

Date : July 4, 2023

CERTIFICATE OF ANALYSIS – GC PROFILING & CHECK FOR CARRIER OIL

SAMPLE IDENTIFICATION

Internal code : 23D20-PTH01

Customer identification : Ylang Ylang Complete - Madagascar - Y10111

Type : Essential oil

Source : *Cananga odorata* var. *genuina* (Ylang-ylang)

Customer : Plant Therapy

ANALYSIS

Method: PC-MAT-014  - Analysis of the composition of an essential oil or other volatile liquid by FAST GC-FID (in French); identifications validated by GC-MS. Carrier oil determination by PC-MAT-010 – GC-FID quantitation of fatty acid methyl esters after derivatization, against an internal standard of tridecanoic acid.

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

Analysis date : April 26, 2023

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Notes: This report may not be published, including online, without the written consent from Laboratoire PhytoChemia.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

This report is an update from the first version issued on April 26, 2023, to update the customer identification.

PYHSICOCHEMICAL DATA

Physical aspect: Light yellow liquid

Refractive index: 1.5002 ± 0.0003 (20 °C; method PC-MAT-016)

CARRIER OIL DETERMINATION

After derivatization, <0.1% m/m of fatty acids, in equivalents of tridecanoic acid, were found.

CONCLUSION

No adulterant, contaminant or diluent has been detected using this method. Absence of fatty acids indicate that the oil is undiluted with a carrier oil.

ANALYSIS SUMMARY – CONSOLIDATED CONTENTS

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

Identification	%	Class
2-Methyl-3-buten-2-ol	0.01	Aliphatic alcohol
Ethyl acetate	0.04	Aliphatic ester
Pentanal	tr	Aliphatic aldehyde
Isobutyl acetate	tr	Aliphatic ester
Prenol	0.01	Aliphatic alcohol
Hexanal	tr	Aliphatic aldehyde
Octane	0.01	Alkane
Butyl acetate	0.01	Aliphatic ester
(3Z)-Hexenol	tr	Aliphatic alcohol
Hexanol	tr	Aliphatic alcohol
Isoamyl acetate	0.02	Aliphatic ester
2-Methylbutyl acetate	0.02	Aliphatic ester
3-Methyl-3-butenyl acetate	0.06	Aliphatic ester
Prenyl acetate	0.20	Aliphatic ester
α-Pinene	0.22	Monoterpene
Campheine	0.01	Monoterpene
Benzaldehyde	0.03	Simple phenolic
Sabinene	0.01	Monoterpene
β-Pinene	0.08	Monoterpene
6-Methyl-5-hepten-2-one	0.01	Aliphatic ketone
Myrcene	0.10	Monoterpene
(3Z)-Hexenyl acetate	0.07	Aliphatic ester
para-Methylanisole	4.29	Simple phenolic
Limonene	0.04	Monoterpene
1,8-Cineole	0.25	Monoterpenic ether
Benzyl alcohol	0.07	Simple phenolic
(Z)-β-Ocimene	0.01	Monoterpene
(E)-β-Ocimene	0.02	Monoterpene
cis-Sabinene hydrate	0.01	Monoterpenic alcohol
cis-Linalool oxide (fur.)	0.02	Monoterpenic alcohol
trans-Linalool oxide (fur.)	0.03	Monoterpenic alcohol
Methyl benzoate	2.24	Phenolic ester
para-Cresol	0.07	Simple phenolic
Linalool	10.25	Monoterpenic alcohol
Camphor	0.02	Monoterpenic ketone
ortho-Dimethoxybenzene	0.02	Simple phenolic
Benzyl acetate	2.32	Phenolic ester
para-Cresyl acetate	0.04	Phenolic ester
Ethyl benzoate	0.07	Phenolic ester
Terpinen-4-ol	0.01	Monoterpenic alcohol
α-Terpineol	0.21	Monoterpenic alcohol
Methyl salicylate	0.02	Phenolic ester
Methylchavicol	0.16	Phenylpropanoid
Nerol	0.04	Monoterpenic alcohol
Neral	0.04	Monoterpenic aldehyde

