

## Supporting Information

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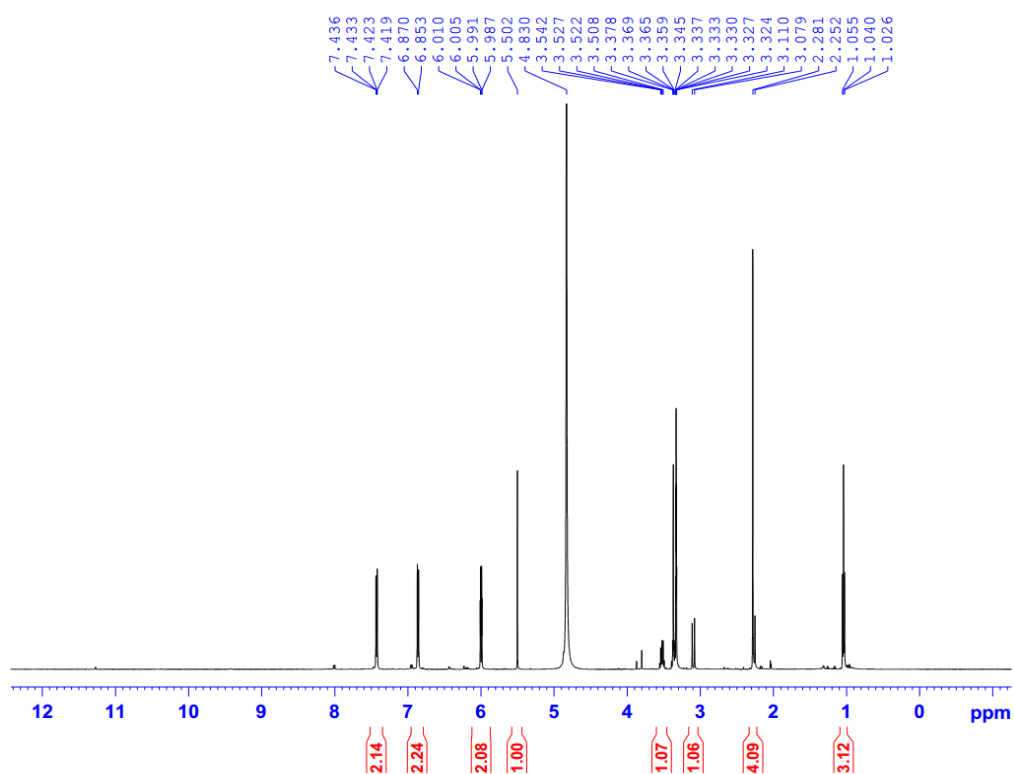
### A New 2,3-Dioxygenated Flavanone and Other Constituents from *Dyosma difformis*

