**Date :** January 28, 2022

# *Certificate of Analysis – GC Profiling*

## *Sample identification*

**Internal code :** 22A14-PTH08
**Customer identification :** Laurel Leaf - Greece - L10108217R
**Type :** Essential oil
**Source :** *Laurus nobilis*
**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, Ph. D.
**Analysis date :** January 26, 2022

Checked and approved by :

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

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*PHYSICOCHEMICAL DATA*

**Physical aspect:** Faintly yellow liquid
**Refractive index:** 1.4661 ± 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 |
|---|---|---|
| Methylcyclopentadiene isomer I | tr | Alkene |
| Isovaleral | tr | Aliphatic aldehyde |
| 3-Methyl-1-penten-3-ol | 0.01 | Aliphatic alcohol |
| 1-Methylpyrrole | tr | Pyrrole |
| Ethyl isobutyrate | tr | Aliphatic ester |
| Toluene | 0.01 | Simple phenolic |
| Hexanal | 0.01 | Aliphatic aldehyde |
| Isopropyl isobutyrate | 0.01 | Aliphatic ester |
| Ethyl 2-methylbutyrate | 0.01 | Aliphatic ester |
| Ethyl isovalerate | 0.01 | Aliphatic ester |
| (3*Z*)-Hexenol | 0.08 | Aliphatic alcohol |
| Hexanol | 0.01 | Aliphatic alcohol |
| Isopropyl 2-methylbutyrate | 0.01 | Aliphatic ester |
| Tricyclene | 0.04 | Monoterpene |
| α-Thujene | 0.46 | Monoterpene |
| α-Pinene | 6.72 | Monoterpene |
| Camphene | 0.78 | Monoterpene |
| Thujadiene isomer | 0.06 | Monoterpene |
| α-Fenchene | 0.03 | Monoterpene |
| Sabinene | 7.20 | Monoterpene |
| β-Pinene | 4.94 | Monoterpene |
| 6-Methyl-5-hepten-2-one | tr | Aliphatic ketone |
| Dehydro-1,8-cineole | 0.15 | Monoterpenic ether |
| Myrcene | 0.81 | Monoterpene |
| α-Phellandrene | 0.18 | Monoterpene |
| Pseudolimonene | 0.03 | Monoterpene |
| Isobutyl 2-methylbutyrate | 0.02 | Aliphatic ester |
| Δ3-Carene | 0.07 | Monoterpene |
| (3*Z*)-Hexenyl acetate | 0.01 | Aliphatic ester |
| α-Terpinene | 0.58 | Monoterpene |
| Unknown | 0.01 | Unknown |
| para-Cymene | 1.30 | Monoterpene |
| Limonene | 1.70 | Monoterpene |
| 1,8-Cineole | 47.43 | Monoterpenic ether |
| (*Z*)-β-Ocimene | 0.02 | Monoterpene |
| Unknown | 0.01 | Unknown |
| (*E*)-β-Ocimene | 0.08 | Monoterpene |
| γ-Terpinene | 0.90 | Monoterpene |
| *cis*-Sabinene hydrate | 0.14 | Monoterpenic alcohol |
| para-Mentha-3,8-diene | 0.01 | Monoterpene |
| *cis*-Linalool oxide (fur.) | 0.01 | Monoterpenic alcohol |
| Terpinolene | 0.27 | Monoterpene |
| para-Cymenene | 0.03 | Monoterpene |
| *trans*-Linalool oxide (fur.) | tr | Monoterpenic alcohol |
| 2-Nonanone | 0.02 | Aliphatic ketone |

