**Date :** 2025-04-22

# *Certificate of Analysis - GC Profiling*

## *Sample Identification*

**Internal code :** 25D11-PTH04
**Customer Identification :** Organic Frankincense Frereana - Somalia - FE0112R
**Type :** Essential Oil
**Source :** *Boswellia frereana*
**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*Notes: This report may not be published, including online, without the written consent from Laboratoire PhytoChemia.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 :** 2025-04-15

## *Physicochemical data*

**Refractive index :** 1.4622 ± 0.0003 (20 °C)
**Method :** PC-MAT-016 - Measure of the refractive index of a liquid.
**Analyst :** Cindy Caron B. Sc.
**Date :** 2025-04-11

## *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 |
|---|---|---|
| 3-Methyl-2-butanone | 0.03 | Aliphatic ketone |
| Unknown | tr | Unknown |
| 4-Methyl-2-pentanone (Methyl isobutyl ketone) | 0.01 | Aliphatic ketone |
| Toluene | 0.03 | Simple phenolic |
| Hexadiene isomer? | 0.01 | Alkene |
| (*Z*)-Salvene | 0.07 | Normonoterpene |
| Unknown | 0.01 | Unknown |
| (*E*)-Salvene | 0.03 | Normonoterpene |
| Unknown | 0.02 | Unknown |
| Unknown | 0.08 | Monoterpene |
| Unknown | 0.21 | Unknown |
| Hashishene | 0.51 | Monoterpene |
| α-Thujene | 59.84 | Monoterpene |
| α-Pinene | 2.38 | Monoterpene |
| Unknown | 2.52 | Monoterpene |
| Camphene | 0.03 | Monoterpene |
| α-Fenchene | tr | Monoterpene |
| Thuja-2,4(10)-diene | 0.02 | Monoterpene |
| Sabinene | 6.35 | Monoterpene |
| β-Pinene | 0.27 | Monoterpene |
| Pseudolimonene isomer | 0.05 | Monoterpene |
| 6-Methyl-5-hepten-2-one | 0.06 | Aliphatic ketone |
| Dehydro-1,8-cineole | 0.02 | Monoterpenic ether |
| Myrcene | 0.23 | Monoterpene |
| 6-Methyl-5-hepten-2-ol | 0.13 | Aliphatic alcohol |
| Pseudolimonene | 0.03 | Monoterpene |
| α-Phellandrene | 0.90 | Monoterpene |
| *ortho*-Methylanisole | 0.02 | Simple phenolic |
| Δ3-Carene | 0.05 | Monoterpene |
| α-Terpinene | 0.44 | Monoterpene |
| Carvomenthene | 0.05 | Aliphatic alcohol |
| *meta*-Cymene | 0.46 | Monoterpene |
| *para*-Cymene | 9.95 | Monoterpene |
| Limonene | 0.49 | Monoterpene |
| β-Phellandrene | [3.34] | Monoterpene |
| 1,8-Cineole | [3.34] | Monoterpenic ether |
| Unknown | 0.04 | Unknown |
| (*Z*)-β-Ocimene | 0.04 | Monoterpene |
| Unknown | 0.25 | Unknown |

