

Analytical Report	
Title	Bergamot Calabrian Essential Oil Profile by GC-MS
Report No.	SE-37270-3
Issue Date	August 31, 2015
Notebook reference	III-29-75
Contributors:	
Quote No.	
Requester	World Naturals LLC

Primary Aim

To identify GC amendable volatile organic compounds present in submitted **Bergamot Calabrian** essential oil sample.

Samples

The sample arrived as clear liquid with characteristic odor labeled as "Bergamot Calabrian lot 11009-a44".

Experimental:

1. Oil was dissolved in methanol to concentration of ~0.1%, 1 ul injected into the GC injector port.
2. GC conditions:
Injector temperature: 250 C
Initial oven temperature: 80 C
Ramp 10 C/min
Final temperature 220 C
Final temperature hold 5 min

Report SE-37270-3 Bergamot Calabrian

3. MS parameters

Ionization and ion polarity	EI+
Scan rate	2 scans/sec
Mass range	35-350 Da
Ion source temperature	150C
Transfer line temperature	280C

4. GC-MS analysis. Waters/Micromass Quatro GC mass spectrometer interfaced to a ThermoElectron Trace gas chromatograph was utilized for the analysis. 30m 0.25 mm ID DB-5 column was used to separate components. Carrier gas was helium at 1.1 ml/min with split ratio of 50.

5. Data treatment.

For each sample, a set of target components was identified with the aid of the AMDIS software¹. The components were identified using the NIST mass spectral library².

Deliverables

1. GC-MS chromatogram. GC-MS chromatogram is shown in Appendix I.
2. Appendix II lists library search results.
 - RT Retention Time, time in minutes at which the compound elutes out of column
 - CAS. CAS registry number or EPA number.
 - Name. IUPAC or common name of identified compound.
 - Area. Peak area of a component in %% to total ion count

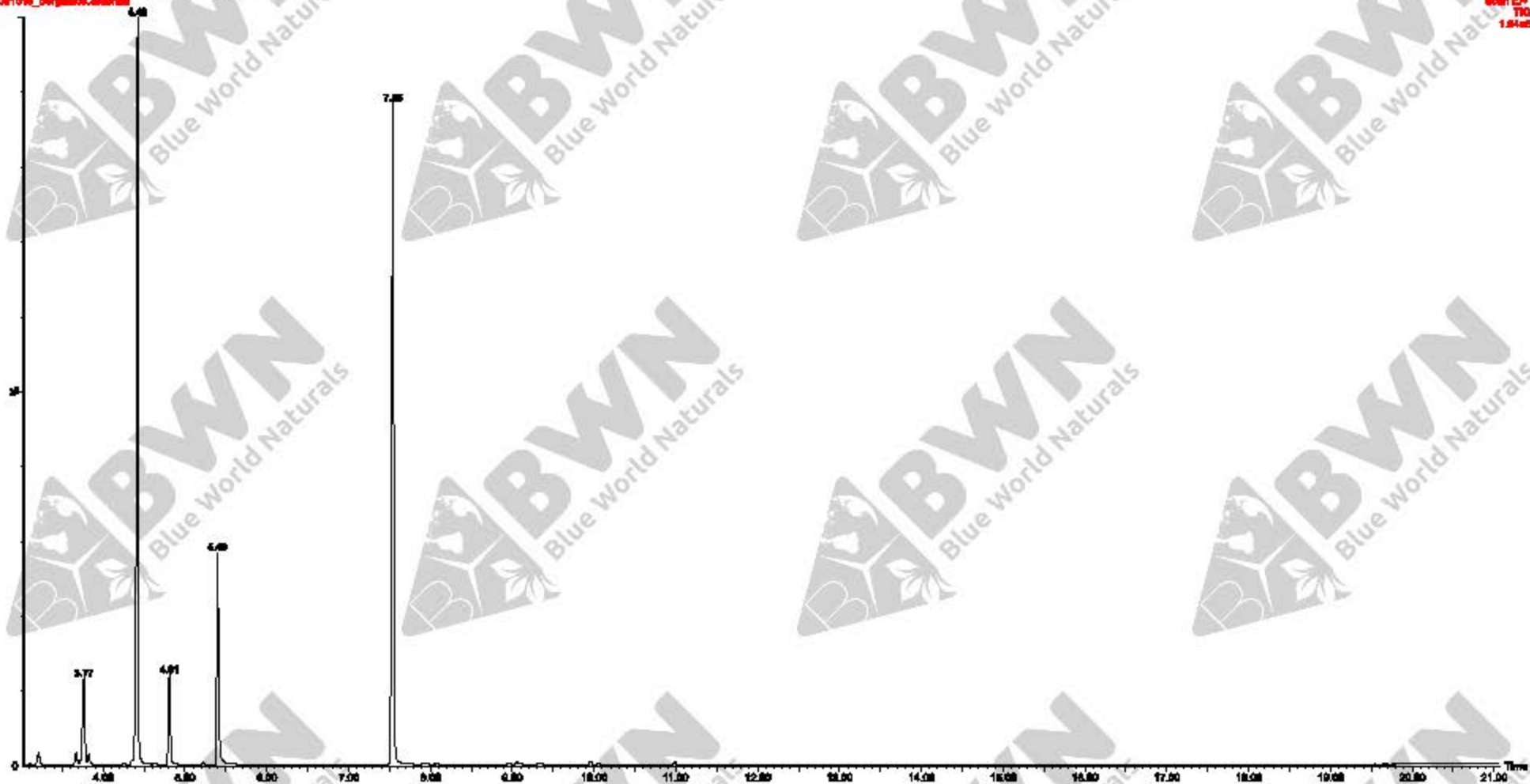
1 <http://chemdata.nist.gov/mass-spc/amdis/>