| | | |
|---|---|---|
| *trans*-Sabinene hydrate | 0.11 | Monoterpenic alcohol |
| Linalool | 1.98 | Monoterpenic alcohol |
| Unknown | 0.06 | Unknown |
| Hotrienol | 0.06 | Monoterpenic alcohol |
| endo-Fenchol | 0.02 | Monoterpenic alcohol |
| Unknown | 0.01 | Unknown |
| *cis*-para-Menth-2-en-1-ol | 0.13 | Monoterpenic alcohol |
| Limona ketone | 0.01 | Normonoterpenic ketone |
| *trans*-Pinocarveol | 0.17 | Monoterpenic alcohol |
| *trans*-para-Menth-2-en-1-ol | 0.08 | Monoterpenic alcohol |
| *trans*-Verbenol | 0.05 | Monoterpenic alcohol |
| Sabinaketone | 0.02 | Normonoterpenic ketone |
| Nerol oxide | 0.03 | Aliphatic ether |
| (*E*)-2,6-Dimethyl-1,5,7-octatrien-3-ol | 0.05 | Monoterpenic alcohol |
| Pinocarvone | 0.09 | Monoterpenic ketone |
| Borneol | 0.22 | Monoterpenic alcohol |
| δ-Terpineol | 0.32 | Monoterpenic alcohol |
| Rosefuran oxide | 3.01* | Monoterpenic ether |
| Terpinen-4-ol | 3.01* | Monoterpenic alcohol |
| Thuj-3-en-10-al | 0.02 | Monoterpenic aldehyde |
| para-Cymen-8-ol | 0.01 | Monoterpenic alcohol |
| α-Terpineol | 1.45 | Monoterpenic alcohol |
| Myrtenal | 0.07 | Monoterpenic aldehyde |
| Myrtenol | 0.09 | Monoterpenic alcohol |
| *cis*-Piperitol | 0.04 | Monoterpenic alcohol |
| Methylchavicol | 0.01 | Phenylpropanoid |
| *trans*-Piperitol | 0.03 | Monoterpenic alcohol |
| *trans*-Carveol | 0.01 | Monoterpenic alcohol |
| 1-para-Menthen-9-al | 0.01 | Monoterpenic aldehyde |
| *cis*-para-Mentha-1(7),8-dien-2-ol | 0.03 | Monoterpenic alcohol |
| Nerol | 0.12 | Monoterpenic alcohol |
| Citronellol | 0.02 | Monoterpenic alcohol |
| Carvone | 0.03 | Monoterpenic ketone |
| Geraniol | 0.03 | Monoterpenic alcohol |
| Linalyl acetate | 0.05 | Monoterpenic ester |
| Geranial | 0.02 | Monoterpenic aldehyde |
| 4-Thujen-2α-yl acetate | 0.11 | Monoterpenic ester |
| Bornyl acetate | 0.52 | Monoterpenic ester |
| para-Cymen-7-ol | 0.03 | Monoterpenic alcohol |
| 2-Undecanone | 0.19 | Aliphatic ketone |
| *trans*-Pinocarvyl acetate | 0.01 | Monoterpenic ester |
| δ-Terpinyl acetate | 0.68 | Monoterpenic ester |
| Unknown | 0.05 | Unknown |
| para-Mentha-1,4-dien-7-ol | 0.01 | Monoterpenic alcohol |
| exo-2-Hydroxycineole acetate | 0.13 | Monoterpenic ester |
| α-Terpinyl acetate | 10.44 | Monoterpenic ester |
| α-Cubebene | 0.09 | Sesquiterpene |
| Eugenol | 0.64 | Phenylpropanoid |
| α-Ylangene | 0.07 | Sesquiterpene |
| Neryl acetate | 0.06 | Monoterpenic ester |
| α-Copaene | 0.05 | Sesquiterpene |
| β-Bourbonene | 0.02 | Sesquiterpene |

| | | |
|---|---|---|
| Geranyl acetate | 0.01 | Monoterpenic ester |
| β-Cubebene | 0.05 | Sesquiterpene |
| β-Elemene | 0.30 | Sesquiterpene |
| Methyleugenol | 0.65 | Phenylpropanoid |
| β-Caryophyllene | 0.32 | Sesquiterpene |
| β-Copaene | 0.01 | Sesquiterpene |
| Aromadendrene | 0.02 | Sesquiterpene |
| α-Guaiene | 0.04 | Sesquiterpene |
| 6,9-Guaiadiene | 0.04 | Sesquiterpene |
| (*E*)-Cinnamyl acetate | 0.06 | Phenylpropanoid ester |
| α-Humulene | 0.05 | Sesquiterpene |
| Selina-4(15),7-diene | 0.05 | Sesquiterpene |
| allo-Aromadendrene | 0.04 | Sesquiterpene |
| *cis*-Muurola-4(15),5-diene | 0.02 | Sesquiterpene |
| Unknown | 0.07 | Unknown |
| Germacrene D | 0.08 | Sesquiterpene |
| β-Selinene | 0.07 | Sesquiterpene |
| Viridiflorene | 0.05 | Sesquiterpene |
| α-Selinene | 0.05 | Sesquiterpene |
| Bicyclogermacrene | 0.05 | Sesquiterpene |
| (3*Z*,6*E*)-α-Farnesene | 0.05 | Sesquiterpene |
| Germacrene A | 0.04 | Sesquiterpene |
| δ-Amorphene | 0.02 | Sesquiterpene |
| γ-Cadinene | 0.13 | Sesquiterpene |
| δ-Cadinene | 0.24 | Sesquiterpene |
| *trans*-Cadina-1,4-diene | 0.01 | Sesquiterpene |
| α-Calacorene | 0.03 | Sesquiterpene |
| (*E*)-α-Bisabolene | 0.06 | Sesquiterpene |
| Elemicin | 0.02 | Phenylpropanoid |
| Germacrene D-4-ol | 0.01 | Sesquiterpenic alcohol |
| Spathulenol | 0.10 | Sesquiterpenic alcohol |
| Caryophyllene oxide | 0.10 | Sesquiterpenic ether |
| Globulol | 0.02 | Sesquiterpenic alcohol |
| Viridiflorol | 0.01 | Sesquiterpenic alcohol |
| Unknown | 0.02 | Oxygenated sesquiterpene |
| Ledol | 0.02 | Sesquiterpenic alcohol |
| Humulene epoxide II | 0.02 | Sesquiterpenic ether |
| Junenol | 0.03 | Sesquiterpenic alcohol |
| Eremoligenol? | 0.03 | Sesquiterpenic alcohol |
| Caryophylladienol II | 0.02 | Sesquiterpenic alcohol |
| τ-Cadinol | 0.02 | Sesquiterpenic alcohol |
| τ-Muurolol | 0.01 | Sesquiterpenic alcohol |
| β-Eudesmol | 0.03 | Sesquiterpenic alcohol |
| α-Eudesmol | 0.01 | Sesquiterpenic alcohol |
| (*E*)-Isoelemicin | 0.03 | Phenylpropanoid |
| Unknown | 0.01 | Unknown |
| Unknown | 0.03 | Aliphatic ester |
| Germacra-4(15),5,10(14)-trien-1-ol isomer | 0.02 | Sesquiterpenic alcohol |
| Gazaniolide | 0.03 | Sesquiterpenic lactone |
| **Consolidated total** | **98.94%** | |

