**Date :** 2025-02-04

# *Certificate of Analysis - GC Profiling*

## *Sample Identification*

**Internal code :** 25A21-PTH02
**Customer Identification :** Laurel Leaf - Greece - L10110R
**Type :** Essential Oil
**Source :** *Laurus nobilis*
**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*Notes: This report may not be published, including online, without the written consent from Laboratoire PhytoChemia.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 :** Alexis St-Gelais, Ph. D., Chimiste 2013-174
**Date :** 2025-02-04

## *Physicochemical data*

**Refractive index :** 1.4708 ± 0.0003 (20 °C)
**Method :** PC-MAT-016 - Measure of the refractive index of a liquid.
**Analyst :** Cindy Caron B. Sc.
**Date :** 2025-01-21

## *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 |
|---|---|---|
| 2-Methylbutyral | 0.01 | Aliphatic aldehyde |
| 2-Ethylfuran | tr | Furan |
| 3-Methyl-1-penten-3-ol | tr | Aliphatic alcohol |
| 1-Methylpyrrole | 0.02 | Pyrrole |
| Ethyl isobutyrate | 0.01 | Aliphatic ester |
| Toluene | tr | Simple phenolic |
| Hexanal | 0.05 | Aliphatic aldehyde |
| Isopropyl butyrate | tr | Aliphatic ester |
| (2*E*)-Hexenal | 0.01 | Aliphatic aldehyde |
| Ethyl 2-methylbutyrate | 0.03 | Aliphatic ester |
| Ethyl isovalerate | 0.02 | Aliphatic ester |
| (3*Z*)-Hexenol | 0.03 | Aliphatic alcohol |
| Isopropyl 2-methylbutyrate | 0.01 | Aliphatic ester |
| Tricyclene | tr | Monoterpene |
| Isobutyl isobutyrate | 0.02 | Aliphatic ester |
| α-Thujene | 0.18 | Monoterpene |
| α-Pinene | 6.29 | Monoterpene |
| α-Fenchene | 0.02 | Monoterpene |
| Camphene | 0.30 | Monoterpene |
| β-Pinene | 3.40 | Monoterpene |
| Sabinene | 7.50 | Monoterpene |
| Dehydro-1,8-cineole | 0.04 | Monoterpenic ether |
| Myrcene | 0.42 | Monoterpene |
| Pseudolimonene | 0.02 | Monoterpene |
| α-Phellandrene | 0.27 | Monoterpene |
| Isobutyl 2-methylbutyrate | 0.05 | Aliphatic ester |
| (3*Z*)-Hexenyl acetate | tr | Aliphatic ester |
| Δ3-Carene | 0.02 | Monoterpene |
| α-Terpinene | 0.36 | Monoterpene |
| *para*-Cymene | 1.41 | Monoterpene |
| Limonene | 2.63 | Monoterpene |
| 1,8-Cineole | 37.16 | Monoterpenic ether |
| (*Z*)-β-Ocimene | 0.10 | Monoterpene |
| (*E*)-β-Ocimene | 0.14 | Monoterpene |
| γ-Terpinene | 1.24 | Monoterpene |
| *cis*-Sabinene hydrate | 0.03 | Monoterpenic alcohol |
| *cis*-Linalool oxide (fur.) | 0.02 | Monoterpenic alcohol |
| *para*-Cymenene | 0.02 | Monoterpene |
| Terpinolene | 0.09 | Monoterpene |
| 2-Nonanone | 0.04 | Aliphatic ketone |



