

CONSOLIDATED TEST RESULTS SUMMARY

Please see the following pages for full test results.

BULK SKU	BATCH #	LOQ: Limit Of Quantitation	
PRODUCT NAME	SERVING SIZE	LOD: Limit Of Detection	
LABORATORY :	OREGON ACCREDITATION: OR100028	1 g = 10 ⁻³ kg = 10 ³ mg = 10 ⁶ µg 1 mg/kg = 1 ppm = 1000 ppb	
POTENCY	PER SERVING	PER GRAM	Percent
Cannabidiol (CBD)	mg/serving	mg/g	%
Total THC (d9-THC, THCA)	mg/serving	mg/g	%
Cannabigerol (CBG)	mg/serving	mg/g	%
Cannabinol (CBN)	mg/serving	mg/g	%
Cannabichromene (CBC)	mg/serving	mg/g	%
Tetrahydrocannabinolic Acid (THCA)	mg/serving	mg/g	%
Delta-9-THC (d9-THC)	mg/serving	mg/g	%
Delta-8-THC (d8-THC)	mg/serving	mg/g	%
HEAVY METALS	PER SERVING	PER GRAM	REGULATORY ACTION LEVEL
Arsenic	µg/serving	µg/g	10 µg/day ^[1]
Cadmium	µg/serving	µg/g	4.1 µg/day ^[1]
Lead	µg/serving	µg/g	6 µg/day ^[1]
Mercury	µg/serving	µg/g	2 µg/day ^[1]
PESTICIDES	REGULATORY ACTION LEVEL		
None of the other 59 pesticides tested found above limit of detection in the sample.			10 ppb ^[1]
RESIDUAL SOLVENTS	Results	REGULATORY ACTION LEVEL	
Ethanol	µg/g	50,000 mg/day	
Heptane	µg/g	50,000 mg/day	
None of the 34 residual solvents tested found above limit of quantitation in the sample.			
MICROBIAL	PASS/FAIL		
Yeast & Mold	Pass		
Coliform	Pass		
Water Activity	Pass		
TERPENES	% OF SAMPLE		
Farnesene	%		
β-Caryophyllene	%		
α-Bisabolol	%		
Guaiol	%		
Humulene	%		
Caryophyllene Oxide	%		



1. American Herbal Pharmacopoeia. (2014). Cannabis Inflorescence: Standards of Identity, Analysis, and Quality Control. Washington DC: AHP.



12423 NE Whitaker Way
Portland, OR 97230
503-254-1794



Report Number: 23-009990/D004.R000
Report Date: 09/09/2023
ORELAP#: OR100028
Purchase Order: 2606808
Received: 08/22/23 16:37

Customer: Etz Hayim Holdings
Product identity: FORM-TN.O.CBGCBD.FS50-FH74
Client/Metric ID: .
Laboratory ID: 23-009990-0001

Summary

Potency:

Analyte per 1g	Result	Limits	Units	Status	
CBC per 1g	2.44		mg/1g		CBD-Total per Serving Size 28.8 mg/1g
CBD per 1g	28.8		mg/1g		
CBDV per 1g	0.337		mg/1g		THC-Total per Serving Size 1.07 mg/1g
CBE per 1g	0.450		mg/1g		(Reported in milligrams per serving)
CBG per 1g	29.6		mg/1g		
CBL per 1g	0.217		mg/1g		
CBN per 1g	0.0519		mg/1g		
CBT per 1g	1.06		mg/1g		
Δ9-THC per 1g	1.07		mg/1g		

Residual Solvents:

All analytes passing and less than LOQ.

Pesticides:

All analytes passing and less than LOQ.

Terpenes:

Analyte	Percent by weight	Percent of Total	Analyte	Percent by weight	Percent of Total
β-Caryophyllene	0.0599	33.46%	(-)-caryophyllene oxide	0.0436	24.36%
(-)-Guaiol	0.0324	18.10%	Humulene	0.0233	13.02%
α-Bisabolol	0.0202	11.28%	Total Terpenes	0.179	100.00%

Metals:

Less than LOQ for all analytes.

Microbiology:

Less than LOQ for all analytes.



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Purchase Order: 2606808
Received: 08/22/23 16:37

Customer: Etz Hayim Holdings
 16427 NE Airport Way
 PORTLAND 97230
 United States of America (USA)

Product identity: FORM-TN.O.CBGCBBD.FS50-FH74

Client/Metric ID: .

Sample Date:

Laboratory ID: 23-009990-0001

Evidence of Cooling: No

Temp: 22.3 °C

Serving Size #1: 1 g

Sample Results

Potency per 1g		Method: J AOAC 2015 V98-6 (mod) ^b		Units mg/se	Batch: 2310408	Analyze: 8/26/23 1:39:00 AM
Analyte	Result	Limits	Units	LOQ	Notes	
CBC per 1g	2.44		mg/1g	0.0328		
CBC-A per 1g	< LOQ		mg/1g	0.0328		
CBC-Total per 1g	2.44		mg/1g	0.0616		
CBD per 1g	28.8		mg/1g	0.328		
CBD-A per 1g	< LOQ		mg/1g	0.0328		
CBD-Total per 1g	28.8		mg/1g	0.357		
CBDV per 1g	0.337		mg/1g	0.0328		
CBDV-A per 1g	< LOQ		mg/1g	0.0328		
CBDV-Total per 1g	0.337		mg/1g	0.0612		
CBE per 1g	0.450		mg/1g	0.0328		
CBG per 1g	29.6		mg/1g	0.328		
CBG-A per 1g	< LOQ		mg/1g	0.0328		
CBG-Total per 1g	29.6		mg/1g	0.357		
CBL per 1g	0.217		mg/1g	0.0328		
CBL-A per 1g	< LOQ		mg/1g	0.0328		
CBL-Total per 1g	0.217		mg/1g	0.0616		
CBN per 1g	0.0519		mg/1g	0.0328		
CBT per 1g	1.06		mg/1g	0.0328		
Δ8-THCV per 1g	< LOQ		mg/1g	0.0328		
Δ10-THC-9R per 1g	< LOQ		mg/1g	0.0328		
Δ10-THC-9S per 1g	< LOQ		mg/1g	0.0328		
Δ10-THC-Total per 1g	< LOQ		mg/1g	0.0656		
Δ8-THC per 1g	< LOQ		mg/1g	0.0328		
Δ9-THC per 1g	1.07		mg/1g	0.0328		
delta-9-THCP per 1g	< LOQ		mg/1g	0.0328		
exo-THC per 1g	< LOQ		mg/1g	0.0328		
THC-A per 1g	< LOQ		mg/1g	0.0328		
THC-Total per 1g	1.07		mg/1g	0.0616		
THCV per 1g	< LOQ		mg/1g	0.0328		
THCV-A per 1g	< LOQ		mg/1g	0.0328		
THCV-Total per 1g	< LOQ		mg/1g	0.0616		



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Report Number: 23-009990/D004.R000
Report Date: 09/09/2023
ORELAP#: OR100028
Purchase Order: 2606808
Received: 08/22/23 16:37

