

Date : 2024-07-15

CERTIFICATE OF ANALYSIS - GC PROFILING

SAMPLE IDENTIFICATION

**Internal code** : 24F28-PTH02

**Customer Identification** : Davana - India - D80106R

**Type** : Essential Oil

**Source** : *Artemisia pallens*

**Customer** : Plant Therapy

Checked and approved by:

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

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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 :** 2024-07-09

## PHYSICOCHEMICAL DATA

**Refractive index :**  $1.4908 \pm 0.0003$  (20 °C)

**Method :** PC-MAT-016 - Measure of the refractive index of a liquid.

**Analyst :** Cindy Caron B. Sc.

**Date :** 2024-06-28

## CONCLUSION

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

## ANALYSIS SUMMARY - CONSOLIDATED CONTENTS

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

Identification	%	Class
2-Methylbutyral	tr	Aliphatic aldehyde
Unknown	0.01	Unknown
Ethyl 2-methylbutyrate	0.15	Aliphatic ester
Ethyl isovalerate	0.10	Aliphatic ester
Propyl isobutyrate	0.01	Aliphatic ester
Nonene	tr	Alkene
( <i>cis?</i> )-5-Ethyl-2-methyl-2-vinyltetrahydrofuran	tr	Furan
( <i>trans?</i> )-5-Ethyl-2-methyl-2-vinyltetrahydrofuran	0.08	Furan
( <i>trans?</i> )-2,5-Divinyl-2-methyltetrahydrofuran	0.06	Furan
$\alpha$ -Pinene	0.02	Monoterpene
Unknown	0.38	Furan
5,5-Dimethyl-2(5H)-furanone	0.12	Aliphatic lactone
Propyl 2-methylbutyrate	0.22	Aliphatic ester
Propyl isovalerate	0.16	Aliphatic ester
Benzaldehyde	0.02	Simple phenolic
Sabinene	0.01	Monoterpene
$\beta$ -Pinene	0.02	Monoterpene
6-Methyl-5-hepten-2-one	0.01	Aliphatic ketone
<i>trans</i> -Dehydroxylinalool oxide	0.02	Monoterpenic ether
Myrcene	0.01	Monoterpene
2-Pentylfuran	0.02	Furan
Unknown	0.01	Furan
Octanal	0.01	Aliphatic aldehyde
<i>cis</i> -Dehydroxylinalool oxide	0.02	Monoterpenic ether
Unknown	0.14	Furan
Isobutyl 2-methylbutyrate	0.02	Aliphatic ester
Isobutyl isovalerate	0.02	Aliphatic ester
$\alpha$ -Terpinene	0.05	Monoterpene
Unknown	0.23	Furan
<i>para</i> -Cymene	0.19	Monoterpene
Limonene	0.02	Monoterpene
1,8-Cineole	0.04	Monoterpenic ether
Lavender lactone	0.06	Aliphatic lactone
<i>cis</i> -Arbusculone	0.04	Furan
( <i>E</i> )- $\beta$ -Ocimene	0.02	Monoterpene
$\gamma$ -Terpinene	0.11	Monoterpene
<i>cis</i> -Arbusculol	0.02	Furan
<i>cis</i> -Sabinene hydrate	0.12	Monoterpenic alcohol
<i>trans</i> -Arbusculone	0.03	Furan