*: 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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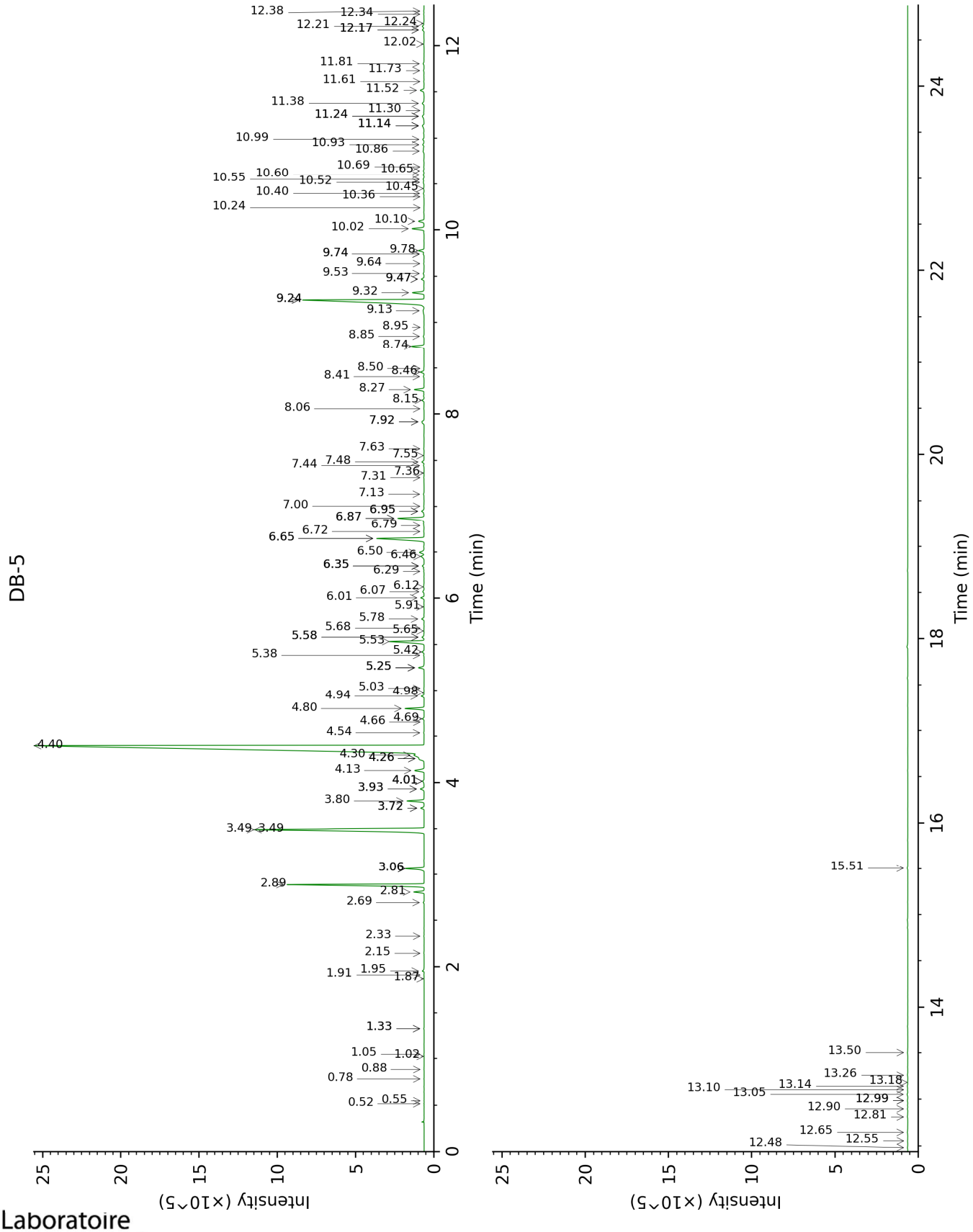
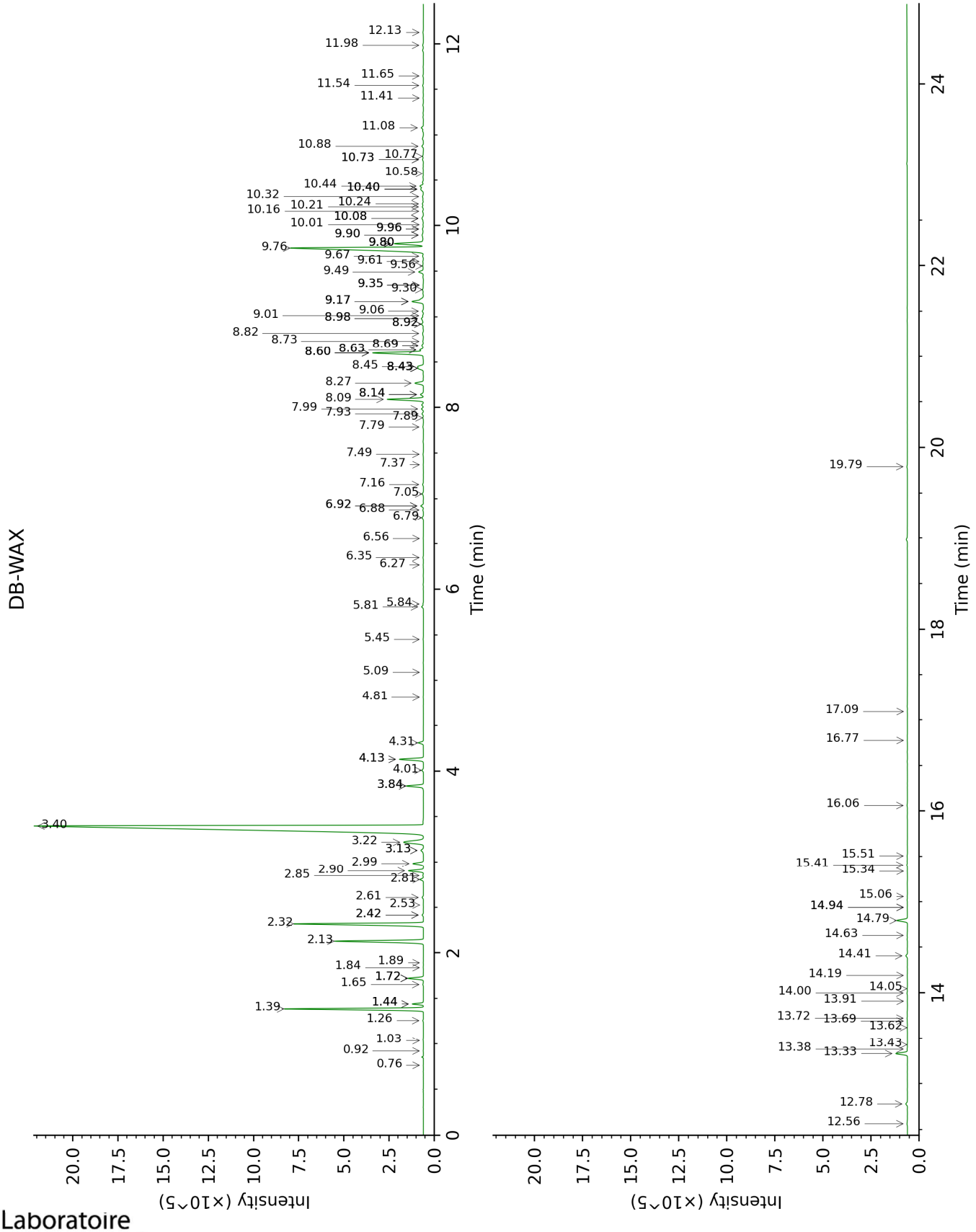