Potency per 1g **Method:** J AOAC 2015 V98-6 (mod)^P **Units** mg/se **Batch:** 2310408 **Analyze:** 8/26/23 1:39:00 AM

Analyte	Result	Limits	Units	LOQ	Notes
Total Cannabinoids per 1g	64.0		mg/1g		

Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2310543	09/03/23 AOAC 991.14 (Petrifilm) ^P		
Total Coliforms	< LOQ		cfu/g	10	2310543	09/03/23 AOAC 991.14 (Petrifilm) ^P		
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2310544	09/03/23 AOAC 2014.05 (RAPID) ^P		
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2310544	09/03/23 AOAC 2014.05 (RAPID) ^P		

Solvents **Method:** Residual Solvents by GC/MS^P **Units** µg/g **Batch** 2310644 **Analyze** 09/06/23 10:31 AM

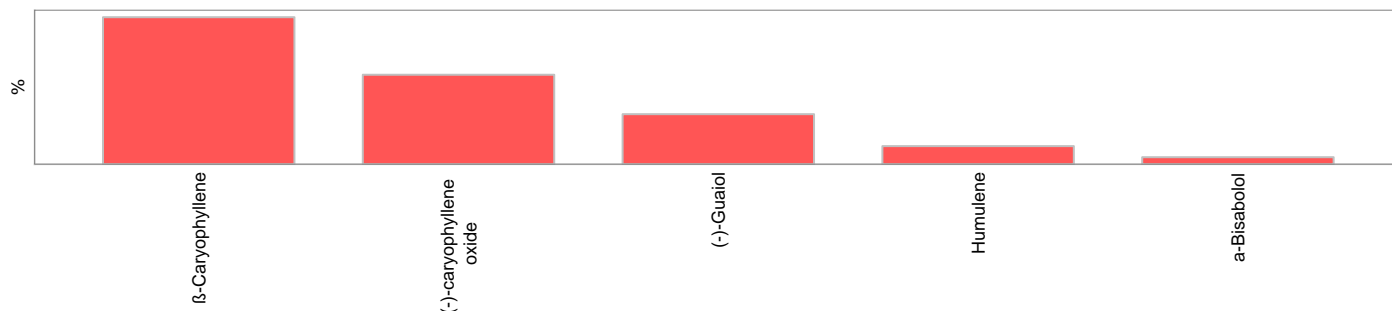
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
1,4-Dioxane	< LOQ	380	100	pass		2-Butanol	< LOQ	5000	200	pass	
2-Ethoxyethanol	< LOQ	160	30.0	pass		2-Methylbutane (Isopentane)	< LOQ		200		
2-Methylpentane	< LOQ		30.0			2-Propanol (IPA)	< LOQ	5000	200	pass	
2,2-Dimethylbutane	< LOQ		30.0			2,2-Dimethylpropane (neo-pentane)	< LOQ		200		
2,3-Dimethylbutane	< LOQ		30.0			3-Methylpentane	< LOQ		30.0		
Acetone	< LOQ	5000	200	pass		Acetonitrile	< LOQ	410	100	pass	
Benzene	< LOQ	2.00	1.00	pass		Butanes (sum)	< LOQ	5000	400	pass	
Cyclohexane	< LOQ	3880	200	pass		Ethanol	< LOQ		200		
Ethyl acetate	< LOQ	5000	200	pass		Ethyl benzene	< LOQ		200		
Ethyl ether	< LOQ	5000	200	pass		Ethylene glycol	< LOQ	620	200	pass	
Ethylene oxide	< LOQ	50.0	20.0	pass		Hexanes (sum)	< LOQ	290	150	pass	
Isopropyl acetate	< LOQ	5000	200	pass		Isopropylbenzene (Cumene)	< LOQ	70.0	30.0	pass	
m,p-Xylene	< LOQ		200			Methanol	< LOQ	3000	200	pass	
Methylene chloride	< LOQ	600	60.0	pass		Methylpropane (Isobutane)	< LOQ		200		
n-Butane	< LOQ		200			n-Heptane	< LOQ	5000	200	pass	
n-Hexane	< LOQ		30.0			n-Pentane	< LOQ		200		
o-Xylene	< LOQ		200			Pentanes (sum)	< LOQ	5000	600	pass	
Propane	< LOQ	5000	200	pass		Tetrahydrofuran	< LOQ	720	100	pass	
Toluene	< LOQ	890	100	pass		Total Xylenes	< LOQ		400		
Total Xylenes and Ethyl benzene	< LOQ	2170	600	pass							



Pesticides											
Method: AOAC 2007.01 & EN 15662 (mod) ^b											
Units mg/kg Batch 2310619 Analyze 09/05/23 01:24 PM											
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
Abamectin [‡]	< LOQ	0.50	0.250	pass		Acephate [‡]	< LOQ	0.40	0.200	pass	
Acequinocyl [‡]	< LOQ	2.0	1.00	pass		Acetamiprid [‡]	< LOQ	0.20	0.100	pass	
Aldicarb [‡]	< LOQ	0.40	0.200	pass		Azoxystrobin [‡]	< LOQ	0.20	0.100	pass	
Bifentazate [‡]	< LOQ	0.20	0.100	pass		Bifenthrin [‡]	< LOQ	0.20	0.100	pass	
Boscalid [‡]	< LOQ	0.40	0.200	pass		Carbaryl [‡]	< LOQ	0.20	0.100	pass	
Carbofuran [‡]	< LOQ	0.20	0.100	pass		Chlorantraniliprole [‡]	< LOQ	0.20	0.100	pass	
Chlorfenapyr [‡]	< LOQ	1.0	0.500	pass		Chlorpyrifos [‡]	< LOQ	0.20	0.100	pass	
Clofentezine [‡]	< LOQ	0.20	0.100	pass		Cyfluthrin [‡]	< LOQ	1.0	0.500	pass	
Cypermethrin [‡]	< LOQ	1.0	0.500	pass		Daminozide [‡]	< LOQ	1.0	0.500	pass	
Diazinon [‡]	< LOQ	0.20	0.100	pass		Dichlorvos [‡]	< LOQ	1.0	0.500	pass	
Dimethoate [‡]	< LOQ	0.20	0.100	pass		Ethoprophos [‡]	< LOQ	0.20	0.100	pass	
Etofenprox [‡]	< LOQ	0.40	0.200	pass		Etoazole [‡]	< LOQ	0.20	0.100	pass	
Fenoxycarb [‡]	< LOQ	0.20	0.100	pass		Fenpyroximate [‡]	< LOQ	0.40	0.200	pass	
Fipronil [‡]	< LOQ	0.40	0.200	pass		Fonicamid [‡]	< LOQ	1.0	0.400	pass	
Fludioxonil [‡]	< LOQ	0.40	0.200	pass		Hexythiazox [‡]	< LOQ	1.0	0.400	pass	
Imazali [‡]	< LOQ	0.20	0.100	pass		Imidacloprid [‡]	< LOQ	0.40	0.200	pass	
Kresoxim-methyl [‡]	< LOQ	0.40	0.200	pass		Malathion [‡]	< LOQ	0.20	0.100	pass	
Metalaxyl [‡]	< LOQ	0.20	0.100	pass		Methiocarb [‡]	< LOQ	0.20	0.100	pass	
Methomyl [‡]	< LOQ	0.40	0.200	pass		MGK-264 [‡]	< LOQ	0.20	0.100	pass	
Myclobutanil [‡]	< LOQ	0.20	0.100	pass		Naled [‡]	< LOQ	0.50	0.250	pass	
Oxamyl [‡]	< LOQ	1.0	0.500	pass		Paclotubrazole [‡]	< LOQ	0.40	0.200	pass	
Parathion-Methyl [‡]	< LOQ	0.20	0.100	pass		Permethrin [‡]	< LOQ	0.20	0.100	pass	
Phosmet [‡]	< LOQ	0.20	0.100	pass		Piperonyl butoxide [‡]	< LOQ	2.0	1.00	pass	
Prallethrin [‡]	< LOQ	0.20	0.100	pass		Propiconazole [‡]	< LOQ	0.40	0.200	pass	
Propoxur [‡]	< LOQ	0.20	0.100	pass		Pyrethrin I (total) [‡]	< LOQ	1.0	0.500	pass	
Pyridaben [‡]	< LOQ	0.20	0.100	pass		Spinosad [‡]	< LOQ	0.20	0.100	pass	
Spiromesifen [‡]	< LOQ	0.20	0.100	pass		Spirotetramat [‡]	< LOQ	0.20	0.100	pass	
Spiroxamine [‡]	< LOQ	0.40	0.200	pass		Tebuconazole [‡]	< LOQ	0.40	0.200	pass	
Thiacloprid [‡]	< LOQ	0.20	0.100	pass		Thiamethoxam [‡]	< LOQ	0.20	0.100	pass	
Trifloxystrobin [‡]	< LOQ	0.20	0.100	pass							