2 <http://www.nist.gov/srd/nist1a.cfm>

APPENDIX I
Bergamot Calabrian
GC-MS Chromatogram

Sample "Bergamot Calabrian"

BergamotCalabrian on DB-5 Int 11890-046
001090_BergamotCalabrian



APPENDIX II
Bergamot Calabrian
Identified Compounds

Bergamot Calabrian

CAS	Name	R.T.	Area
80-56-8	alpha-pinene	3.082	0.1
2437968	Bicyclo[3.1.1]hept-2-ene, 2,6,6-trimethyl-, (±)-	3.19	0.8
3387415	Bicyclo[3.1.0]hexane, 4-methylene-1-(1-methylethyl)-	3.651	0.7
127913	β-Pinene	3.742	5.9
123353	β-Myrcene	3.803	0.6
99865	1,3-Cyclohexadiene, 1-methyl-4-(1-methylethyl)-	4.229	0.1
527844	Benzene, 1-methyl-2-(1-methylethyl)-	4.335	0.2
138863	Limonene	4.399	36.0
99854	1,4-Cyclohexadiene, 1-methyl-4-(1-methylethyl)-	4.789	5.3
586629	Cyclohexene, 1-methyl-4-(1-methylethylidene)-	5.201	0.1
78706	1,β-Octadien-3-ol, 3,7-dimethyl-	5.38	11.3
5986389	5,7-Octadien-2-ol, 2,6-dimethyl-	6.888	0.0
115-95-7	Bergamiol	7.521	37.7
691383	2-Pentene, 4-methyl-, (Z)-	7.903	0.0
598038	Propane, 1,1'-sulfonylbis-	8.06	0.0
54845306	Cyclohexanol, 2-methylene-3-(1-methylethyl)-, acetate, cis-	8.06	0.0
20777495	Cyclohexanol, 2-methyl-5-(1-methylethenyl)-, acetate, (1α,2β,5α)-	8.929	0.1
141128	2,β-Octadien-1-ol, 3,7-dimethyl-, acetate, (Z)-	9.04	0.2
141128	2,β-Octadien-1-ol, 3,7-dimethyl-, acetate, (Z)-	9.307	0.2
118650	Bicyclo[7.2.0]undec-4-ene, 4,11,11-trimethyl-8-methylene-, [1R-(1R*,4Z,9S*)]-	9.936	0.2
26560145	1,3,β,10-Dodecatetraene, 3,7,11-trimethyl-, (Z,E)-	10.041	0.1
495614	Cyclohexene, 1-methyl-4-(5-methyl-1-methylene-4-hexenyl)-, (S)-	10.969	0.2