

## **Supplementary Material**

### **For**

#### **Novel c-Jun N-terminal Kinase (JNK) Inhibitors with an 11*H*-Indeno[1,2-*b*]quinoxalin-11-one Scaffold**

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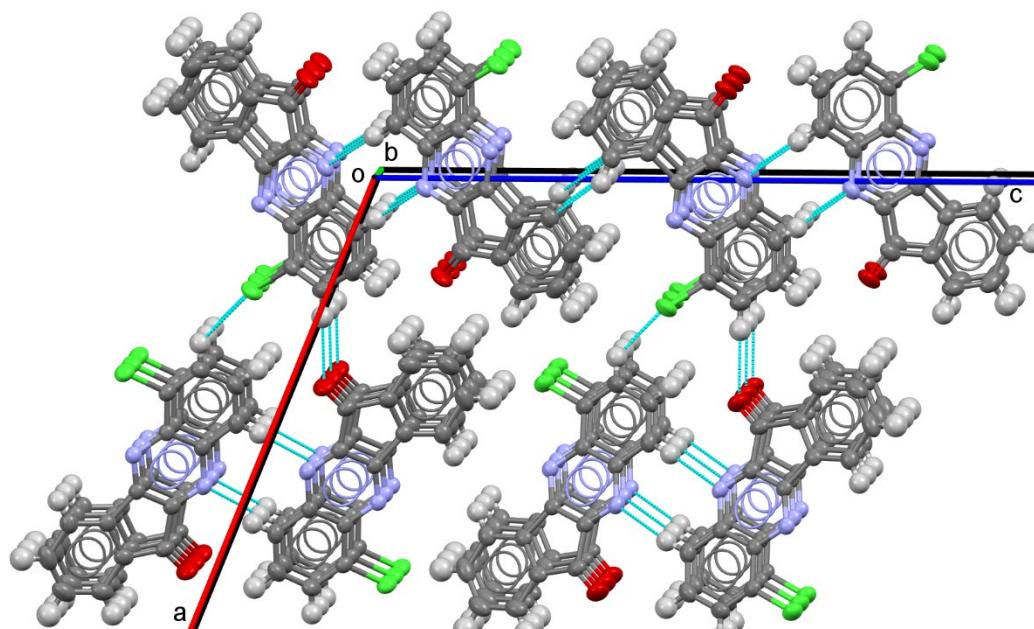
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**Supplementary Table S1.** Bond lengths and angles of H-bonds for compound **3b** in the crystal<sup>a</sup>.

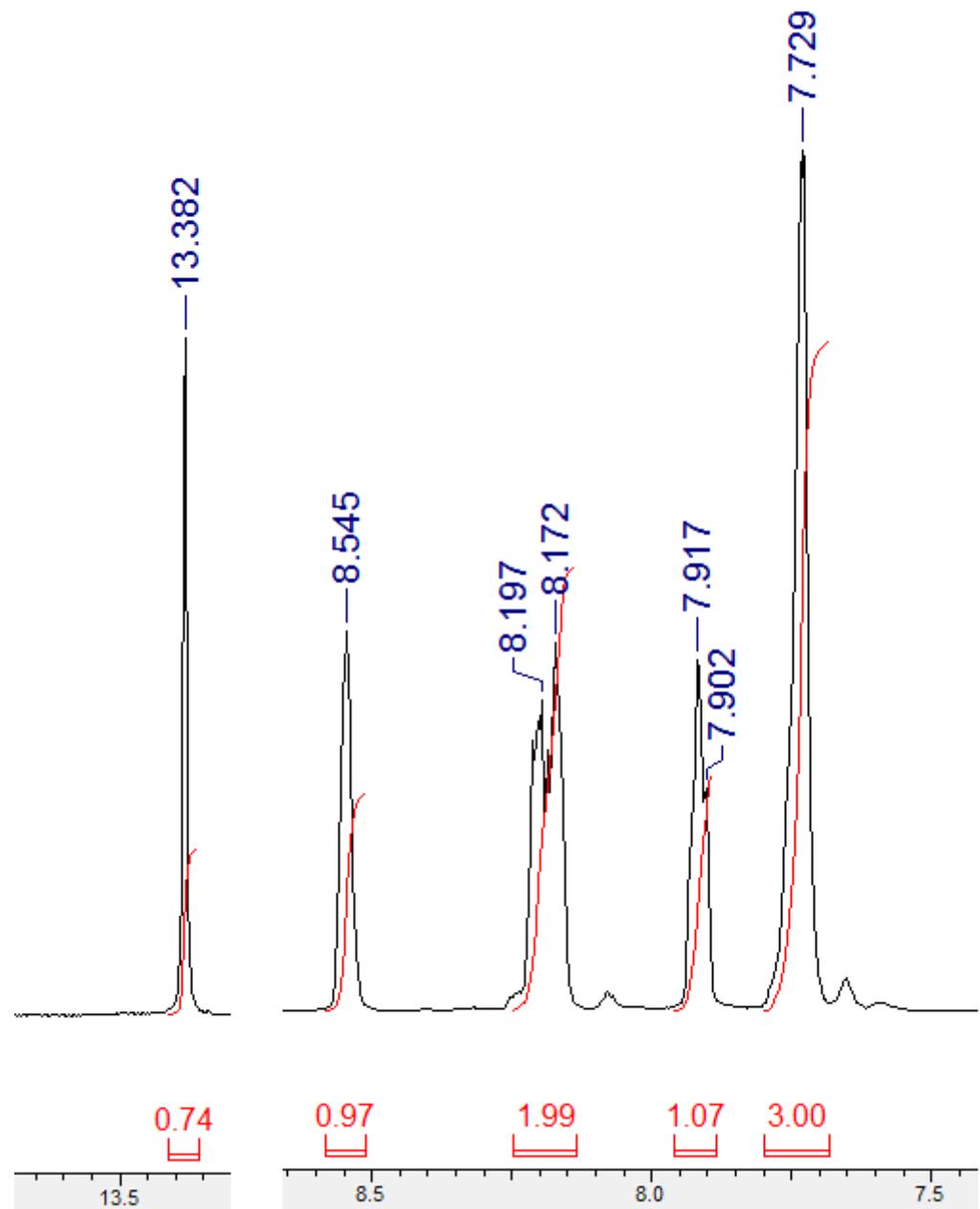
H-bond	D-H	Distance (Å)		Angle <sup>a</sup>
		H...A	D...A	D-H...A
C9-H...N10	0.93	2.58	3.497(3)	168
C7'-H...O13	0.93	2.55	3.451(4)	163
C9'-H...N10'	0.93	2.58	3.493(4)	168

