Date : April 06, 2021

Certificate of Analysis – GC Profiling

Sample identification

Internal code : 21C19-PTH11
Customer identification : Spearmint ORGANIC - India - S40105205R
Type : Essential oil
Source : *Mentha spicata*
Customer : Plant Therapy

Analysis

Method: PC-MAT-014 ISO - Analysis of the composition of an essential oil or other volatile liquid by FAST GC-FID (in French); identifications validated by GC-MS.
Analyst : Seydou Ka, M. Sc.
Analysis date : March 31, 2021

Checked and approved by :

Sylvain Mercier, M. Sc., chimiste 2014-005

Notes: This report may not be published, including online, without the written consent from Laboratoire PhytoChemia. This report is digitally signed, it is only considered valid if the digital signature is intact. The results only describe the samples that were submitted to the assays.

Méthode certifiée ISO 17025:2017 Certified method

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PHYSICOCHEMICAL DATA

Physical aspect: Light yellow liquid
Refractive index: 1.4888 ± 0.0003 (20 °C; method PC-MAT-016)

CONCLUSION

No adulterant, contaminant or diluent has been detected using this method.

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Analysis summary – Consolidated contents

New readers of similar reports are encouraged to read table footnotes at least once.

Identification	%	Class
Isobutyral	tr	Aliphatic aldehyde
3-Buten-2-one	tr	Aliphatic ketone
Isobutanol	tr	Aliphatic alcohol
Isovaleral	0.01	Aliphatic aldehyde
2-Methylbutyral	0.01	Aliphatic aldehyde
2-Ethylfuran	tr	Furan
Isoamyl alcohol	0.01	Aliphatic alcohol
2-Methylbutanol	0.01	Aliphatic alcohol
Toluene	tr	Simple phenolic
Methyl 2-methylbutyrate	0.01	Aliphatic ester
Hexanal	tr	Aliphatic aldehyde
Ethyl 2-methylbutyrate	0.01	Aliphatic ester
Ethyl isovalerate	0.01	Aliphatic ester
(3*Z*)-Hexenol	0.01	Aliphatic alcohol
(2*E*)-Hexenol	tr	Aliphatic alcohol
Hexanol	tr	Aliphatic alcohol
trans-2,5-Diethyltetrahydrofuran	0.07	Furan
Hashishene	0.10	Monoterpene
α-Thujene	0.04	Monoterpene
α-Pinene	0.63	Monoterpene
Camphene	0.02	Monoterpene
3-Methylcyclohexanone	0.02	Aliphatic ketone
Thuja-2,4(10)-diene	0.01	Monoterpene
Benzaldehyde	0.01	Simple phenolic
β-Pinene	0.80	Monoterpene
Sabinene	0.48	Monoterpene
Octen-3-one	tr	Aliphatic ketone
Octen-3-ol	0.01	Aliphatic alcohol
Octan-3-one	0.04	Aliphatic ketone
Myrcene	1.71	Monoterpene
Octan-3-ol	0.37	Aliphatic alcohol
α-Phellandrene	0.01	Monoterpene
Pseudolimonene	0.03	Monoterpene
Octanal	0.06	Aliphatic aldehyde
Δ3-Carene	tr	Monoterpene
α-Terpinene	0.14	Monoterpene
para-Cymene	0.19	Monoterpene
1,8-Cineole	1.47*	Monoterpenic ether
β-Phellandrene	[1.47]*	Monoterpene
Limonene	17.45	Monoterpene
2-Ethylhexanol	0.01	Aliphatic alcohol
(*Z*)-β-Ocimene	0.07	Monoterpene
Unknown	0.09	Unknown
(*E*)-β-Ocimene	0.05	Monoterpene
γ-Terpinene	0.27	Monoterpene

