Supporting Information

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Comprehensive phytochemical profiling and cell-free in vitro bioactivity assessment of the endemic *Campanula kirikkaleensis* Dönmez & Güner

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Table S1: Chemical Composition of the Essential Oils from the Aerial Part and Root of *C. kirikkaleensis* Analyzed by GC-FID/MS

RI	Compound	Aerial Part (%)	Root (%)
1093	Hexanal	-	0.5
1225	(*Z*)-3-Hexenal	0.2	0.1
1244	Amyl furan	-	0.2
1282	Hexyl acetate	-	0.2
1360	Hexanol	0.1	0.4
1399	Nonanal	0.1	1.5
1405	(*Z*)-2-Hexen-1-ol	tr	-
1424	Hexyl butyrate	tr	-
1497	α-Copaene	-	tr
1507	(*E,E*)-2,4-Heptadienal	tr	tr
1516	2-Methoxy-3-sec-butylpyrazine[#]	-	0.1
1541	Benzaldehyde	tr	tr
1557	3,5,5-Trimethyl-3-butylcyclohexanone[#]	0.5	-
1598	2-Undecanone	-	0.1
1599	β-Elemene	0.1	-
1600	Hexadecane	0.1	0.1
1612	β-Caryophyllene	0.1	-
1638	Menthol	-	0.1
1655	(*E*)-2-Decenal	-	0.1
1659	Safranal	0.2	-
1664	Nonanol	-	0.1
1667	β-Barbatene	0.2	-
1682	(*Z*)-3-Nonen-1-ol	-	0.1
1700	Heptadecane	0.1	0.1
1723	3-Ethylbenzaldehyde[#]	0.4	2.9
1726	Germacrene D	1.0	tr
1730	α-Muurolene	0.1	-
1734	Car-2-en-4-one[#]	0.1	-
1738	1-Azabicyclo[3.3.0]octan-2-one	-	0.1
1761	δ-Cadinene	0.5	-
1778	Tridecene[#Δ]	-	3.4
1800	Octadecane	0.1	-
1827	(*E,E*)-2,4-Decadienal	-	1.6
1845	Tridecene[#Δ]	-	0.5
1860	2-Propylphenol methyl ether[#]	-	0.2
1868	(*E*)-Geranyl acetone	0.7	0.7
1870	Tridecene[#Δ]	-	0.7
1878	2-Methyl-3-hydroxy-2,4,4-trimethylpentyl propanoate[#]	0.3	0.4
1889	2,2,4-Trimethyl-3-carboxyisopropyl-isobutyl pentanoate[#]	0.6	0.5
1893	2-Methyl-2,2-dimethyl-1-(2-hydroxymethyl) propyl pentanoate[#]	0.2	0.3
1900	Nonadecane	0.1	-
1933	Tetradecanal	e	-
1958	(*E*)-β-Ionone	0.9	0.1
1970	*trans*-β-Ionone-5,6-epoxide (*isomer-1*) [#Δ]	0.8	-
1973	Dodecanol	0.1	0.1
1996	*trans*-β-Ionone-5,6-epoxide	0.3	-
2001	4,7-Dimethyl-2-benzofuran-1(3H)-one[#]	-	0.2
2000	Eicosane	0.3	-
2037	3-Ethyl-3,4-dihydro-2(1H)-quinoxalinone[#]	2.5	20.6
2041	Pentadecanal	1.5	0.8
2077	Tridecanol	0.1	-