| | | |
|---|---|---|
| (*E*)-β-Ocimene | 0.02 | Monoterpene |
| Unknown | 0.03 | Unknown |
| γ-Terpinene | 0.61 | Monoterpene |
| *cis*-Sabinene hydrate | 0.05 | Monoterpenic alcohol |
| Unknown | 0.05 | Oxygenated monoterpene |
| Terpinolene | 0.15 | Monoterpene |
| *para*-Cymenene | 0.06 | Monoterpene |
| *trans*-Sabinene hydrate | 0.04 | Monoterpenic alcohol |
| Unknown | 0.04 | Unknown |
| α-Thujone | 0.08 | Monoterpenic ketone |
| Linalool | 0.08 | Monoterpenic alcohol |
| β-Thujone | 0.12 | Monoterpenic ketone |
| Unknown | 0.60 | Oxygenated monoterpene |
| Unknown | 0.05 | Oxygenated monoterpene |
| *cis-para*-Menth-2-en-1-ol | 0.14 | Monoterpenic alcohol |
| *trans-para*-Mentha-2,8-dien-1-ol | 0.16 | Monoterpenic alcohol |
| Unknown | 0.06 | Unknown |
| *trans*-Pinocarveol | 0.06 | Monoterpenic alcohol |
| *cis*-Verbenol | 0.05 | Monoterpenic alcohol |
| *trans*-Sabinol | 0.24 | Monoterpenic alcohol |
| *trans*-Verbenol | 0.03 | Monoterpenic alcohol |
| *meta*-Mentha-4,6-dien-8-ol | 0.02 | Monoterpenic alcohol |
| Sabinaketone | 0.05 | Normonoterpenic ketone |
| Unknown | 0.03 | Oxygenated monoterpene |
| Unknown | 0.06 | Oxygenated monoterpene |
| α-Phellandren-8-ol | 0.03 | Monoterpenic alcohol |
| Umbellulone | 0.10 | Monoterpenic ketone |
| Unknown | 0.02 | Oxygenated monoterpene |
| *cis*-Sabinol | 0.23 | Monoterpenic alcohol |
| Terpinen-4-ol | 3.02 | Monoterpenic alcohol |
| Cryptone | 0.02 | Normonoterpenic ketone |
| Thuj-3-en-10-al | 0.16 | Monoterpenic aldehyde |
| *para*-Cymen-8-ol | 0.19 | Monoterpenic alcohol |
| α-Terpineol | 0.03 | Monoterpenic alcohol |
| Myrtenal | 0.02 | Monoterpenic aldehyde |
| *cis*-Piperitol | 0.03 | Monoterpenic alcohol |
| Myrtenol | 0.05 | Monoterpenic alcohol |
| *cis*-α-Phellandrene epoxide (iPr vs Me) | 0.10 | Monoterpenic ether |
| *trans*-Piperitol | 0.06 | Monoterpenic alcohol |
| *trans*-Carveol | 0.02 | Monoterpenic alcohol |
| Citronellol | 0.04 | Monoterpenic alcohol |
| Cuminal | 0.06 | Monoterpenic aldehyde |
| Carvone | 0.01 | Monoterpenic ketone |
| Carvotanacetone | 0.07 | Monoterpenic ketone |
| Unknown | 0.05 | Unknown |

| | | |
|---|---|---|
| Unknown | 0.05 | Oxygenated monoterpene |
| Piperitone | 0.03 | Monoterpenic ketone |
| Unknown | 0.02 | Unknown |
| 3,5-Dimethoxytoluene | 0.01 | Simple phenolic |
| Unknown | 0.03 | Unknown |
| Bornyl acetate | 0.07 | Monoterpenic ester |
| Cuminol | 0.04 | Monoterpenic alcohol |
| Thymol | 0.08 | Monoterpenic alcohol |
| Carvacrol | 0.06 | Monoterpenic alcohol |
| exo-2-Hydroxycineole acetate | 0.02 | Monoterpenic ester |
| α-Cubebene | 0.02 | Sesquiterpene |
| Citronellyl acetate | 0.05 | Monoterpenic ester |
| α-Copaene | 0.05 | Sesquiterpene |
| β-Bourbonene | 1.13 | Sesquiterpene |
| β-Elemene | 0.06 | Sesquiterpene |
| β-Caryophyllene | 0.06 | Sesquiterpene |
| β-Ylangene | 0.08 | Sesquiterpene |
| β-Copaene | 0.13 | Sesquiterpene |
| Isogermacrene D | 0.08 | Sesquiterpene |
| α-Humulene | 0.02 | Sesquiterpene |
| allo-Aromadendrene | 0.02 | Sesquiterpene |
| Germacrene D | 0.03 | Sesquiterpene |
| γ-Cadinene | 0.02 | Sesquiterpene |
| δ-Cadinene | 0.02 | Sesquiterpene |
| Caryophyllene oxide | 0.01 | Sesquiterpenic ether |
| τ-Cadinol | 0.03 | Sesquiterpenic alcohol |
| α-Phellandrene dimer I | 0.03 | Diterpene |
| α-Phellandrene dimer II | 0.44 | Diterpene |
| α-Phellandrene dimer III | 0.06 | Diterpene |
| α-Phellandrene dimer IV | 0.06 | Diterpene |
| α-Phellandrene dimer VI | 0.02 | Diterpene |
| α-Phellandrene dimer VII | 0.02 | Diterpene |
| **Consolidated total** | **99.03** | |

