Date : 2025-09-26

Certificate of Analysis - GC Profiling

Sample Identification

Internal code : 25H20-PTH18
Customer Identification : Wintergreen - China - W10112R
Type : Essential Oil
Source : *Gaultheria procumbens*
Customer : Plant Therapy

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Notes: This report may not be published, including online, without the written consent from Laboratoire PhytoChemia.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

This report is an update from the first version issued on 2025-08-27 to correct the customer identification.

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 : Sylvain Mercier, M. Sc., Chimiste 2014-005
Date : 2025-08-22

Physicochemical data

Refractive index : 1.5359 ± 0.0003 (20 °C)
Method : PC-MAT-016 - Measure of the refractive index of a liquid.
Analyst : Kassandra Baker
Date : 2025-08-21

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
(3*Z*)-Hexenol	0.02	Aliphatic alcohol
Hexanol	0.01	Aliphatic alcohol
α-Pinene	0.01	Monoterpene
Benzaldehyde	0.01	Simple phenolic
β-Pinene	0.01	Monoterpene
Myrcene	0.01	Monoterpene
para-Cymene	0.02	Monoterpene
Limonene	0.03	Monoterpene
1,8-Cineole	0.04	Monoterpenic ether
Benzyl alcohol	0.02	Simple phenolic
γ-Terpinene	0.02	Monoterpene
Octanol	0.01	Aliphatic alcohol
Terpinolene	0.01	Monoterpene
Linalool	0.05	Monoterpenic alcohol
Nonanal	0.01	Aliphatic aldehyde
Menthol	0.01	Monoterpenic alcohol
2,4-Nonanedione?	0.10	Aliphatic ketone
Methyl salicylate	98.35	Phenolic ester
Nerol	0.01	Monoterpenic alcohol
Geraniol	0.04	Monoterpenic alcohol
Ethyl salicylate	0.07	Phenolic ester
Vitispirane	0.03	Terpenic ether
Dehydro elsholtzia ketone	0.02	Monoterpenic ketone
Eugenol	0.03	Phenylpropanoid
α-Copaene	0.01	Sesquiterpene
α-Gurjunene	0.02	Sesquiterpene
β-Caryophyllene	0.03	Sesquiterpene
β-Gurjunene	0.01	Sesquiterpene
α-Maaliene	0.02	Sesquiterpene
Aromadendrene	0.24	Sesquiterpene
α-Humulene	0.02	Sesquiterpene
allo-Aromadendrene	0.06	Sesquiterpene
β-Selinene	0.01	Sesquiterpene
allo-Aromadendr-9-ene	0.01	Sesquiterpene
Viridiflorene	0.04	Sesquiterpene
γ-Cadinene	0.01	Sesquiterpene
δ-Cadinene	0.01	Sesquiterpene
Consolidated total	**99.45**	