cis-Sabinene hydrate	0.32	Monoterpenic alcohol
Octanol	0.06	Aliphatic alcohol
Terpinolene	0.11	Monoterpene
para-Cymenene	0.03	Monoterpene
trans-Sabinene hydrate	0.04	Monoterpenic alcohol
2-Methylbutyl isovalerate?	0.01	Aliphatic ester
Linalool	0.05	Monoterpenic alcohol
Isoamyl isovalerate	tr	Aliphatic ester
2-Methylbutyl 2-methylbutyrate	0.02	Aliphatic ester
Nonanal	0.02	Aliphatic aldehyde
trans-para-Mentha-2,8-dien-1-ol	0.09	Monoterpenic alcohol
cis-Limonene oxide	0.02	Monoterpenic ether
allo-Ocimene	0.01	Monoterpene
Octan-3-yl acetate	0.10	Aliphatic ester
trans-Pinocarveol	0.05	Monoterpenic alcohol
cis-para-Mentha-2,8-dien-1-ol	0.07	Monoterpenic alcohol
trans-Limonene oxide	0.04	Monoterpenic ether
Camphor	0.01	Monoterpenic ketone
cis-Verbenol	0.01	Monoterpenic alcohol
trans-para-Menth-2-en-1-ol	0.01	Monoterpenic alcohol
Isopulegol	0.03	Monoterpenic alcohol
Menthone	0.11	Monoterpenic ketone
Menthofuran	0.01	Monoterpenic ether
Isomenthone	0.04	Monoterpenic ketone
Borneol	0.02	Monoterpenic alcohol
neo-Menthol	0.16	Monoterpenic alcohol
Menthol	0.84	Monoterpenic alcohol
Terpinen-4-ol	0.79	Monoterpenic alcohol
Isomenthol	tr	Monoterpenic alcohol
para-Cymen-8-ol	tr	Monoterpenic alcohol
α-Terpineol	0.02	Monoterpenic alcohol
cis-Dihydrocarvone	1.59	Monoterpenic ketone
Myrtenal	0.01	Monoterpenic aldehyde
neo-Dihydrocarveol	0.65	Monoterpenic alcohol
Methylchavicol	0.01	Phenylpropanoid
trans-Dihydrocarvone	0.25	Monoterpenic ketone
Dihydrocarveol	0.22	Monoterpenic alcohol
trans-Piperitol	tr	Monoterpenic alcohol
Decanal	0.01	Aliphatic aldehyde
iso-Dihydrocarveol ?	0.04	Monoterpenic alcohol
trans-Carveol	0.51	Monoterpenic alcohol
(3*Z*)-Hexenyl 2-methylbutyrate	0.01	Aliphatic ester
Carvone	61.18	Monoterpenic ketone
Piperitone	0.44	Monoterpenic ketone
cis-Carvone oxide	0.01	Monoterpenic ketone
Isopiperitenone	0.06	Monoterpenic ketone
Unknown	0.02	Unknown
trans-Carvone oxide	0.08	Monoterpenic ketone
Decanol	0.03	Aliphatic alcohol
2-Ethylmenthone?	0.01	Aliphatic ketone
Dihydroedulan I	0.02	Terpenic ether
Menthyl acetate	0.10	Monoterpenic ester

Dihydroedulan II	0.01	Terpenic ether
Isomenthyl acetate	0.01	Monoterpenic alcohol
Bicycloelemene analog	tr	Sesquiterpene
Dihydrocarvyl acetate	0.23	Monoterpenic ester
Bicycloelemene	0.03	Sesquiterpene
trans-Carvyl acetate	0.02	Monoterpenic ester
α-Cubebene	0.01	Sesquiterpene
Menthofurolactone	0.03	Aliphatic alcohol
iso-Dihydrocarvyl acetate	0.03	Monoterpenic ester
cis-Carvyl acetate	0.21	Monoterpenic ester
α-Copaene	0.04	Sesquiterpene
β-Bourbonene	1.15	Sesquiterpene
1,5-diepi-β-Bourbonene	0.11	Sesquiterpene
β-Elemene	0.11	Sesquiterpene
(*Z*)-Jasmone	0.16	Jasmonate
Unknown	0.01	Sesquiterpene
Isocaryophyllene	0.03	Sesquiterpene
β-Caryophyllene	0.72	Sesquiterpene
β-Ylangene	0.15	Sesquiterpene
β-Copaene	0.18	Sesquiterpene
Isogermacrene D	0.13	Sesquiterpene
α-Humulene	0.07	Sesquiterpene
allo-Aromadendrene	0.02	Sesquiterpene
Unknown	0.04	Sesquiterpene
(*E*)-β-Farnesene	0.49	Sesquiterpene
trans-Cadina-1(6),4-diene	0.01	Sesquiterpene
Germacrene D	0.57	Sesquiterpene
Menthalactone	0.03	Monoterpenic lactone
Bicyclogermacrene	0.05	Sesquiterpene
α-Muurolene	0.06	Sesquiterpene
γ-Cadinene	0.02	Sesquiterpene
δ-Cadinene	0.07	Sesquiterpene
α-Cadinene	0.01	Sesquiterpene
1,5-Epoxysalvial-4(14)-ene	0.01	Sesquiterpenic ether
(*E*)-Nerolidol	0.01	Sesquiterpenic alcohol
Spathulenol	0.02	Sesquiterpenic alcohol
Caryophyllene oxide	0.04	Sesquiterpenic ether
Caryophyllene oxide isomer	0.01	Sesquiterpenic ether
Viridiflorol	0.08	Sesquiterpenic alcohol
Isospathulenol	0.02	Sesquiterpenic alcohol
τ-Cadinol	0.03	Sesquiterpenic alcohol
τ-Muurolol	0.01	Sesquiterpenic alcohol
α-Muurolol	0.02	Sesquiterpenic alcohol
α-Cadinol	0.03	Sesquiterpenic alcohol
Unknown	0.05	Unknown
meta-Camphorene	0.02	Diterpene
para-Camphorene	0.03	Diterpene
Consolidated total	**97.73%**	

