Supplementary Materials:

Structure - antioxidant - antiproliferative activity relationships of natural C7 and C7-C8 hydroxylated flavones and flavanones

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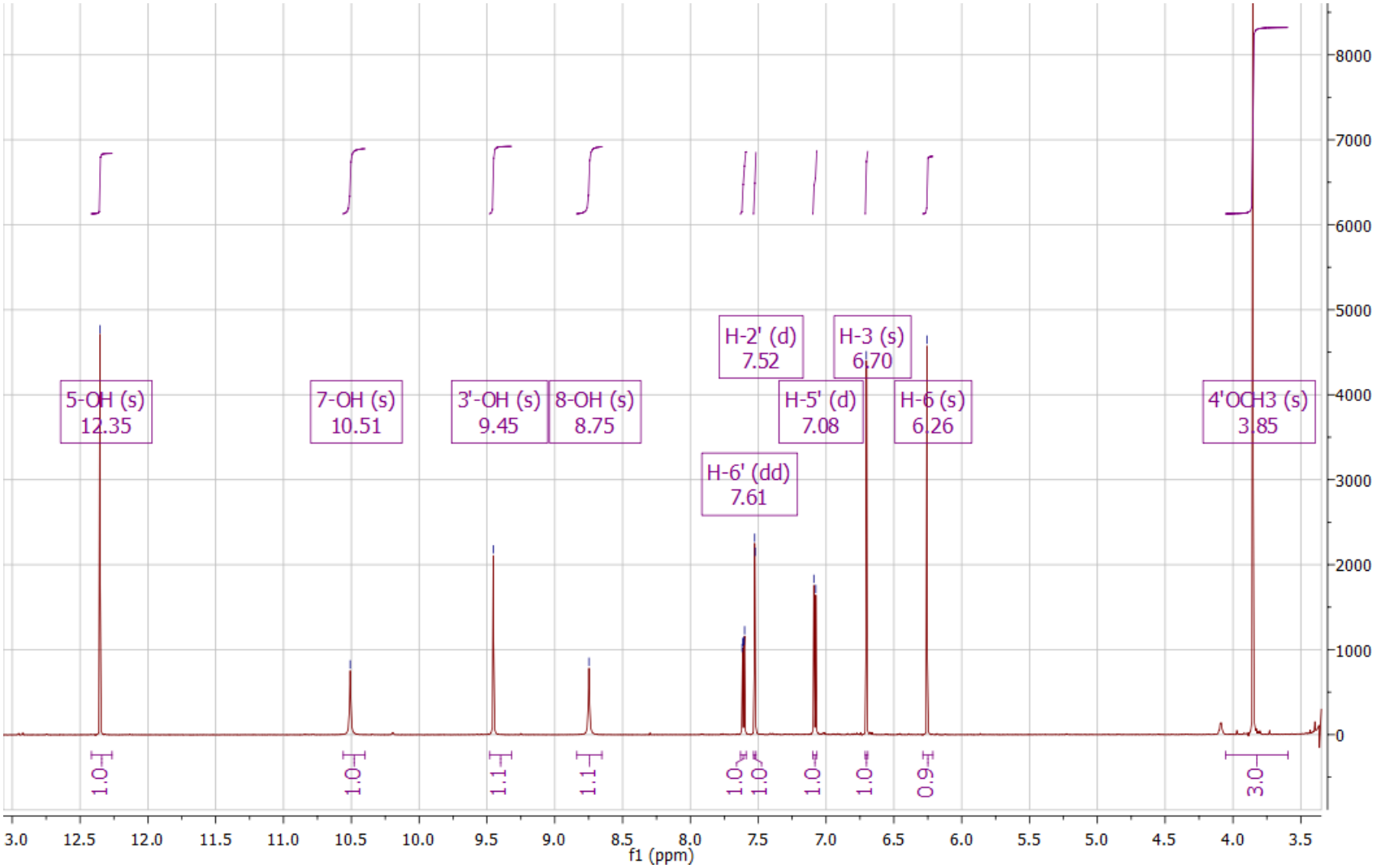

Figure S1. ^{1}H-NMR (600 MHz, DMSO-d_6) spectrum of 8-hydroxydiosmetin (**12**).

