

Client/Code

Date

No.

Source 09 Nov 17 2:07p

W136927

Type of Sampleabis - wax

No. of Samples other

1

Comments

Arrival temp.: 23.0C

PD E-transfer Batch 834

Sample: SJ THI

CANNABINOLS

<u>Compounds</u>	<u>Sample</u>	<u>Lab Blank</u>	<u>S_o</u>	<u>Units</u>	<u>reference recovery(%)</u>
Delta-9 THC	0.510	ND	0.001	%	100
Delta-9 THC Acid	77.4	ND	0.001	%	98.2
Delta-8 THC	ND	ND	0.001	%	99.3
Delta-8 THC Acid	ND	ND	0.001		
Cannabichromene (CBC)	ND	ND	0.001	%	99.4
Cannabichromene-Acid	0.700	ND	0.001	%	99.6
Cannabidiol (CBD)	0.070	ND	0.001	%	98.1
Cannabidiol-Acid	0.160	ND	0.001		99.7
Cannabigerol (CBG)	0.660	ND	0.001	%	99.1
Cannabigerol-Acid	1.68	ND	0.001		97.5
Cannabicyclol (CBL)	ND	ND	0.001	%	94.8
Cannabicyclol-Acid	ND	ND	0.001	%	
Cannabidivarin (CBDV)	ND	ND	0.001	%	97.7
Cannabidivarin-Acid	ND	ND	0.001	%	98.1
Delta-9 THCV	ND	ND	0.001	%	99.3
Delta-9 THCV Acid	0.300	ND	0.001		
Cannabinol (CBN)	ND	ND	0.001	%	98.5
Moisture	4.77			%	

Methods: solvent extraction; measured by LC-ESI-MSMS and UPLC-UV.

Pharma.Intern 1.14 & based on USP monograph 29

S_o = standard deviation at zero analyte concentration; method detection limit is generally considered to be 3x S_o value

ND = none detected n/a = not applicable

ug/g = micrograms per gram (ppm), ug/Kg = micrograms per kilogram (ppb)

% = percent (10mg/g = 1.0 %)

^9 -THC = delta 9-tetrahydrocannabinol, ^8 -THC = delta 8-tetrahydrocannabinol

Material will be held for upto 3 weeks unless alternative arrangements have been made. Sample holding time may vary and is dependant upon MBL licence restrictions.



H. Hartmann
H. Hartmann
Sr. Analytical Chemist

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Date

No.

Source 09 Nov 17 2:07p

W136927 P2

Type of Sample Tablets - wax

No. of Samples other

1

Comments

Arrival temp.: 23.0C

PD E-transfer Batch 834

Sample: SJ TH1

Solvent Residue

Solvent	Class	Sample	S _d	Units	Acceptance Criteria
Benzene	I	ND	0.05	ug/g	2 ppm
Methanol	II	ND	3.00	ug/g	3000 ppm
Hexane	II	ND	3.00	ug/g	290 ppm
Toluene	II	ND	3.00	ug/g	890 ppm
Butane	III	ND	3.00	ug/g	5000 ppm
Ethanol	III	ND	3.00	ug/g	5000 ppm
Isopropanol	III	ND	3.00	ug/g	5000 ppm
Pentanes	III	ND	3.00	ug/g	5000 ppm

S_d = standard deviation at zero analyte concentration; method detection limit is generally considered to be 3x S_d value

ND = none detected n/a = not applicable

ref: US Pharmacopeia National Formulary USP 38 NF 33 2015.

Class I to be avoided

Class II to be limited

Class III limit by GMP-low toxicity potential

R. Bilodeau
Analytical Chemist



H. Hartmann
Sr. Analytical Chemist



Client/Code

Date

No.

