

## Supporting Information

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### **Monoterpene Flavonoid from Aerial Parts of *Satureja khuzistanica***

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*NMR data of compounds 2-6*

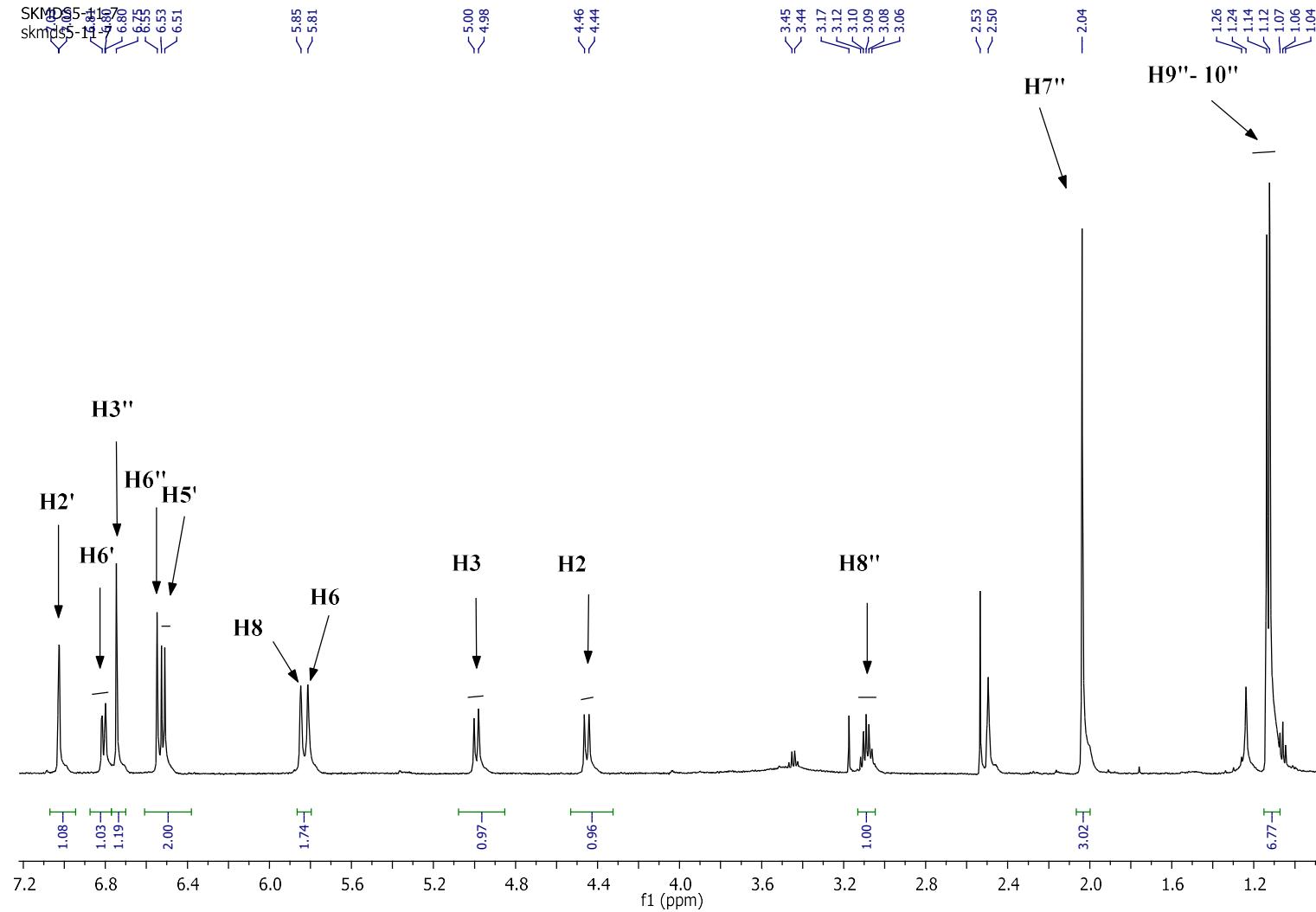
*Keshonin (2)*:  $^1\text{H}$  NMR (500 MHz, DMSO)  $\delta$ : 7.27 (1H, d,  $J = 2.2$  Hz, H-2'), 7.68 (1H, dd,  $J = 8.5, 2.2$  Hz, H-6'), 7.07 (1H, d,  $J = 8.5$  Hz, H-5'), 6.79 (1H, s, H-3''), 6.66 (1H, d,  $J = 2.0$  Hz, H-8), 6.66 (1H, s, H-3), 6.53 (1H, s, H-6''), 6.47 (1H, d,  $J = 2.1$  Hz, H-6), 5.21 (1H, d,  $J = 7.3$  Hz, H-1''), 3.98 (1H, d,  $J = 9.6$  Hz, H-5''), 3.56 – 3.05 (3H, m, H-2'', H-3'', H-4''), 3.16 – 3.08 (1H, m, H-8''), 2.04 (3H, s, H-7''), 1.19 (3H, d,  $J = 1.8$  Hz, H-9''), 1.18 (3H, d,  $J = 1.8$  Hz, H-10'').

*Saturejin (3'-(2,5-dihydroxy-p-cymene) 5,7,4'-trihydroxy flavone) (3)*:  $^1\text{H}$  NMR (500 MHz, DMSO)  $\delta$ : 6.61 (1H, s, H-3), 6.16 (1H, d,  $J = 2.0$ , H-6), 6.30 (1H, d,  $J = 2.0$ , H-8), 7.20 (1H, d,  $J = 2.2$ , H-2'), 7.05 (1H, d,  $J = 8.5$ , H-5'), 7.64 (1H, dd,  $J = 8.5, 2.2$ , H-6'), 6.78 (1H, s, H-1''), 6.57 (1H, s, H-6''), 2.04 (3H, s, H-7''), 3.11 (1H, m, H-8''), 1.15 (1H, d,  $J = 6.8$ , H-9''), 1.15 (1H, d,  $J = 6.8$ , H-10'').

