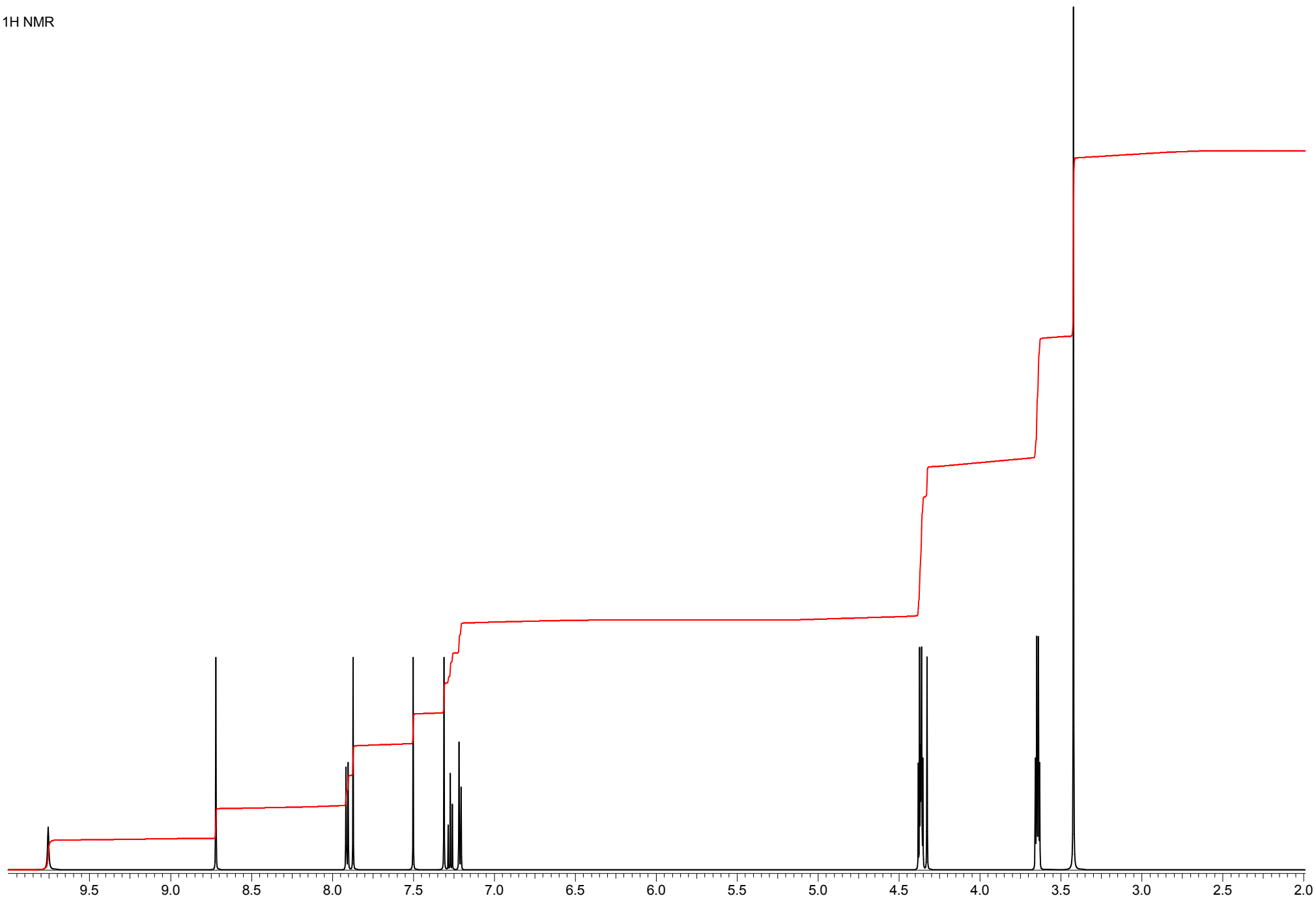
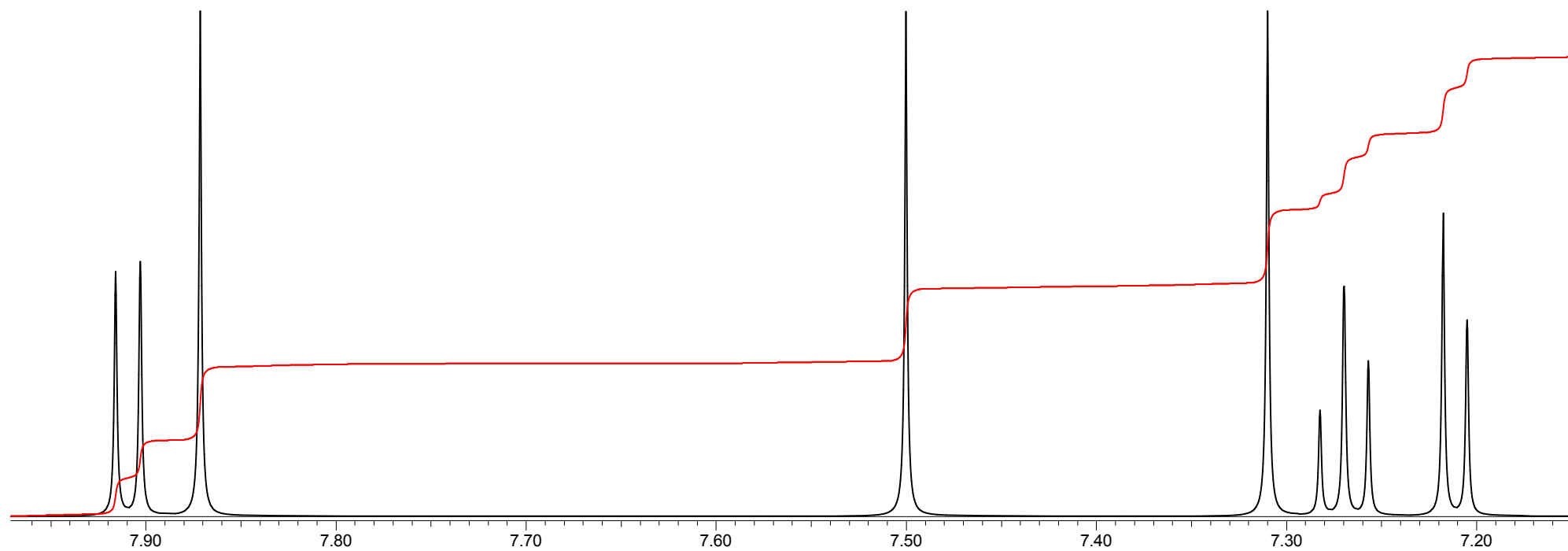