<i>trans</i> -Arbusculol	0.02	Furan
Terpinolene	0.03	Monoterpene
<i>trans</i> -Sabinene hydrate	0.04	Monoterpenic alcohol
Linalool	0.46	Monoterpenic alcohol
Hotrienol	0.08	Monoterpenic alcohol
Nonanal	0.03	Aliphatic aldehyde
2-Methylbutyl 2-methylbutyrate	0.27	Aliphatic ester
2-Methylbutyl isovalerate	0.13	Aliphatic ester
3-Methyl-3-butenyl 2-methylbutyrate	0.02	Aliphatic ester
<i>cis-para</i> -Menth-2-en-1-ol	0.04	Monoterpenic alcohol
<i>trans-para</i> -Menth-2-en-1-ol	0.02	Monoterpenic alcohol
Prenyl 2-methylbutyrate	0.03	Aliphatic ester
Prenyl isovalerate	0.04	Aliphatic ester
Nerol oxide	0.01	Aliphatic ether
Terpinen-4-ol	0.25	Monoterpenic alcohol
<i>para</i> -Cymen-8-ol	0.01	Monoterpenic alcohol
$\alpha$ -Terpineol	0.02	Monoterpenic alcohol
<i>cis</i> -Piperitol	0.03	Monoterpenic alcohol
Ethyl octanoate	0.01	Aliphatic ester
<i>cis</i> -nor-Davanone?	0.03	Furan
nor-Davanone	0.20	Furan
(3 <i>Z</i> )-Hexenyl 2-methylbutyrate	0.02	Aliphatic ester
(3 <i>Z</i> )-Hexenyl isovalerate	0.06	Aliphatic ester
Neral	0.04	Monoterpenic aldehyde
Geraniol	0.01	Monoterpenic alcohol
Geranial	0.03	Monoterpenic aldehyde
Methyl hydrocinnamate	0.01	Phenylpropanoid ester
Cogeijerene	0.05	Terpene derivative
Methyl ( <i>Z</i> )-cinnamate	0.12	Phenylpropanoid ester
Bicycloelemene analog	0.02	Sesquiterpene
Bicycloelemene	0.23	Sesquiterpene
Ethyl hydrocinnamate	0.13	Phenylpropanoid ester
Eugenol	0.10	Phenylpropanoid
Isoledene	0.04	Sesquiterpene
$\alpha$ -Copaene	0.04	Sesquiterpene
Ethyl ( <i>Z</i> )-cinnamate	1.10	Phenylpropanoid ester
Modhephene	0.04	Sesquiterpene
Methyl ( <i>E</i> )-cinnamate	0.64	Phenylpropanoid ester
<i>cis</i> - $\beta$ -Elemene	0.04	Sesquiterpene
Geranyl acetate	0.71	Monoterpenic ester
$\beta$ -Cubebene	0.01	Sesquiterpene
$\beta$ -Elemene	0.29	Sesquiterpene
Benzyl isovalerate	0.06	Phenolic ester
<i>trans</i> -erythro-Davanafuran?	0.07	Furan
<i>cis</i> -erythro-Davanafuran?	0.06	Furan

$\alpha$ -Gurjunene	0.07	Sesquiterpene
<i>trans</i> -threo-Davana-furan?	0.04	Furan
$\beta$ -Maaliene	0.02	Sesquiterpene
<i>cis</i> -threo-Davana-furan	0.45	Furan
$\beta$ -Caryophyllene	0.10	Sesquiterpene
$\beta$ -Copaene	0.05	Sesquiterpene
$\alpha$ -Maaliene	0.01	Sesquiterpene
Aromadendrene	0.34	Sesquiterpene
Selina-5,11-diene	0.04	Sesquiterpene
$\alpha$ -Humulene	0.14	Sesquiterpene
allo-Aromadendrene	0.60	Sesquiterpene
Unknown	0.13	Unknown
Ethyl ( <i>E</i> )-cinnamate	3.80	Phenylpropanoid ester
Selina-4,11-diene	0.05	Sesquiterpene
$\gamma$ -Murolene	0.17	Sesquiterpene
Germacrene D	0.91	Sesquiterpene
$\beta$ -Selinene	1.44	Sesquiterpene
$\alpha$ -Selinene	0.06	Sesquiterpene
Davana ether isomer I	1.56	Sesquiterpenic ether
Bicyclogermacrene	9.93	Sesquiterpene
Viridiflorene	0.62	Sesquiterpene
$\alpha$ -Murolene	0.03	Sesquiterpene
$\gamma$ -Cadinene	0.37	Sesquiterpene
Davana ether isomer II	4.13	Sesquiterpenic ether
Davana ether isomer III	1.11	Sesquiterpenic ether
Artedouglasia oxide C	0.26	Sesquiterpenic ketone
Zonarene	0.07	Sesquiterpene
$\delta$ -Cadinene	0.15	Sesquiterpene
Laciniata furanone G?	0.05	Sesquiterpenic ketone
Laciniata furanone F?	0.05	Sesquiterpenic ketone
Artedouglasia oxide A	0.34	Sesquiterpenic ketone
Davana ether isomer IV	2.58	Sesquiterpenic ether
$\alpha$ -Calacorene	0.05	Sesquiterpene
Laciniata furanone <i>E</i> ?	0.07	Sesquiterpenic ketone
Laciniata furanone H	0.10	Sesquiterpenic ketone
Davanone A	0.52	Sesquiterpenic ketone
Artedouglasia oxide D	0.23	Sesquiterpenic ketone
Geranyl butyrate	0.02	Monoterpenic ester
Davanone B	1.42	Sesquiterpenic ketone
( <i>E</i> )-Nerolidol	0.37	Sesquiterpenic alcohol
Davanone C	0.28	Sesquiterpenic ketone
Spathulenol	1.13	Sesquiterpenic alcohol
Artedouglasia oxide B	0.19	Sesquiterpenic ketone
Globulol	0.40	Sesquiterpenic alcohol
Cubeban-11-ol	0.17	Sesquiterpenic alcohol