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| | | |
|---|---|---|
| *trans*-Sabinene hydrate | 0.03 | Monoterpenic alcohol |
| Linalool | 8.36 | Monoterpenic alcohol |
| Unknown | 0.02 | Unknown |
| Hotrienol | 0.02 | Monoterpenic alcohol |
| endo-Fenchol | 0.01 | Monoterpenic alcohol |
| *cis-para*-Menth-2-en-1-ol | 0.02 | Monoterpenic alcohol |
| *trans*-Pinocarveol | 0.05 | Monoterpenic alcohol |
| *trans-para*-Menth-2-en-1-ol | 0.02 | Monoterpenic alcohol |
| *trans*-Verbenol | 0.02 | Monoterpenic alcohol |
| Nerol oxide | tr | Aliphatic ether |
| Pinocarvone | 0.03 | Monoterpenic ketone |
| (*E*)-2,6-Dimethyl-1,5,7-octatrien-3-ol | 0.01 | Monoterpenic alcohol |
| Borneol | 0.03 | Monoterpenic alcohol |
| δ-Terpineol | 0.09 | Monoterpenic alcohol |
| Isopinocamphone | tr | Monoterpenic ketone |
| Terpinen-4-ol | 2.34 | Monoterpenic alcohol |
| Thuj-3-en-10-al | 0.01 | Monoterpenic aldehyde |
| *para*-Cymen-8-ol | 0.01 | Monoterpenic alcohol |
| Myrtenal | 0.03 | Monoterpenic aldehyde |
| α-Terpineol | 2.13 | Monoterpenic alcohol |
| Myrtenol | 0.05 | Monoterpenic alcohol |
| Methylchavicol | 0.12 | Phenylpropanoid |
| Nerol | 0.17 | Monoterpenic alcohol |
| Linalyl acetate | 0.07 | Monoterpenic ester |
| Geraniol | 0.05 | Monoterpenic alcohol |
| Geranial | 0.01 | Monoterpenic aldehyde |
| 4-Thujen-2α-yl acetate | 0.03 | Monoterpenic ester |
| Bornyl acetate | 0.62 | Monoterpenic ester |
| *para*-Cymen-7-ol | 0.01 | Monoterpenic alcohol |
| 2-Undecanone | 0.07 | Aliphatic ketone |
| *para*-Menth-5-en-1,2-diol isomer III | 0.01 | Monoterpenic alcohol |
| δ-Terpinyl acetate | 0.15 | Monoterpenic ester |
| Unknown | 0.02 | Unknown |
| exo-2-Hydroxycineole acetate | 0.02 | Monoterpenic ester |
| α-Cubebene | 0.02 | Sesquiterpene |
| α-Terpinyl acetate | 13.13 | Monoterpenic ester |
| Eugenol | 2.43 | Phenylpropanoid |
| Neryl acetate | 0.09 | Monoterpenic ester |
| Unknown | 0.02 | Sesquiterpene |
| α-Ylangene | 0.03 | Sesquiterpene |
| α-Copaene | 0.02 | Sesquiterpene |
| β-Elemene | 0.09 | Sesquiterpene |
| Methyleugenol | 5.80 | Phenylpropanoid |
| β-Caryophyllene | 0.33 | Sesquiterpene |
| Aromadendrene | 0.01 | Sesquiterpene |

| | | |
|---|---|---|
| α-Guaiene | 0.01 | Sesquiterpene |
| 6,9-Guaiadiene | 0.01 | Sesquiterpene |
| (*E*)-Cinnamyl acetate | 0.06 | Phenylpropanoid ester |
| α-Humulene | 0.05 | Sesquiterpene |
| Selina-4(15),7-diene | 0.02 | Sesquiterpene |
| *cis*-Muurola-4(15),5-diene | 0.01 | Sesquiterpene |
| Germacrene D | 0.02 | Sesquiterpene |
| β-Selinene | 0.03 | Sesquiterpene |
| Viridiflorene | 0.01 | Sesquiterpene |
| α-Selinene | 0.03 | Sesquiterpene |
| Bicyclogermacrene | 0.01 | Sesquiterpene |
| (3*Z*,6*E*)-α-Farnesene | 0.02 | Sesquiterpene |
| Methyl (*E*)-isoeugenol | 0.10 | Phenylpropanoid |
| γ-Cadinene | 0.04 | Sesquiterpene |
| δ-Cadinene | 0.06 | Sesquiterpene |
| α-Calacorene | 0.01 | Sesquiterpene |
| (*E*)-α-Bisabolene | 0.03 | Sesquiterpene |
| Elemicin | 0.02 | Phenylpropanoid |
| Spathulenol | 0.04 | Sesquiterpenic alcohol |
| Caryophyllene oxide | 0.06 | Sesquiterpenic ether |
| Viridiflorol | 0.01 | Sesquiterpenic alcohol |
| Ledol | 0.01 | Sesquiterpenic alcohol |
| Humulene epoxide II | 0.01 | Sesquiterpenic ether |
| Junenol | 0.02 | Sesquiterpenic alcohol |
| Eremoligenol? | 0.01 | Sesquiterpenic alcohol |
| Caryophylladienol II | 0.02 | Sesquiterpenic alcohol |
| τ-Cadinol | 0.01 | Sesquiterpenic alcohol |
| β-Eudesmol | 0.03 | Sesquiterpenic alcohol |
| α-Eudesmol | 0.01 | Sesquiterpenic alcohol |
| (*E*)-Isoelemicin | 0.01 | Phenylpropanoid |
| α-Cadinol | 0.01 | Sesquiterpenic alcohol |
| Unknown | 0.01 | Aliphatic ester |
| **Consolidated total** | **99.41** | |

