Date : 2024-08-23

Certificate of Analysis - GC Profiling

Sample Identification

Internal code : 24H09-NSO03
Customer Identification : Yuzu Oil - Lot No: K0H0611Z
Type : Essential Oil
Source : *Citrus junos*
Customer : Natural Sourcing LLC

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Notes: This report may not be published, including online, without the written consent from Laboratoire PhytoChemia.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 : Benoit Roger, Ph. D.
Date : 2024-08-20

Physicochemical data

Refractive index : 1.4735 ± 0.0003 (20 °C)
Method : PC-MAT-016 - Measure of the refractive index of a liquid.
Analyst : Cindy Caron B. Sc.
Date : 2024-08-09

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
α-Thujene	0.29	Monoterpene
α-Pinene	1.09	Monoterpene
Camphene	0.01	Monoterpene
β-Pinene	0.60	Monoterpene
Sabinene	0.17	Monoterpene
Myrcene	1.39	Monoterpene
Pseudolimonene	0.02	Monoterpene
α-Phellandrene	0.32	Monoterpene
Octanal	tr	Aliphatic aldehyde
Δ3-Carene	0.01	Monoterpene
α-Terpinene	0.15	Monoterpene
para-Cymene	0.98	Monoterpene
Limonene	79.93	Monoterpene
β-Phellandrene	2.50	Monoterpene
(*E*)-β-Ocimene	0.15	Monoterpene
γ-Terpinene	8.08	Monoterpene
para-Cymenene	0.04	Monoterpene
Terpinolene	0.32	Monoterpene
Linalool	1.72	Monoterpenic alcohol
trans-para-Mentha-2,8-dien-1-ol	0.02	Monoterpenic alcohol
cis-Limonene oxide	0.04	Monoterpenic ether
trans-Limonene oxide	0.04	Monoterpenic ether
trans-Sabinol	0.01	Monoterpenic alcohol
Epoxyterpinolene	0.01	Monoterpenic ether
Terpinen-4-ol	0.12	Monoterpenic alcohol
Cryptone	0.01	Normonoterpenic ketone
α-Terpineol	0.10	Monoterpenic alcohol
Decanal	0.02	Aliphatic aldehyde
trans-Carveol	0.02	Monoterpenic alcohol
cis-Carveol	0.02	Monoterpenic alcohol
Carvone	0.02	Monoterpenic ketone
Thymol	0.06	Monoterpenic alcohol
δ-Elemene	0.03	Sesquiterpene
α-Copaene	0.02	Sesquiterpene
β-Cubebene	0.02	Sesquiterpene
β-Elemene	0.02	Sesquiterpene
β-Caryophyllene	0.11	Sesquiterpene
γ-Elemene	0.01	Sesquiterpene
α-Humulene	0.03	Sesquiterpene
(*E*)-β-Farnesene	0.23	Sesquiterpene

Germacrene D	0.09	Sesquiterpene
Bicyclogermacrene	0.22	Sesquiterpene
α-Muurolene	0.02	Sesquiterpene
δ-Cadinene	0.04	Sesquiterpene
β-Sesquiphellandrene	0.05	Sesquiterpene
Germacrene B	0.07	Sesquiterpene
Spathulenol	0.20	Sesquiterpenic alcohol
Alismol	0.02	Sesquiterpenic alcohol
Isospathulenol	0.02	Sesquiterpenic alcohol
τ-Muurolol	0.01	Sesquiterpenic alcohol
τ-Cadinol	tr	Sesquiterpenic alcohol
α-Cadinol	0.02	Sesquiterpenic alcohol
Consolidated total	**99.48**	