<sup>a</sup>Bond angles in degrees.

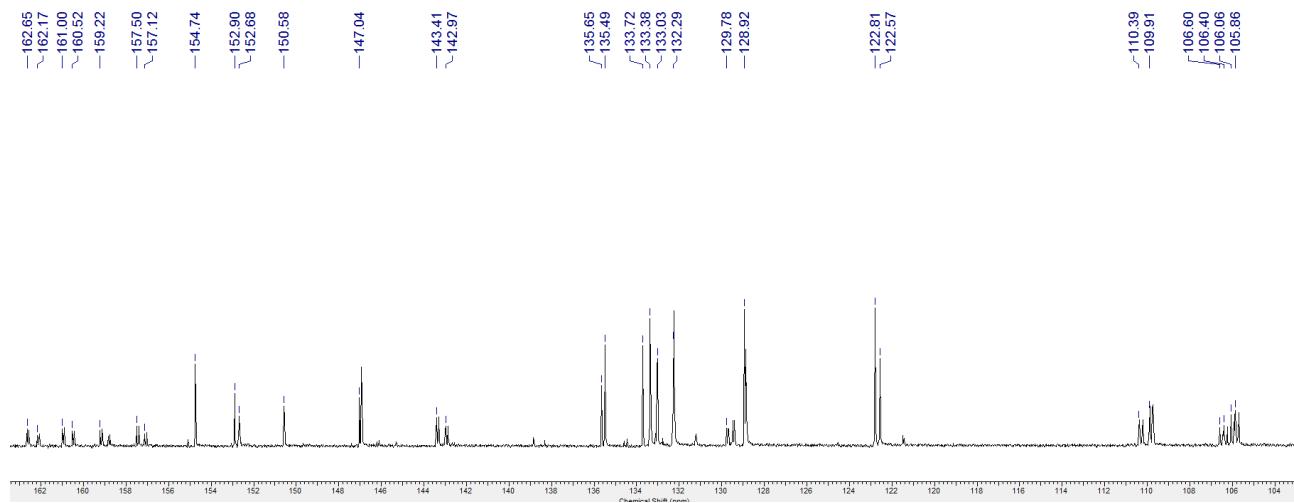


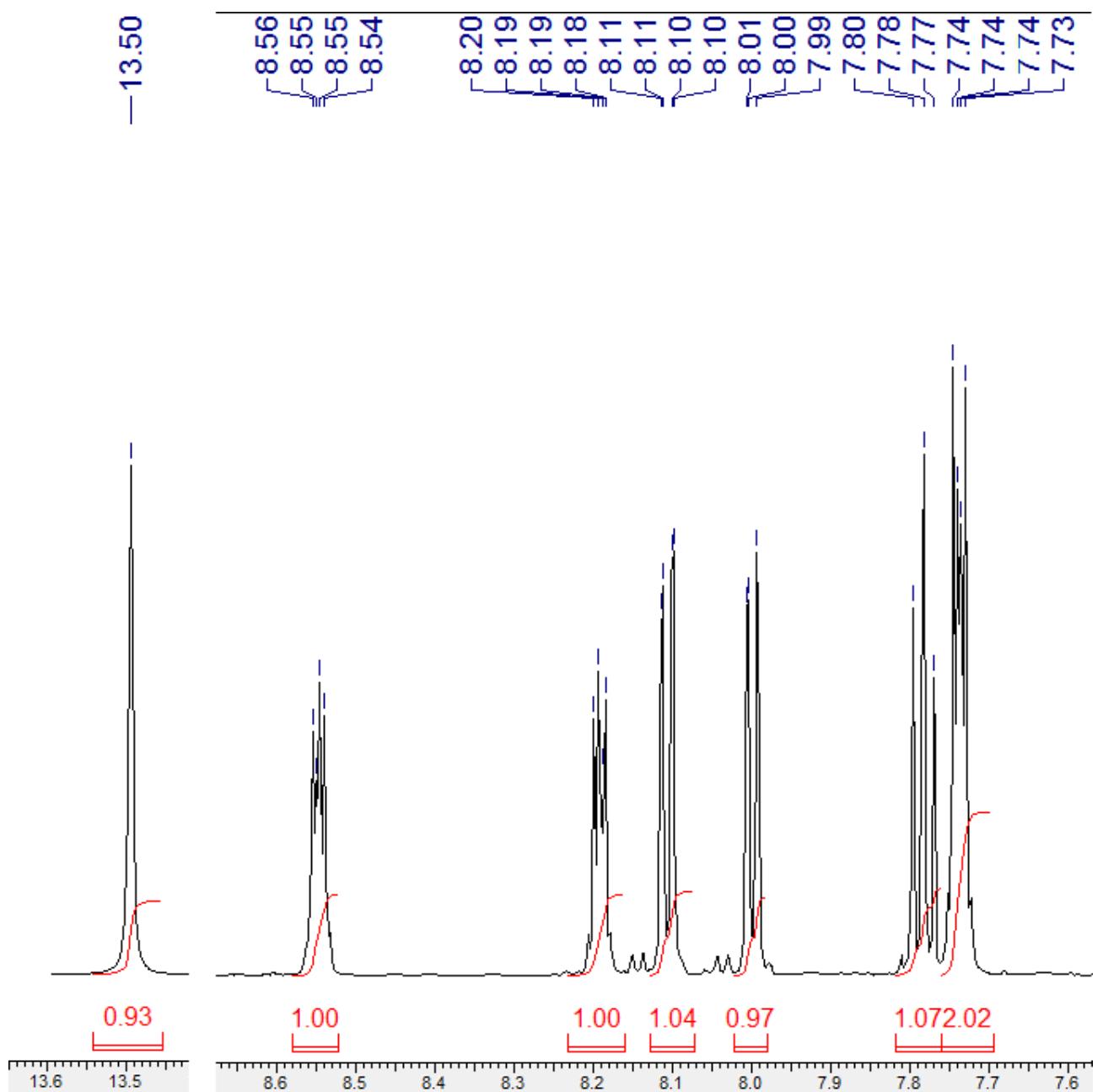
**Supplementary Figure S1.** Crystal packing of compound **3b** along crystallographic axis *b*. Oxygen and chlorine