DB-5
Time (min)
Intensity (×10^5)
12.38
12.34
12.21
12.17
12.24
12.02
11.81
11.73
11.61
11.52
11.38
11.24
11.30
11.14
10.99
10.93
10.86
10.69
10.65
10.55
10.60
10.52
10.45
10.40
10.36
10.24
10.10
10.02
9.74
9.78
9.64
9.53
9.47
9.32
9.24
9.13
8.95
8.85
8.74
8.50
8.41
8.46
8.27
8.15
8.06
7.92
7.63
7.55
7.44
7.48
7.31
7.36
7.13
7.00
6.95
6.87
6.79
6.72
6.65
6.50
6.46
6.35
6.29
6.12
6.07
6.01
5.91
5.78
5.68
5.65
5.58
5.53
5.42
5.38
5.25
5.03
4.98
4.94
4.80
4.69
4.66
4.54
4.40
4.30
4.26
4.13
4.01
3.93
3.80
3.72
3.49
3.49
3.06
2.89
2.81
2.69
2.33
2.15
1.91
1.95
1.87
1.33
1.05
1.02
0.88
0.78
0.52
0.55
15.51
13.50
13.26
13.18
13.10
13.05
13.14
12.99
12.90
12.81
12.65
12.55
12.48


DB-WAX
Time (min)
Intensity (×10^5)
12.13
11.98
11.65
11.54
11.41
11.08
10.88
10.73
10.77
10.58
10.44
10.40
10.32
10.24
10.21
10.16
10.08
10.01
9.96
9.90
9.80
9.76
9.67
9.61
9.56
9.49
9.35
9.30
9.17
9.06
9.01
8.98
8.92
8.82
8.73
8.69
8.63
8.60
8.45
8.43
8.27
8.14
8.09
7.99
7.93
7.89
7.79
7.49
7.37
7.16
7.05
6.92
6.88
6.79
6.56
6.35
6.27
5.84
5.81
5.45
5.09
4.81
4.31
4.13
4.01
3.84
3.40
3.22
3.13
2.99
2.90
2.85
2.81
2.61
2.53
2.42
2.32
2.13
1.89
1.84
1.72
1.65
1.44
1.39
1.26
1.03
0.92
0.76
19.79
17.09
16.77
16.06
15.51
15.41
15.34
15.06
14.94
14.79
14.63
14.41
14.19
14.05
14.00
13.91
13.72
13.69
13.62
13.43
13.38
13.33
12.78
12.56

*FULL ANALYSIS DATA*

| Identification | Column DB-5 | | | Column DB-WAX | | |
|---|---|---|---|---|---|---|
| | R.T | R.I | % | R.T | R.I | % |
| Methylcyclopentadiene isomer I | 0.52 | 628 | tr | | | |
| Isovaleral | 0.55 | 640 | tr | 0.76 | 887 | tr |
| 3-Methyl-1-penten-3-ol | 0.78 | 716 | 0.01 | 2.53 | 1104 | 0.01 |
| 1-Methylpyrrole | 0.88 | 732 | tr | 2.42* | 1096 | 0.06 |
| Ethyl isobutyrate | 1.02 | 753 | tr | 1.03 | 936 | tr |
| Toluene | 1.05 | 757 | 0.01 | 1.44* | 999 | 0.47 |
| Hexanal | 1.33* | 801 | 0.01 | 1.89 | 1044 | 0.01 |
| Isopropyl isobutyrate | 1.33* | 801 | [0.01] | 0.92 | 917 | 0.01 |
| Ethyl 2-methylbutyrate | 1.87 | 850 | 0.01 | 1.72* | 1027 | 0.80 |
| Ethyl isovalerate | 1.91 | 853 | 0.01 | 1.84 | 1038 | 0.01 |
| (3*Z*)-Hexenol | 1.95 | 857 | 0.08 | 5.81 | 1348 | 0.09 |
| Hexanol | 2.15 | 874 | 0.01 | 5.45 | 1322 | 0.01 |
| Isopropyl 2-methylbutyrate | 2.33 | 890 | 0.01 | 1.72* | 1027 | [0.80] |
| Tricyclene | 2.70 | 918 | 0.04 | 1.26 | 972 | 0.03 |
| α-Thujene | 2.81 | 926 | 0.46 | 1.44* | 999 | [0.47] |
| α-Pinene | 2.89 | 931 | 6.72 | 1.39 | 993 | 6.69 |
| Camphene | 3.06* | 943 | 0.88 | 1.72* | 1027 | [0.80] |
| Thujadiene isomer | 3.06* | 943 | [0.88] | 2.42* | 1096 | [0.06] |
| α-Fenchene | 3.06* | 943 | [0.88] | 1.65 | 1020 | 0.03 |
| Sabinene | 3.49* | 972 | 12.16 | 2.32 | 1086 | 7.20 |
| β-Pinene | 3.49* | 972 | [12.16] | 2.13 | 1067 | 4.94 |
| 6-Methyl-5-hepten-2-one | 3.72* | 988 | 0.16 | 5.09 | 1301 | tr |
| Dehydro-1,8-cineole | 3.72* | 988 | [0.16] | 3.13* | 1152 | 0.17 |
| Myrcene | 3.80 | 993 | 0.81 | 2.90 | 1134 | 0.79 |
| α-Phellandrene | 3.93* | 1002 | 0.21 | 2.81 | 1126 | 0.18 |
| Pseudolimonene | 3.93* | 1002 | [0.21] | 2.85 | 1130 | 0.03 |
| Isobutyl 2-methylbutyrate | 4.02* | 1007 | 0.10 | 3.13* | 1152 | [0.17] |
| Δ3-Carene | 4.02* | 1007 | [0.10] | 2.61 | 1111 | 0.07 |
| (3*Z*)-Hexenyl acetate | 4.02* | 1007 | [0.10] | 4.81 | 1280 | 0.01 |
| α-Terpinene | 4.13 | 1015 | 0.58 | 2.99 | 1141 | 0.59 |
| Unknown [m/z 109, 43 (58), 95 (26)... 137 (15)...] | 4.26*† | 1023 | 50.47 | 6.27 | 1381 | 0.01 |
| para-Cymene | 4.26*† | 1023 | [50.47] | 4.13* | 1229 | 1.31 |
| Limonene | 4.30† | 1025 | [50.47] | 3.22 | 1159 | 1.70 |
| 1,8-Cineole | 4.40† | 1032 | [50.47] | 3.40 | 1173 | 47.43 |
| (*Z*)-β-Ocimene | 4.54 | 1040 | 0.02 | 3.84* | 1208 | 0.94 |
| Unknown [m/z 109, 43 (57), 91 (28), 67 (25), 93 (24), 95 (22), 77 (21), 137 (21), 41 (17), 79 (14)...] | 4.66 | 1048 | 0.01 | 7.37 | 1462 | 0.01 |
| (*E*)-β-Ocimene | 4.69 | 1050 | 0.08 | 4.01 | 1220 | 0.08 |
| γ-Terpinene | 4.80 | 1057 | 0.90 | 3.84* | 1208 | [0.94] |

