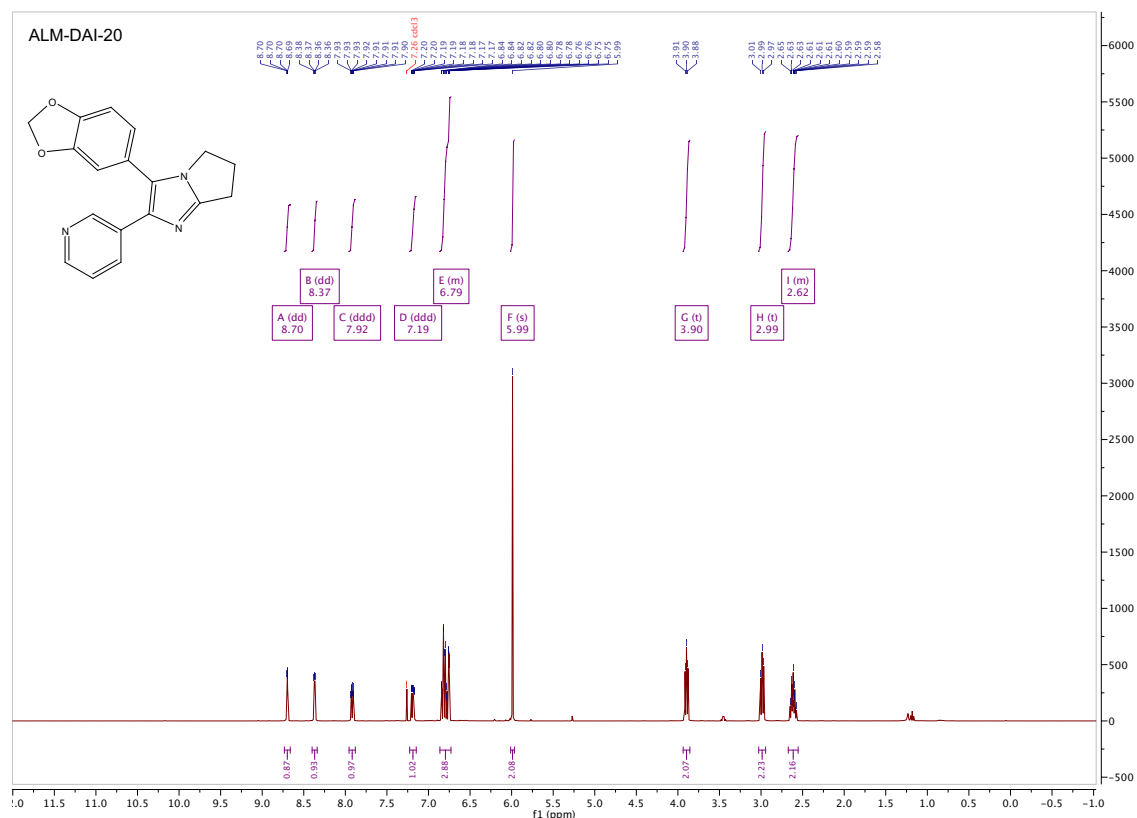
ALM_DAI_23 #520-879 RT: 4.21-4.50 AV: 60 NL: 5.92E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



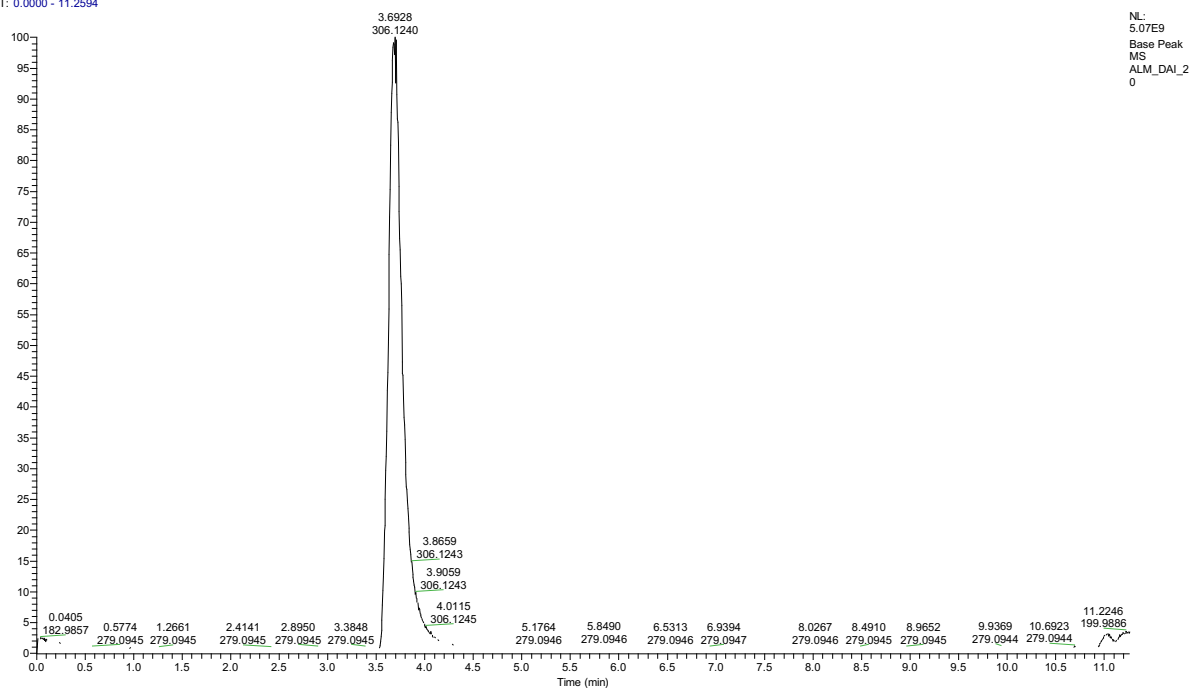
ALM_DAI_23 #820-879 RT: 4.21-4.50 AV: 60 NL: 5.92E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



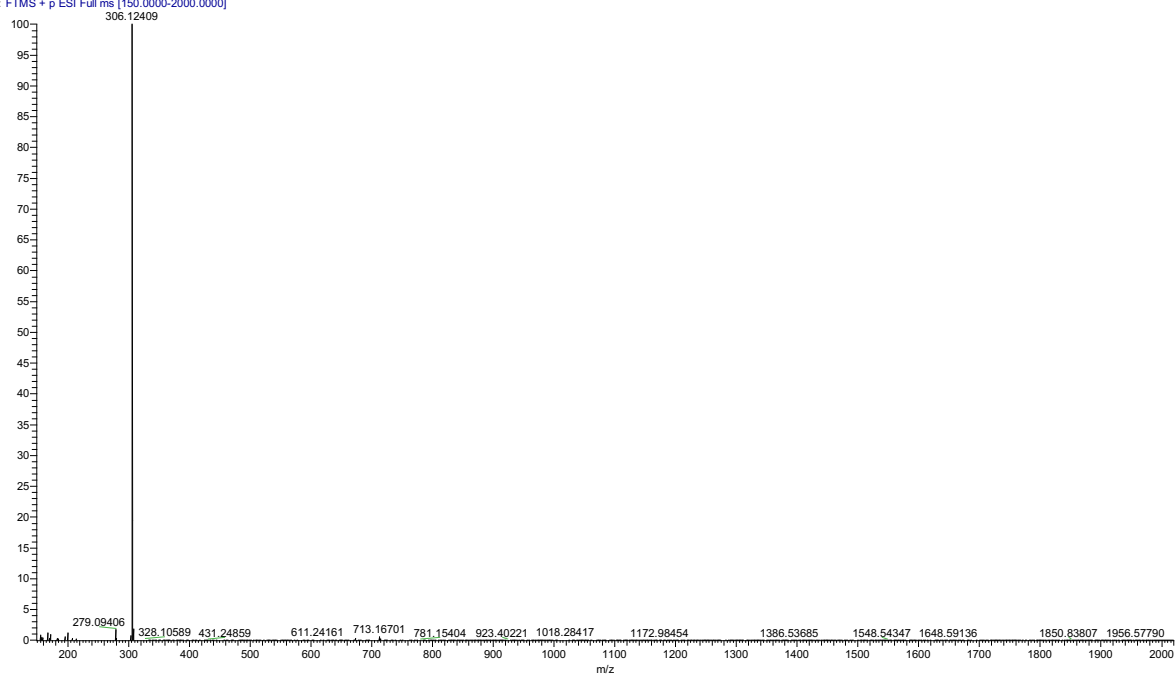
3-(benzo[d][1,3]dioxol-5-yl)-2-(pyridin-3-yl)-6,7-dihydro-5H-pyrrolo[1,2-a]imidazole (ALM-DAI-20).



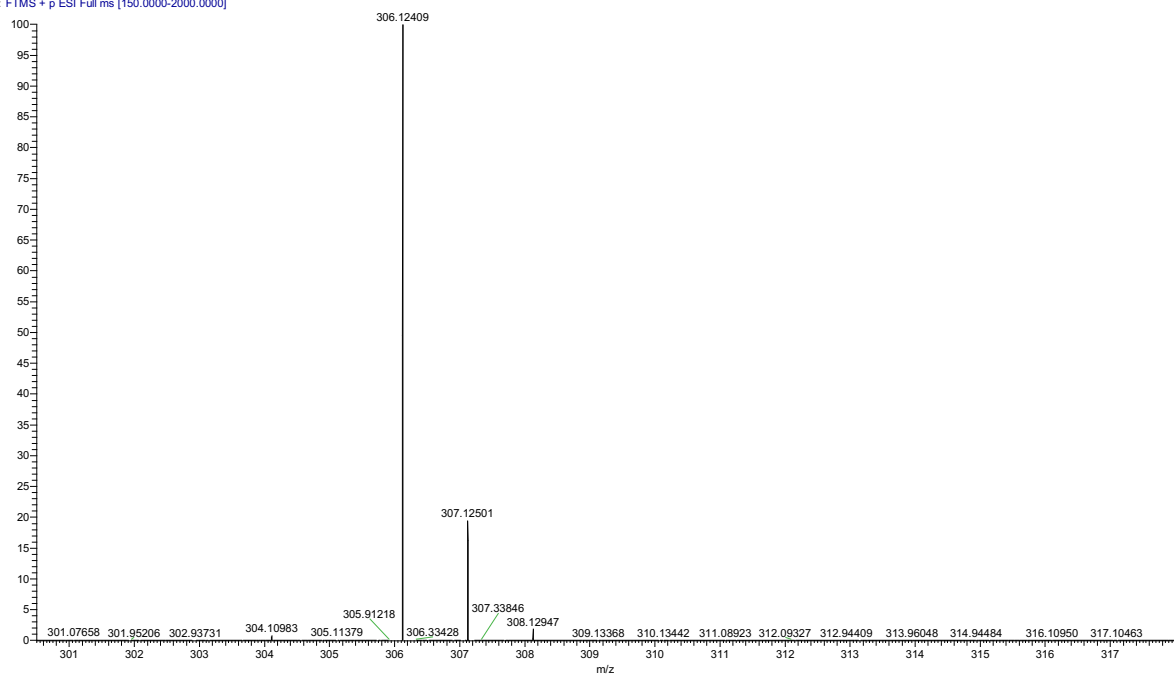
RT: 0.0000 - 11.2594



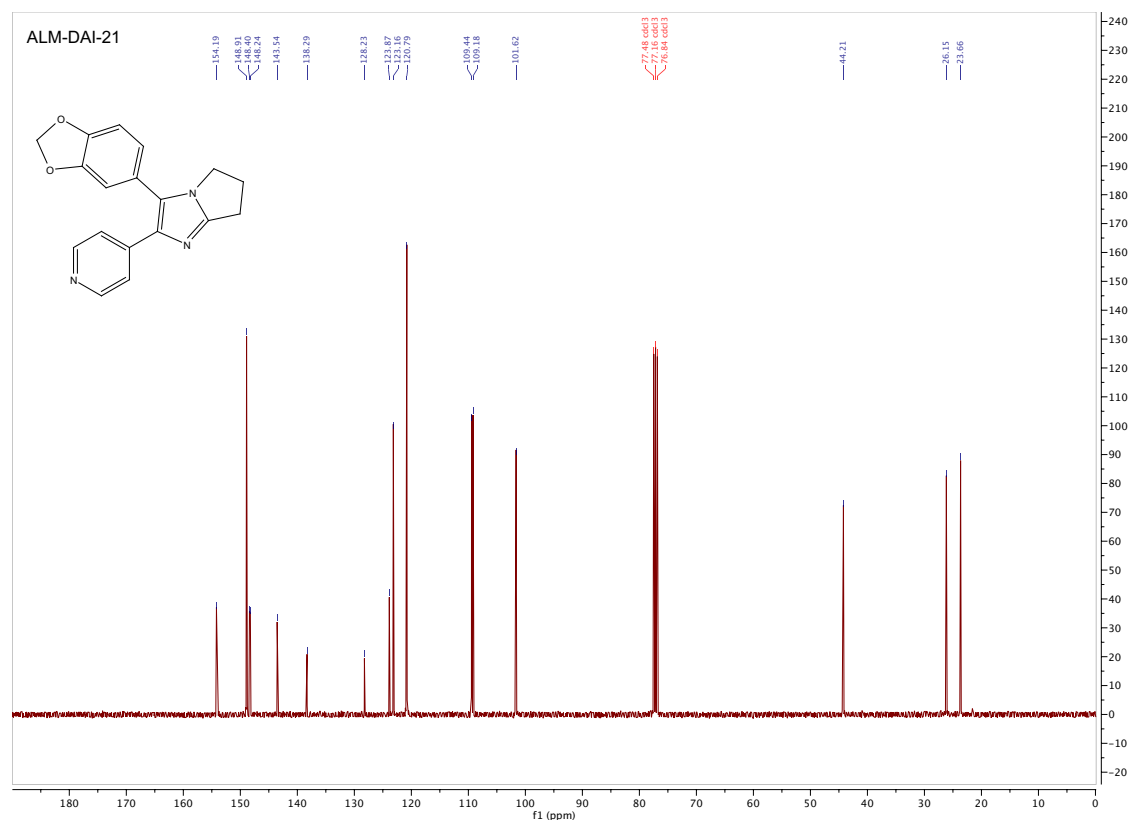
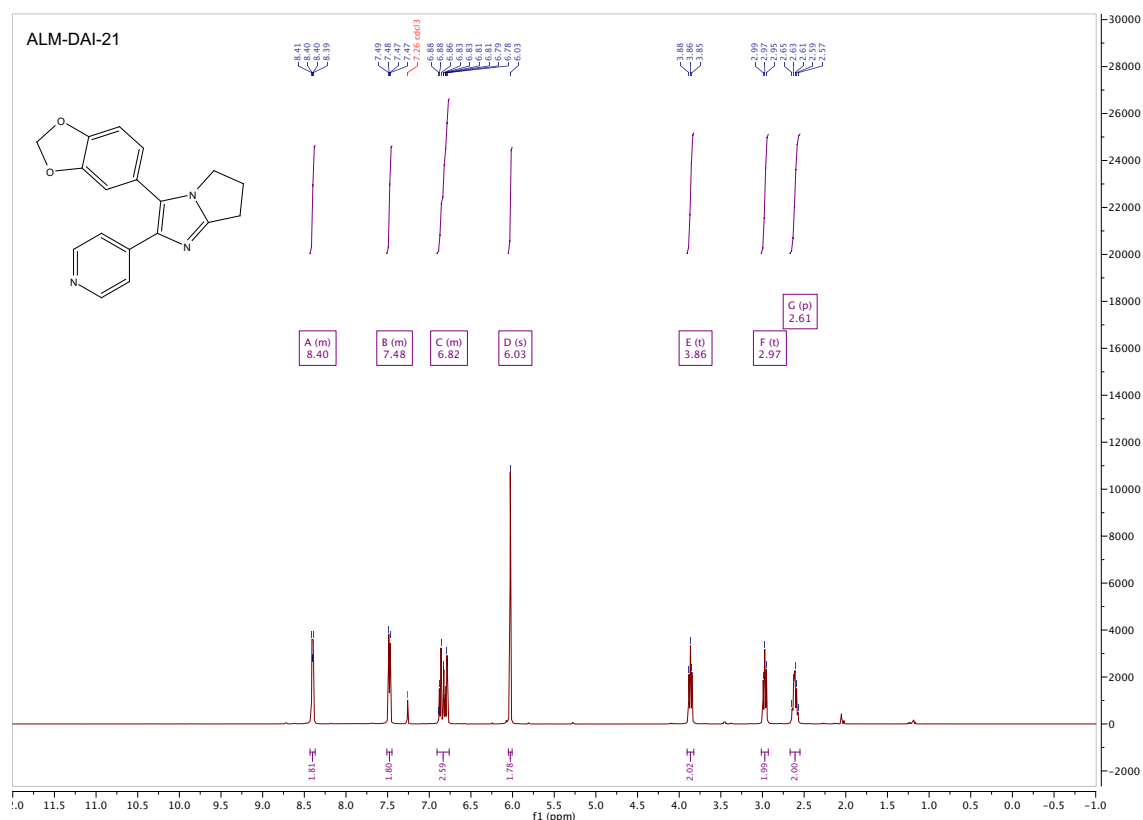
ALM_DAI_20 #596-753 RT: 3.58-3.86 AV: 58 NL: 2.63E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



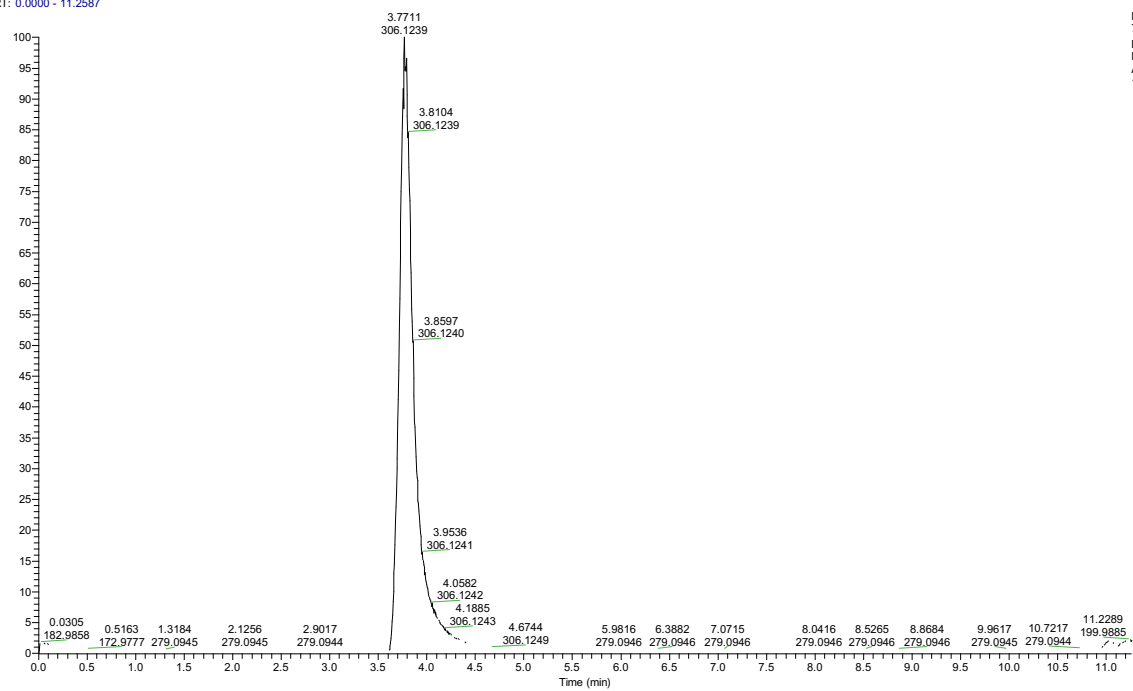
ALM_DAI_20 #696-753 RT: 3.58-3.86 AV: 58 NL: 2.63E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



3-(benzo[d][1,3]dioxol-5-yl)-2-(pyridin-4-yl)-6,7-dihydro-5H-pyrrolo[1,2-a]imidazole (ALM-DAI-21).