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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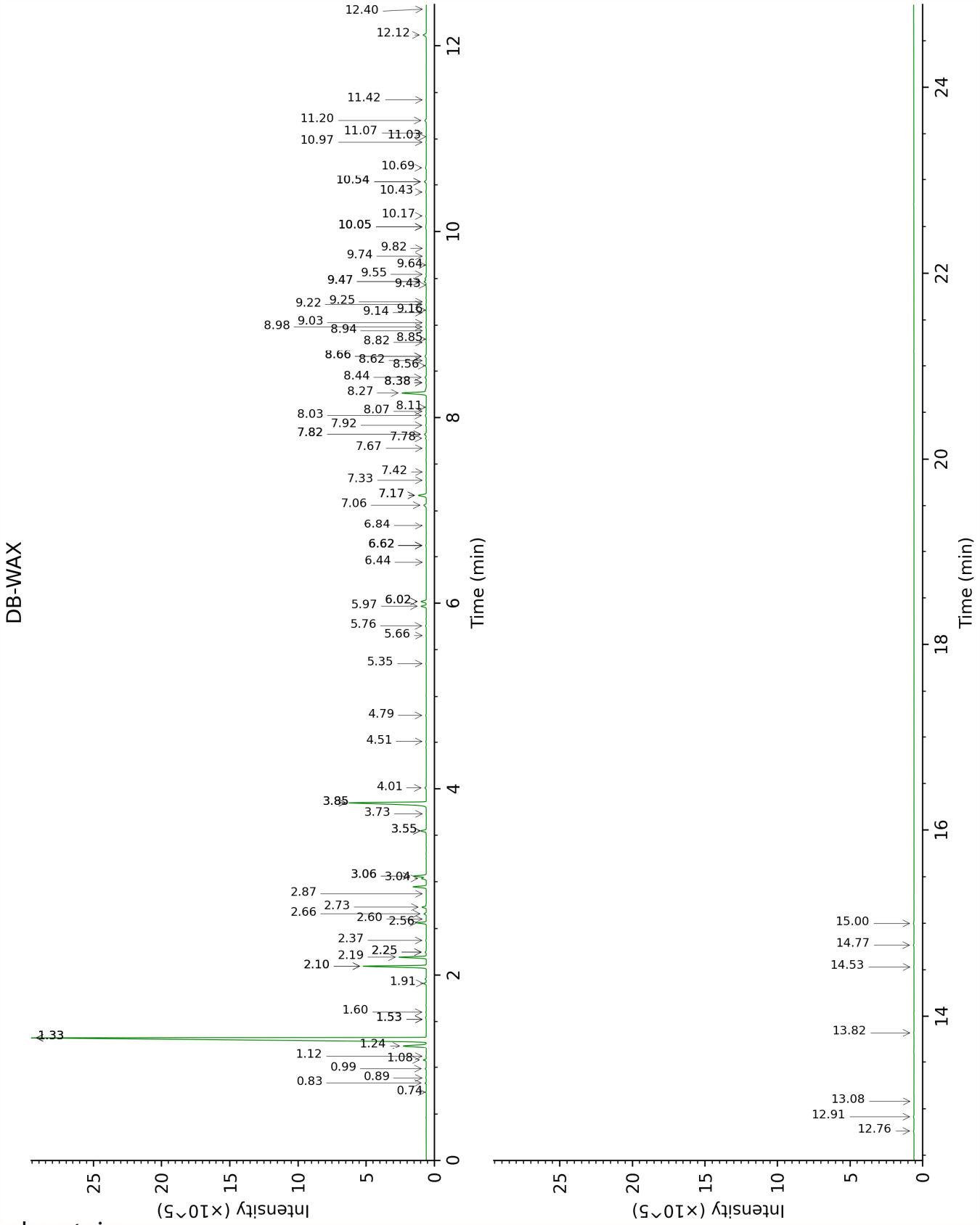
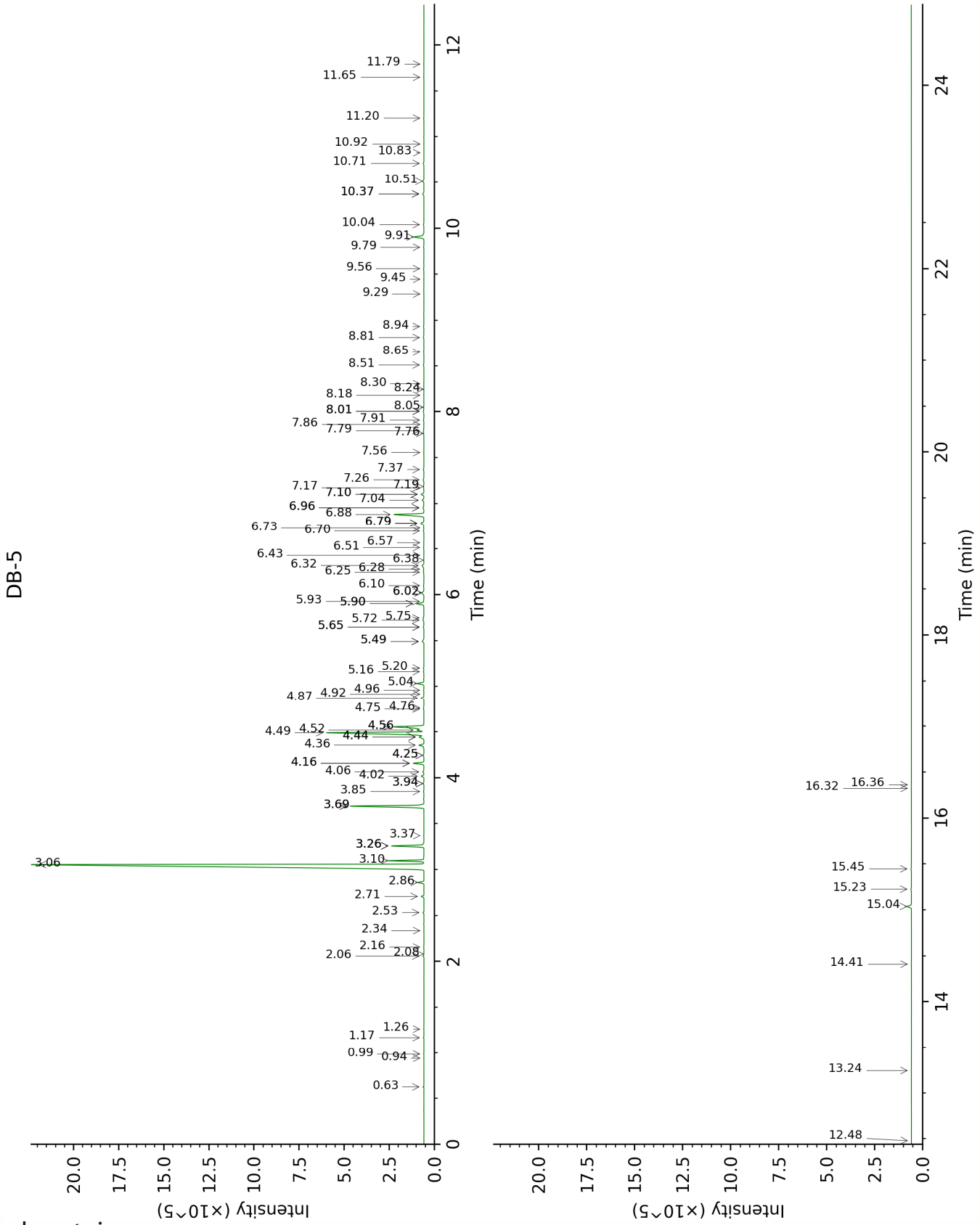