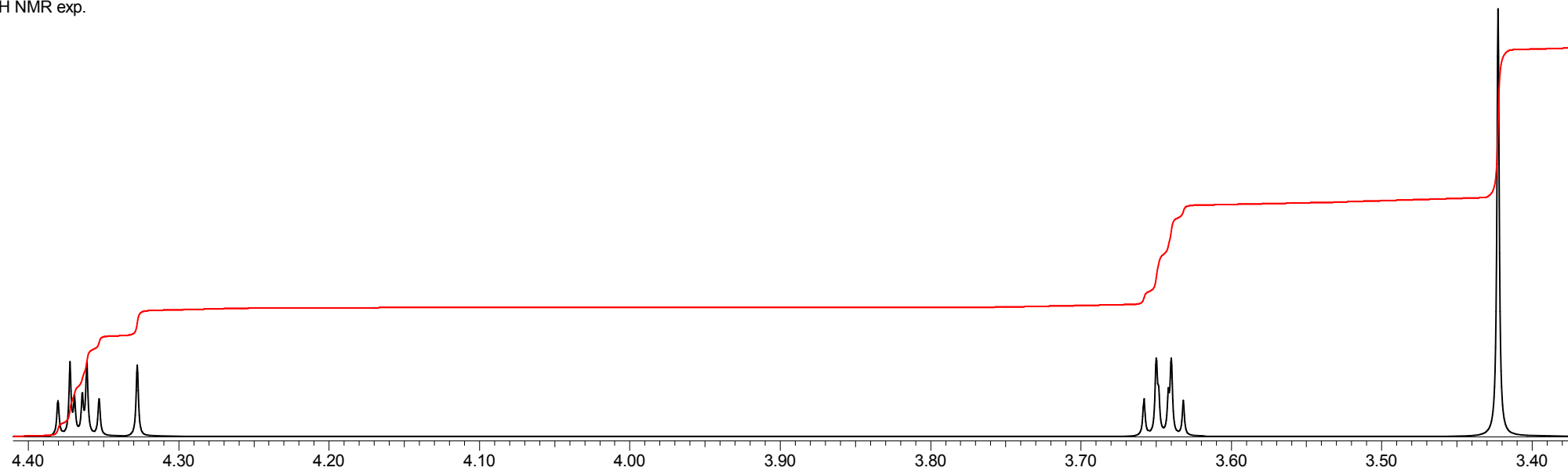


