Date : 2025-06-02

Certificate of Analysis - GC Profiling

Sample Identification

Internal code : 25E16-PTH05
Customer Identification : Organic Spearmint - India - S40110R
Type : Essential Oil
Source : *Mentha spicata*
Customer : Plant Therapy

Checked an approved by:

Alexis St-Gelais, Ph. D., Chimiste 2013-174

Notes: This report may not be published, including online, without the written consent from Laboratoire PhytoChemia.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

Laboratoire PhytoChemia

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Gas Chromatographic Analysis

Method : PC-MAT-014 - Analysis of the composition of an essential oil or other volatile liquide by FAST GC-FID
ISO
Results : See analysis summary (next page)
Analyst : Sylvain Mercier, M. Sc., Chimiste 2014-005
Date : 2025-05-27

Physicochemical data

Refractive index : 1.4908 ± 0.0003 (20 °C)
Method : PC-MAT-016 - Measure of the refractive index of a liquid.
Analyst : Cindy Caron B. Sc.
Date : 2025-05-16

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
Isovaleral	tr	Aliphatic aldehyde
2-Methylbutyral	tr	Aliphatic aldehyde
Isoamyl alcohol	tr	Aliphatic alcohol
2-Methylbutanol	tr	Aliphatic alcohol
Methyl 2-methylbutyrate	tr	Aliphatic ester
Ethyl 2-methylbutyrate	0.01	Aliphatic ester
(3*Z*)-Hexenol	0.01	Aliphatic alcohol
Hexanol	tr	Aliphatic alcohol
trans-2,5-Diethyltetrahydrofuran	0.05	Furan
Hashishene	0.08	Monoterpene
α-Thujene	0.04	Monoterpene
α-Pinene	0.58	Monoterpene
trans-3-Methylcyclohexanol	0.01	Aliphatic alcohol
Camphene	0.01	Monoterpene
3-Methylcyclohexanone	0.03	Aliphatic ketone
α-Fenchene	tr	Monoterpene
Propyl 2-methylbutyrate	tr	Aliphatic ester
Thuja-2,4(10)-diene	0.01	Monoterpene
Sabinene	0.42	Monoterpene
β-Pinene	0.77	Monoterpene
Octen-3-ol	0.02	Aliphatic alcohol
cis-Carane	0.01	Monoterpene
Octan-3-one	0.03	Aliphatic ketone
Myrcene	1.33	Monoterpene
Octan-3-ol	0.39	Aliphatic alcohol
α-Phellandrene	0.02	Monoterpene
Octanal	0.02	Aliphatic aldehyde
Pseudolimonene	0.04	Monoterpene
Δ3-Carene	tr	Monoterpene
α-Terpinene	0.10	Monoterpene
Carvomenthene	0.02	Aliphatic alcohol
para-Cymene	0.21	Monoterpene
β-Phellandrene	[1.34]	Monoterpene
1,8-Cineole	[1.34]	Monoterpenic ether
Limonene	15.63	Monoterpene
2-Ethylhexanol	0.01	Aliphatic alcohol
(*Z*)-β-Ocimene	0.07	Monoterpene
(*E*)-β-Ocimene	0.05	Monoterpene
γ-Terpinene	0.19	Monoterpene
cis-Sabinene hydrate	0.24	Monoterpenic alcohol