Davanone D	40.99	Sesquiterpenic ketone
Viridiflorol	0.27	Sesquiterpenic alcohol
Unknown	0.50	Oxygenated sesquiterpene
Eudesm-5-en-11-ol	0.34	Sesquiterpenic alcohol
Unknown	0.20	Oxygenated sesquiterpene
Davanol D isomer I	0.56	Sesquiterpenic alcohol
Unknown	0.14	Unknown
Isospathulenol	0.61	Sesquiterpenic alcohol
$\tau$ -Cadinol	1.11	Sesquiterpenic alcohol
Unknown	0.50	Unknown
Methyl <i>cis</i> -jasmonate	0.23	Jasmonate
$\beta$ -Eudesmol	0.57	Sesquiterpenic alcohol
5-Hydroxy-6-methyl-2-(5-methyl-5-vinyltetrahydrofuran-2-yl)hepta-4,6-dien-3-one, isomer II	0.19	Sesquiterpenic alcohol
$\alpha$ -Cadinol	0.21	Sesquiterpenic alcohol
Unknown	0.22	Oxygenated sesquiterpene
Unknown	2.50	Oxygenated sesquiterpene
Davanyl acetate	0.14	Sesquiterpenic ester
Davanonol isomer	0.08	Sesquiterpenic alcohol
Unknown	0.26	Oxygenated sesquiterpene
$\beta$ -Davanon-2-ol	0.32	Sesquiterpenic alcohol
Phytone	0.13	Terpenic ketone
Nonadecane	0.04	Alkane
Heneicosane	0.11	Alkane
Phytol	0.13	Diterpenic alcohol
<b>Consolidated total</b>	<b>93.19</b>	

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

Note: no correction factor was applied

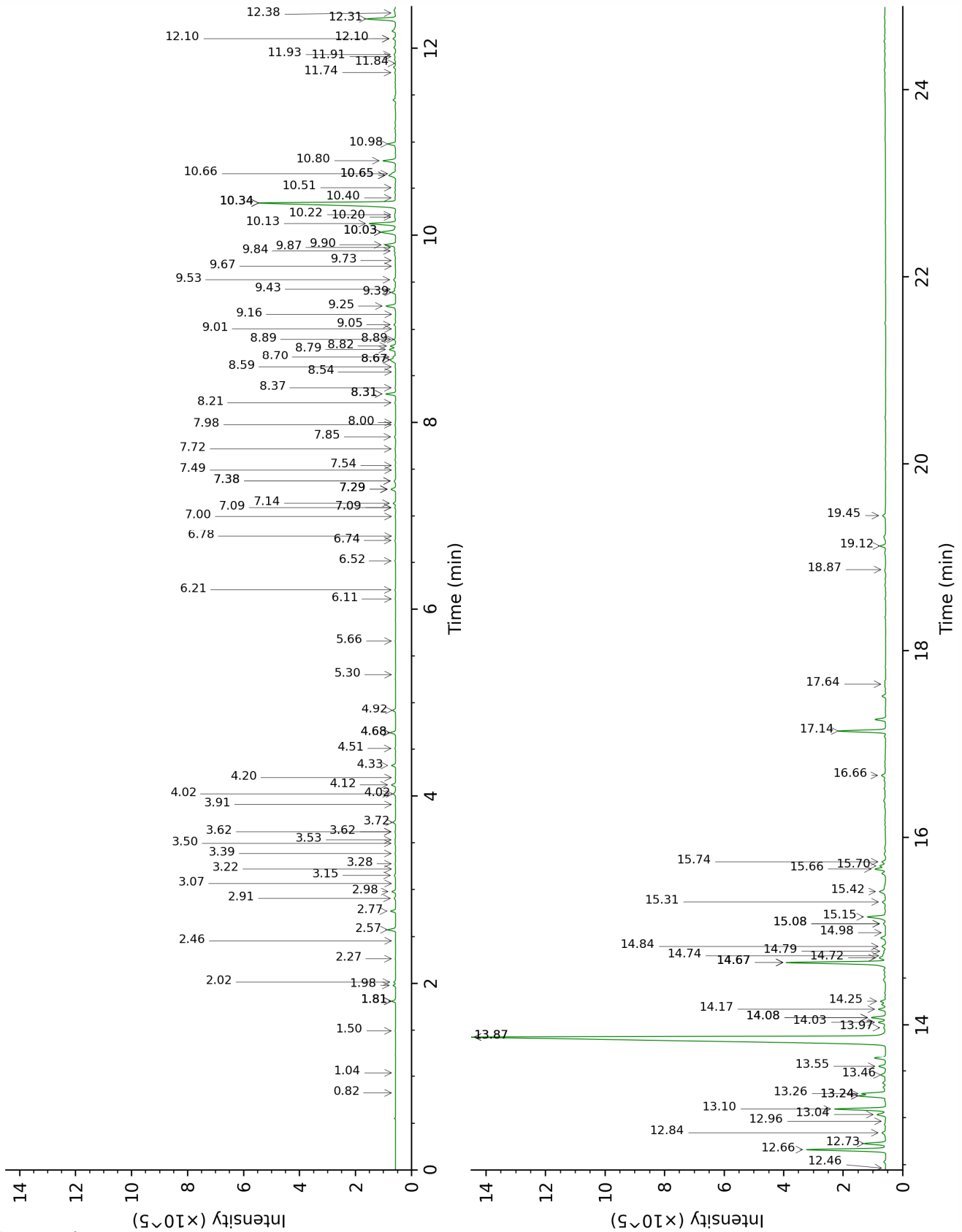
**About "consolidated" data:** The table above presents the breakdown of the sample volatile constituents after applying an algorithm to collapse data acquired from the multi-columns system of PhytoChemia into a single set of consolidated contents. In case of discrepancies between columns, the algorithm 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.