Terpenes				Method: J AOAC 2015 V98-6	Units %	Batch 2310728	Analyze 09/08/23 05:39 AM		
Analyte	Result	LOQ	% of Total	Notes	Analyte	Result	LOQ	% of Total	Notes
β-Caryophyllene	0.0599	0.019	33.4637%		(-)-caryophyllene oxide	0.0436	0.019	24.3575%	
(-)-Guaiol	0.0324	0.019	18.1006%		Humulene	0.0233	0.019	13.0168%	
α-Bisabolol	0.0202	0.019	11.2849%		farnesene	< LOQ	0.019	0.00%	
valencene	< LOQ	0.019	0.00%		β-Myrcene	< LOQ	0.019	0.00%	
(+)-Cedrol	< LOQ	0.019	0.00%		Geraniol	< LOQ	0.019	0.00%	
(-)-β-Pinene	< LOQ	0.019	0.00%		α-Terpinene	< LOQ	0.019	0.00%	
Terpinolene	< LOQ	0.019	0.00%		(R)-(+)-Limonene	< LOQ	0.019	0.00%	
(±)-trans-Nerolidol	< LOQ	0.019	0.00%		Eucalyptol	< LOQ	0.019	0.00%	
Linalool	< LOQ	0.019	0.00%		(±)-fenchone	< LOQ	0.019	0.00%	
nerol	< LOQ	0.019	0.00%		p-Cymene	< LOQ	0.019	0.00%	
(+)-Pulegone	< LOQ	0.019	0.00%		(±)-cis-Nerolidol	< LOQ	0.019	0.00%	
(±)-Camphor	< LOQ	0.019	0.00%		(+)-fenchol	< LOQ	0.019	0.00%	
Isoborneol	< LOQ	0.019	0.00%		(+)-Borneol	< LOQ	0.019	0.00%	
Geranyl acetate	< LOQ	0.019	0.00%		(-)-α-Terpineol	< LOQ	0.019	0.00%	
(-)-Isopulegol	< LOQ	0.019	0.00%		α-cedrene	< LOQ	0.019	0.00%	
α-phellandrene	< LOQ	0.019	0.00%		α-pinene	< LOQ	0.019	0.00%	
Camphene	< LOQ	0.019	0.00%		cis-β-Ocimene	< LOQ	0.006	0.00%	
d-3-Carene	< LOQ	0.019	0.00%		γ-Terpinene	< LOQ	0.019	0.00%	
Menthol	< LOQ	0.019	0.00%		Sabinene	< LOQ	0.019	0.00%	
Sabinene hydrate	< LOQ	0.019	0.00%		trans-β-Ocimene	< LOQ	0.013	0.00%	
Total Terpenes	0.179								



Metals									
Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method		Status	Notes
Arsenic [‡]	< LOQ	0.200	mg/kg	0.0652	2310661	09/06/23	AOAC 2013.06 (mod.) ^p	pass	
Cadmium [‡]	< LOQ	0.200	mg/kg	0.0652	2310661	09/06/23	AOAC 2013.06 (mod.) ^p	pass	
Lead [‡]	< LOQ	0.500	mg/kg	0.0652	2310661	09/06/23	AOAC 2013.06 (mod.) ^p	pass	
Mercury [‡]	< LOQ	0.100	mg/kg	0.0326	2310661	09/06/23	AOAC 2013.06 (mod.) ^p	pass	



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Abbreviations

Limits: Action Levels per OAR-333-007-0400, OAR-333-007-0210, OAR-333-007-0220, CCR title 16-division 42. BCC-section 5723

Limit(s) of Quantitation (LOQ): The minimum levels, concentrations, or quantities of a target variable (e.g., target analyte) that can be reported with a specified degree of confidence.

Ⓟ = ISO/IEC 17025:2017 accredited method.

Ⓜ = TNI accredited analyte.

Units of Measure

cfu/g = Colony forming units per gram

g = g

µg/g = Microgram per gram

mg/kg = Milligram per kilogram = parts per million (ppm)

mg/1g = Milligram per 1g

% = Percentage of sample

% wt = µg/g divided by 10,000

Approved Signatory

Derrick Tanner
General Manager



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8/23/23, 9:50 AM

Columbia Laboratories: ECOC

<https://submissions.us.tentamus.com/cannabis/Lazarus-Naturals-1692743367>

2/2



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Revision: 4 Document ID: 7148
 Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