*: Individual compounds concentration could not be found due to overlapping coelutions on columns considered
[xx]: Duplicate percentage due to coelutions, not taken into account in the consolidated total
tr: The compound has been detected below 0.005% of total signal.

Note: no correction factor was applied

About "consolidated" data: The table above presents the breakdown of the sample volatile constituents after applying an algorithm to collapse data acquired from the multi-columns system of PhytoChemia into a single set of consolidated contents. In case of discrepancies between columns, the algorithm is set to prioritize data from the most standard DB-5 column, and smallest values so as to avoid overestimating individual content. This process is semi-automatic. Advanced users are invited to consult the "Full analysis data" table after the chromatograms in this report to access the full untreated data and perform their own calculations if needed.

Unknowns: Unknown compounds' mass spectral data is presented in the "Full analysis data" table. The occurrence of unknown compounds is to be expected in many samples, and does not denote particular problems unless noted otherwise in the conclusion.

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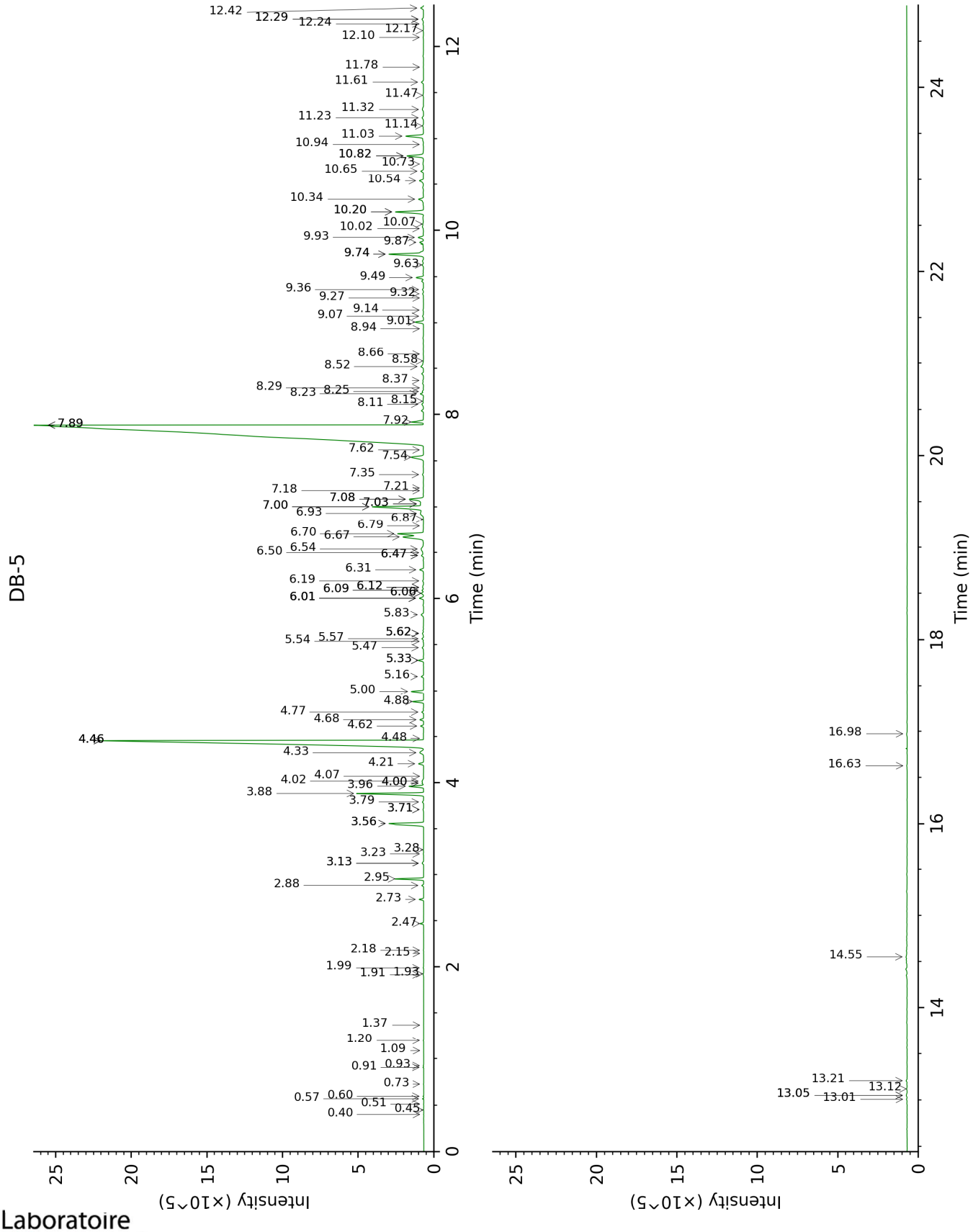
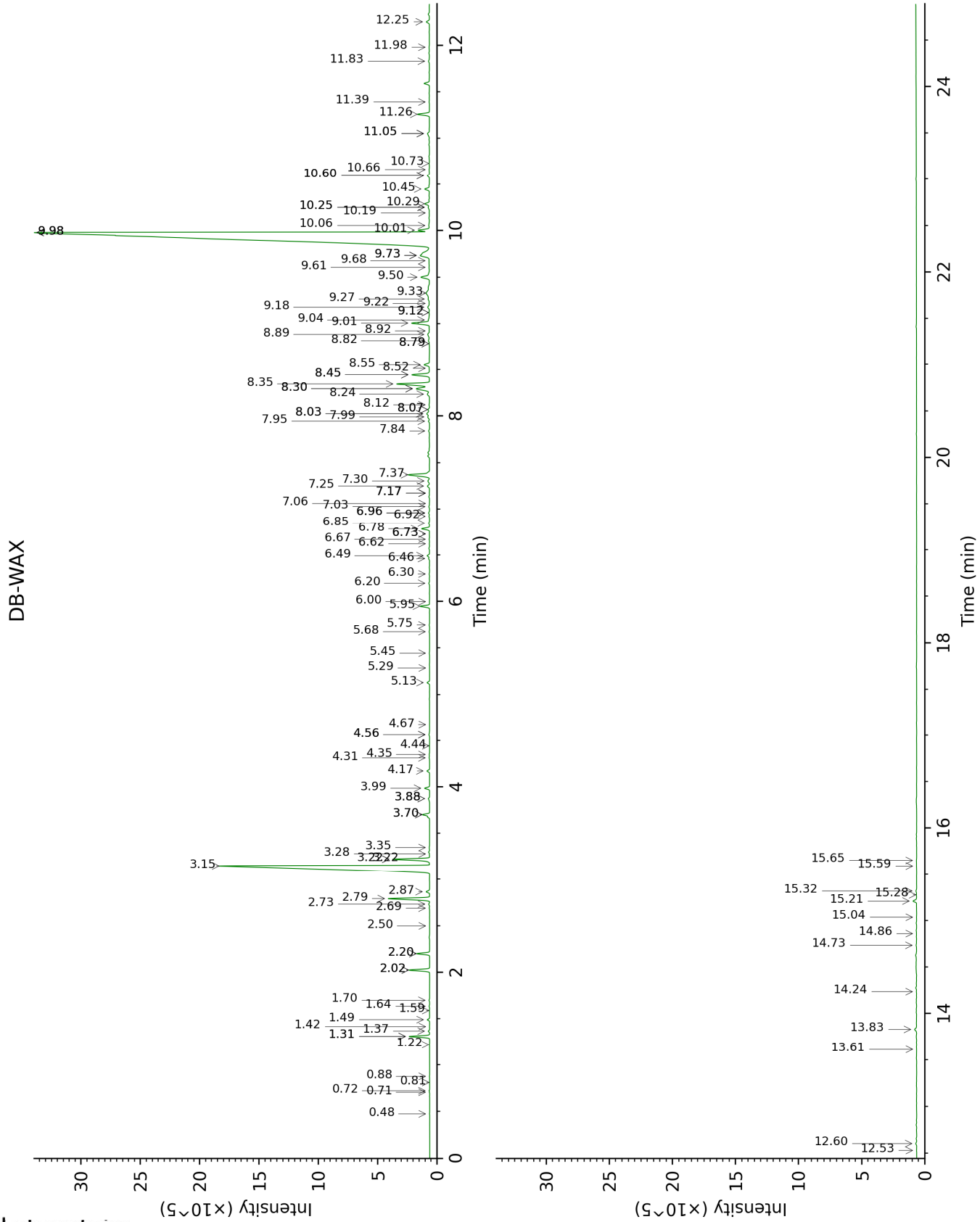