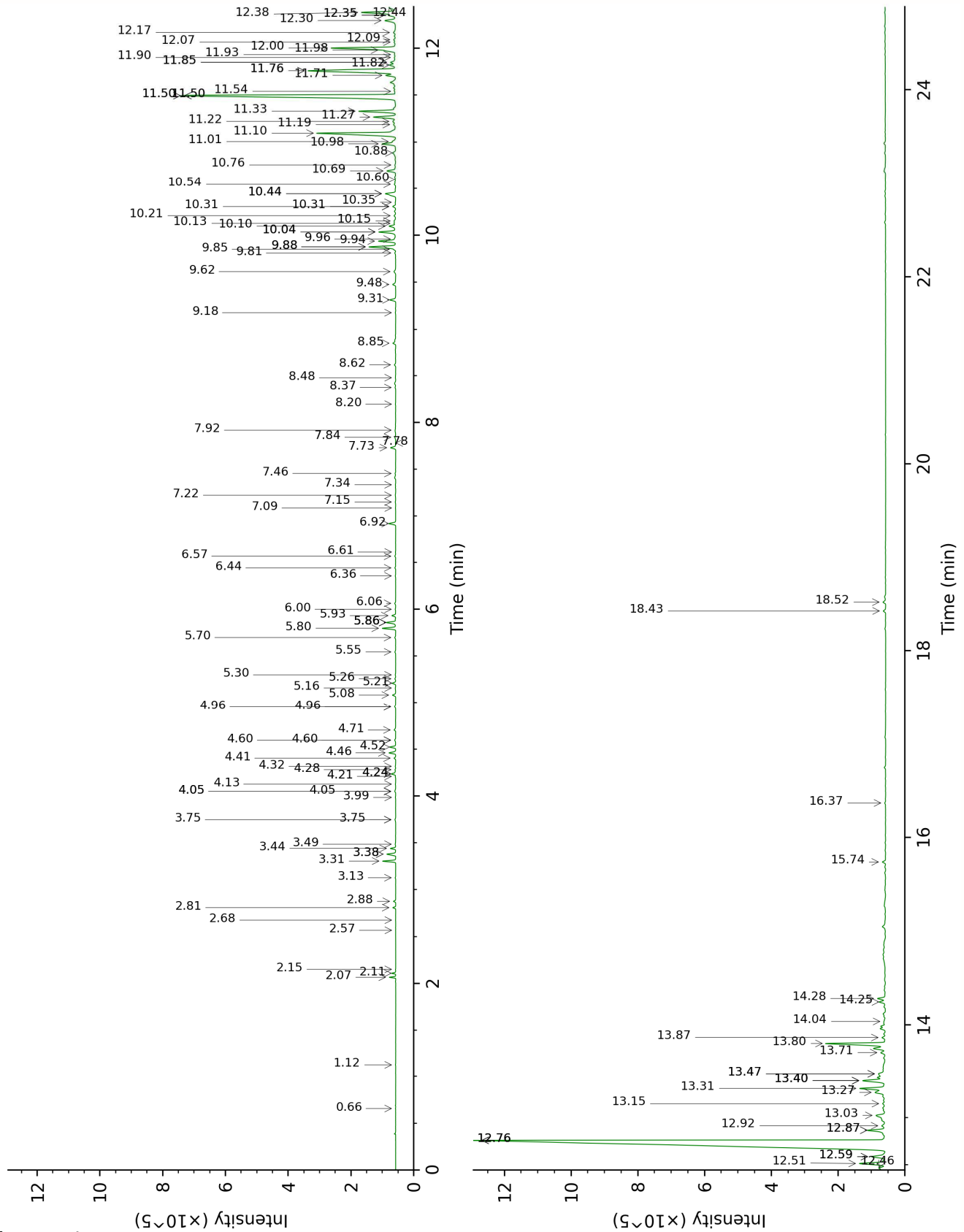
This page was intentionally left blank. The following pages present the complete data of the analysis.