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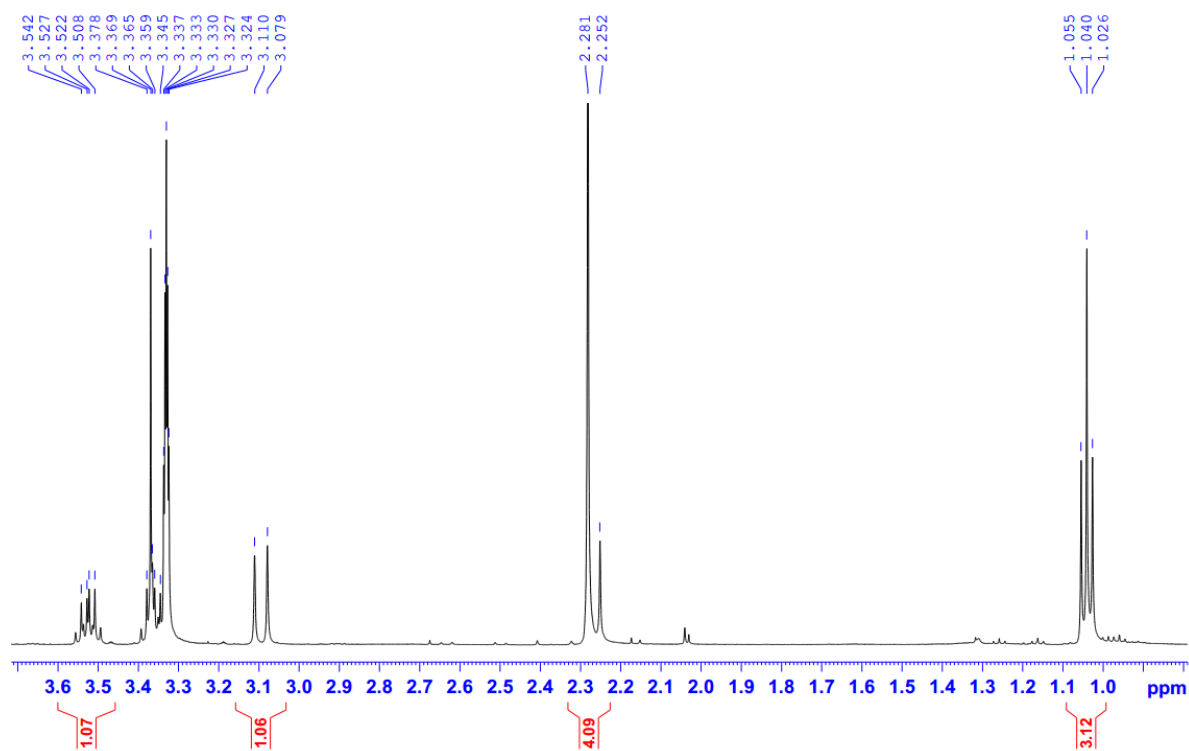
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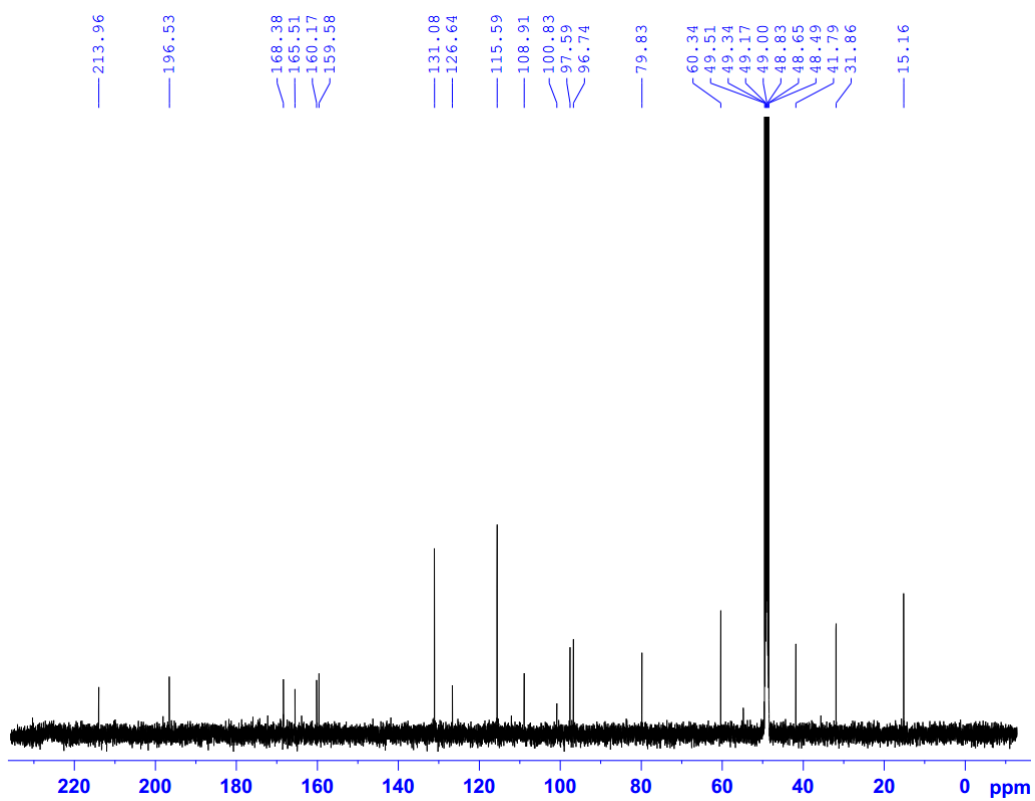
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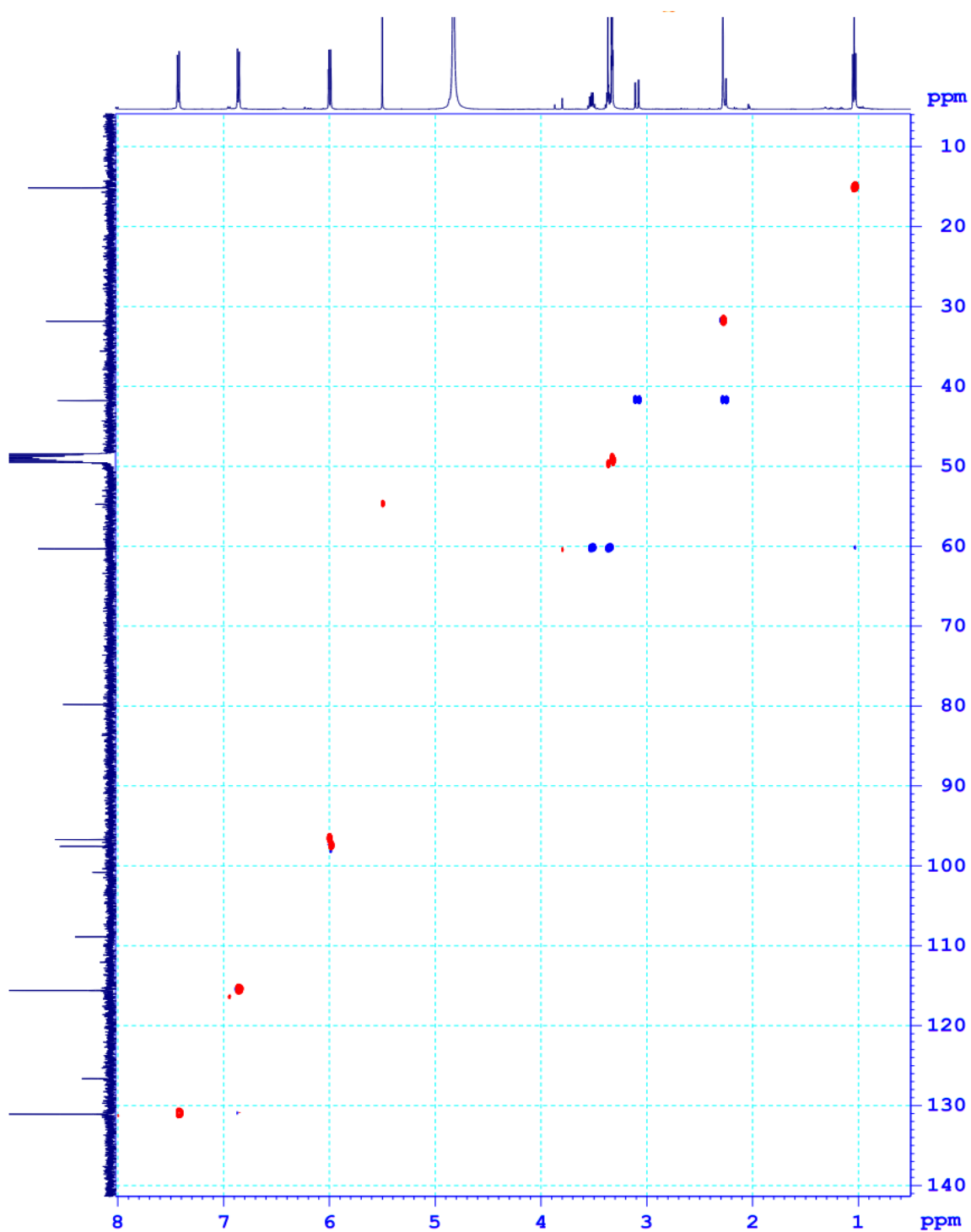
**Figure S1:** The  $^1\text{H}$ -NMR spectrum of the compound (**2**) in  $\text{CD}_3\text{OD}$