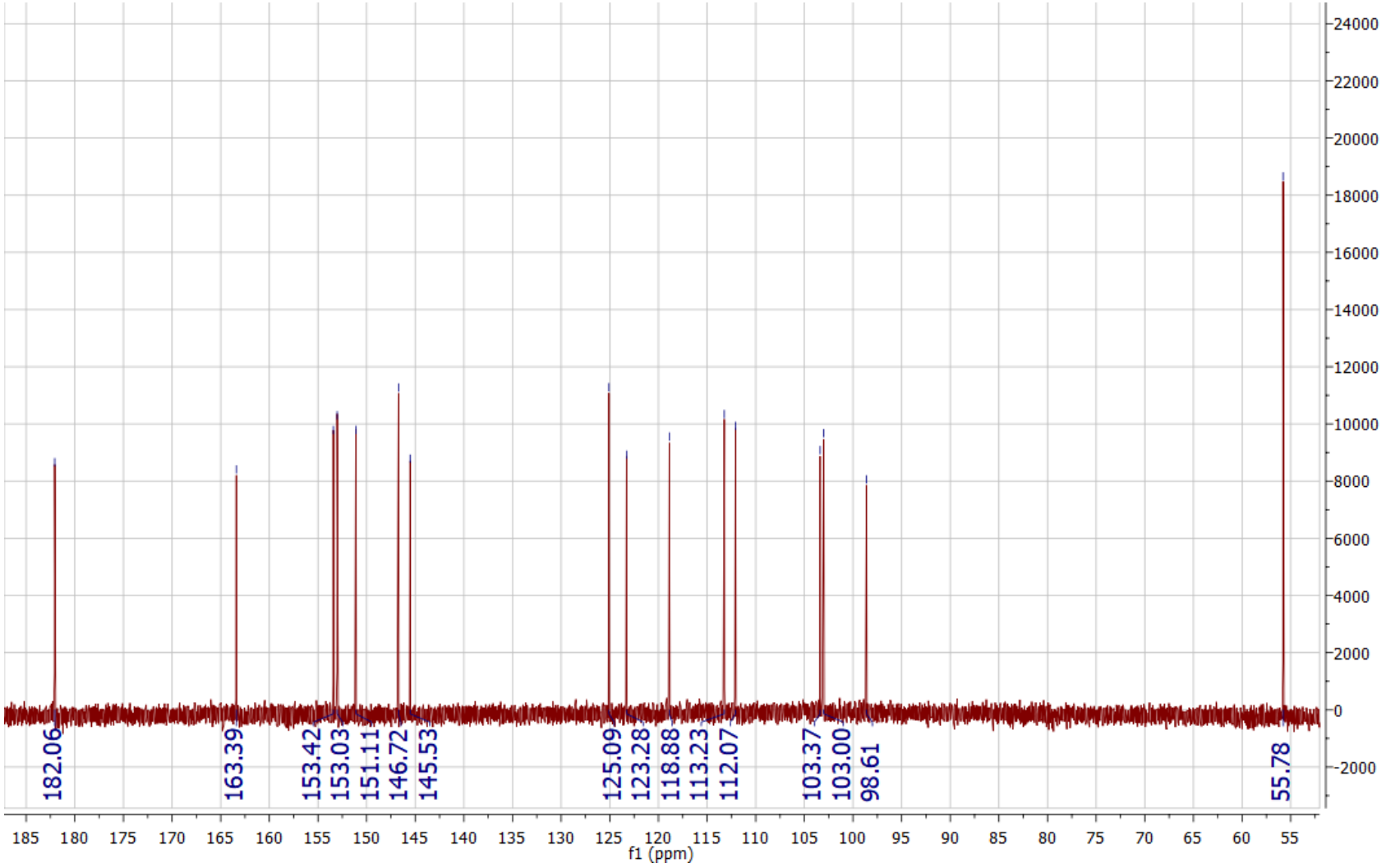

Figure S2. ^{13}C-NMR (150 MHz, DMSO-d_6) spectrum of 8-hydroxydiosmetin (**12**).