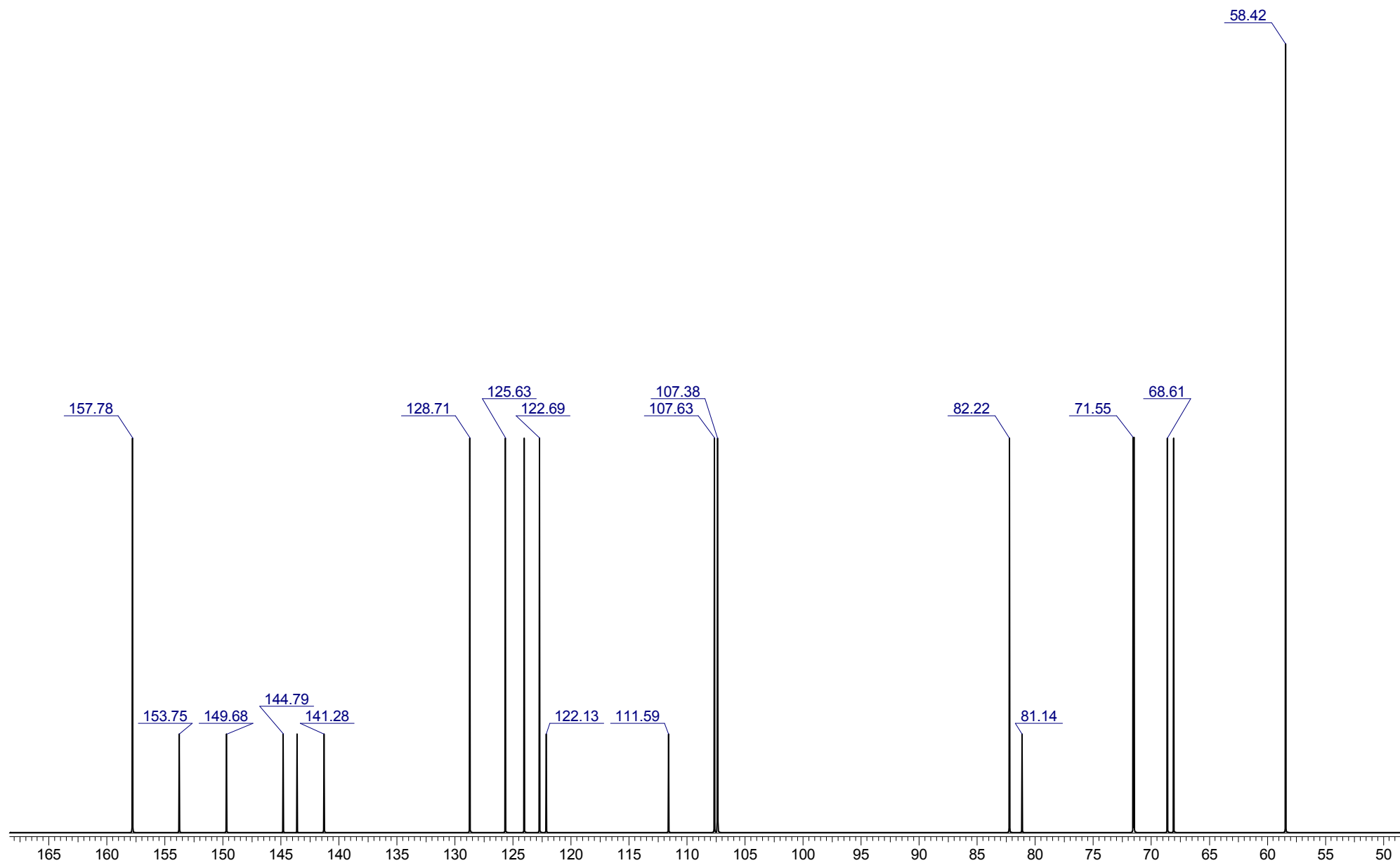
¹H NMR



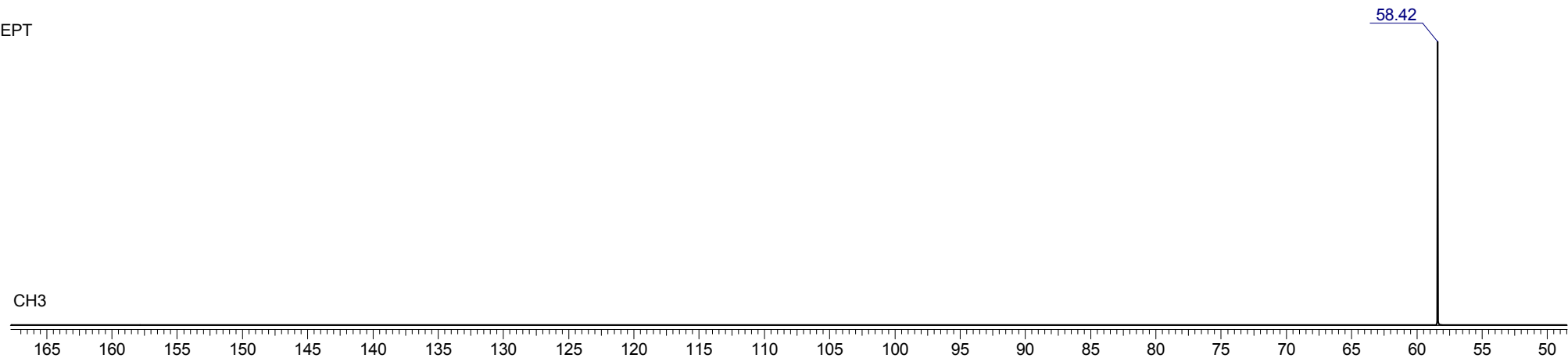
^1H NMR exp.



