

Date : March 06, 2018

CERTIFICATE OF ANALYSIS - GC PROFILING

SAMPLE IDENTIFICATION

Internal code : 18B20-PLG13-1-CC
Customer identification : Cinnamon Cassia
Type : Essential oil
Source : *Cinnamomum cassia*
Customer : Plant Guru

ANALYSIS

Method: PC-PA-014-17J19 - Analysis of the composition of an essential oil, or other volatile liquid, by FAST GC-FID (in French); identifications validated by GC-MS.

Analyst : Sarah-Eve Tremblay, M. Sc. A., Chimiste

Analysis date : February 28, 2018

Checked and approved by :

Alexis St-Gelais, M. Sc., chimiste 2013-174

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PHYSICOCHEMICAL DATA

Physical aspect: Bright yellow liquid

Refractive index: 1.6068 ± 0.0003 (20 °C)

CONCLUSION

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

ANALYSIS SUMMARY

Identification	DB-5 (%)	DB-WAX (%)	Classe
Hexanal	0.01	0.01	Aliphatic aldehyde
Furfural	0.01	0.01	Aliphatic alcohol
(2E)-Hexenal	tr	tr	Aliphatic aldehyde
Ethylbenzene	tr	tr*	Simple phenolic
Styrene	0.12	0.12	Simple phenolic
α -Thujene	0.01	0.01	Monoterpene
α -Pinene	0.06	0.06	Monoterpene
Camphene	0.04	0.04	Monoterpene
Benzaldehyde	0.81	0.84	Simple phenolic
β -Pinene	0.03*	0.03	Monoterpene
Sabinene	[0.03]*	[tr]*	Monoterpene
6-Methyl-5-hepten-2-one	0.01	0.01	Aliphatic ketone
Benzofuran	0.01		Simple phenolic
Octanal	tr	tr	Aliphatic aldehyde
Δ^3 -Carene	0.01	tr	Monoterpene
para-Cymene	0.03	0.03	Monoterpene
Limonene	0.05*	0.04	Monoterpene
1,8-Cineole	[0.05]*	0.01*	Monoterpenic ether
β -Phellandrene	[0.05]*	[0.01]*	Monoterpene
Benzyl alcohol	0.31*	0.32*	Simple phenolic
Salicylaldehyde	[0.31]*	0.29	Simple phenolic
(E)- β -Ocimene	tr	tr	Monoterpene
γ -Terpinene	0.01	tr	Monoterpene
Acetophenone	0.03	0.06*	Simple phenolic
Terpinolene	tr	tr	Monoterpene
ortho-Guaiacol	0.03	tr	Simple phenolic
Linalool	0.01	0.01	Monoterpenic alcohol
Nonanal	0.01	0.01	Aliphatic aldehyde
Phenylethyl alcohol	0.63	0.69*	Simple phenolic
ortho-Vinylanisole	0.02	0.04	Simple phenolic
trans-Pinocarveol	0.01	0.01	Monoterpenic alcohol
2-Methylbenzofuran	0.03	0.02	Phenylpropanoid
Hydrocinnamal	0.48*	0.42*	Phenylpropanoid
Unknown	[0.48]*	tr	Phenylpropanoid
Borneol	0.10	0.12*	Monoterpenic alcohol
3-Methylbenzofuran?	0.08	0.08	Phenylpropanoid
Terpinen-4-ol	0.01	0.01	Monoterpenic alcohol
para-Cymen-8-ol	0.04	0.03	Monoterpenic alcohol
α -Terpineol	0.01	[0.12]*	Monoterpenic alcohol
Methyl salicylate	0.01	[0.42]*	Phenolic ester
(Z)-Cinnamal	0.33	[0.32]*	Phenylpropanoid
Hydrocinnamyl alcohol	0.08	0.15	Phenylpropanoid
ortho-Anisaldehyde	0.46	0.47	Simple phenolic
Phenylethyl acetate	0.07	0.09	Phenolic ester
(E)-Cinnamal	76.59	77.04	Phenylpropanoid
(E)-Cinnamyl alcohol	0.19	0.19	Phenylpropanoid
Eugenol	0.18*	0.14*	Phenylpropanoid
Cyclosativene I	[0.18]*	0.03	Sesquiterpene

