

Article

Comparative Study on the Phenolic Fingerprint and Antioxidant Activity of Strawberry Tree (*Arbutus unedo* L.) Leaves and Fruits

Irena Brčić Karačonji ^{1,2,*,#}, Karlo Jurica ^{3,#}, Uroš Gašić ^{4,*}, Aleksandra Dramićanin ⁵, Živoslav Tešić ⁵ and Dušanka Milojković Opsenica ⁵

¹ Analytical Toxicology and Mineral Metabolism Unit, Institute for Medical Research and Occupational Health, Ksaverska cesta 2, 10000 Zagreb, Croatia; ibrcic@imi.hr (I.B.K.)

² Faculty of Health Studies, University of Rijeka, Viktora Cara Emina 5, 51000 Rijeka, Croatia

³ Special Security Operations Directorate, Ministry of the Interior, Ulica grada Vukovara 33, 10000 Zagreb, Croatia; juricakarlo@gmail.com (K.J.)

⁴ Institute for Biological Research "Siniša Stanković"–National Institute of Republic of Serbia, University of Belgrade, Bulevar despota Stefana 142, 11060 Belgrade, Serbia; uros.gasic@ibiss.bg.ac.rs (U.G.)

⁵ Faculty of Chemistry, University of Belgrade, Studentski trg 12-16, 11158 Belgrade, Serbia; akosovic@chem.bg.ac.rs (A.D.); ztesic@chem.bg.ac.rs (Ž.T.); dusankam@chem.bg.ac.rs (D.M.O.)

* Correspondence: ibrcic@imi.hr (I.B.K.); uros.gasic@ibiss.bg.ac.rs (U.G.)

Equally contributed (I.B.K. and K.J.)

Figure S1. Proposed fragmentation pathway of compound 42 (kaempferol 3-O-(6''-galloyl)hexoside).

Table S1. The list of phenolic compounds with regression equation parameters, correlation coefficients, limits of detection (LOD) and quantification (LOQ), obtained from the calibration curves.