Phenylethyl acetate	0.05	Phenolic ester
Geraniol	1.77	Monoterpenic alcohol
Chavicol	0.01	Phenylpropanoid
Geranal	0.11	Monoterpenic aldehyde
(E)-Anethole	0.07	Phenylpropanoid
1-Nitro-2-phenylethane	0.09	Simple phenolic
(E)-Cinnamyl alcohol	0.02	Phenylpropanoid
4-Vinylguaiacol	0.03	Simple phenolic
δ-Elemene	0.04	Sesquiterpene
Bicycloelemene	0.02	Sesquiterpene
Benzyl butyrate	0.05	Phenolic ester
α-Cubebene	0.14	Sesquiterpene
Eugenol	0.70	Phenylpropanoid
Neryl acetate	0.04	Monoterpenic ester
α-Ylangene	0.14	Sesquiterpene
Hydrocinnamyl acetate	0.04	Phenylpropanoid ester
α-Copaene	0.89	Sesquiterpene
β-Bourbonene	0.03	Sesquiterpene
Geranyl acetate	8.48	Monoterpenic ester
β-Cubebene	0.13	Sesquiterpene
β-Elemene	0.35	Sesquiterpene
Cyperene	0.04	Sesquiterpene
Isocaryophyllene	0.01	Sesquiterpene
Methyleugenol	0.04	Phenylpropanoid
β-Caryophyllene	11.43	Sesquiterpene
Caryophylla-4(12),8(13)-diene	0.07	Sesquiterpene
β-Copaene	0.43	Sesquiterpene
Aromadendrene	0.02	Sesquiterpene
α-Guaiene	0.04	Sesquiterpene
Isogermacrene D	0.08	Sesquiterpene
(E)-Cinnamyl acetate	0.69	Phenylpropanoid ester
trans-Muurola-3,5-diene	0.11	Sesquiterpene
α-Humulene	3.12	Sesquiterpene
ε-Muurolene?	0.08	Sesquiterpene
(E)-Isoeugenol	0.18	Phenylpropanoid
allo-Aromadendrene	0.05	Sesquiterpene
cis-Cadina-1(6),4-diene	0.01	Sesquiterpene
cis-Muurola-4(15),5-diene	0.14	Sesquiterpene
trans-Cadina-1(6),4-diene	0.21	Sesquiterpene
γ-Muurolene	1.88	Sesquiterpene
Germacrene D	12.46	Sesquiterpene
trans-Muurola-4(15),5-diene	0.50	Sesquiterpene
γ-Amorphene	0.04	Sesquiterpene
Prenyl benzoate	0.41	Phenolic ester
Bicyclogermacrene	0.15	Sesquiterpene
Viridiflorene	0.07	Sesquiterpene
epi-Cubebol	0.06	Sesquiterpenic alcohol
Methyl (E)-iseugenol	0.06	Phenylpropanoid
α-Muurolene	0.89	Sesquiterpene
(3Z,6E)-α-Farnesene	0.03	Sesquiterpene
δ-Amorphene	0.31	Sesquiterpene
Unknown	1.01	Sesquiterpene