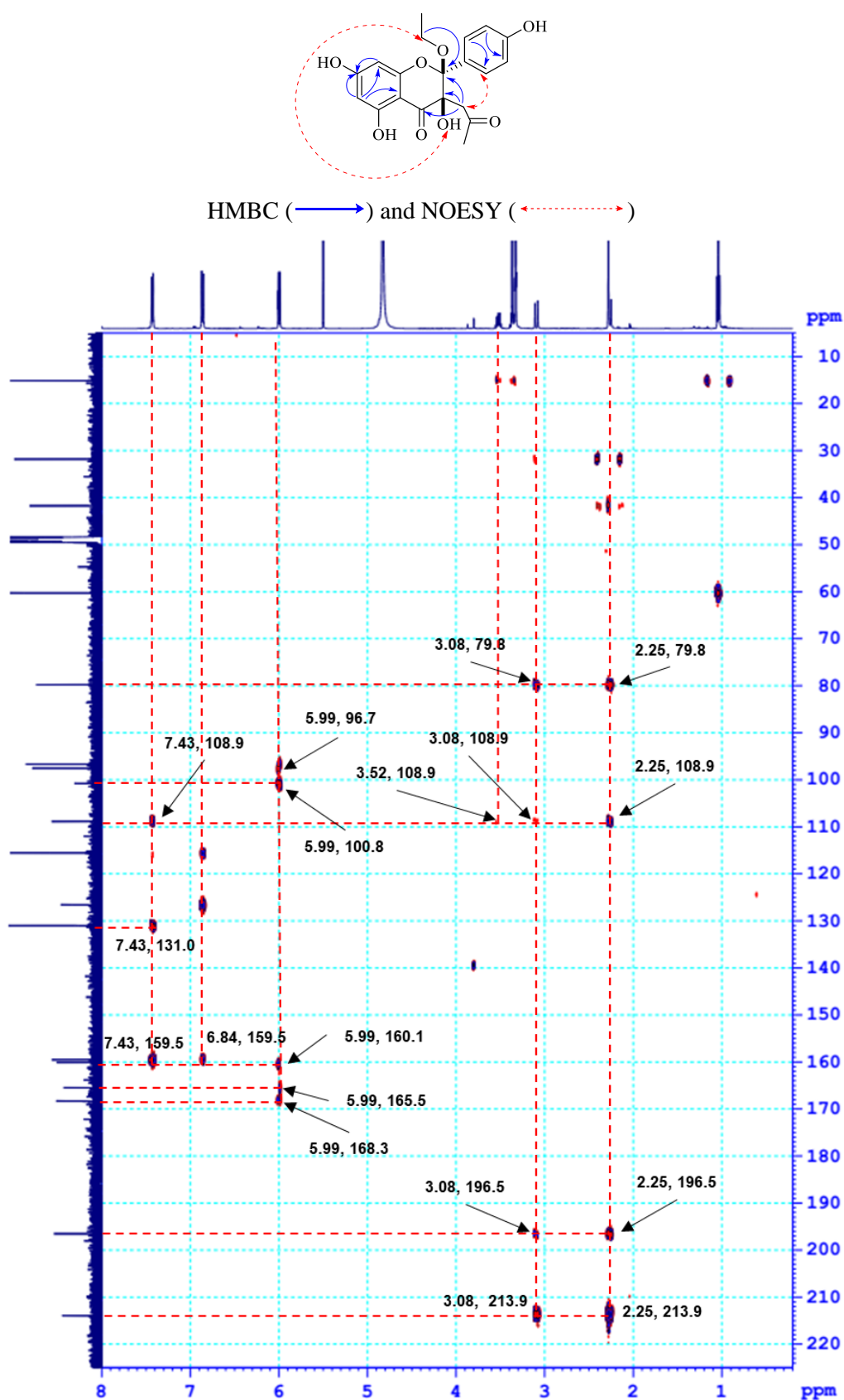
**Figure S2:** The extended  $^1\text{H}$ -NMR spectrum of the compound (**2**) in  $\text{CD}_3\text{OD}$



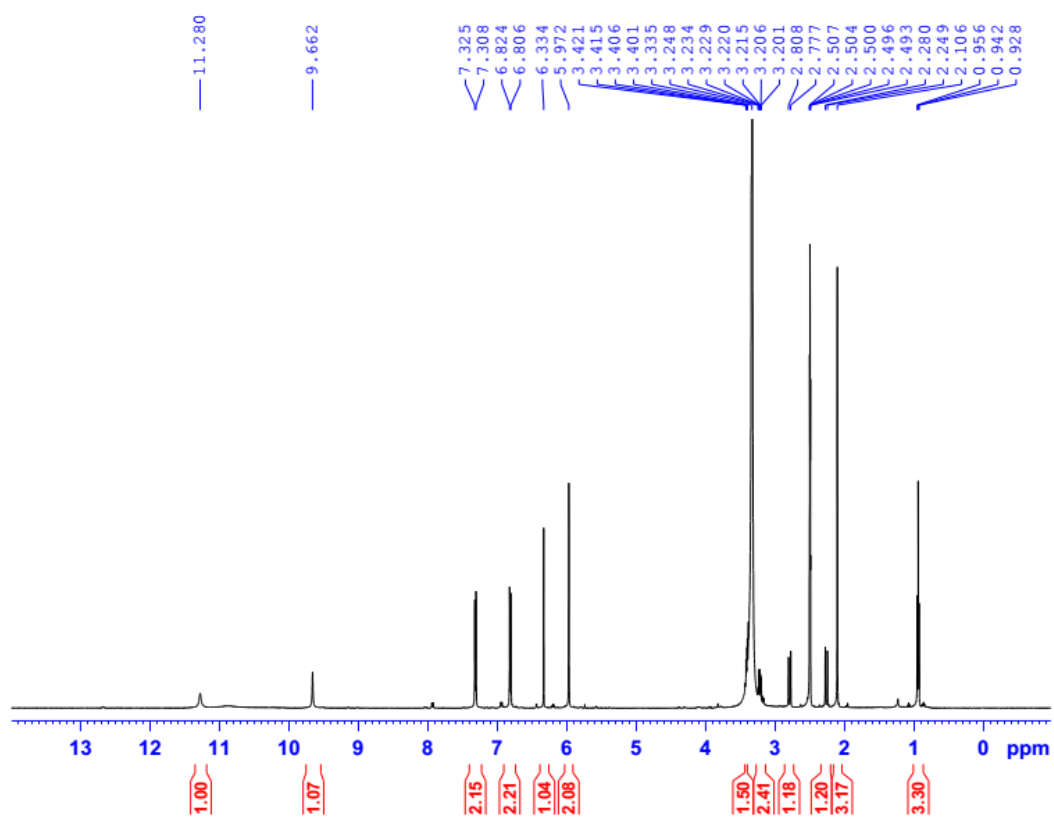
**Figure S3:** The  $^{13}\text{C}$ -NMR spectrum of the compound (2) in  $\text{CD}_3\text{OD}$