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.

This page was intentionally left blank. The following pages present the complete data of the analysis.

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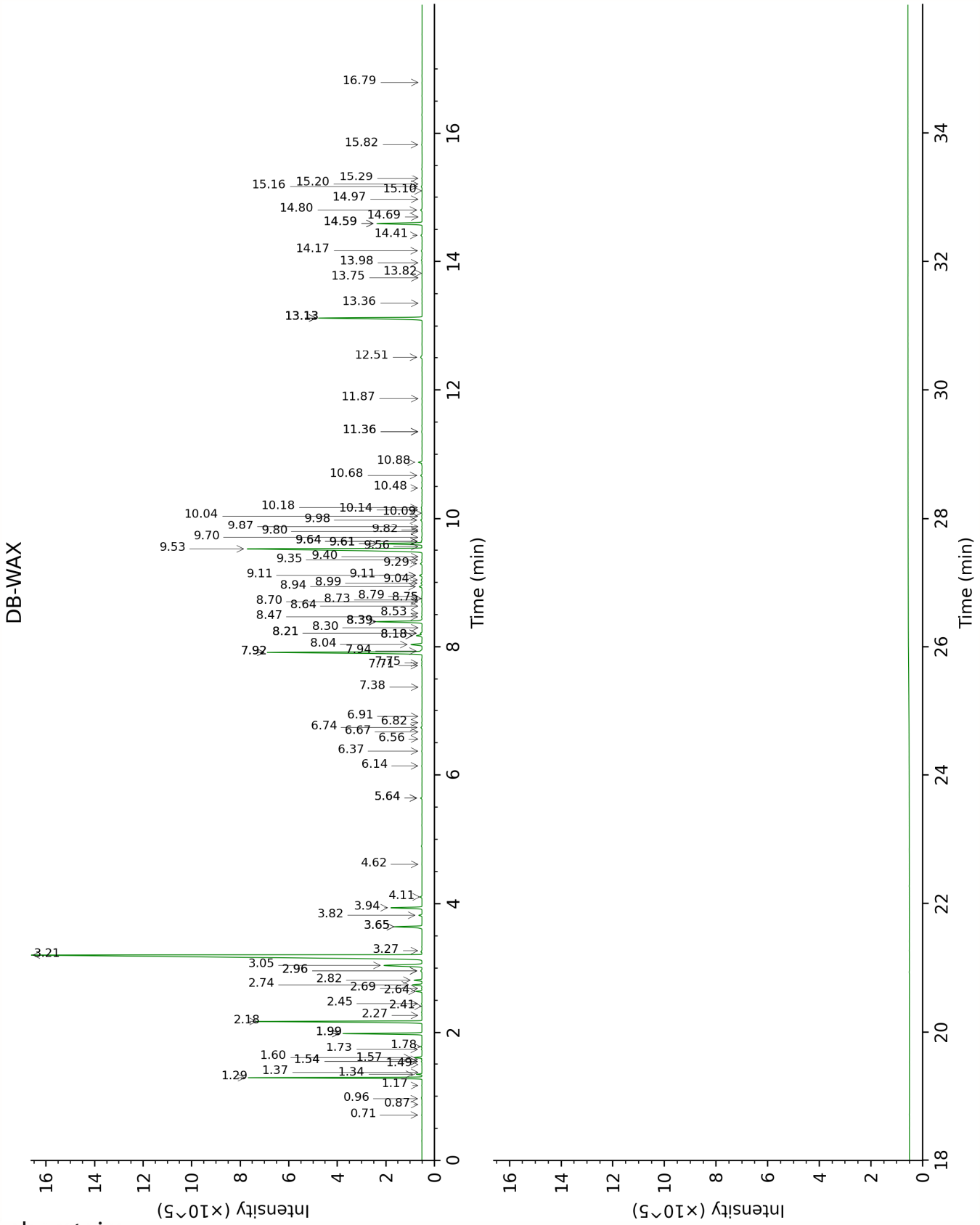