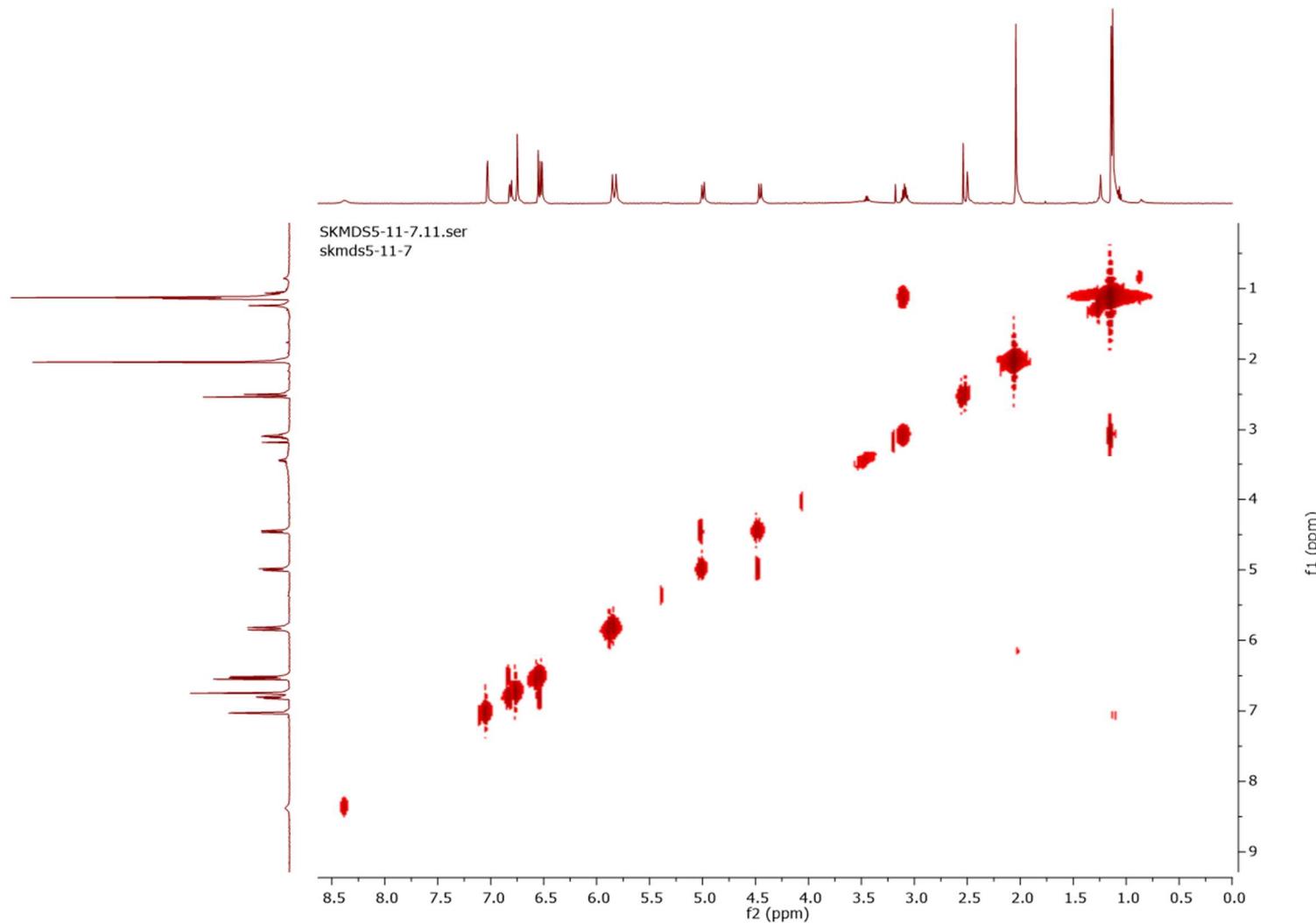
*Ponciretin (4)*:  $^1\text{H}$  NMR (500 MHz, DMSO)  $\delta$ : 7.39 (2H, d,  $J = 8.7$  Hz, H-2', H-6'), 6.93 (2H, d,  $J = 8.7$  Hz, H-3', H-5'), 5.80 (2H, d,  $J = 1.9$  Hz, H-6, H-8), 5.44 (1H, dd,  $J = 12.4, 3.0$  Hz, H-2), 3.74 (3H, s, OMe), 3.17 (1H, dd,  $J = 16.9, 12.3$  Hz, H-3ax), 2.67 (1H, dd,  $J = 17.1, 3.2$  Hz, H-3eq).

*5,6-dihydroxy-3',4',7-trimethoxyflavone (5)*:  $^1\text{H}$  NMR (500 MHz, DMSO)  $\delta$  7.60 (1H, s, H-3), 7.59 (1H, s, H-8), 6.96 (2H, d,  $J = 8.1$  Hz,), 6.91 (1H, d,  $J = 10.0$  Hz,), 3.94 (3H,s, OMe), 3.91 (3H,s, OMe), 3.75 (3H,s, OMe).