DB-WAX
Time (min)
Intensity (×10^5)
12.40
12.12
11.42
11.20
11.07
11.03
10.97
10.69
10.54
10.43
10.17
10.05
9.82
9.74
9.64
9.55
9.47
9.43
9.25
9.22
9.14
9.16
9.03
8.98
8.94
8.85
8.82
8.66
8.62
8.56
8.44
8.38
8.27
8.11
8.07
8.03
7.92
7.82
7.78
7.67
7.42
7.33
7.17
7.06
6.84
6.62
6.44
6.02
5.97
5.76
5.66
5.35
4.79
4.51
4.01
3.85
3.73
3.55
3.06
3.04
2.87
2.73
2.66
2.60
2.56
2.37
2.25
2.19
2.10
1.91
1.60
1.53
1.33
1.24
1.12
1.08
0.99
0.89
0.83
0.74
15.00
14.77
14.53
13.82
13.08
12.91
12.76


DB-5
Time (min)
Intensity (×10^5)
11.79
11.65
11.20
10.92
10.83
10.71
10.51
10.37
10.04
9.91
9.79
9.56
9.45
9.29
8.94
8.81
8.65
8.51
8.30
8.24
8.18
8.05
8.01
7.91
7.86
7.79
7.76
7.56
7.37
7.26
7.19
7.17
7.10
7.04
6.96
6.88
6.79
6.73
6.70
6.57
6.51
6.43
6.38
6.32
6.28
6.25
6.10
6.02
5.93
5.90
5.75
5.72
5.65
5.49
5.20
5.16
5.04
4.96
4.92
4.87
4.76
4.75
4.56
4.52
4.49
4.44
4.36
4.25
4.16
4.06
4.02
3.94
3.85
3.69
3.37
3.26
3.10
3.06
2.86
2.71
2.53
2.34
2.16
2.08
2.06
1.26
1.17
0.99
0.94
0.63
16.36
16.32
15.45
15.23
15.04
14.41
13.24
12.48

*FULL ANALYSIS DATA*

| | Column DB-WAX | | | Column DB-5 | | |
|---|---|---|---|---|---|---|
| **3-Methyl-2-butanone** | 0.74 | 894.9 | 0.04 | 0.63 | 640.8 | 0.03 |
| Unknown BODA I [m/z 93, 91 (70), 77 (48), 108 (42)] | | | | 0.94 | 722.1 | tr |
| 4-Methyl-2-pentanone (Methyl isobutyl ketone) | 1.12 | 965.2 | 0.02 | 0.99 | 728.8 | 0.01 |
| Toluene | 1.33* | 999.4 | [59.86] | 1.17 | 755.2 | 0.03 |
| Hexadiene isomer? | | | | 1.26 | 768.7 | 0.01 |
| (*Z*)-Salvene | 0.84 | 913.0 | 0.07 | 2.06 | 849.6 | 0.07 |
| Unknown PRME II [m/z 109, 43 (28), 124 (28), 41 (14), 55 (11), 79 (9), 81 (8)...] | 1.53* | 1020.3 | [0.02] | 2.08 | 851.6 | 0.01 |
| (*E*)-Salvene | 0.89 | 922.6 | 0.03 | 2.16 | 857.9 | 0.03 |
| Unknown BOCA II [m/z 79, 78 (45), 91 (28), 77 (28), 41 (13), 80 (12), 107 (11)... 122 (1)] | 1.08 | 957.9 | 0.23 | 2.34 | 873.0 | 0.02 |
| Unknown BOFR I [m/z 93, 91 (75), 121 (61), 77 (58), 79 (38), 92 (26), 43 (24), 41 (23), 105 (22),107 (19), 136 (16)] | 0.99 | 941.0 | 0.08 | 2.53 | 889.6 | 0.08 |
| Unknown BOFR II [m/z 93, 91 (72), 121 (58), 77 (49), 79 (41), 43 (22), 105 (20),107 (20), 41 (18), 136 (17), 92 (17)] | | | | 2.71 | 904.1 | 0.21 |
| Hashishene | 1.33* | 999.4 | [59.86] | 2.86 | 914.3 | 0.51 |
| α-Thujene | 1.33* | 999.4 | [59.86] | 3.06 | 927.3 | 59.84 |
| α-Pinene | 1.24 | 986.6 | 2.88 | 3.10 | 930.1 | 2.38 |
| Unknown SAOF I [m/z 91, 92 (47), 65 (11)... 134 (1)] | 2.19 | 1090.5 | 2.52 | 3.26* | 940.9 | [2.54] |