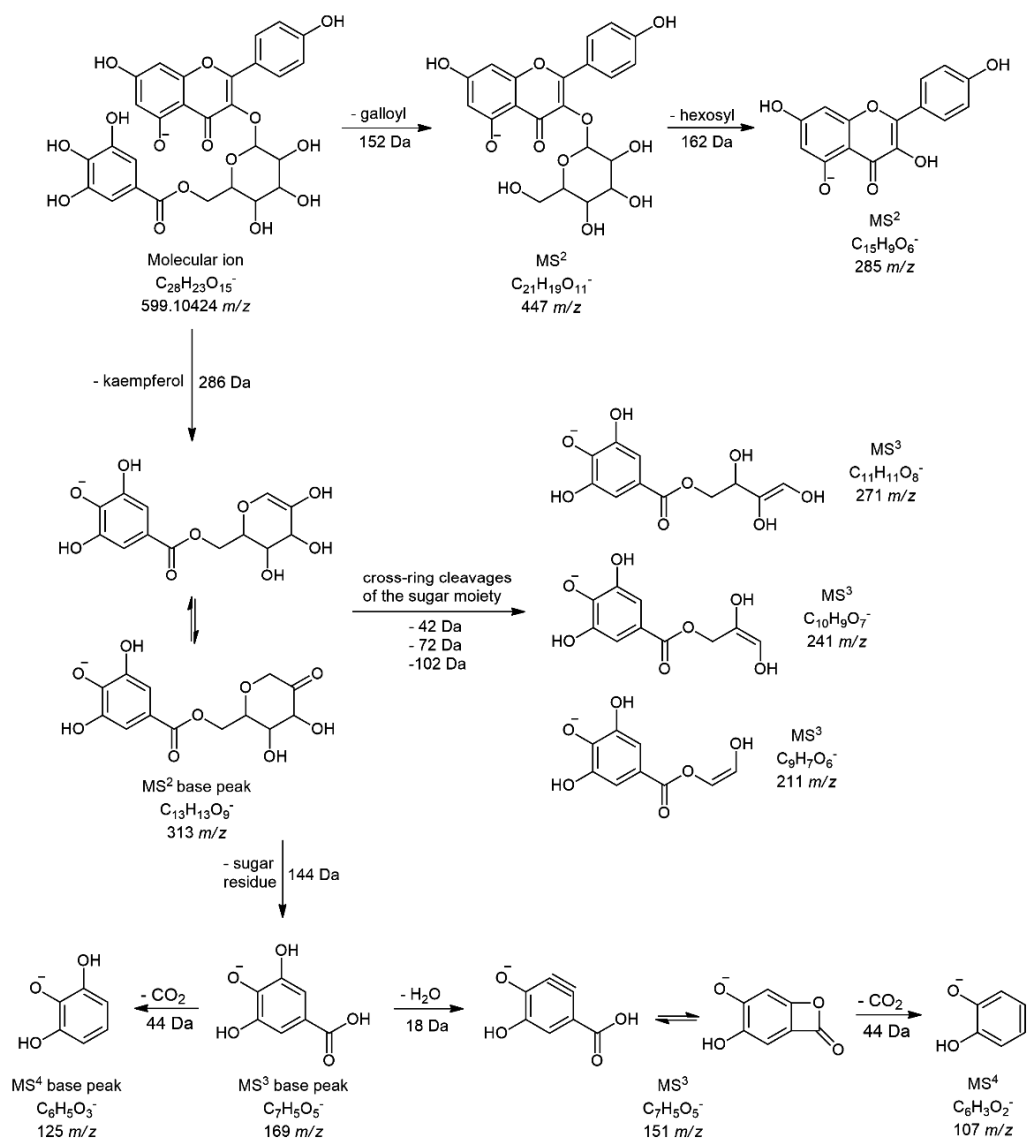


Figure S1

Table S1

No	Compounds	Regression equation (Y = A + BX)		R ²	LOD, mg/L	LOQ, mg/L
		(A ± SE) × 10 ⁵	(B ± SE) × 10 ⁵			
1	Gallic acid	- 3.22 ± 1.11	72.56 ± 2.64	0.9961	0.06	0.21
2	Gallocatechin	- 5.41 ± 0.81	32.05 ± 1.57	0.9952	0.06	0.21
3	Protocatechuic acid	2.66 ± 1.36	223.63 ± 2.63	0.9993	0.03	0.10
4	Aesculin	- 0.16 ± 1.27	156.22 ± 2.66	0.9985	0.04	0.15
5	Chlorogenic acid	- 4.01 ± 1.69	94.49 ± 2.88	0.9963	0.08	0.27
6	Catechin	- 1.17 ± 1.43	112.93 ± 3.71	0.9957	0.06	0.21
7	<i>p</i> -Hydroxybenzoic acid	8.35 ± 4.69	359.63 ± 10.99	0.9963	0.07	0.22
8	Caffeic acid	- 0.12 ± 1.20	223.33 ± 2.45	0.9994	0.03	0.10
9	Syringic acid	2.42 ± 3.34	236.76 ± 6.12	0.9967	0.08	0.25
10	Vanillic acid	2.84 ± 2.62	223.66 ± 5.28	0.9972	0.06	0.21
11	Rutin	- 3.43 ± 1.19	155.02 ± 2.10	0.9991	0.04	0.14
12	<i>p</i> -Hydroxyphenylacetic acid	5.66 ± 2.58	180.38 ± 4.77	0.9965	0.08	0.26
13	Hyperoside	- 3.96 ± 1.43	129.05 ± 2.76	0.9977	0.06	0.20
14	<i>p</i> -Coumaric acid	8.11 ± 4.37	313.16 ± 8.62	0.9962	0.08	0.25
15	Catechin gallate	- 9.18 ± 2.99	50.19 ± 3.11	0.9924	0.20	0.68
16	Ferulic acid	6.04 ± 5.97	251.15 ± 5.74	0.9974	0.13	0.43
17	Myricetin	0.11 ± 0.28	28.11 ± 0.73	0.9980	0.04	0.14
18	Quercetin	- 8.99 ± 7.31	183.45 ± 8.87	0.9927	0.18	0.60
19	Naringenin	- 8.30 ± 5.74	409.51 ± 9.70	0.9972	0.08	0.25
20	Kaempferol	- 6.94 ± 2.31	119.60 ± 7.62	0.9920	0.07	0.24

SE – standard error; LOD–limit of detection; LOQ–limit of quantification.