RI	Compound	Aerial Part (%)	Root (%)
2100	Heneicosane	1.1	2.1
2131	Hexahydrofarnesyl acetone	6.6	3.4
2178	3,4-Dimethyl-5-pentylidene-2(5H)-furanone	1.1	-
2180	Tetradecanol	0.2	-
2187	T-Cadinol	0.2	-
2195	3,4-Dimethyl-5-pentyl-5H-furan-2-one	0.9	-
2196	4-Vinyl guaiacol	0.5	0.2
2200	Docosane	0.8	-
2208	Megastigmatrienone#	0.3	-
2226	Methyl hexadecanoate (=*Methyl palmitate*)	0.2	-
2255	α-Cadinol	0.4	-
2294	Myristicin	3.8	5.5
2298	Pentadecanol	0.2	-
2300	Tricosane	7.5	0.6
2305	9-Hydroxy-thymol methyl ether	-	2.3
2330	Methyl heptadecanoate (=*Methyl margarate*)	0.8	-
2367	(*Z*)-9-Tricosene	0.5	-
2384	Farnesyl acetone	0.9	-
2390	6,9,12-Octadecatrienoic acid methyl ester#	2.4	0.4
2400	Tetracosane	3.8	0.7
2452	9-Tetracosene	0.9	-
2500	Pentacosane	12.8	1.6
2538	9-Pentacosene	1.5	-
2590	Phytol	14.2	4.9
2600	Hexacosane	3.0	-
2648	Nonadecanol	-	0.7
2700	Heptacosane	9.2	1.4
2734	Tetradecanoic acid (=*Myristic acid*)	-	1.4
2800	Octacosane	-	0.7
2815	1-Eicosanol	1.4	1.1
2900	Nonacosane	1.1	-
2931	Hexadecanoic acid (=*Palmitic acid*)	4.5	22.5
3000	Triacontane	3.4	8.0
3100	Hentriacontane	-	3.2
		96.5	**97.5**

Note: RRI: Relative Retention Index was calculated for the polar column with *n*-alkanes (C_7-C_{40}); %: Calculated from the FID chromatogram; tr: Trace (< 0.1%); #: Identified based on spectral similarity with the Wiley-NIST library; Δ: Correct isomer not identified.

Table S2: GC-FID/MS Analysis of the Volatile Compounds in *C. kirikkaleensis* Aerial Part and Root using the MSD-SPME Technique.

RRI	Compound	Aerial Part %	Root %
1093	Hexanal	-	4.3
1129	Thuja-2,4(10)-diene	-	0.1
1194	Heptanal	-	0.3
1225	(*Z*)-3-Hexenal	1.0	-
1232	(*E*)-2-Hexenal	-	1.2
1244	Amyl furan	-	0.3
1260	Pentanol	-	0.2
1282	Hexyl acetate	-	1.0

RRI	Compound	Aerial Part %	Root %
1296	Octanal	-	0.3
1304	1-Octen-3-one	-	0.1
1327	(*Z*)-3-Hexenyl acetate	-	0.1
1332	4-Nonanone	-	0.1
1335	(*E*)-2-Heptenal	-	0.6
1348	6-Methyl-5-hepten-2-one	-	0.2
1360	Hexanol	0.2	10.7
1391	(*Z*)-3-Hexen-1-ol	-	0.5
1399	Nonanal	0.9	3.5
1412	(*E*)-2-Hexen-1-ol	-	1.0
1416	3-Octen-2-one	-	1.0
1441	(*E*)-2-Octenal	0.1	2.5
1452	1-Octen-3-ol	-	0.9
1479	(*E,Z*)-2,4-Heptadienal	0.6	2.0
1481	3,4-Dihydro-2H-1,4-benzoxazine	0.2	-
1493	Tetramethyl pyrazine[#]	0.8	-
1496	3-Nonanol	0.2	-
1507	(*E,E*)-2,4-Heptadienal	0.7	1.9
1516	2-Methoxy-3-sec-butylpyrazine[#]	-	0.6
1523	(*E,Z*)-3,5-Octadien-2-one	-	4.7
1532	Camphor	1.2	-
1541	Benzaldehyde	0.3	0.6
1548	(*E*)-2-Nonenal	0.2	0.7
1553	Linalool	1.7	0.4
1562	Octanol	0.2	0.6
1573	(*E,E*)-3,5-Octadien-2-one	0.5	1.0
1600	Hexadecane	0.9	-
1604	2-Undecanone	0.4	-
1606	6-Methyl-3,5-heptadien-2-one	0.4	0.2
1617	Hexyl hexanoate	-	0.2
1638	β-Cyclocitral	0.9	-
1642	*neoiso*-Menthol	0.6	0.4
1655	(*E*)-2-Decenal	0.6	0.7
1660	Benzeneacetaldehyde	1.8	-
1664	Nonanol	1.5	0.4
1671	Acetophenone	-	0.2
1681	2-Butyl-2-octenal	-	0.8
1696	*trans*-Chrysanthenol	-	0.9
1704	10-Methyl-2-undecanone	0.7	-
1706	α-Terpineol	0.2	-
1714	2-Dodecanone	0.2	-
1715	(*E,E*)-2,4-Nonadienal	-	0.5
1720	Dodecanal	1.0	-
1723	3-Ethylbenzaldehyde [#]	0.4	2.6
1730	Veratrole	-	0.1
1751	Carvone	0.5	-
1763	Naphthalene	-	0.5
1764	(*E*)-2-Undecenal	0.6	-
1766	Decanol	0.8	-
1778	Tridecene [#Δ]	1.3	-
1779	(*E,Z*)-2,4-Decadienal	-	7.1
1800	Octadecane	0.4	-
1815	2-Tridecanone	0.2	-
1827	(*E,E*)-2,4-Decadienal	1.5	10.1
1865	Hexanoic acid	-	0.7