DB-WAX





DB-5



FULL ANALYSIS DATA

2-Methylbutyral	Column DB-WAX			Column DB-5		
	0.82	879.1	tr	0.66	651.2	tr
Unknown ARPA I [m/z 85, 41 (62), 43 (47), 60 (22)... 95 (t)]	3.22	1144.8	0.01	1.12	748.4	0.01
Ethyl 2-methylbutyrate	1.81*	1023.0	[0.18]	2.07	849.7	0.15
Ethyl isovalerate	1.98	1039.2	0.10	2.11	853.2	0.10
Propyl isobutyrate	1.81*	1023.0	[0.18]	2.15	856.7	0.01
Nonene	1.04	920.5	tr	2.57	891.3	tr
( <i>cis?</i> )-5-Ethyl-2-methyl-2-vinyltetrahydrofuran	1.81*	1023.0	[0.18]	2.68	900.3	tr
( <i>trans?</i> )-5-Ethyl-2-methyl-2-vinyltetrahydrofuran	2.02	1042.4	0.08	2.81	910.0	0.08
( <i>trans?</i> )-2,5-Divinyl-2-methyltetrahydrofuran	2.91	1121.1	0.07	2.88	914.4	0.06
$\alpha$ -Pinene	1.50	991.1	0.02	3.13	931.1	0.02
Unknown ARPA II [m/z 111, 43 (87), 82 (72), 55 (70), 67 (63), 83 (56), 41 (44)... 125 (17), 140 (5)]	2.57	1095.6	0.38	3.31	942.8	0.38
5,5-Dimethyl-2(5H)-furanone	8.70	1546.8	0.12	3.38*	947.7	[0.32]
Propyl 2-methylbutyrate	2.77	1110.8	0.22	3.38*	947.7	[0.32]
Propyl isovalerate	2.98	1126.6	0.18	3.44	951.8	0.16
Benzaldehyde	7.54	1460.0	0.02	3.49	954.7	0.02
Sabinene	2.46	1084.5	0.01	3.75*	972.1	[0.03]
$\beta$ -Pinene	2.26	1066.3	0.02	3.75*	972.1	[0.03]
6-Methyl-5-hepten-2-one	5.30	1300.1	0.01	3.99	987.8	0.01
<i>trans</i> -Dehydroxylinalool oxide	3.62*	1174.9	[0.03]	4.05*	992.0	[0.05]
Myrcene	3.07	1133.1	0.01	4.05*	992.0	[0.05]
2-Pentylfuran	3.91	1196.6	0.02	4.05*	992.0	[0.05]
Unknown ARPA IV [m/z 43, 71 (48), 41 (46), 55 (39), 69 (38), 70 (38), 82 (35)... 140 (23)]	3.62*	1174.9	[0.03]	4.13	997.2	0.01
Octanal	4.68*	1250.5	[0.28]	4.21	1002.6	0.01
<i>cis</i> -Dehydroxylinalool oxide	4.02*	1204.6	[0.13]	4.24*	1004.4	[0.16]