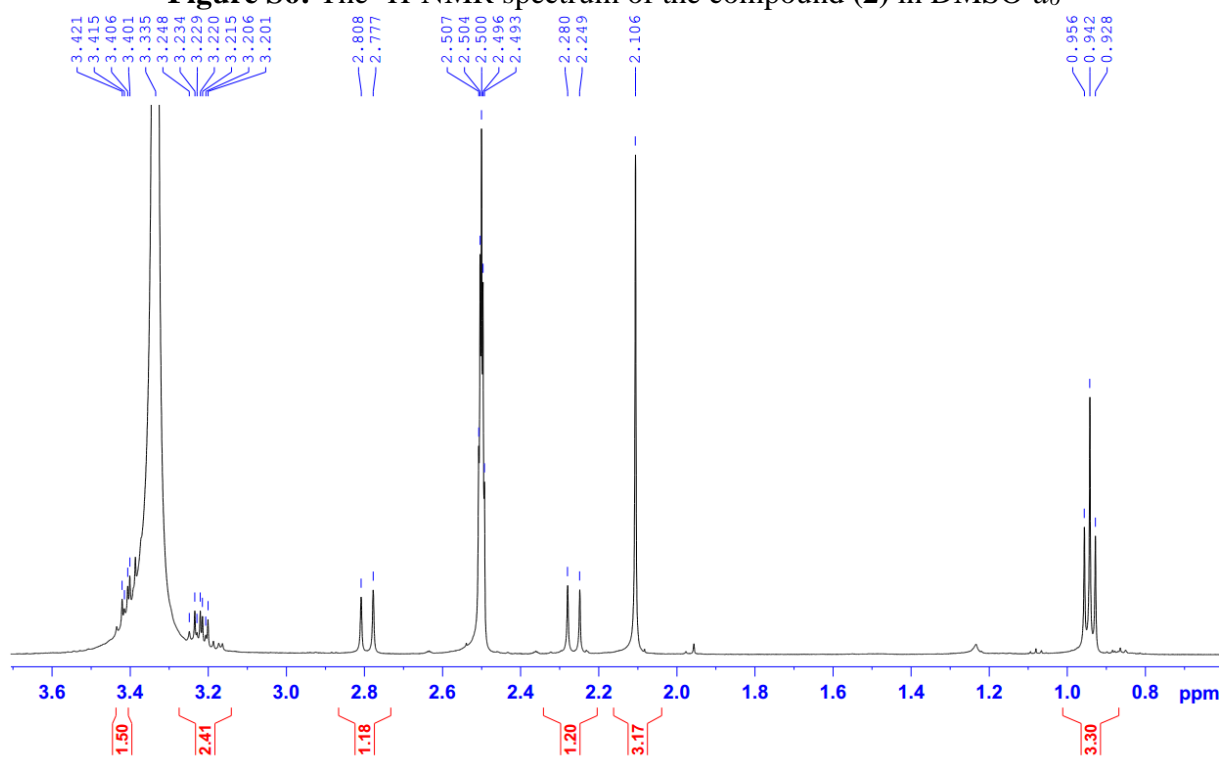
**Figure S4:** HSQC spectrum of the compound (2) in CD<sub>3</sub>OD



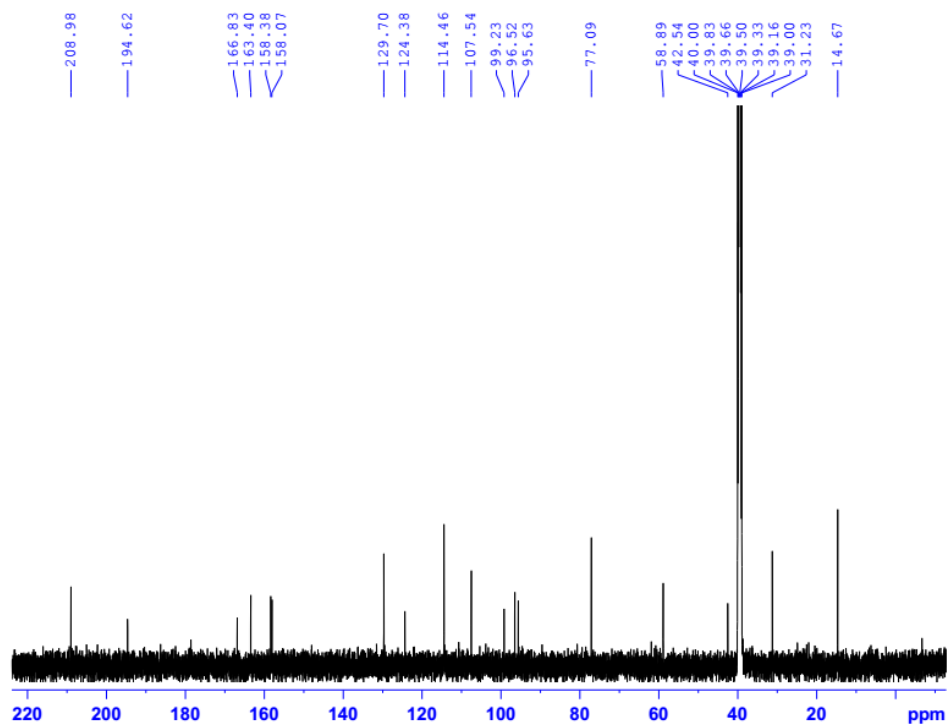
**Figure S5:** HMBC spectrum of the compound (2) in CD<sub>3</sub>OD