| | | | | | | |
|---|---|---|---|---|---|---|
| *cis*-Sabinene hydrate | 4.94 | 1066 | 0.14 | 6.92* | 1429 | 0.15 |
| para-Mentha-3,8-diene | 4.98 | 1068 | 0.01 | 4.13* | 1229 | [1.31] |
| *cis*-Linalool oxide (fur.) | 5.03 | 1071 | 0.01 | 6.56 | 1402 | 0.01 |
| Terpinolene | 5.25* | 1086 | 0.30 | 4.31 | 1243 | 0.27 |
| para-Cymenene | 5.25* | 1086 | [0.30] | 6.35 | 1387 | 0.03 |
| *trans*-Linalool oxide (fur.) | 5.25* | 1086 | [0.30] | 6.92* | 1429 | [0.15] |
| 2-Nonanone | 5.38 | 1094 | 0.02 | 5.84 | 1350 | 0.02 |
| *trans*-Sabinene hydrate | 5.42 | 1096 | 0.11 | 7.98 | 1508 | 0.11 |
| Linalool | 5.53 | 1104 | 1.98 | 8.09 | 1517 | 1.98 |
| Unknown [m/z 43, 59 (37), 79 (33), 91 (32), 119 (31)...] | 5.58* | 1107 | 0.13 | 9.06 | 1592 | 0.06 |
| Hotrienol | 5.58* | 1107 | [0.13] | 8.82 | 1573 | 0.06 |
| endo-Fenchol | 5.65 | 1111 | 0.02 | 8.43* | 1543 | 0.32 |
| Unknown [m/z 93, 121 (64), 91 (58), 77 (58), 94 (53), 79 (47)...] | 5.68 | 1113 | 0.01 | | | |
| *cis*-para-Menth-2-en-1-ol | 5.78 | 1119 | 0.13 | 8.14* | 1521 | 0.18 |
| Limona ketone | 5.91 | 1128 | 0.01 | 7.89 | 1501 | 0.03 |
| *trans*-Pinocarveol | 6.01 | 1134 | 0.17 | 9.17* | 1600 | 0.88 |
| *trans*-para-Menth-2-en-1-ol | 6.07 | 1138 | 0.08 | 8.98* | 1586 | 0.17 |
| *trans*-Verbenol | 6.12 | 1142 | 0.05 | 9.56 | 1632 | 0.04 |
| Sabinaketone | 6.29 | 1152 | 0.02 | 8.73 | 1566 | 0.04 |
| Nerol oxide | 6.35* | 1156 | 0.10 | 6.88 | 1426 | 0.03 |
| (*E*)-2,6-Dimethyl-1,5,7-octatrien-3-ol | 6.35* | 1156 | [0.10] | 10.32 | 1694 | 0.05 |
| Pinocarvone | 6.35* | 1156 | [0.10] | 7.93 | 1504 | 0.09 |
| Borneol | 6.46 | 1163 | 0.22 | 9.80* | 1652 | 1.75 |
| δ-Terpineol | 6.50 | 1166 | 0.32 | 9.49 | 1627 | 0.31 |
| Rosefuran oxide | 6.65* | 1175 | 3.01 | 8.60* | 1556 | 3.04 |
| Terpinen-4-ol | 6.65* | 1175 | [3.01] | 8.60* | 1556 | [3.04] |
| Thuj-3-en-10-al | 6.72 | 1180 | 0.02 | 8.69* | 1563 | 0.10 |
| para-Cymen-8-ol | 6.79 | 1184 | 0.01 | 11.54 | 1798 | 0.04 |
| α-Terpineol | 6.87* | 1190 | 1.58 | 9.80* | 1652 | [1.75] |
| Myrtenal | 6.87* | 1190 | [1.58] | 8.69* | 1563 | [0.10] |
| Myrtenol | 6.95* | 1195 | 0.15 | 10.88 | 1741 | 0.09 |
| *cis*-Piperitol | 6.95* | 1195 | [0.15] | 9.61 | 1636 | 0.04 |
| Methylchavicol | 7.00 | 1198 | 0.01 | 9.35* | 1615 | 0.02 |
| *trans*-Piperitol | 7.14 | 1207 | 0.03 | 10.40* | 1701 | 0.20 |
| *trans*-Carveol | 7.31 | 1219 | 0.01 | 11.41 | 1786 | 0.02 |
| 1-para-Menthen-9-al | 7.36 | 1222 | 0.01 | 8.92* | 1581 | 0.11 |
| *cis*-para-Mentha-1(7),8-dien-2-ol | 7.44 | 1227 | 0.03 | 11.98 | 1837 | 0.05 |
| Nerol | 7.48 | 1230 | 0.12 | 11.08 | 1758 | 0.13 |
| Citronellol | 7.55 | 1235 | 0.02 | 10.77 | 1732 | 0.05 |
| Carvone | 7.63 | 1240 | 0.03 | 10.01 | 1669 | 0.02 |
| Geraniol | 7.92* | 1259 | 0.19 | 11.65 | 1807 | 0.03 |
| Linalyl acetate | 7.92* | 1259 | [0.19] | 8.14* | 1521 | [0.18] |
| Geranial | 8.06 | 1269 | 0.02 | 10.16 | 1681 | 0.04 |