atoms are shown in red and green, respectively. Intermolecular H-bonds are shown as thin blue lines.

### NMR Spectra Data

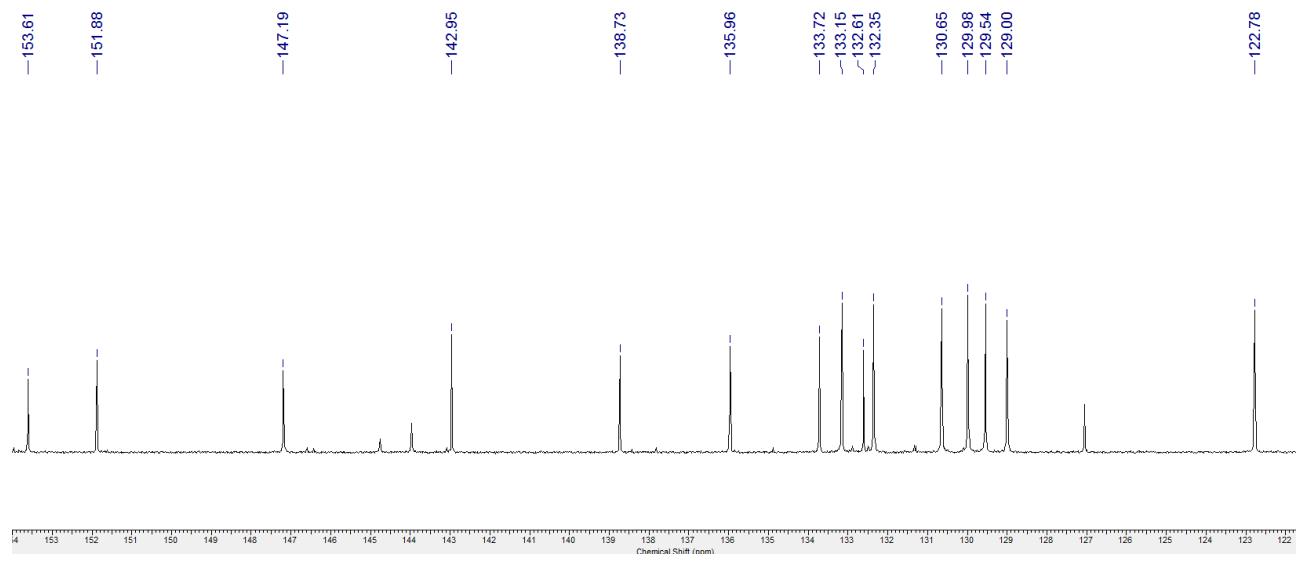


NMR <sup>1</sup>H spectrum (600 MHz, DMSO-D<sub>6</sub>) of compound 4a.