Laboratoire
PhytoChemia

Plus que des analyses... des conseils

δ -Guaiene	0.48	Sesquiterpene
(3E,6E)- α -Farnesene	5.43	Sesquiterpene
Cubebol	0.10	Sesquiterpenic alcohol
γ -Cadinene	1.08	Sesquiterpene
δ -Cadinene	2.86	Sesquiterpene
<i>trans</i> -Calamenene	0.13	Sesquiterpene
<i>trans</i> -Cadina-1,4-diene	0.18	Sesquiterpene
α -Cadinene	0.26	Sesquiterpene
α -Calacorene	0.05	Sesquiterpene
<i>cis</i> -Dracunculifoliol	0.02	Sesquiterpenic alcohol
α -Elemol	0.10	Sesquiterpenic alcohol
Germacrene B	0.01	Sesquiterpene
Salviadienol?	0.02	Sesquiterpenic alcohol
Sesquirosefuran?	0.03	Sesquiterpenic ether
(E)-Nerolidol	0.14	Sesquiterpenic alcohol
(3Z)-Hexenyl benzoate	0.03	Phenolic ester
Germacrene D-4-ol	0.06	Sesquiterpenic alcohol
Spathulenol	0.07	Sesquiterpenic alcohol
<i>trans</i> -Dracunculifoliol	0.10*	Sesquiterpenic alcohol
Caryophyllene oxide	0.25	Sesquiterpenic ether
10-epi-Junenol	0.10*	Sesquiterpenic alcohol
Unknown	0.05	Sesquiterpenic alcohol
Globulol	0.05	Sesquiterpenic alcohol
Unknown	0.09	Oxygenated sesquiterpene
Viridiflorol	0.03	Sesquiterpenic alcohol
Guaiol	0.07	Sesquiterpenic alcohol
Copaborneol	0.06	Sesquiterpenic alcohol
Humulene epoxide II	0.08	Sesquiterpenic ether
10-epi-Cubenol	0.05	Sesquiterpenic alcohol
Junenol	0.30	Sesquiterpenic alcohol
(E)-Isoeugenyl acetate	0.02	Phenylpropanoid ester
Unknown	0.06	Oxygenated sesquiterpene
1-epi-Cubenol	0.25	Sesquiterpenic alcohol
γ -Eudesmol	0.14	Sesquiterpenic alcohol
Caryophylladienol II	0.02	Sesquiterpenic alcohol
τ -Cadinol	0.30	Sesquiterpenic alcohol
τ -Muurolol	0.74	Sesquiterpenic alcohol
Cubenol	0.09	Sesquiterpenic alcohol
α -Muurolol	0.40	Sesquiterpenic alcohol
Unknown	0.24	Sesquiterpenic alcohol
α -Cadinol	1.47	Sesquiterpenic alcohol
<i>cis</i> -Calamenen-10-ol	0.05	Sesquiterpenic alcohol
<i>trans</i> -Calamenen-10-ol	0.03	Sesquiterpenic alcohol
Bulnesol	0.04	Sesquiterpenic alcohol
Unknown	0.18	Oxygenated sesquiterpene
(2Z,6E)-Farnesol	0.02	Sesquiterpenic alcohol
(2E,6E)-Farnesol	1.86	Sesquiterpenic alcohol
(2E,6E)-Farnesal	0.04	Sesquiterpenic aldehyde
Benzyl benzoate	6.78	Phenolic ester
Unknown	0.03	Unknown
(2E,6E)-Farnesyl acetate	1.49	Sesquiterpenic ester
Unknown	0.02	Oxygenated sesquiterpene

Laboratoire
PhytoChemia

Plus que des analyses... des conseils

Benzyl salicylate	1.79	Phenolic ester
Unknown	0.01	Unknown
Unknown	0.02	Unknown
Geranyl benzoate	0.15	Phenolic ester
Unknown	0.01	Unknown
Unknown	0.11	Unknown
Unknown	0.01	Unknown
Unknown	0.02	Unknown
Unknown	0.05	Unknown
Consolidated total	97.58%	