¹³C



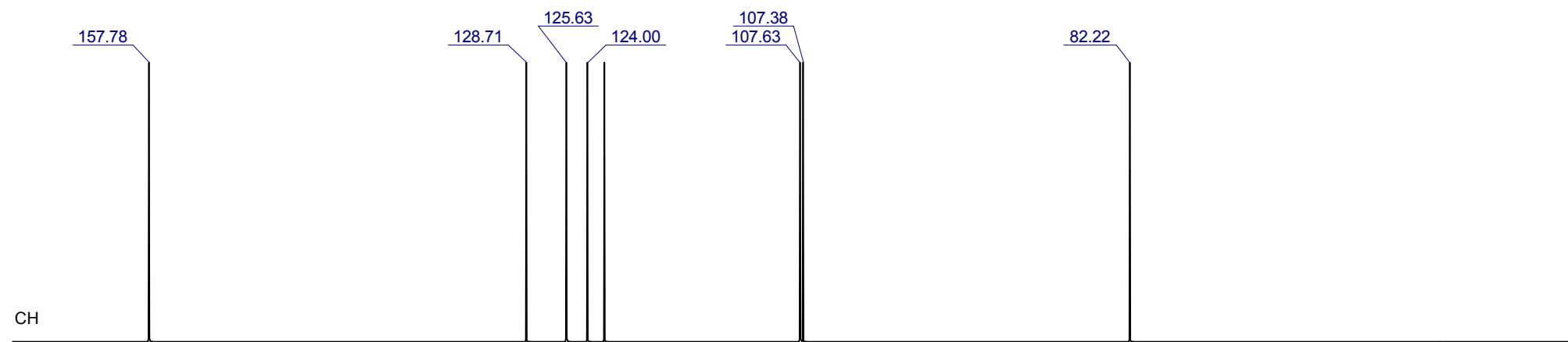
DEPT



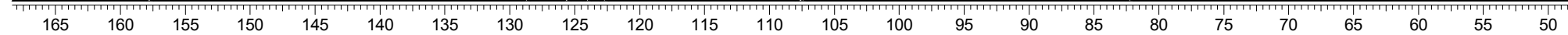
CH3