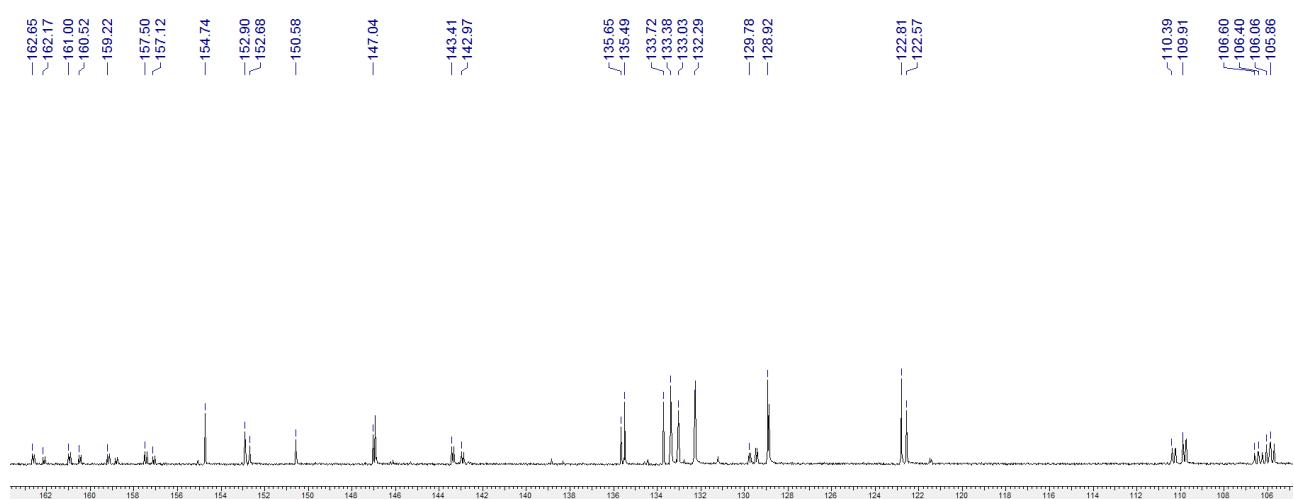
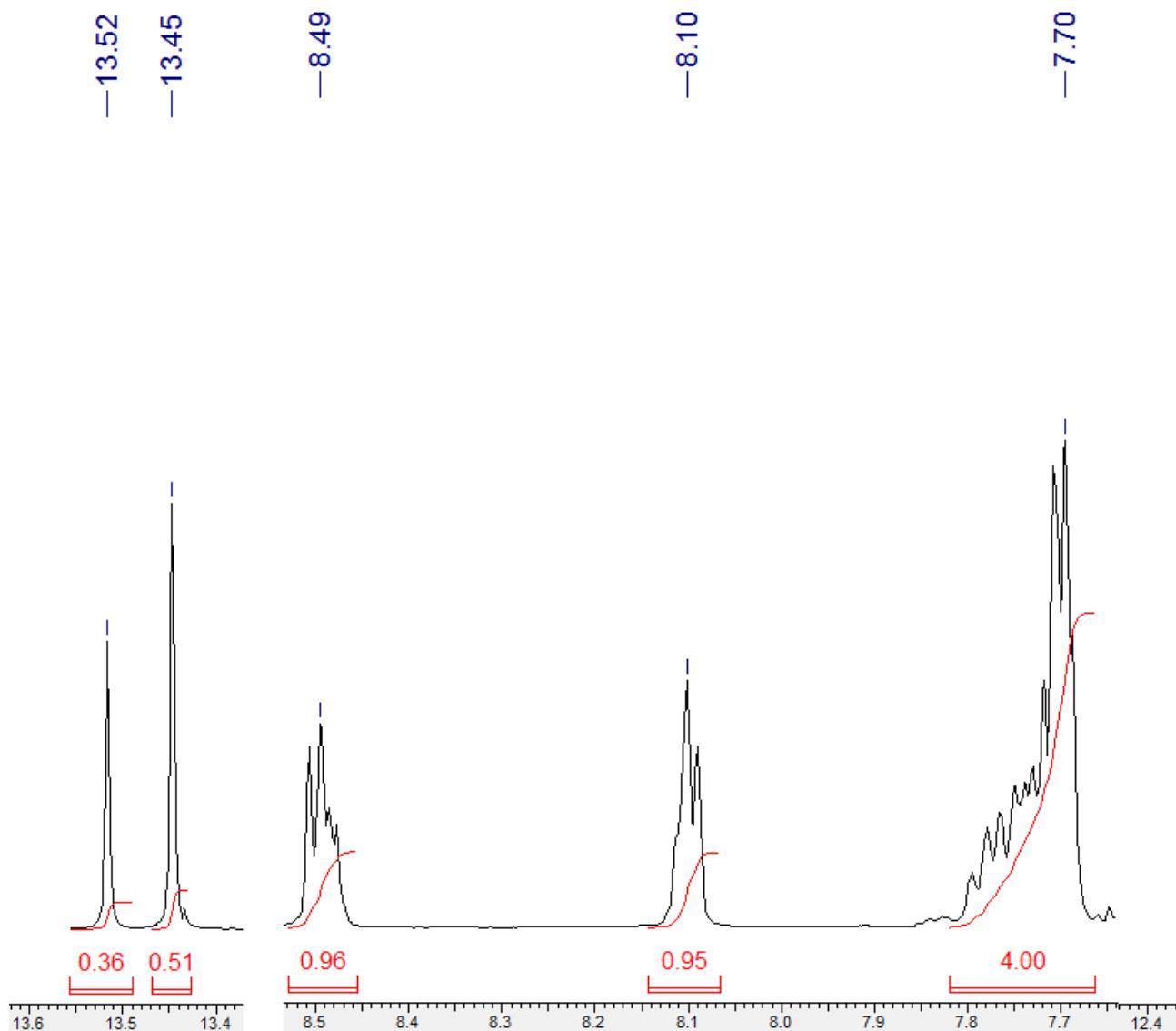