*: 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

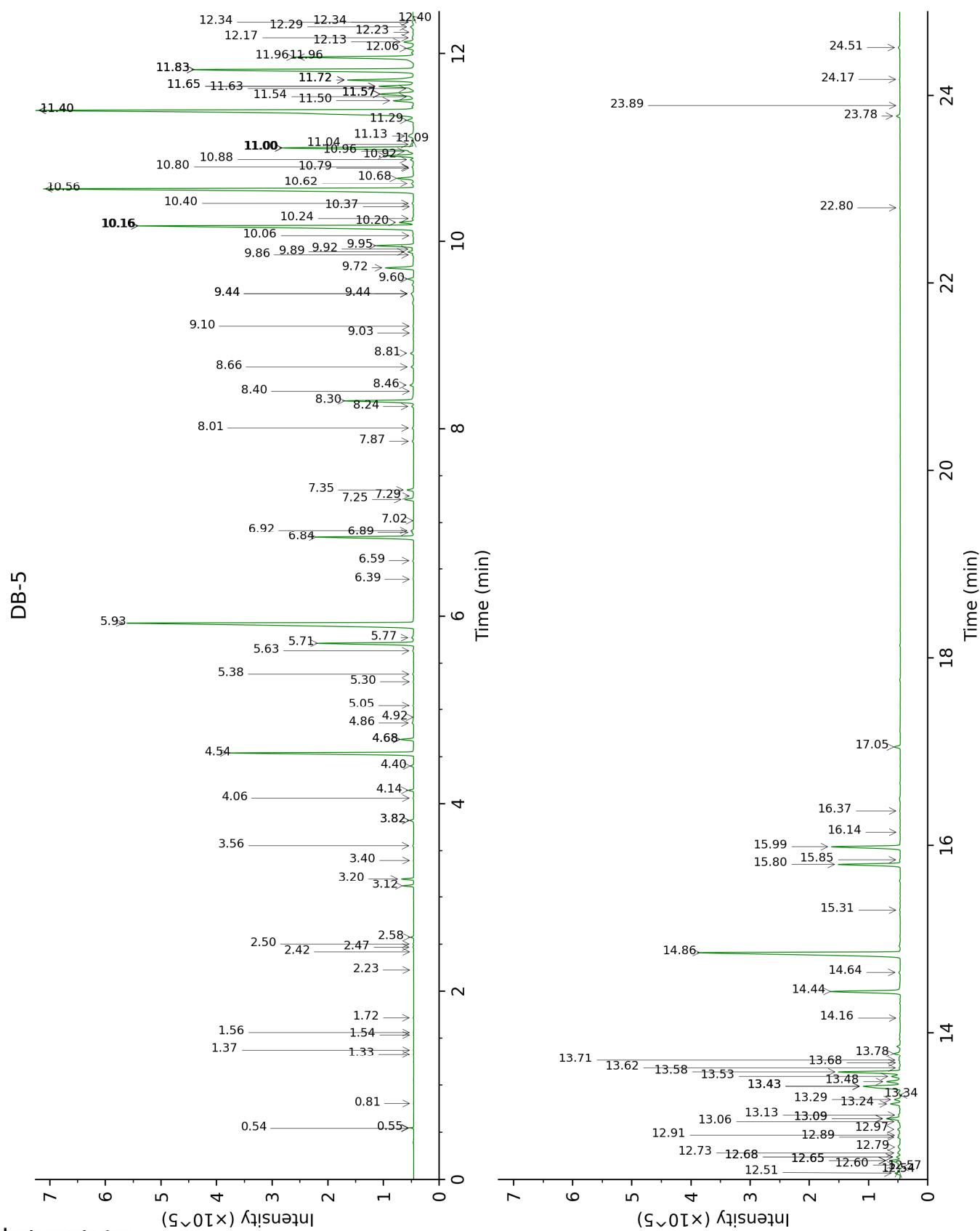
tr: The compound has been detected below 0.005% of total signal.