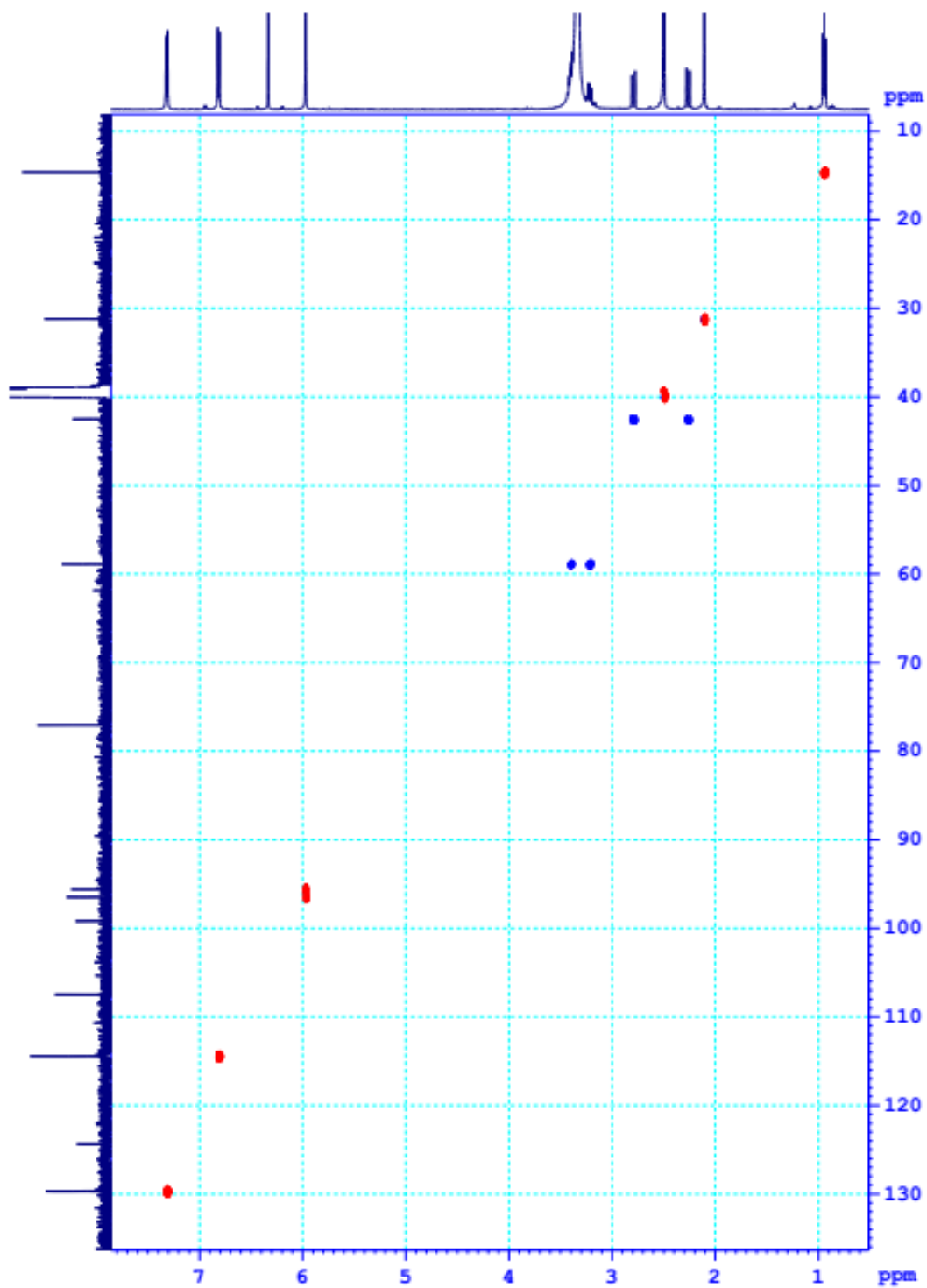
**Figure S6:** The  $^1\text{H}$ -NMR spectrum of the compound (2) in  $\text{DMSO-}d_6$