DB-5
Time (min)
Intensity (×10^5)
12.42
12.29
12.24
12.17
12.10
11.78
11.61
11.47
11.32
11.23
11.14
11.03
10.94
10.82
10.73
10.65
10.54
10.34
10.20
10.07
10.02
9.93
9.87
9.74
9.63
9.49
9.36
9.32
9.27
9.14
9.07
9.01
8.94
8.66
8.58
8.52
8.37
8.29
8.25
8.23
8.15
8.11
7.92
7.89
7.62
7.54
7.35
7.21
7.18
7.08
7.03
7.00
6.93
6.87
6.79
6.70
6.67
6.54
6.50
6.47
6.31
6.19
6.12
6.09
6.08
6.01
5.83
5.62
5.57
5.54
5.47
5.33
5.16
5.00
4.88
4.77
4.68
4.62
4.48
4.46
4.33
4.21
4.07
4.02
4.00
3.96
3.88
3.79
3.71
3.56
3.28
3.23
3.13
2.95
2.88
2.73
2.47
2.18
2.15
1.99
1.93
1.91
1.37
1.20
1.09
0.93
0.91
0.73
0.60
0.57
0.51
0.45
0.40
16.98
16.63
14.55
13.21
13.12
13.05
13.01


DB-WAX
Time (min)
Intensity (×10^5)
12.25
11.98
11.83
11.39
11.26
11.05
10.73
10.66
10.60
10.45
10.29
10.25
10.19
10.06
10.01
9.98
9.73
9.68
9.61
9.50
9.33
9.27
9.22
9.18
9.12
9.04
9.01
8.92
8.89
8.82
8.79
8.55
8.52
8.45
8.35
8.30
8.24
8.12
8.07
8.03
7.99
7.95
7.84
7.37
7.30
7.25
7.17
7.06
7.03
6.96
6.92
6.85
6.78
6.73
6.67
6.62
6.49
6.46
6.30
6.20
6.00
5.95
5.75
5.68
5.45
5.29
5.13
4.67
4.56
4.44
4.35
4.31
4.17
3.99
3.88
3.70
3.35
3.28
3.22
3.15
2.87
2.79
2.73
2.69
2.50
2.20
2.02
1.70
1.64
1.59
1.49
1.42
1.37
1.31
1.22
0.88
0.81
0.72
0.71
0.48
15.65
15.59
15.32
15.28
15.21
15.04
14.86
14.73
14.24
13.83
13.61
12.60
12.53