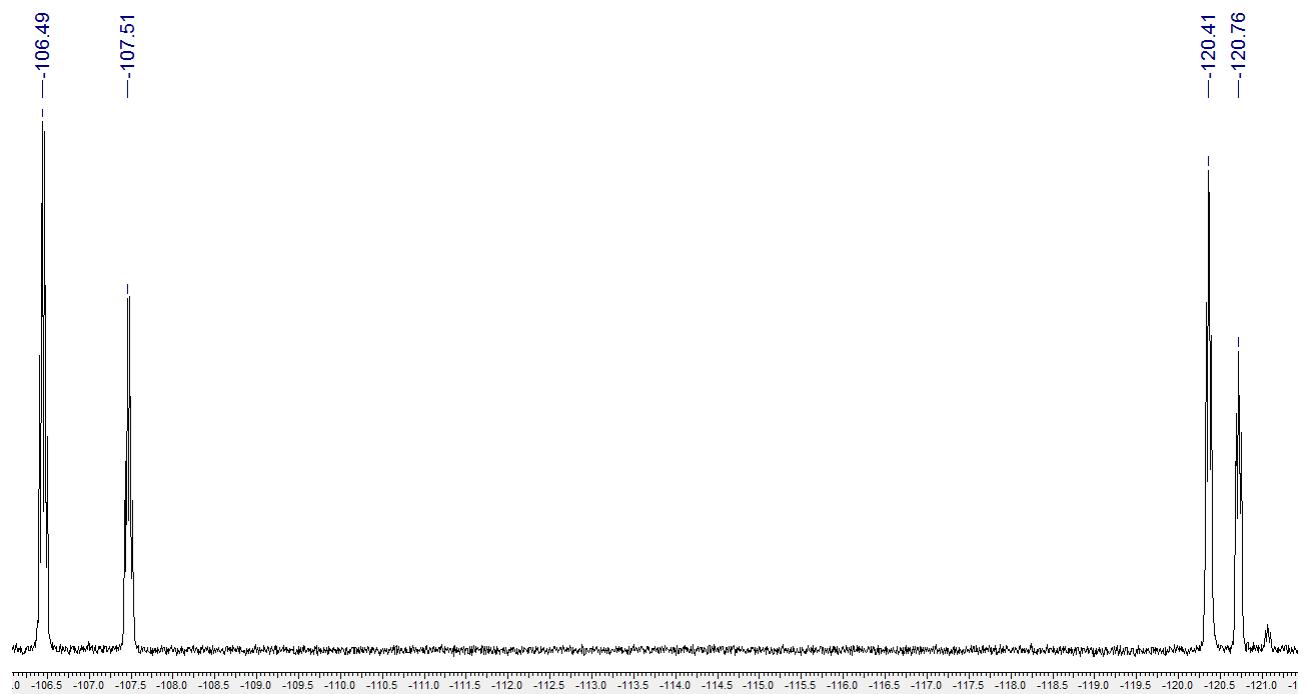


NMR  $^1\text{H}$  spectrum (600 MHz, DMSO-D6) of compound **4b**.