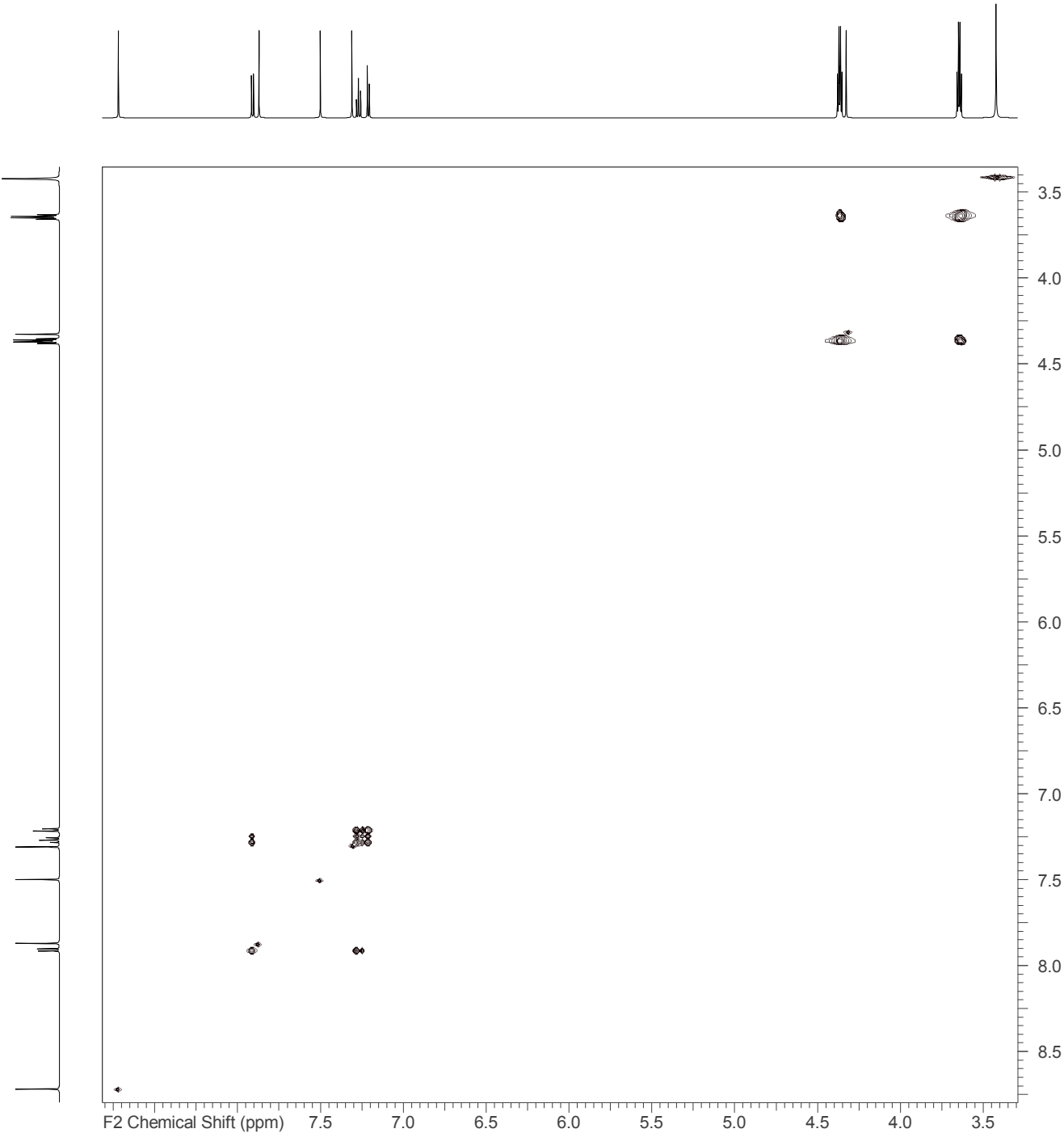
CH2



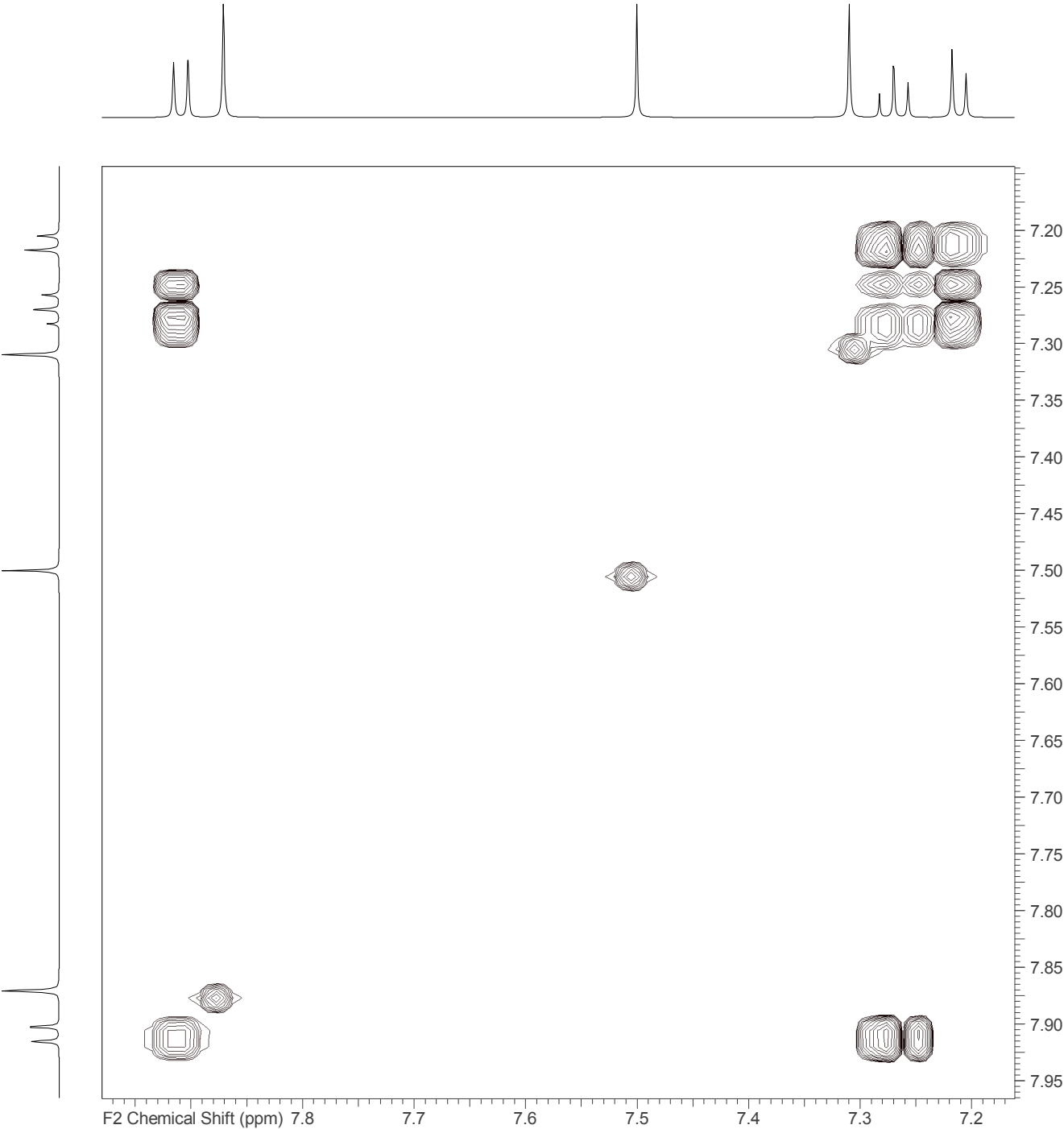
CH