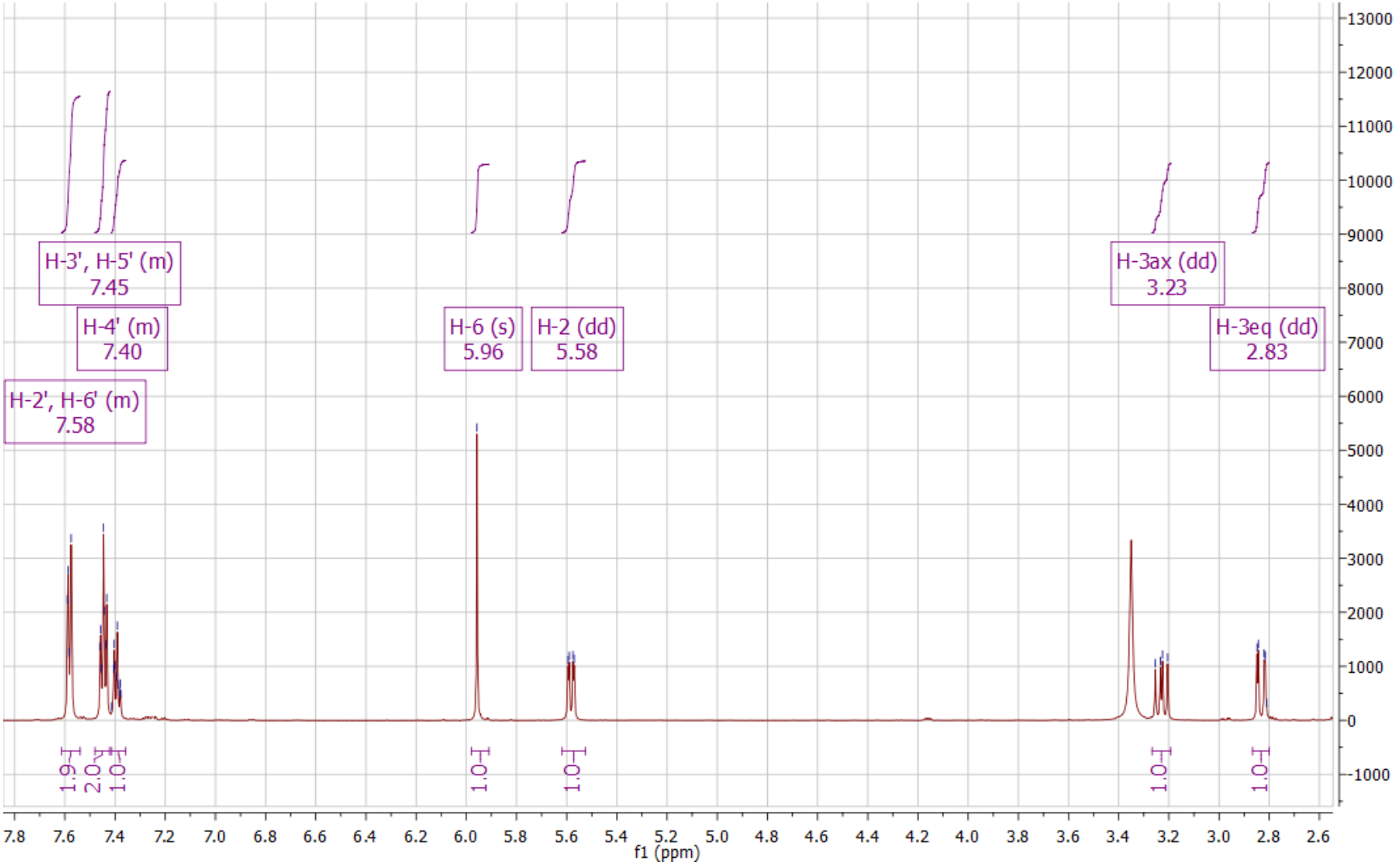

Figure S3. ^{1}H-NMR (600 MHz, DMSO-d_6) spectrum of 8-hydroxypinocembrin (**13**).