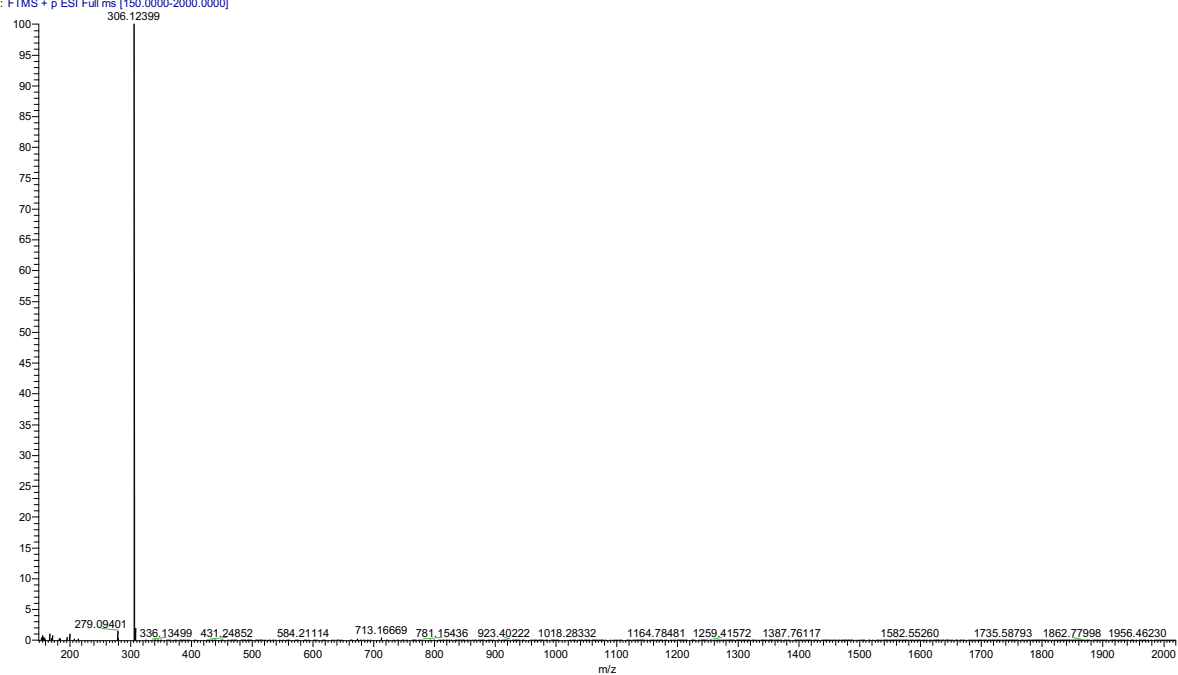


RT: 0.0000 - 11.2587

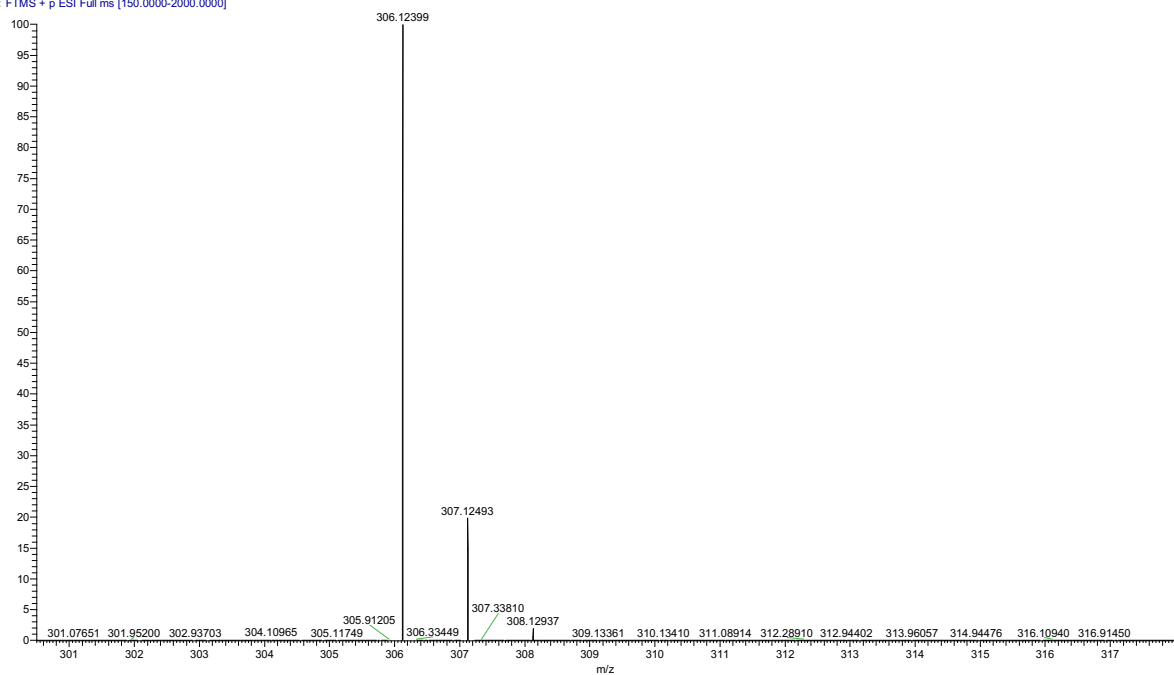


NL:
7.80E9
Base Peak
MS
ALM_DAI_2
1

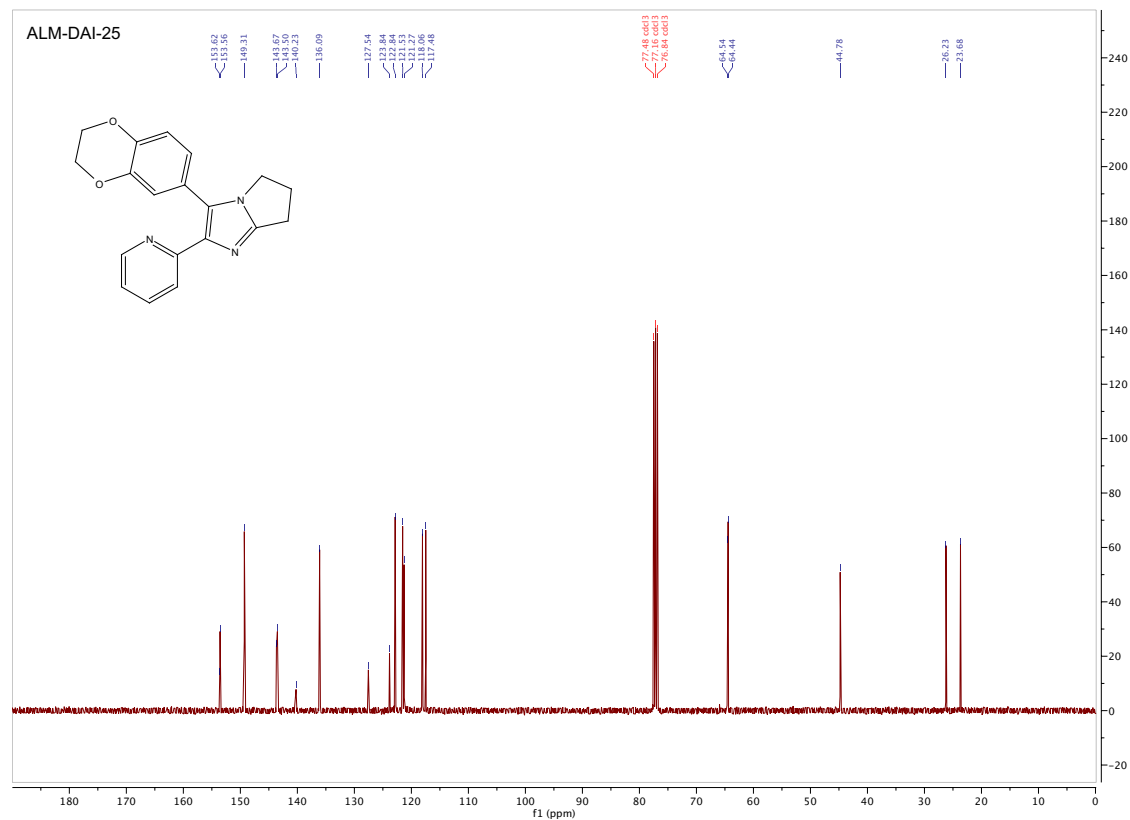
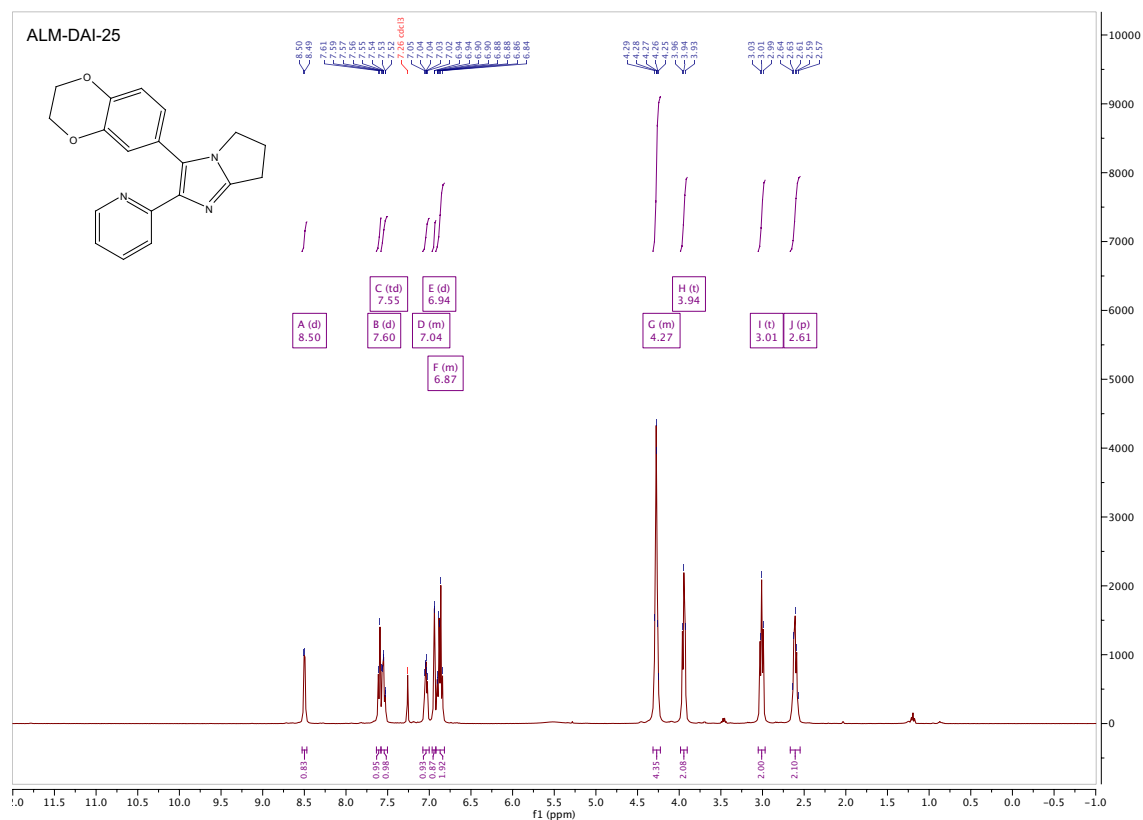
ALM_DAI_21 #709-788 RT: 3.65-4.04 AV: 80 NL: 2.97E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



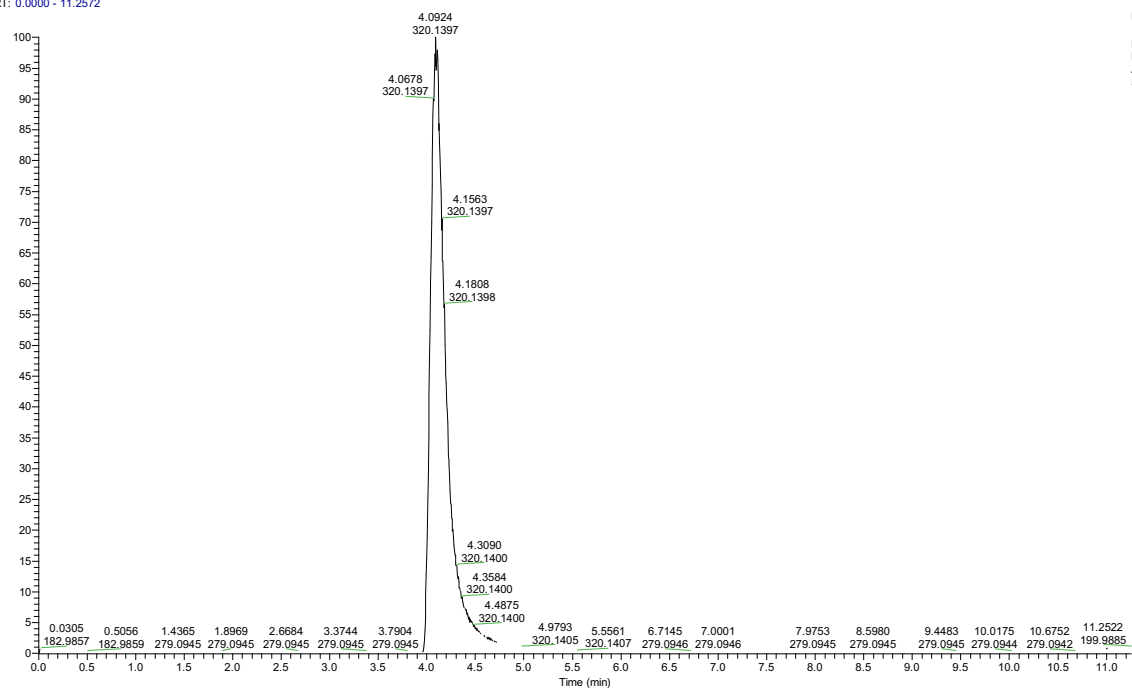
ALM_DAI_21 #709-788 RT: 3.65-4.04 AV: 80 NL: 2.97E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



3-(2,3-dihydrobenzo[*b*][1,4]dioxin-6-yl)-2-(pyridin-2-yl)-6,7-dihydro-5*H*-pyrrolo[1,2-*a*]imidazole (ALM-DAI-25).

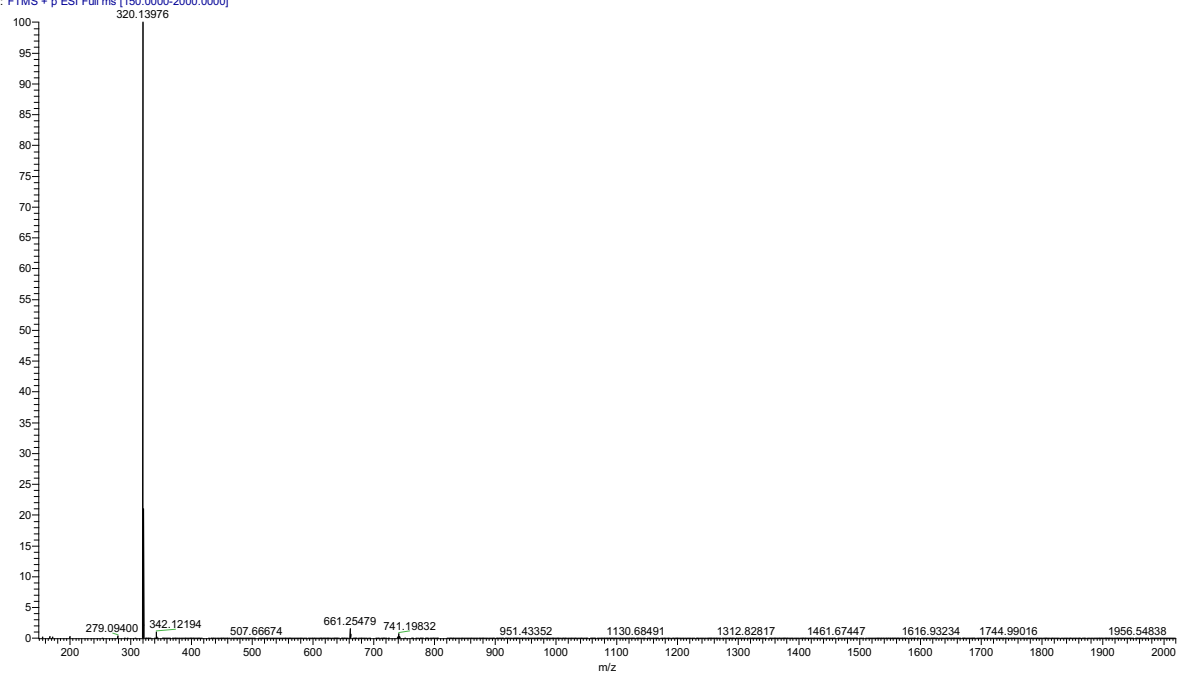


RT: 0.0000 - 11.2572

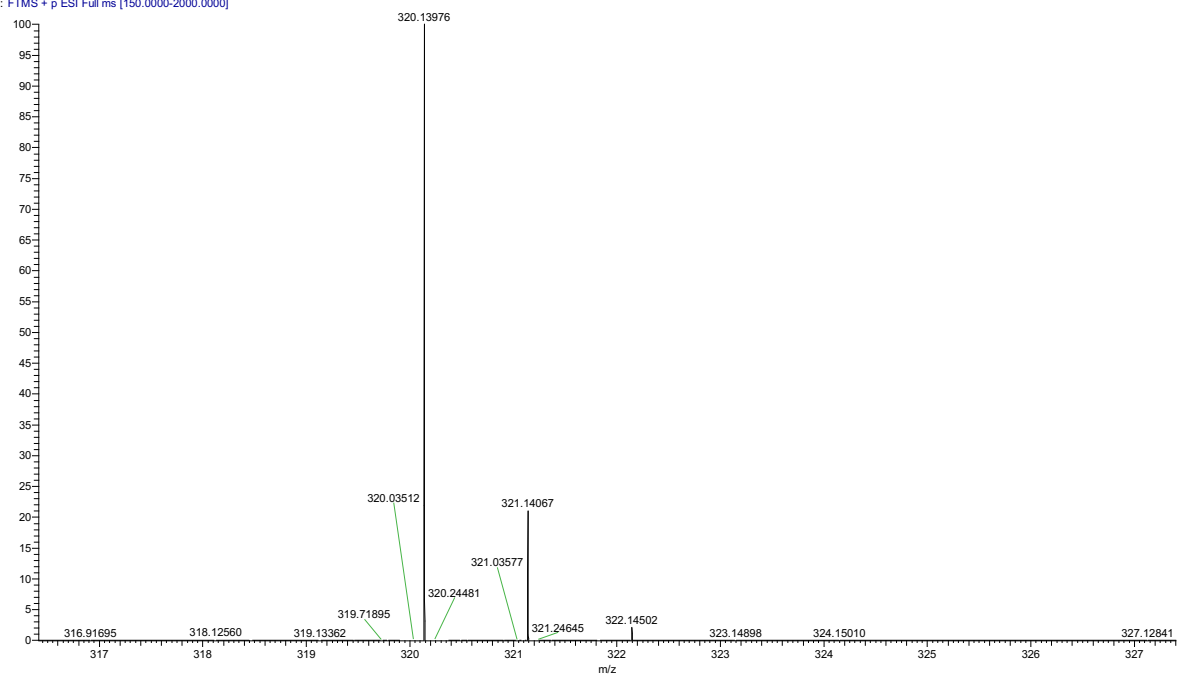


NL:
1.70E10
Base Peak
MS
ALM_DAI_2
5

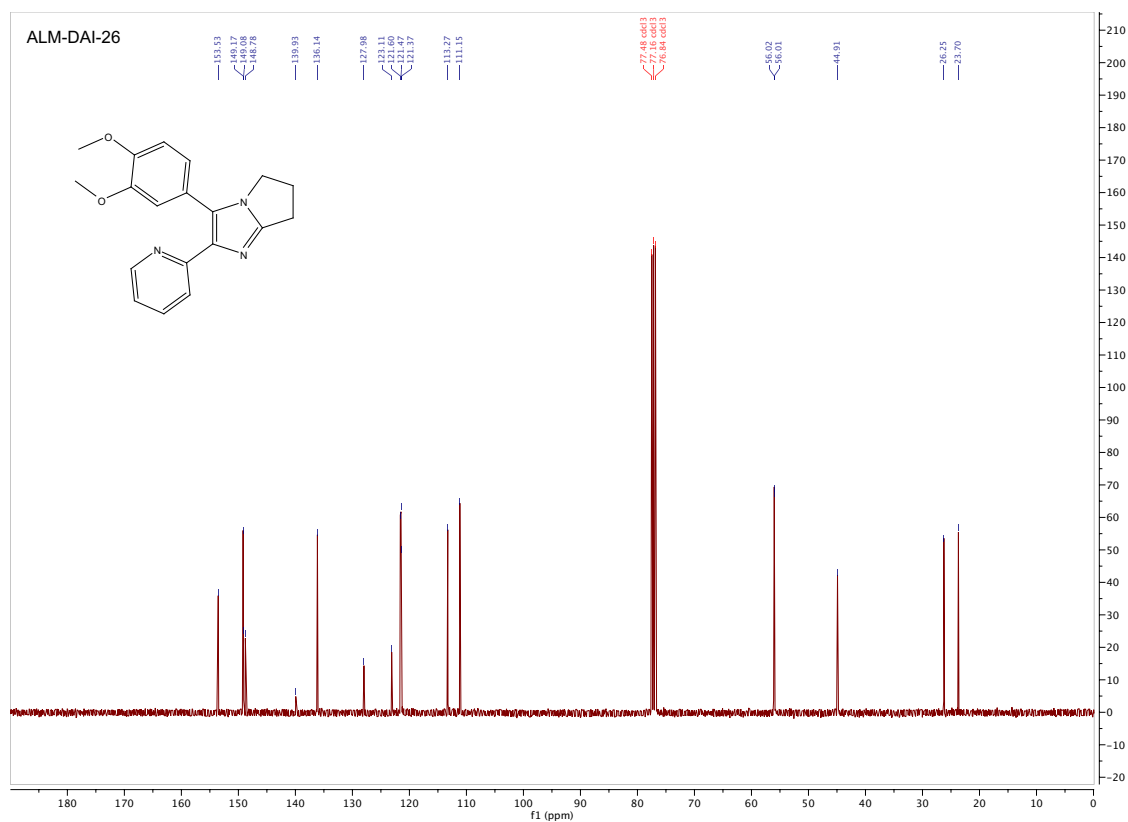
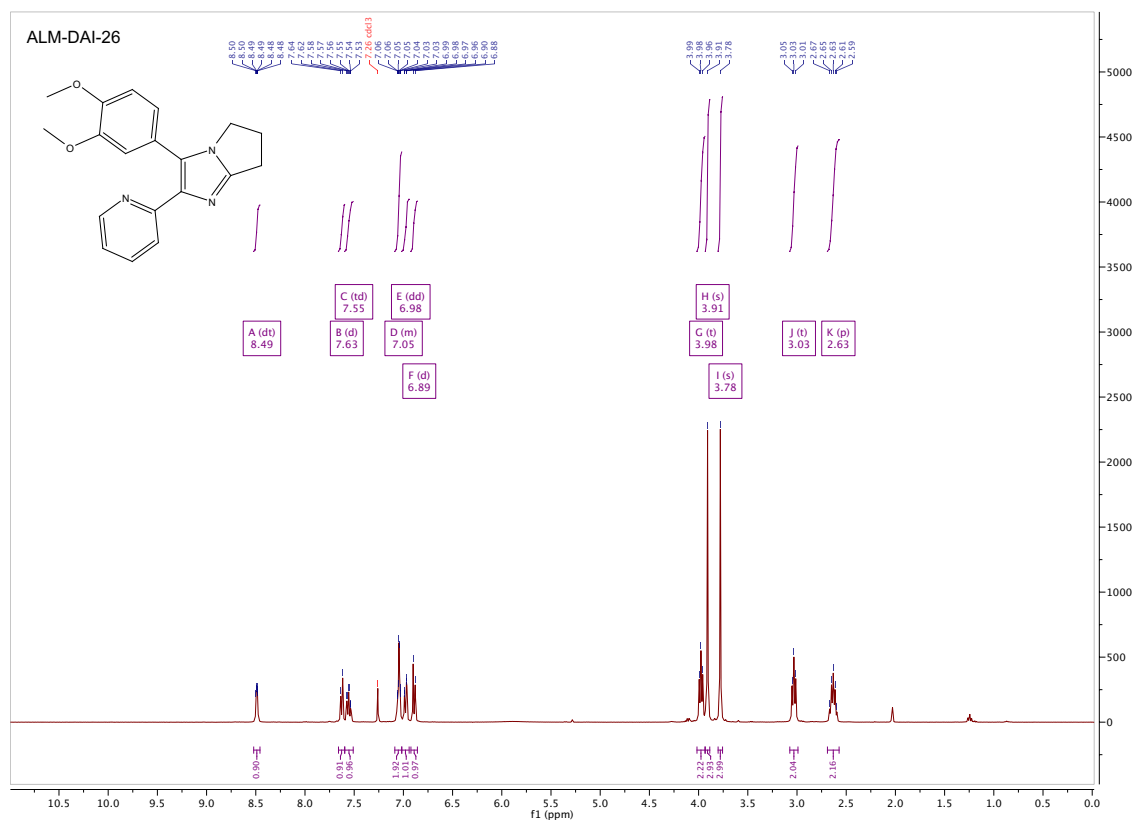
ALM_DAI_25 #778-839 RT: 4.00-4.30 AV: 62 NL: 8.69E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



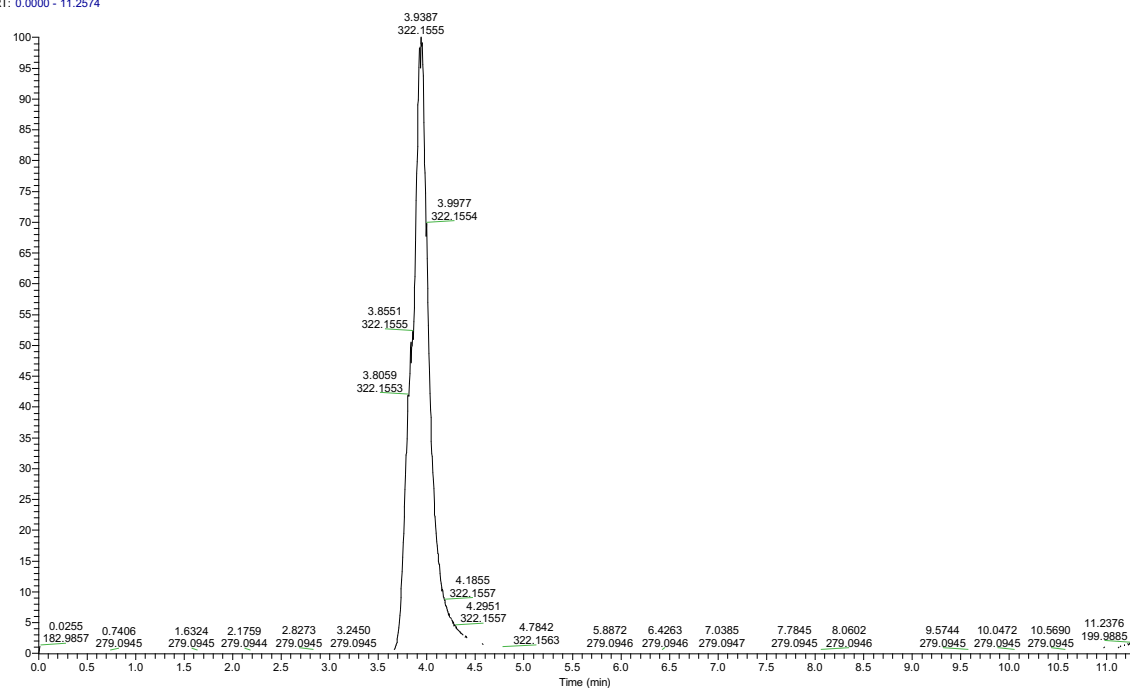
ALM_DAI_25 #778-839 RT: 4.00-4.30 AV: 62 NL: 8.69E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



3-(3,4-dimethoxyphenyl)-2-(pyridin-2-yl)-6,7-dihydro-5H-pyrrolo[1,2-a]imidazole (ALM-DAI-26).