FULL ANALYSIS DATA

Identification	Column DB-5			Column DB-WAX		
	R.T	R.I	%	R.T	R.I	%
Isobutyral	0.40	536	tr	0.48	785	tr
3-Buten-2-one	0.45	580	tr	0.81	908	tr
Isobutanol	0.51	619	tr	2.02*	1068	0.81
Isovaleral	0.57	640	0.01	0.72	886	0.01
2-Methylbutyral	0.60	651	0.01	0.71	880	0.01
2-Ethylfuran	0.73	701	tr	0.88	919	0.01
Isoamyl alcohol	0.91	730	0.01	3.28	1174	0.01
2-Methylbutanol	0.93	733	0.01	3.35	1179	0.02
Toluene	1.09	758	tr	1.42	1006	tr
Methyl 2-methylbutyrate	1.20	774	0.01	1.22	977	0.01
Hexanal	1.37	800	tr			
Ethyl 2-methylbutyrate	1.91	849	0.01	1.59	1024	0.01
Ethyl isovalerate	1.92	850	0.01	1.70	1035	0.02
(3*Z*)-Hexenol	1.99	856	0.01	5.68	1350	0.02
(2*E*)-Hexenol	2.15	869	tr	6.00	1373	0.01
Hexanol	2.18	872	tr	5.29	1321	tr
trans-2,5-Diethyltetrahydrofuran	2.47	897	0.07	1.49	1014	0.08
Hashishene	2.73	916	0.10	1.31*	993	0.73
α-Thujene	2.88	926	0.04	1.37	1002	0.05
α-Pinene	2.95	931	0.63	1.31*	993	[0.73]
Camphene	3.13*	943	0.05	1.64	1028	0.02
3-Methylcyclohexanone	3.13*	943	[0.05]	4.56*	1269	0.03
Thuja-2,4(10)-diene	3.23	950	0.01	2.20*	1086	0.48
Benzaldehyde	3.28	953	0.01	7.17*	1460	0.01
β-Pinene	3.56*	972	1.28	2.02*	1068	[0.81]
Sabinene	3.56*	972	[1.28]	2.20*	1086	[0.48]
Octen-3-one	3.71*	982	0.02	4.44	1260	tr
Octen-3-ol	3.71*	982	[0.02]	6.62	1419	0.01
Octan-3-one	3.79	988	0.04	3.88*	1219	0.07
Myrcene	3.88	994	1.71	2.79	1135	1.69
Octan-3-ol	3.96	999	0.37	5.95	1369	0.37
α-Phellandrene	4.00*	1002	0.04	2.69	1127	0.01
Pseudolimonene	4.00*	1002	[0.04]	2.73	1131	0.03
Octanal	4.02	1003	0.06	4.31	1251	0.02
Δ3-Carene	4.07	1006	tr	2.50	1112	tr
α-Terpinene	4.21	1015	0.14	2.87	1141	0.14
para-Cymene	4.33	1023	0.19	3.99	1227	0.20
1,8-Cineole	4.46*	1031	18.92	3.22*	1169	1.49
β-Phellandrene	4.46*	1031	[18.92]	3.22*	1169	[1.49]
Limonene	4.46*	1031	[18.92]	3.15	1164	17.45
2-Ethylhexanol	4.48	1032	0.01	7.17*	1460	[0.01]
(*Z*)-β-Ocimene	4.62	1041	0.07	3.70*	1206	0.29
Unknown [m/z 57, 73 (44), 43 (26), 115 (25), 55 (14), 97 (8)...]	4.68	1045	0.09			