JAOAC2015 V986 Batch ID: 2310408

Laboratory Control Sample									
Analyte	LCS	Result	Spike	Units	% Rec	Limits		Evaluation	Notes
CBDVA	2	0.0310	0.0309	%	100	80.0	- 120	Acceptable	
CBDV	2	0.0315	0.0313	%	101	80.0	- 120	Acceptable	
CBE	2	0.0327	0.0329	%	99.1	80.0	- 120	Acceptable	
CBDA	1	0.0334	0.0338	%	98.7	90.0	- 110	Acceptable	
CBGA	1	0.0342	0.0343	%	99.8	80.0	- 120	Acceptable	
CBG	1	0.0361	0.0363	%	99.4	80.0	- 120	Acceptable	
CBD	1	0.0349	0.0351	%	99.4	90.0	- 110	Acceptable	
THCV	2	0.0201	0.0200	%	101	80.0	- 120	Acceptable	
d8THCV	2	0.0275	0.0276	%	99.7	80.0	- 120	Acceptable	
THCVA	2	0.0307	0.0307	%	100	80.0	- 120	Acceptable	
CBN	1	0.0344	0.0343	%	100	80.0	- 120	Acceptable	
exo-THC	2	0.0301	0.0302	%	99.6	80.0	- 120	Acceptable	
d9THC	1	0.0358	0.0355	%	101	90.0	- 110	Acceptable	
d8THC	1	0.0367	0.0364	%	101	90.0	- 110	Acceptable	
9S-d10THC	1	0.0357	0.0354	%	101	80.0	- 120	Acceptable	
CBL	2	0.0320	0.0311	%	103	80.0	- 120	Acceptable	
9R-d10THC	1	0.0113	0.0115	%	98.0	80.0	- 120	Acceptable	
CBG	2	0.0332	0.0335	%	99.1	80.0	- 120	Acceptable	
THCA	1	0.0343	0.0344	%	99.8	90.0	- 110	Acceptable	
CBGA	2	0.0320	0.0319	%	100	80.0	- 120	Acceptable	
CBLA	2	0.0650	0.0647	%	101	80.0	- 120	Acceptable	
d9THCP	2	0.0311	0.0316	%	98.3	80.0	- 120	Acceptable	
CBT	2	0.0312	0.0308	%	101	80.0	- 120	Acceptable	

Method Blank							
Analyte	Result	LOQ	Units	Limits		Evaluation	Notes
CBDVA	<LOQ	0.00330	%	< 0.00330		Acceptable	
CBDV	<LOQ	0.00330	%	< 0.00330		Acceptable	
CBE	<LOQ	0.00330	%	< 0.00330		Acceptable	
CBDA	<LOQ	0.00330	%	< 0.00330		Acceptable	
CBGA	<LOQ	0.00330	%	< 0.00330		Acceptable	
CBG	<LOQ	0.00330	%	< 0.00330		Acceptable	
CBD	<LOQ	0.00330	%	< 0.00330		Acceptable	
THCV	<LOQ	0.00330	%	< 0.00330		Acceptable	
d8THCV	<LOQ	0.00330	%	< 0.00330		Acceptable	
THCVA	<LOQ	0.00330	%	< 0.00330		Acceptable	
CBN	<LOQ	0.00330	%	< 0.00330		Acceptable	
exo-THC	<LOQ	0.00330	%	< 0.00330		Acceptable	
d9THC	<LOQ	0.00330	%	< 0.00330		Acceptable	
d8THC	<LOQ	0.00330	%	< 0.00330		Acceptable	
9S-d10THC	<LOQ	0.00330	%	< 0.00330		Acceptable	
CBL	<LOQ	0.00330	%	< 0.00330		Acceptable	
9R-d10THC	<LOQ	0.00330	%	< 0.00330		Acceptable	
CBG	<LOQ	0.00330	%	< 0.00330		Acceptable	
THCA	<LOQ	0.00330	%	< 0.00330		Acceptable	
CBGA	<LOQ	0.00330	%	< 0.00330		Acceptable	
CBLA	<LOQ	0.00330	%	< 0.00330		Acceptable	
d9THCP	<LOQ	0.00330	%	< 0.00330		Acceptable	
CBT	<LOQ	0.00330	%	< 0.00330		Acceptable	

Abbreviations
 ND - None Detected at or above MRL
 RPD - Relative Percent Difference
 LOQ - Limit of Quantitation

Units of Measure:
 % - Percent



12423 NE Whitaker Way
Portland, OR 97230
503-254-1794



Report Number: 23-009990/D004.R000
Report Date: 09/09/2023
ORELAP#: OR100028
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Received: 08/22/23 16:37

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Laboratory Quality Control Results

JAOAC2015 V986		Batch ID: 2310408						
Sample Duplicate		Sample ID: 23-0099510002						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.00328	%	NA	< 20	Acceptable	
CBDV	0.0300	0.0300	0.00328	%	0.0581	< 20	Acceptable	
CBE	<LOQ	<LOQ	0.00328	%	NA	< 20	Acceptable	
CBD	0.0945	0.0941	0.00328	%	0.449	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.00328	%	NA	< 20	Acceptable	
CBG	0.138	0.137	0.00328	%	0.259	< 20	Acceptable	
CBD	9.76	10.0	0.00328	%	2.59	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.00328	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.00328	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.00328	%	NA	< 20	Acceptable	
CBN	0.00653	0.00659	0.00328	%	0.976	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.00328	%	NA	< 20	Acceptable	
d9THC	0.229	0.229	0.00328	%	0.0408	< 20	Acceptable	
d8THC	<LOQ	<LOQ	0.00328	%	NA	< 20	Acceptable	
9S-d10THC	<LOQ	<LOQ	0.00328	%	NA	< 20	Acceptable	
CBL	0.0123	0.0130	0.00328	%	5.84	< 20	Acceptable	
9R-d10THC	<LOQ	<LOQ	0.00328	%	NA	< 20	Acceptable	
CBC	0.276	0.276	0.00328	%	0.0349	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.00328	%	NA	< 20	Acceptable	
CBCA	<LOQ	<LOQ	0.00328	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.00328	%	NA	< 20	Acceptable	
d9THCP	<LOQ	<LOQ	0.00328	%	NA	< 20	Acceptable	
CBT	0.0372	0.0360	0.00328	%	3.43	< 20	Acceptable	

Abbreviations

ND - None Detected at or above MRL
RPD - Relative Percent Difference
LOQ - Limit of Quantitation

Units of Measure:

% - Percent



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Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662		Units: mg/Kg			Batch ID: 2310619			
Method Blank		Laboratory Control Sample						
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes
Abamectin	0.000	< 0.250		1.022	1.000	102.2	50.0	150
Acephate	0.000	< 0.200		0.769	0.800	96.1	60.0	120
Acetamiprid	0.000	< 1.000		3.988	4.000	99.7	40.0	160
Acetamiprid	0.000	< 0.100		0.398	0.400	99.6	60.0	120
Aldicarb	0.000	< 0.200		0.823	0.800	102.8	60.0	120
Azoxystrobin	0.000	< 0.100		0.402	0.400	100.5	60.0	120
Bifenazate	0.000	< 0.100		0.393	0.400	98.3	60.0	120
Bifenthrin	0.000	< 0.100		0.392	0.400	98.1	50.0	150
Boscalid	0.000	< 0.200		0.790	0.800	98.7	60.0	120
Carbaryl	0.000	< 0.100		0.398	0.400	99.6	60.0	120
Carbofuran	0.000	< 0.100		0.412	0.400	103.0	60.0	120
Chlorantraniliprole	0.000	< 0.100		0.385	0.400	96.2	60.0	120
Chlorfenapyr	0.000	< 0.500		1.546	2.000	77.3	60.0	120
Chlorpyrifos	0.004	< 0.100		0.388	0.400	97.0	60.0	120
Clofentazine	0.000	< 0.100		0.364	0.400	91.0	60.0	120
Cyfluthrin	0.000	< 0.500		1.974	2.000	98.7	50.0	150
Cypermethrin	0.000	< 0.500		1.980	2.000	99.0	50.0	150
Daminozide	0.012	< 0.500		0.784	2.000	39.2	60.0	120
Diazinon	0.000	< 0.100		0.429	0.400	107.3	60.0	120
Dichlorvos	0.000	< 0.500		2.034	2.000	101.7	60.0	120
Dimethoate	0.000	< 0.100		0.405	0.400	101.3	60.0	120
Ethoprophos	0.000	< 0.100		0.405	0.400	101.3	60.0	120
Etofenprox	0.000	< 0.200		0.821	0.800	102.6	50.0	150
Etoxazole	0.000	< 0.100		0.429	0.400	107.3	60.0	120
Fenoxycarb	0.000	< 0.100		0.407	0.400	101.8	60.0	120
Fenpyroximate	0.000	< 0.200		0.816	0.800	102.0	60.0	120
Fipronil	0.000	< 0.200		0.796	0.800	99.5	60.0	120
Fonicamid	0.000	< 0.250		1.000	1.000	100.0	60.0	120
Fludioxonil	0.000	< 0.200		0.768	0.800	96.0	50.0	150
Hexythiazox	0.000	< 0.250		1.002	1.000	100.2	60.0	120
Imazalil	0.000	< 0.100		0.387	0.400	96.6	60.0	120
Imidacloprid	0.000	< 0.200		0.814	0.800	101.8	60.0	120
Kresoxim-methyl	0.000	< 0.200		0.782	0.800	97.7	60.0	120
Malathion	0.000	< 0.100		0.400	0.400	99.9	60.0	120
Metaxalyl	0.000	< 0.100		0.405	0.400	101.2	60.0	120
Methiocarb	0.000	< 0.100		0.398	0.400	99.6	60.0	120
Methomyl	0.000	< 0.200		0.797	0.800	99.6	60.0	120
MGK-264	0.000	< 0.100		0.392	0.400	97.9	50.0	150
Myclobutanil	0.000	< 0.100		0.390	0.400	97.6	60.0	120
Naled	0.000	< 0.250		0.993	1.000	99.3	50.0	150
Oxamyl	0.000	< 0.500		2.031	2.000	101.6	60.0	120
Pacllobutrazole	0.000	< 0.200		0.823	0.800	102.9	60.0	120
Parathion-Methyl	0.002	< 0.100		0.428	0.400	107.0	50.0	150
Permethrin	0.000	< 0.100		0.394	0.400	98.5	50.0	150
Phosmet	0.000	< 0.100		0.400	0.400	99.9	50.0	150
Piperonyl butoxide	0.000	< 0.500		2.174	2.000	108.7	60.0	120
Prallethrin	0.000	< 0.100		0.391	0.400	97.8	60.0	120
Propiconazole	0.000	< 0.200		0.774	0.800	96.8	60.0	120
Propoxur	0.000	< 0.100		0.410	0.400	102.6	60.0	120
Pyrethrin (Summe)	0.002	< 0.100		0.493	0.488	101.0	60.0	120
Pyridaben	0.000	< 0.100		0.409	0.400	102.3	50.0	150
Spirosad	0.000	< 0.100		0.408	0.388	105.1	50.0	150
Spiromesifen	0.000	< 0.100		0.384	0.400	96.0	60.0	120
Spirotetramat	0.000	< 0.100		0.397	0.400	99.4	60.0	120
Spiroxamine	0.000	< 0.200		0.820	0.800	102.5	60.0	120
Tebuconazole	0.000	< 0.200		0.812	0.800	101.6	60.0	120
Thiacloprid	0.000	< 0.100		0.400	0.400	100.0	60.0	120
Thiamethoxam	0.000	< 0.100		0.403	0.400	100.8	60.0	120
Trifloxystrobin	0.000	< 0.100		0.409	0.400	102.2	60.0	120