α-Ylangene	0.02	0.02	Sesquiterpene
α-Copaene	0.52*	0.32	Sesquiterpene
ortho-Methoxyhydrocinnamal?	[0.52]*	0.07	Phenylpropanoid
β-Elemene	0.02	0.15*	Sesquiterpene
β-Caryophyllene	0.12	[0.15]*	Sesquiterpene
Coumarin	2.25	2.21	Coumarin
trans-α-Bergamotene	tr	tr	Sesquiterpene
(E)-Cinnamyl acetate	2.93	2.91	Phenylpropanoid ester
allo-Aromadendrene	0.19*	[0.06]*	Sesquiterpene
(Z)-ortho-Methoxycinnamal	[0.19]*	0.06	Phenylpropanoid
(E)-Cinnamic acid	0.40	0.02	Phenylpropanoid
γ-Muurolene	0.10	0.12	Sesquiterpene
ar-Curcumene	0.06	0.07	Sesquiterpene
Viridiflorene	0.05	0.04	Sesquiterpene
α-Muurolene	0.06	0.07	Sesquiterpene
β-Bisabolene	0.09	0.09	Sesquiterpene
γ-Cadinene	0.08	0.17*	Sesquiterpene
δ-Cadinene	0.14*	[0.17]*	Sesquiterpene
trans-Calamenene	[0.14]*	0.03	Sesquiterpene
α-Calacorene	9.58*	[0.69]*	Sesquiterpene
(E)-ortho-Methoxycinnamal	[9.58]*	9.46	Phenylpropanoid
(E)-Nerolidol	0.18	0.19	Sesquiterpenic alcohol
Spathulenol	0.11	0.12	Sesquiterpenic alcohol
Caryophyllene oxide	0.07	0.08	Sesquiterpenic ether
Humulene epoxide II	0.01		Sesquiterpenic ether
Tetradecanal?	0.03	0.01	Aliphatic aldehyde
1-epi-Cubenol	0.03	0.01	Sesquiterpenic alcohol
Caryophylladienol II	0.02	0.02	Sesquiterpenic alcohol
τ-Cadinol	0.05	0.03	Sesquiterpenic alcohol
α-Cadinol	0.03	0.02	Sesquiterpenic alcohol
(3Z)-Caryophylla-3,8(13)-dien-5β-ol	0.03	0.02	Sesquiterpenic alcohol
Mustakone?	0.14*		Sesquiterpenic ketone
Cadalene	[0.14]*	0.05*	Sesquiterpene
α-Bisabolol	0.05	[0.05]*	Sesquiterpenic alcohol
Benzyl benzoate	0.06	0.05	Phenolic ester
Phenylethyl benzoate	0.05	0.03	Phenolic ester
Benzyl salicylate	0.01	0.01	Phenolic ester
Rimuene	tr	[0.14]*	Diterpene
Dolabradiene	0.06	0.06	Diterpene
Manoyl oxide	0.05	0.04	Diterpenic ether
Kaurene?	0.01		Diterpene
Phenylethyl (E)-cinnamate	0.04	0.01	Phenylpropanoid ester
Total identified	98.54%	98.02%	