**Figure S7:** The extended  $^1\text{H}$ -NMR spectrum of the compound (2) in  $\text{DMSO-}d_6$



**Figure S8:** The  $^{13}\text{C}$ -NMR spectrum of the compound (2) in  $\text{DMSO}-d_6$

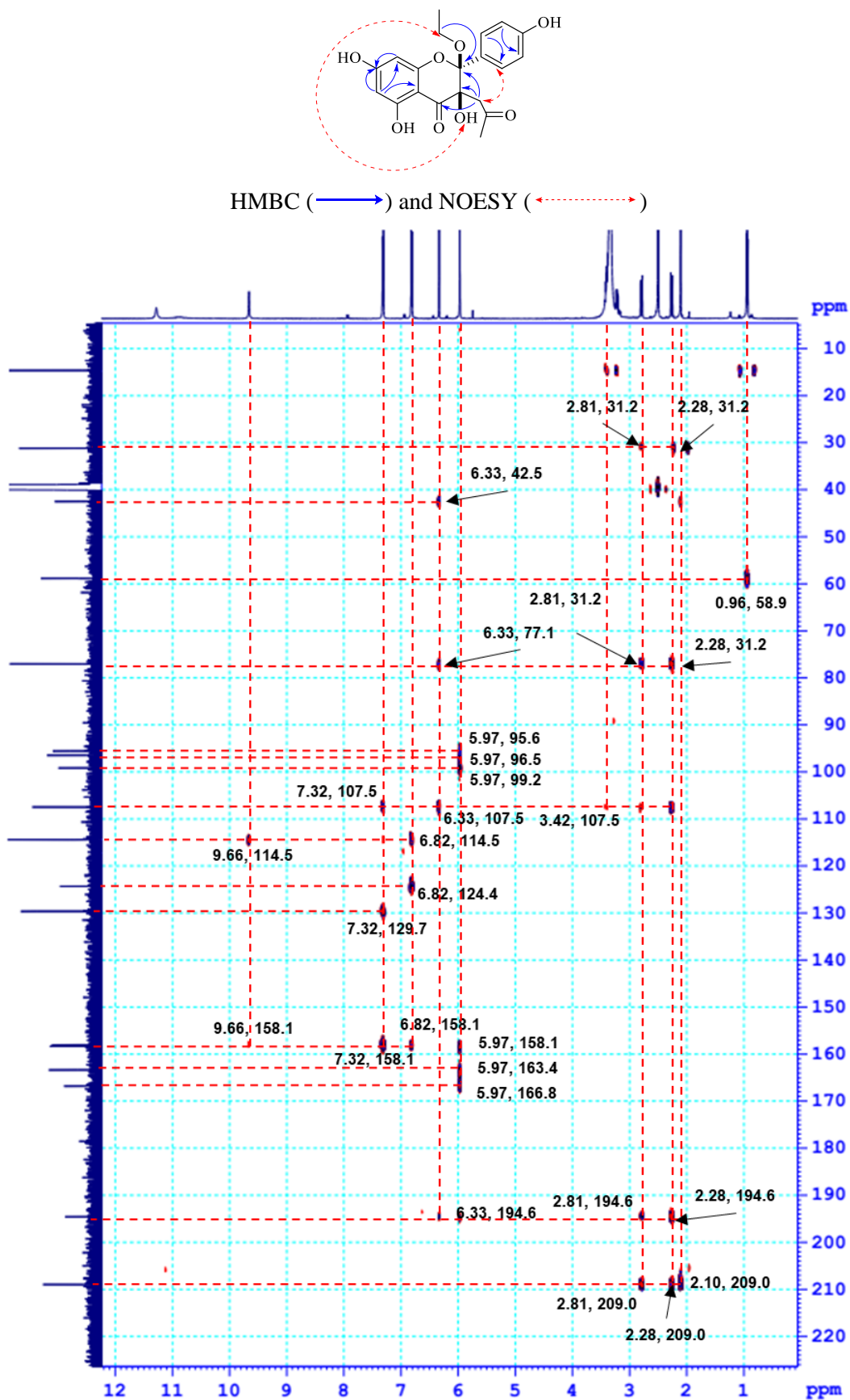


**Figure S9:** HSQC spectrum of the compound (**2**) in