(*E*)-β-Ocimene	4.77	1051	0.05	3.88*	1219	[0.07]
γ-Terpinene	4.88	1058	0.27	3.70*	1206	[0.29]
cis-Sabinene hydrate	5.00	1065	0.32	6.78	1431	0.33
Octanol	5.16	1075	0.06	8.07*	1528	0.07
Terpinolene	5.33*	1086	0.13	4.17	1241	0.11
para-Cymenene	5.33*	1086	[0.13]	6.20	1387	0.03
trans-Sabinene hydrate	5.47	1095	0.04	7.84	1510	0.04
2-Methylbutyl isovalerate?	5.54	1100	0.01	4.56*	1269	[0.03]
Linalool	5.57	1101	0.05	7.95	1518	0.04
Isoamyl isovalerate	5.62*	1105	0.04	4.67	1277	tr
2-Methylbutyl 2-methylbutyrate	5.62*	1105	[0.04]	4.35	1253	0.02
Nonanal	5.62*	1105	[0.04]	5.75	1355	0.02
trans-para-Mentha-2,8-dien-1-ol	5.83	1118	0.09	8.89	1592	0.12
cis-Limonene oxide	6.01*	1130	0.12	6.30	1394	0.02
allo-Ocimene	6.01*	1130	[0.12]	5.45	1333	0.01
Octan-3-yl acetate	6.01*	1130	[0.12]	5.13	1310	0.10
trans-Pinocarveol	6.06*	1133	0.07	9.04	1604	0.05
cis-para-Mentha-2,8-dien-1-ol	6.06*	1133	[0.07]	9.27	1623	0.07
trans-Limonene oxide	6.09*	1135	0.04	6.46	1407	0.04
Camphor	6.09*	1135	[0.04]	7.06	1452	0.01
cis-Verbenol	6.12*	1137	0.05	9.12*	1611	0.02
trans-para-Menth-2-en-1-ol	6.12*	1137	[0.05]	8.79*	1584	0.01
Isopulegol	6.19	1142	0.03	8.03*	1525	0.18
Menthone	6.31	1149	0.11	6.49	1409	0.10
Menthofuran	6.47*	1159	0.07	6.73*	1427	0.01
Isomenthone	6.47*	1159	[0.07]	6.85	1435	0.04
Borneol	6.50	1161	0.02	9.61	1650	0.02
neo-Menthol	6.54	1164	0.16	8.45*	1558	0.79
Menthol	6.67	1172	0.84	9.01	1602	0.86
Terpinen-4-ol	6.70	1174	0.79	8.45*	1558	[0.79]
Isomenthol	6.79	1180	tr	8.79*	1584	[0.01]
para-Cymen-8-ol	6.86	1185	tr	11.39	1801	0.02
α-Terpineol	6.93†	1189	2.74	9.68	1656	0.02
cis-Dihydrocarvone	7.00*†	1194	[2.74]	8.35	1550	1.59
Myrtenal	7.00*†	1194	[2.74]	8.52	1563	0.01
neo-Dihydrocarveol	7.03*†	1196	[2.74]	10.01†	1683	[61.94]
Methylchavicol	7.03*†	1196	[2.74]	9.22	1619	0.01
trans-Dihydrocarvone	7.08*†	1199	[2.74]	8.55	1566	0.25
Dihydrocarveol	7.08*†	1199	[2.74]	10.29	1706	0.22
trans-Piperitol	7.18	1205	tr	10.19	1698	0.01
Decanal	7.21	1207	0.01	7.17*	1460	[0.01]
iso-Dihydrocarveol ?	7.35	1217	0.04	10.60*	1733	0.09
trans-Carveol	7.54	1229	0.51	11.26	1789	0.53
(3*Z*)-Hexenyl 2-methylbutyrate	7.62	1234	0.01	6.92	1441	0.02
Carvone	7.89*†	1252	62.40	9.98*†	1681	61.94