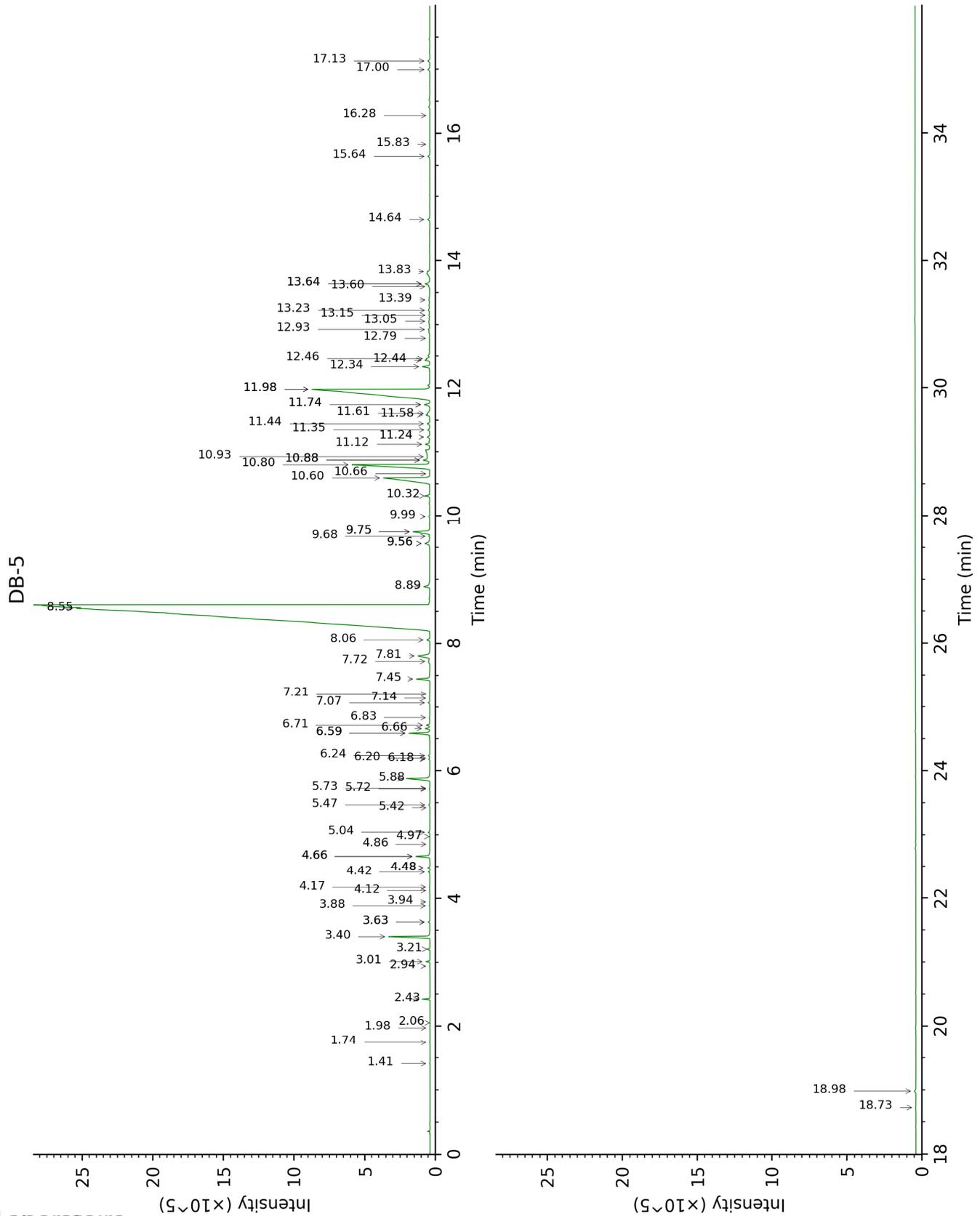
*: Two or more compounds are coeluting on this column

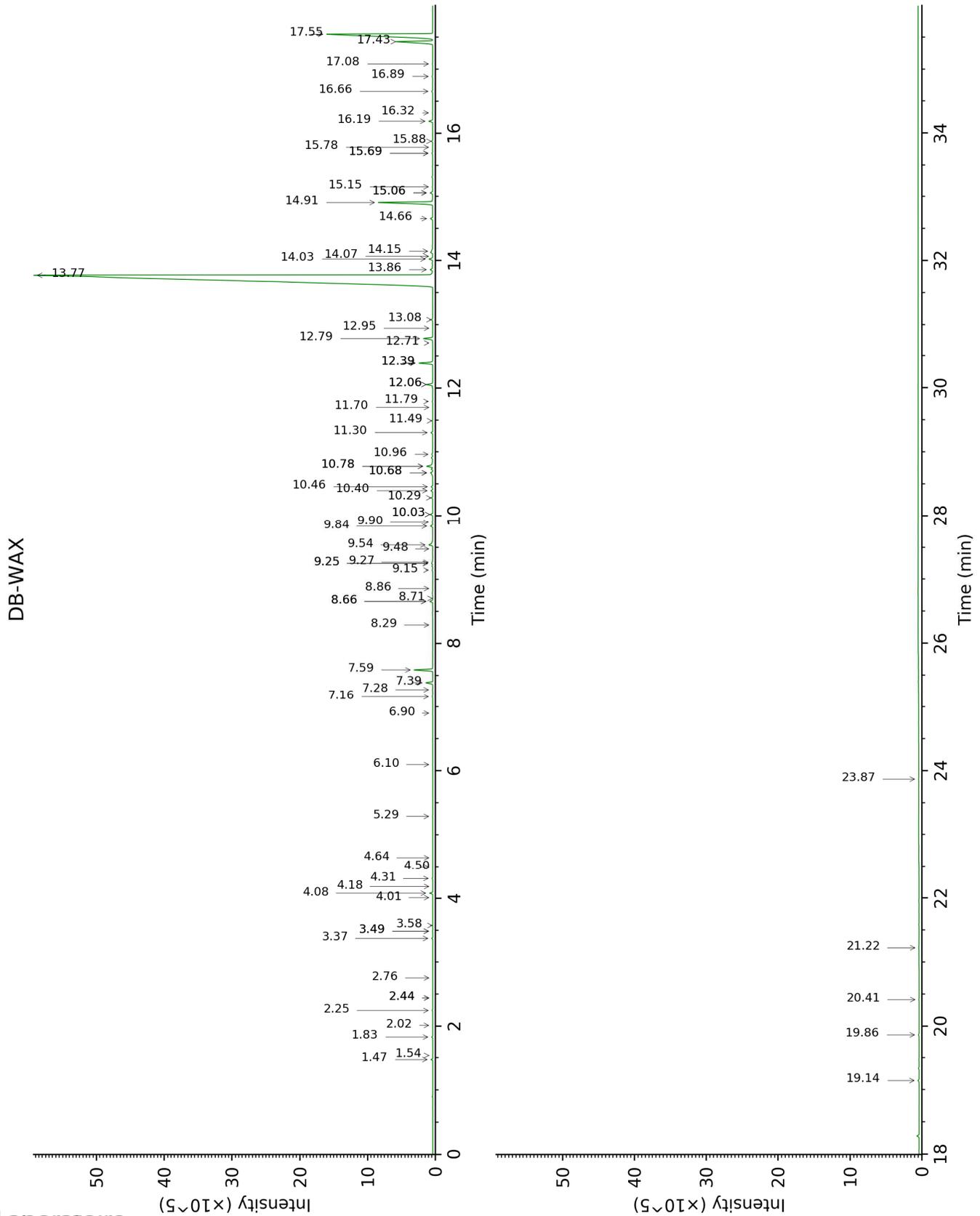
[xx]: Duplicate percentage due to coelutions, not taken account in the identified total