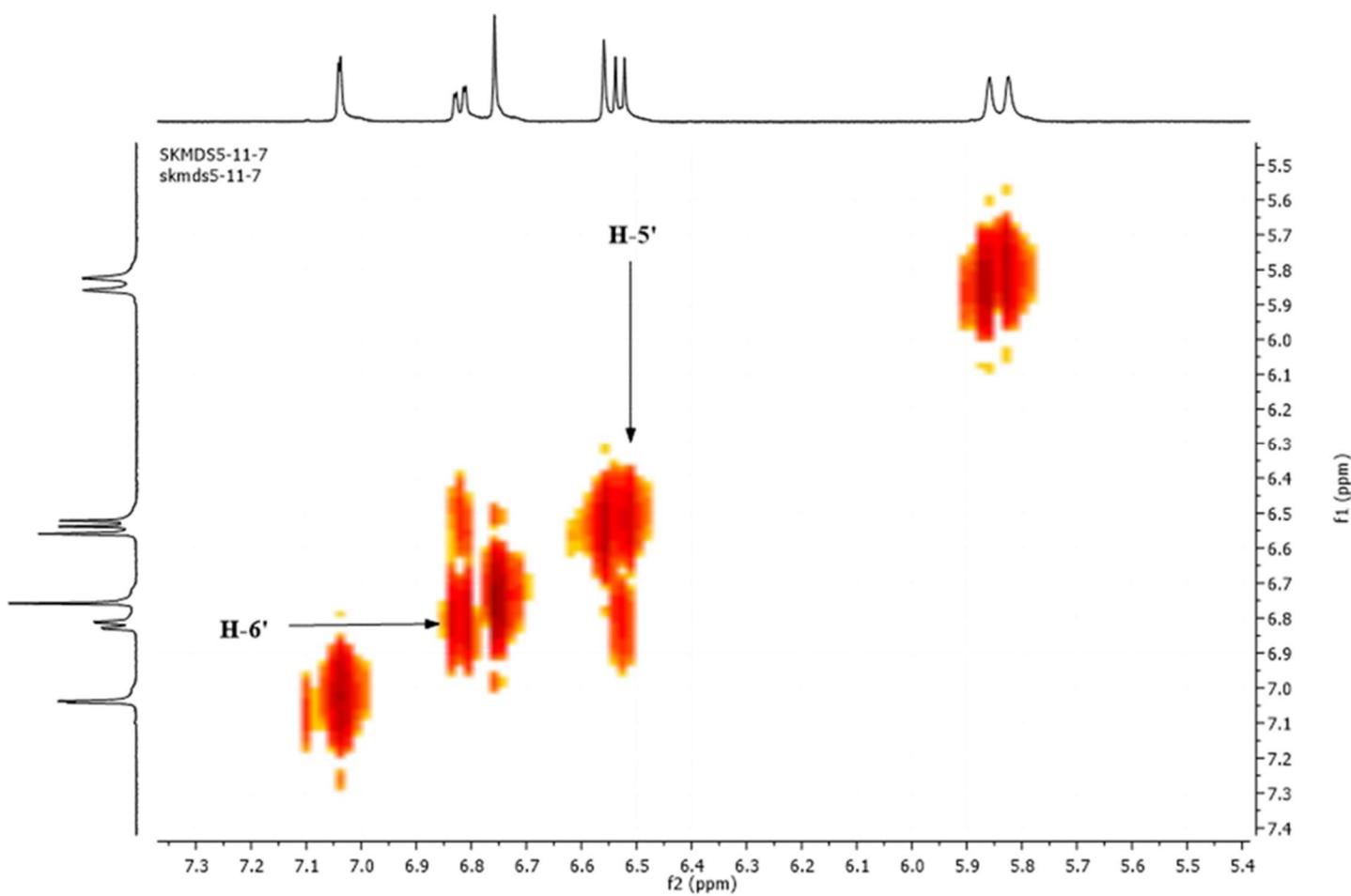
*5,6-dihydroxy-4',7-dimethoxyflavone (6)*:  $^1\text{H}$  NMR (500 MHz, DMSO)  $\delta$ : 7.95 (2H, d,  $J = 8.8$  Hz,H-2', H-6'), 6.94 (2H, d,  $J = 8.7$  Hz, H-3', H-5'), 6.90 (1H, s, H-3), 6.80 (1H, s, H-8), 3.94 (3H,s, OMe), 3.75 (3H,s,OMe).



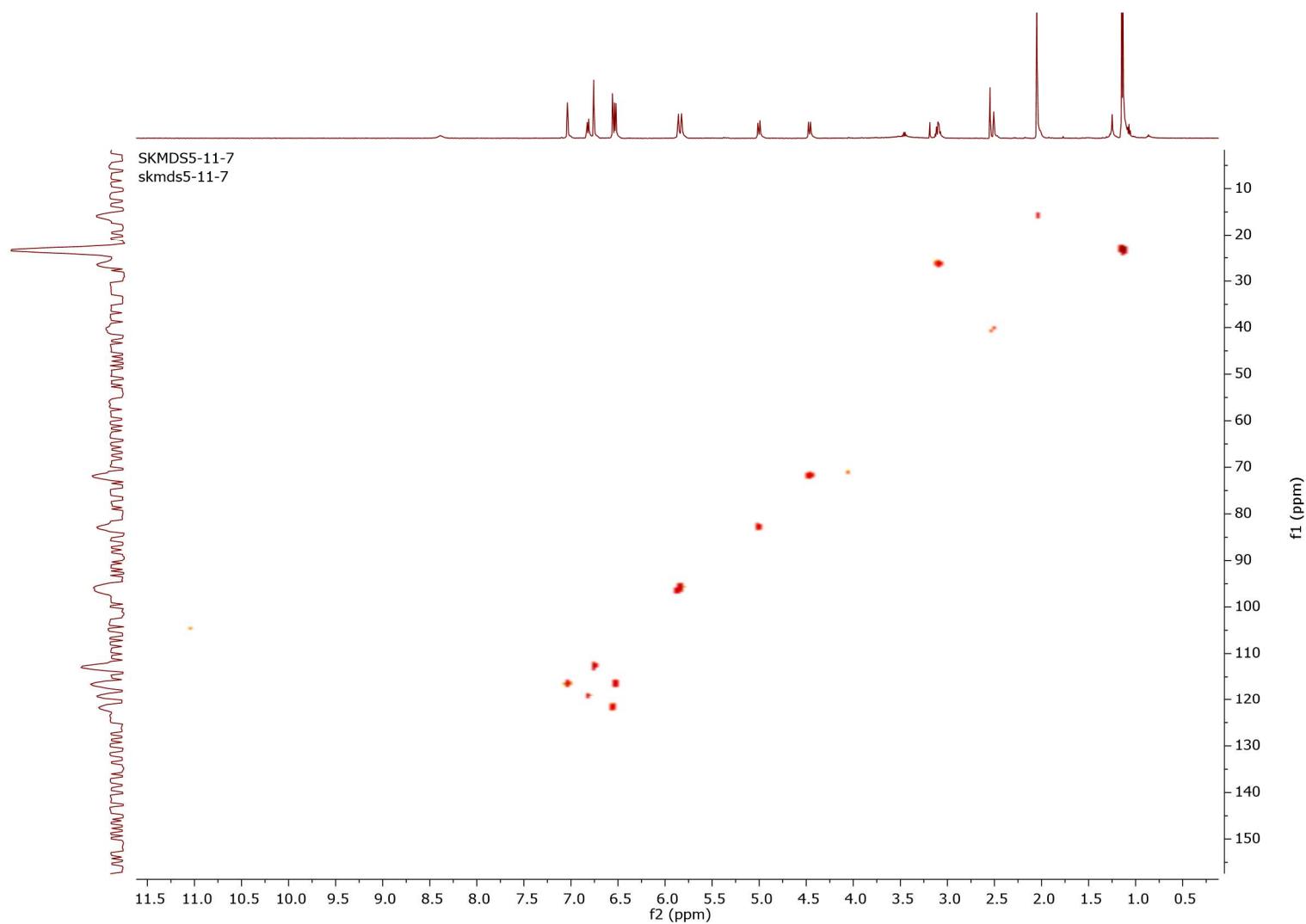
**Figure S1:**  $^1\text{H}$ NMR (500 MHz, DMSO- $d_6$ ) spectrum of **1**