RRI	Compound	Aerial Part %	Root %
1868	(*E*)-Geranyl acetone	4.8	1.6
1871	Undecanol	0.3	-
1878	2-Methyl-3-hydroxy-2,4,4-trimethylpentyl propanoate [#]	1.8	-
1889	2,2,4-Trimethyl-3-carboxyisopropyl-isobutyl pentanoate [#]	2.6	0.2
1893	2-Methyl-2,2-dimethyl-1-(2-hydroxy methyl) propyl pentanoate [#]	1.5	-
1933	Tetradecanal	0.6	-
1958	(*E*)-β-Ionone	5.5	-
1973	Dodecanol	1.8	0.4
1980	Benzothiazole[#]	-	0.2
1986	5,7-Dimethyl-2-benzofuran-1(3H)-one[#]	-	0.9
1996	*trans*-β-Ionone-5,6-epoxide	1.4	-
2025	1-(4-tert-Butylphenyl) propan-2-one [#]	0.6	-
2037	3-Ethyl-3,4-dihydro-2(1H)-quinoxalinone [#]	9.8	14.5
2053	Anisaldehyde	1.4	0.2
2084	Octanoic acid	-	0.2
2131	Hexahydrofarnesyl acetone	4.7	0.3
2192	Nonanoic acid	0.9	-
2178	3,4-Dimethyl-5-pentylidene-2(5H)-furanone	5.2	-
2196	4-Vinyl guaiacol	0.9	-
2239	Carvacrol	4.6	0.6
2264	Hexyl salicylate	0.7	-
2332	Methylpropyl salicylate [#Δ]	1.2	-
2343	2-Ethylhexyl salicylate	-	7.0
2371	Dihydroactinidiolide #	1.8	-
2382	2-Methoxy-α-5-dimethylbenzeneacetaldehyde[#]	-	0.7
2500	Pentacosane	1.5	-
2629	*MW:276, BP:121*	5.5	-
2670	Tetradecanoic acid	3.8	-
2935	Hexadecanoic acid (=*Palmitic acid*)	7.8	1.1
		94.4	**94.7**

Note: RRI: Relative Retention Index was calculated for the polar column with *n*-alkanes (C_7-C_{40}); %: Calculated from the FID chromatogram; #: Identified based on spectral similarity with the Wiley-NIST library; Δ: Correct isomer not identified; MW: Molecular weight; BP: Base peak

Table S3: GC-FID/MS Analysis of TMS Derivatives of *C. kirikkaleensis* Aerial Part Different Extracts

RRI	Compound*	HE %	CE %	ME %	HCE %	HCME %
1281	Glycerol - 3 TMS	-	-	0.7	-	-
1506	Malic acid - 3TMS	0.6	-	-	-	-
1582	1,4-Dideoxy-1,4-imino-D-arabinitol - 3TMS[#]	-	-	0.5	-	-
1622	3-Hydroxy-3-methylglutaric acid - 3TMS	-	-	0.8	-	-
1797	Tagatofuranose - 5TMS (*isomer-2*)	-	-	0.8	-	-
1843	Neophytadiene	0.7	1.2	2.8	2.2	-
1843	α-Fructofuranose - 5TMS	-	-	-	-	8.7
1849	D-Psicofuranose - 5TMS	-	-		-	1.2
1852	β-D-Fructofuranose, pentakis (trimethylsilyl) ether (*isomer 2*)	-	-	2.9	-	-
1855	Tatradecanoic acid (=*Myristic acid*) - TMS	0.6	-	2.5	1.2	-
1885	Heptadecanal	-	-	-	0.7	-
1898	Quinic acid - 5TMS	-	-	4.9	-	2.2
1907	Crotonic acid	-	-	-	0.9	-

