

Date : 2023-10-19

CERTIFICATE OF ANALYSIS - GC PROFILING

SAMPLE IDENTIFICATION

Internal code : 23J12-PTH04

Customer Identification : Organic Wintergreen - Nepal - W20107R

Type : Essential Oil

Source : *Gaultheria fragrantissima*

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 liquide by FAST GC-FID

✖ISO

Results : See analysis summary (next page)

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

Date : 2023-10-16

PHYSICOCHEMICAL DATA

Refractive index : 1.5361 ± 0.0003 (20 °C)

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

Analyst : Cindy Caron B. Sc.

Date : 2023-10-12

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
Isovaleral	0.01	Aliphatic aldehyde
2-Methylbutyral	tr	Aliphatic aldehyde
(3Z)-Hexenol	0.01	Aliphatic alcohol
α -Pinene	0.01	Monoterpene
Camphene	0.01	Monoterpene
β -Pinene	0.02	Monoterpene
α -Phellandrene	tr	Monoterpene
<i>para</i> -Cymene	0.01	Monoterpene
1,8-Cineole	tr	Monoterpenic ether
Limonene	0.06	Monoterpene
Linalool	0.02	Monoterpenic alcohol
Methyl salicylate	99.33	Phenolic ester
Ethyl salicylate	0.06	Phenolic ester
Eugenol	0.08	Phenylpropanoid
Consolidated total	99.63	