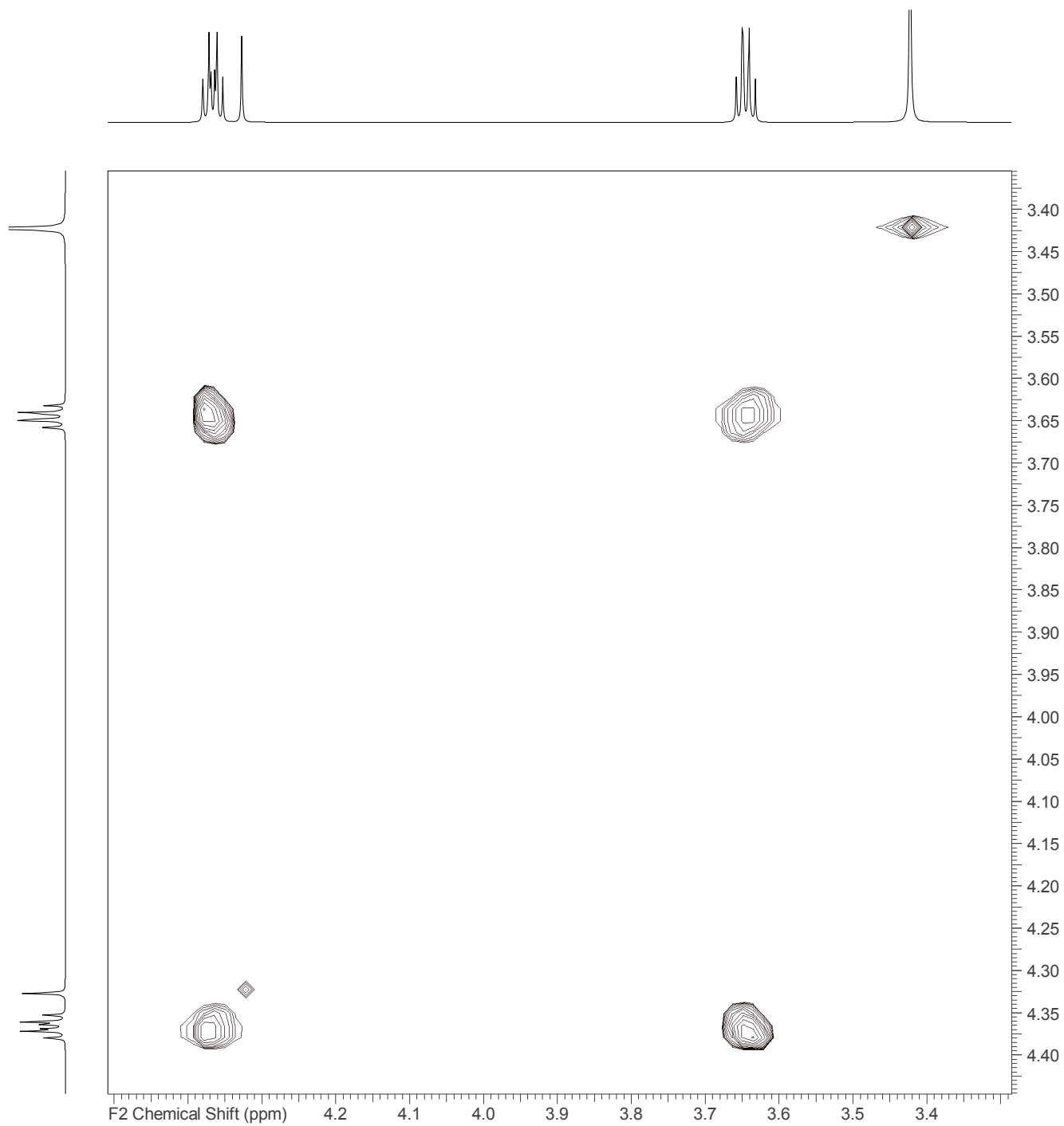
COSY



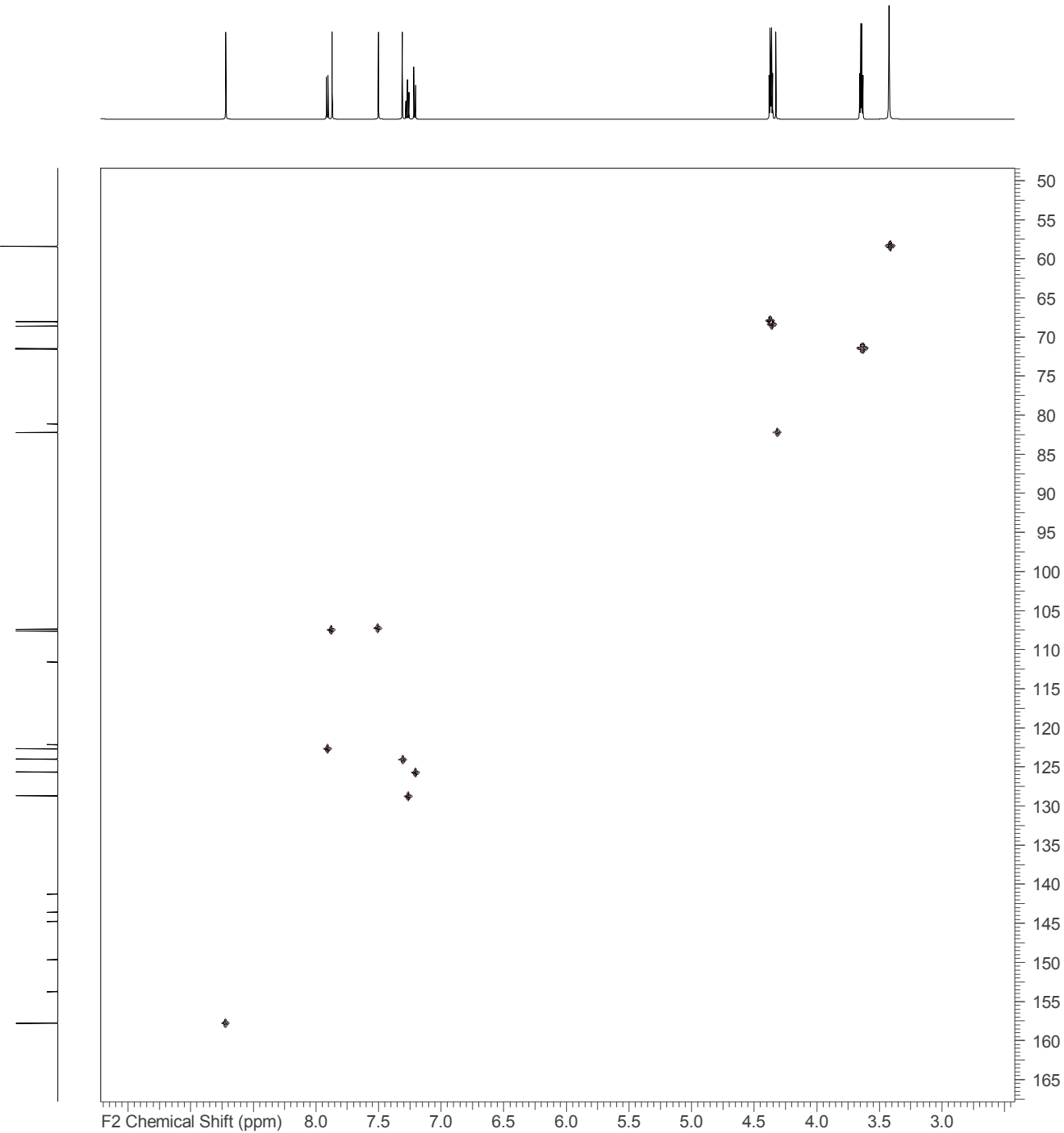
COSY exp.