Octanol	0.04	Aliphatic alcohol
Terpinolene	0.08	Monoterpene
para-Cymenene	0.03	Monoterpene
trans-Sabinene hydrate	0.04	Monoterpenic alcohol
Linalool	0.07	Monoterpenic alcohol
2-Methylbutyl 2-methylbutyrate	0.01	Aliphatic ester
Nonanal	0.01	Aliphatic aldehyde
endo-Fenchol	0.01	Monoterpenic alcohol
trans-para-Mentha-2,8-dien-1-ol	0.07	Monoterpenic alcohol
cis-Limonene oxide	0.02	Monoterpenic ether
Octan-3-yl acetate	0.09	Aliphatic ester
cis-para-Mentha-2,8-dien-1-ol	0.05	Monoterpenic alcohol
trans-Limonene oxide	0.04	Monoterpenic ether
Camphor	0.02	Monoterpenic ketone
trans-Sabinol	0.01	Monoterpenic alcohol
Isopulegol	0.03	Monoterpenic alcohol
Menthone	0.19	Monoterpenic ketone
Menthofuran	0.01	Monoterpenic ether
Isomenthone	0.06	Monoterpenic ketone
neo-Menthol	0.10	Monoterpenic alcohol
δ-Terpineol	0.05	Monoterpenic alcohol
Menthol	0.77	Monoterpenic alcohol
Terpinen-4-ol	0.44	Monoterpenic alcohol
Isomenthol	0.02	Monoterpenic alcohol
α-Terpineol	0.20	Monoterpenic alcohol
neoiso-Menthol	0.03	Monoterpenic alcohol
cis-Piperitol	0.01	Monoterpenic alcohol
cis-Dihydrocarvone	1.20	Monoterpenic ketone
Myrtenol	0.01	Monoterpenic alcohol
Dihydrocarveol	0.16	Monoterpenic alcohol
neo-Dihydrocarveol	0.24	Monoterpenic alcohol
Methylchavicol	0.14	Phenylpropanoid
trans-Dihydrocarvone	0.22	Monoterpenic ketone
iso-Dihydrocarveol ?	0.03	Monoterpenic alcohol
trans-Carveol	0.29	Monoterpenic alcohol
Pulegone	0.03	Monoterpenic ketone
cis-Carveol	0.16	Monoterpenic alcohol
Carvone	67.50	Monoterpenic ketone
Piperitone	0.75	Monoterpenic ketone
Isopiperitenone	0.04	Monoterpenic ketone
trans-Carvone oxide	0.07	Monoterpenic ketone
Decanol	0.03	Aliphatic alcohol
Dihydroedulan I	0.01	Terpenic ether
Dihydroedulan II	0.05	Terpenic ether
Menthyl acetate	0.10	Monoterpenic ester

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Thymol	0.01	Monoterpenic alcohol
Isomenthyl acetate	0.01	Monoterpenic alcohol
neo-Dihydrocarvyl acetate	0.01	Monoterpenic ester
Dihydrocarvyl acetate	0.23	Monoterpenic ester
Bicycloelemene	0.02	Sesquiterpene
trans-Carvyl acetate	0.01	Monoterpenic ester
α-Cubebene	0.01	Sesquiterpene
iso-Dihydrocarvyl acetate	0.02	Monoterpenic ester
cis-Carvyl acetate	0.16	Monoterpenic ester
α-Copaene	0.05	Sesquiterpene
β-Bourbonene	1.06	Sesquiterpene
1,5-diepi-β-Bourbonene	0.05	Sesquiterpene
β-Elemene	0.12	Sesquiterpene
(*Z*)-Jasmone	0.14	Jasmonate
Longifolene	0.01	Sesquiterpene
Unknown	0.01	Sesquiterpene
Isocaryophyllene	0.04	Sesquiterpene
β-Caryophyllene	0.69	Sesquiterpene
β-Ylangene	0.13	Sesquiterpene
β-Copaene	0.10	Sesquiterpene
Aromadendrene	0.02	Sesquiterpene
Isogermacrene D	0.09	Sesquiterpene
α-Humulene	0.06	Sesquiterpene
allo-Aromadendrene	0.01	Sesquiterpene
(*E*)-β-Farnesene	0.40	Sesquiterpene
trans-Cadina-1(6),4-diene	0.01	Sesquiterpene
γ-Muurolene	0.01	Sesquiterpene
Germacrene D	0.40	Sesquiterpene
Viridiflorene	0.03	Sesquiterpene
α-Muurolene	0.02	Sesquiterpene
γ-Cadinene	0.01	Sesquiterpene
δ-Cadinene	0.04	Sesquiterpene
Caryophyllene oxide	0.04	Sesquiterpenic ether
Caryophyllene oxide isomer	0.01	Sesquiterpenic ether
Viridiflorol	0.05	Sesquiterpenic alcohol
Isospathulenol	0.01	Sesquiterpenic alcohol
Consolidated total	**99.40**	

tr: The compound has been detected below 0.005% of the 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.

Bracketed value ([xx]): A bracketed percent value indicate that two or more compound percentage could not be solved due to coelution.