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

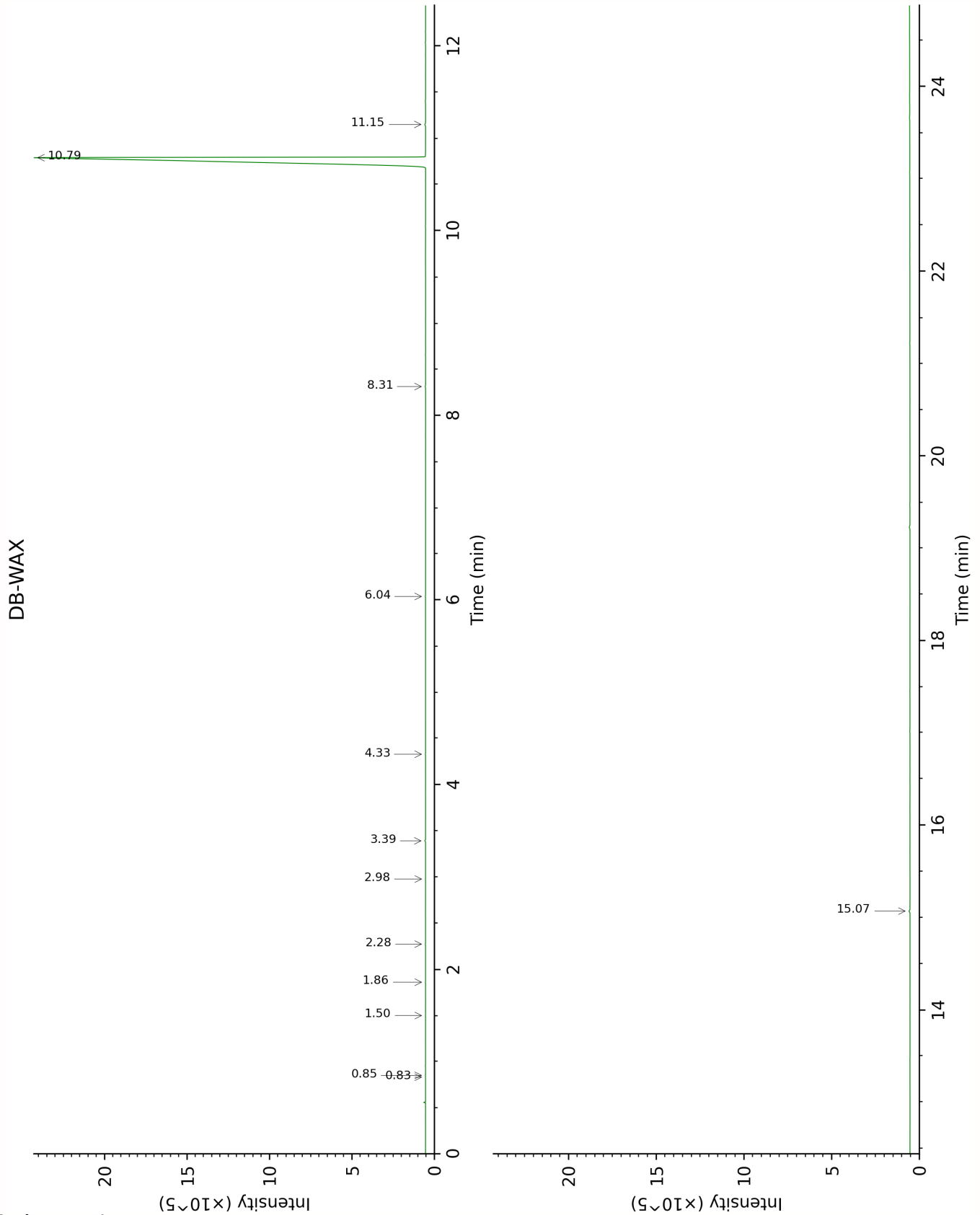
Note: no correction factor was applied

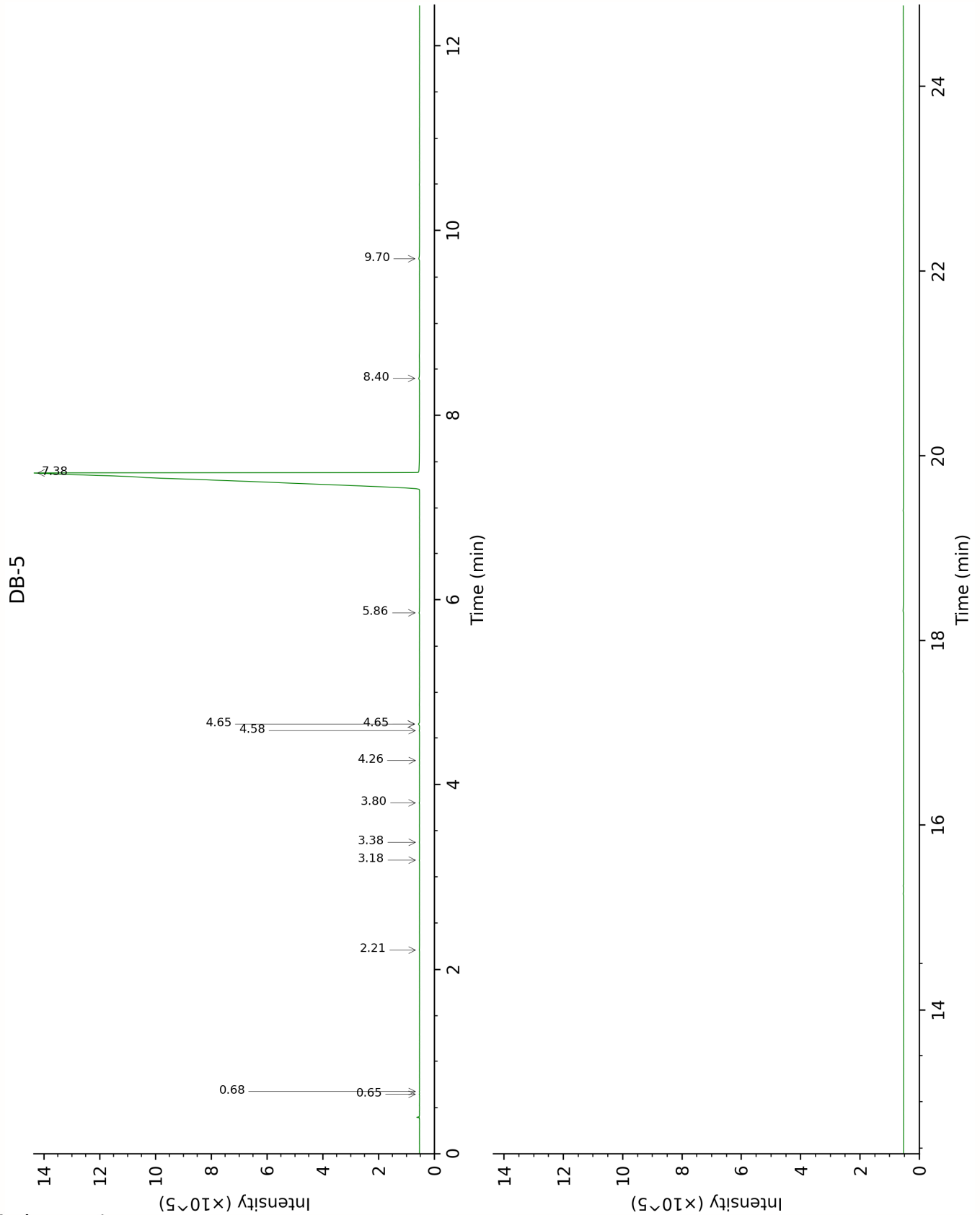
About "consolidated" data: The table above presents the breakdown of the sample volatile constituents after applying an algorithm to collapse data acquired from the multi-columns system of PhytoChemia into a single set of consolidated contents. In case of discrepancies between columns, the algorithm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

Isovaleral	Column DB-WAX			Column DB-5		
	0.85	889.7	0.01	0.64	643.0	0.01
2-Methylbutyral	0.83	883.0	0.01	0.68	653.1	tr
(3Z)-Hexenol	6.04	1350.9	0.02	2.21	859.0	0.01
α-Pinene	1.50	993.5	0.01	3.18	931.9	0.01
Camphene	1.86	1029.2	0.01	3.38	944.6	0.01
β-Pinene	2.28	1068.0	0.02	3.80	972.5	0.02
α-Phellandrene	2.98	1126.9	0.01	4.26	1002.6	tr
para-Cymene	4.33	1226.7	0.02	4.58	1022.8	0.01
1,8-Cineole				4.65*	1027.3	[0.06]
Limonene	3.39	1158.1	0.06	4.65*	1027.3	[0.06]
Linalool	8.31	1517.0	0.04	5.86	1102.7	0.02
Methyl salicylate	10.79	1713.1	98.79	7.38	1199.3	99.33
Ethyl salicylate	11.15	1743.3	0.08	8.40	1267.2	0.06
Eugenol	15.07	2098.5	0.10	9.70	1356.3	0.08
Total reported		99.17%			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

†: 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