RRI	Compound*	HE %	CE %	ME %	HCE %	HCME %
1929	α-Glucopyranose - 5TMS	-	-	2.6	-	2.6
1952	Pentadecanoic acid	-	-	-	0.7	-
1955	*MW: 612 BP:305*	-	-	2.6	-	-
2023	β-Glucopyranose - 5TMS	-	-	2.1	0.5	3.4
2049	Hexadecanoic acid (=*Palmitic acid*) - TMS	4.8	4.0	-	9.0	-
2076	*Unidentified*	-	-	-	-	3.5
2083	*MW: 612 Inositol derivative*	-	-	4.4	-	
2127	*myo*-Inositol - 6TMS	-	-	3.0	-	11.7
2161	*meso*-Inositol - 6TMS[#]	-	-	-	-	2.8
2168	*MW: 612 Inositol derivative*	-	-	1.3	-	-
2179	Phytol - TMS	0.6	-	-	-	-
2182	*MW: 612 Inositol derivative*	-	-	0.8	-	-
2186	Phytol	-	-	-	1.4	-
2189	*Unidentified*	-	-	0.9	-	-
2217	4-Methylvaleric acid	-	-	-	2.3	-
2217	(*Z,Z*)-9,12-Octadecadienoic acid (=*Linoleic acid*) - TMS	4.0	2.4	-	-	-
2221	(*Z*)-9-Octadecenoic acid (=*Oleic acid*) - TMS	5.6	3.8	-	-	-
2225	α-Linolenic acid	-	-	-	4.3	-
2250	Octadecanoic acid (=*Stearic acid*) - TMS	0.5	-	-	1.0	-
2531	2-Methyl-6-nitroindolizine	-	-	-	2.6	-
2591	*Sucrose derivative - 7TMS*	-	-	-	-	0.8
2608	Monopalmitin - 2TMS	-	-	0.9	-	-
2646	Docosanoic acid - TMS	0.7	-	-	-	-
2687	*Unidentified*	-	-	-	-	1.6
2704	*Sucrose derivative - 8TMS*	-	-	-	-	5.0
2713	Sucrose - 8TMS	-	-	20.8	1.0	46.3
2761	*Unidentified*	-	-	-	-	2.1
2826	*Unidentified*	-	-	-	-	2.1
2844	Tetracosanoic acid - TMS	2.7	2.4	-	-	-
2893	*Unidentified*	-	-	1.2	-	-
2900	Nonacosane	0.5	-	-	-	-
2944	Pentacosanoic acid - TMS	0.6	-	-	-	-
2953	Hexacosanol - TMS	1.0	1.1	-	-	-
3011	*Unidentified*	0.8	-	-	-	-
3014	γ-Tocopherol - TMS	0.5	-	-	-	-
3023	*MW: 740*	0.6	-	1.1	-	-
3042	Benzene-1,2,4-triol	-	-	-	9.4	-
3042	Hexacosanoic acid - TMS	8.8	8.9	-	-	-
3071	*BP: 204*	-	-	0.7	-	-
3085	*MP: 361, 412*	-	-	2.0	-	-
3095	*BP: 203 MW:740*	-	-	2.9	-	-
3100	Hentriacontane	0.5	-	-	-	-
3104	*BP: 361*	-	-	0.7	-	-
3144	*MP: 361, 217, 382*	-	-	1.5	-	-
3144	Heptacosanoic acid - TMS	0.1	1.4	-	-	-
3149	1-Octacosanol - TMS	2.7	5.0	-	4.2	-
3159	α-Tocopherol (=*Vitamin E*) - TMS	1.8	1.6	-	-	-
3184	*crypto*-Chlorogenic acid - 6TMS	-	-	3.0	-	-
3214	*Unidentified*	-	-	-	-	1.0
3219	*BP: 575*	-	-	0.7	-	-
3224	*BP: 249*	-	-	1.2	-	-
3226	*BP: 204*	-	-	0.7	-	-
3239	*BP: 202 MP:289*	-	-	1.9	-	-
3244	Octacosanoic acid - TMS	16.8	20.0	-	18.0	-