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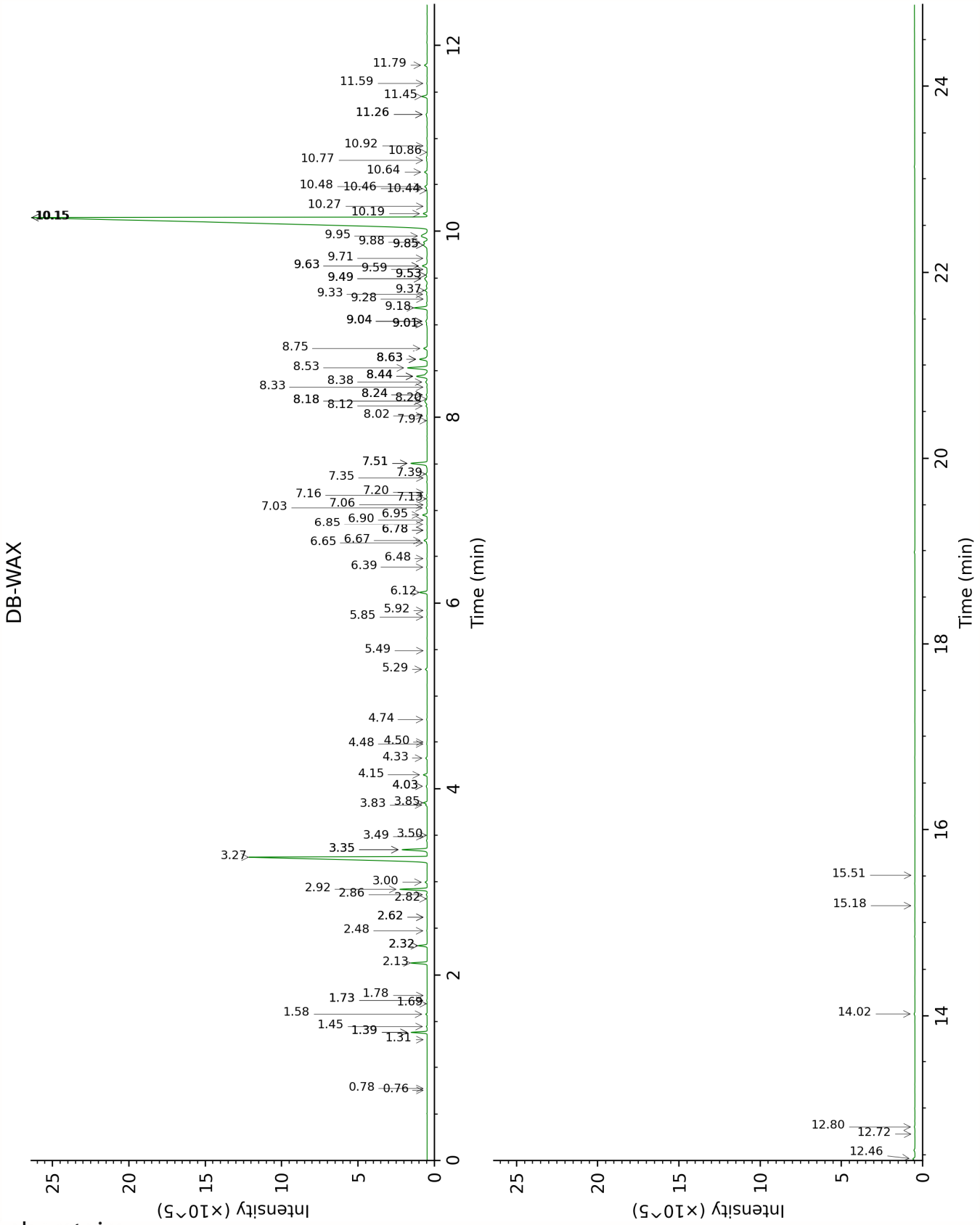
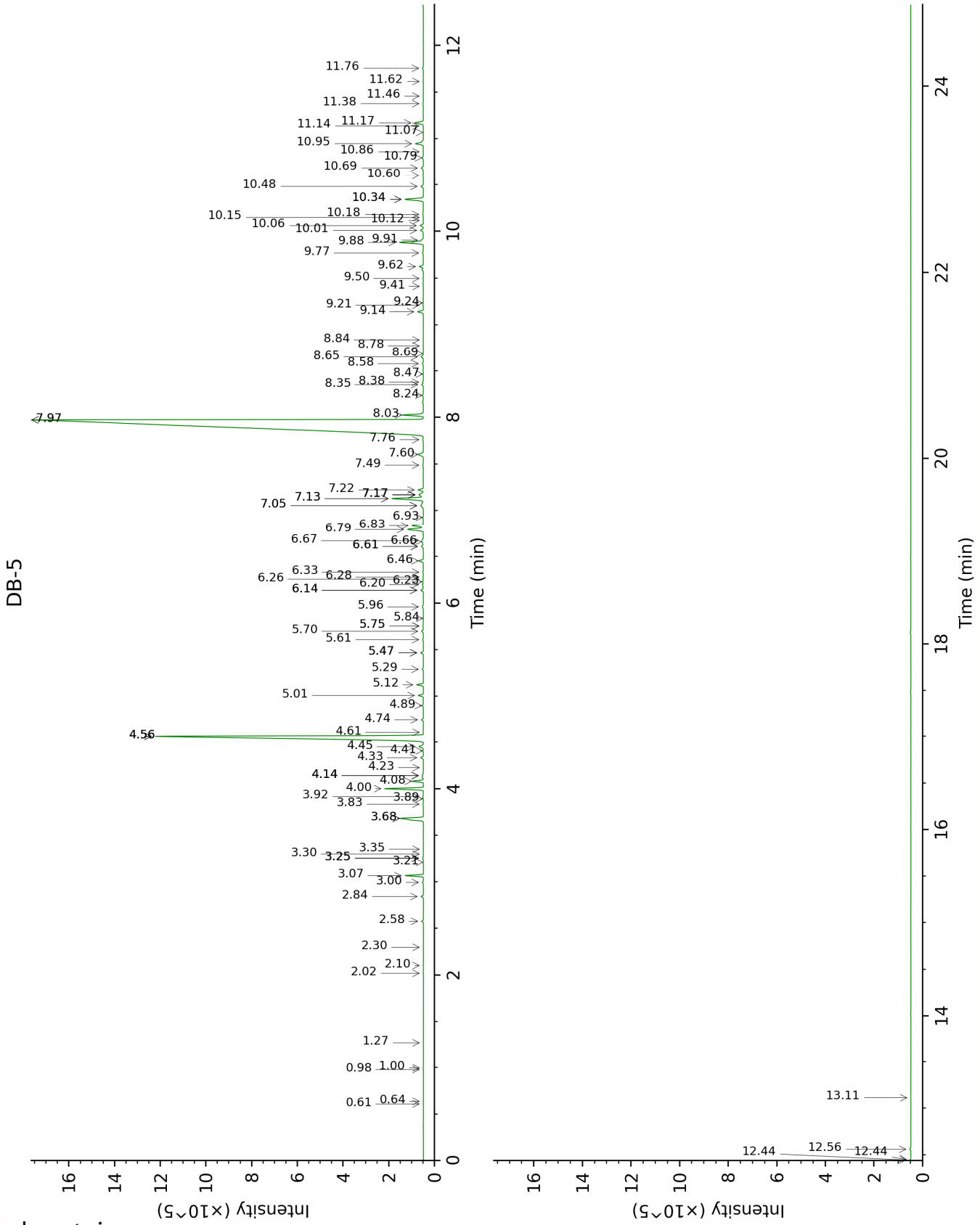