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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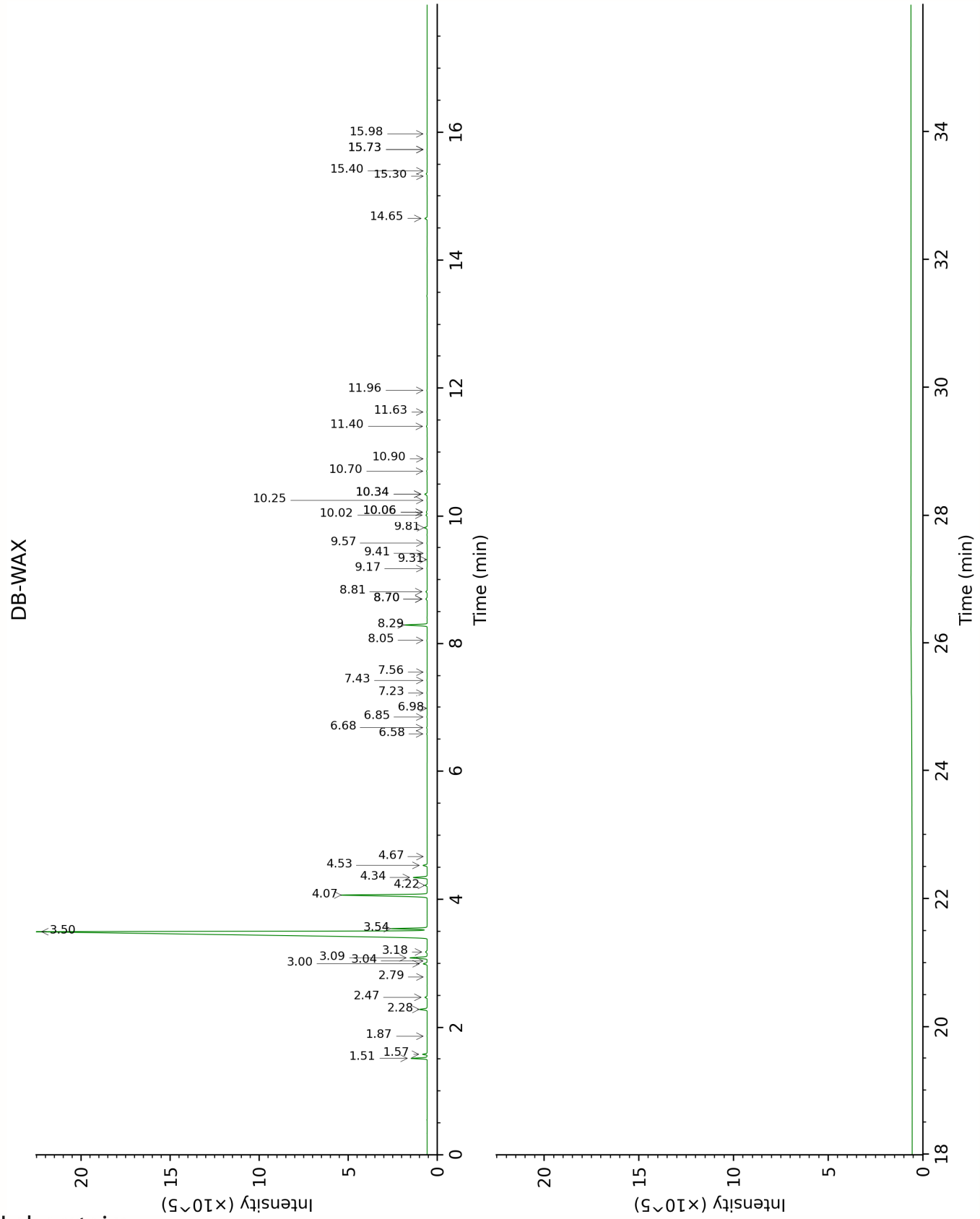
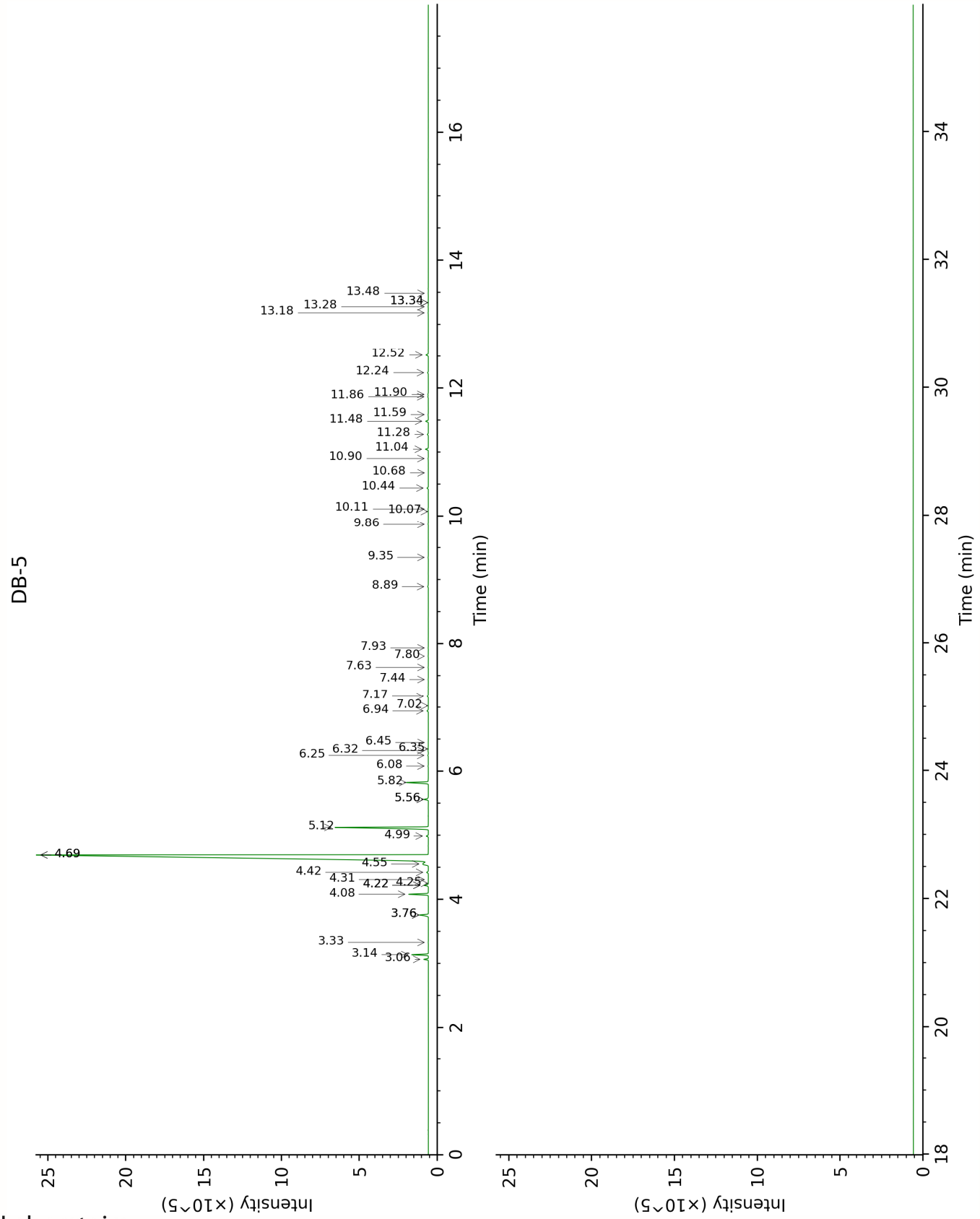