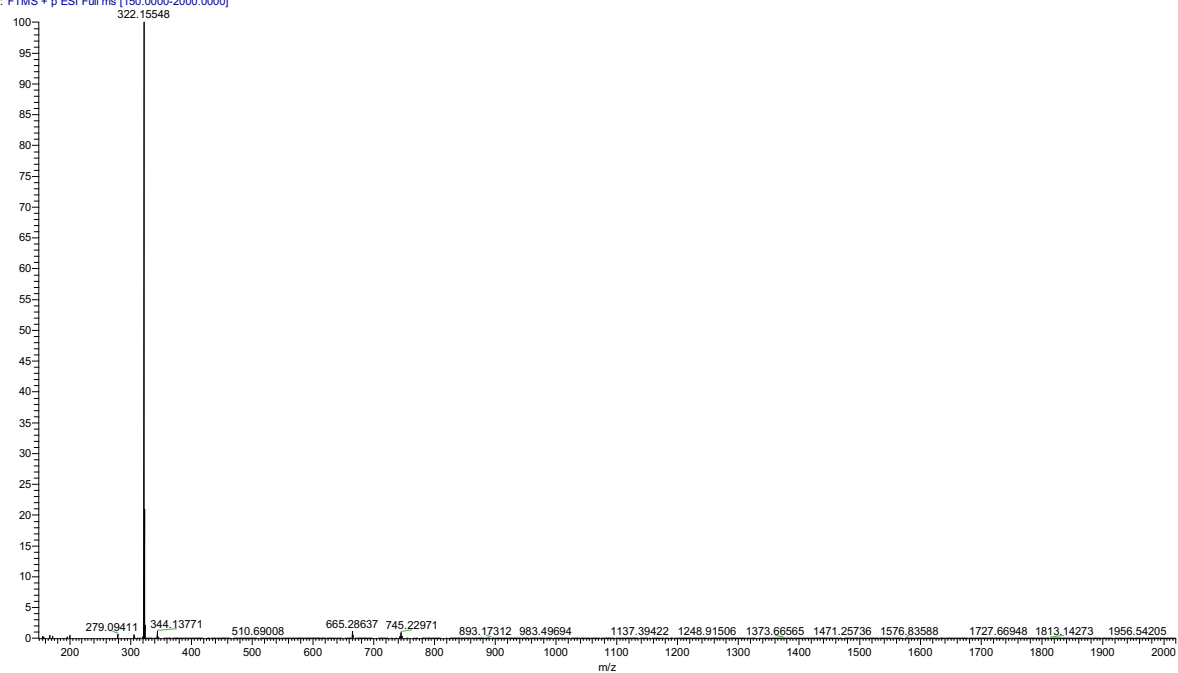


RT: 0.0000 - 11.2574

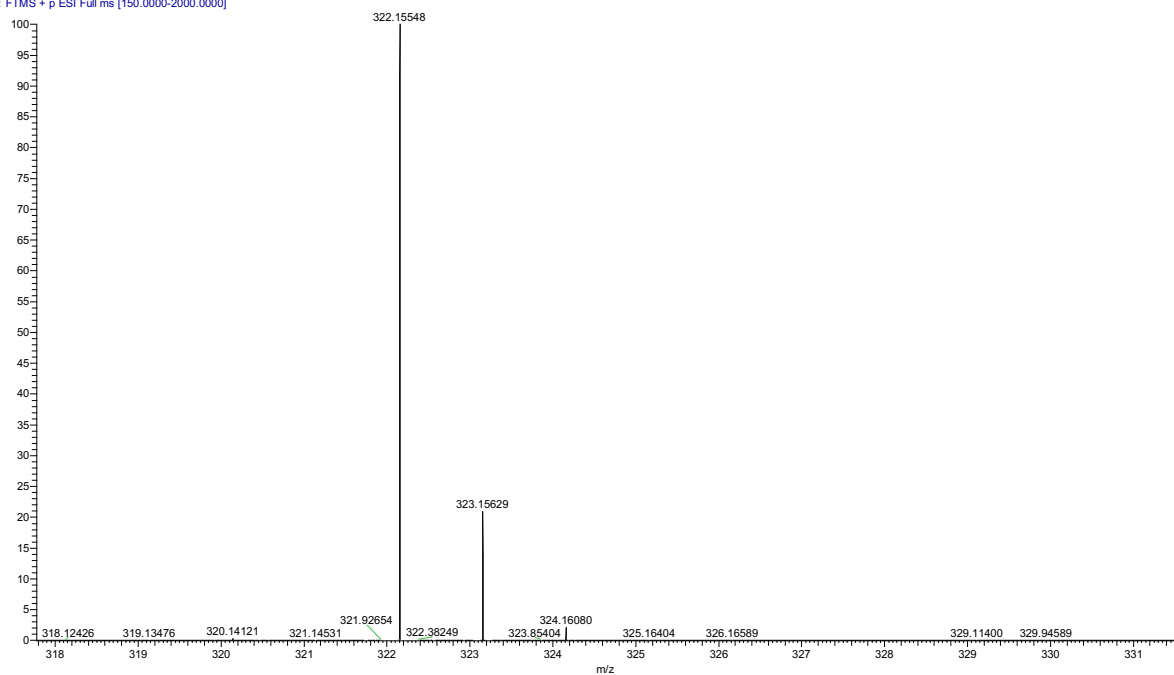


NL:
1.11E10
Base Peak
MS
ALM_DAI_2
6

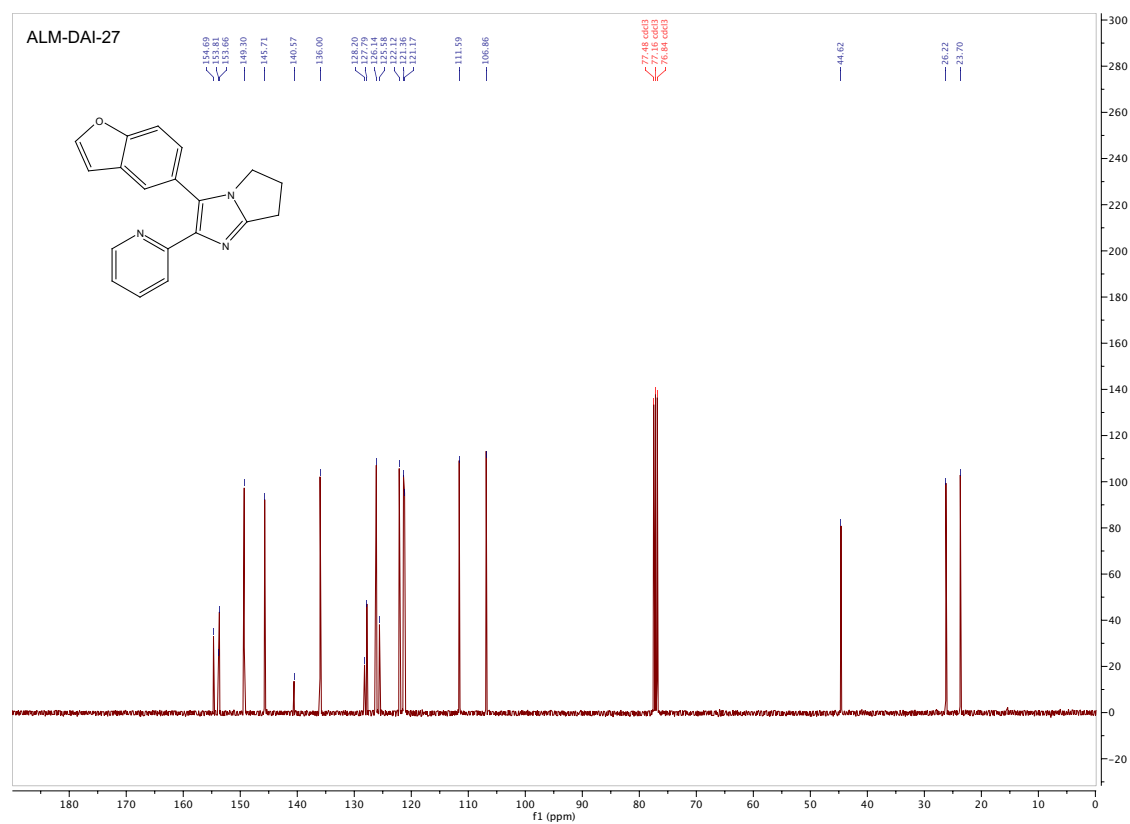
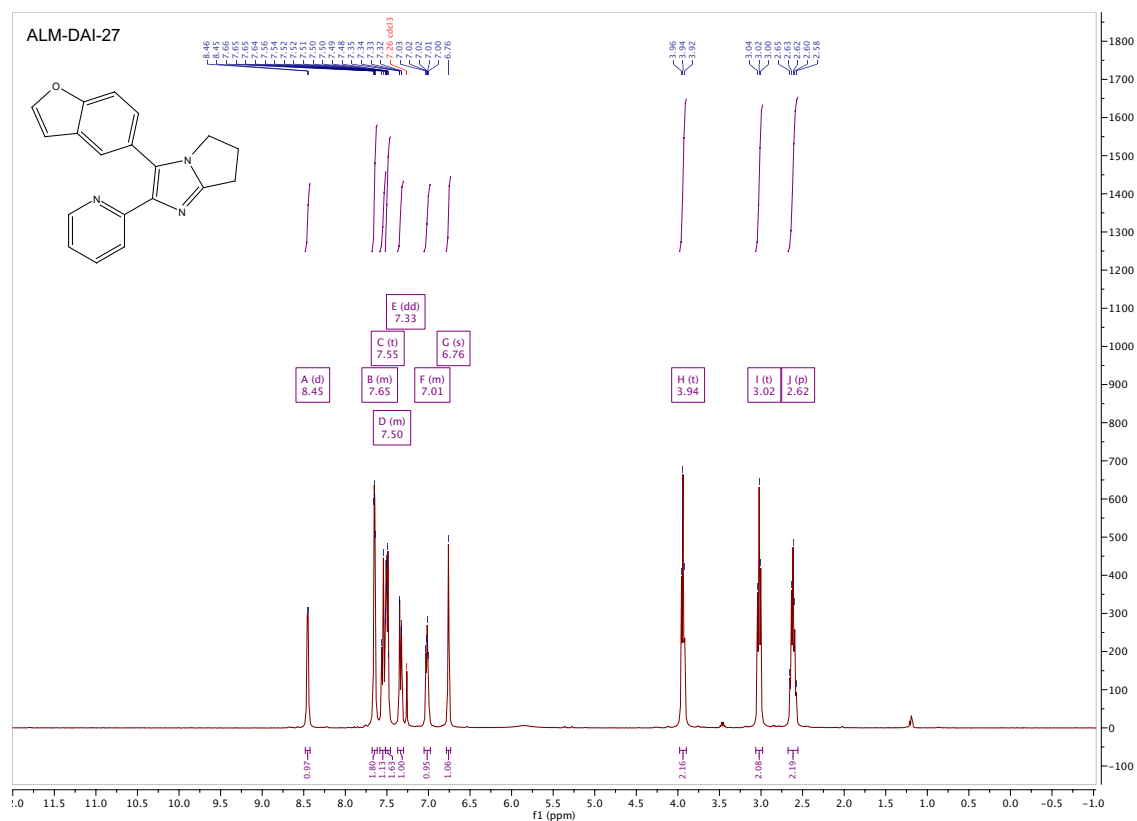
ALM_DAI_26 #731-796 RT: 3.76-4.08 AV: 66 NL: 5.85E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



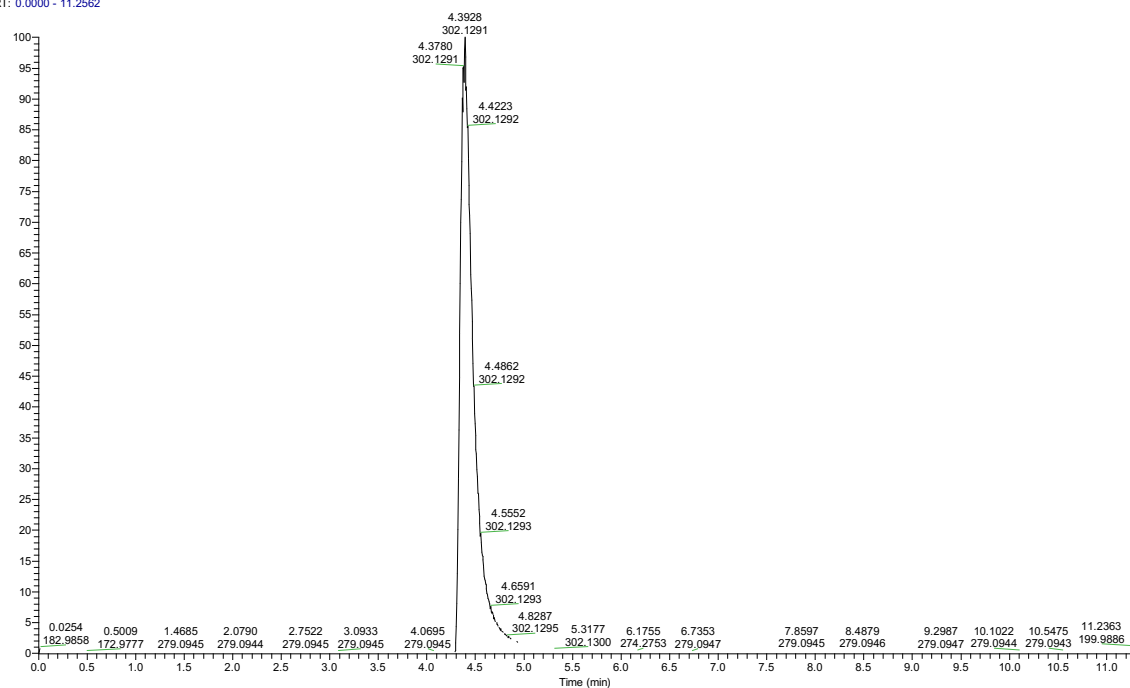
ALM_DAI_26 #731-796 RT: 3.76-4.08 AV: 66 NL: 5.85E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



3-(benzofuran-5-yl)-2-(pyridin-2-yl)-6,7-dihydro-5H-pyrrolo[1,2-a]imidazole (ALM-DAI-27).