RRI	Compound*	HE %	CE %	ME %	HCE %	HCME %
3264	5-Caffeoylquinic acid - 6TMS	-	-	0.5	-	-
3275	Campesterol - TMS	1.8	1.5	-	-	-
3305	Stigmasterol - TMS	2.2	2.1	-	-	-
3347	Triacontanol - TMS	0.9	1.5	-	-	-
3368	β-Sitosterol - TMS	12.0	11.3	0.9	4.9	-
3378	5α-Stigmasta-3β-ol - TMS	1.7	2.4	-	-	-
3383	*MW: 414*	-	1.1	-	-	-
3426	*MW: 486*	-	3.1	-	-	-
3431	Cycloartenol - TMS	3.2	-	-	-	-
3442	Triacontanoic acid - TMS	1.0	1.9	-	-	-
3444	*Unidentified*	0.6	-	-	-	-
3480	*MW: 512*	3.3	2.6	-	-	-
3491	*Unidentified*	1.0	-	-	-	-
3507	*Unidentified*	-	-	-	-	3.6
3507	Raffinose - 11TMS	-	-	4.2	-	-
3535	*MW: 498*	1.2	1.1	-	-	-
3563	*Unidentified*	1.0	1.4	-	-	-
3630	*Unidentified*	0.7	-	-	-	-
4093	Docosyl ferulate	-	-	-	8.2	-
4093	Unidentified	0.7	-	-	-	-
4218	*MW: 408*	-	3.1	-	-	-
4646	*MW: 1050, BP: 217*	-	-	3.2	-	-
5665	*α-D-Mannopyranoside methyl-butyl derivative*	-	-	-	26.3	-
	Total	**86.8**	**84.9**	**81.7**	**98.8**	**98.6**

Note: RRI: Relative Retention Index was calculated for the polar column with *n*-alkanes (C_7-C_{40}); *: ≥ 0.5%; #: Identified based on spectral similarity with the Wiley-NIST library; HE: Hexane extract; CE: Chloroform extract; ME: Methanol extract; HCE: Hexane+Chloroform extract; Hexane+Chloroform+Methanol extract; MW: Molecular weight; BP: Base peak

Table S4: GC-FID/MS Analysis of TMS Derivatives of *C. kirikkaleensis* Different Root Extracts

RRI	Compound*	HE %	CE %	ME %	HCE %	HCME %
1281	Glycerol - 3TMS	-	-	0.1	-	-
1319	5-Hydroxymethyl furfural - TMS	-	-	-	-	9.7
1325	*Unidentified*	-	-	-	-	2.5
1506	Malic acid - 3TMS	-	-	0.5	-	-
1512	*Unidentified*	-	-	-	-	3.3
1518	*Unidentified*	-	-	-	-	5.5
1520	Trimethyl ((4-methylcyclohexyl) methoxy) silane#	-	-	-	-	7.9
1532	Pyroglutamic acid - 2TMS	-	-	0.1	-	-
1565	*Unidentified*	-	-	-	-	3.4
1582	1,4-Dideoxy-1,4-imino-D-arabinitol - 3TMS#	-	-	0.1	-	-
1585	*Unidentified*	-	-	-	-	11.0
1585	Threonic acid - 4TMS (*isomer-2*)	-	-	0.1	-	-
1602	*Unidentified*	-	-	-	-	4.7
1622	3-Hydroxy-3-methylglutaric acid - 3TMS	-	-	0.1	-	-
1628	*Unidentified*	-	-	-	-	1.3
1636	*Unidentified*	-	-	-	-	1.0
1738	Xylopyranose - 4TMS	-	-	0.1	-	-
1753	Xylitol - 5TMS	-	-	0.1	-	-
1767	*Unidentified*	-	-	-	-	1.5
1782	*Unidentified*	-	-	-	-	1.2