Source 09 Nov 17 2:07p

W136927 P3

Type of Sample Cannabis - wax

No. of Samples other 1

Comments

Arrival temp.: 23.0C
PD E-transfer Batch 834

Sample: SJ THI

PESTICIDE ANALYSIS SCREEN & TARGET COMPOUNDS -Cannabis

Test Compound Groups*	Sample	Lab	
		Blank	Units
Organophosphates	ND	<0.10	ug/g
Organochlorines	ND	<0.10	ug/g
Carbamates	ND	<0.10	ug/g
Organonitrogens	ND	<0.10	ug/g
Non-Ionic Herbicides	ND	<0.10	ug/g
Ionic Herbicides	ND	<0.10	ug/g
Botanicals	ND	<0.10	ug/g
Carbamates	ND	ND	0.2-37 ng/g
Other	ND	ND	0.2-37 ng/g
-Pyrethroids	ND	ND	0.2-37 ng/g
-Avermectins	ND	ND	0.2-37 ng/g

Method 1: Analysis by GC/MS-MS. Data is analyzed using Agilent RTL Pesticide and Endocrine Disruptor Library with DRS (Deconvolution Reporting Software). The software uses the National Institute of Standards and Technology (NIST) Mass Spectral Search Program with NIST 2011 MS Library. Detection of compounds in the library are 10 ng/g (ppb) level or better. Procedure Ref. AOAC Method 2007.01; USP

Method 2: Analysis is carried out by using UPLC-ESI-MS/MS UV. Sample is solvent extracted then cleaned using dispersive SPE (QuEChERS) methods. Detection of compounds are 0.2-37 ng/g (ppb) level or better. Procedure Ref. AOAC Method 2007.01; USP

* list of total compounds attached


Notes: Some of the compounds analyzed in this method may be naturally occurring & have biological activity. These compounds may not be regulated.

ND = none detected

n/a = not applicable

Acceptance Criteria: none present or are on Health Canada Approved List

R. Bilodeau
Analytical Chemist


H. Hartmann
Sr. Analytical Chemist




Terpene Profile

W136927-1 SJ THI

Compound Name	mg/g	% of Total	Standard Recovery (%)
a-Pinene	6.02	15.55	97.7
Camphene	0.00	0.00	91.0
Sabinene	0.00	0.00	97.1
B-Pinene	2.10	5.44	94.5
Myrcene	24.77	64.00	99.1
a-Phellandrene	0.00	0.00	87.2
3-Carene	0.00	0.00	97.5
Cymene	0.00	0.00	102
D-Limonene	0.94	2.43	105
Eucalyptol	0.00	0.00	108
Ocimene	0.00	0.00	108
Sabinene hydrate	0.00	0.00	82.2
Fenchone	0.00	0.00	101
Terpinolene	0.00	0.00	94.4
Linalol	0.06	0.17	99.1
Fenchol	0.18	0.48	107
Camphor	0.00	0.00	96.3
Isopulegol	0.00	0.00	114
Isoborneol	0.00	0.00	106
Menthofuran	0.00	0.00	94.2
Borneol	0.18	0.47	102
Menthol	0.00	0.00	105
a-Terpineol	0.19	0.48	98.5
Nerol	0.00	0.00	115
Geraniol	0.00	0.00	112
Nerol acetate	0.00	0.00	103
Dihydrojasnone	0.00	0.00	106
Cedrene	0.00	0.00	100
Caryophyllene	2.47	6.38	102
Humulene	1.03	2.66	110
Valencene	0.00	0.00	95.2
Caryophyllene oxide	0.50	1.29	106
Guaiol	0.00	0.00	107
a-Bisabolol	0.26	0.66	104
Total	38.71		

M.B. LABS LTD PO Box 2103 2062 West Henry Ave. Sidney, BC V8L 3S6
T: 250-656-1334 E: mblabs@pacificcoast.net W: www.mblabs.com

R. Bilodeau
Analytical Chemist



H. Hartmann
Sr. Analytical Chemist

Terpene Profile

W136927-1 SJ THI