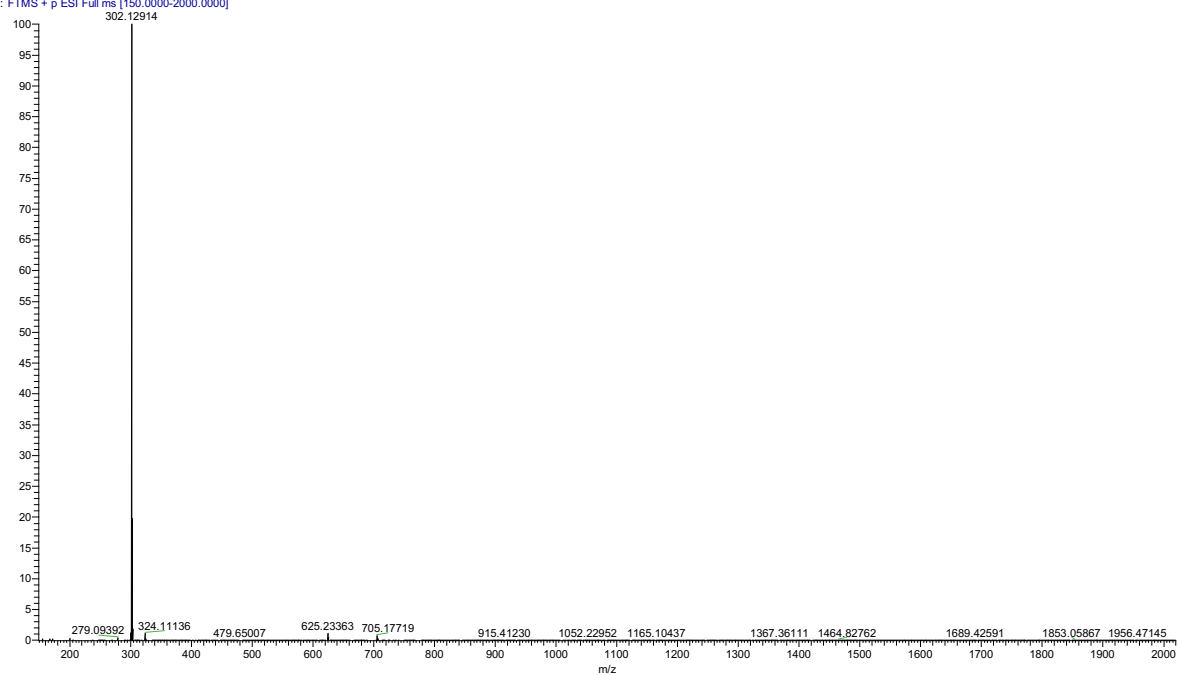


RT: 0.0000 - 11.2562

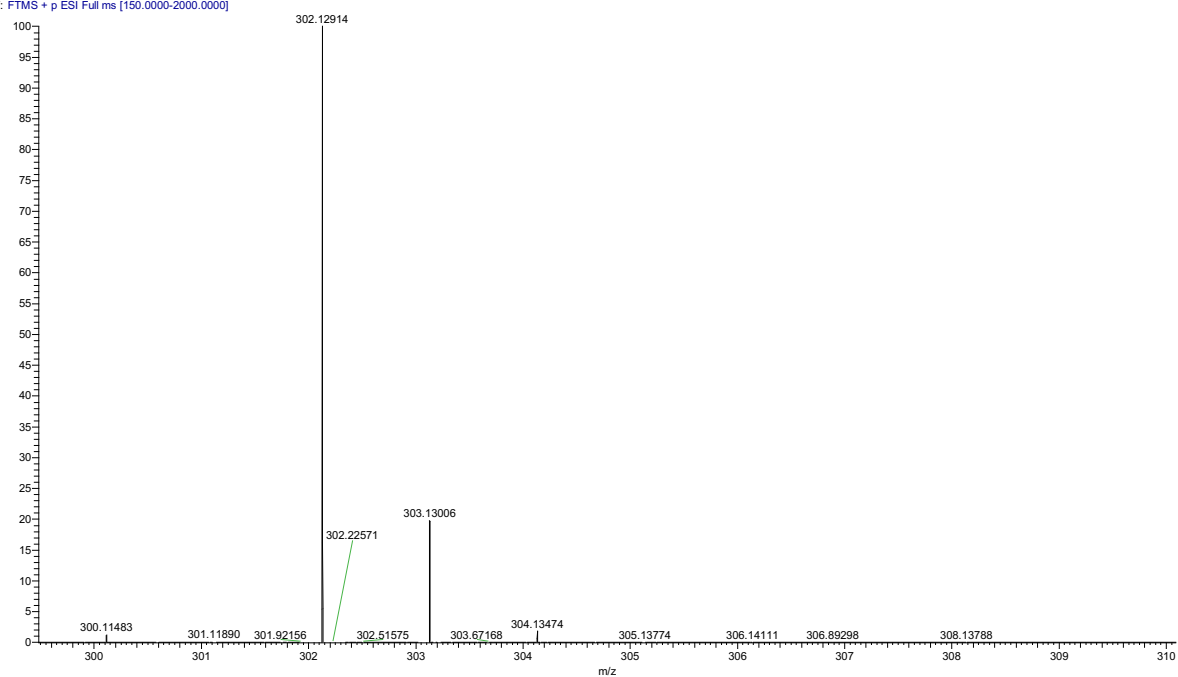


NL:
1.46E10
Base Peak
MS
ALM_DAI_2
7

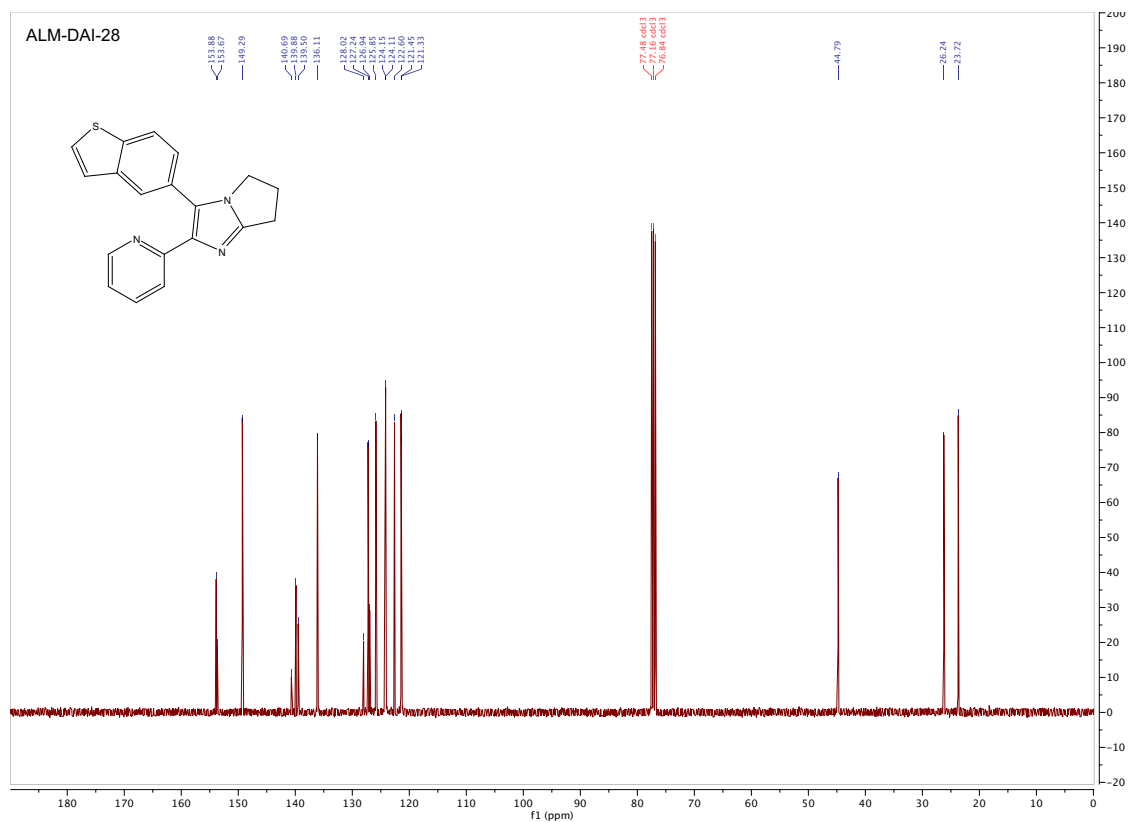
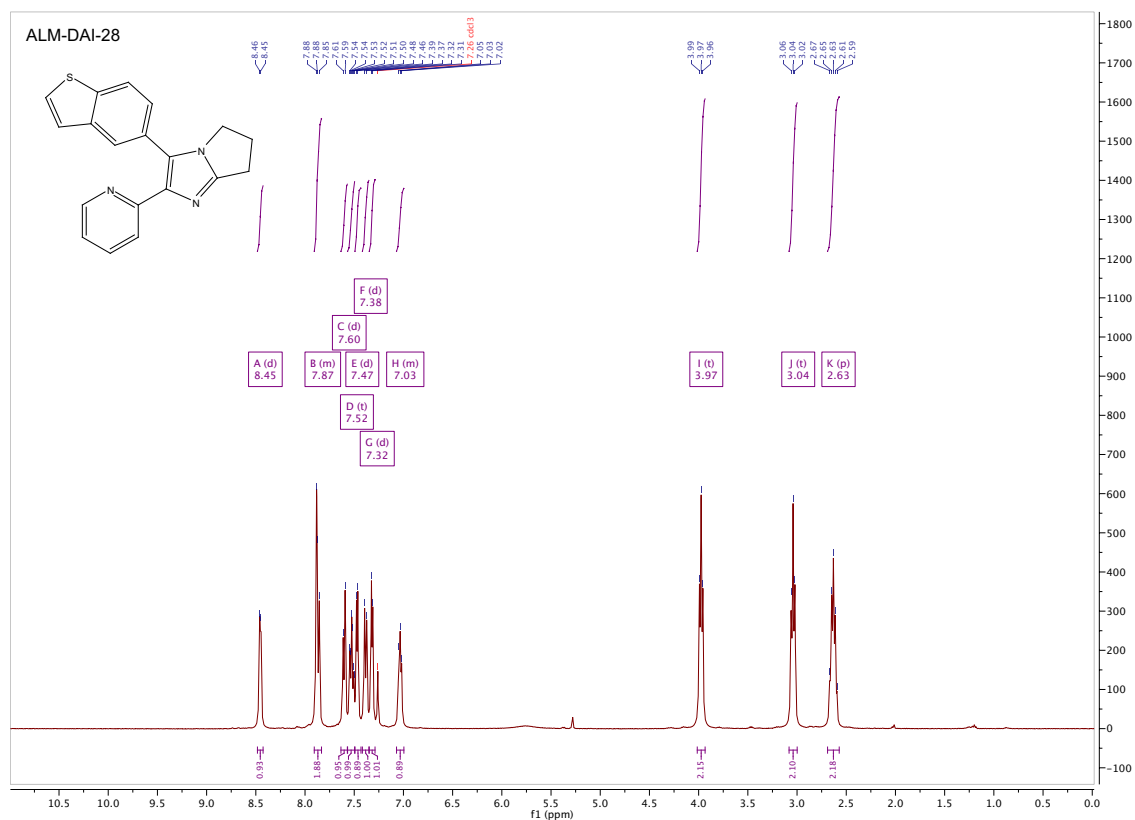
ALM_DAI_27 #940-877 RT: 4.32-4.50 AV: 38 NL: 9.39E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



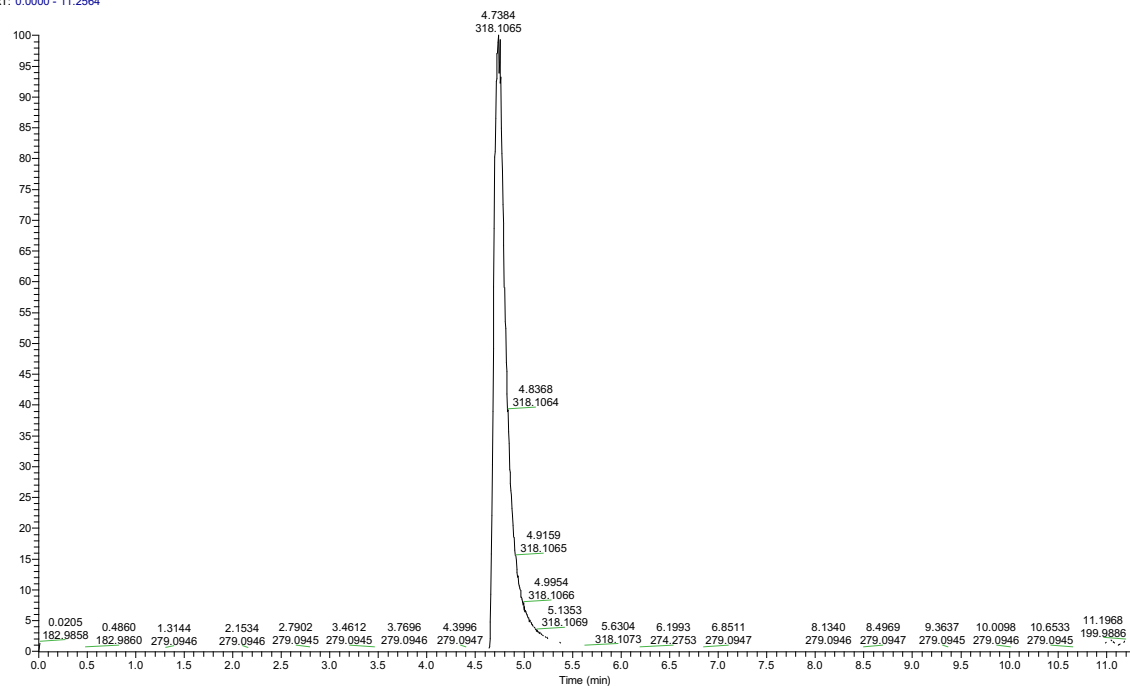
ALM_DAI_27 #840-877 RT: 4.32-4.50 AV: 38 NL: 9.39E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



3-(benzo[*b*]thiophen-5-yl)-2-(pyridin-2-yl)-6,7-dihydro-5*H*-pyrrolo[1,2-*a*]imidazole (ALM-DAI-28).

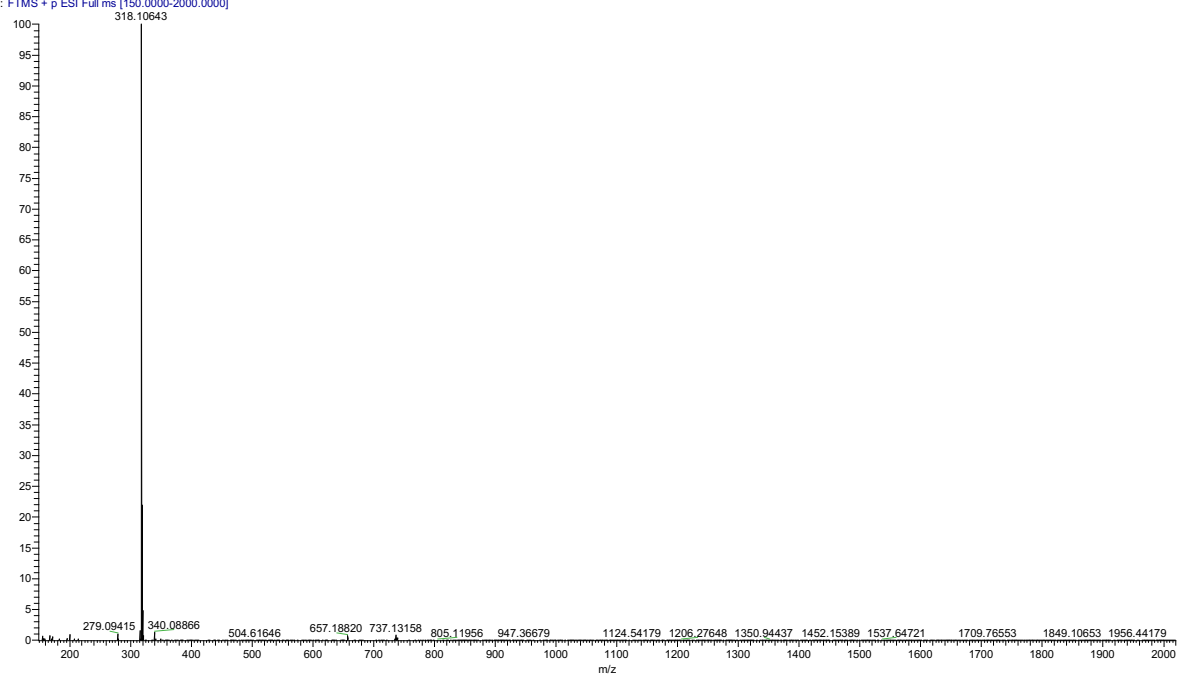


RT: 0.0000 - 11.2564

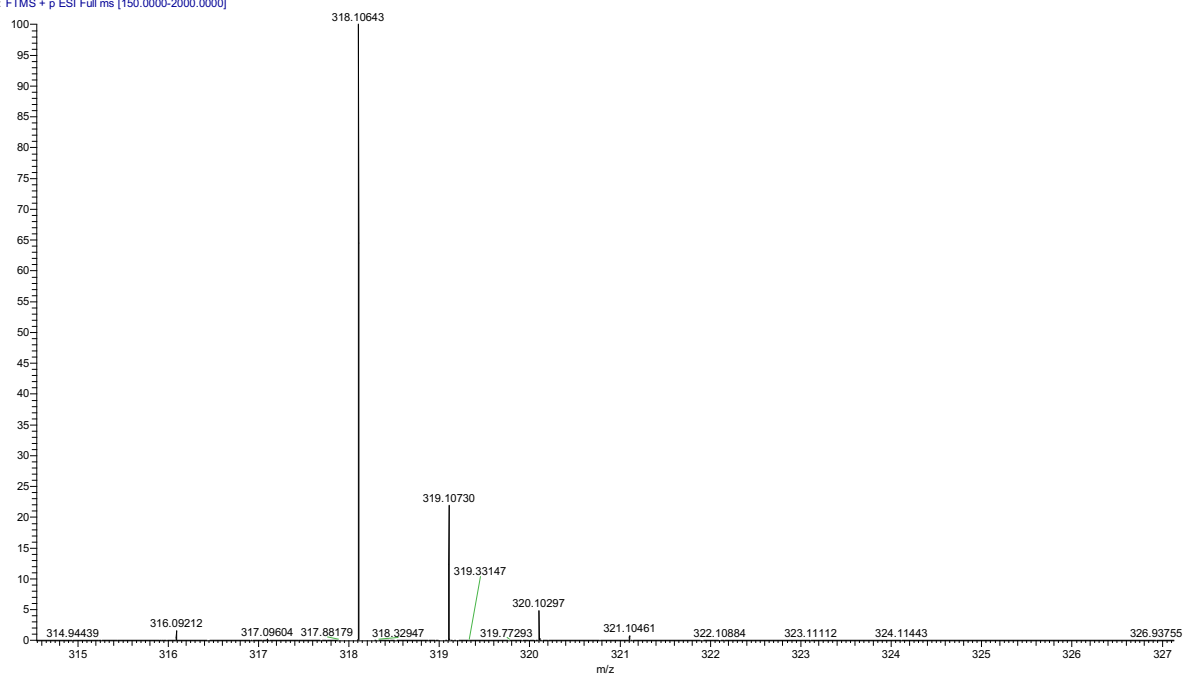


NL:
9.09E9
Base Peak
MS
ALM_DAI_2
8

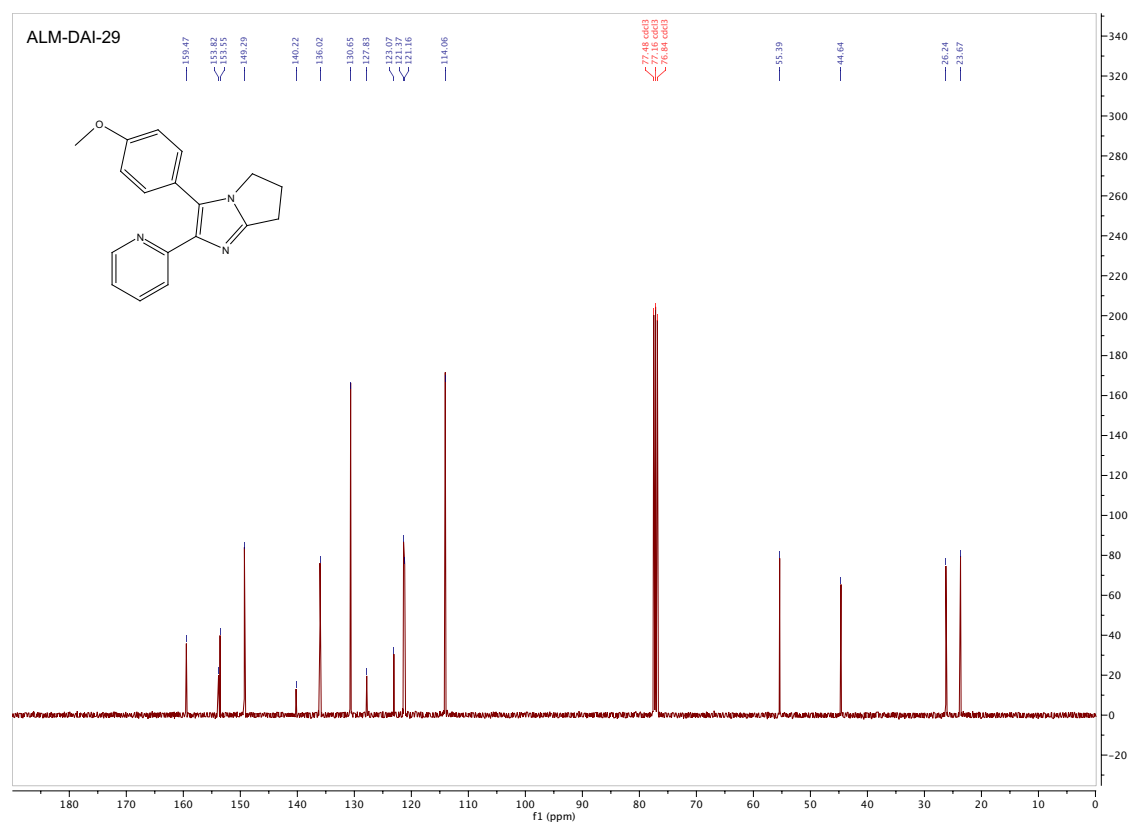
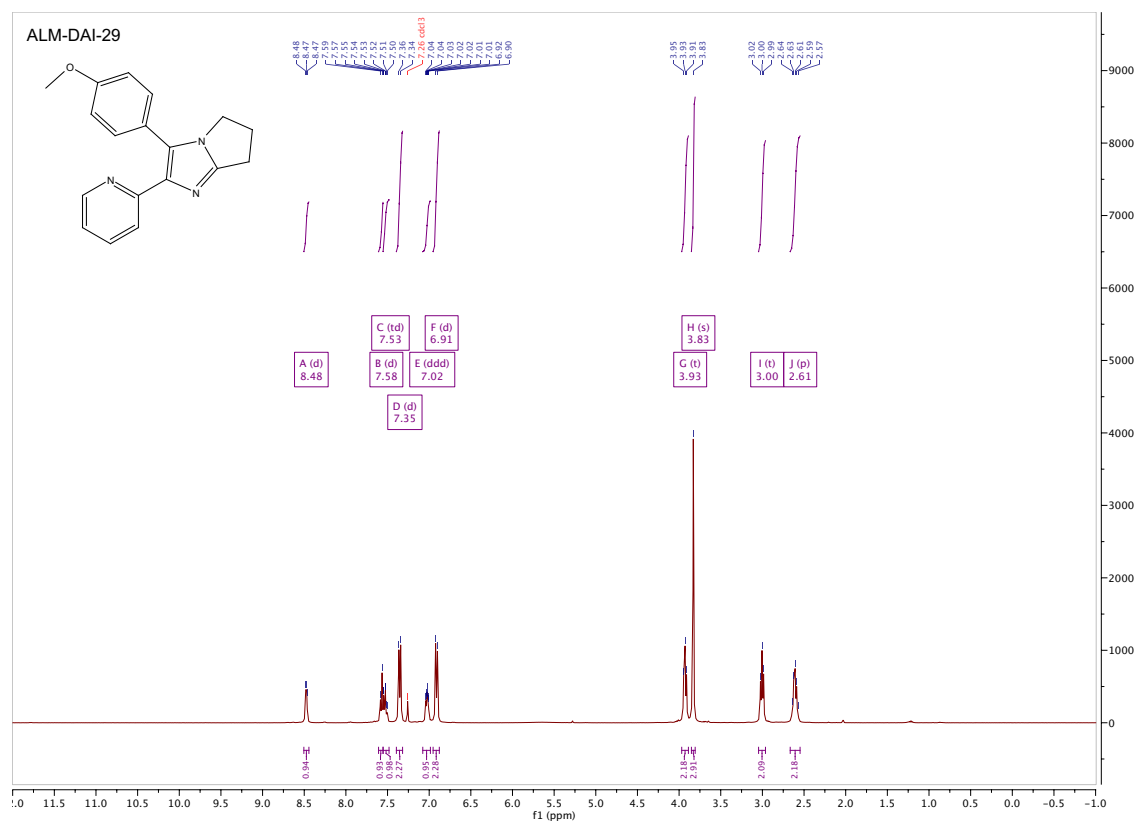
ALM_DAI_28 #905-963 RT: 4.65-4.94 AV: 59 NL: 4.08E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



ALM_DAI_28 #905-963 RT: 4.65-4.94 AV: 59 NL: 4.08E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



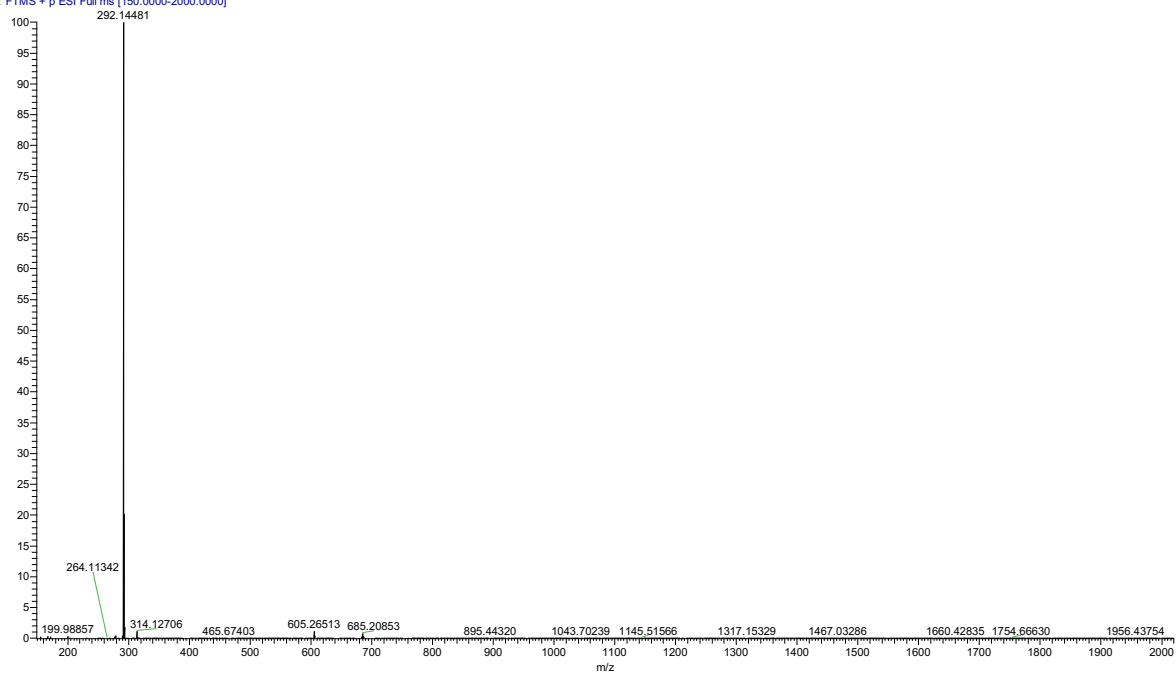
3-(4-methoxyphenyl)-2-(pyridin-2-yl)-6,7-dihydro-5H-pyrrolo[1,2-a]imidazole (ALM-DAI-29).