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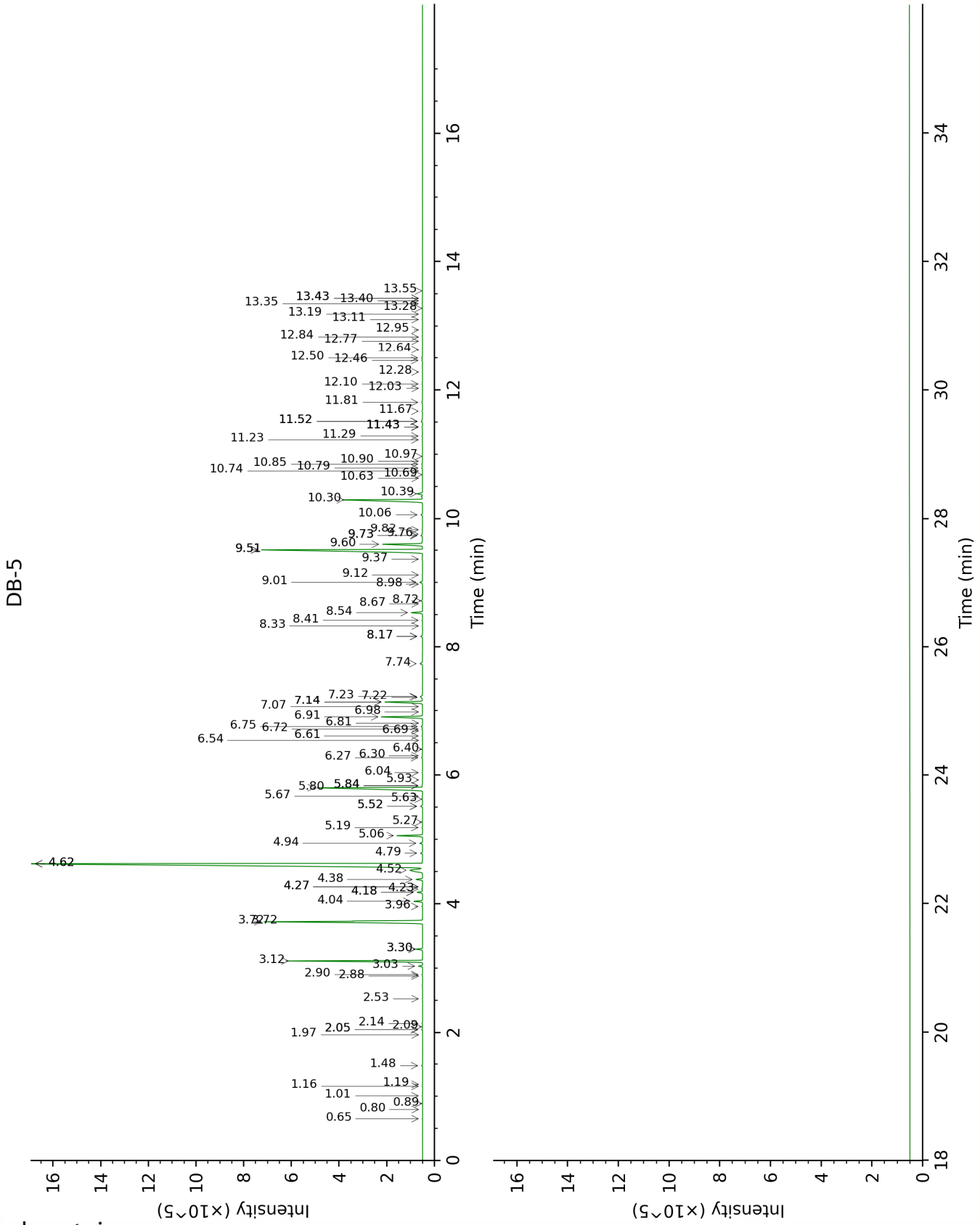


