

Date : 2023-07-05

CERTIFICATE OF ANALYSIS - GC PROFILING

SAMPLE IDENTIFICATION

Internal code : 23F27-PTH02

Customer Identification : Organic Clove Bud - Sri Lanka - CH0113R

Type : Essential Oil

Source : *Syzygium aromaticum*

Customer : Plant Therapy

Checked and approved by:

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 liquid by FAST GC-FID



Results : See analysis summary (next page)

Analyst : Sylvain Mercier, M. Sc., Chimiste 2014-005

Date : 2023-07-04

PHYSICOCHEMICAL DATA

Physical aspect : Yellow liquid

Analyst : Cindy Caron B. Sc.

Date : 2023-06-28

Refractive index : 1.535 ± 0.0003 (20 °C)

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

Analyst : Cindy Caron B. Sc.

Date : 2023-06-28

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
Furfural	0.02	Furan
Linalool	0.01	Monoterpnic alcohol
Methyl salicylate	0.02	Phenolic ester
Chavicol	0.14	Phenylpropanoid
Chavicyl acetate	0.02	Phenylpropanoid ester
Eugenol	80.32	Phenylpropanoid
Dihydroeugenol	0.13	Phenylpropanoid
α -Copaene	0.13	Sesquiterpene
Vanillin	0.01	Simple phenolic
Isocaryophyllene	0.03	Sesquiterpene
Methyleugenol	0.04	Phenylpropanoid
β -Caryophyllene	4.95	Sesquiterpene
Caryophylla-4(12),8(13)-diene	0.04	Sesquiterpene
(E)-Isoeugenol	0.05	Phenylpropanoid
α -Humulene	0.64	Sesquiterpene
allo-Aromadendrene	0.01	Sesquiterpene
trans-Cadina-1(6),4-diene	0.05	Sesquiterpene
γ -Muurolene	0.03	Sesquiterpene
β -Selinene	0.01	Sesquiterpene
α -Selinene	0.02	Sesquiterpene
α -Muurolene	0.01	Sesquiterpene
(3Z,6E)- α -Farnesene	0.02	Sesquiterpene
γ -Cadinene	0.03	Sesquiterpene
Cubebol	0.02	Sesquiterpenic alcohol
trans-Calamenene	0.05	Sesquiterpene
δ -Cadinene	0.21	Sesquiterpene
Eugenyl acetate	11.66	Phenylpropanoid ester
Unknown	0.08	Unknown
Unknown	0.02	Phenylpropanoid
Caryophyllenyl alcohol	0.10	Sesquiterpenic alcohol
Unknown	0.01	Oxygenated sesquiterpene
Caryophyllene oxide	0.13	Sesquiterpenic ether
Unknown	0.01	Unknown
Humulene epoxide I	0.02	Sesquiterpenic ether
Unknown	0.05	Unknown
Humulene epoxide II	0.03	Sesquiterpenic ether
(E)-Isoeugenyl acetate	0.07	Phenylpropanoid ester
1-epi-Cubenol	0.05	Sesquiterpenic alcohol
Caryophylladienol I	0.03	Sesquiterpenic alcohol
Caryophylladienol II	0.04	Sesquiterpenic alcohol

τ -Cadinol	0.02	Sesquiterpenic alcohol
Cubenol	0.02	Sesquiterpenic alcohol
τ -Muurolol	0.01	Sesquiterpenic alcohol
α -Muurolol	0.01	Sesquiterpenic alcohol
α -Cadinol	0.01	Sesquiterpenic alcohol
14-Hydroxy-(Z)-caryophyllene	0.07	Sesquiterpenic alcohol
14-Hydroxy-9-epi-(E)-caryophyllene	0.02	Sesquiterpenic alcohol
(3Z)-Caryophylla-3,8(13)-dien-5 β -ol	0.04	Sesquiterpenic alcohol
Trimethoxypropylbenzene analog	0.03	Phenylpropanoid
Unknown	0.11	Lignan
Consolidated total	99.67	