DB-WAX
Time (min)
Intensity (×10^5)
11.79
11.59
11.45
11.26
10.92
10.86
10.77
10.64
10.48
10.46
10.44
10.27
10.19
10.15
9.95
9.88
9.85
9.71
9.63
9.59
9.53
9.49
9.37
9.33
9.28
9.18
9.04
9.01
8.75
8.63
8.53
8.44
8.38
8.33
8.24
8.20
8.18
8.12
8.02
7.97
7.51
7.39
7.35
7.20
7.16
7.13
7.06
7.03
6.95
6.90
6.85
6.78
6.67
6.65
6.48
6.39
6.12
5.92
5.85
5.49
5.29
4.74
4.50
4.48
4.33
4.15
4.03
3.85
3.83
3.50
3.49
3.35
3.27
3.00
2.92
2.86
2.82
2.62
2.48
2.32
2.13
1.78
1.73
1.69
1.58
1.45
1.39
1.31
0.78
0.76
15.51
15.18
14.02
12.80
12.72
12.46


DB-5
Time (min)
Intensity (×10^5)
11.76
11.62
11.46
11.38
11.14
11.17
11.07
10.95
10.86
10.79
10.69
10.60
10.48
10.34
10.15
10.18
10.12
10.06
10.01
9.88
9.91
9.77
9.62
9.50
9.41
9.21
9.24
9.14
8.84
8.78
8.65
8.69
8.58
8.47
8.35
8.38
8.24
8.03
7.97
7.76
7.60
7.49
7.22
7.17
7.13
7.05
6.93
6.79
6.83
6.67
6.61
6.66
6.46
6.33
6.28
6.26
6.20
6.23
6.14
5.96
5.84
5.75
5.70
5.61
5.47
5.29
5.12
5.01
4.89
4.74
4.61
4.56
4.45
4.41
4.33
4.23
4.14
4.08
4.00
3.92
3.89
3.83
3.68
3.35
3.30
3.25
3.21
3.07
3.00
2.84
2.58
2.30
2.10
2.02
1.27
1.00
0.98
0.64
0.61
13.11
12.56
12.44
12.44

FULL ANALYSIS DATA

	Column DB-WAX			Column DB-5		
Isovaleral	0.78	886.2	tr	0.61	642.2	tr
2-Methylbutyral	0.76	879.0	tr	0.64	652.2	tr
Isoamyl alcohol	3.49	1177.6	tr	0.98	732.2	tr
2-Methylbutanol	3.50	1178.7	tr	1.00	734.9	tr
Methyl 2-methylbutyrate	1.31	976.8	tr	1.27	774.4	tr
Ethyl 2-methylbutyrate	1.73*	1025.1	[0.02]	2.02	849.3	0.01
(3*Z*)-Hexenol	5.85	1350.6	0.02	2.10	856.3	0.01
Hexanol	5.49	1324.9	0.01	2.30	872.8	tr
trans-2,5-Diethyltetrahydrofuran	1.58	1010.5	0.05	2.58	896.3	0.05
Hashishene	1.39*	989.6	[0.65]	2.84	915.6	0.08
α-Thujene	1.45	997.3	0.03	3.00	925.7	0.04
α-Pinene	1.39*	989.6	[0.65]	3.07	930.6	0.58
trans-3-Methylcyclohexanol	6.85	1422.3	0.02	3.21	940.1	0.01
Camphene	1.73*	1025.1	[0.02]	3.26*	943.1	[0.04]
3-Methylcyclohexanone	4.74	1270.8	0.03	3.26*	943.1	[0.04]
α-Fenchene	1.69	1021.6	tr	3.26*	943.1	[0.04]
Propyl 2-methylbutyrate	2.62*	1109.6	[0.01]	3.30	946.0	tr
Thuja-2,4(10)-diene	2.32*	1083.4	[0.42]	3.35	949.6	0.01
Sabinene	2.32*	1083.4	[0.42]	3.68*	971.7	[1.19]
β-Pinene	2.13	1065.0	0.77	3.68*	971.7	[1.19]
Octen-3-ol	6.78*	1417.3	[0.01]	3.84	981.9	0.02
cis-Carane	1.78	1030.3	0.01	3.89	985.8	0.01
Octan-3-one	4.03*	1218.4	[0.06]	3.92	987.4	0.03
Myrcene	2.92	1133.2	1.33	4.00	993.1	1.33
Octan-3-ol	6.12	1369.3	0.40	4.08	998.5	0.39
α-Phellandrene	2.82	1125.1	0.02	4.14*†	1002.5	[0.06]
Octanal	4.48	1251.6	0.02	4.14*†	1002.5	[0.06]
Pseudolimonene	2.86	1128.6	0.04	4.14*†	1002.5	[0.06]
Δ3-Carene	2.62*	1109.6	[0.01]	4.23	1008.0	tr
α-Terpinene	3.00	1139.1	0.10	4.33	1014.8	0.10
Carvomenthene	2.48	1098.2	0.02	4.41	1019.5	0.02
para-Cymene	4.15	1227.4	0.20	4.45	1022.3	0.21
β-Phellandrene	3.35*	1166.6	[1.31]	4.56*	1029.3	[16.97]
1,8-Cineole	3.35*	1166.6	[1.31]	4.56*	1029.3	[16.97]
Limonene	3.27	1160.2	15.63	4.56*	1029.3	[16.97]
2-Ethylhexanol	7.35	1459.1	0.01	4.61	1032.0	0.01
(*Z*)-β-Ocimene	3.83	1203.7	0.07	4.74	1040.6	0.07
(*E*)-β-Ocimene	4.03*	1218.4	[0.06]	4.89	1050.3	0.05