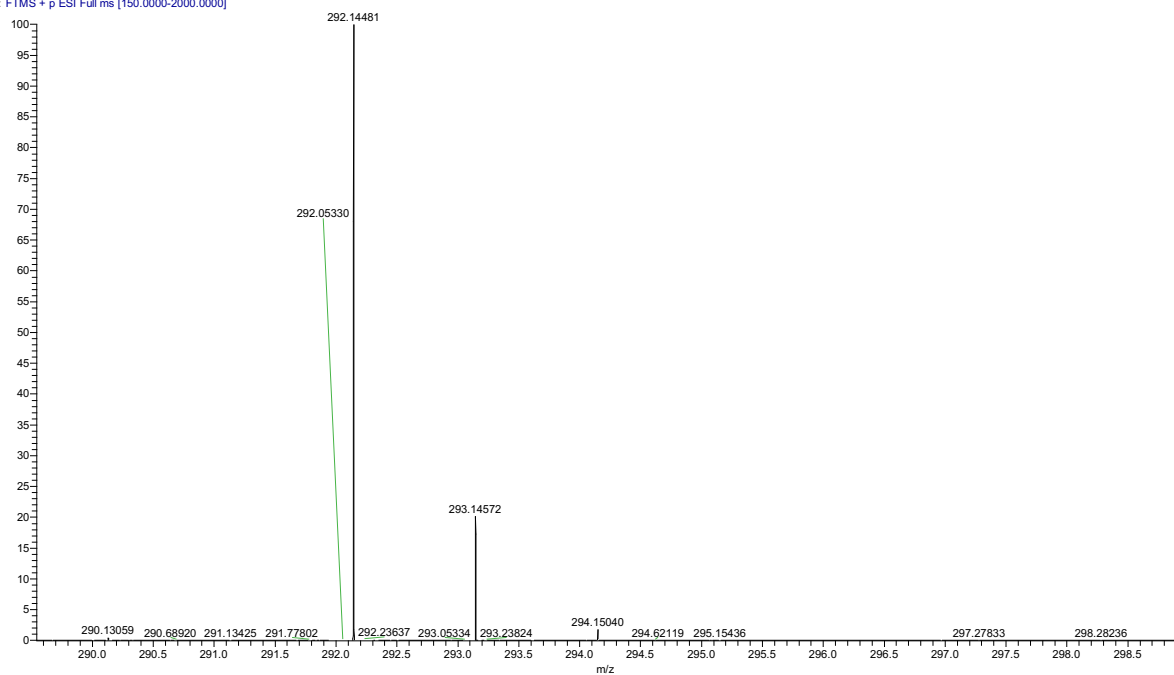
Chromatogram showing a major peak at 4.1823 minutes. The y-axis represents intensity from 0 to 100, and the x-axis represents time in minutes from 0.0 to 11.0. Numerous small peaks are labeled with retention times and molecular weights.

Retention Time (min)	Molecular Weight
0.0205	162.9658
0.5871	279.0946
1.3290	279.0945
2.0592	279.0947
3.0417	279.0946
3.8285	279.0945
4.1823	292.1449
4.1922	292.1449
4.2118	292.1449
4.4438	292.1452
4.5733	292.1451
5.0673	292.1456
5.8495	279.0946
6.8227	279.0946
7.7838	279.0946
8.3458	279.0946
8.7750	279.0946
9.8801	279.0945
10.6658	279.0945
11.2487	199.9855

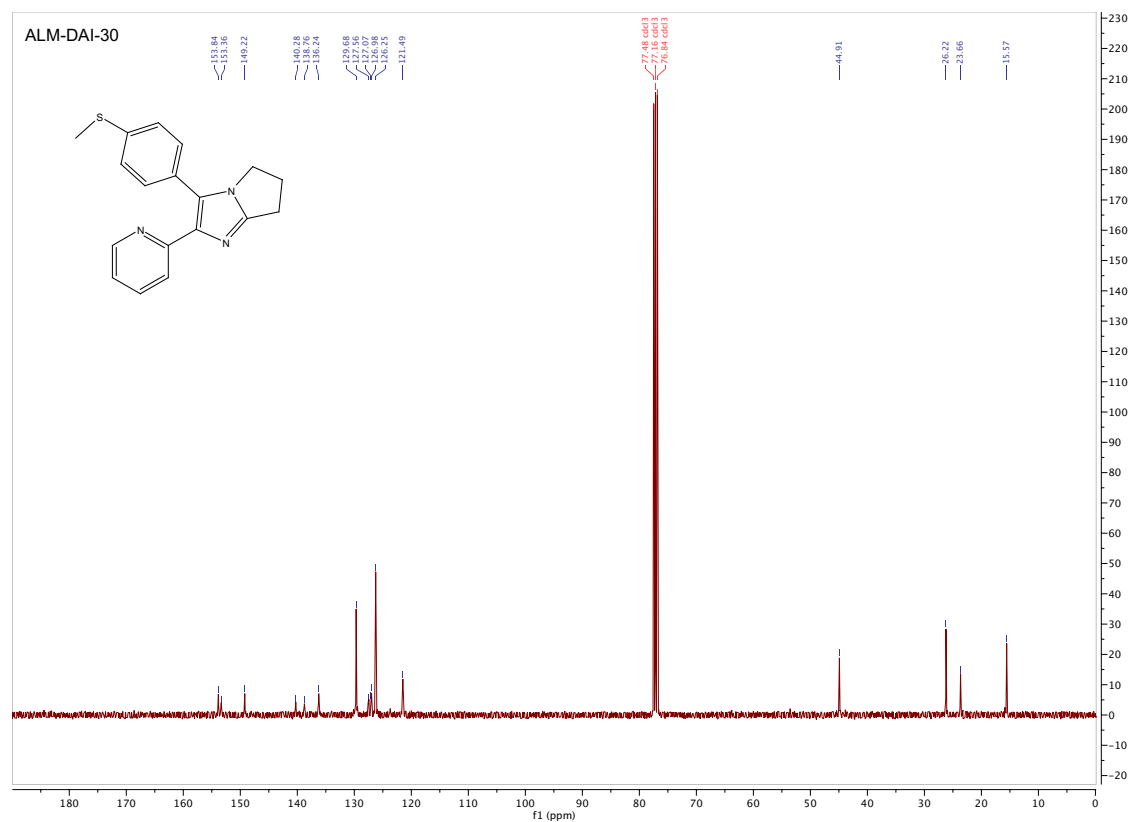
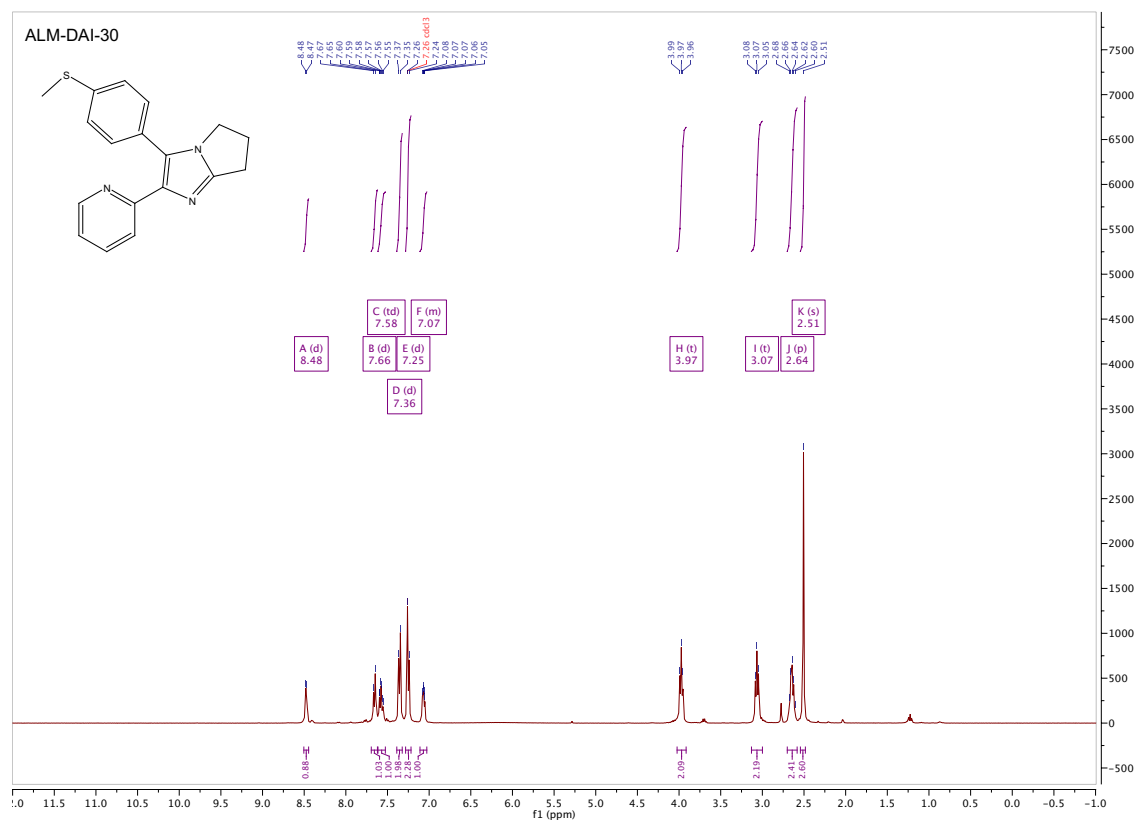
ALM_DAI_29 #797-838 RT: 4.10-4.30 AV: 42 NL: 8.46E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



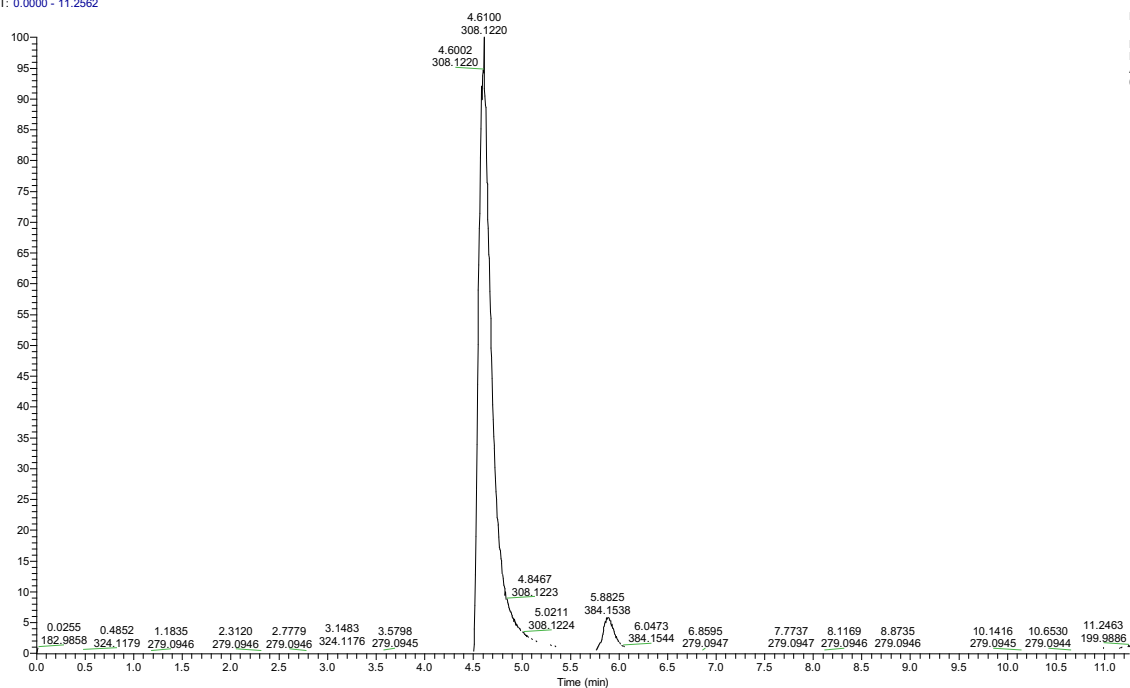
ALM_DAI_29 #797-838 RT: 4.10-4.30 AV: 42 NL: 8.46E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



3-(4-(methylthio)phenyl)-2-(pyridin-2-yl)-6,7-dihydro-5H-pyrrolo[1,2-a]imidazole (ALM-DAI-30).

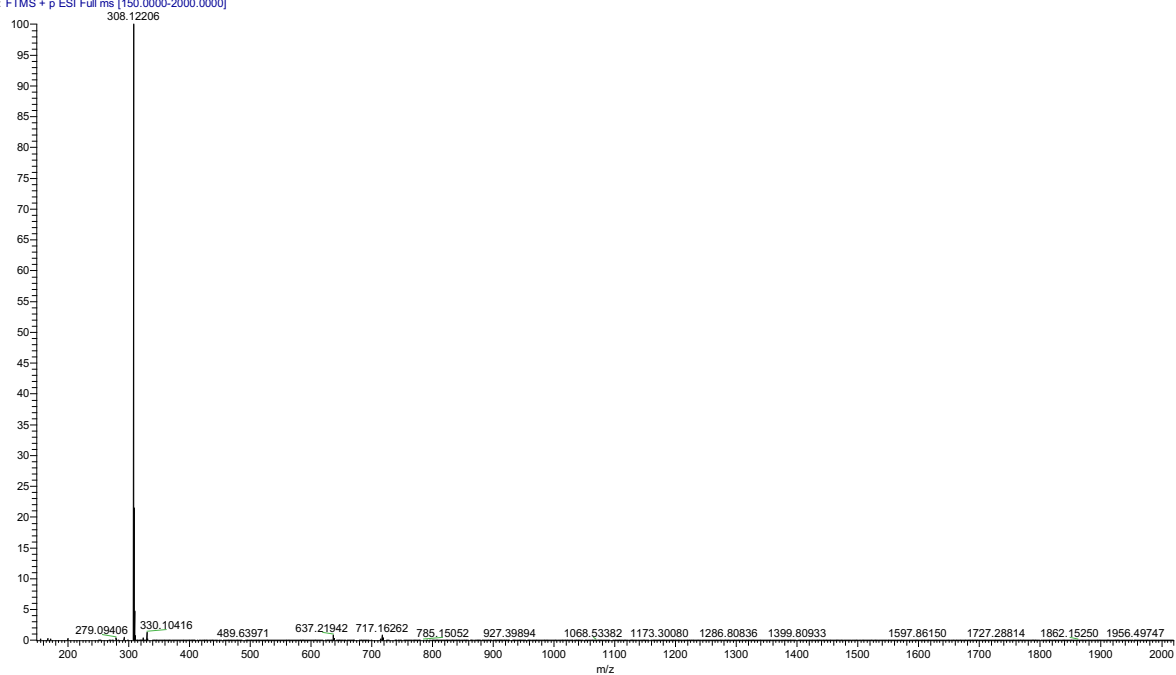


RT: 0.0000 - 11.2562

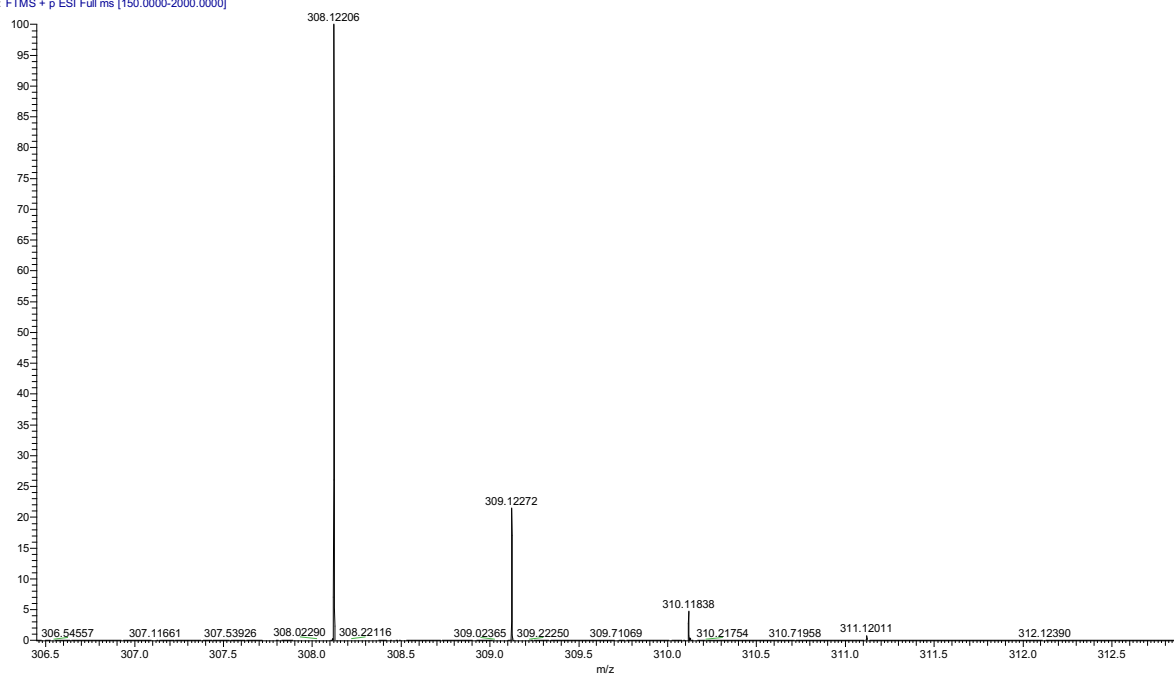


NL:
1.36E10
Base Peak
MS
ALM_DAI_3
0

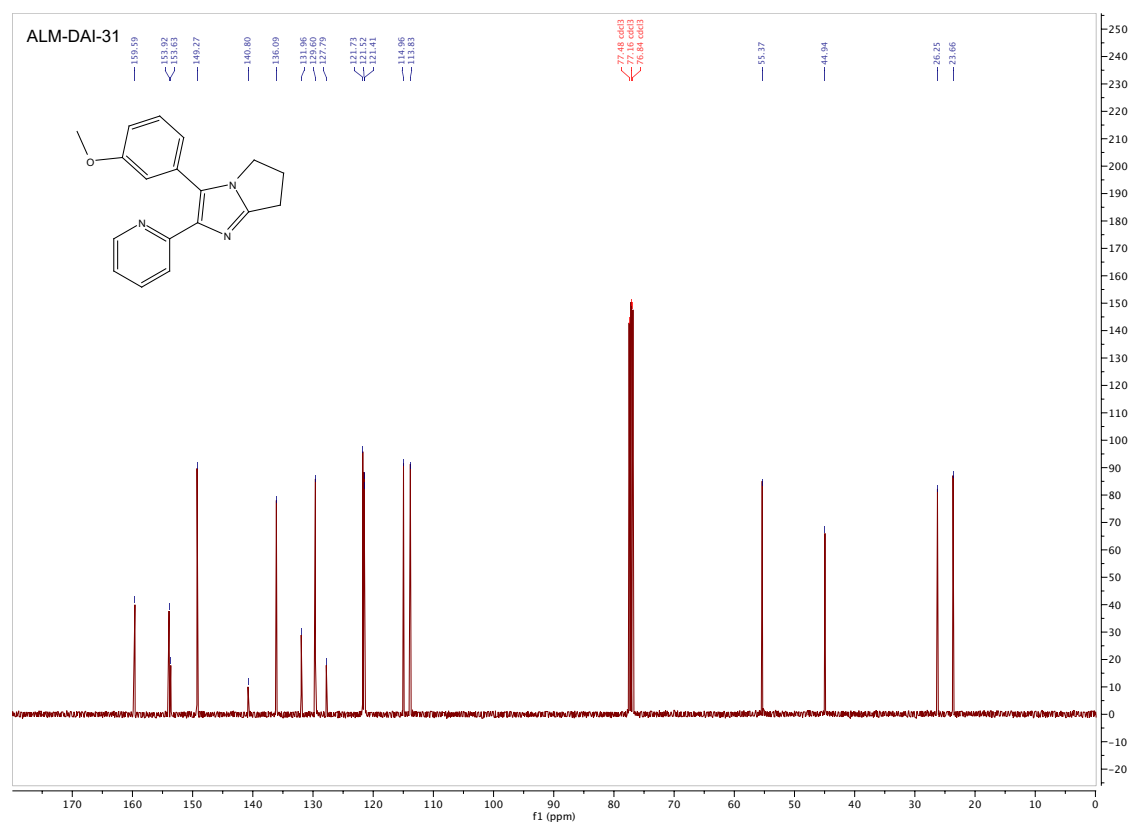
ALM_DAI_30 #884-921 RT: 4.54-4.72 AV: 38 NL: 8.43E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



ALM_DAI_30 #884-921 RT: 4.54-4.72 AV: 38 NL: 8.43E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]

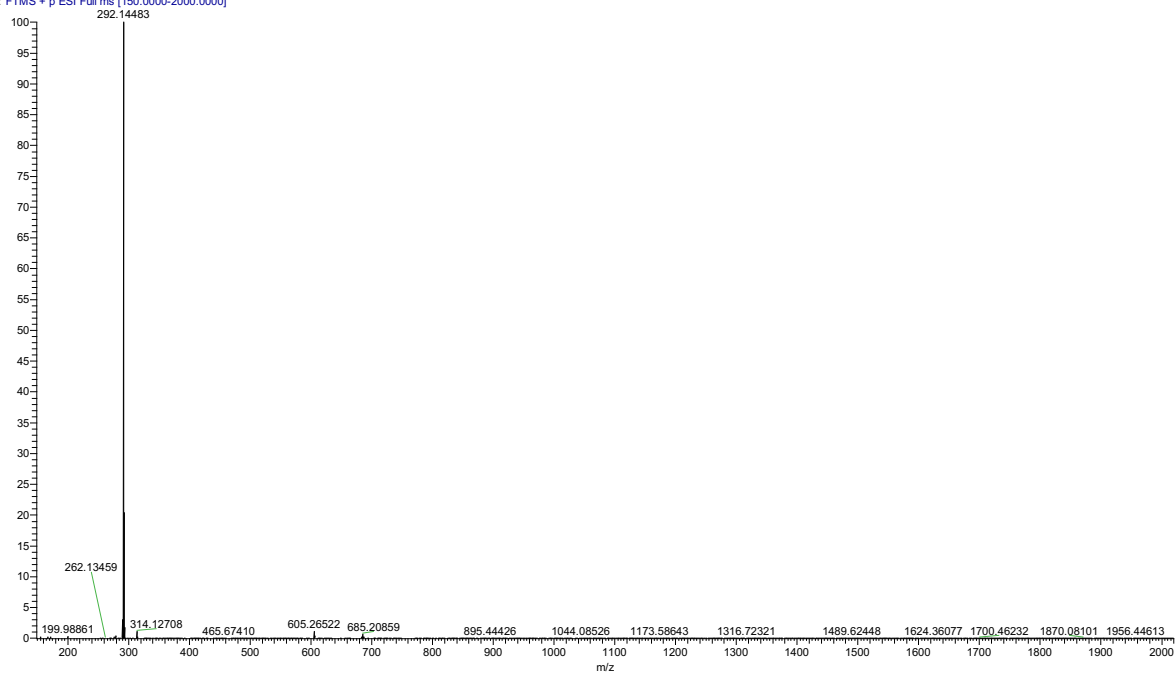


3-(3-methoxyphenyl)-2-(pyridin-2-yl)-6,7-dihydro-5H-pyrrolo[1,2-a]imidazole (ALM-DAI-31).