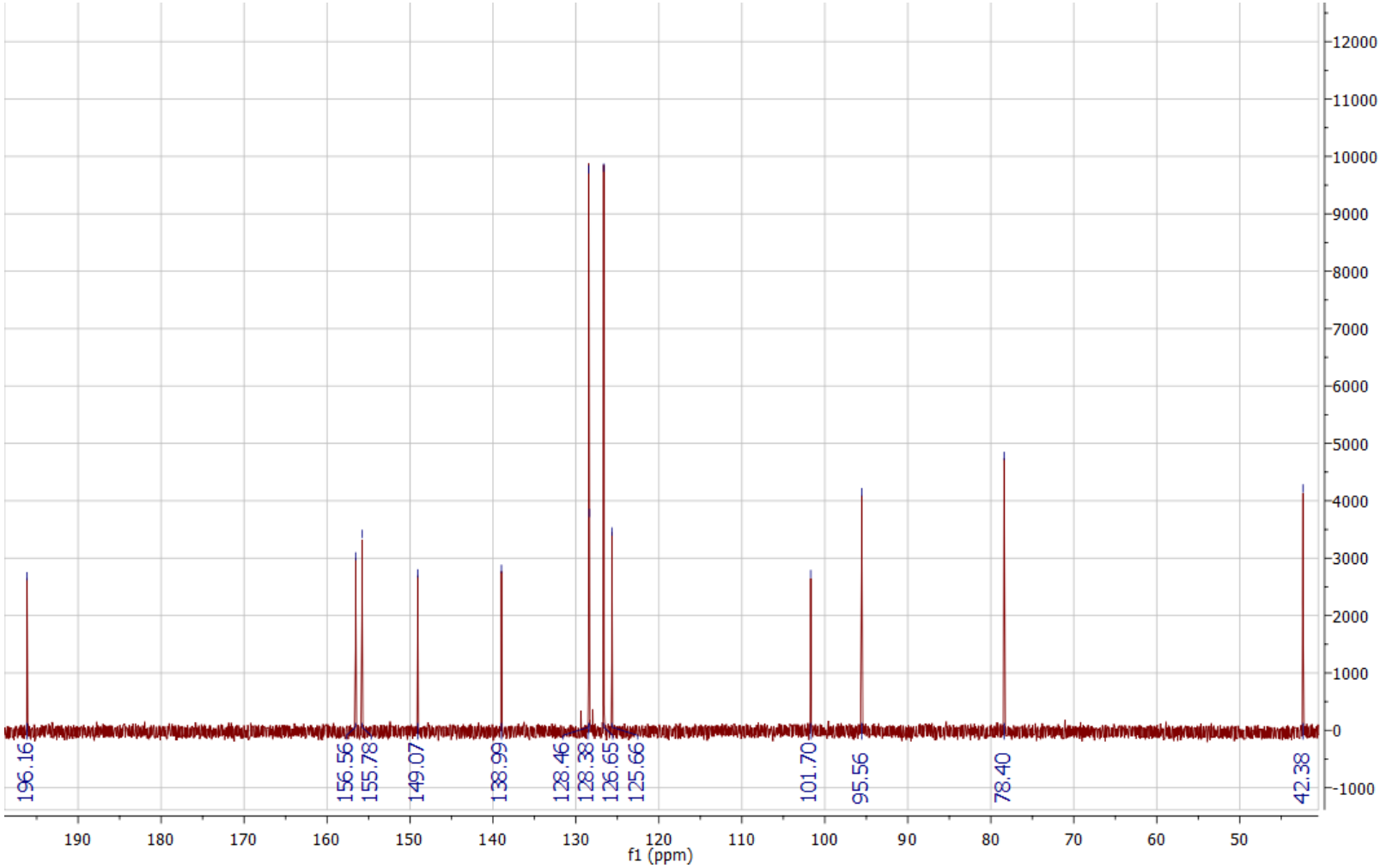

Figure S4. ^{13}C-NMR (150 MHz, DMSO-d_6) spectrum of 8-hydroxypinocembrin (**13**).