Unknown ARPA V [m/z 43, 111 (99), 55 (82), 125 (75), 82 (73), 67 (68), 83 (62)... 140 (5)]	3.72	1182.5	0.14	4.24*	1004.4	[0.16]
Isobutyl 2-methylbutyrate	3.28	1149.1	0.03	4.28	1007.2	0.02
Isobutyl isovalerate	3.50	1165.5	0.04	4.32	1009.5	0.02
$\alpha$ -Terpinene	3.15	1139.7	0.08	4.40	1015.0	0.05
Unknown ARPA VI [m/z 111, 43 (92), 55 (75), 82 (70), 67 (65), 83 (58), 41 (48), 81 (41)... 125 (15), 140 (4)]	4.12	1211.4	0.22	4.46	1018.5	0.23
<i>para</i> -Cymene	4.33	1225.9	0.20	4.52	1022.4	0.19
Limonene	3.39	1157.4	0.02	4.60*	1027.0	[0.05]
1,8-Cineole	3.53	1168.4	0.04	4.60*	1027.0	[0.05]
Lavender lactone	9.43	1602.8	0.08	4.71	1033.8	0.06
<i>cis</i> -Arbusculone	6.74	1401.1	0.04	4.96*	1049.5	[0.05]
( <i>E</i> )- $\beta$ -Ocimene	4.20	1216.9	0.02	4.96*	1049.5	[0.05]
$\gamma$ -Terpinene	4.02*	1204.6	[0.13]	5.08	1057.6	0.11
<i>cis</i> -Arbusculol				5.16	1062.4	0.02
<i>cis</i> -Sabinene hydrate	7.14	1430.7	0.13	5.21	1065.6	0.12
<i>trans</i> -Arbusculone	7.29*	1441.5	[0.27]	5.26	1068.5	0.03
<i>trans</i> -Arbusculol	7.49	1456.4	0.02	5.30	1071.1	0.02
Terpinolene	4.51	1238.7	0.03	5.55	1086.5	0.03
<i>trans</i> -Sabinene hydrate	8.21	1509.5	0.04	5.70	1096.1	0.04
Linalool	8.31*	1516.7	[0.47]	5.80	1102.4	0.46
Hotrienol	9.05	1573.6	0.08	5.86*	1106.3	[0.35]
Nonanal	6.11	1357.0	0.03	5.86*	1106.3	[0.35]
2-Methylbutyl 2-methylbutyrate	4.68*	1250.5	[0.28]	5.86*	1106.3	[0.35]
2-Methylbutyl isovalerate	4.92	1267.1	0.14	5.93	1111.0	0.13
3-Methyl-3-butenyl 2-methylbutyrate	5.66	1325.3	0.02	6.00	1115.3	0.02
<i>cis-para</i> -Menth-2-en-1-ol	8.37	1521.7	0.04	6.06	1119.3	0.04
<i>trans-para</i> -Menth-2-en-1-ol	9.16	1582.0	0.01	6.36	1138.0	0.02
Prenyl 2-methylbutyrate	6.21	1363.8	0.03	6.44	1143.5	0.03
Prenyl isovalerate	6.52	1385.6	0.04	6.57	1151.4	0.04
Nerol oxide	7.09*	1427.3	[0.04]	6.61	1154.3	0.01
Terpinen-4-ol	8.82	1556.3	0.27	6.92	1174.2	0.25