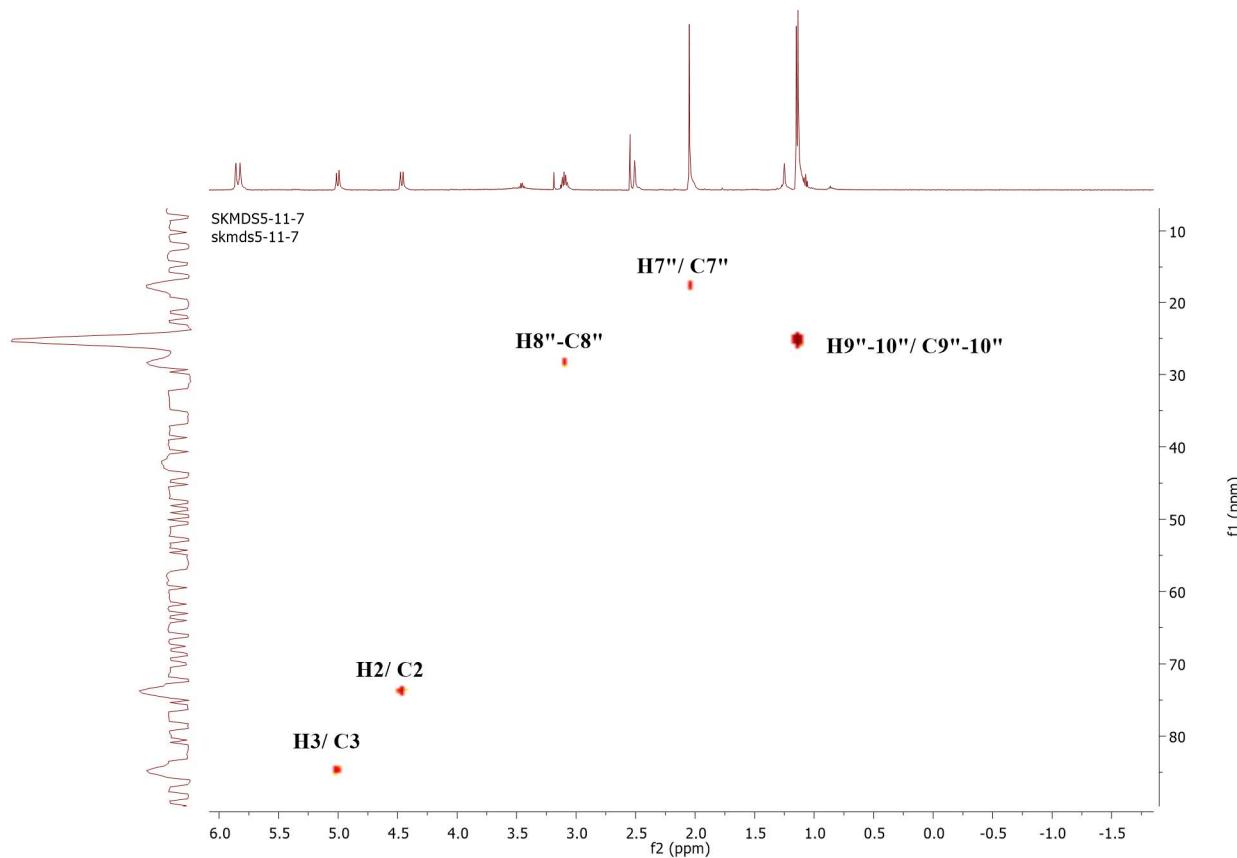
**Figure S2:**  $^1\text{H}$   $^1\text{H}$  COSY (500 MHz, DMSO-d<sub>6</sub>) spectrum of **1**



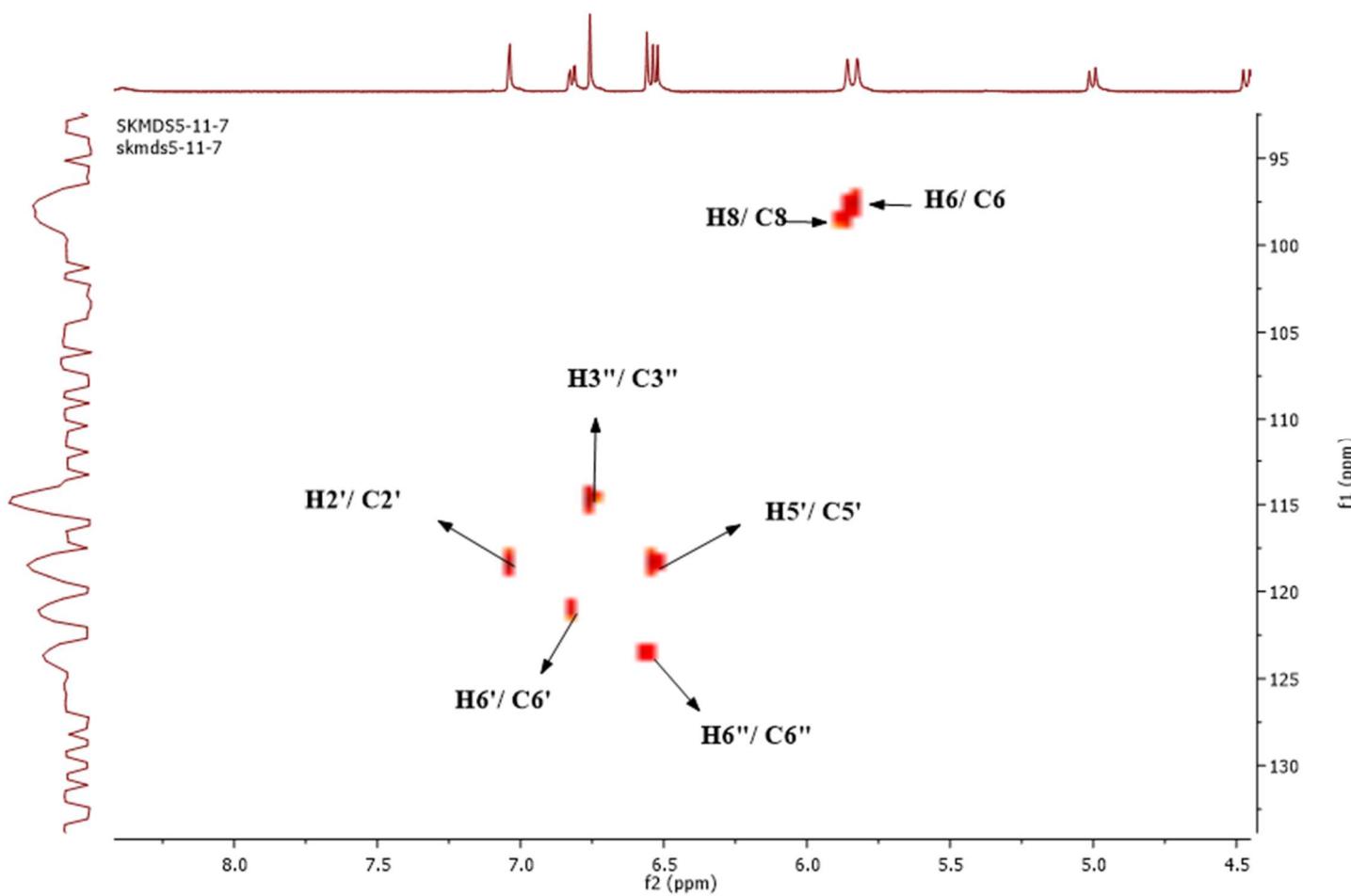
**Figure S3:** Expanded  $^1\text{H}$   $^1\text{H}$  COSY (500 MHz, DMSO-d<sub>6</sub>) spectrum of **1**