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*FULL ANALYSIS DATA*

| | Column DB-WAX | | | Column DB-5 | | |
|---|---|---|---|---|---|---|
| **2-Methylbutyral** | 0.71 | 885.0 | 0.01 | 0.65 | 653.3 | 0.01 |
| 2-Ethylfuran | 0.87 | 922.6 | 0.01 | 0.80 | 703.7 | tr |
| 3-Methyl-1-penten-3-ol | 2.41 | 1108.9 | tr | 0.89 | 718.4 | tr |
| 1-Methylpyrrole | 2.27 | 1096.4 | 0.01 | 1.01 | 735.4 | 0.02 |
| Ethyl isobutyrate | 0.96 | 937.9 | 0.01 | 1.16 | 756.4 | 0.01 |
| Toluene | 1.37 | 1005.3 | tr | 1.19 | 760.8 | tr |
| Hexanal | 1.78 | 1047.0 | 0.05 | 1.48 | 801.9 | 0.05 |
| Isopropyl butyrate | 1.49 | 1017.6 | 0.01 | 1.97 | 844.5 | tr |
| (2*E*)-Hexenal | 3.27 | 1177.6 | 0.01 | 2.05* | 851.2 | [0.04] |
| Ethyl 2-methylbutyrate | 1.54* | 1022.5 | [0.02] | 2.05* | 851.2 | [0.04] |
| Ethyl isovalerate | 1.73 | 1041.6 | 0.02 | 2.09 | 854.9 | 0.02 |
| (3*Z*)-Hexenol | 5.64* | 1352.8 | [0.07] | 2.14 | 859.0 | 0.03 |
| Isopropyl 2-methylbutyrate | 1.58 | 1025.9 | 0.02 | 2.53 | 891.1 | 0.01 |
| Tricyclene | 1.17 | 973.1 | 0.01 | 2.88 | 917.5 | tr |
| Isobutyl isobutyrate | 1.99* | 1067.8 | [3.37] | 2.90 | 919.0 | 0.02 |
| α-Thujene | 1.34 | 1002.5 | 0.18 | 3.03 | 927.7 | 0.18 |
| α-Pinene | 1.29 | 993.9 | 6.23 | 3.12 | 933.1 | 6.29 |
| α-Fenchene | 1.54* | 1022.5 | [0.02] | 3.30* | 945.2 | [0.33] |
| Camphene | 1.60 | 1028.8 | 0.30 | 3.30* | 945.2 | [0.33] |
| β-Pinene | 1.99* | 1067.8 | [3.37] | 3.72* | 973.4 | [10.90] |
| Sabinene | 2.18 | 1086.9 | 7.50 | 3.72* | 973.4 | [10.90] |
| Dehydro-1,8-cineole | 2.96* | 1152.5 | [0.07] | 3.96 | 989.0 | 0.04 |
| Myrcene | 2.74 | 1135.3 | 0.41 | 4.04 | 994.5 | 0.42 |
| Pseudolimonene | 2.69 | 1130.8 | 0.02 | 4.18* | 1003.7 | [0.28] |
| α-Phellandrene | 2.64 | 1127.3 | 0.27 | 4.18* | 1003.7 | [0.28] |
| Isobutyl 2-methylbutyrate | 2.96* | 1152.5 | [0.07] | 4.23 | 1006.9 | 0.05 |
| (3*Z*)-Hexenyl acetate | 4.62 | 1278.9 | tr | 4.27* | 1009.1 | [0.03] |
| Δ3-Carene | 2.45 | 1112.1 | 0.02 | 4.27* | 1009.1 | [0.03] |
| α-Terpinene | 2.82 | 1141.1 | 0.36 | 4.38 | 1016.2 | 0.36 |
| *para*-Cymene | 3.94 | 1228.8 | 1.41 | 4.52 | 1025.2 | 1.41 |
| Limonene | 3.05 | 1159.4 | 2.63 | 4.62* | 1031.5 | [39.94] |
| 1,8-Cineole | 3.21 | 1172.2 | 37.16 | 4.62* | 1031.5 | [39.94] |
| (*Z*)-β-Ocimene | 3.65* | 1207.0 | [1.35] | 4.79 | 1041.8 | 0.10 |
| (*E*)-β-Ocimene | 3.82 | 1220.2 | 0.14 | 4.94 | 1051.5 | 0.14 |
| γ-Terpinene | 3.65* | 1207.0 | [1.35] | 5.06 | 1058.9 | 1.24 |
| *cis*-Sabinene | 6.74 | 1432.7 | 0.06 | 5.19 | 1066.8 | 0.03 |