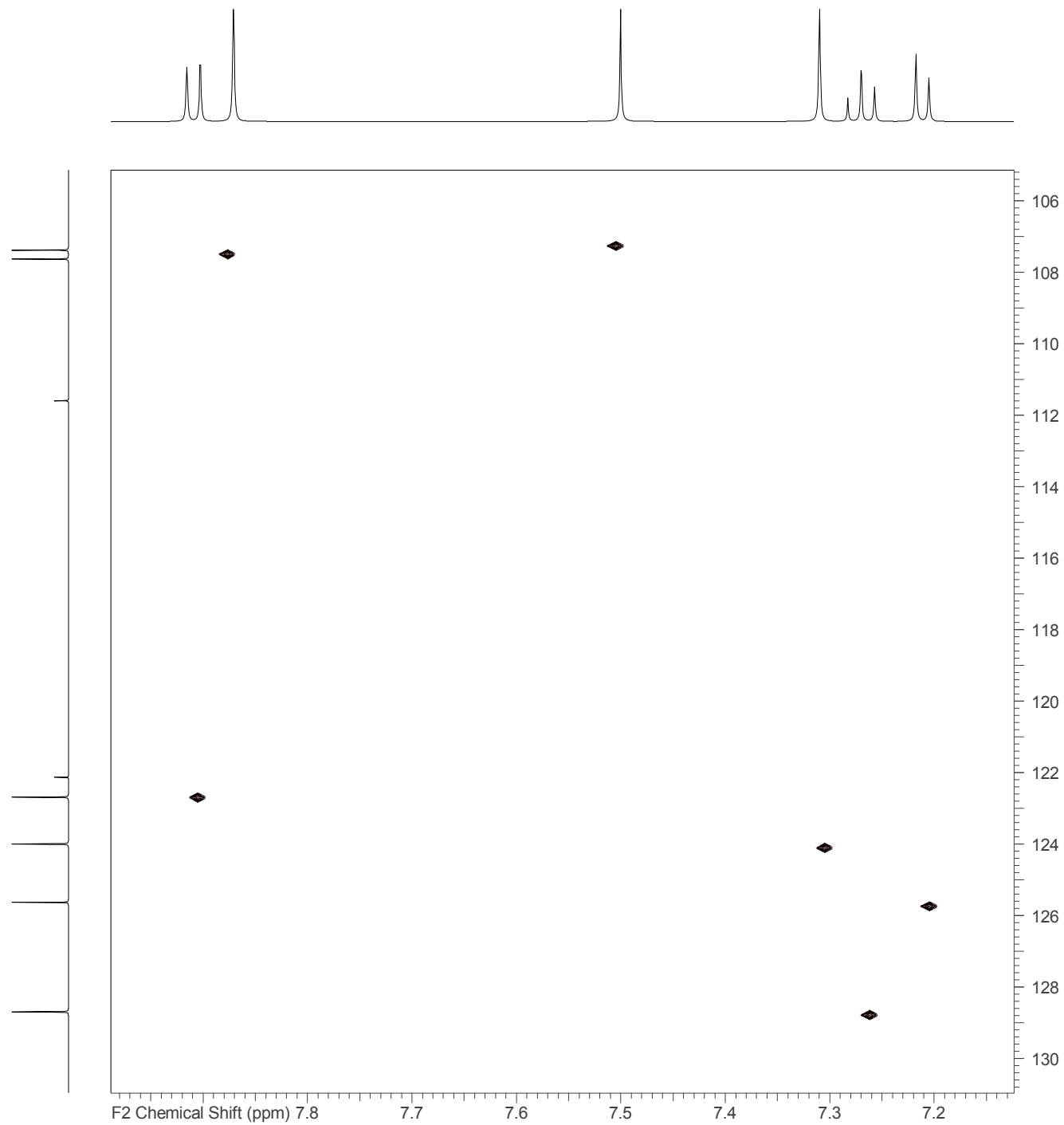
COSY exp.