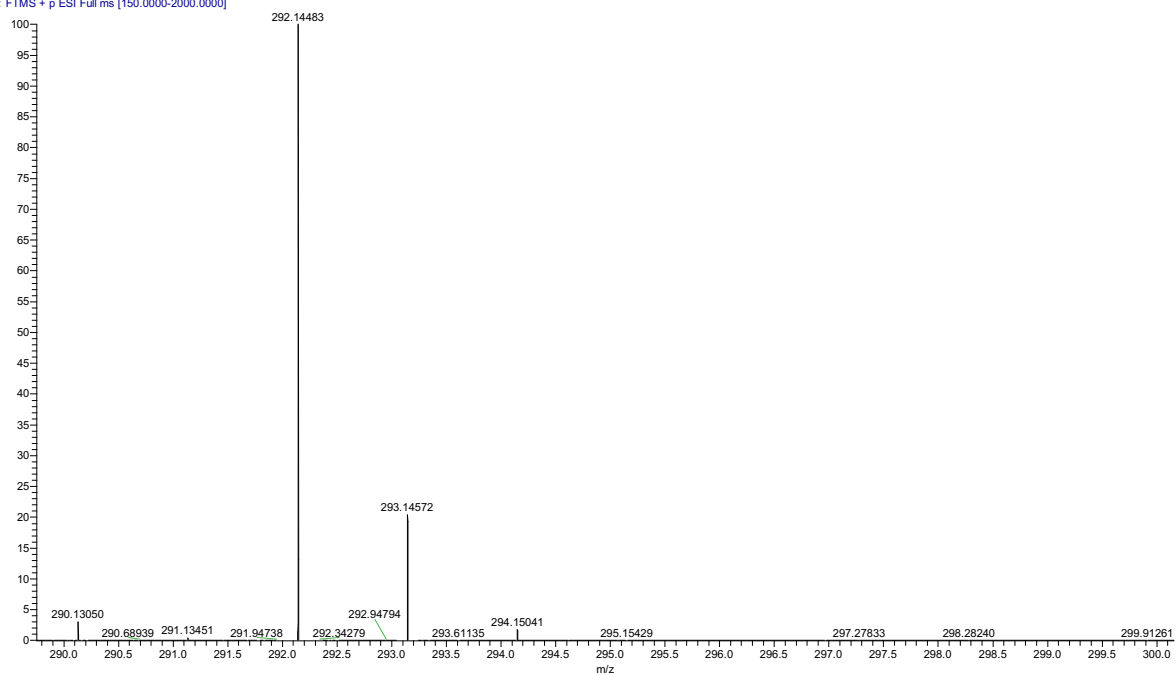


Chromatogram showing a major peak at 4.1472 minutes and several minor peaks. The x-axis is Time (min) from 0.0 to 11.0, and the y-axis is relative intensity from 0 to 100. The major peak is labeled with retention times 4.1472 and 292.1447. Other labeled peaks include 4.1816, 292.1449, 4.3835, 292.1451, 4.4578, 292.1452, 4.6174, 292.1452, 5.1165, 292.1457, 6.1738, 279.0946, 6.8624, 279.0947, 8.0637, 279.0946, 8.3550, 279.0948, 9.5230, 279.0947, 10.0969, 279.0945, 10.6580, 279.0945, and 11.0315, 167.0134.

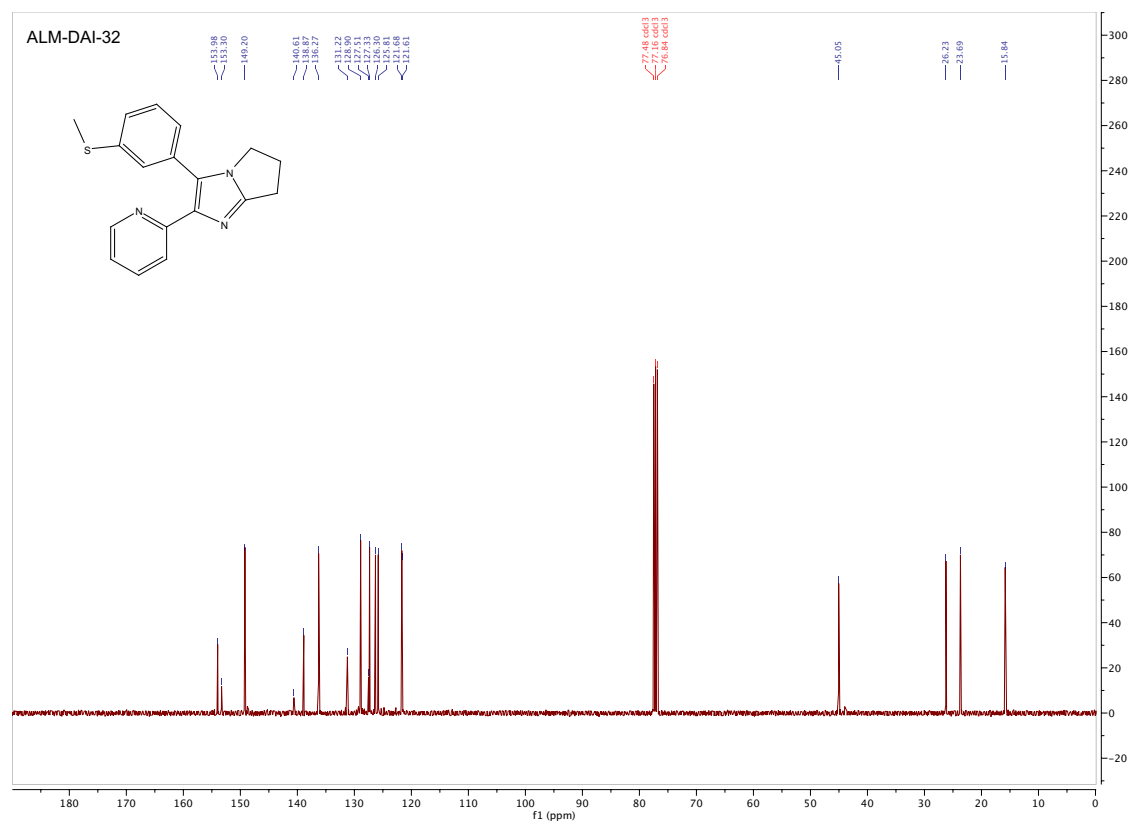
ALM_DAI_31 #793-840 RT: 4.08-4.31 AV: 48 NL: 8.59E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]
292 14483



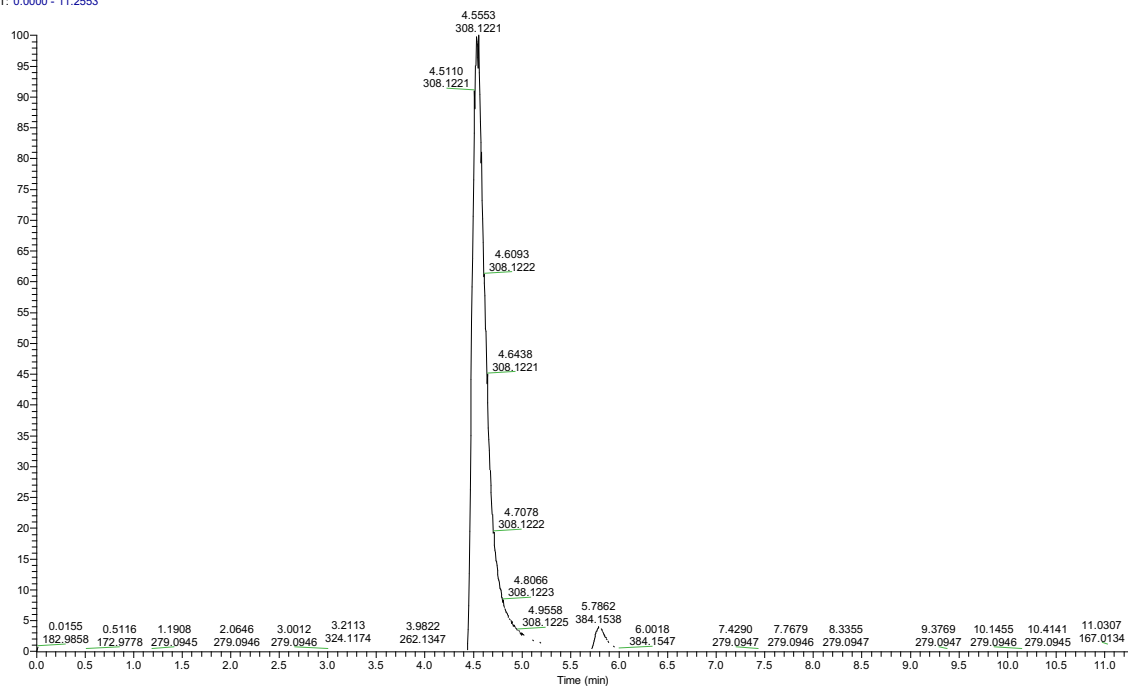
ALM_DAI_31 #793-840 RT: 4.08-4.31 AV: 48 NL: 8.59E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



3-(3-(methylthio)phenyl)-2-(pyridin-2-yl)-6,7-dihydro-5H-pyrrolo[1,2-a]imidazole (ALM-DAI-32).

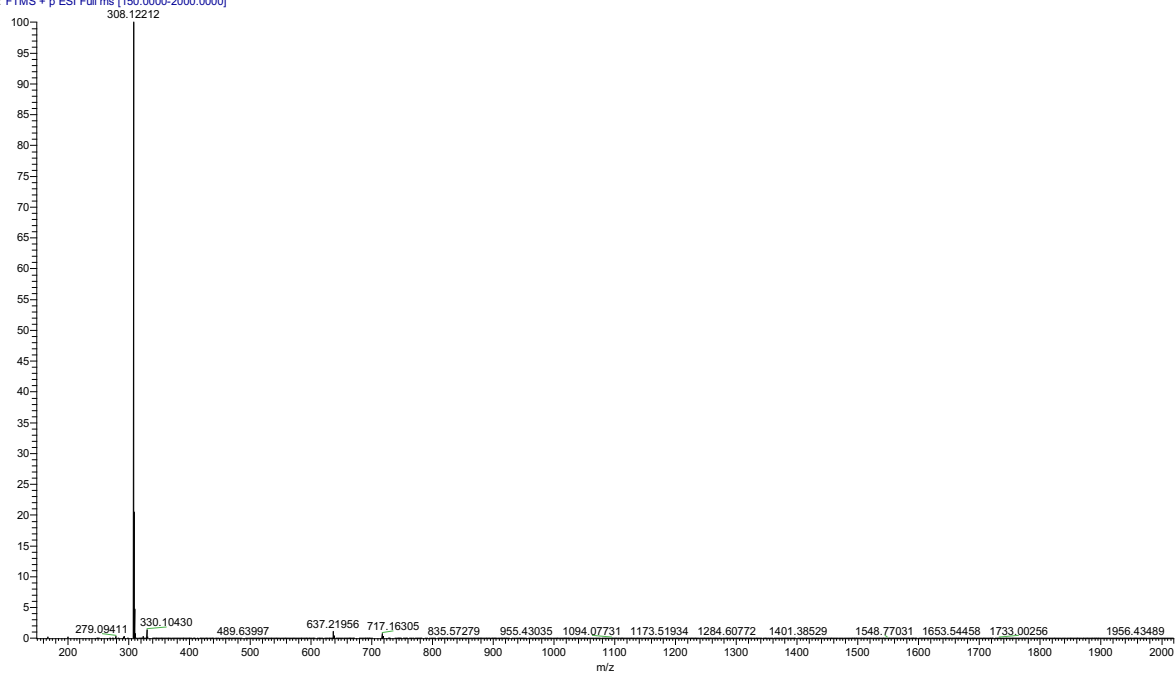


RT: 0.0000 - 11.2553

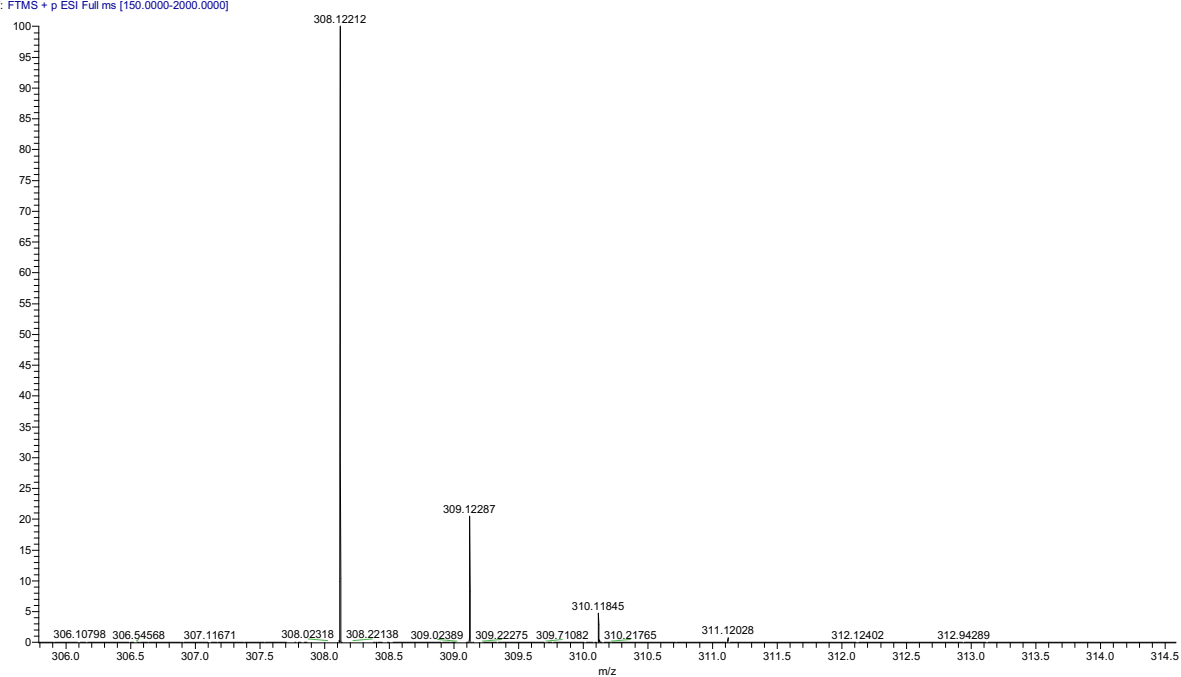


NL:
1.67E10
Base Peak
MS
ALM_DAI_3
2

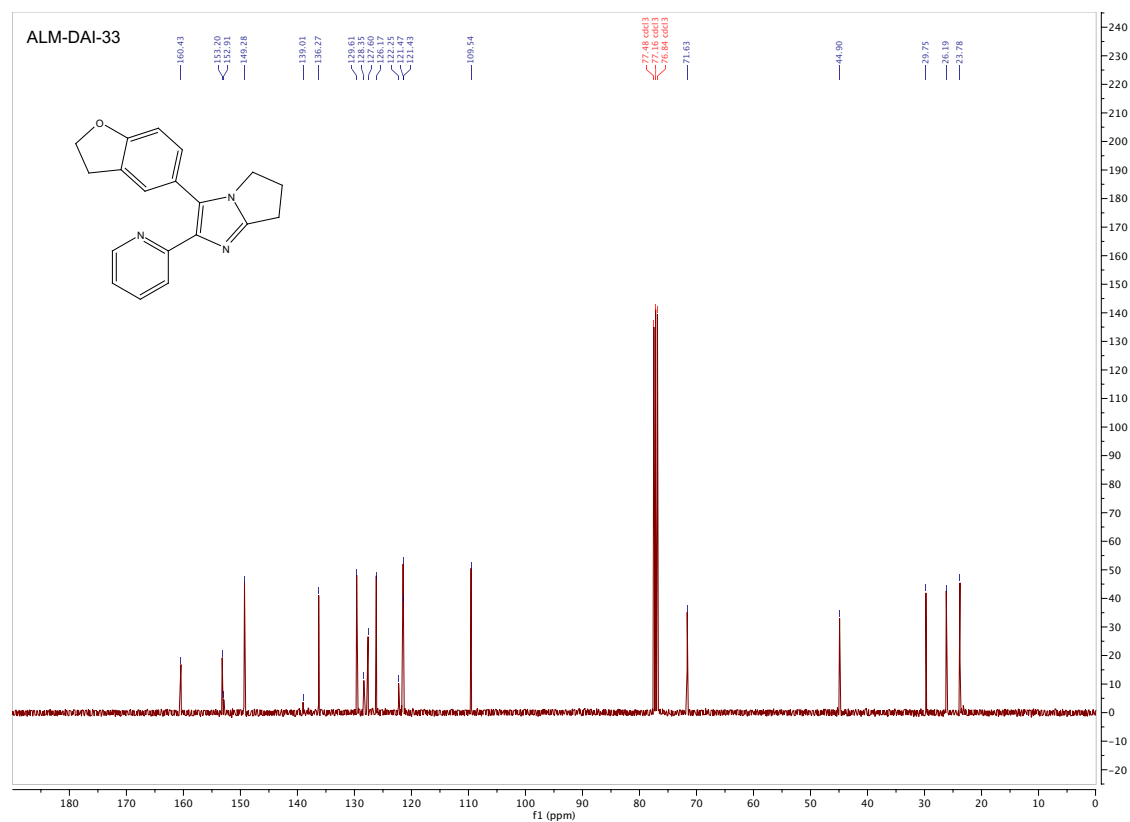
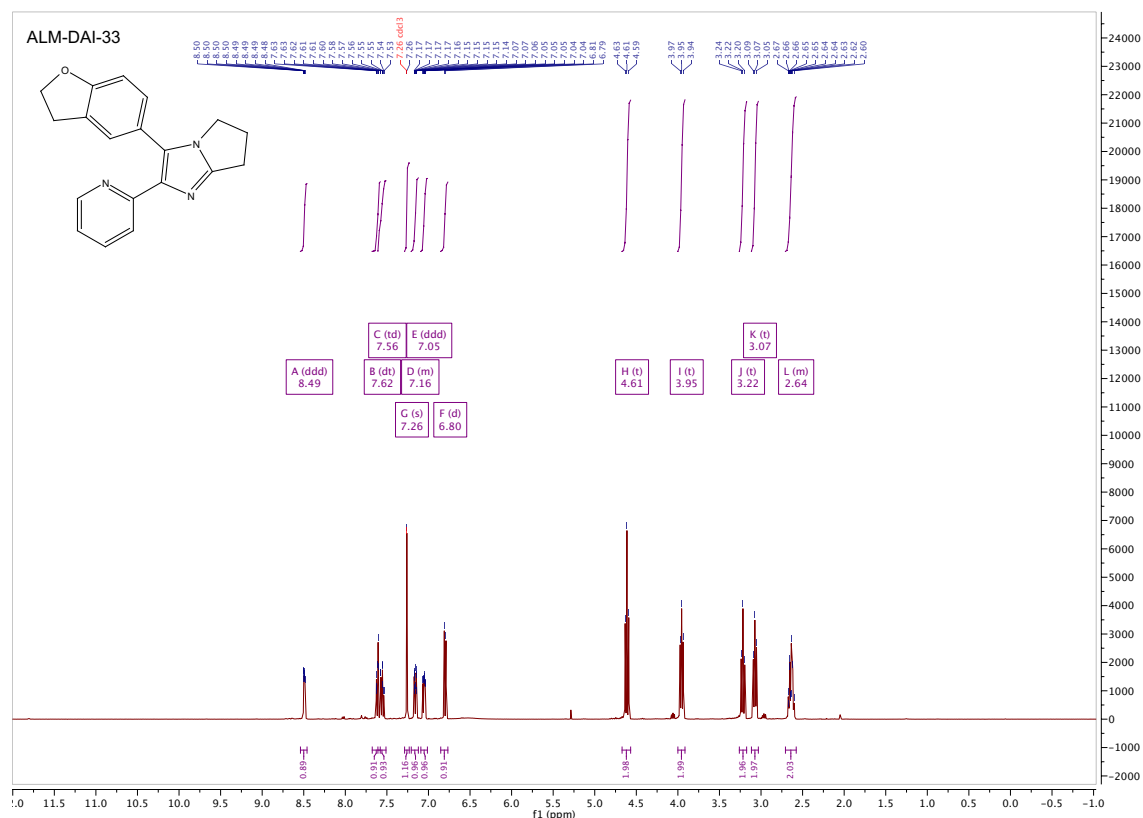
ALM_DAI_32 #670-912 RT: 4.47-4.68 AV: 43 NL: 1.04E10
T: FTMS + p ESI Full ms [150.0000-2000.0000]



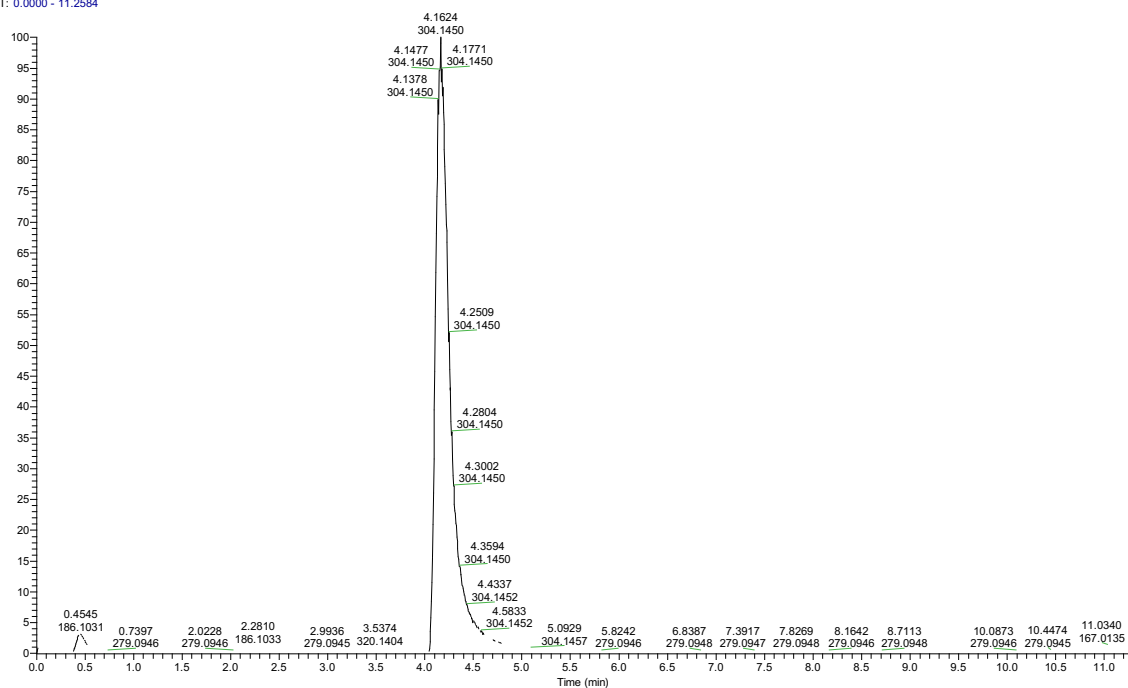
ALM_DAI_32 #870-912 RT: 4.47-4.68 AV: 43 NL: 1.04E10
T: FTMS + p ESI Full ms [150.0000-2000.0000]



3-(2,3-dihydrobenzofuran-5-yl)-2-(pyridin-2-yl)-6,7-dihydro-5*H*-pyrrolo[1,2-*a*]imidazole (ALM-DAI-33).