| | | | | | | |
|---|---|---|---|---|---|---|
| 4-Thujen-2α-yl acetate | 8.15 | 1275 | 0.11 | 8.92* | 1581 | [0.11] |
| Bornyl acetate | 8.27 | 1282 | 0.52 | 8.27 | 1530 | 0.52 |
| para-Cymen-7-ol | 8.41 | 1292 | 0.03 | 14.19 | 2040 | 0.02 |
| 2-Undecanone | 8.46 | 1295 | 0.19 | 8.63* | 1559 | 0.19 |
| *trans*-Pinocarvyl acetate | 8.50 | 1298 | 0.01 | 9.17* | 1600 | [0.88] |
| δ-Terpinyl acetate | 8.74 | 1315 | 0.68 | 9.17* | 1600 | [0.88] |
| Unknown [m/z 119, 43 (99), 93 (52), 59 (44), 91 (41), 134 (34)...] | 8.85 | 1323 | 0.05 | 9.90* | 1660 | 0.15 |
| para-Mentha-1,4-dien-7-ol | 8.95 | 1330 | 0.01 | 13.72 | 1995 | 0.03 |
| exo-2-Hydroxycineole acetate | 9.13 | 1342 | 0.13 | 10.08* | 1674 | 0.12 |
| α-Terpinyl acetate | 9.24* | 1350 | 10.49 | 9.76 | 1648 | 10.44 |
| α-Cubebene | 9.24* | 1350 | [10.49] | 6.79 | 1419 | 0.09 |
| Eugenol | 9.32 | 1356 | 0.64 | 14.79 | 2098 | 0.61 |
| α-Ylangene | 9.47* | 1366 | 0.16 | 7.05 | 1439 | 0.07 |
| Neryl acetate | 9.47* | 1366 | [0.16] | 10.21 | 1685 | 0.06 |
| α-Copaene | 9.53 | 1371 | 0.05 | 7.16 | 1446 | 0.05 |
| β-Bourbonene | 9.64 | 1378 | 0.02 | 7.49 | 1471 | 0.03 |
| Geranyl acetate | 9.74* | 1386 | 0.05 | 10.58 | 1715 | 0.01 |
| β-Cubebene | 9.74* | 1386 | [0.05] | 7.79 | 1493 | 0.05 |
| β-Elemene | 9.78 | 1388 | 0.30 | 8.43* | 1543 | [0.32] |
| Methyleugenol | 10.02 | 1405 | 0.65 | 13.33 | 1958 | 0.66 |
| β-Caryophyllene | 10.10 | 1411 | 0.32 | 8.45 | 1544 | 0.33 |
| β-Copaene | 10.24 | 1422 | 0.01 | 8.43* | 1543 | [0.32] |
| Aromadendrene | 10.36 | 1431 | 0.02 | 8.60* | 1556 | [3.04] |
| α-Guaiene | 10.40 | 1434 | 0.04 | 8.43* | 1543 | [0.32] |
| 6,9-Guaiadiene | 10.45 | 1438 | 0.04 | 8.63* | 1559 | [0.19] |
| (*E*)-Cinnamyl acetate | 10.52 | 1443 | 0.06 | 14.63 | 2082 | 0.02 |
| α-Humulene | 10.55 | 1445 | 0.05 | 9.30 | 1611 | 0.06 |
| Selina-4(15),7-diene | 10.60 | 1449 | 0.05 | 8.98* | 1586 | [0.17] |
| allo-Aromadendrene | 10.65 | 1452 | 0.04 | 9.01 | 1588 | 0.03 |
| *cis*-Muurola-4(15),5-diene | 10.69 | 1456 | 0.02 | 9.35* | 1615 | [0.02] |
| Unknown [m/z 43, 67 (61), 79 (57), 81 (44), 54 (44)...] | 10.86 | 1468 | 0.07 | | | |
| Germacrene D | 10.93 | 1474 | 0.08 | 9.80* | 1652 | [1.75] |
| β-Selinene | 10.99 | 1478 | 0.07 | 9.90* | 1660 | [0.15] |
| Viridiflorene | 11.14* | 1489 | 0.15 | 9.67 | 1641 | 0.05 |
| α-Selinene | 11.14* | 1489 | [0.15] | 9.96* | 1665 | 0.07 |
| Bicyclogermacrene | 11.14* | 1489 | [0.15] | 10.08* | 1674 | [0.12] |
| (3*Z*,6*E*)-α-Farnesene | 11.24* | 1497 | 0.16 | 10.24 | 1687 | 0.05 |
| Germacrene A | 11.24* | 1497 | [0.16] | 10.40* | 1701 | [0.20] |
| δ-Amorphene | 11.30 | 1501 | 0.02 | 9.96* | 1665 | [0.07] |
| γ-Cadinene | 11.38 | 1507 | 0.13 | 10.40* | 1701 | [0.20] |
| δ-Cadinene | 11.52 | 1518 | 0.24 | 10.44 | 1704 | 0.21 |
| *trans*-Cadina-1,4-diene | 11.61 | 1526 | 0.01 | 10.73* | 1729 | 0.05 |
| α-Calacorene | 11.73 | 1535 | 0.03 | 12.13 | 1849 | 0.04 |
| (*E*)-α-Bisabolene | 11.81 | 1541 | 0.06 | 10.73* | 1729 | [0.05] |