| | | | | | | |
|---|---|---|---|---|---|---|
| Camphene | 1.60 | 1028.4 | 0.03 | 3.26* | 940.9 | [2.54] |
| α-Fenchene | 1.53* | 1020.3 | [0.02] | 3.26* | 940.9 | [2.54] |
| Thuja-2,4(10)-diene | 2.10* | 1080.3 | [6.35] | 3.37 | 948.4 | 0.02 |
| Sabinene | 2.10* | 1080.3 | [6.35] | 3.69* | 970.0 | [6.62] |
| β-Pinene | 1.91 | 1060.9 | 0.27 | 3.69* | 970.0 | [6.62] |
| Pseudolimonene isomer | 2.25* | 1096.1 | [0.10] | 3.85 | 980.6 | 0.05 |
| 6-Methyl-5-hepten-2-one | 4.79 | 1292.4 | 0.06 | 3.94* | 986.4 | [0.07] |
| Dehydro-1,8-cineole | 2.87 | 1146.7 | 0.02 | 3.94* | 986.4 | [0.07] |
| Myrcene | 2.66 | 1129.5 | 0.23 | 4.02 | 991.8 | 0.23 |
| 6-Methyl-5-hepten-2-ol | 6.62* | 1427.9 | [0.06] | 4.06 | 994.9 | 0.13 |
| Pseudolimonene | 2.60 | 1124.9 | 0.03 | 4.16* | 1001.2 | [0.93] |
| α-Phellandrene | 2.56 | 1121.6 | 0.90 | 4.16* | 1001.2 | [0.93] |
| *ortho*-Methylanisole | 5.66 | 1356.0 | 0.02 | 4.24* | 1006.7 | [0.09] |
| Δ3-Carene | 2.37 | 1106.4 | 0.05 | 4.24* | 1006.7 | [0.09] |
| α-Terpinene | 2.73 | 1135.2 | 0.43 | 4.36 | 1013.7 | 0.44 |
| Carvomenthene | 2.25* | 1096.1 | [0.10] | 4.44* | 1019.3 | [0.56] |
| *meta*-Cymene | 3.85* | 1223.2 | [10.41] | 4.44* | 1019.3 | [0.56] |
| *para*-Cymene | 3.85* | 1223.2 | [10.41] | 4.49 | 1022.3 | 9.95 |
| Limonene | 3.04 | 1160.1 | 0.49 | 4.52*† | 1024.1 | [0.60] |
| β-Phellandrene | 3.06* | 1162.1 | [1.36] | 4.56*† | 1026.4 | [3.23] |
| 1,8-Cineole | 3.06* | 1162.1 | [1.36] | 4.56*† | 1026.4 | [3.23] |
| Unknown ARAN I [m/z 43, 55 (65), 41 (34), 67 (32), 107 (30), 122 (26)... 125 (10)] | 5.35 | 1333.9 | 0.04 | 4.75 | 1038.7 | 0.04 |
| (*Z*)-β-Ocimene | 3.55* | 1201.0 | [0.64] | 4.76 | 1039.5 | 0.04 |
| Unknown BOFR III [m/z 109, 43 (57), 91 (28), 67 (25), 93 (24), 95 (22), 77 (21), 137 (21), 41 (17), 79 (14)...] | 7.06 | 1461.2 | 0.34 | 4.87 | 1046.5 | 0.25 |
| (*E*)-β-Ocimene | 3.73 | 1214.4 | 0.02 | 4.92 | 1049.3 | 0.02 |
| Unknown BOFR IV [m/z 109 , 45 (67), 41 (40), 67 (39), 81 (33), 79 (27), 95 (24), 91 (23), 82 | 6.62* | 1427.9 | [0.06] | 4.96 | 1052.0 | 0.03 |

| | | | | | | |
|---|---|---|---|---|---|---|
| (21), 55 (21), 93 (20)...] | | | | | | |
| γ-Terpinene | 3.55* | 1201.0 | [0.64] | 5.04 | 1056.7 | 0.61 |
| *cis*-Sabinene hydrate | 6.62* | 1427.9 | [0.06] | 5.16 | 1064.8 | 0.05 |
| Unknown PIMA I [m/z 79, 93 (60), 43 (40), 94 (35), 137 (33), 77 (26), 91 (20), 152 (18)] | 4.51 | 1271.8 | 0.04 | 5.20 | 1067.3 | 0.05 |
| Terpinolene | 4.01 | 1235.0 | 0.15 | 5.49* | 1085.5 | [0.21] |
| *para*-Cymenene | 6.02* | 1382.9 | [0.64] | 5.49* | 1085.5 | [0.21] |
| *trans*-Sabinene hydrate | 7.67 | 1507.7 | 0.04 | 5.65* | 1095.4 | [0.07] |
| Unknown BOFR VII [m/z 109, 43 (48), 91 (39), 81 (33), 95 (32), 93 (32), 67 (31)... 152 (9)] | | | | 5.65* | 1095.4 | [0.07] |
| α-Thujone | 5.76 | 1363.7 | 0.08 | 5.72*† | 1100.2 | [0.10] |
| Linalool | 7.78 | 1516.2 | 0.08 | 5.75*† | 1101.8 | [0.08] |
| β-Thujone | 6.02* | 1382.9 | [0.64] | 5.90* | 1111.8 | [0.72] |
| Unknown BOSE I [m/z 109, 81 (54), 91 (32), 79 (22)...] | 5.97 | 1379.2 | 0.60 | 5.90* | 1111.8 | [0.72] |
| Unknown BOSE II [m/z 109, 91 (57), 93 (47), 81 (44), 77 (40)... 154 (1)] | | | | 5.93 | 1113.3 | 0.05 |
| *cis-para*-Menth-2-en-1-ol | 7.82* | 1519.6 | [0.23] | 6.02* | 1119.4 | [0.29] |
| *trans-para*-Mentha-2,8-dien-1-ol | 8.66* | 1585.5 | [0.18] | 6.02* | 1119.4 | [0.29] |
| Unknown BOSE III [m/z 111, 43 (22), 55 (14), 41 (12), 110 (11)...] | | | | 6.10 | 1124.3 | 0.06 |
| *trans*-Pinocarveol | 8.82 | 1597.8 | 0.04 | 6.25 | 1133.7 | 0.06 |
| *cis*-Verbenol | 8.98 | 1611.4 | 0.06 | 6.28 | 1135.8 | 0.05 |
| *trans*-Sabinol | 9.47* | 1651.3 | [0.27] | 6.32 | 1138.4 | 0.24 |
| *trans*-Verbenol | 9.16 | 1626.2 | 0.03 | 6.38 | 1142.1 | 0.03 |
| *meta*-Mentha-4,6-dien-8-ol | 9.03 | 1615.1 | 0.04 | 6.43 | 1145.5 | 0.02 |
| Sabinaketone | 8.38* | 1563.3 | [0.07] | 6.52 | 1151.0 | 0.05 |