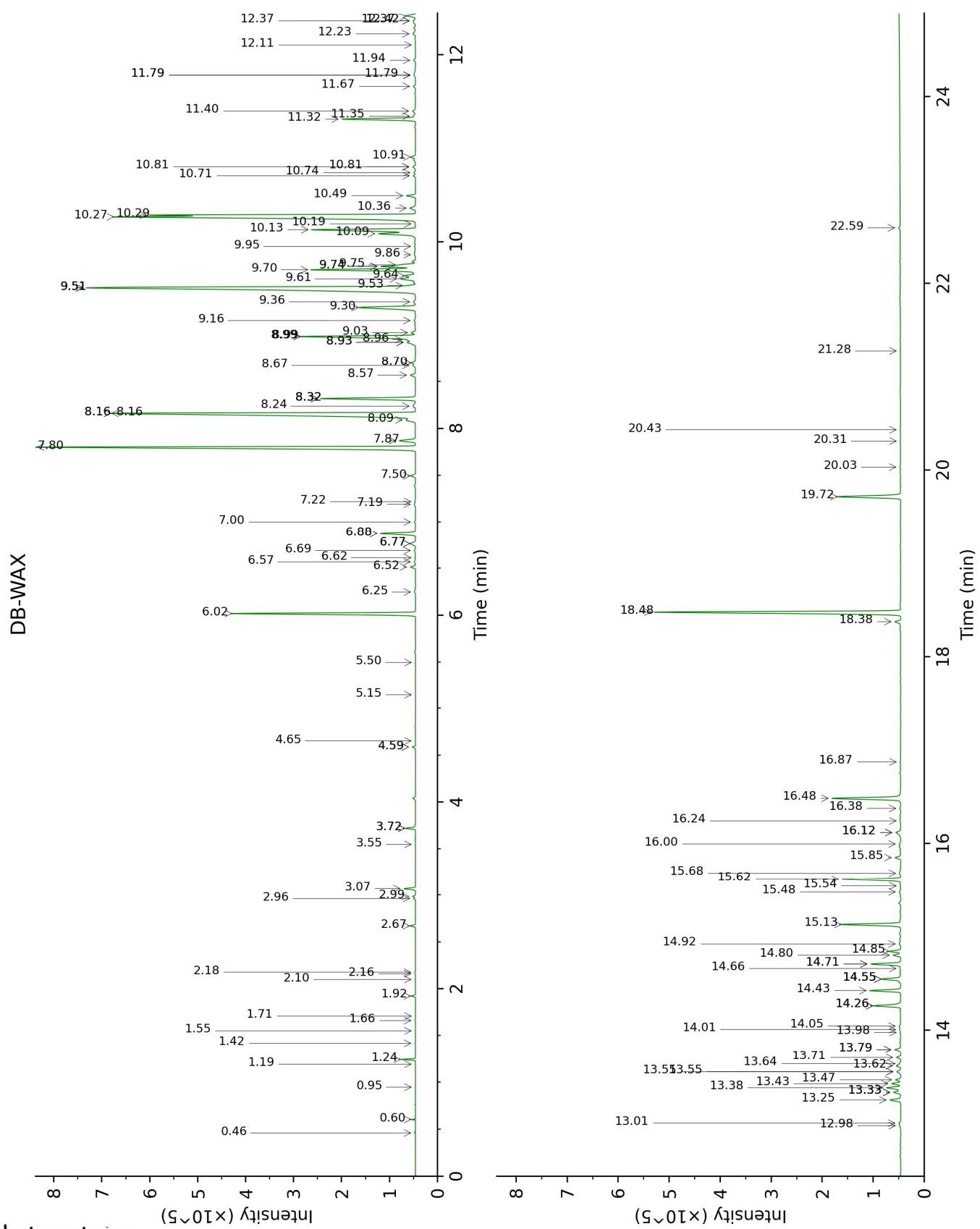
Note: no correction factor was applied

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.

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





FULL ANALYSIS DATA

Identification	Column DB-5			Column DB-WAX		
	R.T	R.I	%	R.T	R.I	%
2-Methyl-3-buten-2-ol	0.54	606	0.01	1.42	1014	0.01
Ethyl acetate	0.55	608	0.04	0.60	854	0.03
Pentanal	0.80	693	tr	0.95	941	tr
Isobutyl acetate	1.33	771	tr	1.19	983	tr
Prenol	1.37	777	0.01	4.59*	1288	0.07
Hexanal	1.54	799	tr	1.71	1044	tr
Octane	1.56	803	0.01	0.46	786	0.01
Butyl acetate	1.72	816	0.01	1.66	1039	0.01
(3Z)-Hexenol	2.23	858	tr	5.50	1347	tr
Hexanol	2.42	874	tr	5.15	1321	tr
Isoamyl acetate	2.47	878	0.02	2.18	1092	0.02
2-Methylbutyl acetate	2.50	880	0.02	2.16	1091	0.01
3-Methyl-3-butenyl acetate	2.58	886	0.06	2.99	1160	0.06
Prenyl acetate	3.12	926	0.20	3.72*	1220	0.22
α-Pinene	3.20	931	0.22	1.24	992	0.22
Camphepane	3.40	944	0.01	1.55	1027	0.01
Benzaldehyde	3.56	954	0.03	7.00	1458	0.03
Sabinene	3.82*	972	0.09	2.10	1084	0.01
β-Pinene	3.82*	972	[0.09]	1.92	1066	0.08
6-Methyl-5-hepten-2-one	4.06	987	0.01	4.65	1293	0.01
Myrcene	4.14	993	0.10	2.67	1135	0.11
(3Z)-Hexenyl acetate	4.40	1010	0.07	4.59*	1288	[0.07]
para-Methylanisole	4.54	1018	4.29	6.02	1385	4.26
Limonene	4.68*	1027	0.29	2.96	1159	0.04
1,8-Cineole	4.68*	1027	[0.29]	3.07	1167	0.25
Benzyl alcohol	4.86	1038	0.07	11.40	1816	0.10
(Z)-β-Ocimene	4.92	1042	0.01	3.55	1207	0.01
(E)-β-Ocimene	5.05	1050	0.02	3.72*	1220	[0.22]
cis-Sabinene hydrate	5.30	1066	0.01	6.62	1429	0.02
cis-Linalool oxide (fur.)	5.38	1071	0.02	6.25	1401	0.04
trans-Linalool oxide (fur.)	5.63	1086	0.03	6.57	1426	0.04
Methyl benzoate	5.71	1091	2.24	8.32*	1560	2.29
para-Cresol	5.77	1095	0.07	13.55*	2014	0.14
Linalool	5.93	1105	10.25	7.80	1519	10.20
Camphor	6.39	1134	0.02	6.88*	1449	0.88
ortho-Dimethoxybenzene	6.59	1147	0.02			
Benzyl acetate	6.84	1163	2.32	9.70	1671	2.65
para-Cresyl acetate	6.89	1166	0.04			
Ethyl benzoate	6.92	1168	0.07	8.99*	1613	3.19