<i>para</i> -Cymen-8-ol	11.84	1801.2	0.02	7.09	1184.9	0.01
$\alpha$ -Terpineol	10.03*	1651.3	[1.03]	7.15	1189.0	0.02
<i>cis</i> -Piperitol	9.87	1638.3	0.05	7.22	1193.5	0.03
Ethyl octanoate	6.78	1404.5	0.02	7.34	1200.7	0.01
<i>cis</i> -nor-Davanone?	8.89*	1561.6	[0.08]	7.46	1208.6	0.03
nor-Davanone	9.39	1600.1	0.21	7.73	1227.0	0.20
(3 <i>Z</i> )-Hexenyl 2-methylbutyrate	7.29*	1441.5	[0.27]	7.78	1230.4	0.02
(3 <i>Z</i> )-Hexenyl isovalerate	7.38*	1447.9	[0.09]	7.84	1234.4	0.06
Neral	9.73	1627.2	0.03	7.92	1239.5	0.04
Geraniol	11.91	1807.5	0.02	8.20	1258.1	0.01
Geranial	10.40	1680.4	0.05	8.37	1270.0	0.03
Methyl hydrocinnamate	11.74	1792.6	0.06	8.48	1277.0	0.01
Cogeijerene	8.31*	1516.7	[0.47]	8.62	1286.2	0.05
Methyl ( <i>Z</i> )-cinnamate	12.84	1889.0	0.26	8.85	1302.1	0.12
Bicycloelemene analog	7.00	1420.3	0.03	9.18	1324.9	0.02
Bicycloelemene	7.29*	1441.5	[0.27]	9.32	1334.5	0.23
Ethyl hydrocinnamate	12.10*	1824.3	[0.14]	9.48	1346.1	0.13
Eugenol	15.08*	2099.4	[0.13]	9.62	1355.7	0.10
Isoledene	7.09*	1427.3	[0.04]	9.81	1369.7	0.04
$\alpha$ -Copaene	7.38*	1447.9	[0.09]	9.85	1372.5	0.04
Ethyl ( <i>Z</i> )-cinnamate	13.26	1927.4	1.10	9.88*	1374.3	[1.07]
Modhephene	7.72	1473.0	0.04	9.88*	1374.3	[1.07]
Methyl ( <i>E</i> )-cinnamate	14.08*	2003.1	[0.76]	9.94	1378.6	0.64
<i>cis</i> - $\beta$ -Elemene	8.54	1534.6	0.01	9.96	1380.1	0.04
Geranyl acetate	10.80	1713.7	0.71	10.04*	1385.5	[0.71]
$\beta$ -Cubebene	8.00	1493.5	0.01	10.04*	1385.5	[0.71]
$\beta$ -Elemene	8.67*	1544.6	[0.45]	10.10	1389.8	0.29
Benzyl isovalerate	12.10*	1824.3	[0.14]	10.13	1391.9	0.06
<i>trans</i> -erythro-Davanafuran?	10.22	1666.0	0.07	10.15	1393.7	0.07
<i>cis</i> -erythro-Davanafuran?	10.34*	1676.0	[10.02]	10.21	1397.7	0.06
$\alpha$ -Gurjunene	7.85	1482.3	0.07	10.31*	1404.6	[0.11]
<i>trans</i> -threo-Davanafuran?	10.51	1689.0	0.04	10.31*	1404.6	[0.11]
$\beta$ -Maaliene	7.98	1491.7	0.04	10.35	1407.9	0.02
<i>cis</i> -threo-Davanafuran	10.98	1728.7	0.45	10.44*	1414.7	[0.55]
$\beta$ -Caryophyllene	8.67*	1544.6	[0.45]	10.44*	1414.7	[0.55]
$\beta$ -Copaene	8.59	1538.7	0.04	10.54	1422.3	0.05
$\alpha$ -Maaliene	8.89*	1561.6	[0.08]	10.60	1426.2	0.01
Aromadendrene	8.79	1553.5	0.36	10.69	1433.3	0.34

Selina-5,11-diene	9.01	1570.2	0.02	10.76	1438.0	0.04
$\alpha$ -Humulene	9.53	1610.8	0.14	10.88	1447.6	0.14
allo-Aromadendrene	9.25	1588.7	0.54	10.98	1454.8	0.60
Unknown ARPA VII [m/z 153, 43 (78), 111 (62), 96 (58), 95 (50), 67 (44), 109 (42)...]	11.93	1809.4	0.09	11.01	1456.6	0.13
Ethyl ( <i>E</i> )-cinnamate	14.67*	2059.3	[5.27]	11.10	1463.4	3.80
Selina-4,11-diene	9.67†	1622.3	0.03	11.19*†	1470.2	[0.10]
$\gamma$ -Muurolole	9.84†	1635.3	0.09	11.22*†	1472.7	[0.14]
Germacrene D	10.03*	1651.3	[1.03]	11.27	1476.1	0.91
$\beta$ -Selinene	10.13	1658.6	1.48	11.33	1480.8	1.44
$\alpha$ -Selinene	10.20	1664.1	0.06	11.50*	1493.2	[11.83]
Davana ether isomer I	12.32	1842.8	1.56	11.50*	1493.2	[11.83]
Bicyclogermacrene	10.34*	1676.0	[10.02]	11.50*	1493.2	[11.83]
Viridiflorene	9.90	1640.5	0.62	11.50*	1493.2	[11.83]
$\alpha$ -Muurolole	10.34*	1676.0	[10.02]	11.54	1496.6	0.03
$\gamma$ -Cadinene	10.65*	1700.6	[0.45]	11.71	1509.6	0.37
Davana ether isomer II	12.66	1873.2	4.13	11.76*	1513.4	[5.08]
Davana ether isomer III	12.73	1878.9	1.11	11.76*	1513.4	[5.08]
Artedouglasia oxide C	13.24*	1925.2	[1.46]	11.82	1518.2	0.26
Zonarene	10.65*	1700.6	[0.45]	11.85*	1520.5	[0.22]
$\delta$ -Cadinene	10.66	1702.0	0.15	11.85*	1520.5	[0.22]
Laciniata furanone G?				11.90	1524.6	0.05
Laciniata furanone F?				11.93	1527.0	0.05
Artedouglasia oxide A	13.55	1954.1	0.40	11.98	1530.7	0.34
Davana ether isomer IV	13.10	1912.3	2.57	12.00	1532.5	2.58
$\alpha$ -Calacorene	12.38	1848.4	0.04	12.07	1537.4	0.05
Laciniata furanone <i>E</i> ?				12.09	1539.4	0.07
Laciniata furanone H				12.17	1545.4	0.10
Davanone A	13.04	1906.8	0.52	12.30	1555.5	0.52
Artedouglasia oxide D	14.08*	2003.1	[0.76]	12.35*	1560.0	[0.25]
Geranyl butyrate	12.46	1855.5	0.02	12.35*	1560.0	[0.25]
Davanone B	13.24*	1925.2	[1.46]	12.38	1562.3	1.42
( <i>E</i> )-Nerolidol	14.03	1998.3	0.40	12.44	1566.5	0.37
Davanone C	13.46	1945.6	0.32	12.46	1568.6	0.28
Spathulenol	14.67*	2059.3	[5.27]	12.51	1572.4	1.13
Artedouglasia oxide B	14.74	2066.2	0.19	12.59*	1578.2	[0.70]
Globulol	14.17	2011.7	0.40	12.59*	1578.2	[0.70]
Cubeban-11-ol	13.97	1992.8	0.17	12.76*	1591.9	[41.43]
Davanone D	13.87*	1983.3	[41.59]	12.76*	1591.9	[41.43]
Viridiflorol	14.25	2019.8	0.27	12.76*	1591.9	[41.43]
Unknown ARPA VIII [m/z 43, 219 (54), 41	13.87*	1983.3	[41.59]	12.87*	1600.4	[0.84]