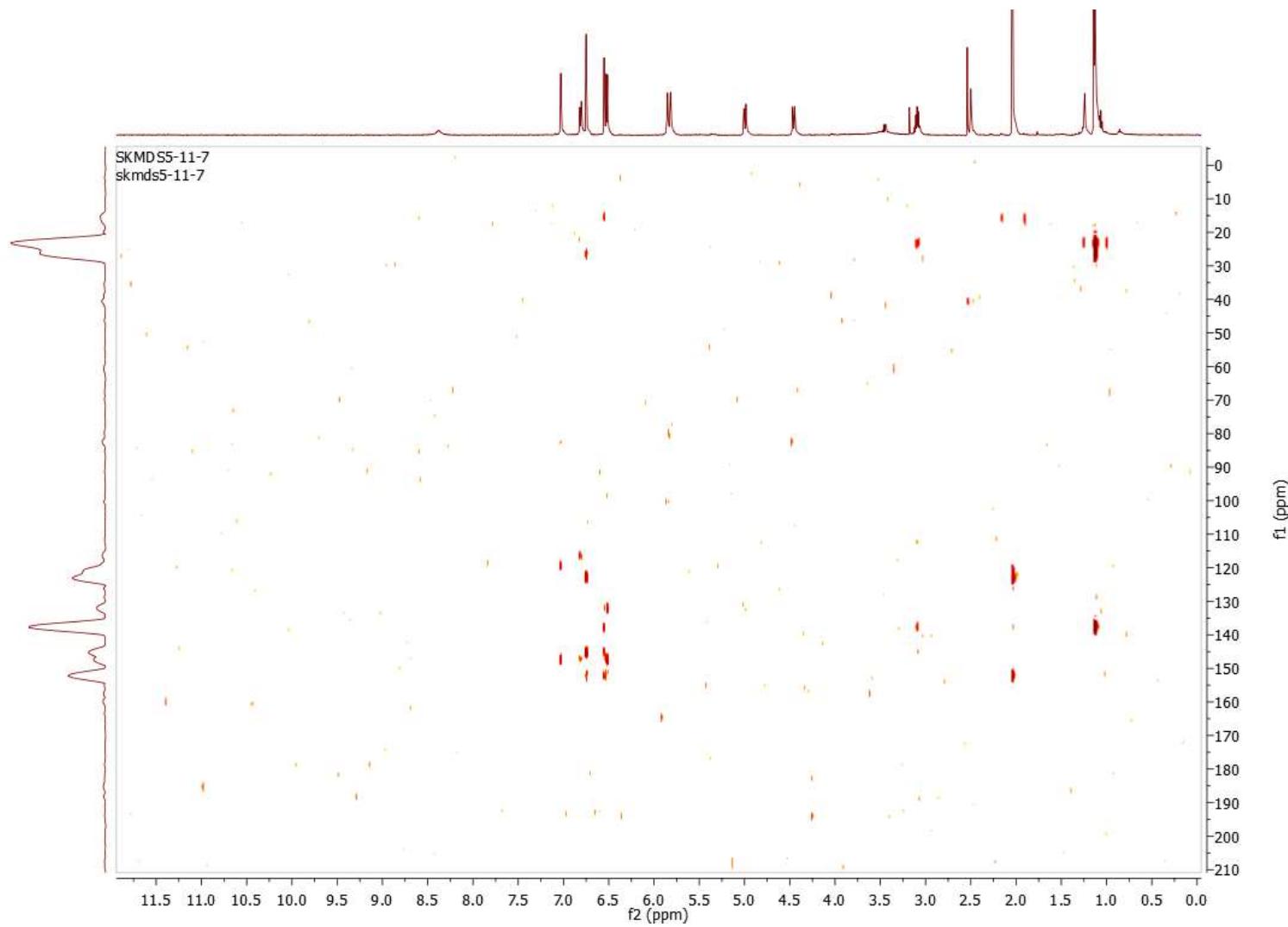
**Figure S4:** HSQC (500 MHz, DMSO-d<sub>6</sub>) spectrum of **1**