HSQC

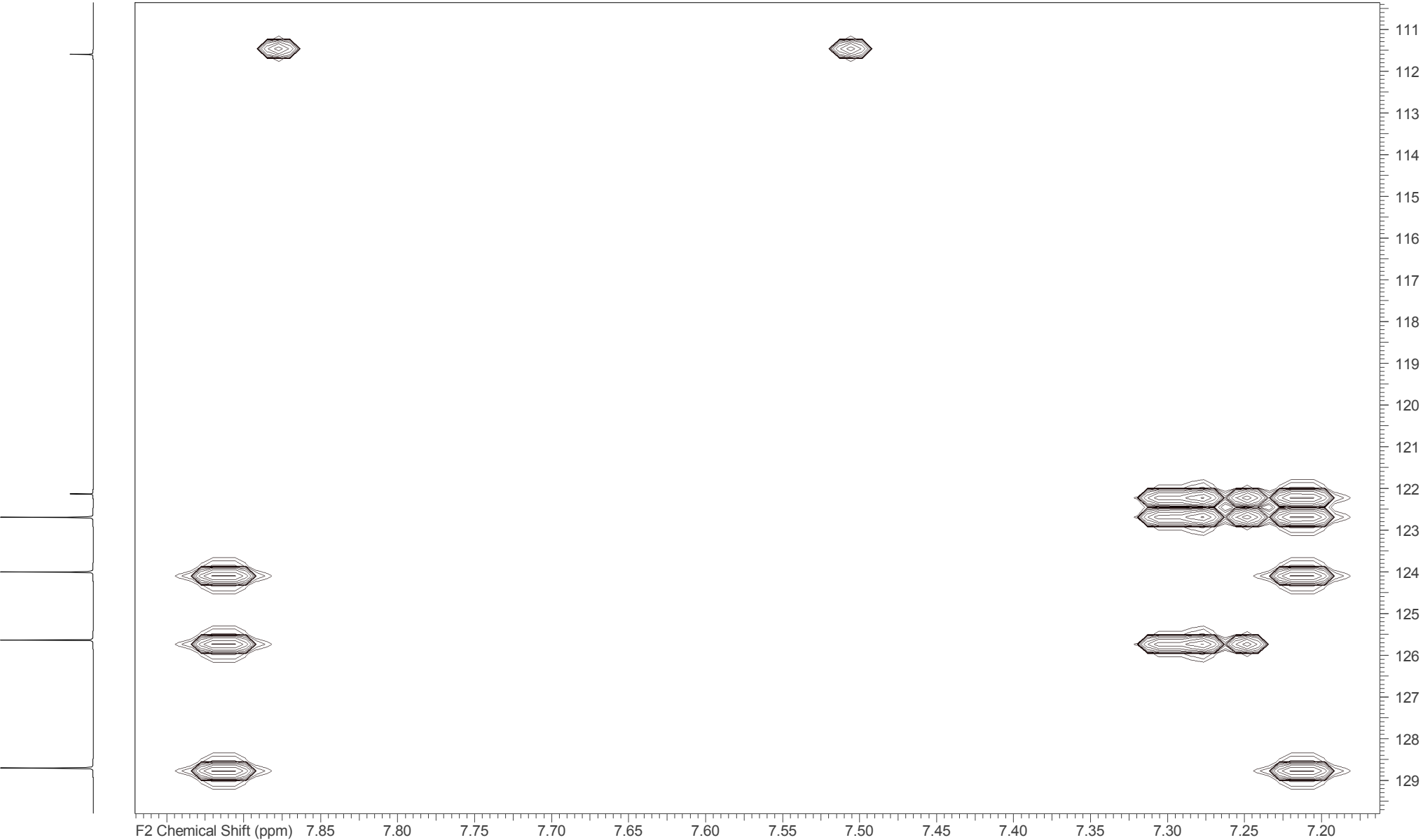
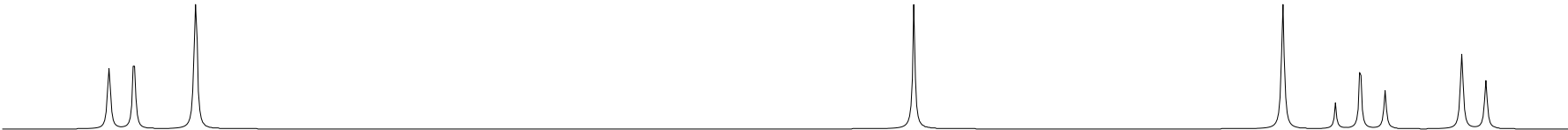


HSQC exp.



2D NMR spectrum (HSQC) of compound 1. The x-axis is F2 Chemical Shift (ppm) from 8.0 to 3.5. The y-axis is F1 Chemical Shift (ppm) from 60 to 155. The spectrum shows correlations between ^1H and ^{13}C chemical shifts. Key features include a cluster of peaks in the aromatic region (7.0-8.0 ppm ^1H , 120-145 ppm ^{13}C), a cluster in the aliphatic region (3.5-4.5 ppm ^1H , 60-75 ppm ^{13}C), and a cluster in the carbonyl region (7.5-8.0 ppm ^1H , 145-155 ppm ^{13}C).

HMBC exp.



HMBC exp.