RRI	Compound*	HE %	CE %	ME %	HCE %	HCME %
1794	*Unidentified*	-	-	-	-	2.8
1797	Tagatofuranose (isomer-2) - 5TMS	-	-	0.1	-	-
1800	Ribonic acid - 5TMS	-	-	0.1	-	-
1803	*Unidentified*	-	-	-	-	1.8
1809	α-Methylfuranoside - 4TMS	-	-	0.1	-	-
1824	*Unidentified*	-	-	-	-	3.6
1843	*Unidentified*	-	-	-	-	2.3
1843	α-Fructofuranose - 5TMS	-	-	1.8	-	-
1852	β-Fructofuranose - 5TMS	-	-	2.0	-	-
1858	Fructopyranose - 5TMS	-	-	0.5	-	-
1858	*Unidentified*	-	-	-	-	2.6
1901	Quinic acid - 5TMS	-	-	2.7	-	-
1907	Lidocaine[#]	-	-	-	-	3.0
1917	Gluconolactone - 4TMS	-	-	0.2	-	-
1927	Sorbose - 5TMS[#]	-	-	1.1	-	-
1929	*Unidentified*	-	-	-	-	1.1
1929	α-Glucopyranose - 5TMS	-	-	0.9	-	-
1937	Unidentified	-	-	-	-	0.9
1958	*MW:612, BP:305*	-	-	1.4	-	-
1971	Mannitol - 6TMS	-	-	0.2	-	-
1980	Sorbitol - 6TMS	-	-	0.2	-	-
1987	*Unidentified*	-	-	-	-	0.5
2023	β-Glucopyranose - 5TMS	-	-	1.8	-	-
2046	Gluconic acid - 6TMS	-	-	0.1	-	-
2049	Hexadecanoic acid (=*Palmitic acid*) - TMS	1.8	-	0.1	2.1	0.6
2083	*Unidentified*	-	-	1.8	-	-
2127	*myo*-Inositol - 6TMS	-	-	3.1	-	-
2151	(*E*)-Caffeic acid - 6TMS	-	-	0.1	-	-
2151	Heptadecanoic acid - TMS	-	1.5	-	-	-
2165	*Unidentified*	-	-	0.5	-	-
2182	*Unidentified*	-	-	0.7	-	-
2207	*Unidentified*	-	-	0.3	-	-
2217	(Z,Z)-9,12-Octadecadienoic acid (=*Linoleic acid*) - TMS	3.4	1.5	0.1	0.6	-
2221	(Z)-9-Octadecenoic acid (=*Oleic acid*) -TMS	1.9	1.0	0.1	-	-
2250	Octadecanoic acid (=*Stearic acid*) - TMS	-	-	-	0.4	-
2364	2-O-Glycerol-α-D-galactopyranoside – 6TMS	-	-	0.1	-	-
2429	*Unidentified*	-	-	0.7	-	-
2519	*Unidentified*	-	-	-	-	2.3
2555	*Unidentified*	-	-	-	-	4.3
2555	Docosanol - TMS	-	-	-	0.4	-
2591	*Unidentified*	-	-	-	-	3.3
2646	Docosanoic acid (=*Behenic acid*)-TMS	1.0	1.2	-	1.6	-
2662	*Unidentified*	-	-	-	-	2.8
2692	*Unidentified*	-	-	0.2	-	5.0
2709	*Unidentified*	-	-	-	-	2.6
2713	Sucrose - 8TMS	-	-	52.9	-	-
2717	*Unidentified*	-	-	-	-	1.4
2730	*Unidentified*	-	-	-	-	2.6
2765	*Unidentified*	-	-	0.2	-	-
2769	*Unidentified*	-	-	-	-	2.9
2778	1-Monooleoylglycerol - 2TMS	-	-	0.1	-	-
2817	Trehalose - 8TMS	-	-	1.1	-	-
2844	Tetracosanoic acid (=*Lignoceric acid*) - TMS	-	-	-	2.9	-
2844	*neo*-Trehalose - 8TMS	-	-	0.1	-	-
2844	Tetracosanoic acid - TMS	1.8	2.9	-	-	-