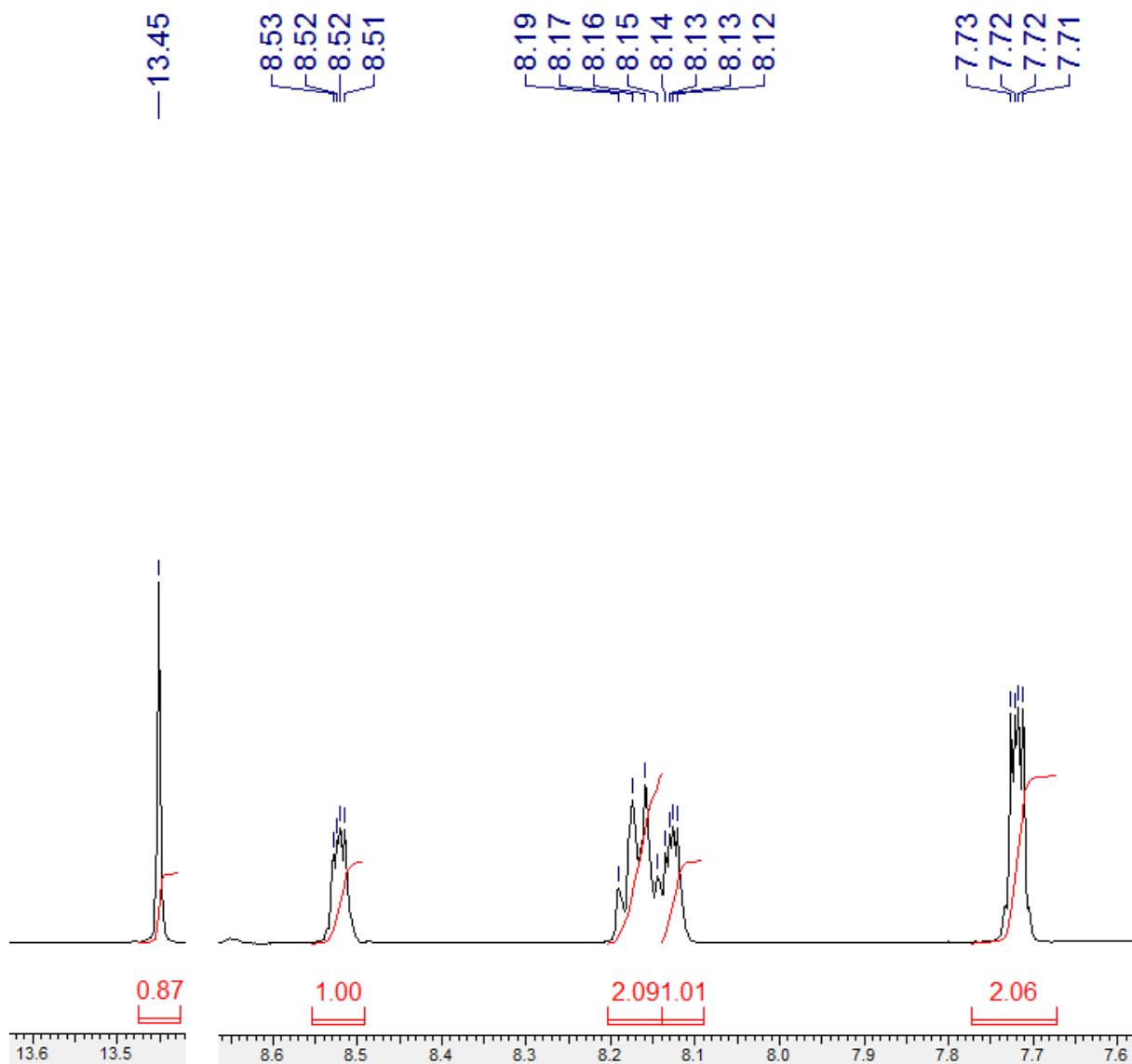


NMR  $^{13}\text{C}$  spectrum (150 MHz, DMSO-D6) of compound **4b**.