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

Note: no correction factor was applied

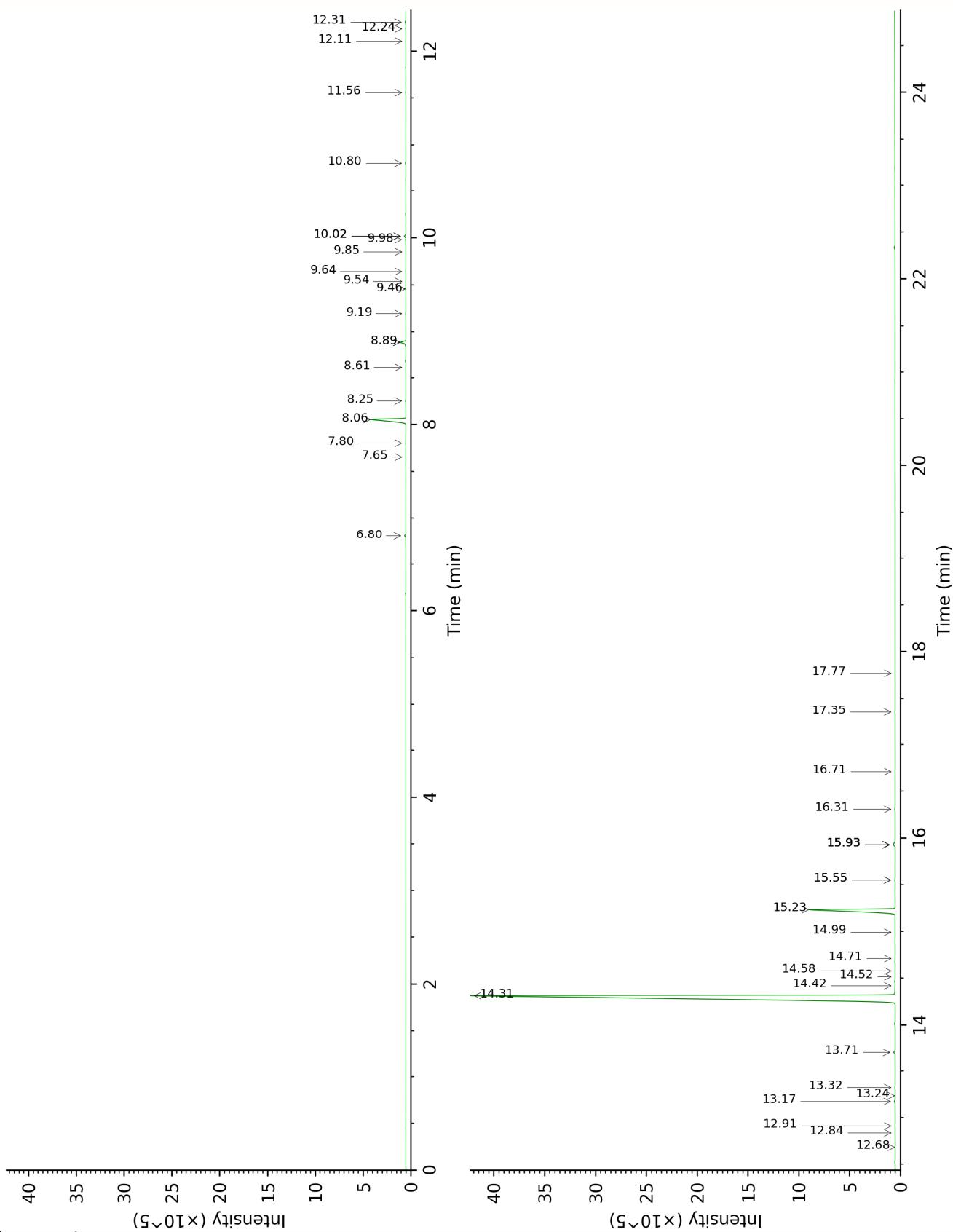
About "consolidated" data: The table above presents the breakdown of the sample volatile constituents after applying an algorithm to collapse data acquired from the multi-columns system of PhytoChemia into a single set of consolidated contents. In case of discrepancies between columns, the algorithm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