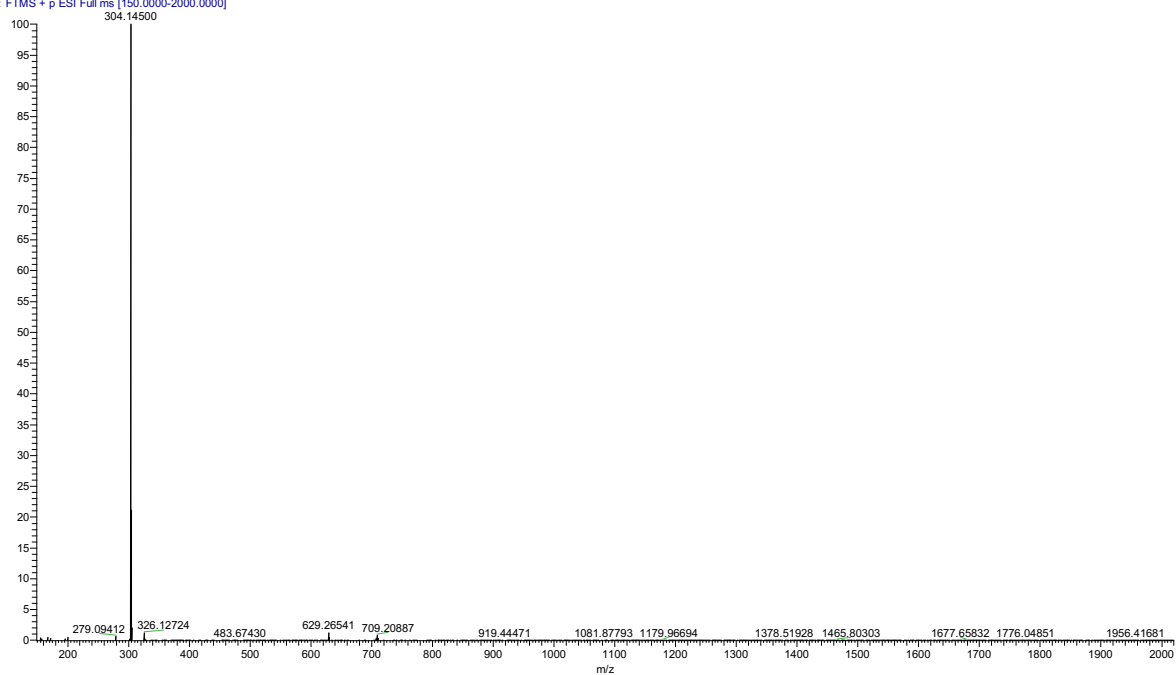


RT: 0.0000 - 11.2584

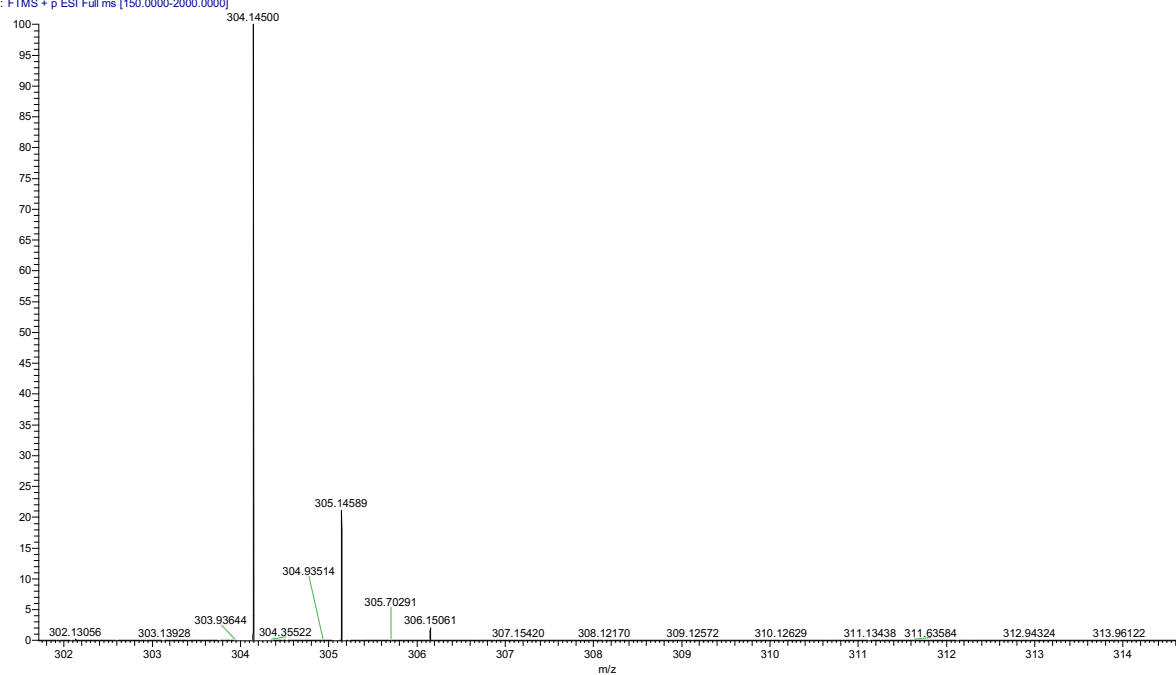


NL:
1.26E10
Base Peak
MS
ALM_DAI_3
3

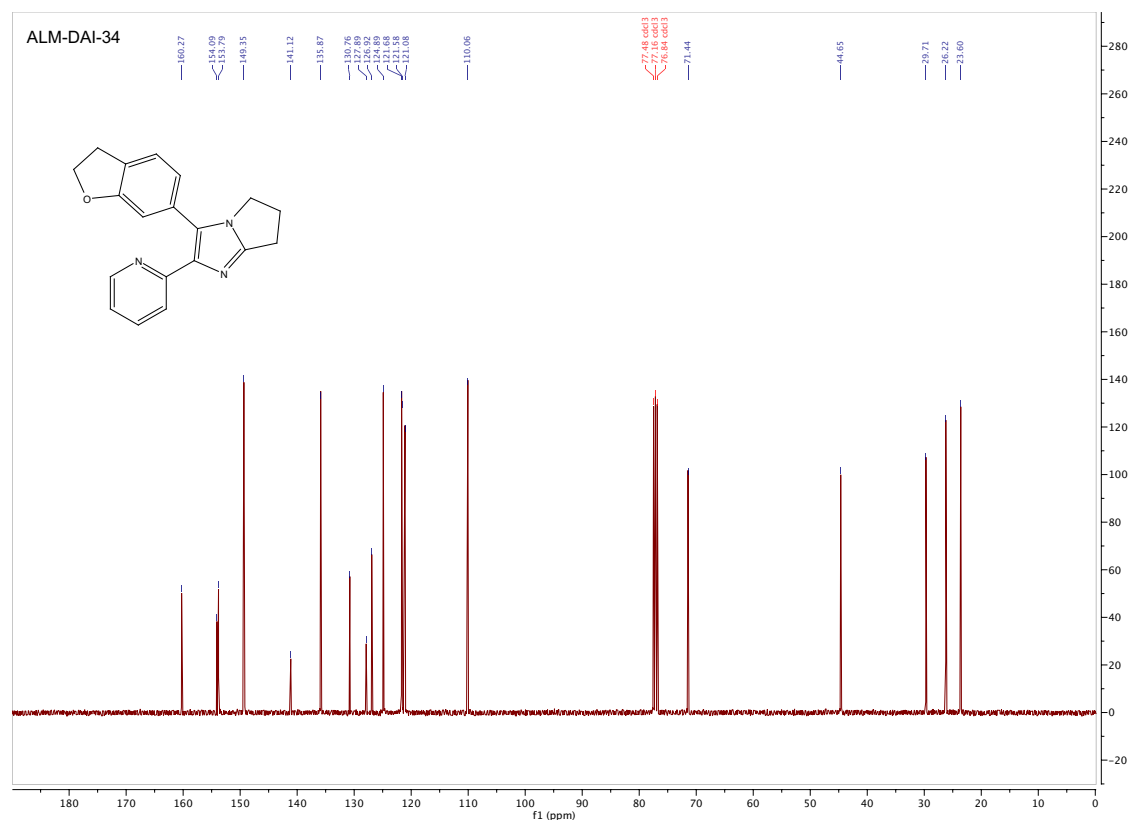
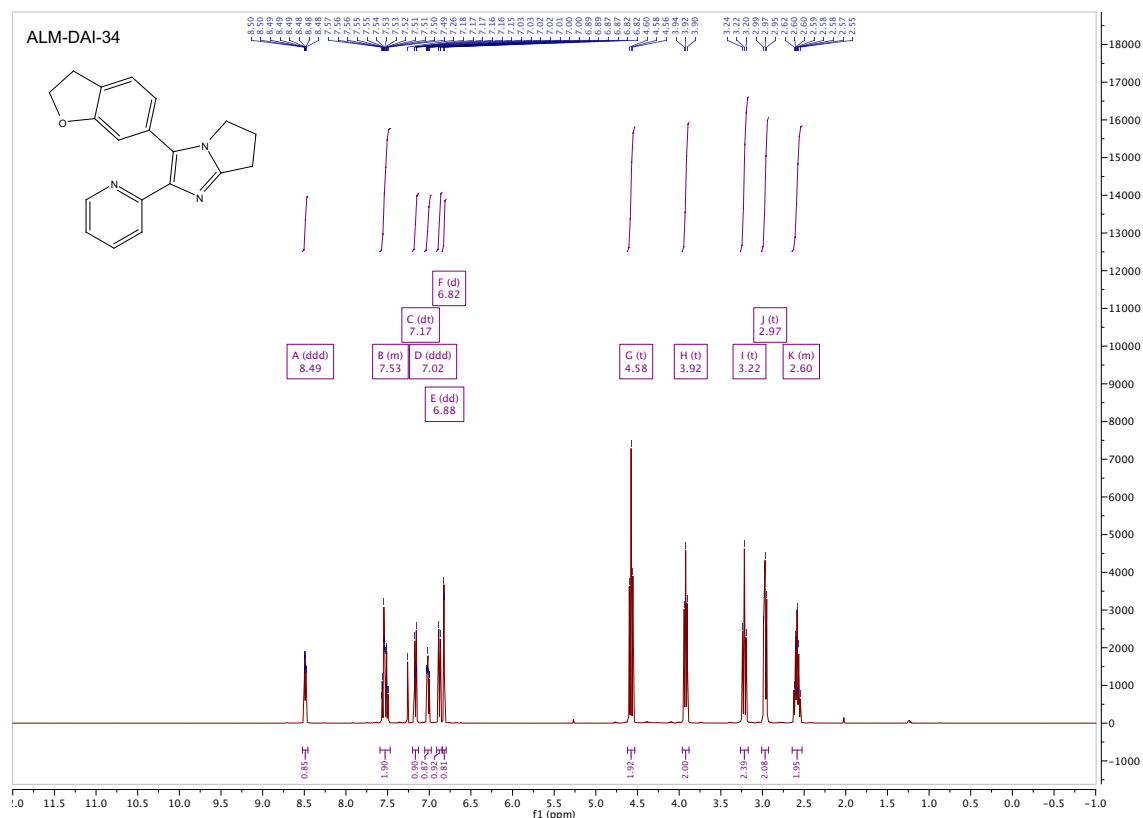
ALM_DAI_33 #791-854 RT: 4.06-4.37 AV: 64 NL: 5.75E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



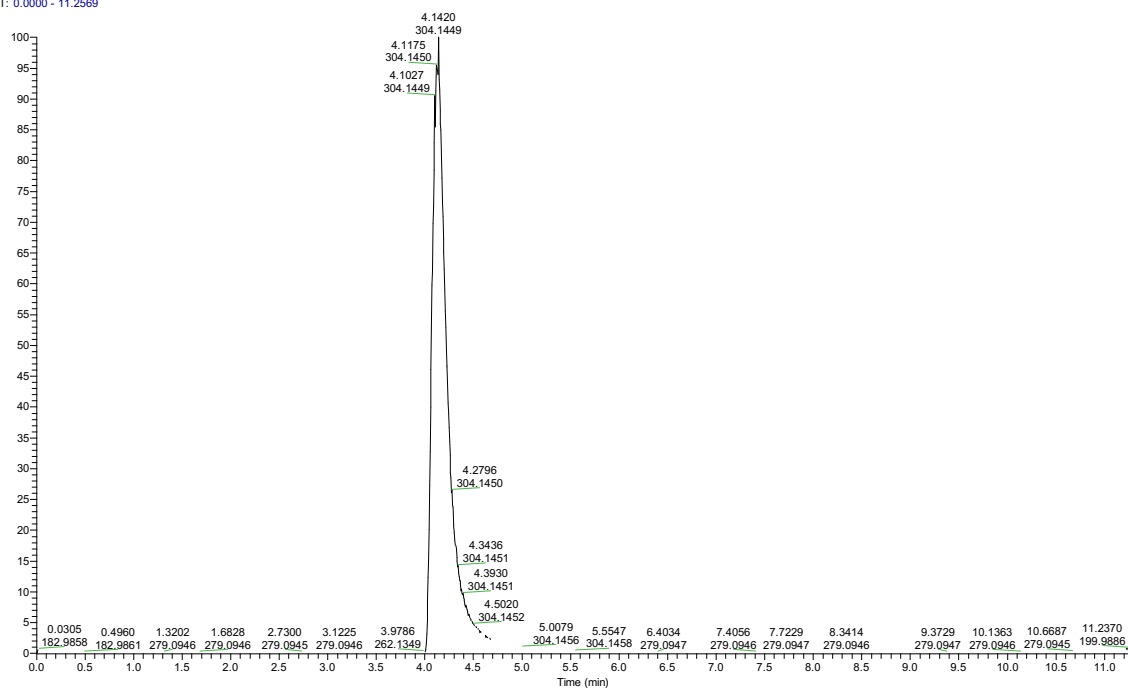
ALM_DAI_33 #791-854 RT: 4.06-4.37 AV: 64 NL: 5.75E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



3-(2,3-dihydrobenzofuran-6-yl)-2-(pyridin-2-yl)-6,7-dihydro-5H-pyrrolo[1,2-*a*]imidazole (ALM-DAI-34).

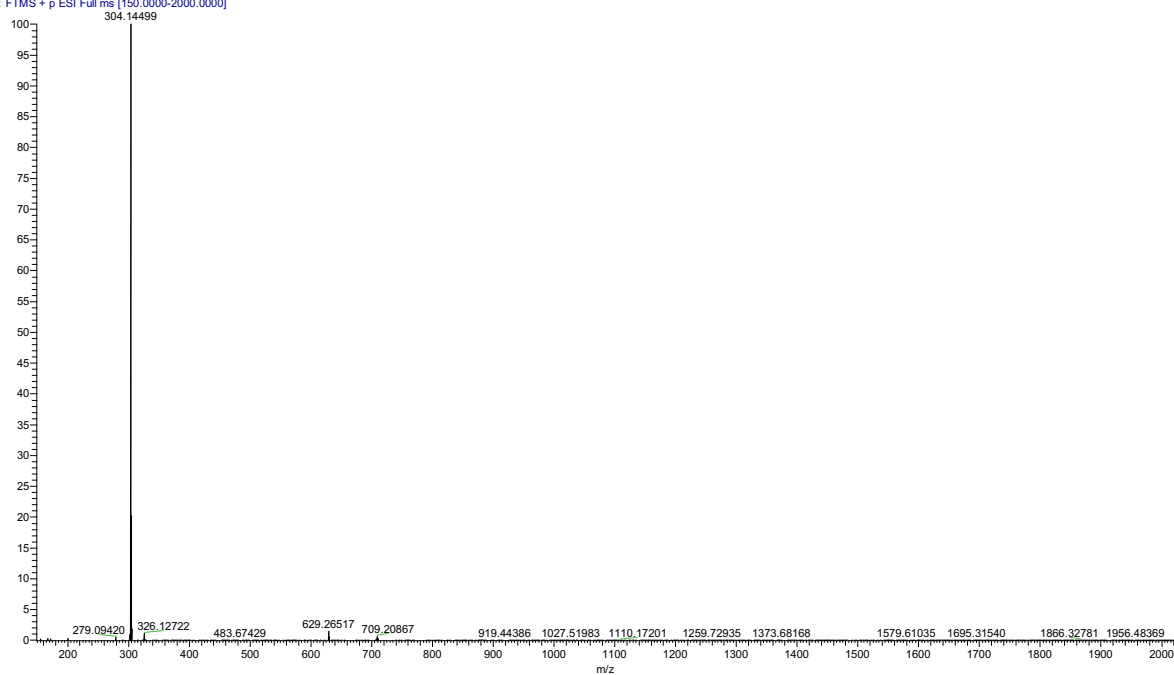


RT: 0.0000 - 11.2569

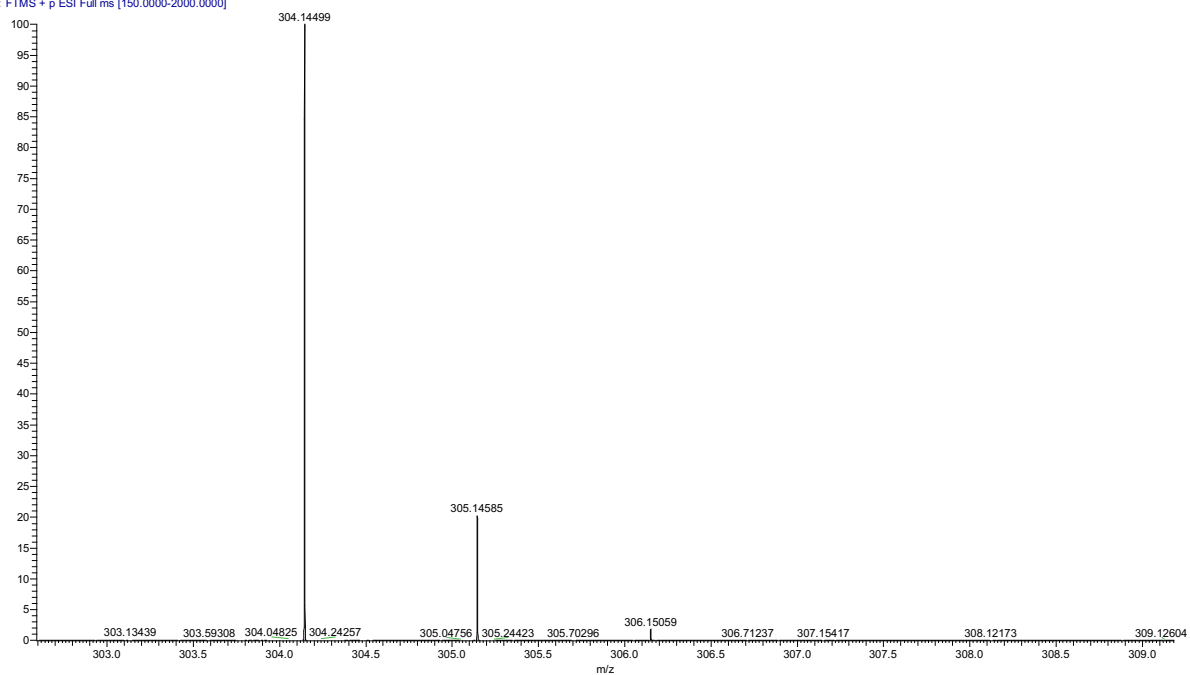


NL:
1.99E10
Base Peak
MS
ALM_DAI_3
4

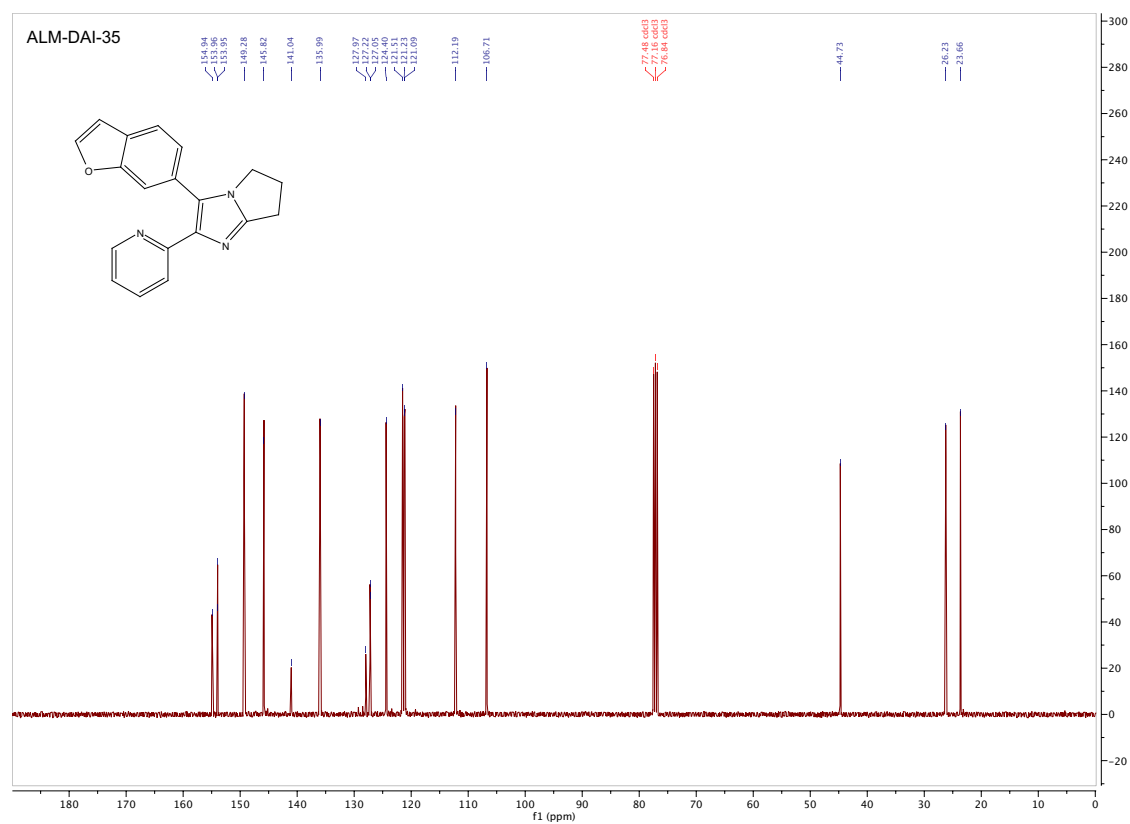
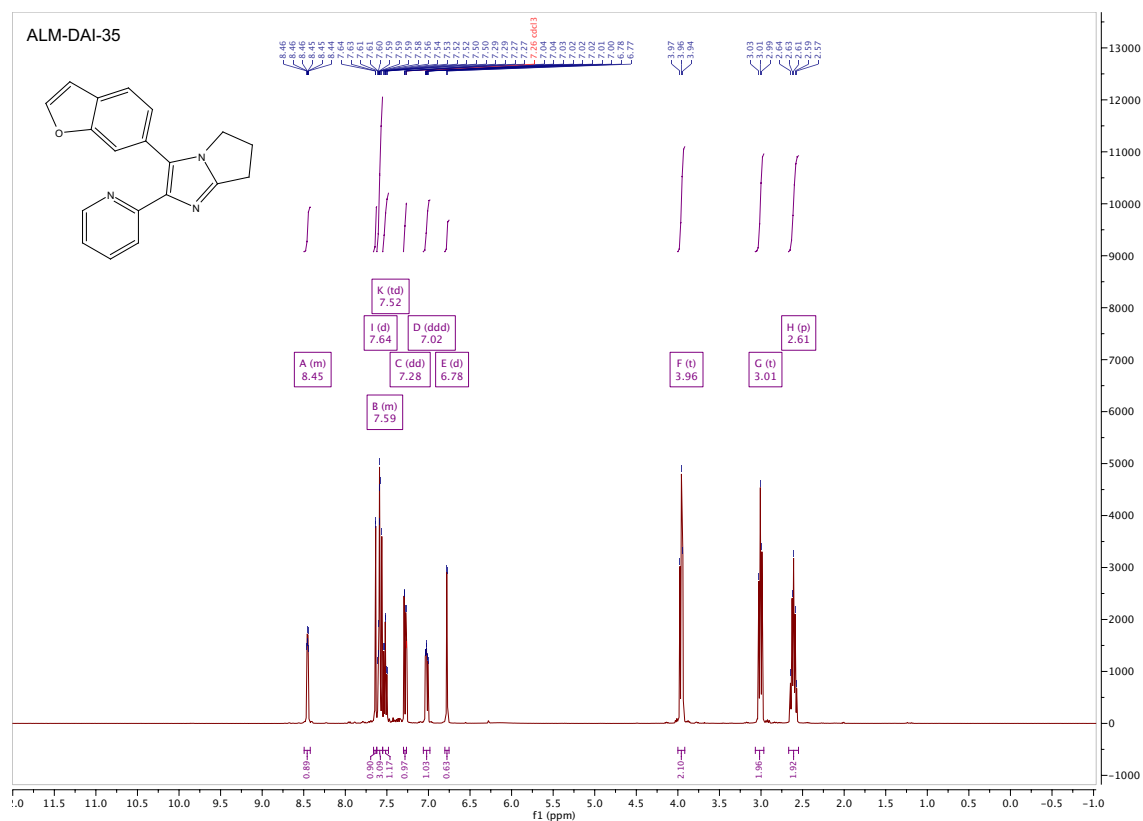
ALM_DAI_34 #785-865 RT: 4.04-4.43 AV: 81 NL: 8.04E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