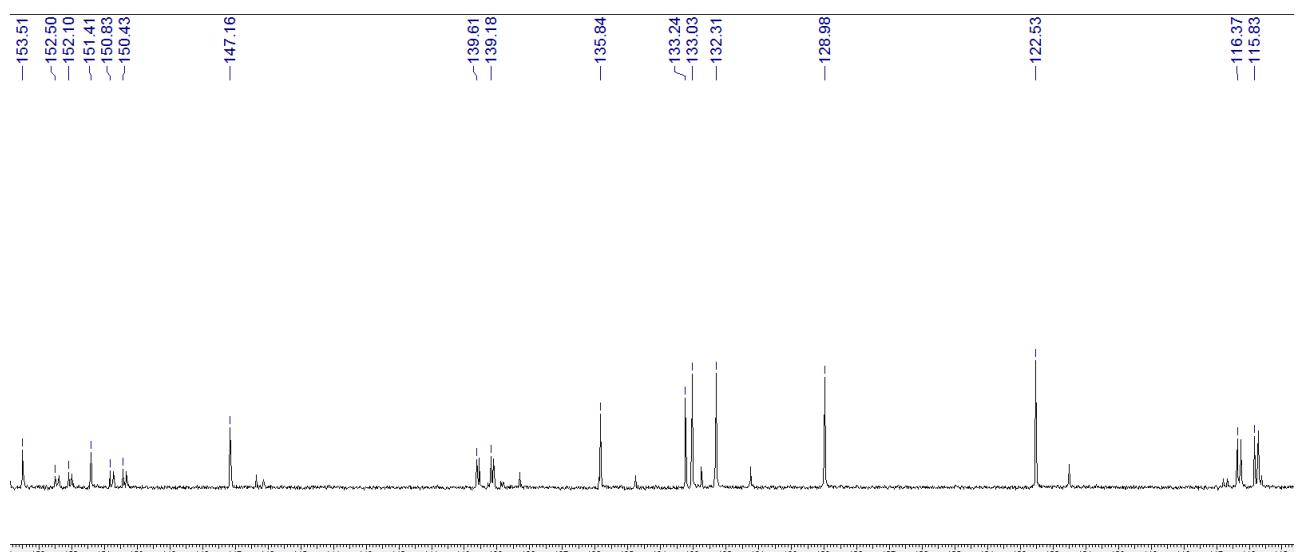




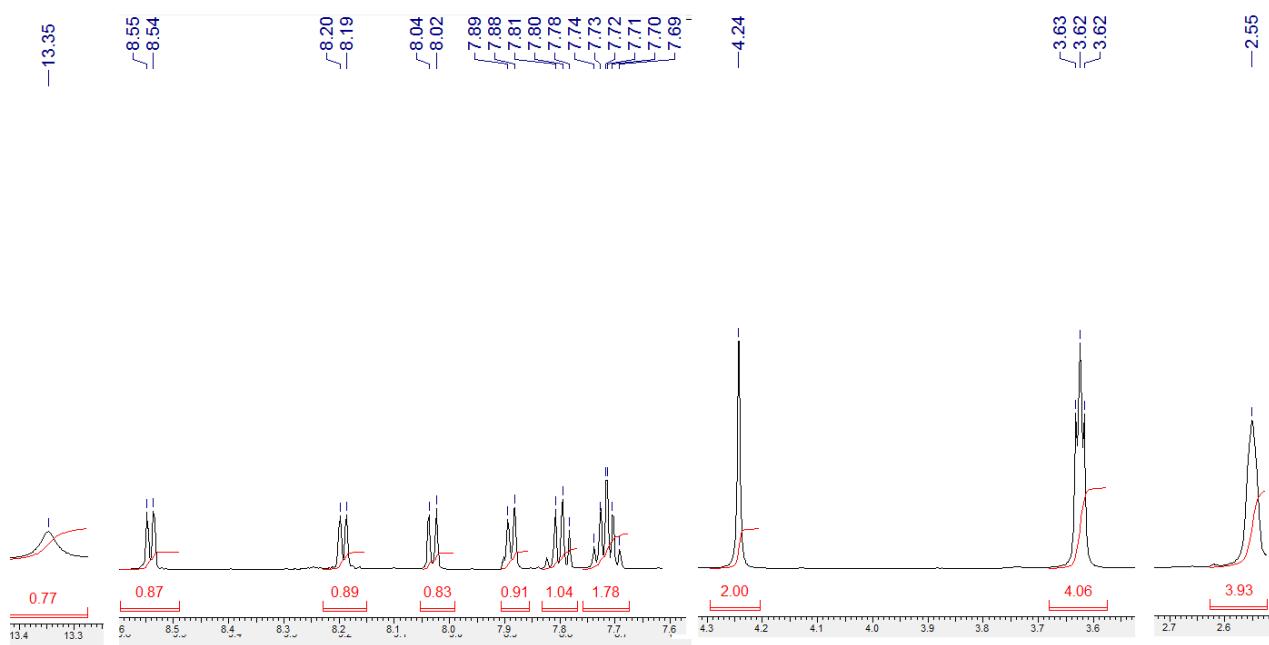
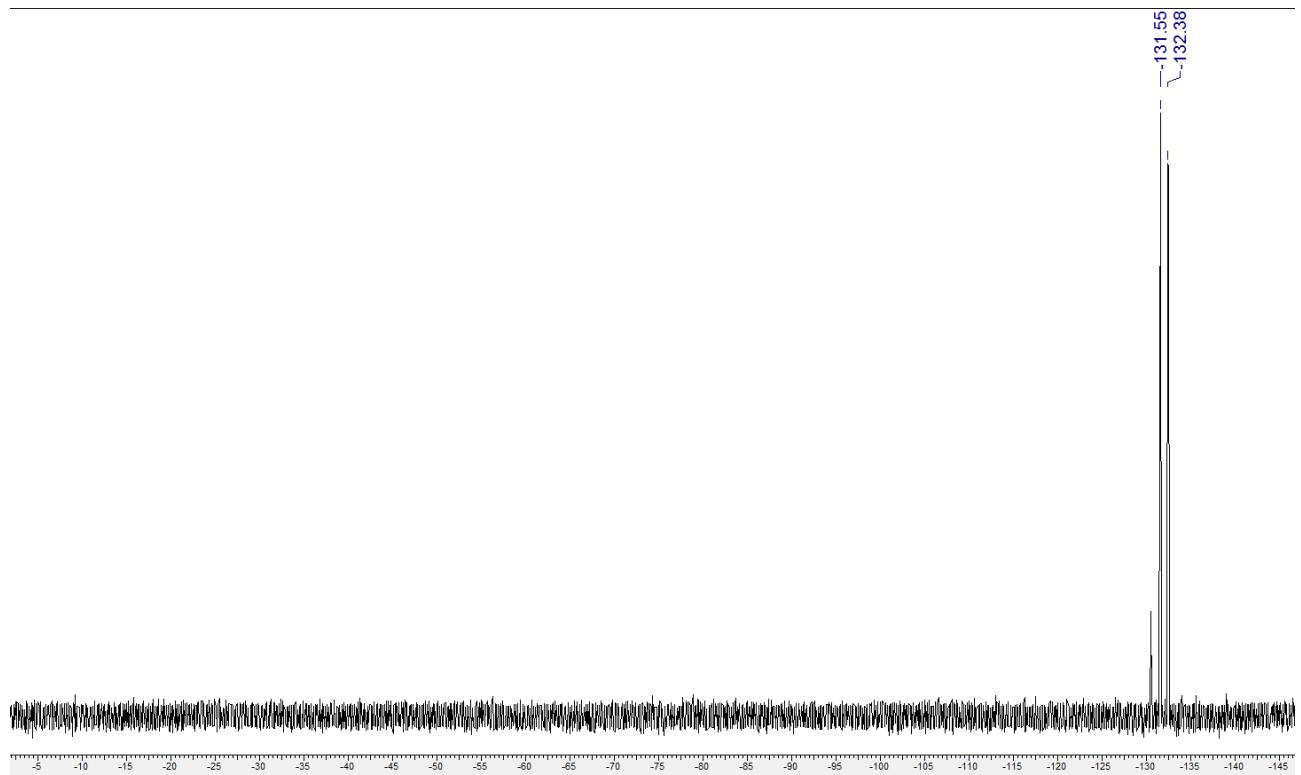
NMR  ${}^{19}\text{F}$  spectrum (280 MHz, DMSO-D6) of compound **4e**.