| | | | | | | |
|---|---|---|---|---|---|---|
| hydrate | | | | | | |
| *cis*-Linalool oxide (fur.) | 6.37 | 1405.4 | 0.02 | 5.27 | 1072.0 | 0.02 |
| *para*-Cymenene | 6.14 | 1388.8 | 0.02 | 5.52* | 1087.4 | [0.13] |
| Terpinolene | 4.11 | 1241.3 | 0.09 | 5.52* | 1087.4 | [0.13] |
| 2-Nonanone | 5.64* | 1352.8 | [0.07] | 5.63 | 1094.4 | 0.04 |
| *trans*-Sabinene hydrate | 7.75 | 1508.2 | 0.02 | 5.67 | 1097.2 | 0.03 |
| Linalool | 7.92* | 1521.2 | [8.35] | 5.80 | 1105.4 | 8.36 |
| Unknown CASA I [m/z 43, 59 (37), 79 (33), 91 (32), 119 (31)...] | 8.79 | 1588.7 | 0.02 | 5.84* | 1107.5 | [0.05] |
| Hotrienol | 8.64 | 1576.9 | 0.02 | 5.84* | 1107.5 | [0.05] |
| endo-Fenchol | 8.18* | 1541.4 | [0.32] | 5.93 | 1113.2 | 0.01 |
| *cis-para*-Menth-2-en-1-ol | 7.92* | 1521.2 | [8.35] | 6.04 | 1120.4 | 0.02 |
| *trans*-Pinocarveol | 8.99 | 1604.8 | 0.06 | 6.27 | 1135.2 | 0.05 |
| *trans-para*-Menth-2-en-1-ol | 8.73 | 1584.3 | 0.02 | 6.30 | 1137.2 | 0.02 |
| *trans*-Verbenol | 9.35 | 1634.1 | 0.03 | 6.40 | 1143.6 | 0.02 |
| Nerol oxide | 6.67 | 1427.5 | 0.01 | 6.54 | 1152.3 | tr |
| Pinocarvone | 7.71 | 1505.2 | 0.03 | 6.61 | 1156.7 | 0.03 |
| (*E*)-2,6-Dimethyl-1,5,7-octatrien-3-ol | 10.09 | 1694.2 | 0.02 | 6.69 | 1161.8 | 0.01 |
| Borneol | 9.61* | 1654.8 | [2.15] | 6.72 | 1163.6 | 0.03 |
| δ-Terpineol | 9.29 | 1629.4 | 0.09 | 6.75 | 1166.0 | 0.09 |
| Isopinocamphone | 7.38 | 1480.1 | 0.01 | 6.81 | 1169.4 | tr |
| Terpinen-4-ol | 8.40* | 1558.2 | [2.40] | 6.91 | 1175.7 | 2.34 |
| Thuj-3-en-10-al | 8.52 | 1568.3 | 0.02 | 6.98 | 1180.5 | 0.01 |
| *para*-Cymen-8-ol | 11.36* | 1801.6 | [0.03] | 7.06 | 1185.9 | 0.01 |
| Myrtenal | 8.47 | 1564.0 | 0.03 | 7.14* | 1190.4 | [2.16] |
| α-Terpineol | 9.61* | 1654.8 | [2.15] | 7.14* | 1190.4 | [2.16] |
| Myrtenol | 10.68 | 1743.7 | 0.07 | 7.22 | 1195.6 | 0.05 |
| Methylchavicol | 9.11* | 1614.7 | [0.13] | 7.23 | 1196.5 | 0.12 |
| Nerol | 10.88 | 1761.2 | 0.19 | 7.74 | 1230.2 | 0.17 |
| Linalyl acetate | 7.94 | 1522.7 | 0.07 | 8.17* | 1258.6 | [0.11] |
| Geraniol | 11.36* | 1801.6 | [0.03] | 8.17* | 1258.6 | [0.11] |
| Geranial | 9.87 | 1676.2 | 0.02 | 8.33 | 1269.4 | 0.01 |
| 4-Thujen-2α-yl acetate | 8.70 | 1582.2 | 0.03 | 8.42 | 1275.2 | 0.03 |
| Bornyl acetate | 8.04 | 1530.6 | 0.60 | 8.54 | 1283.2 | 0.62 |
| *para*-Cymen-7-ol | 13.98 | 2043.3 | 0.02 | 8.67 | 1292.4 | 0.01 |
| 2-Undecanone | 8.40* | 1558.2 | [2.40] | 8.72 | 1295.6 | 0.07 |
| *para*-Menth-5-en- | 14.97 | 2140.1 | 0.02 | 8.98 | 1313.2 | 0.01 |