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

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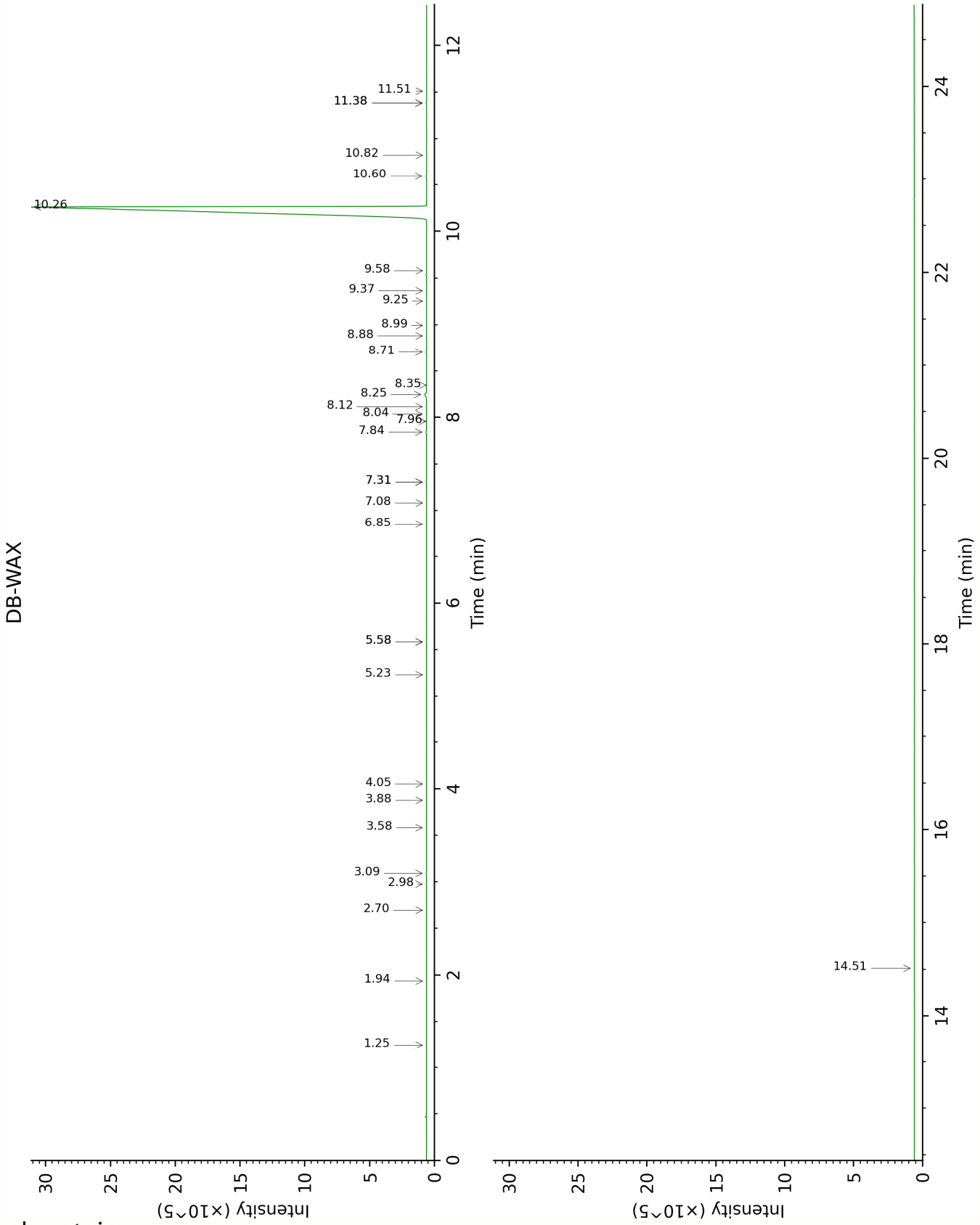
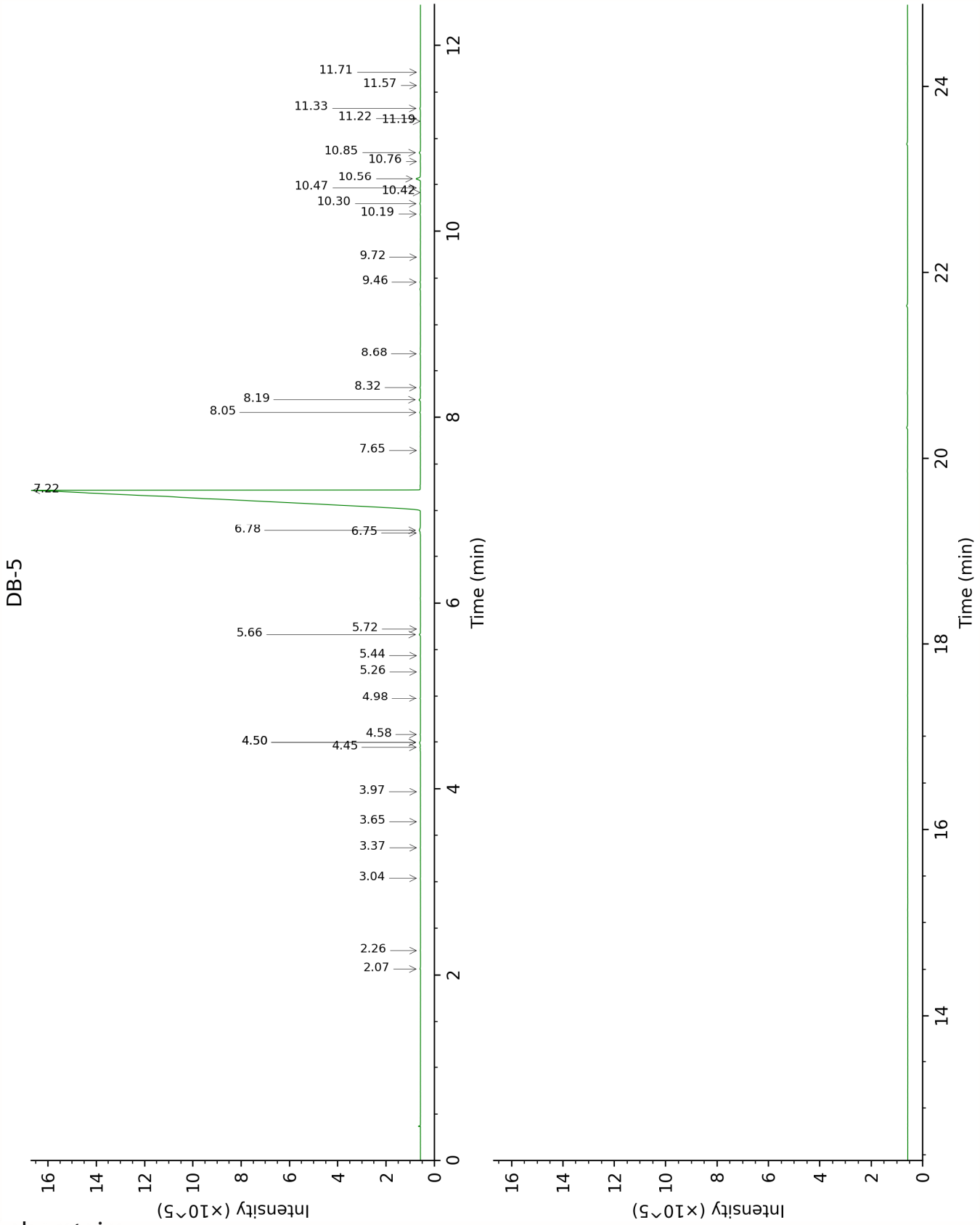