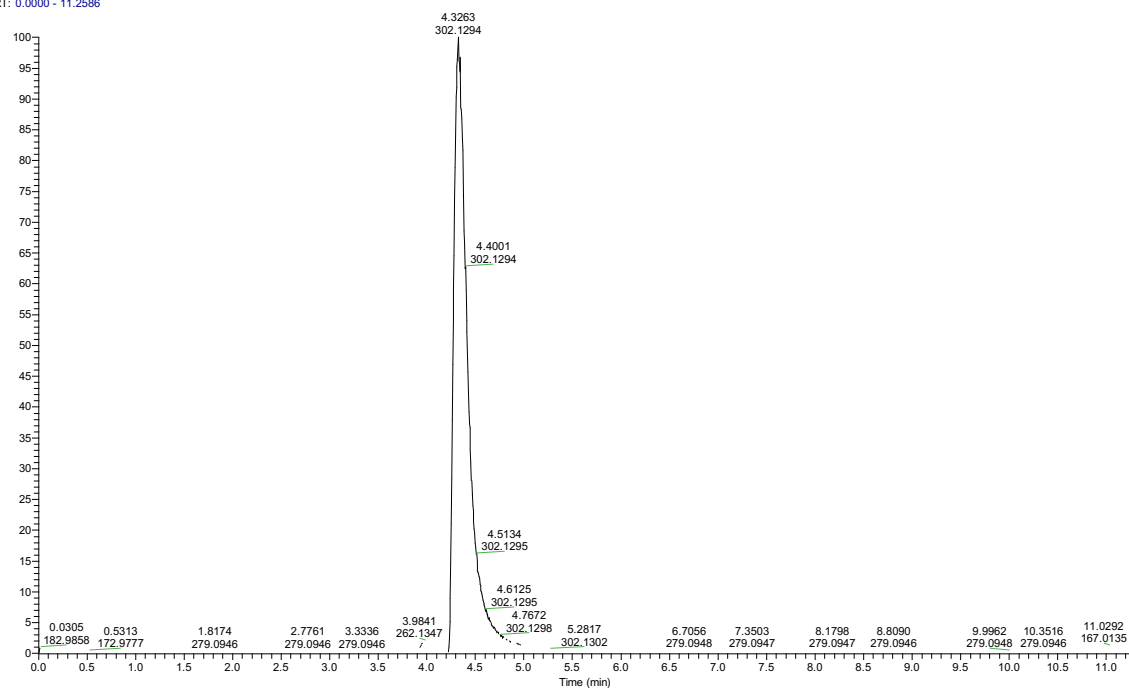
ALM_DAI_34 #785-865 RT: 4.04-4.43 AV: 81 NL: 8.04E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



3-(benzofuran-6-yl)-2-(pyridin-2-yl)-6,7-dihydro-5H-pyrrolo[1,2-a]imidazole (ALM-DAI-35).

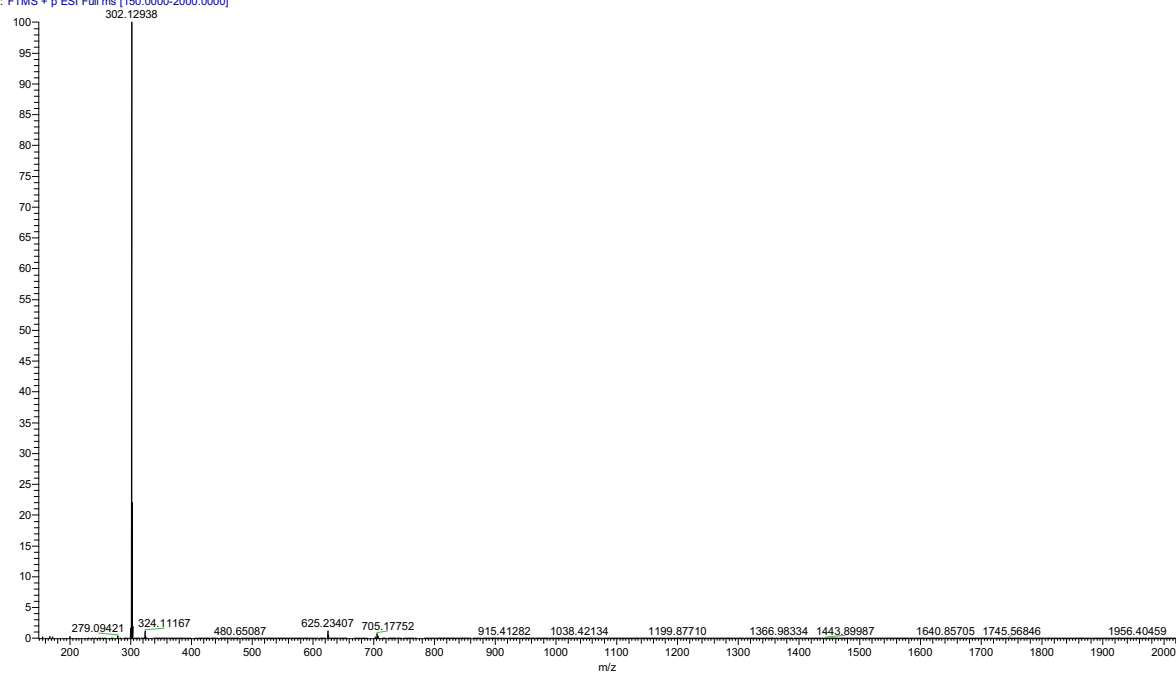


RT: 0.0000 - 11.2586

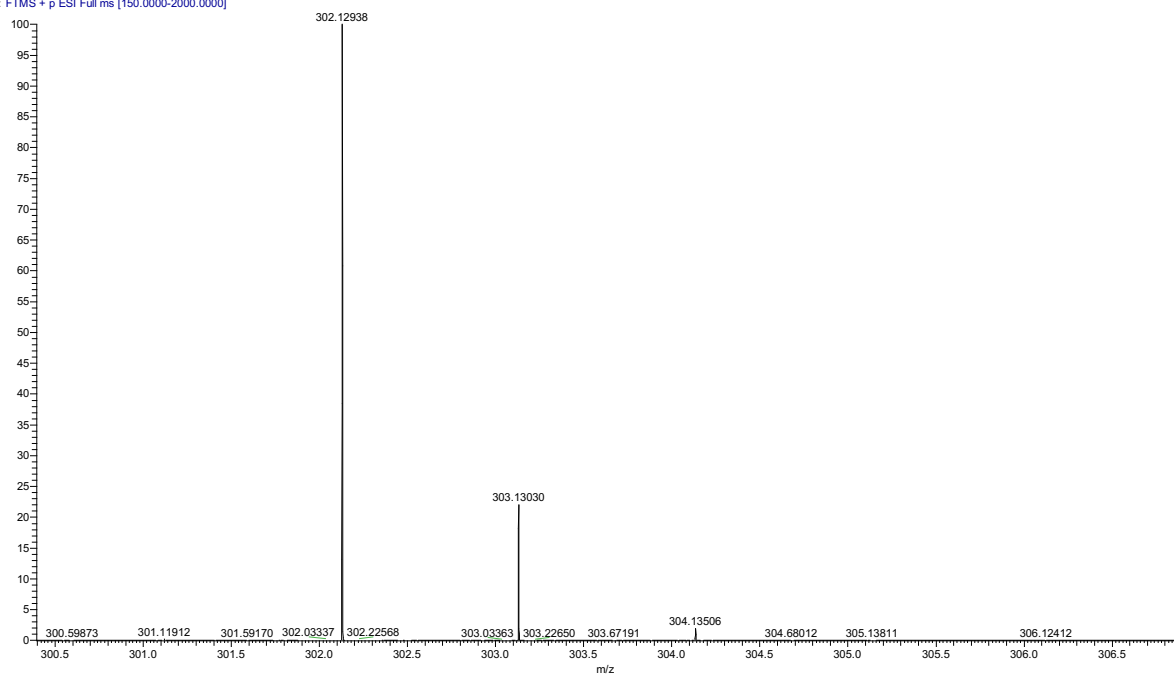


NL:
1.37E10
Base Peak
MS
ALM_DAI_3
5

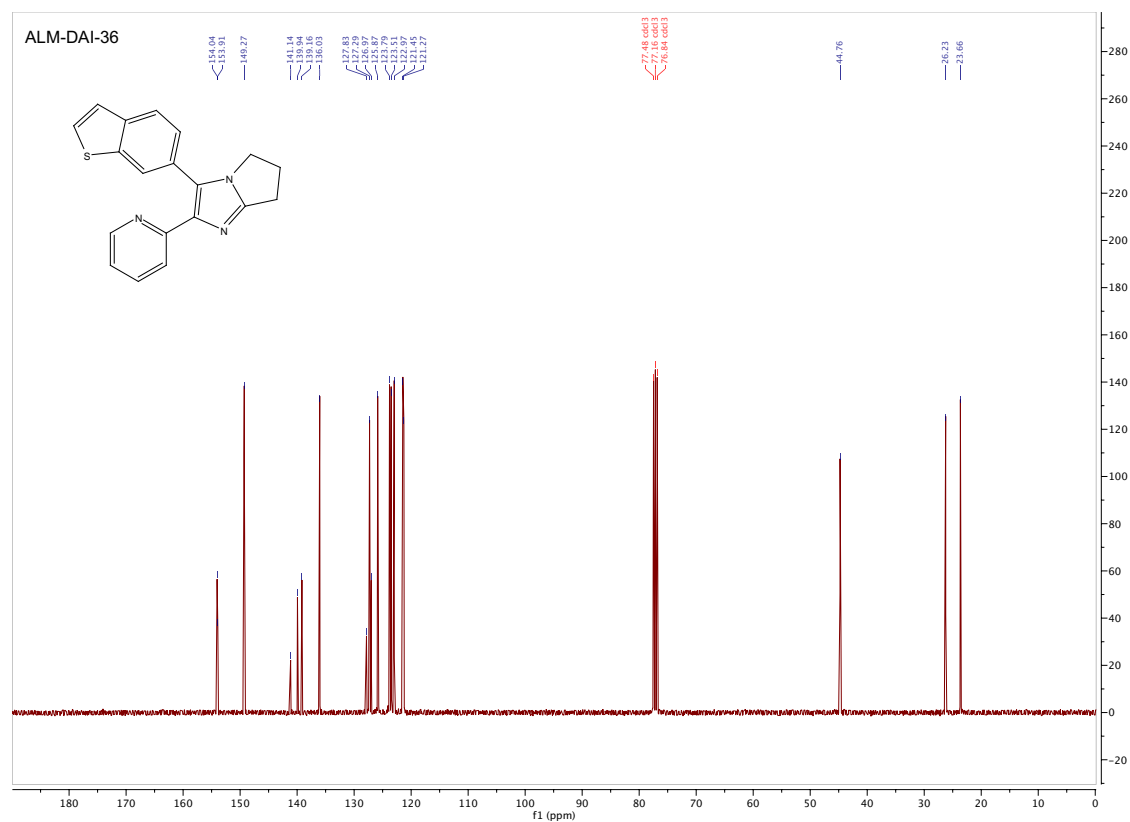
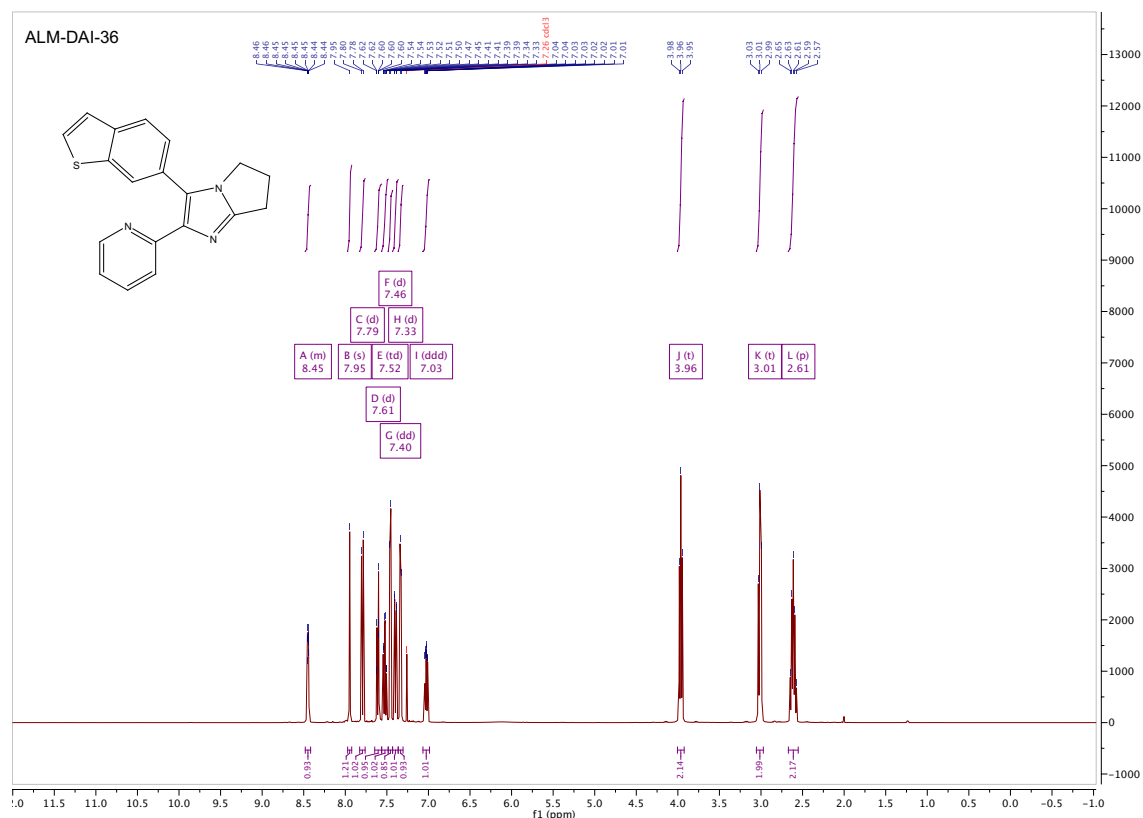
ALM_DAI_35 #928-867 RT: 4.26-4.45 AV: 40 NL: 8.68E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



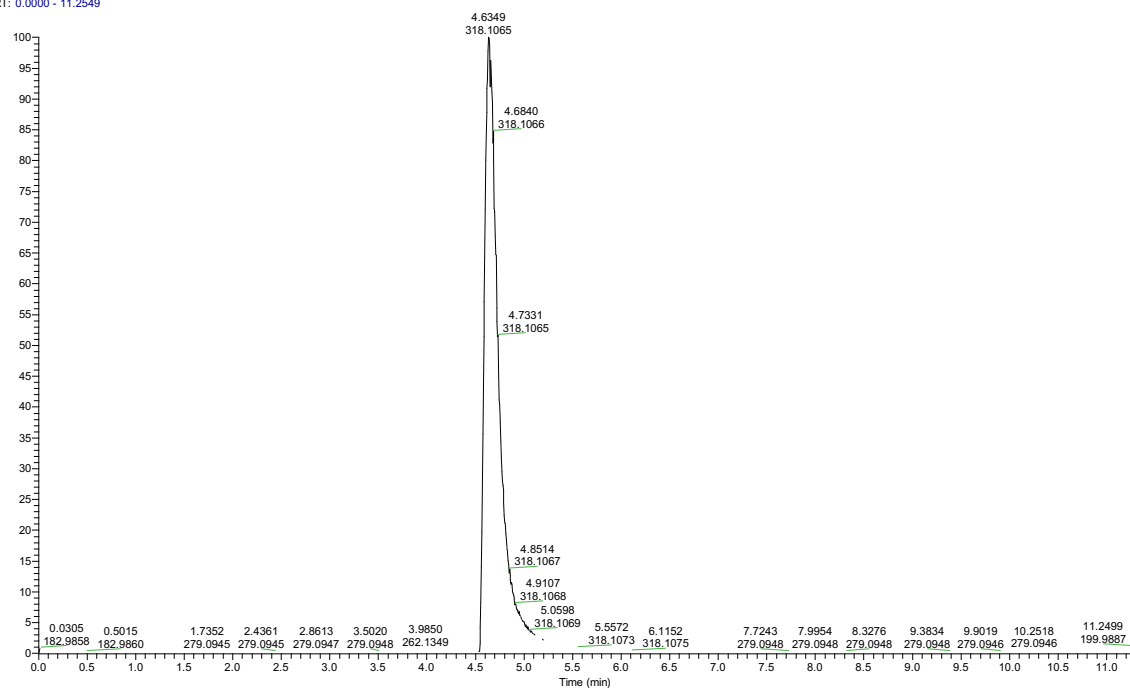
ALM_DAI_35 #828-867 RT: 4.26-4.45 AV: 40 NL: 8.68E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



3-(benzo[*b*]thiophen-6-yl)-2-(pyridin-2-yl)-6,7-dihydro-5*H*-pyrrolo[1,2-*a*]imidazole (ALM-DAI-36).

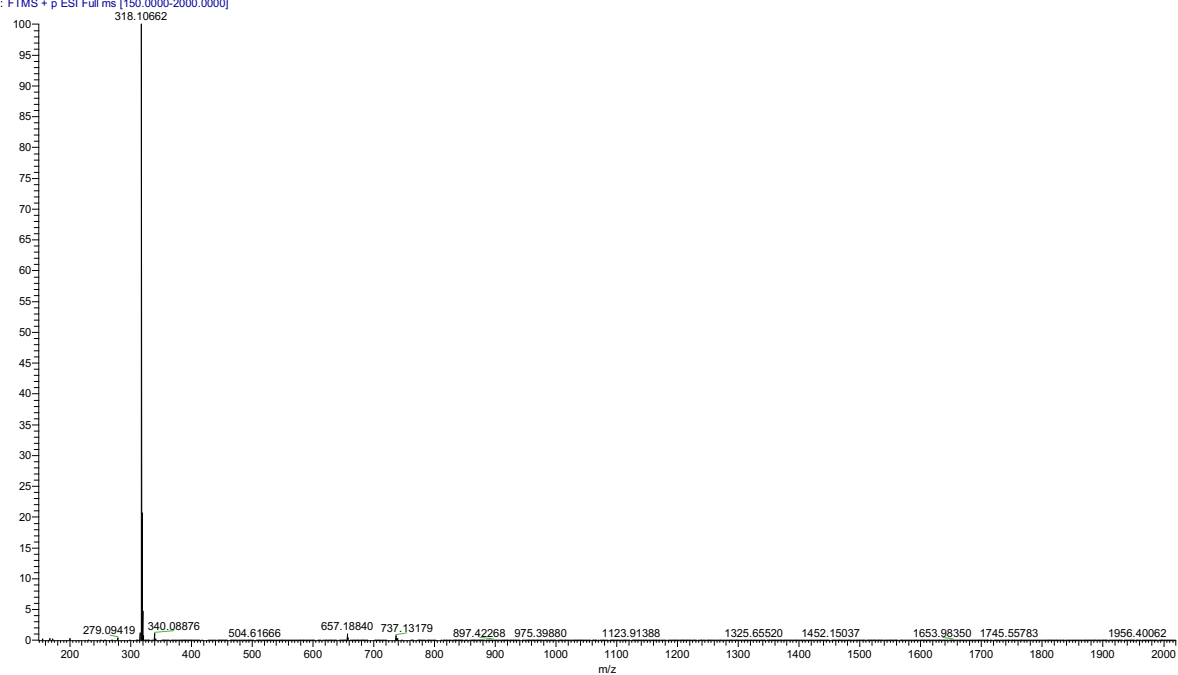


RT: 0.0000 - 11.2549



NL:
1.52E10
Base Peak
MS
ALM_DAI_3
6

ALM_DAI_36 #989-938 RT: 4.58-4.82 AV: 50 NL: 8.92E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]



ALM_DAI_36 #889-938 RT: 4.58-4.82 AV: 50 NL: 8.92E9
T: FTMS + p ESI Full ms [150.0000-2000.0000]