Laboratoire
PhytoChemia

Plus que des analyses... des conseils

Terpinen-4-ol	7.02	1174	0.01	8.32*	1560	[2.29]
α-Terpineol	7.25	1189	0.21	9.51*	1655	14.14
Methyl salicylate	7.29	1191	0.02	10.19	1712	0.02
Methylchavicol	7.35	1195	0.16	8.96	1610	0.18
Nerol	7.87	1229	0.04	10.74	1759	0.04
Neral	8.01	1239	0.04	9.16	1627	0.04
Phenylethyl acetate	8.24	1254	0.05	10.71	1756	0.06
Geraniol	8.30	1258	1.77	11.32	1809	1.79
Chavicol	8.40	1265	0.01	16.12*	2273	0.19
Geranal	8.46	1269	0.11	9.75†	1676	[1.15]
(E)-Anethole	8.66	1282	0.07	10.81*	1765	0.08
1-Nitro-2-phenylethane	8.81	1292	0.09	13.79*	2038	0.18
(E)-Cinnamyl alcohol	9.03	1307	0.02	15.54	2212	0.02
4-Vinylguaiacol	9.10	1312	0.03	14.71*	2128	0.77
δ-Elemene	9.44*	1336	0.11	6.69	1435	0.04
Bicycloelemene	9.44*	1336	[0.11]	6.77*	1440	0.19
Benzyl butyrate	9.44*	1336	[0.11]	11.35	1812	0.05
α-Cubebene	9.60	1347	0.14	6.52	1422	0.13
Eugenol	9.72	1355	0.70	14.43	2099	0.74
Neryl acetate	9.86	1365	0.04	9.86	1684	0.03
α-Ylangene	9.89	1367	0.14	6.77*	1440	[0.19]
Hydrocinnamyl acetate	9.92	1369	0.04	12.11	1879	0.01
α-Copaene	9.95	1372	0.89	6.88*	1449	[0.88]
β-Bourbonene	10.06	1379	0.03	7.22	1475	0.02
Geranyl acetate	10.16*	1387	8.61	10.28†	1720	[14.16]
β-Cubebene	10.16*	1387	[8.61]	7.50	1495	0.13
β-Elemene	10.20	1389	0.35	8.09	1542	0.32
Cyperene	10.24	1392	0.04	7.19	1472	0.05
Isocaryophyllene	10.37	1401	0.01	7.87	1524	0.48
Methyleugenol	10.40	1404	0.04	12.98	1960	0.05
β-Caryophyllene	10.56	1415	11.43	8.16*	1547	11.36
Caryophylla-4(12),8(13)-diene	10.62	1420	0.07	8.32*	1560	[2.29]
β-Copaene	10.68	1424	0.43	8.16*	1547	[11.36]
Aromadendrene	10.79	1432	0.02	8.24	1553	0.12
α-Guaiene	10.80	1433	0.04	8.16*	1547	[11.36]
Isogermacrene D	10.88	1439	0.08	8.67	1588	0.07
(E)-Cinnamyl acetate	10.92	1442	0.69	14.26*	2083	0.76
trans-Muurola-3,5-diene	10.96	1445	0.11	8.57	1579	0.15
α-Humulene	11.00*	1448	3.40	8.99*	1613	[3.19]
ε-Muurolene?	11.00*	1448	[3.40]	8.93*	1608	0.29
(E)-Isoeugenol	11.00*	1448	[3.40]	16.12*	2273	[0.19]
allo-Aromadendrene	11.04	1451	0.05	8.70*	1590	0.12
cis-Cadin-1(6),4-diene	11.09	1454	0.01	8.70*	1590	[0.12]