RRI	Compound*	HE %	CE %	ME %	HCE %	HCME %
2853	Leucrose – 6TMS	-	-	0.1	-	-
2944	Pentacosanoic acid (=*Pentadecylic acid*) - TMS	-	-	-	0.5	-
2953	Hexacosan-1-ol (=*Ceryl alcohol*) - TMS	2.5	2.6	-	1.4	-
2976	*Unidentified*	-	-	0.1	-	-
2986	*Unidentified*	-	-	0.2	-	
3042	Hexacosanoic acid (=*Cerotic acid*) - TMS	6.1	11.7	-	14.9	-
3080	*Unidentified*	-	-	0.1	-	-
3085	*Unidentified*	-	-	0.2	-	-
3090	*Unidentified*	-	-	0.1	-	-
3095	*Unidentified*	-	-	0.3	-	-
3104	*Unidentified*	-	-	0.2	-	-
3144	*Unidentified*	-	-	0.2	-	-
3144	Heptacosanoic acid (=*Carboceric acid*) - TMS	-	1.1	-	1.3	-
3149	Octacosan-1-ol (=*Stearyl alcohol*) - TMS	12.6	15.6	-	8.4	-
3159	α-Tocopherol (=*Vitamin E*) - TMS	1.5	-	-	-	-
3184	Chlorogenic acid - 6TMS	-	-	1.6	-	-
3224	*Unidentified*	-	-	0.4	-	-
3244	Octacosanoic acid (=*Montanic acid*) - TMS	6.9	18.9	-	16.6	-
3249	*Unidentified*	-	-	0.4	-	-
3275	Campesterol - TMS	2.9	1.5	-	-	-
3305	Stigmasterol - TMS	2.7	1.8	-	-	-
3347	1-Triacontanol (=*Triacontoxy*) - TMS	2.3	2.6	-	-	-
3347	1,1-Dimethylethyl-dimethyl-triacontyloxy silane#	-	-	-	2.0	-
3368	β-Sitosterol - TMS	10.5	6.8	0.1	2.1	-
3378	5α-Stigmasta-3β-ol (=*Stigmastanol*) -TMS	2.2	1.7	-	-	-
3383	*MW:414*	1.4	-	-	-	-
3389	*MW:424*	1.1	-	-	-	-
3431	Cycloartenol - TMS	2.6	1.9	-	-	-
3442	*MW:524*	2.7	1.6	-	-	-
3442	*Unidentified*	-	-	-	1.4	-
3480	*MA:512*	5.2	2.1	-	-	-
3513	1-Kestose - 11TMS	-	-	9.1	-	-
3591	Palatinose methyloxime - 7TMS#	-	-	0.1	-	-
3602	Oleanolic acid - 2TMS	-	-	0.1	-	-
3778	*Unidentified*	-	-	-	1.4	-
3935	Docosyl-*p*-coumarate - TMS	1.6	1.5	-	6.3	-
4170	*Unidentified*	-	-	-	4.1	-
4489	*Unidentified*	-	-	-	3.1	-
4675	*MW:1050, BP:217*	-	-	7.6	-	-
4939	*Unidentified*	-	-	-	4.9	-
4993	*Unidentified*	-	-	-	4.4	-
5070	*Unidentified*	-	-	-	2.9	-
5251	*Unidentified*	-	-	-	16.3	-
	Total	**74.7**	**79.5**	**97.5**	**100.0**	**99.4**

Note: RRI: Relative Retention Index was calculated for the polar column with *n*-alkanes (C_7-C_{40}); *: ≥ 0.5%**;** #: Identified based on spectral similarity with the Wiley-NIST library; HE: Hexane extract; CE: Chloroform extract; ME: Methanol extract; HCE: Hexane+Chloroform extract; Hexane+Chloroform+Methanol extract; MW: Molecular weight; BP: Base peak