**Figure S5:** Expanded HSQC (500 MHz, DMSO-d6) spectrum of **1**



**Figure S6:** Expanded HSQC (500 MHz, DMSO-d6) spectrum of **1**



**Figure S7:** HMBC (500 MHz, DMSO-d6) spectrum of **1**

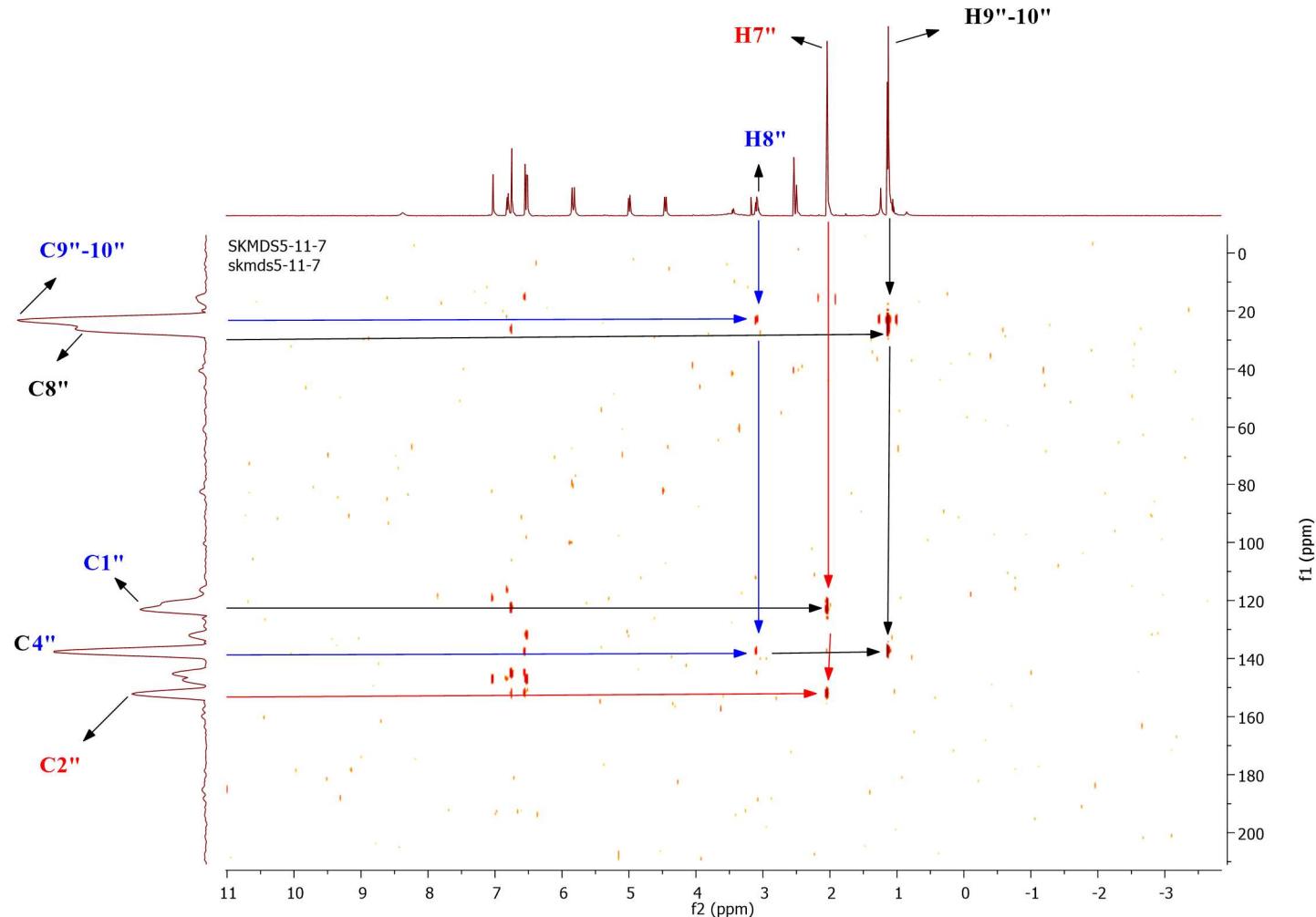
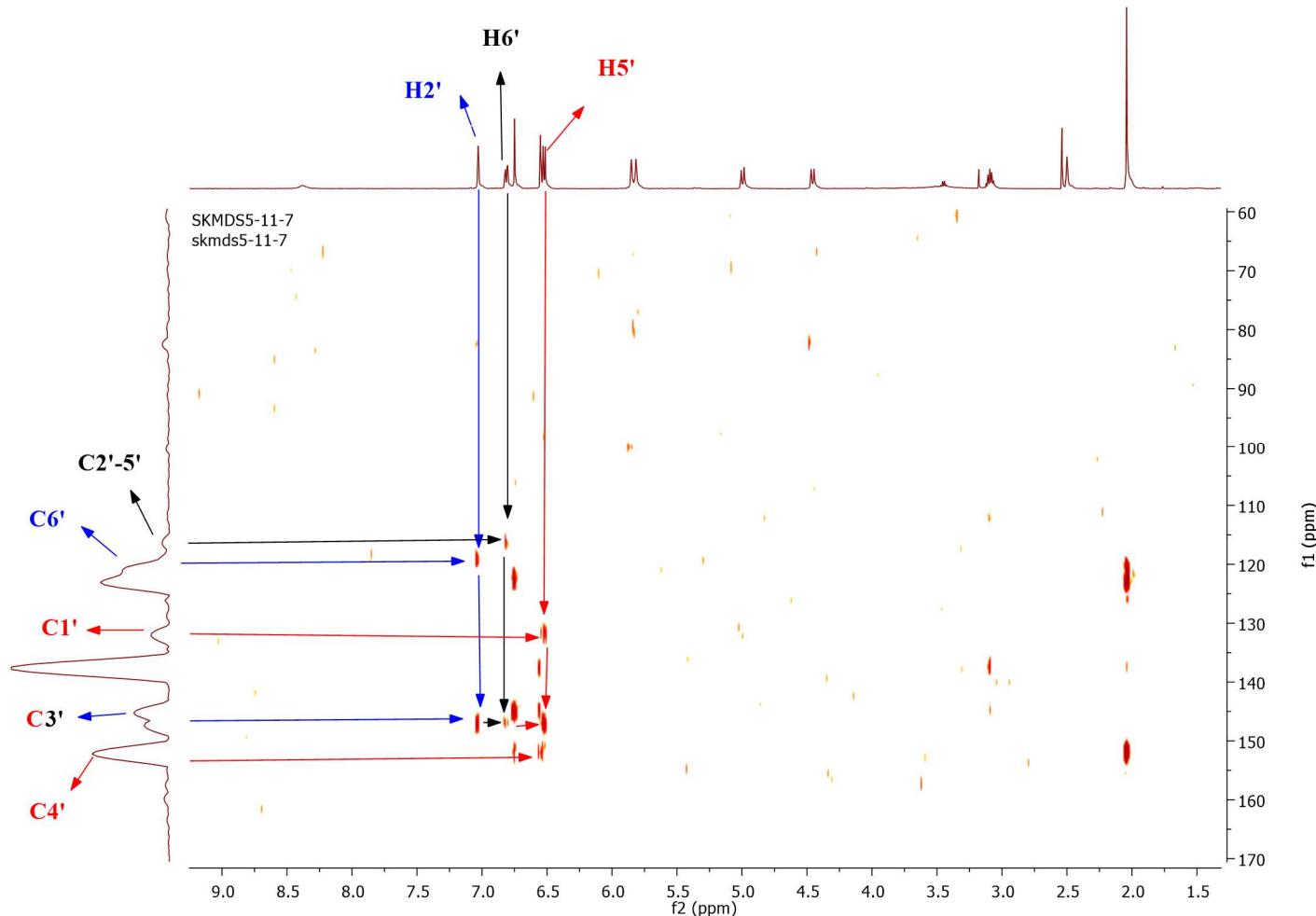