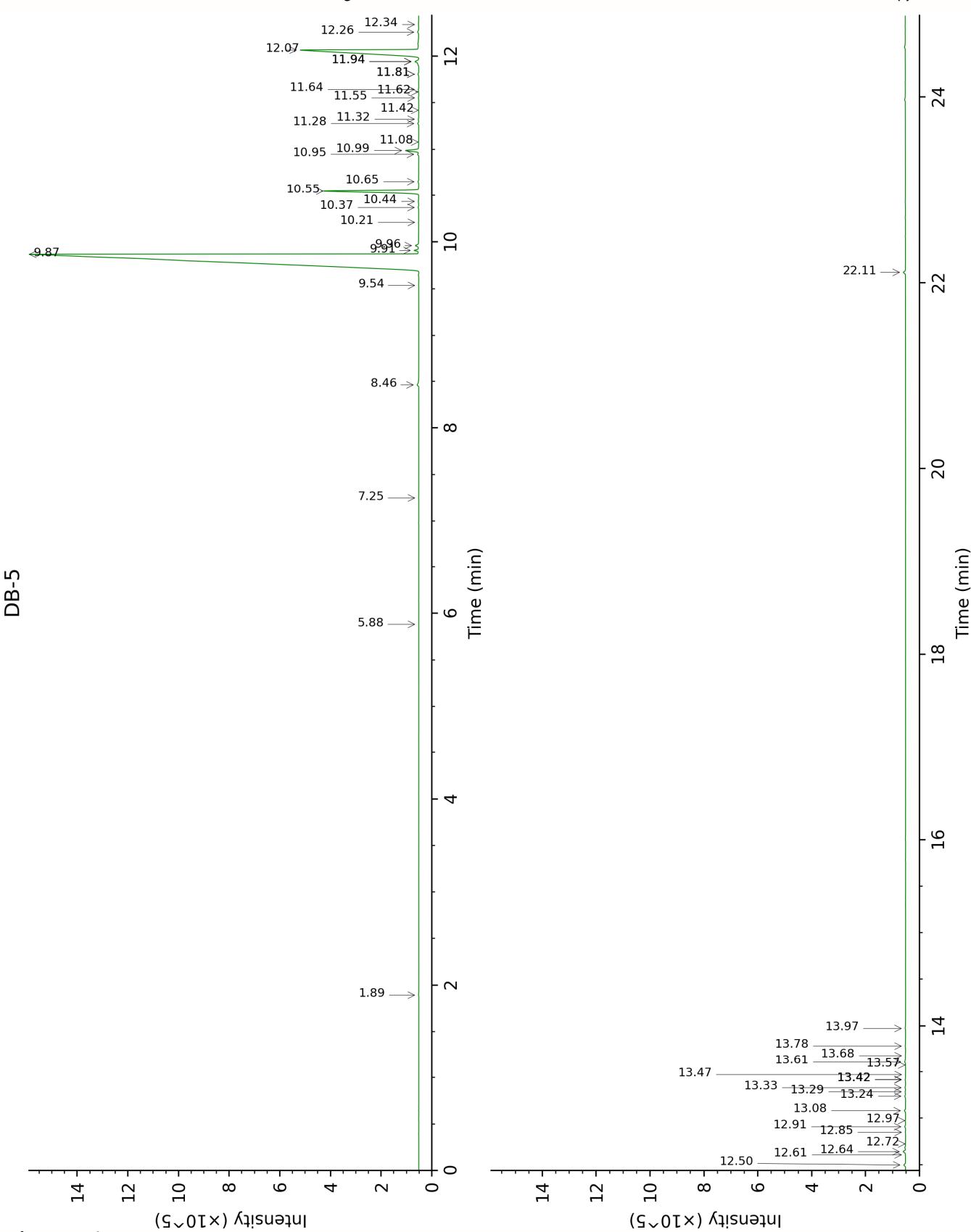
Unknowns: Unknown compounds' mass spectral data is presented in the "Full analysis data" table. The occurrence of unknown compounds is to be expected in many samples, and does not denote particular problems unless noted otherwise in the conclusion.