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γ-Terpinene	3.85	1205.4	0.20	5.01	1057.6	0.19
cis-Sabinene hydrate	6.95	1429.9	0.25	5.12	1064.9	0.24
Octanol	8.24*	1525.9	[0.05]	5.29	1075.5	0.04
Terpinolene	4.33	1240.5	0.08	5.47*	1086.7	[0.10]
para-Cymenene	6.39	1388.5	0.03	5.47*	1086.7	[0.10]
trans-Sabinene hydrate	8.02	1508.6	0.03	5.61	1095.7	0.04
Linalool	8.12	1516.7	0.07	5.70	1101.4	0.07
2-Methylbutyl 2-methylbutyrate	4.50	1253.2	0.01	5.76*	1105.0	[0.03]
Nonanal	5.92	1355.5	0.01	5.76*	1105.0	[0.03]
endo-Fenchol	8.44*	1541.1	[0.82]	5.84	1110.3	0.01
trans-para-Mentha-2,8-dien-1-ol	9.01*	1584.3	[0.06]	5.96	1118.4	0.07
cis-Limonene oxide	6.48	1395.1	0.02	6.14*	1129.9	[0.11]
Octan-3-yl acetate	5.29	1310.7	0.09	6.14*	1129.9	[0.11]
cis-para-Mentha-2,8-dien-1-ol	9.53*	1626.3	[0.09]	6.20	1133.8	0.05
trans-Limonene oxide	6.65	1407.4	0.03	6.23	1135.8	0.04
Camphor	7.20	1447.8	0.01	6.26	1137.6	0.02
trans-Sabinol	9.88†	1654.4	0.26	6.28	1139.1	0.01
Isopulegol	8.20	1522.6	0.02	6.33	1142.5	0.03
Menthone	6.67	1409.3	0.19	6.46	1150.4	0.19
Menthofuran	6.90	1425.9	0.01	6.61*	1160.4	[0.09]
Isomenthone	7.03	1435.5	0.06	6.61*	1160.4	[0.09]
neo-Menthol	8.63*	1555.2	[0.49]	6.66	1163.7	0.10
δ-Terpineol	9.50*	1623.2	[0.25]	6.67	1164.5	0.05
Menthol	9.18	1598.2	0.77	6.79*†	1172.4	[0.74]
Terpinen-4-ol	8.63*	1555.2	[0.49]	6.83*†	1175.1	[0.46]
Isomenthol	9.01*	1584.3	[0.06]	6.93	1181.1	0.02
α-Terpineol	9.85*†	1652.0	[0.36]	7.05*	1189.4	[0.24]
neoiso-Menthol	9.53*	1626.3	[0.09]	7.05*	1189.4	[0.24]
cis-Piperitol	9.59	1631.2	0.01	7.13*†	1194.3	[1.37]
cis-Dihydrocarvone	8.53	1548.1	1.20	7.13*†	1194.3	[1.37]
Myrtenol	10.92	1740.2	0.01	7.17*†	1197.0	[0.28]
Dihydrocarveol	10.48	1702.8	0.16	7.17*†	1197.0	[0.28]
neo-Dihydrocarveol	10.19	1679.0	0.24	7.17*†	1197.0	[0.28]
Methylchavicol	9.37	1613.2	0.14	7.17*†	1197.0	[0.28]
trans-Dihydrocarvone	8.75	1564.6	0.22	7.22*†	1200.4	[0.21]
iso-Dihydrocarveol ?	10.86	1734.4	0.01	7.49	1218.2	0.03
trans-Carveol	11.45	1784.7	0.30	7.60	1226.1	0.29
Pulegone	9.01*	1584.3	[0.06]	7.76	1236.9	0.03
cis-Carveol	11.79	1814.0	0.16	7.98*	1251.4	[67.66]
Carvone	10.15*	1675.6	[67.36]	7.98*	1251.4	[67.66]
Piperitone	9.95	1659.9	0.67	8.03	1255.0	0.75