Figure S8: Expanded HMBC (500 MHz, DMSO-d<sub>6</sub>) spectrum of **1**



**Figure S9:** Expanded HMBC (500 MHz, DMSO-d<sub>6</sub>) spectrum of **1**

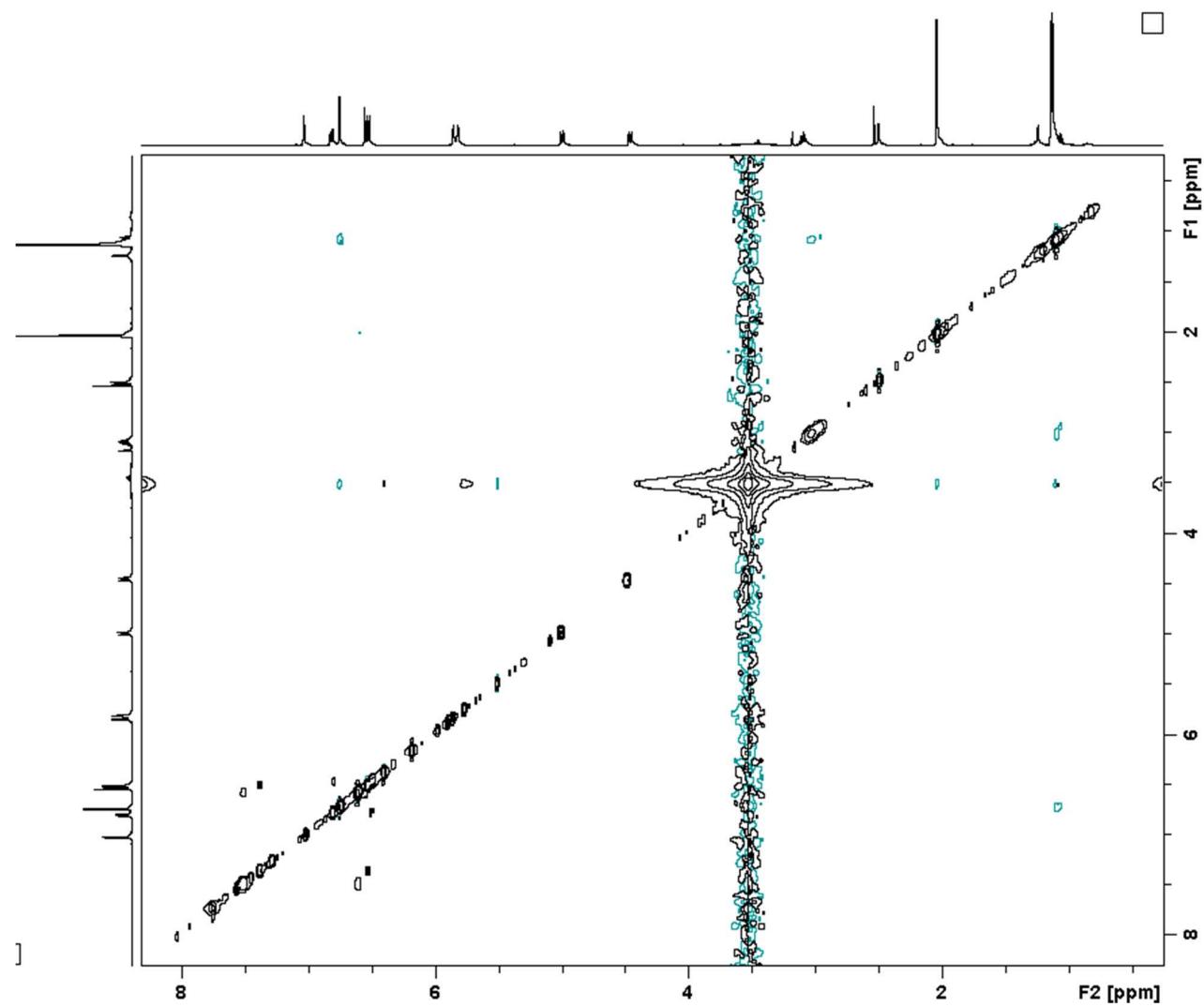
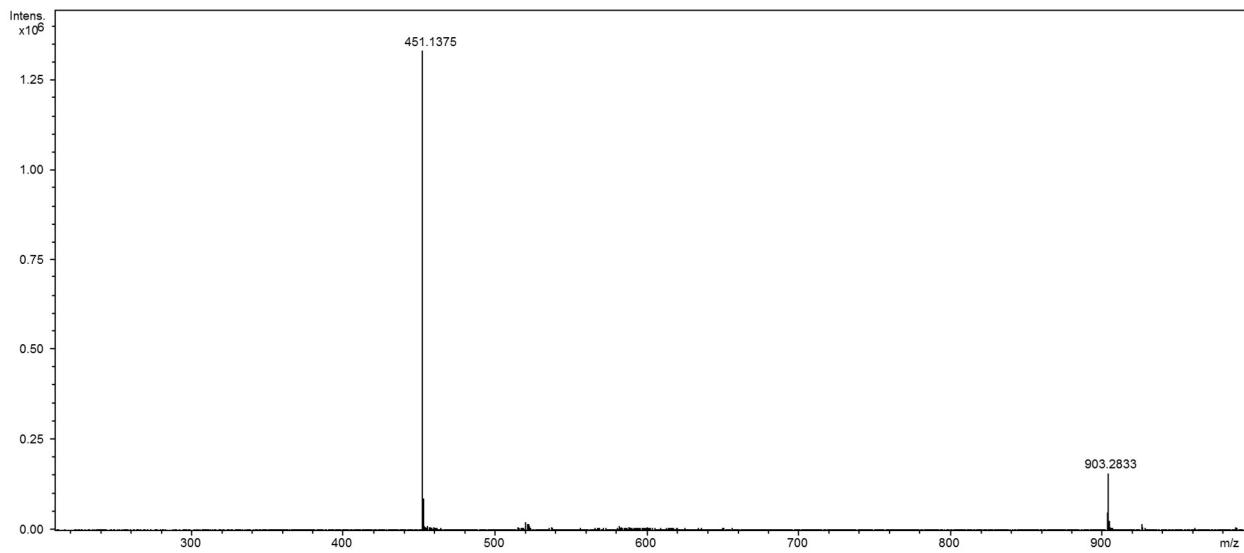
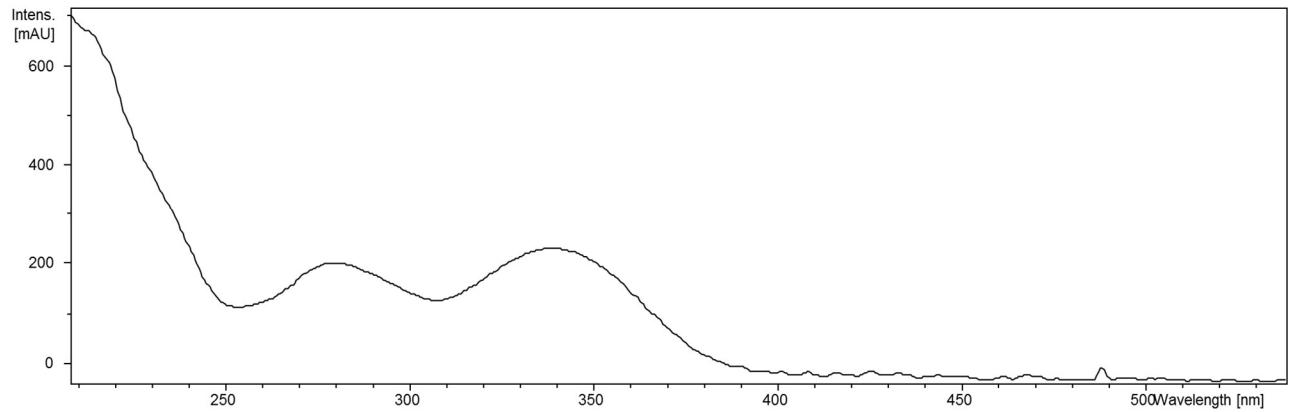