DMSO-*d*<sub>6</sub>

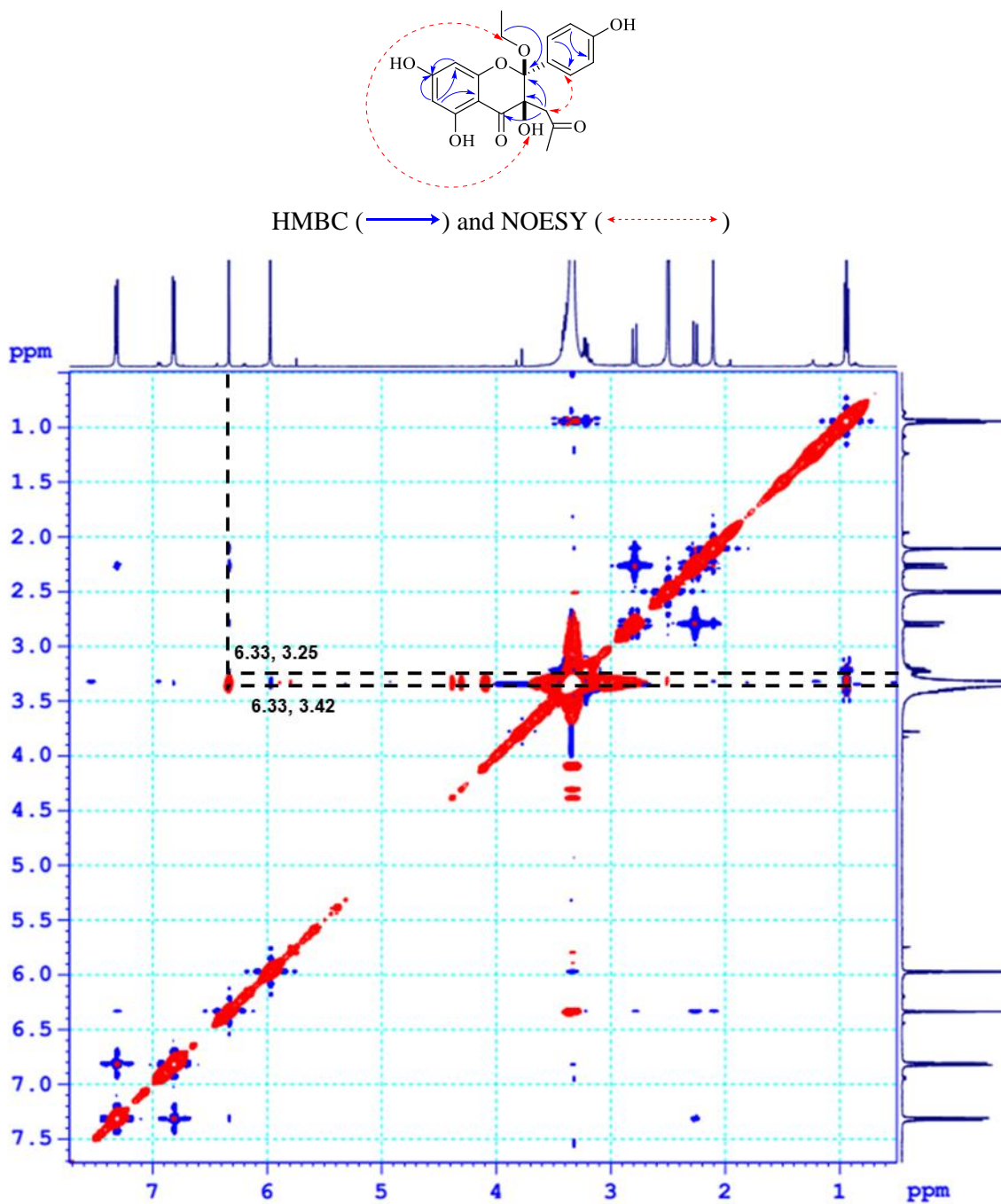




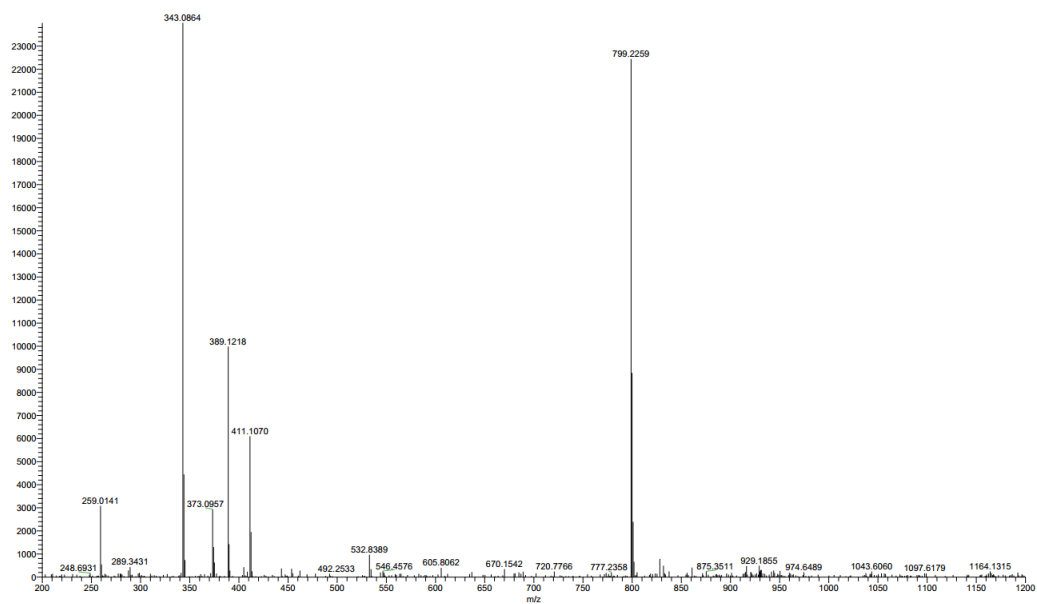
**Figure S10:** COSY spectrum of the compound (2) in DMSO-*d*<sub>6</sub>



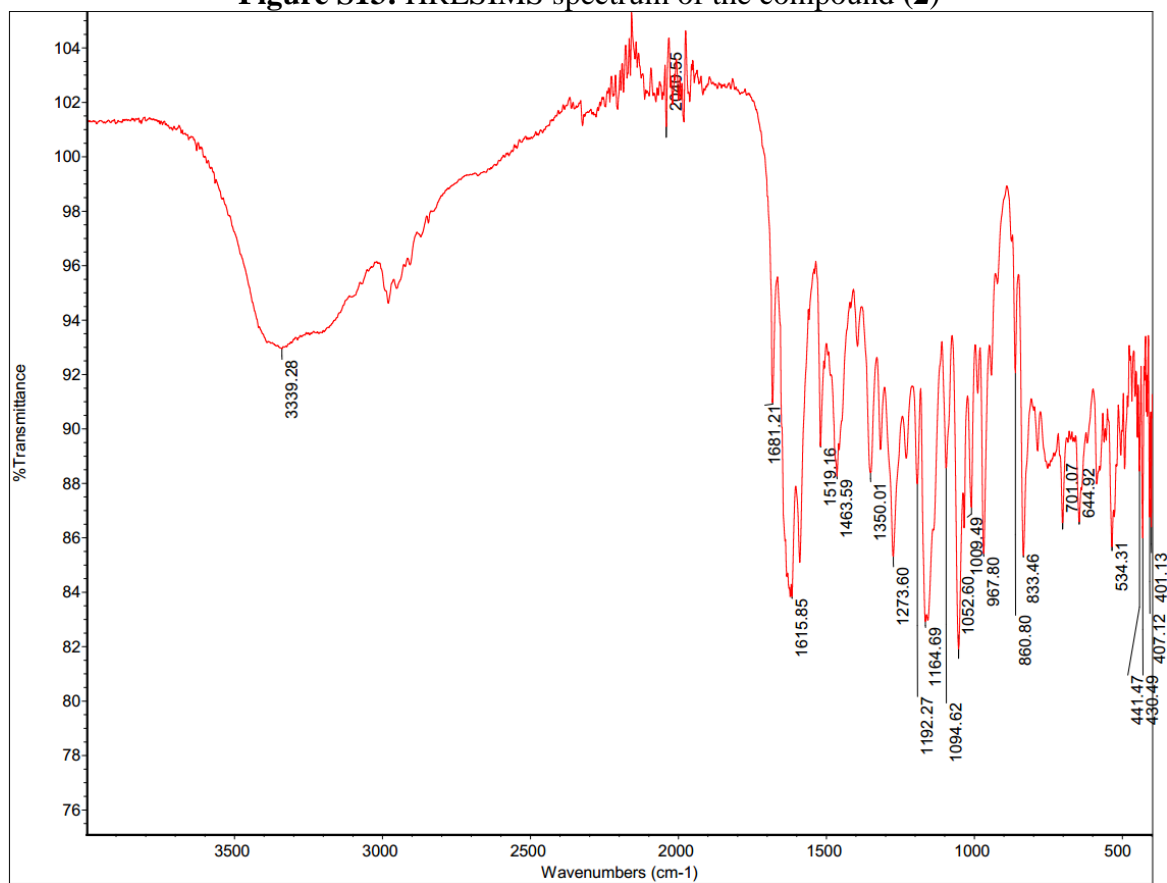
**Figure S11:** HMBC spectrum of the compound (2) in DMSO-*d*<sub>6</sub>