| | | | | | | |
|---|---|---|---|---|---|---|
| 1,2-diol isomer III | | | | | | |
| δ-Terpinyl acetate | 8.94 | 1600.5 | 0.15 | 9.01 | 1315.3 | 0.15 |
| Unknown LANO II [m/z 119, 43 (99), 93 (52), 59 (44), 91 (41), 134 (34)...] | 9.64* | 1657.6 | [0.05] | 9.12 | 1323.3 | 0.02 |
| exo-2-Hydroxycineole acetate | 9.82 | 1672.0 | 0.02 | 9.36 | 1340.6 | 0.02 |
| α-Cubebene | 6.56 | 1419.3 | 0.02 | 9.51* | 1350.8 | [13.32] |
| α-Terpinyl acetate | 9.52 | 1648.1 | 13.13 | 9.51* | 1350.8 | [13.32] |
| Eugenol | 14.59* | 2102.3 | [2.43] | 9.60 | 1356.9 | 2.43 |
| Neryl acetate | 9.98 | 1685.4 | 0.09 | 9.73* | 1366.5 | [0.10] |
| Unknown ENKR II [m/z 189, 91 (89), 82 (78), 109 (72), 93 (68), 67 (63), 161 (59)... 204 (tr)] | | | | 9.73* | 1366.5 | [0.10] |
| α-Ylangene | 6.82 | 1438.3 | 0.02 | 9.76 | 1368.1 | 0.03 |
| α-Copaene | 6.91 | 1445.5 | 0.02 | 9.82 | 1372.8 | 0.02 |
| β-Elemene | 8.22* | 1544.2 | [0.10] | 10.06 | 1389.9 | 0.09 |
| Methyleugenol | 13.13* | 1962.3 | [5.73] | 10.30 | 1406.5 | 5.80 |
| β-Caryophyllene | 8.18* | 1541.4 | [0.32] | 10.39 | 1413.5 | 0.33 |
| Aromadendrene | 8.30 | 1550.8 | 0.01 | 10.63 | 1431.3 | 0.01 |
| α-Guaiene | 8.22* | 1544.2 | [0.10] | 10.69 | 1435.4 | 0.01 |
| 6,9-Guaiadiene | 8.40* | 1558.2 | [2.40] | 10.74 | 1439.7 | 0.01 |
| (*E*)-Cinnamyl acetate | 14.41 | 2084.4 | 0.07 | 10.79 | 1443.2 | 0.06 |
| α-Humulene | 9.04 | 1608.9 | 0.04 | 10.85 | 1447.6 | 0.05 |
| Selina-4(15),7-diene | 8.75 | 1586.1 | 0.01 | 10.90 | 1451.1 | 0.02 |
| *cis*-Muurola-4(15),5-diene | 9.11* | 1614.7 | [0.13] | 10.97 | 1456.7 | 0.01 |
| Germacrene D | 9.56 | 1651.1 | 0.01 | 11.23 | 1475.8 | 0.02 |
| β-Selinene | 9.64* | 1657.6 | [0.05] | 11.29 | 1480.2 | 0.03 |
| Viridiflorene | 9.40 | 1638.3 | 0.01 | 11.43* | 1490.6 | [0.04] |
| α-Selinene | 9.70 | 1662.6 | 0.03 | 11.43* | 1490.6 | [0.04] |
| Bicyclogermacrene | 9.80 | 1670.2 | 0.01 | 11.43* | 1490.6 | [0.04] |
| (3*Z*,6*E*)-α-Farnesene | 10.04 | 1690.3 | 0.02 | 11.52* | 1497.1 | [0.12] |
| Methyl (*E*)-isoeugenol | 14.80 | 2123.2 | 0.10 | 11.52* | 1497.1 | [0.12] |
| γ-Cadinene | 10.14 | 1698.2 | 0.05 | 11.67 | 1509.1 | 0.04 |
| δ-Cadinene | 10.18 | 1701.3 | 0.05 | 11.81 | 1519.9 | 0.06 |
| α-Calacorene | 11.87 | 1847.3 | 0.01 | 12.03 | 1537.0 | 0.01 |