| | | | | | | |
|---|---|---|---|---|---|---|
| Elemicin | 12.02 | 1558 | 0.02 | 15.51 | 2170 | 0.02 |
| Germacrene D-4-ol | 12.17* | 1570 | 0.10 | 13.69 | 1992 | 0.01 |
| Spathulenol | 12.17* | 1570 | [0.10] | 14.41 | 2061 | 0.10 |
| Caryophyllene oxide | 12.21 | 1573 | 0.10 | 12.78 | 1908 | 0.10 |
| Globulol | 12.24 | 1575 | 0.02 | 13.91 | 2013 | 0.01 |
| Viridiflorol | 12.34 | 1583 | 0.01 | 14.00 | 2021 | 0.03 |
| Unknown [m/z 133, 93 (64), 43 (64), 177 (60), 107 (59), 91 (55)...220 (7)] | 12.38 | 1586 | 0.02 | 12.56 | 1888 | 0.03 |
| Ledol | 12.48 | 1594 | 0.02 | 13.38 | 1963 | 0.02 |
| Humulene epoxide II | 12.55 | 1600 | 0.02 | 13.43 | 1968 | 0.01 |
| Junenol | 12.65 | 1607 | 0.03 | 13.62 | 1985 | 0.01 |
| Eremoligenol? | 12.81 | 1621 | 0.03 | 14.94* | 2112 | 0.01 |
| Caryophylladienol II | 12.90 | 1628 | 0.02 | 16.06 | 2226 | 0.02 |
| τ-Cadinol | 12.99* | 1635 | 0.03 | 14.94* | 2112 | [0.01] |
| τ-Muurolol | 12.99* | 1635 | [0.03] | 15.06 | 2124 | 0.01 |
| β-Eudesmol | 13.05 | 1641 | 0.03 | 15.41 | 2159 | 0.04 |
| α-Eudesmol | 13.10 | 1645 | 0.01 | 15.34 | 2152 | 0.01 |
| (*E*)-Isoelemicin | 13.14 | 1648 | 0.03 | 17.09 | 2335 | 0.01 |
| Unknown [m/z 41, 79 (99), 91 (88), 69 (82), 93 (81), 55 (74), 95 (65)...] | 13.18 | 1652 | 0.01 | | | |
| Unknown [m/z 79, 43 (66), 67 (59), 80 (56), 41 (41), 81 (37), 55 (29)...] | 13.26 | 1658 | 0.03 | 14.05 | 2026 | 0.02 |
| Germacra-4(15),5,10(14)-trien-1-ol isomer | 13.50 | 1678 | 0.02 | 16.78 | 2301 | tr |
| Gazaniolide | 15.51 | 1853 | 0.03 | 19.79 | 2641 | 0.03 |
| **Total identified** | | **98.98%** | | | **98.60%** | |
| **Total reported** | | **99.17%** | | | **98.73%** | |

*: 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