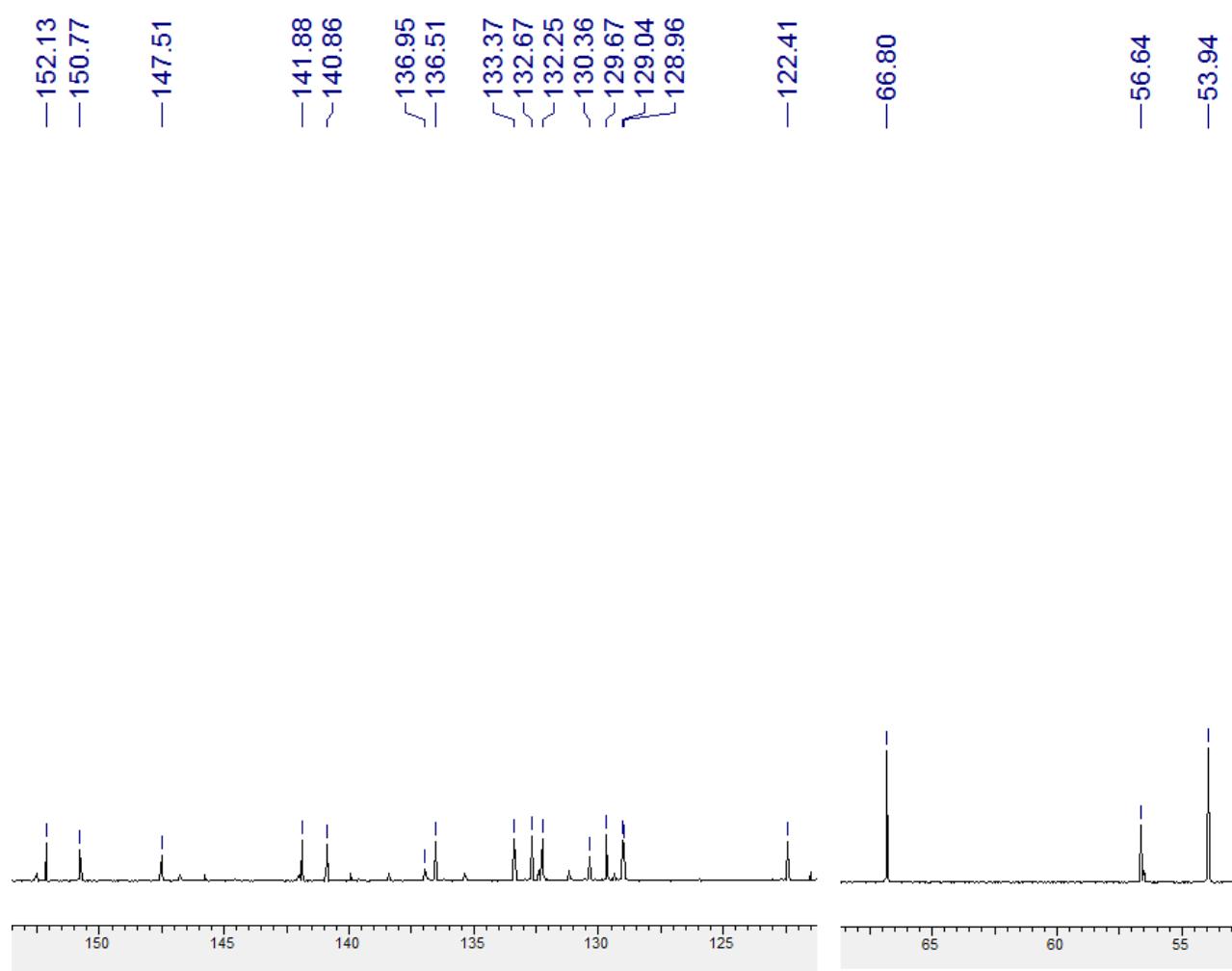


NMR  $^1\text{H}$  spectrum (600 MHz, DMSO-D<sub>6</sub>) of compound **4f**.



NMR  $^{13}\text{C}$  spectrum (150 MHz, DMSO-D<sub>6</sub>) of compound **4f**.





NMR  $^{13}\text{C}$  spectrum (150 MHz, DMSO-D6) of compound **4m**.