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

Note: no correction factor was applied

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FULL ANALYSIS DATA

Identification	Column DB-5			Column DB-WAX		
	R.T	R.I	%	R.T	R.I	%
Hexanal	1.41	795	0.01	2.02	1047	0.01
Furfural	1.74	825	0.01	6.90	1413	0.01
(2E)-Hexenal	1.98	844	tr	3.58	1174	tr
Ethylbenzene	2.06	851	tr	2.44*	1087	tr
Styrene	2.43	882	0.12	4.08	1211	0.12
α-Thujene	2.94	920	0.01	1.54	1002	0.01
α-Pinene	3.01	925	0.06	1.47	995	0.06
Camphene	3.21	938	0.04	1.83	1030	0.04
Benzaldehyde	3.40	951	0.81	7.59	1463	0.84
β-Pinene	3.63*	966	0.03	2.25	1069	0.03
Sabinene	3.63*	966	[0.03]	2.44*	1087	[tr]
6-Methyl-5-hepten-2-one	3.88	983	0.01	5.29	1297	0.01
Benzofuran	3.94	987	0.01			
Octanal	4.12	999	tr	4.64	1251	tr
Δ3-Carene	4.18	1002	0.01	2.76	1112	tr
para-Cymene	4.42	1017	0.03	4.31	1228	0.03
Limonene	4.48*	1022	0.05	3.37	1159	0.04
1,8-Cineole	4.48*	1022	[0.05]	3.49*	1167	0.01
β-Phellandrene	4.48*	1022	[0.05]	3.49*	1167	[0.01]
Benzyl alcohol	4.66*	1033	0.31	12.06*	1823	0.32
Salicylaldehyde	4.66*	1033	[0.31]	9.54	1613	0.29
(E)-β-Ocimene	4.86	1045	tr	4.18	1219	tr
γ-Terpinene	4.98	1053	0.01	4.01	1206	tr
Acetophenone	5.04	1057	0.03	9.25*	1590	0.06
Terpinolene	5.42	1081	tr	4.50	1241	tr
ortho-Guaiacol	5.47	1084	0.03	11.70	1791	tr
Linalool	5.72	1099	0.01	8.29	1516	0.01
Nonanal	5.73	1100	0.01	6.10	1355	0.01
Phenylethyl alcohol	5.88	1110	0.63	12.39*	1852	0.69
ortho-Vinylanisole	6.18	1129	0.02	9.15	1582	0.04
trans-Pinocarveol	6.20	1130	0.01	9.48	1608	0.01
2-Methylbenzofuran	6.24	1132	0.03	9.27	1591	0.02
Hydrocinnamal	6.59*	1155	0.48	10.78*	1714	0.42
Unknown [m/z 133, 77 (86), 105 (75), 79 (68), 134 (48)]	6.59*	1155	[0.48]	12.95	1902	tr
Borneol	6.66	1159	0.10	10.02*	1652	0.12
3-Methylbenzofuran?	6.71	1163	0.08	10.46	1687	0.08
Terpinen-4-ol	6.83	1170	0.01	8.86	1560	0.01
para-Cymen-8-ol	7.07	1185	0.04	11.79	1799	0.03
α-Terpineol	7.14	1190	0.01	10.02*	1652	[0.12]
Methyl salicylate	7.21	1194	0.01	10.78*	1714	[0.42]
(Z)-Cinnamal	7.45	1210	0.33	12.06*	1823	[0.32]
Hydrocinnamyl alcohol	7.72	1228	0.08	13.86	1985	0.15
ortho-Anisaldehyde	7.81	1234	0.46	12.78	1887	0.47
Phenylethyl acetate	8.06	1251	0.07	11.30	1758	0.09
(E)-Cinnamal	8.55	1284	76.59	13.77	1977	77.04