<i>cis</i> -Muurola-4(15),5-diene	11.13	1457	0.14	9.03	1616	0.12
<i>trans</i> -Cadina-1(6),4-diene	11.29	1470	0.21	8.93*	1608	[0.29]
γ -Murolene	11.40*	1478	14.34	9.30	1638	1.88
Germacrene D	11.40*	1478	[14.34]	9.51*	1655	[14.14]
<i>trans</i> -Muurola-4(15),5-diene	11.50	1485	0.50	9.51*	1655	[14.14]
γ -Amorphene	11.54	1488	0.04	9.53	1657	0.08
Prenyl benzoate	11.57*	1490	1.10	13.38	1998	0.41
Bicyclogermacrene	11.57*	1490	[1.10]	9.74*†	1674	1.15
Viridiflorene	11.57*	1490	[1.10]	9.36	1643	0.07
epi-Cubebol	11.57*	1490	[1.10]	11.67	1840	0.06
Methyl (<i>E</i>)-isoeugenol	11.63	1495	0.06	14.66	2123	0.03
α -Murolene	11.66*	1496	0.92	9.74*†	1674	[1.15]
(3 <i>Z</i> ,6 <i>E</i>)- α -Farnesene	11.66*	1496	[0.92]	9.95	1692	0.03
δ -Amorphene	11.72*	1502	1.80	9.64	1666	0.31
Unknown [m/z 119, 41 (95), 123 (53), 80 (49), 161 (44), 105 (42)... 204 (2)]	11.72*	1502	[1.80]			
δ -Guaiene	11.72*	1502	[1.80]	9.60	1663	0.48
(3 <i>E</i> ,6 <i>E</i>)- α -Farnesene	11.83*	1510	6.62	10.27†	1718	14.16
Cubebol	11.83*	1510	[6.62]	12.23	1890	0.10
γ -Cadinene	11.83*	1510	[6.62]	10.09	1703	1.08
δ -Cadinene	11.96*	1520	3.04	10.13	1706	2.86
<i>trans</i> -Calamenene	11.96*	1520	[3.04]	10.91	1774	0.13
<i>trans</i> -Cadina-1,4-diene	12.06	1528	0.18	10.36	1726	0.19
α -Cadinene	12.13	1533	0.26	10.49	1738	0.23
α -Calacorene	12.17	1536	0.05	11.78*	1851	0.09
<i>cis</i> -Dracunculifoliol	12.23	1541	0.02	11.78*	1851	[0.09]
α -Elemol	12.29	1546	0.10	13.71	2029	0.11
Germacrene B	12.34*	1550	0.04	10.81*	1765	[0.08]
Salviadienol?	12.34*	1550	[0.04]	13.98	2055	0.02
Sesquirosefuran?	12.40	1555	0.03	11.78*	1851	[0.09]
(<i>E</i>)-Nerolidol	12.51	1563	0.14	13.47	2006	0.13
(3 <i>Z</i>)-Hexenyl benzoate	12.54	1566	0.03	14.01	2058	0.04
Germacrene D-4-ol	12.57	1568	0.06	13.33*	1993	0.20
Spathulenol	12.60	1570	0.07	14.05	2062	0.06
<i>trans</i> -Dracunculifoliol	12.65*	1574	0.35	12.36*	1902	0.11
Caryophyllene oxide	12.65*	1574	[0.35]	12.42	1908	0.25
10-epi-Junenol	12.65*	1574	[0.35]	12.36*	1902	[0.11]
Unknown cadinol or murolol analog [m/z 161, 119 (77),	12.68*	1577	0.10	11.94	1865	0.05