Isopiperitenone	11.26*	1768.5	[0.08]	8.24	1269.4	0.04
trans-Carvone oxide	11.26*	1768.5	[0.08]	8.35	1277.2	0.07
Decanol	10.77	1727.1	0.04	8.38	1279.0	0.03
Dihydroedulan I	7.13	1442.7	0.04	8.47	1285.0	0.01
Dihydroedulan II	7.51*	1470.5	[1.08]	8.58	1292.7	0.05
Menthyl acetate	8.18*	1520.7	[0.24]	8.65	1297.6	0.10
Thymol	15.18	2130.5	0.01	8.69	1300.1	0.01
Isomenthyl acetate	8.33	1532.2	0.02	8.78	1306.1	0.01
neo-Dihydrocarvyl acetate	9.04*	1586.9	[0.11]	8.84	1306.8	0.01
Dihydrocarvyl acetate	9.50*	1623.2	[0.25]	9.14	1328.3	0.23
Bicycloelemene	7.06	1437.9	0.01	9.21	1333.3	0.02
trans-Carvyl acetate	10.27	1685.3	0.01	9.24	1335.1	0.01
α-Cubebene	6.78*	1417.3	[0.01]	9.41	1347.7	0.01
iso-Dihydrocarvyl acetate				9.50	1353.7	0.02
cis-Carvyl acetate	10.64	1716.5	0.16	9.62	1362.7	0.16
α-Copaene	7.16	1445.3	0.04	9.77	1373.1	0.05
β-Bourbonene	7.51*	1470.5	[1.08]	9.88	1381.2	1.06
1,5-diepi-β-Bourbonene	7.39	1462.2	0.09	9.91	1382.9	0.05
β-Elemene	8.44*	1541.1	[0.82]	10.01	1390.3	0.12
(*Z*)-Jasmone	12.46	1872.7	0.12	10.06	1394.0	0.14
Longifolene	7.97	1504.6	0.01	10.12	1397.8	0.01
Unknown MEPI VIII [m/z 106, 119 (99), 43 (78), 91 (74), 105 (60), 134 (55)... 204 (19)]	11.59	1796.9	0.02	10.15	1400.1	0.01
Isocaryophyllene	8.24*	1525.9	[0.05]	10.18	1402.4	0.04
β-Caryophyllene	8.44*	1541.1	[0.82]	10.34*	1414.5	[0.84]
β-Ylangene	8.18*	1520.7	[0.24]	10.34*	1414.5	[0.84]
β-Copaene	8.38	1536.4	0.11	10.48	1424.9	0.10
Aromadendrene	8.63*	1555.2	[0.49]	10.60	1433.9	0.02
Isogermacrene D	9.04*	1586.9	[0.11]	10.68	1440.1	0.09
α-Humulene	9.33	1609.8	0.05	10.80	1448.4	0.06
allo-Aromadendrene	9.04*	1586.9	[0.11]	10.86	1453.2	0.01
(*E*)-β-Farnesene	9.63*	1634.2	[0.35]	10.95	1459.8	0.40
trans-Cadina-1(6),4-diene	9.28	1605.7	0.01	11.07	1468.9	0.01
γ-Muurolene	9.63*	1634.2	[0.35]	11.14	1474.0	0.01
Germacrene D	9.85*†	1652.0	[0.36]	11.17	1476.6	0.40
Viridiflorene	9.71	1640.7	0.04	11.38	1492.0	0.03
α-Muurolene	10.15*	1675.6	[67.36]	11.46	1498.1	0.02
γ-Cadinene	10.44	1699.2	0.01	11.62	1510.1	0.01
δ-Cadinene	10.46	1700.9	0.04	11.76	1521.3	0.04

Caryophyllene oxide	12.80	1903.6	0.04	12.44*	1575.6	[0.04]
Caryophyllene oxide isomer	12.72	1896.5	0.01	12.44*	1575.6	[0.04]
Viridiflorol	14.02	2017.1	0.05	12.56	1584.9	0.05
Isospathulenol	15.51	2163.0	0.01	13.11	1629.8	0.01
Total reported		98.82%			99.33%	

*: Two or more compounds are coeluting on this column

[xx]: Duplicate percentage due to coelutions, only the first one is 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