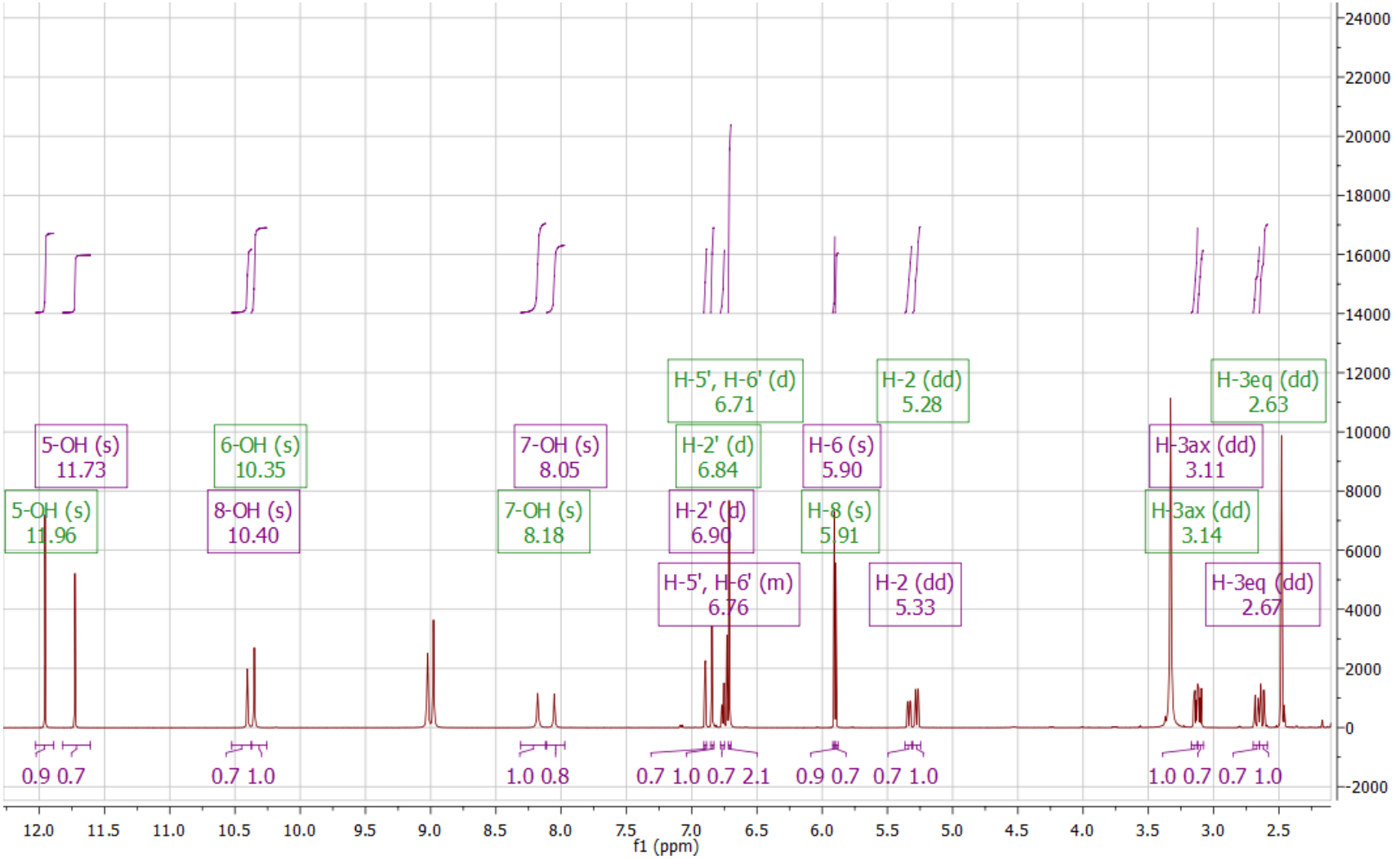

Figure S5. ^{1}H-NMR (600 MHz, DMSO-d_6) spectrum of 8-hydroxyeriodictyol (**16**) (violet) and 6-hydroxyeriodictyol (**17**) (green).