| | | | | | | |
|---|---|---|---|---|---|---|
| Unknown BOCA III [m/z 97, 81 (96), 109 (80), 43 (53), 53 (40), 41 (36), 56 (29), 95 (25)... 152 (1)] | 7.17* | 1469.2 | [1.04] | 6.57 | 1154.4 | 0.03 |
| Unknown CALU II [m/z 95, 110 (38), 81 (21), 79 (16)... 152 (7)] | 7.33 | 1481.4 | 0.04 | 6.70 | 1162.7 | 0.06 |
| α-Phellandren-8-ol | 9.82 | 1680.5 | 0.02 | 6.73 | 1164.7 | 0.03 |
| Umbellulone | 8.56 | 1577.6 | 0.10 | 6.78* | 1168.3 | [0.36] |
| Unknown CALU III [m/z 95, 110 (43), 81 (28), 41 (15)... 152 (8)] | 7.42 | 1488.2 | 0.02 | 6.78* | 1168.3 | [0.36] |
| *cis*-Sabinol | 10.54* | 1740.9 | [0.25] | 6.78* | 1168.3 | [0.36] |
| Terpinen-4-ol | 8.27 | 1554.5 | 3.04 | 6.88 | 1174.5 | 3.02 |
| Cryptone | 8.85 | 1600.6 | 0.02 | 6.96* | 1179.2 | [0.12] |
| Thuj-3-en-10-al | 8.44 | 1567.7 | 0.16 | 6.96* | 1179.2 | [0.12] |
| *para*-Cymen-8-ol | 11.20 | 1798.0 | 0.19 | 7.04 | 1184.5 | 0.19 |
| α-Terpineol | 9.47* | 1651.3 | [0.27] | 7.10* | 1188.7 | [0.27] |
| Myrtenal | 8.38* | 1563.3 | [0.07] | 7.10* | 1188.7 | [0.27] |
| *cis*-Piperitol | 9.25 | 1633.6 | 0.02 | 7.17 | 1193.2 | 0.03 |
| Myrtenol | 10.54* | 1740.9 | [0.25] | 7.19 | 1194.1 | 0.05 |
| *cis*-α-Phellandrene epoxide (iPr vs Me) | 10.69 | 1754.3 | 0.13 | 7.26 | 1198.8 | 0.10 |
| *trans*-Piperitol | 10.05* | 1699.4 | [0.15] | 7.37 | 1206.1 | 0.06 |
| *trans*-Carveol | 11.07 | 1786.4 | 0.02 | 7.56 | 1218.5 | 0.02 |
| Citronellol | 10.42 | 1731.3 | 0.07 | 7.76 | 1232.5 | 0.04 |
| Cuminal | 10.17 | 1709.4 | 0.06 | 7.79 | 1234.5 | 0.06 |
| Carvone | 9.64 | 1665.9 | 0.02 | 7.86 | 1239.1 | 0.01 |
| Carvotanacetone | 9.22 | 1631.2 | 0.05 | 7.91 | 1242.2 | 0.07 |
| Unknown CALU IV [m/z 43, 97 (69), 107 (46), 41 (28), 55 (21), 109 (20)...] | | | | 8.01* | 1248.7 | [0.10] |
| Unknown BODA V [m/z 109, 119 (84), 91 (81), 134 (55)... 137 (27)...] | 10.97 | 1777.9 | 0.05 | 8.01* | 1248.7 | [0.10] |
| Piperitone | 9.55 | 1657.7 | 0.04 | 8.05 | 1251.6 | 0.03 |
| Unknown CIGL III [m/z 43, 97 (78), 41 | 13.08 | 1969.0 | 0.02 | 8.18 | 1260.4 | 0.02 |