DB-WAX
Time (min)
Intensity (×10^5)
11.51
11.38
10.82
10.60
10.26
9.58
9.37
9.25
8.99
8.88
8.71
8.35
8.25
8.12
8.04
7.96
7.84
7.31
7.08
6.85
5.58
5.23
4.05
3.88
3.58
3.09
2.98
2.70
1.94
1.25
14.51


DB-5
11.71
11.57
11.33
11.22
11.19
10.85
10.76
10.56
10.47
10.42
10.30
10.19
9.72
9.46
8.68
8.32
8.19
8.05
7.65
7.22
6.78
6.75
5.72
5.66
5.44
5.26
4.98
4.58
4.50
4.45
3.97
3.65
3.37
3.04
2.26
2.07
Time (min)
Intensity (×10^5)

FULL ANALYSIS DATA

	Column DB-WAX			Column DB-5		
(3*Z*)-Hexenol	5.58*	1351.0	[0.03]	2.07	854.7	0.02
Hexanol	5.23	1325.4	0.01	2.26	871.5	0.01
α-Pinene	1.25	989.0	0.01	3.04	930.3	0.01
Benzaldehyde	7.08	1461.2	0.01	3.37	952.3	0.01
β-Pinene	1.94	1065.0	0.01	3.65	971.0	0.01
Myrcene	2.70	1133.9	0.01	3.97	992.7	0.01
para-Cymene	3.88	1225.9	0.02	4.45	1023.6	0.02
Limonene	2.98	1156.4	0.03	4.50*	1026.9	[0.06]
1,8-Cineole	3.09	1165.6	0.04	4.50*	1026.9	[0.06]
Benzyl alcohol	11.51	1820.3	0.02	4.58	1032.3	0.02
γ-Terpinene	3.58	1204.3	0.02	4.98	1057.2	0.02
Octanol	7.96	1527.9	0.01	5.26	1075.4	0.01
Terpinolene	4.05	1238.7	0.01	5.44	1086.4	0.01
Linalool	7.84	1518.9	0.05	5.66	1100.8	0.05
Nonanal	5.58*	1351.0	[0.03]	5.72	1104.6	0.01
Menthol	8.88	1599.9	0.02	6.75	1171.2	0.01
2,4-Nonanedione?				6.78	1173.1	0.10
Methyl salicylate	10.26	1712.9	98.62	7.22	1201.4	98.35
Nerol	10.82	1760.6	0.01	7.65	1230.1	0.01
Geraniol	11.38*	1809.0	[0.06]	8.05	1257.8	0.04
Ethyl salicylate	10.60	1741.3	0.07	8.19	1267.1	0.07
Vitispirane	7.31*	1478.0	[0.05]	8.32	1275.9	0.03
Dehydro elsholtzia ketone	11.38*	1809.0	[0.06]	8.68	1300.5	0.02
Eugenol	14.51	2100.3	0.02	9.46	1352.1	0.03
α-Copaene	6.85	1444.1	0.01	9.72	1371.1	0.01
α-Gurjunene	7.31*	1478.0	[0.05]	10.19	1403.9	0.02
β-Caryophyllene	8.12	1540.1	0.05	10.30	1412.2	0.03
β-Gurjunene	8.04	1534.0	0.02	10.42	1421.2	0.01
α-Maaliene	8.35	1558.2	0.01	10.47	1424.9	0.02
Aromadendrene	8.25	1550.3	0.26	10.56	1432.2	0.24
α-Humulene	8.99	1608.9	0.02	10.76	1446.7	0.02
allo-Aromadendrene	8.71	1586.0	0.06	10.85	1453.8	0.06
β-Selinene	9.58	1656.7	0.01	11.19	1479.1	0.01
allo-Aromadendr-9-ene	9.25	1630.2	0.02	11.22	1481.2	0.01
Viridiflorene	9.37	1639.2	0.03	11.33	1489.4	0.04
γ-Cadinene				11.57	1508.1	0.01
δ-Cadinene				11.71	1519.3	0.01

Total reported	99.60%			99.44%		

*: 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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