DB-WAX
Time (min)
Intensity (×10^5)
15.98
15.73
15.40
15.30
14.65
11.96
11.63
11.40
10.90
10.70
10.34
10.25
10.02
10.06
9.81
9.57
9.41
9.31
9.17
8.81
8.70
8.29
8.05
7.56
7.43
7.23
6.98
6.85
6.68
6.58
4.67
4.53
4.34
4.22
4.07
3.54
3.50
3.18
3.09
3.04
3.00
2.79
2.47
2.28
1.87
1.57
1.51


DB-5
Time (min)
Intensity (×10^5)
13.48
13.34
13.28
13.18
12.52
12.24
11.90
11.86
11.59
11.48
11.28
11.04
10.90
10.68
10.44
10.11
10.07
9.86
9.35
8.89
7.93
7.80
7.63
7.44
7.17
7.02
6.94
6.45
6.35
6.32
6.25
6.08
5.82
5.56
5.12
4.99
4.69
4.55
4.42
4.31
4.25
4.22
4.08
3.76
3.33
3.14
3.06

FULL ANALYSIS DATA

α-Thujene	**Column DB-WAX**			**Column DB-5**		
	1.57	999.8	0.30	3.06	925.7	0.29
α-Pinene	1.51	993.1	1.08	3.14	930.5	1.09
Camphene	1.87	1027.7	0.01	3.33	943.1	0.01
β-Pinene	2.28	1067.6	0.60	3.76*	971.3	[0.77]
Sabinene	2.47	1085.5	0.17	3.76*	971.3	[0.77]
Myrcene	3.09	1134.4	1.40	4.08	992.7	1.39
Pseudolimonene	3.04	1130.6	0.02	4.22*	1002.0	[0.34]
α-Phellandrene	3.00	1127.3	0.32	4.22*	1002.0	[0.34]
Octanal	4.67	1251.1	0.01	4.25	1003.9	tr
Δ3-Carene	2.79	1111.6	0.01	4.31	1007.5	0.01
α-Terpinene	3.18	1141.4	0.15	4.42	1014.6	0.15
para-Cymene	4.34	1227.8	1.10	4.55	1022.8	0.98
Limonene	3.50	1165.5	79.93	4.69*	1031.6	[82.85]
β-Phellandrene	3.54	1169.2	2.50	4.69*	1031.6	[82.85]
(*E*)-β-Ocimene	4.22	1219.0	0.16	4.99	1050.0	0.15
γ-Terpinene	4.07	1208.3	8.07	5.12	1058.5	8.08
para-Cymenene	6.58	1387.6	0.04	5.56*	1085.8	[0.36]
Terpinolene	4.53	1241.3	0.32	5.56*	1085.8	[0.36]
Linalool	8.29	1514.3	1.72	5.82	1102.3	1.72
trans-para-Mentha-2,8-dien-1-ol	9.17	1582.0	0.02	6.08	1118.6	0.02
cis-Limonene oxide	6.68	1394.6	0.04	6.25	1129.2	0.04
trans-Limonene oxide	6.85	1406.9	0.03	6.32	1134.1	0.04
trans-Sabinol	10.06*	1653.2	[0.09]	6.35	1135.7	0.01
Epoxyterpinolene	6.98	1416.8	0.01	6.45	1142.0	0.01
Terpinen-4-ol	8.81	1554.3	0.11	6.94	1173.5	0.12
Cryptone	9.41	1600.7	0.01	7.02	1178.7	0.01
α-Terpineol	10.02	1649.5	0.10	7.17	1188.2	0.10
Decanal	7.56	1459.2	0.01	7.44	1205.5	0.02
trans-Carveol	11.63	1783.0	0.02	7.63	1218.1	0.02
cis-Carveol	11.96	1812.4	0.02	7.80	1229.8	0.02
Carvone	10.25	1668.0	0.03	7.94	1238.5	0.02
Thymol	15.40	2132.0	0.01	8.89	1302.3	0.06
δ-Elemene	7.23	1435.1	0.03	9.35	1334.4	0.03
α-Copaene	7.43	1449.7	0.02	9.86	1370.8	0.02
β-Cubebene	8.05	1495.9	0.03	10.07	1385.5	0.02
β-Elemene	8.70*	1545.4	[0.13]	10.11	1388.2	0.02
β-Caryophyllene	8.70*	1545.4	[0.13]	10.44	1411.8	0.11
γ-Elemene	9.31	1592.8	0.01	10.68	1429.5	0.01
α-Humulene	9.57	1613.6	0.03	10.90	1446.1	0.03
(*E*)-β-Farnesene	9.81	1632.9	0.24	11.04	1456.8	0.23
Germacrene D	10.06*	1653.2	[0.09]	11.28	1474.1	0.09

Bicyclogermacrene	10.34*	1675.8	[0.23]	11.48	1489.3	0.22
α-Muurolene	10.34*	1675.8	[0.23]	11.59	1497.0	0.02
δ-Cadinene	10.70	1705.3	0.05	11.86	1518.5	0.04
β-Sesquiphellandrene	10.90	1721.6	0.02	11.90	1521.4	0.05
Germacrene B	11.40	1764.2	0.07	12.24	1548.0	0.07
Spathulenol	14.65	2058.5	0.21	12.52	1569.7	0.20
Alismol	15.98	2189.6	0.02	13.18	1622.8	0.02
Isospathulenol	15.73*	2165.1	[0.01]	13.28	1630.7	0.02
τ-Muurolol	15.30	2122.6	0.01	13.34*	1636.0	[0.01]
τ-Cadinol				13.34*	1636.0	[0.01]
α-Cadinol	15.73*	2165.1	[0.01]	13.48	1647.8	0.02
Total reported		99.52%			99.89%	

*: 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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