| | | | | | | |
|---|---|---|---|---|---|---|
| (*E*)-α-Bisabolene | 10.48 | 1727.1 | 0.03 | 12.10 | 1542.3 | 0.03 |
| Elemicin | 15.29 | 2172.5 | 0.02 | 12.28 | 1556.9 | 0.02 |
| Spathulenol | 14.17 | 2061.4 | 0.04 | 12.46 | 1571.0 | 0.04 |
| Caryophyllene oxide | 12.51 | 1904.9 | 0.12 | 12.50 | 1574.2 | 0.06 |
| Viridiflorol | 13.75 | 2020.8 | 0.01 | 12.64 | 1584.7 | 0.01 |
| Ledol | 13.13* | 1962.3 | [5.73] | 12.77 | 1595.2 | 0.01 |
| Humulene epoxide II | 13.13* | 1962.3 | [5.73] | 12.84 | 1600.3 | 0.01 |
| Junenol | 13.36 | 1983.4 | 0.01 | 12.95 | 1609.1 | 0.02 |
| Eremoligenol? | 14.59* | 2102.3 | [2.43] | 13.11 | 1622.4 | 0.01 |
| Caryophylladienol II | 15.82 | 2227.5 | 0.03 | 13.19 | 1629.2 | 0.02 |
| τ-Cadinol | 14.69 | 2112.5 | 0.01 | 13.28 | 1636.8 | 0.01 |
| β-Eudesmol | 15.16 | 2159.7 | 0.05 | 13.35 | 1642.6 | 0.03 |
| α-Eudesmol | 15.10 | 2153.1 | 0.01 | 13.40 | 1646.5 | 0.01 |
| (*E*)-Isoelemicin | 16.79 | 2329.9 | 0.01 | 13.43* | 1649.5 | [0.02] |
| α-Cadinol | 15.20 | 2163.7 | 0.01 | 13.43* | 1649.5 | [0.02] |
| Unknown HULU VI [m/z 79, 43 (66), 67 (59), 80 (56), 41 (41), 81 (37), 55 (29)...] | 13.82 | 2027.7 | 0.01 | 13.55 | 1659.0 | 0.01 |
| Total reported | | 99.19% | | | 99.74% | |

*: 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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