| | | | | | | |
|---|---|---|---|---|---|---|
| (45), 55 (35), 69 (28), 107 (24), 83 (23)...] | | | | | | |
| 3,5-Dimethoxytoluene | 11.03 | 1783.0 | 0.01 | 8.24 | 1264.8 | 0.01 |
| Unknown PECR X [m/z 109, 43 (83), 95 (70), 110 (70), 99 (53), 119 (48)...] | | | | 8.30 | 1268.8 | 0.03 |
| Bornyl acetate | 7.92 | 1527.2 | 0.07 | 8.51 | 1282.6 | 0.07 |
| Cuminol | 13.82 | 2040.0 | 0.03 | 8.65 | 1292.2 | 0.04 |
| Thymol | 14.77 | 2132.8 | 0.07 | 8.82 | 1303.1 | 0.08 |
| Carvacrol | 15.00 | 2156.3 | 0.07 | 8.94 | 1311.5 | 0.06 |
| exo-2-Hydroxycineole acetate | 9.74 | 1673.6 | 0.04 | 9.29 | 1336.5 | 0.02 |
| α-Cubebene | 6.44 | 1414.3 | 0.01 | 9.45 | 1347.8 | 0.02 |
| Citronellyl acetate | 9.14 | 1624.0 | 0.03 | 9.56 | 1355.9 | 0.05 |
| α-Copaene | 6.84 | 1444.1 | 0.03 | 9.79 | 1372.3 | 0.05 |
| β-Bourbonene | 7.17* | 1469.2 | [1.04] | 9.91 | 1380.2 | 1.13 |
| β-Elemene | 8.07 | 1539.0 | 0.06 | 10.04 | 1389.7 | 0.06 |
| β-Caryophyllene | 8.11 | 1542.4 | 0.06 | 10.37* | 1413.7 | [0.14] |
| β-Ylangene | 7.82* | 1519.6 | [0.23] | 10.37* | 1413.7 | [0.14] |
| β-Copaene | 8.02 | 1535.4 | 0.14 | 10.51 | 1424.1 | 0.13 |
| Isogermacrene D | 8.62 | 1581.9 | 0.03 | 10.71 | 1439.0 | 0.08 |
| α-Humulene | 8.94 | 1608.0 | 0.02 | 10.83 | 1447.7 | 0.02 |
| allo-Aromadendrene | 8.66* | 1585.5 | [0.18] | 10.92 | 1454.8 | 0.02 |
| Germacrene D | 9.43 | 1648.3 | 0.05 | 11.20 | 1475.8 | 0.03 |
| γ-Cadinene | 10.05* | 1699.4 | [0.15] | 11.65 | 1509.2 | 0.02 |
| δ-Cadinene | 10.05* | 1699.4 | [0.15] | 11.79 | 1520.5 | 0.02 |
| Caryophyllene oxide | 12.40 | 1905.4 | 0.02 | 12.48 | 1574.2 | 0.01 |
| τ-Cadinol | 14.53 | 2109.1 | 0.03 | 13.24 | 1636.1 | 0.03 |
| α-Phellandrene dimer I | 11.42 | 1817.7 | 0.02 | 14.41 | 1733.7 | 0.03 |
| α-Phellandrene dimer II | 12.12 | 1880.1 | 0.44 | 15.04 | 1788.1 | 0.44 |
| α-Phellandrene dimer III | 12.76 | 1939.3 | 0.06 | 15.23 | 1804.4 | 0.06 |
| α-Phellandrene dimer IV | 12.92 | 1953.6 | 0.07 | 15.45 | 1824.5 | 0.06 |
| α-Phellandrene dimer VI | | | | 16.32 | 1904.0 | 0.02 |
| α-Phellandrene | | | | 16.36 | 1907.8 | 0.02 |

| | | | | | | |
|---|---|---|---|---|---|---|
| dimer VII | | | | | | |
| Total reported | 97.01% | | | 99.25% | | |
| | | | | | | |

*: 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

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