Q6



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Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662		Units: mg/Kg				Batch ID: 2310619				
Matrix Spike/Matrix Spike Duplicate Recoveries		Sample ID: 23-010390-0001								
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes
Abamectin	0.000	1.034	1.013	1.000	2.0%	< 30	103.4%	101.3%	50 - 150	
Acephate	0.000	0.721	0.730	0.800	1.3%	< 30	90.1%	91.3%	50 - 150	
Acequinocyl	0.000	4.087	4.433	4.000	8.1%	< 30	102.2%	110.8%	50 - 150	
Acetamiprid	0.000	0.396	0.411	0.400	3.7%	< 30	99.0%	102.7%	50 - 150	
Aldicarb	0.000	0.815	0.808	0.800	0.9%	< 30	101.9%	101.0%	50 - 150	
Azoxystrobin	0.000	0.395	0.400	0.400	1.2%	< 30	98.8%	99.9%	50 - 150	
Bifenazate	0.000	0.382	0.406	0.400	6.1%	< 30	95.5%	101.5%	50 - 150	
Bifenthrin	0.000	0.389	0.378	0.400	2.8%	< 30	97.2%	94.5%	50 - 150	
Boscalid	0.000	0.763	0.817	0.800	6.8%	< 30	95.3%	102.1%	50 - 150	
Carbaryl	0.000	0.384	0.403	0.400	4.8%	< 30	95.9%	100.6%	50 - 150	
Carbofuran	0.000	0.389	0.396	0.400	1.8%	< 30	97.2%	99.0%	50 - 150	
Chlorantraniliprole	0.000	0.406	0.386	0.400	4.9%	< 30	101.4%	96.6%	50 - 150	
Chlorfenapyr	0.000	1.554	1.568	2.000	0.8%	< 30	77.7%	78.4%	50 - 150	
Chlorpyrifos	0.000	0.367	0.371	0.400	1.1%	< 30	91.8%	92.8%	50 - 150	
Clofentezine	0.000	0.332	0.339	0.400	2.2%	< 30	82.9%	84.7%	50 - 150	
Cyfluthrin	0.000	1.950	2.006	2.000	2.8%	< 30	97.5%	100.3%	30 - 150	
Cypermethrin	0.000	1.867	1.992	2.000	6.5%	< 30	93.3%	99.6%	50 - 150	
Daminozide	0.010	0.752	0.763	2.000	1.5%	< 30	37.1%	37.7%	30 - 150	
Diazinon	0.000	0.404	0.426	0.400	5.2%	< 30	101.0%	106.4%	50 - 150	
Dichlorvos	0.000	1.906	1.959	2.000	2.7%	< 30	95.3%	98.0%	50 - 150	
Dimethoate	0.000	0.397	0.407	0.400	2.5%	< 30	99.3%	101.7%	50 - 150	
Ethoprophos	0.000	0.386	0.398	0.400	3.1%	< 30	96.4%	99.5%	50 - 150	
Etofenprox	0.000	0.801	0.816	0.800	1.9%	< 30	100.1%	102.0%	50 - 150	
Etoxazole	0.000	0.452	0.453	0.400	0.2%	< 30	112.9%	113.2%	50 - 150	
Fenoxycarb	0.000	0.394	0.388	0.400	1.6%	< 30	98.6%	97.0%	50 - 150	
Fenpyroximate	0.000	0.779	0.804	0.800	3.1%	< 30	97.4%	100.5%	50 - 150	
Fipronil	0.000	0.787	0.780	0.800	0.9%	< 30	98.4%	97.5%	50 - 150	
Fonicamid	0.000	0.956	1.014	1.000	5.9%	< 30	95.6%	101.4%	50 - 150	
Fludioxonil	0.000	0.751	0.767	0.800	2.2%	< 30	93.8%	95.9%	50 - 150	
Hexythiazox	0.000	1.195	1.036	1.000	14.3%	< 30	119.5%	103.6%	50 - 150	
Imazalil	0.000	0.394	0.380	0.400	3.6%	< 30	98.4%	95.0%	50 - 150	
Imidacloprid	0.000	0.769	0.774	0.800	0.6%	< 30	96.2%	96.7%	50 - 150	
Kresoxim-methyl	0.000	0.733	0.750	0.800	2.4%	< 30	91.6%	93.8%	50 - 150	
Malathion	0.000	0.387	0.406	0.400	4.8%	< 30	96.6%	101.4%	50 - 150	
Metaxalaxyl	0.000	0.397	0.397	0.400	0.1%	< 30	99.2%	99.3%	50 - 150	
Methiocarb	0.000	0.404	0.394	0.400	2.7%	< 30	101.1%	98.4%	50 - 150	
Methomyl	0.000	0.780	0.833	0.800	6.6%	< 30	97.5%	104.1%	50 - 150	
MGK-264	0.000	0.370	0.410	0.400	10.2%	< 30	92.6%	102.5%	50 - 150	
Myclobutanil	0.000	0.374	0.374	0.400	0.1%	< 30	93.5%	93.6%	50 - 150	
Naled	0.000	0.987	0.931	1.000	5.8%	< 30	98.7%	93.1%	50 - 150	
Oxamyl	0.000	1.949	1.937	2.000	0.6%	< 30	97.4%	96.8%	50 - 150	
Pacllobutrazole	0.000	0.843	0.827	0.800	2.0%	< 30	105.4%	103.3%	50 - 150	
Parathion-Methyl	0.000	0.433	0.400	0.400	7.9%	< 30	108.3%	100.0%	30 - 150	
Permethrin	0.000	0.359	0.390	0.400	8.2%	< 30	89.8%	97.4%	50 - 150	
Phosmet	0.000	0.385	0.420	0.400	8.7%	< 30	96.3%	105.1%	50 - 150	
Piperonyl butoxide	0.000	2.012	2.133	2.000	5.8%	< 30	100.6%	106.6%	50 - 150	
Prallethrin	0.000	0.388	0.378	0.400	2.6%	< 30	96.9%	94.5%	50 - 150	
Propiconazole	0.000	0.762	0.778	0.800	2.0%	< 30	95.3%	97.2%	50 - 150	
Propoxur	0.000	0.388	0.398	0.400	2.7%	< 30	96.9%	99.6%	50 - 150	
Pyrethrin (Summe)	0.000	0.466	0.481	0.488	3.2%	< 30	95.5%	98.6%	50 - 150	
Pyridaben	0.000	0.401	0.410	0.400	2.4%	< 30	100.1%	102.6%	50 - 150	
Spinosad	0.000	0.392	0.411	0.388	4.8%	< 30	100.9%	105.9%	50 - 150	
Spiromesifen	0.000	0.395	0.386	0.400	2.2%	< 30	98.8%	96.6%	50 - 150	
Spirotetramat	0.000	0.391	0.397	0.400	1.5%	< 30	97.7%	99.2%	50 - 150	
Spiroxamine	0.000	0.791	0.804	0.800	1.6%	< 30	98.9%	100.5%	50 - 150	
Tebuconazole	0.000	0.767	0.772	0.800	0.6%	< 30	95.9%	96.5%	50 - 150	
Thiacloprid	0.000	0.383	0.387	0.400	1.1%	< 30	95.8%	96.8%	50 - 150	
Thiamethoxam	0.000	0.380	0.416	0.400	9.0%	< 30	95.0%	104.0%	50 - 150	
Trifloxystrobin	0.000	0.397	0.416	0.400	4.7%	< 30	99.2%	104.0%	50 - 150	

Test results relate only to the parameters tested and to the samples as received by the laboratory. Test results meet all requirements of NELAP and the Columbia Laboratories quality assurance plan unless otherwise noted. This report shall not be reproduced, except in full, without the written consent of this laboratory. Samples will be retained for a maximum of 30 days from the receipt date unless prior arrangements have been made.

Testing in accordance with: OAR 333-007-0390 OAR 333-007-0400 OAR 333-007-0410 OAR 333-007-0430



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Revision: 2 Document ID: 7087
 Legacy ID: CFL-E33Effective:

Laboratory Quality Control Results

Residual Solvents				Batch ID: 2310644					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		541	584	µg/g	92.6	60 - 120	
Isobutane	ND	< 200		715	767	µg/g	93.2	60 - 120	
Butane	ND	< 200		743	782	µg/g	95.0	60 - 120	
2,2-Dimethylpropane	ND	< 200		913	939	µg/g	97.2	60 - 120	
Methanol	ND	< 200		1640	1670	µg/g	98.2	60 - 120	
Ethylene Oxide	ND	< 30		58.3	57.1	µg/g	102.1	60 - 120	
2-Methylbutane	ND	< 200		1520	1680	µg/g	90.5	60 - 120	
Pentane	ND	< 200		1530	1670	µg/g	91.6	60 - 120	
Ethanol	ND	< 200		1640	1660	µg/g	98.8	70 - 130	
Ethyl Ether	ND	< 200		1570	1670	µg/g	94.0	60 - 120	
2,2-Dimethylbutane	ND	< 30		174	189	µg/g	92.1	60 - 120	
Acetone	ND	< 200		1630	1670	µg/g	97.6	60 - 120	
2-Propanol	ND	< 200		1630	1630	µg/g	100.0	60 - 120	
Ethyl Formate	ND	< 500		1520	1600	µg/g	95.0	70 - 130	
Acetonitrile	ND	< 100		478	492	µg/g	97.2	60 - 120	
Methyl Acetate	ND	< 500		1640	1600	µg/g	102.5	70 - 130	
2,3-Dimethylbutane	ND	< 30		174	180	µg/g	96.7	60 - 120	
Dichloromethane	ND	< 60		477	488	µg/g	97.7	60 - 120	
2-Methylpentane	ND	< 30		170	182	µg/g	93.4	60 - 120	
MTBE	ND	< 500		1640	1610	µg/g	101.9	70 - 130	
3-Methylpentane	ND	< 30		166	177	µg/g	93.8	60 - 120	
Hexane	ND	< 30		170	177	µg/g	96.0	60 - 120	
1-Propanol	ND	< 500		1610	1600	µg/g	100.6	70 - 130	
Methylethylketone	ND	< 500		1650	1610	µg/g	102.5	70 - 130	
Ethyl acetate	ND	< 200		1610	1630	µg/g	98.8	60 - 120	
2-Butanol	ND	< 200		1650	1630	µg/g	101.2	60 - 120	
Tetrahydrofuran	ND	< 100		491	488	µg/g	100.6	60 - 120	
Cyclohexane	ND	< 200		1580	1610	µg/g	98.1	60 - 120	
2-methyl-1-propanol	ND	< 500		1620	1610	µg/g	100.6	70 - 130	
Benzene	ND	< 1		4.99	4.79	µg/g	104.2	60 - 120	
Isopropyl Acetate	ND	< 200		1680	1650	µg/g	101.8	60 - 120	
Heptane	ND	< 200		1630	1630	µg/g	100.0	60 - 120	
1-Butanol	ND	< 500		1540	1600	µg/g	96.3	70 - 130	
Propyl Acetate	ND	< 500		1630	1600	µg/g	101.9	70 - 130	
1,4-Dioxane	ND	< 100		506	523	µg/g	96.7	60 - 120	
2-Ethoxyethanol	ND	< 30		195	179	µg/g	108.9	60 - 120	
Methylisobutylketone	ND	< 500		1630	1600	µg/g	101.9	70 - 130	
3-Methyl-1-butanol	ND	< 500		1510	1600	µg/g	94.4	70 - 130	
Ethylene Glycol	ND	< 200		283	506	µg/g	55.9	60 - 120	Q6
Toluene	ND	< 100		521	496	µg/g	105.0	60 - 120	
Isobutyl Acetate	ND	< 500		1710	1610	µg/g	106.2	70 - 130	
1-Pentanol	ND	< 500		1480	1600	µg/g	92.5	70 - 130	
Butyl Acetate	ND	< 500		1650	1610	µg/g	102.5	70 - 130	
Ethylbenzene	ND	< 200		1000	978	µg/g	102.2	60 - 120	
m,p-Xylene	ND	< 200		1040	994	µg/g	104.6	60 - 120	
o-Xylene	ND	< 200		1010	982	µg/g	102.9	60 - 120	
Cumene	ND	< 30		151	171	µg/g	88.3	60 - 120	
Anisole	ND	< 500		1670	1600	µg/g	104.4	70 - 130	
DMSO	ND	< 500		1730	1620	µg/g	106.8	70 - 130	
1,2-dimethoxyethane	ND	< 50		184	185	µg/g	98.9	70 - 130	
Triethylamine	ND	< 500		1550	1600	µg/g	96.9	70 - 130	
N,N-dimethylformamide	ND	< 150		533	480	µg/g	111.0	70 - 130	
N,N-dimethylacetamide	ND	< 150		481	483	µg/g	99.6	70 - 130	
Pyridine	ND	< 50		172	168	µg/g	102.4	70 - 130	
Sulfone	ND	< 50		139	161	µg/g	86.3	70 - 130	
1,2-Dichloroethane	ND	< 1		1.06	1	µg/g	106.0	70 - 130	
Chloroform	ND	< 1		1.05	1	µg/g	105.0	70 - 130	
Trichloroethylene	ND	< 1		1.02	1	µg/g	102.0	70 - 130	
1,1-Dichloroethane	ND	< 1		1.02	1	µg/g	102.0	70 - 130	



12423 NE Whitaker Way
Portland, OR 97230
503-254-1794



Report Number: 23-009990/D004.R000
Report Date: 09/09/2023
ORELAP#: OR100028
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Legacy ID: CFL-E33Effective:

QC - Sample Duplicate		Sample ID: 23-010144-0001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Accept/Fail	Notes
Propane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Isobutane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Butane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2,2-Dimethylpropane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Methanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethylene Oxide	ND	ND	30	µg/g	0.0	< 20	Acceptable	
2-Methylbutane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Pentane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethyl Ether	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2,2-Dimethylbutane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Acetone	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2-Propanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethyl Formate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Acetonitrile	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Methyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
2,3-Dimethylbutane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Dichloromethane	ND	ND	60	µg/g	0.0	< 20	Acceptable	
2-Methylpentane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
MTBE	ND	ND	500	µg/g	0.0	< 20	Acceptable	
3-Methylpentane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Hexane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
1-Propanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Methyl ethyl ketone	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Ethyl acetate	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2-Butanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Tetrahydrofuran	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Cyclohexane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2-methyl-1-propanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Benzene	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Isopropyl Acetate	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Heptane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
1-Butanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Propyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
1,4-Dioxane	ND	ND	100	µg/g	0.0	< 20	Acceptable	
2-Ethoxyethanol	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Methylisobutylketone	ND	ND	500	µg/g	0.0	< 20	Acceptable	
3-Methyl-1-butanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Ethylene Glycol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Toluene	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Isobutyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
1-Pentanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Butyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Ethylbenzene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
m,p-Xylene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
o-Xylene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Cumene	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Anisole	ND	ND	500	µg/g	0.0	< 20	Acceptable	
DMSO	ND	ND	500	µg/g	0.0	< 20	Acceptable	
1,2-dimethoxyethane	ND	ND	50	µg/g	0.0	< 20	Acceptable	
Triethylamine	ND	ND	500	µg/g	0.0	< 20	Acceptable	
N,N-dimethylformamide	ND	ND	150	µg/g	0.0	< 20	Acceptable	
N,N-dimethylacetamide	ND	ND	150	µg/g	0.0	< 20	Acceptable	
Pyridine	ND	ND	50	µg/g	0.0	< 20	Acceptable	
Sulfolane	ND	ND	50	µg/g	0.0	< 20	Acceptable	
1,2-Dichloroethane	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Chloroform	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Trichloroethylene	ND	ND	1	µg/g	0.0	< 20	Acceptable	
1,1-Dichloroethane	ND	ND	1	µg/g	0.0	< 20	Acceptable	

Abbreviations

ND - None Detected at or above MRL
RPD - Relative Percent Difference
LOQ - Limit of Quantitation
Q6 - Quality control outside QC limits. Data acceptable based on remaining QC.

Units of Measure:

µg/g- Microgram per gram or ppm


 Revision: 1 Document ID: 7086
 Legacy ID: CFL-E57Worksheet Validated 11/04/2020

Terpenes Quality Control Results

Method Reference: EPA5035				Batch ID: 2310728					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	LCS	Units	LCS% Rec	Limits	Notes
a-pinene	<LOQ	< 200		422	500	µg/g	84%	70 - 130	
Camphene	<LOQ	< 200		517	500	µg/g	103%	70 - 130	
Sabinene	<LOQ	< 200		416	500	µg/g	83%	70 - 130	
b-Pinene	<LOQ	< 200		417	500	µg/g	83%	70 - 130	
b-Myrcene	<LOQ	< 200		544	500	µg/g	109%	70 - 130	
a-phellandrene	<LOQ	< 200		555	500	µg/g	111%	70 - 130	
d-3-Carene	<LOQ	< 200		530	500	µg/g	106%	70 - 130	
a-Terpinene	<LOQ	< 200		435	500	µg/g	87%	70 - 130	
p-Cymene	<LOQ	< 200		538	500	µg/g	108%	70 - 130	
D-Limonene	<LOQ	< 200		448	500	µg/g	90%	70 - 130	
Eucalyptol	<LOQ	< 200		552	500	µg/g	110%	70 - 130	
b-cis-Cimene	<LOQ	< 67		186	167	µg/g	111%	70 - 130	
b-trans-Cimene	<LOQ	< 133		375	333	µg/g	113%	70 - 130	
g-Terpinene	<LOQ	< 200		450	500	µg/g	90%	70 - 130	
Sabinene Hydrate	<LOQ	< 200		434	500	µg/g	87%	70 - 130	
Terpinolene	<LOQ	< 200		442	500	µg/g	88%	70 - 130	
D-Fenchone	<LOQ	< 200		426	500	µg/g	85%	70 - 130	
Linalool	<LOQ	< 200		642	500	µg/g	128%	70 - 130	
Fenchol	<LOQ	< 200		461	500	µg/g	92%	70 - 130	
Camphor	<LOQ	< 200		557	500	µg/g	111%	70 - 130	
Isopulego	<LOQ	< 200		591	500	µg/g	118%	70 - 130	
Isoborneol	<LOQ	< 200		592	500	µg/g	118%	70 - 130	
Borneol	<LOQ	< 200		449	500	µg/g	90%	70 - 130	
DL-Menthol	<LOQ	< 200		620	500	µg/g	124%	70 - 130	
Terpineol	<LOQ	< 200		459	500	µg/g	92%	70 - 130	
Nerd	<LOQ	< 200		564	500	µg/g	113%	70 - 130	
Pulegone	<LOQ	< 200		480	500	µg/g	96%	70 - 130	
Geraniol	<LOQ	< 200		483	500	µg/g	97%	70 - 130	
Geranyl_Acetate	<LOQ	< 200		571	500	µg/g	114%	70 - 130	
a-Cedrene	<LOQ	< 200		425	500	µg/g	85%	70 - 130	
b-Caryophyllene	<LOQ	< 200		570	500	µg/g	114%	70 - 130	
a-Humulene	<LOQ	< 200		478	500	µg/g	96%	70 - 130	
Valene	<LOQ	< 200		543	500	µg/g	109%	70 - 130	
cis-Nerolidol	<LOQ	< 200		613	500	µg/g	123%	70 - 130	
a-Farnesene	<LOQ	< 200		640	500	µg/g	128%	70 - 130	
trans-Nerolidol	<LOQ	< 200		479	500	µg/g	96%	70 - 130	
Caryophyllene_Oxide	<LOQ	< 200		585	500	µg/g	117%	70 - 130	
Guaiol	<LOQ	< 200		468	500	µg/g	94%	70 - 130	
Cedrol	<LOQ	< 200		625	500	µg/g	125%	70 - 130	
a-Bisabolol	<LOQ	< 200		626	500	µg/g	125%	70 - 130	

Definitions

LOQ	Limit of Quantitation
LCS	Laboratory Control Sample
% REC	Percent Recovery



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Terpenes Quality Control Results

Method Reference: EPA5035		Batch ID: 2310728					
Sample/ Sample Duplicate		Sample ID: 23-009990-001					
Analyte	Result	Org. Result	LOQ	Units	% RPD	LIMIT	Notes
a-pinene	<LOQ	<LOQ	194	µg/g	0%	< 20	
Camphene	<LOQ	<LOQ	194	µg/g	0%	< 20	
Sabinene	<LOQ	<LOQ	194	µg/g	0%	< 20	
b-Pinene	<LOQ	<LOQ	194	µg/g	0%	< 20	
b-Myrcene	<LOQ	<LOQ	194	µg/g	0%	< 20	
a-phellandrene	<LOQ	<LOQ	194	µg/g	0%	< 20	
d-3-Carene	<LOQ	<LOQ	194	µg/g	0%	< 20	
a-Terpinene	<LOQ	<LOQ	194	µg/g	0%	< 20	
p-Cymene	<LOQ	<LOQ	194	µg/g	0%	< 20	
D-Limonene	<LOQ	<LOQ	194	µg/g	0%	< 20	
Eucalyptol	<LOQ	<LOQ	194	µg/g	0%	< 20	
b-cis-Cimene	<LOQ	<LOQ	64.8	µg/g	0%	< 20	
b-trans-Cimene	<LOQ	<LOQ	130	µg/g	0%	< 20	
g-Terpinene	<LOQ	<LOQ	194	µg/g	0%	< 20	
Sabinene Hydrate	<LOQ	<LOQ	194	µg/g	0%	< 20	
Terpinolene	<LOQ	<LOQ	194	µg/g	0%	< 20	
D-Fenchone	<LOQ	<LOQ	194	µg/g	0%	< 20	
Linalool	<LOQ	<LOQ	194	µg/g	0%	< 20	
Fenchol	<LOQ	<LOQ	194	µg/g	0%	< 20	
Camphor	<LOQ	<LOQ	194	µg/g	0%	< 20	
Isopulego	<LOQ	<LOQ	194	µg/g	0%	< 20	
Isoborneol	<LOQ	<LOQ	194	µg/g	0%	< 20	
Borneol	<LOQ	<LOQ	194	µg/g	0%	< 20	
DL-Menthhol	<LOQ	<LOQ	194	µg/g	0%	< 20	
Terpineol	<LOQ	<LOQ	194	µg/g	0%	< 20	
Nerd	<LOQ	<LOQ	194	µg/g	0%	< 20	
Pulegone	<LOQ	<LOQ	194	µg/g	0%	< 20	
Geraniol	<LOQ	<LOQ	194	µg/g	0%	< 20	
Geranyl Acetate	<LOQ	<LOQ	194	µg/g	0%	< 20	
a-Cedrene	<LOQ	<LOQ	194	µg/g	0%	< 20	
b-Caryophyllene	594	599	194	µg/g	1%	< 20	
a-Humulene	229	233	194	µg/g	2%	< 20	
Valnene	<LOQ	<LOQ	194	µg/g	0%	< 20	
cis-Nerolidol	<LOQ	<LOQ	194	µg/g	0%	< 20	
a-Farnesene	<LOQ	<LOQ	194	µg/g	0%	< 20	
trans-Nerolidol	<LOQ	<LOQ	194	µg/g	0%	< 20	
Caryophyllene Oxide	432	436	194	µg/g	1%	< 20	
Guaiol	318	324	194	µg/g	2%	< 20	
Cedrol	<LOQ	<LOQ	194	µg/g	0%	< 20	
a-Bisabolol	200	202	194	µg/g	1%	< 20	

Definitions

RPD Relative Percent Difference



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Explanation of QC Flag Comments:

Code	Explanation
Q	Matrix interferences affecting spike or surrogate recoveries.
Q1	Quality control result biased high. Only non-detect samples reported.
Q2	Quality control outside QC limits. Data considered estimate.
Q3	Sample concentration greater than four times the amount spiked.
Q4	Non-homogenous sample matrix, affecting RPD result and/or % recoveries.
Q5	Spike results above calibration curve.
Q6	Quality control outside QC limits. Data acceptable based on remaining QC.
R	Relative percent difference (RPD) outside control limit.
R1	RPD non-calculable, as sample or duplicate results are less than five times the LOQ.
R2	Sample replicates RPD non-calculable, as only one replicate is within the analytical range.
LOQ1	Quantitation level raised due to low sample volume and/or dilution.
LOQ2	Quantitation level raised due to matrix interference.
B	Analyte detected in method blank, but not in associated samples.
B1	The sample concentration is greater than 5 times the blank concentration.
B2	The sample concentration is less than 5 times the blank concentration.