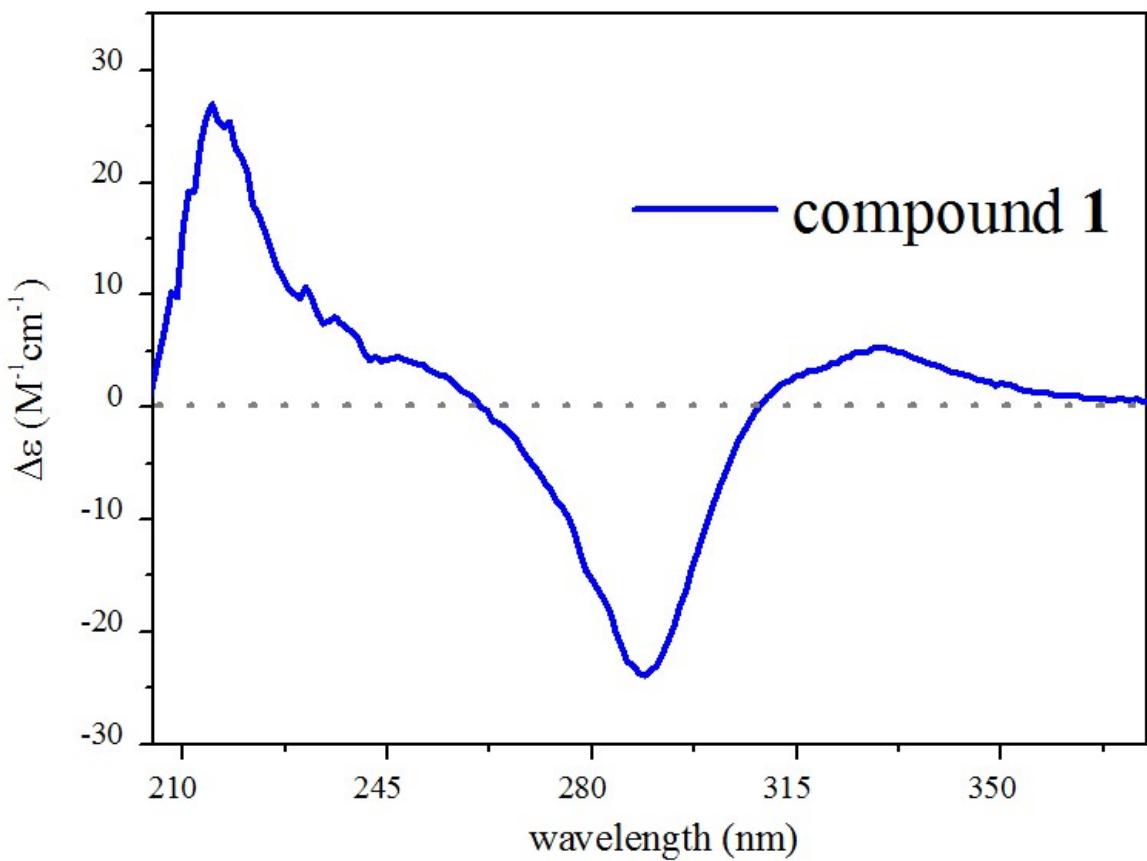
Figure S10: NOESY (500 MHz, DMSO-d<sub>6</sub>) spectrum of compound 1



**Figure S11:** HR-ESIMS spectrum of **1**



**Figure S12:** UV-Vis spectrum of compound **1**



**Figure S13:** Experimental ECD spectrum of compound 1