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





FULL ANALYSIS DATA

Furfural	Column DB-WAX			Column DB-5		
				1.89	830.6	0.02
Linalool	7.65	1514.0	0.01	5.88	1102.3	0.01
Methyl salicylate	10.02*	1704.4	[0.17]	7.25	1189.2	0.01
Chavicol	15.93*	2265.8	[0.29]	8.46	1269.5	0.14
Chavicyl acetate	12.24	1899.9	0.03	9.54	1343.1	0.02
Eugenol	14.31	2098.5	80.19	9.87	1366.6	80.33
Dihydroeugenol	13.71	2038.9	0.15	9.91	1369.5	0.13
α -Copaene	6.80	1449.8	0.14	9.96	1373.1	0.13
Vanillin	17.77	2470.6	0.01	10.21	1390.6	0.01
Isocaryophyllene	7.80	1525.7	0.03	10.37	1401.9	0.03
Methyleugenol	12.84	1955.7	0.01	10.44	1406.5	0.04
β -Caryophyllene	8.06	1545.6	4.91	10.55	1414.8	4.95
Caryophylla-4(12),8(13)-diene	8.25	1561.2	0.05	10.65	1422.7	0.04
(E)-Isoeugenol	15.93*	2265.8	[0.29]	10.95	1444.6	0.05
α -Humulene	8.89*	1611.2	[0.67]	10.99	1447.7	0.64
allo-Aromadendrene	8.61	1589.4	0.01	11.08	1454.5	0.01
trans-Cadina-1(6),4-diene	8.89*	1611.2	[0.67]	11.28	1469.3	0.05
γ -Muurolene	9.19	1636.4	0.03	11.32	1472.7	0.03
β -Selinene	9.46	1658.0	0.01	11.42	1479.9	0.01
α -Selinene	9.54	1664.7	0.02	11.55	1489.8	0.02
α -Muurolene	9.64	1673.4	0.02	11.62	1494.6	0.01
(3Z,6E)- α -Farnesene	9.85	1690.6	0.01	11.64	1496.5	0.02
γ -Cadinene	9.98	1701.3	0.03	11.81*	1508.9	[0.04]
Cubebol	12.11	1888.0	0.02	11.81*	1508.9	[0.04]
trans-Calamenene	10.80	1771.9	0.05	11.94*	1519.5	[0.26]
δ -Cadinene	10.02*	1704.4	[0.17]	11.94*	1519.5	[0.26]
Eugenyl acetate	15.23	2192.4	11.55	12.07	1529.2	11.65
Unknown SYAR II [m/z 164, 135 (98), 93 (86), 107 (83), 79 (69)...]	11.56	1838.3	0.03	12.26	1544.3	0.08
Unknown SYAR III [m/z 180, 93 (70), 55 (62), 77 (55), 164 (55), 103 (50)]				12.34	1550.7	0.02
Caryophyllenyl alcohol	13.17	1987.4	0.09	12.50	1563.0	0.10
Unknown SYAR IV [m/z 161, 187 (32), 105 (30), 205 (24)... 222 (3)]	14.52	2118.6	0.01	12.61	1571.6	0.01
Caryophyllene oxide	12.31	1906.3	0.09	12.64	1574.3	0.13
Unknown SYAR V [m/z 151, 178 (54), 123 (20), 55 (13), 161 (11), 77 (10)...]				12.72	1580.8	0.01
Humulene epoxide I	12.68	1941.1	0.02	12.85	1590.6	0.02

Unknown SYAR XII [m/z 164, 93 (48), 43 (44), 91 (27), 55 (27)...]			12.91	1595.3	0.05
Humulene epoxide II	12.91	1962.6	0.03	12.98	1600.5
(E)-Isoeugenyl acetate	16.71	2350.8	0.02	13.08	1609.0
1-epi-Cubenol	13.32	2001.6	0.04	13.24	1621.9
Caryophylladienol I	15.55*	2225.4	[0.05]	13.28	1625.8
Caryophylladienol II	15.55*	2225.4	[0.05]	13.33	1629.4
τ-Cadinol	14.42	2108.9	0.02	13.42*	1636.6
Cubenol	13.24	1993.2	0.02	13.42*	1636.6
τ-Muurolol	14.58	2125.1	0.01	13.42*	1636.6
α-Muurolol	14.71	2138.7	0.01	13.47	1641.1
α-Cadinol	14.99	2167.6	0.02	13.57	1649.6
14-Hydroxy-(Z)-caryophyllene	15.93*	2265.8	[0.29]	13.61	1652.4
14-Hydroxy-9-epi-(E)-caryophyllene	15.93*	2265.8	[0.29]	13.68	1658.3
(3Z)-Caryophylla-3,8(13)-dien-5β-ol	16.31	2306.5	0.05	13.78	1666.9
Trimethoxypropylbenzene analog	17.35	2422.0	0.03	13.97	1682.6
Unknown OCSA V [m/z 326, 148 (67), 147 (41), 117 (30), 91 (22)...]			22.11	2494.4	0.11
Total reported		98.97%		99.63%	

*: 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

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