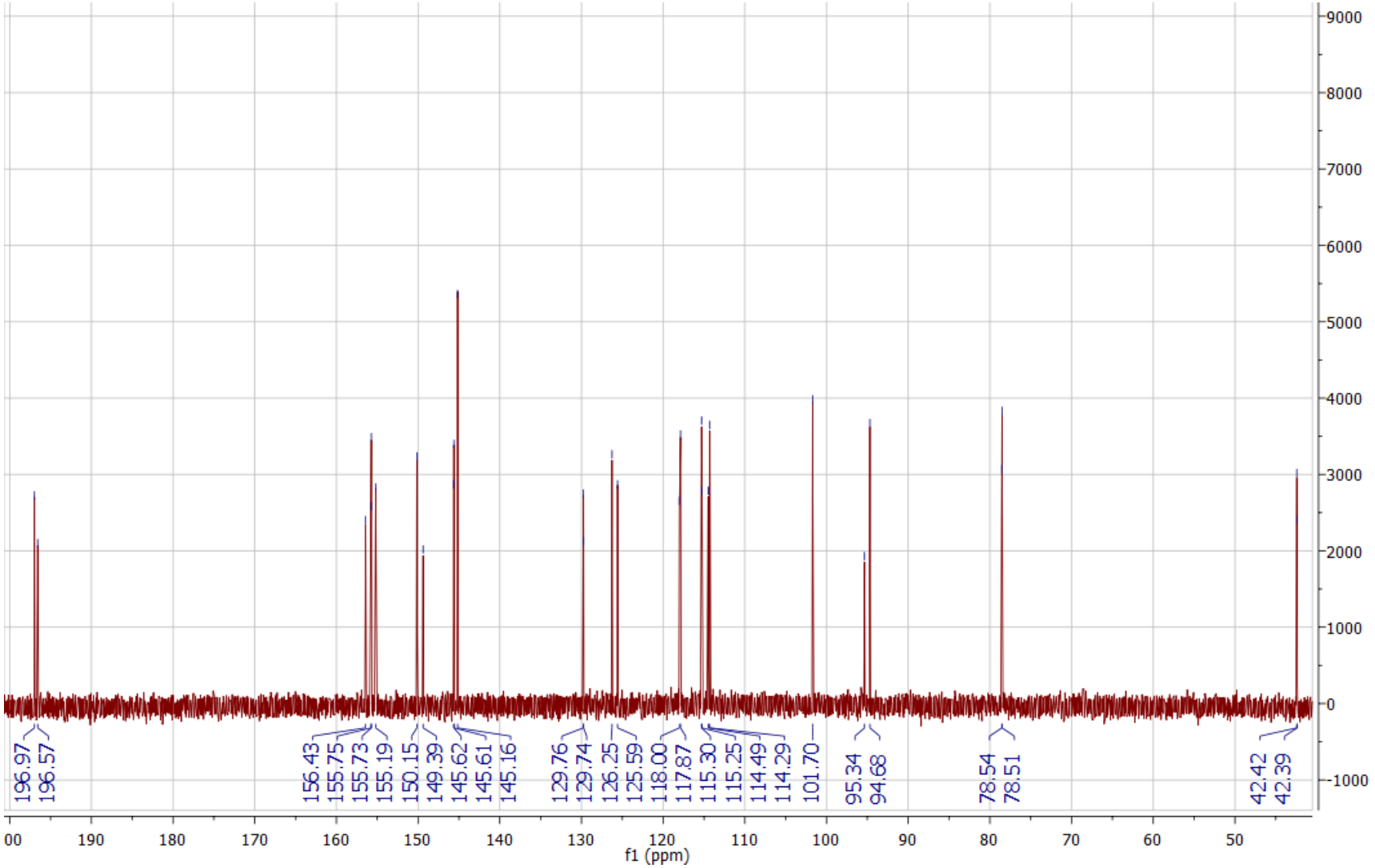

Figure S6. ^{13}C-NMR (150 MHz, DMSO-d_6) spectrum of 8-hydroxyeriodictyol and 6-hydroxyeriodictyol (**16**, **17**).

Table S1. Antioxidant activity of tested flavonoids.[a]

Compound	$TEAC_{ABTS}$	$TEAC_{DPPH}$	$TEAC_{FRAP}$
1	0.084 ± 0.011	0.004 ± 0.001	ND
2	0.250 ± 0.014	0.005 ± 0.002	ND
3	16.645 ± 0.897	10.50 ± 0.117	17.686 ± 0.327
4	1.712 ± 0.041	0.008 ± 0.001	ND
5	0.165 ± 0.009	0.005 ± 0.001	ND
6	0.347 ± 0.020	0.017 ± 0.002	ND
7	13.988 ± 0.251	9.582 ± 0.315	10.619 ± 0.319
8	13.254 ± 0.123	0.098 ± 0.002	4.025 ± 0.175
9	14.173 ± 1.591	8.038 ± 0.061	8.900 ± 0.582
10	10.184 ± 1.768	5.560 ± 0.121	7.824 ± 0.657
11	34.450 ± 1.016	13.208 ± 0.080	36.895 ± 0.432
12	13.405 ± 0.826	8.496 ± 0.040	11.340 ± 0.203
13	20.763 ± 0.939	9.164 ± 0.097	12.861 ± 0.313
14 and **15***	17.370 ± 0.989	9.098 ± 0.045	8.686 ± 0.093
16 and **17****	29.243 ± 1.362	14.649 ± 0.182	22.367 ± 1.801
18	31.187 ± 1.353	14.150 ± 0.229	25.260 ± 0.909

[a] Results are expressed as the equivalent of µmol trolox ($TEAC_{ABTS}$, $TEAC_{DPPH}$, $TEAC_{FRAP}$) per µmol of tested compound; SD – standard deviation; ND – not determined in the range of concentrations tested (50 µg/mL - 10 mg/mL); *ratio 19:1 (by HPLC); **ratio 1:1.3 (by HPLC).