Piperitone	7.89*†	1252	[62.40]	9.74*	1661	1.01
cis-Carvone oxide	7.92†	1255	[62.40]	10.73	1744	0.01
Isopiperitenone	8.11	1268	0.06	11.05*	1772	0.14
Unknown [m/z 82, 109 (35), 135 (22), 127 (19), 54 (16), 43 (14)...]	8.15	1270	0.02			
trans-Carvone oxide	8.23	1275	0.08	11.05*	1772	[0.14]
Decanol	8.25	1277	0.03	10.60*	1733	[0.09]
2-Ethylmenthone?	8.29	1280	0.01			
Dihydroedulan I	8.37	1285	0.02	6.96*	1444	0.04
Menthyl acetate	8.52	1295	0.10	7.99	1522	0.10
Dihydroedulan II	8.58	1299	0.01	7.30	1470	0.10
Isomenthyl acetate	8.66	1304	0.01	8.12	1532	0.02
Bicycloelemene analog	8.94	1324	tr	6.73*	1427	[0.01]
Dihydrocarvyl acetate	9.01	1329	0.23	9.33	1628	0.35
Bicycloelemene	9.07	1334	0.03	6.96*	1444	[0.04]
trans-Carvyl acetate	9.14	1338	0.02	10.06	1687	0.02
α-Cubebene	9.27	1348	0.01	6.67	1422	0.02
Menthofurolactone	9.32	1351	0.03	11.83	1840	0.04
iso-Dihydrocarvyl acetate	9.36	1354	0.03			
cis-Carvyl acetate	9.49	1363	0.21	10.45	1720	0.21
α-Copaene	9.63	1373	0.04	7.03	1449	0.03
β-Bourbonene	9.74*	1381	1.15	7.37	1475	1.15
1,5-diepi-β-Bourbonene	9.74*	1381	[1.15]	7.25	1466	0.11
β-Elemene	9.87	1390	0.11	8.30*	1546	0.83
(*Z*)-Jasmone	9.93	1394	0.16	12.25	1877	0.15
Unknown [m/z 106, 119 (99), 43 (78), 91 (74), 105 (60), 134 (55)... 204 (19)]	10.02	1401	0.01			
Isocaryophyllene	10.07	1404	0.03	8.07*	1528	[0.07]
β-Caryophyllene	10.20*	1414	0.91	8.30*	1546	[0.83]
β-Ylangene	10.20*	1414	[0.91]	8.03*	1525	[0.18]
β-Copaene	10.34	1424	0.18	8.24	1541	0.13
Isogermacrene D	10.54	1440	0.13	8.82	1587	0.06
α-Humulene	10.65	1448	0.07	9.18	1615	0.10
allo-Aromadendrene	10.73	1454	0.02	8.92	1595	0.01
Unknown [m/z 161, 105 (56), 91 (50), 93 (36), 119 (33), 79 (31)...204 (5)]	10.82*	1460	0.53			
(*E*)-β-Farnesene	10.82*	1460	[0.53]	9.50	1642	0.49
trans-Cadina-1(6),4-diene	10.94	1469	0.01	9.12*	1611	[0.02]
Germacrene D	11.03	1476	0.57	9.74*	1661	[1.01]
Menthalactone	11.14	1484	0.03	15.65	2204	0.02
Bicyclogermacrene	11.23	1491	0.05	9.98*†	1681	[61.94]
α-Muurolene	11.32	1498	0.06	9.98*†	1681	[61.94]
γ-Cadinene	11.47	1509	0.02	10.25*	1703	0.07
δ-Cadinene	11.61	1520	0.07	10.25*	1703	[0.07]

α-Cadinene	11.78	1533	0.01	10.66	1738	0.02
1,5-Epoxysalvial-4(14)-ene	12.10	1559	0.01	11.98	1853	0.03
(*E*)-Nerolidol	12.17	1564	0.01	13.61	2003	tr
Spathulenol	12.24	1570	0.02	14.24	2063	0.01
Caryophyllene oxide	12.30*	1574	0.06	12.60	1909	0.04
Caryophyllene oxide isomer	12.30*	1574	[0.06]	12.53	1902	0.01
Viridiflorol	12.42	1584	0.08	13.83	2024	0.11
Isospathulenol	13.01	1632	0.02	15.28	2167	0.01
τ-Cadinol	13.05*	1635	0.03	14.74	2112	0.03
τ-Muurolol	13.05*	1635	[0.03]	14.86	2125	0.01
α-Muurolol	13.12	1641	0.02	15.04	2142	0.01
α-Cadinol	13.21	1648	0.03	15.32	2171	0.03
Unknown [m/z 95, 139 (69), 55 (65), 81 (55), 108 (51)...]	14.55	1762	0.05			
meta-Camphorene	16.63	1952	0.02	15.21	2160	0.15
para-Camphorene	16.98	1985	0.03	15.59	2198	0.01
Total identified		**98.25%**			**97.65%**	
Total reported		**98.42%**			**97.65%**	

*: Two or more compounds are coeluting on this column
[xx]: Duplicate percentage due to coelutions, not taken into account in the consolidated total
†: Peaks apexes were resolved, but peaks overlapped and were summed for analysis
tr: The compound has been detected below 0.005% of total signal.

Note: no correction factor was applied
R.T.: Retention time (minutes)
R.I.: Retention index