(E)-Cinnamyl alcohol	8.89	1306	0.19	16.19	2214	0.19
Eugenol	9.56*	1354	0.18	15.06*	2101	0.14
Cyclosativene I	9.56*	1354	[0.18]	7.16	1432	0.03
α-Ylangene	9.68	1362	0.02	7.28	1440	0.02
α-Copaene	9.75*	1367	0.52	7.39	1448	0.32
ortho-Methoxyhydrocinnamal?	9.75*	1367	[0.52]	14.15	2013	0.07
β-Elemene	9.99	1384	0.02	8.66*	1544	0.15
β-Caryophyllene	10.32	1407	0.12	8.66*	1544	[0.15]
Coumarin	10.60	1428	2.25	17.43	2345	2.21
trans-α-Bergamotene	10.66	1433	tr	8.70	1548	tr
(E)-Cinnamyl acetate	10.80	1444	2.93	14.91	2086	2.91
allo-Aromadendrene	10.88*	1449	0.19	9.25*	1590	[0.06]
(Z)-ortho-Methoxycinnamal	10.88*	1449	[0.19]	15.88	2182	0.06
(E)-Cinnamic acid	10.93	1453	0.40	21.22	2787	0.02
γ-Muurolene	11.12	1467	0.10	9.84	1637	0.12
ar-Curcumene	11.24	1476	0.06	10.96	1729	0.07
Viridiflorene	11.35	1484	0.05	9.90	1642	0.04
α-Muurolene	11.44	1491	0.06	10.29	1673	0.07
β-Bisabolene	11.58	1501	0.09	10.40	1682	0.09
γ-Cadinene	11.61	1503	0.08	10.68*†	1705	0.17
δ-Cadinene	11.74*	1514	0.14	10.68*†	1705	[0.17]
trans-Calamenene	11.74*	1514	[0.14]	11.49	1774	0.03
α-Calacorene	11.98*	1533	9.58	12.39*	1852	[0.69]
(E)-ortho-Methoxycinnamal	11.98*	1533	[9.58]	17.55	2358	9.46
(E)-Nerolidol	12.34	1560	0.18	14.03	2001	0.19
Spathulenol	12.44	1568	0.11	14.66	2062	0.12
Caryophyllene oxide	12.46	1570	0.07	13.08	1914	0.08
Humulene epoxide II	12.79	1596	0.01			
Tetradecanal?	12.93	1607	0.03	12.72	1881	0.01
1-epi-Cubenol	13.05	1618	0.03	14.07	2005	0.01
Caryophylladienol II	13.15	1625	0.02	16.32	2228	0.02
τ-Cadinol	13.23	1632	0.05	15.15	2110	0.03
α-Cadinol	13.39	1645	0.03	15.78	2173	0.02
(3Z)-Caryophylla-3,8(13)-dien-5β-ol	13.60	1662	0.03	17.08	2308	0.02
Mustakone?	13.64*	1666	0.14			
Cadalene	13.64*	1666	[0.14]	15.69*	2163	0.05
α-Bisabolol	13.83	1681	0.05	15.69*	2163	[0.05]
Benzyl benzoate	14.64	1750	0.06	19.14	2537	0.05
Phenylethyl benzoate	15.64	1838	0.05	19.86	2621	0.03
Benzyl salicylate	15.83	1856	0.01	20.42	2687	0.01
Rimuene	16.28	1896	tr	15.06*	2101	[0.14]
Dolabradiene	17.00	1964	0.06	16.66	2263	0.06
Manoyl oxide	17.13	1977	0.05	16.89	2287	0.04
Kaurene?	18.73	2135	0.01			
Phenylethyl (E)-cinnamate	18.98	2161	0.04	23.87	3138	0.01
Total identified		98.54%			98.02%	

Total reported	98.54%	98.02%
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*: Two or more compounds are coeluting on this column

[xx]: Duplicate percentage due to coelutions, not taken account in the identified 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