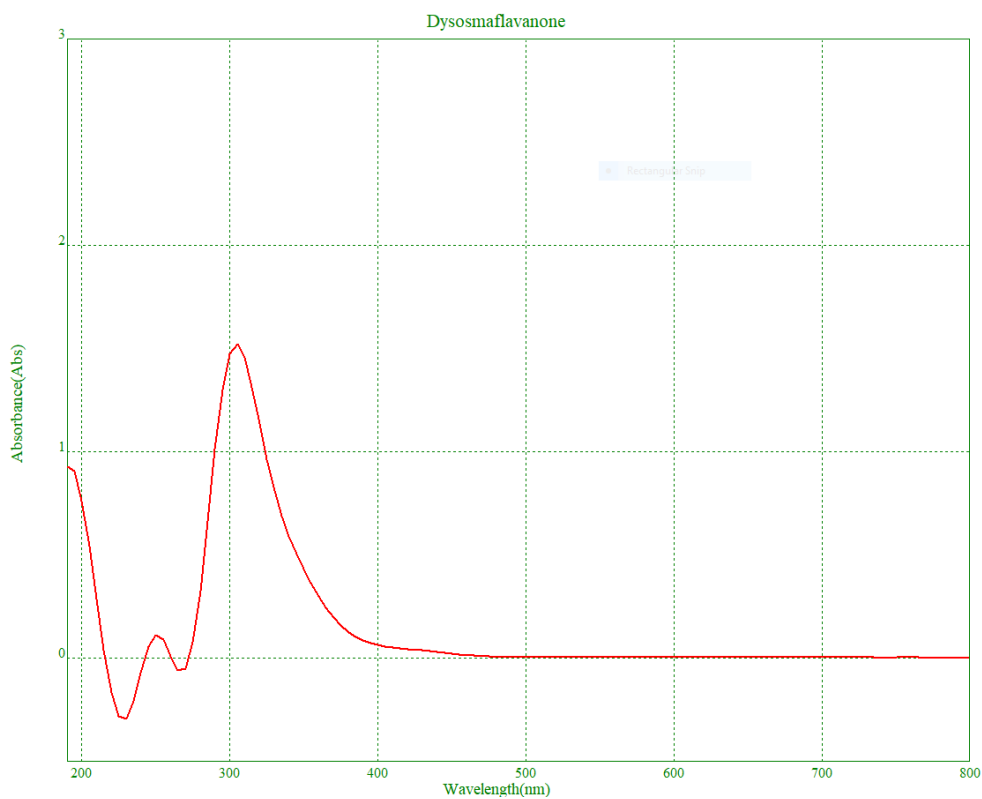
**Figure S12:** NOESY spectrum of the compound (**2**) in DMSO- $d_6$



**Figure S13:** HRESIMS spectrum of the compound (2)



**Figure S14:** IR spectrum of the compound (2)



**Figure S15:** UV-Vis spectrum of the compound (2)

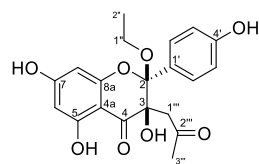
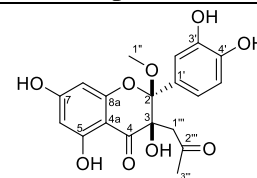
Substances (5)

2212305-01-4

C20H20O9  
 4*H*-1-Benzopyran-4-one, 2-(3,4-dihydroxyphenyl)-2-ethoxy-2,3-dihydro-3,5,7-trihydroxy-3-(2-oxopropyl)-

Key Physical Properties	Value	Condition
Molecular Weight	404.37	-
Boiling Point (Predicted)	716.3±60.0 °C	Press: 760 Torr
Density (Predicted)	1.59±0.1 g/cm <sup>3</sup>	Temp: 20 °C; Press: 760 Torr
pKa (Predicted)	7.17±0.60	Most Acidic Temp: 25 °C

**Figure S16:** The Scifinder search for the new compound (2)

**Table S1:** NMR data in CD<sub>3</sub>OD for the compound (**2**) and the reference Cepaflava B**Dysosmaflavanone****Cepaflava B**

Position	$\delta_C$	$\delta_H$	$\delta_C$	$\delta_H$
2	108.9		109.4	
3	79.8		80.0	
4	196.5		199.0	
4a	100.8		101.9	
5	165.5		164.4	
6	97.5	5.99 (1H, d, $J = 2.5$ Hz)	97.8	5.95 (1H, brs)
7	168.3		168.3	
8	96.7	6.01 (1H, d, $J = 2.5$ Hz)	97.1	6.00 (1H, brs)
8a	160.1		158.8	
1'	126.6		126.3	
2'	131.0	7.43 (1H, d, $J = 8.5$ Hz)	117.1	7.11 (1H, s)
3'	115.5	6.84 (1H, d, $J = 8.5$ Hz)	145.8	
4'	159.5		147.5	
5'	115.5	6.84 (1H, d, $J = 8.5$ Hz)	115.6	6.80 (1H, d, $J = 7.8$ Hz)
6'	131.0	7.43 (1H, d, $J = 8.5$ Hz)	121.6	7.02 (1H, d, $J = 7.8$ Hz)
1''	60.3	3.37 (1H, dq, overlap) 3.52 (1H, dq, $J = 7.5, 2.5$ Hz)	51.4	3.88 (3H, s)
2''	15.1	1.04 (3H, t, $J = 7.5$ Hz)	-	-
1'''	41.7	2.25 (1H, d, $J = 15.5$ Hz) 3.08 (1H, d, $J = 15.5$ Hz)	41.3	2.56 (1H, d, $J = 13.5$ Hz) 2.42 (1H, d, $J = 13.5$ Hz)
2'''	213.9		209.3	
3'''	31.8	2.28 (3H, s)	32.3	1.98 (3H, s)