(51), 234 (49)]						
Eudesm-5-en-11-ol	14.67*	2059.3	[5.27]	12.87*	1600.4	[0.84]
Unknown ARPA IX [m/z 207, 43 (87), 83 (84), 149 (77), 109 (70), 81 (66), 93 (54)... 235 (10)...]	14.84	2075.7	0.23	12.92	1604.1	0.20
Davanol D isomer I	14.72	2064.1	0.42	13.03	1613.2	0.56
Unknown ARPA X [m/z 83, 55 (25), 111 (22), 93 (19), 43 (12)...]	14.79	2070.8	0.10	13.15	1623.6	0.14
Isospathulenol	15.70	2160.9	0.31	13.27	1633.6	0.61
τ-Cadinol	15.15	2106.7	1.11	13.32	1637.1	1.11
Unknown ARPA XI [m/z 95, 67 (56), 93 (41), 41 (37), 55 (34), 109 (30), 111 (24)...]	15.42	2133.6	0.50	13.40*	1644.0	[1.14]
Methyl <i>cis</i> -jasmonate	16.66	2260.0	0.23	13.40*	1644.0	[1.14]
β-Eudesmol	15.66	2157.6	0.57	13.40*	1644.0	[1.14]
5-Hydroxy-6-methyl-2- (5-methyl-5- vinyltetrahydrofuran- 2-yl)hepta-4,6-dien-3- one, isomer II	15.31	2122.6	0.19	13.47*	1650.0	[0.50]
α-Cadinol	15.74	2165.2	0.21	13.47*	1650.0	[0.50]
Unknown ARPA XII [m/z 153, 43 (51), 93 (25), 111 (23), 109 (22), 41 (22), 55 (19), 123 (17)... 250 (9)]				13.70	1669.3	0.22
Unknown ARPA XIV [m/z 163, 43 (56), 121 (14), 105 (13), 164 (13)... 234 (3)...]	17.14	2310.2	2.37	13.80	1677.4	2.50
Davanyl acetate				13.87	1682.6	0.14
Davanonol isomer	18.87	2501.6	0.09	14.04	1696.8	0.08
Unknown ARPA XV [m/z 43, 71 (88), 93 (86), 41 (74), 55 (73), 81 (71), 95 (59), 91 (53), 67 (52)... 220 (13)... 236? (t)]	17.64	2364.8	0.06	14.25	1714.5	0.26
β-Davanon-2-ol	19.12	2530.9	0.29	14.28	1717.7	0.32
Phytone	14.98	2089.9	0.02	15.74	1846.0	0.13
Nonadecane	12.96	1900.3	0.02	16.37	1903.3	0.04

Heneicosane	15.08*	2099.4	[0.13]	18.42	2103.1	0.11
Phytol	19.45	2569.4	0.16	18.52	2112.8	0.13
Total reported		91.95%			92.64%	

\*: 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