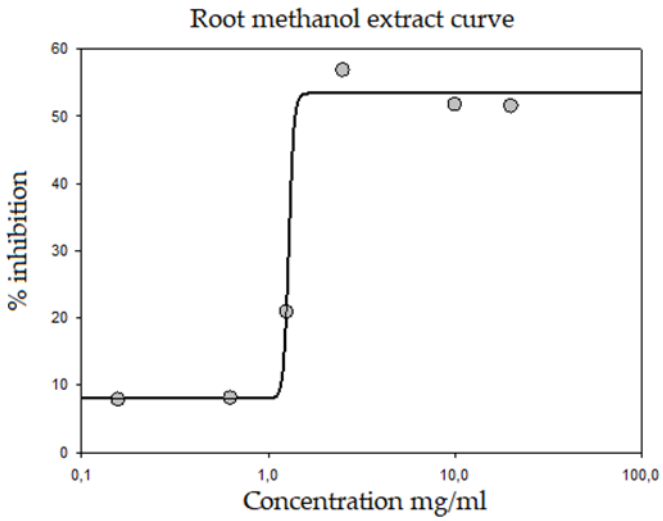

Figure S1: Root Methanol Extract Curve

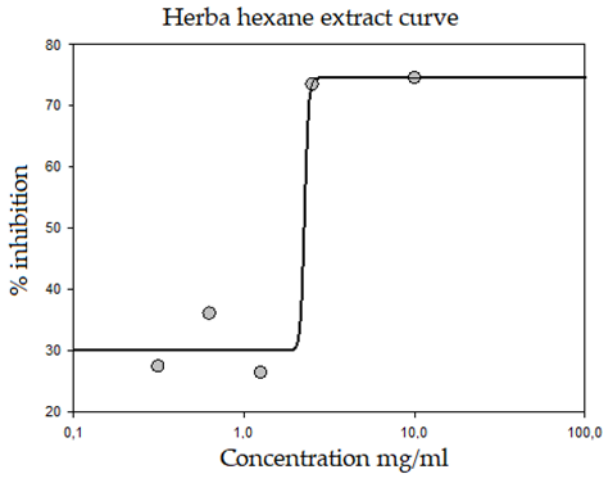

Figure S2: Herba Hexane Extract Curve

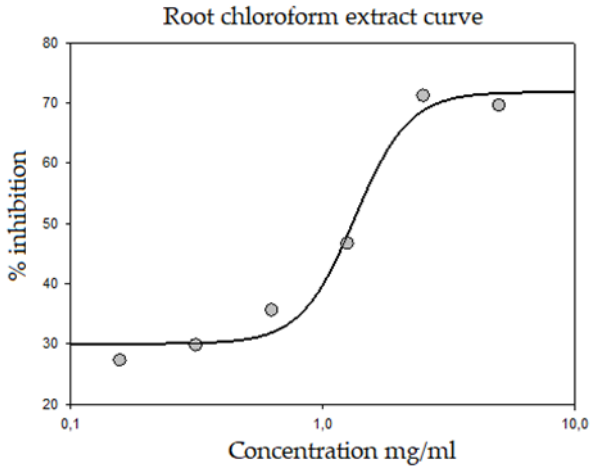

Figure S3: Root Chloroform Extract Curve

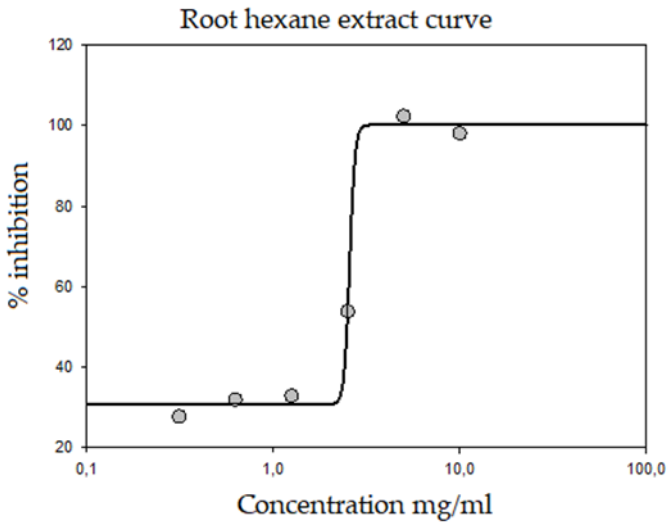

Figure S4: Root Hexane Extract Curve