120 (76), 105 (73), 93 (57)... 204 (36)]						
Globulol	12.68*	1577	[0.10]	13.55*	2014	[0.14]
Unknown [m/z 161, 105 (84), 43 (80), 119 (72), 93 (62), 121 (54)... 204 (38), 222 (2)]	12.73	1580	0.09	13.62	2020	0.13
Viridiflorol	12.79	1585	0.03	13.64	2022	0.04
Guaiol	12.89	1593	0.07	13.79*	2038	[0.18]
Copaborneol	12.91	1595	0.06	14.55*	2111	0.50
Humulene epoxide II	12.98	1600	0.08	13.01	1963	0.06
10-epi-Cubenol	13.06	1606	0.05	13.33*	1993	[0.20]
Junenol	13.09*	1609	0.37	13.25	1986	0.30
(E)-Isoeugenyl acetate	13.09*	1609	[0.37]	16.87	2354	0.02
Unknown [m/z 179, 161 (66), 119 (44), 95 (38), 105 (35)... 204 (24), 222 (1)]	13.13	1612	0.06	14.26*	2083	[0.76]
1-epi-Cubenol	13.24	1622	0.25	13.43	2002	0.23
γ -Eudesmol	13.29	1625	0.14	14.55*	2111	[0.50]
Caryophylladienol II	13.34	1629	0.02	15.68	2227	0.03
τ -Cadinol	13.43*	1637	1.34	14.55*	2111	[0.50]
τ -Murolol	13.43*	1637	[1.34]	14.71*	2128	[0.77]
Cubenol	13.43*	1637	[1.34]	13.33*	1993	[0.20]
α -Murolol	13.48	1641	0.40	14.84	2141	0.29
Unknown cadinol analog II [m/z 95, 121 (73), 43 (57), 79 (43), 161 (43), 109 (40)... 204 (35), 222 (2)]	13.53	1645	0.24	14.80	2137	0.20
α -Cadinol	13.58	1649	1.47	15.13	2170	1.49
cis-Calamenen-10-ol	13.62	1653	0.05	16.00	2260	0.04
trans-Calamenen-10-ol	13.68	1657	0.03	16.38	2300	0.03
Bulnesol	13.71	1660	0.04	14.92	2149	0.06
Unknown [m/z 123, 95 (31), 81 (29), 105 (27)... 222 (5)]	13.78	1666	0.18	15.85	2245	0.16
(2Z,6E)-Farnesol	14.16	1697	0.02	16.24	2286	0.02
(2E,6E)-Farnesol	14.44	1721	1.86	16.48	2312	1.91
(2E,6E)-Farnesal	14.64	1739	0.04	15.48	2206	0.05
Benzyl benzoate	14.86	1757	6.78	18.48	2536	6.72
Unknown [m/z 121, 107 (86), 81 (71), 93 (71), 59 (68), 43 (67)...]	15.31	1796	0.03			

(2E,6E)-Farnesyl acetate	15.80	1840	1.49	15.62	2220	1.48
Unknown [m/z 43, 107 (97), 81 (83), 121 (77), 123 (74), 93 (73)... 220 (26)...]	15.85	1845	0.02	20.03	2722	0.02
Benzyl salicylate	15.99	1857	1.79	19.72	2684	1.76
Unknown [m/z 91, 93 (98), 81 (92), 41 (92), 105 (86), 107 (86)...]	16.14	1871	0.01	20.43	2772	0.02
Unknown [m/z 123, 81 (96), 41 (74), 43 (64), 91 (62), 95 (57)...]	16.37	1892	0.02	20.31	2757	0.01
Geranyl benzoate	17.05	1956	0.15	18.38	2524	0.16
Unknown [m/z 326, 311 (47), 327 (22), 202 (16), 137 (15)...]	22.80	2575	0.01			
Unknown [m/z 326, 327 (22), 311 (17), 137 (8), 202 (7)...]	23.78	2696	0.11			
Unknown [m/z 69, 83 (59), 81 (56), 137 (48), 41 (29), 139 (28)...]	23.89	2710	0.01	21.28	2881	0.01
Unknown [m/z 69, 81 (44), 147 (31), 41 (26), 119 (26), 93 (24)...]	24.17	2746	0.02			
Unknown [m/z 69, 81 (42), 147 (26), 119 (25), 93 (25), 41 (24)...]	24.51	2790	0.05	22.59	3057	0.05
Total identified	97.47%			97.21%		
Total reported	98.33%			97.86%		

*: Two or more compounds are coeluting on this column

[xx]: Duplicate percentage due to coelutions, not taken into account in the consolidated total

t: 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