

## CONSOLIDATED TEST RESULTS SUMMARY

Please see the following pages for full test results.

<b>BULK SKU</b>	<b>BATCH #</b>	<b>LOQ: Limit Of Quantitation</b>	
<b>PRODUCT NAME</b>	<b>SERVING SIZE</b>	<b>LOD: Limit Of Detection</b>	
<b>LABORATORY :</b>	<b>OREGON ACCREDITATION: OR100028</b>	1 g = 10 <sup>-3</sup> kg = 10 <sup>3</sup> mg = 10 <sup>6</sup> 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	1.5 ppm
Cadmium	µg/serving	µg/g	0.5 ppm
Lead	µg/serving	µg/g	0.5 ppm
Mercury	µg/serving	µg/g	3.0 ppm
PESTICIDES	None of the other 59 pesticides tested found above limit of detection in the sample.		REGULATORY ACTION LEVEL
			10 ppb <sup>[1]</sup>
RESIDUAL SOLVENTS	Results	REGULATORY ACTION LEVEL	
Ethanol*	µg/g	5,000 ppm	
Heptane	µg/g	5,000 ppm	
None of the 34 residual solvents tested found above limit of quantitation in the sample.			
MICROBIAL	PASS/FAIL		
Yeast & Mold	Pass		
Coliform	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.

\*Ethanol is a food additive used in some of our ingredients. The FDA has labeled ethanol as Generally Recognized as Safe (GRAS). Many foods contain trace amounts of ethanol, including soy sauce, pasta sauces, fruits and juices, etc. Our products contain safe levels of ethanol and always below pertinent regulatory action levels.



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



**Report Number:** 23-006471/D012.R000  
**Report Date:** 06/22/2023  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 05/31/23 16:23

**Customer:** Etz Hayim Holdings  
**Product identity:** FORM-SG100.V2-  
**Client/Metric ID:** FE32-23081101.  
**Laboratory ID:** 23-006471-0004

### Summary

**Potency:**

Analyte per 1g	Result	Limits	Units	Status	
CBC per 1g	10.3		mg/1g		CBD-Total per Serving Size 259 mg/1g
CBD per 1g	259		mg/1g		
CBDV per 1g	2.01		mg/1g		THC-Total per Serving Size 1.84 mg/1g (Reported in milligrams per serving)
CBE per 1g	21.4		mg/1g		
CBG per 1g	2.30		mg/1g		
CBL per 1g	1.49		mg/1g		
CBN per 1g	2.24		mg/1g		
CBT per 1g	10.8		mg/1g		
Δ9-THC per 1g	1.84		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
(R)-(+)-Limonene	3.18	45.04%	a-pinene	2.76	39.09%
β-Caryophyllene	0.704	9.97%	(-)-caryophyllene oxide	0.138	1.95%
α-Bisabolol	0.112	1.59%	Humulene	0.0827	1.17%
β-Myrcene	0.0429	0.61%	Camphene	0.0395	0.56%
<b>Total Terpenes</b>	<b>7.06</b>	<b>100.00%</b>			

**Metals:**

Less than LOQ for all analytes.

**Microbiology:**

Less than LOQ for all analytes.



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**Microbiology**

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
Aerobic Plate Count	< LOQ		cfu/g	10	2308277	06/18/23 AOAC 990.12 (Petrifilm) <sup>P</sup>		I
E.coli	< LOQ		cfu/g	10	2308275	06/18/23 AOAC 991.14 (Petrifilm) <sup>P</sup>		I
Total Coliforms	< LOQ		cfu/g	10	2308275	06/18/23 AOAC 991.14 (Petrifilm) <sup>P</sup>		I
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2308276	06/18/23 AOAC 2014.05 (RAPID) <sup>P</sup>		I
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2308276	06/18/23 AOAC 2014.05 (RAPID) <sup>P</sup>		I

**Solvents** Method: Residual Solvents by GC/MS<sup>B</sup> Units µg/g Batch 2308405 Analyze 06/21/23 09:20 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							

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



12423 NE Whitaker Way  
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 503-254-1794



**Report Number:** 23-011802/D002.R001  
**Report Date:** 10/19/2023  
**ORELAP#:** OR100028  
**Purchase Order:** 2663180  
**Received:** 10/04/23 16:16

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

**Product identity:** FORM-SG100.V2-23081101-3

**Client/Metric ID:** .

**Sample Date:**

**Laboratory ID:** 23-011802-0003

**Evidence of Cooling:** No

**Temp:** 24 °C

**Relinquished by:** client

**Serving Size #1:** 1 g

### Sample Results

Potency per 1g					
Method: J AOAC 2015 V98-6 (mod) <sup>b</sup>		Units mg/se		Batch: 2311963	
				Analyze: 10/18/23 12:12:00 A	
Analyte	Result	Limits	Units	LOQ	Notes
CBC per 1g	10.3		mg/1g	3.06	
CBC-A per 1g	< LOQ		mg/1g	0.0306	
CBC-Total per 1g	10.3		mg/1g	3.08	
CBD per 1g	259		mg/1g	3.06	
CBD-A per 1g	< LOQ		mg/1g	0.0306	
CBD-Total per 1g	259		mg/1g	3.08	
CBDV per 1g	2.01		mg/1g	0.0306	
CBDV-A per 1g	< LOQ		mg/1g	0.0306	
CBDV-Total per 1g	2.01		mg/1g	0.0571	
CBE per 1g	21.4		mg/1g	3.06	
CBG per 1g	2.30		mg/1g	0.0306	
CBG-A per 1g	< LOQ		mg/1g	0.0306	
CBG-Total per 1g	2.30		mg/1g	0.0571	
CBL per 1g	1.49		mg/1g	0.0306	
CBL-A per 1g	< LOQ		mg/1g	0.0306	
CBL-Total per 1g	1.49		mg/1g	0.0574	
CBN per 1g	2.24		mg/1g	0.0306	
CBT per 1g	10.8		mg/1g	3.06	
Δ8-THCV per 1g	< LOQ		mg/1g	0.0306	
Δ10-THC-9R per 1g	< LOQ		mg/1g	0.0306	
Δ10-THC-9S per 1g	< LOQ		mg/1g	0.0306	
Δ10-THC-Total per 1g	< LOQ		mg/1g	0.0611	
Δ8-THC per 1g	< LOQ		mg/1g	0.0306	
Δ9-THC per 1g	1.84		mg/1g	0.0306	
delta-9-THCP per 1g	< LOQ		mg/1g	0.0306	
exo-THC per 1g	< LOQ		mg/1g	0.0306	
THC-A per 1g	< LOQ		mg/1g	0.0306	
THC-Total per 1g	1.84		mg/1g	0.0574	
THCV per 1g	< LOQ		mg/1g	0.0306	
THCV-A per 1g	< LOQ		mg/1g	0.0306	



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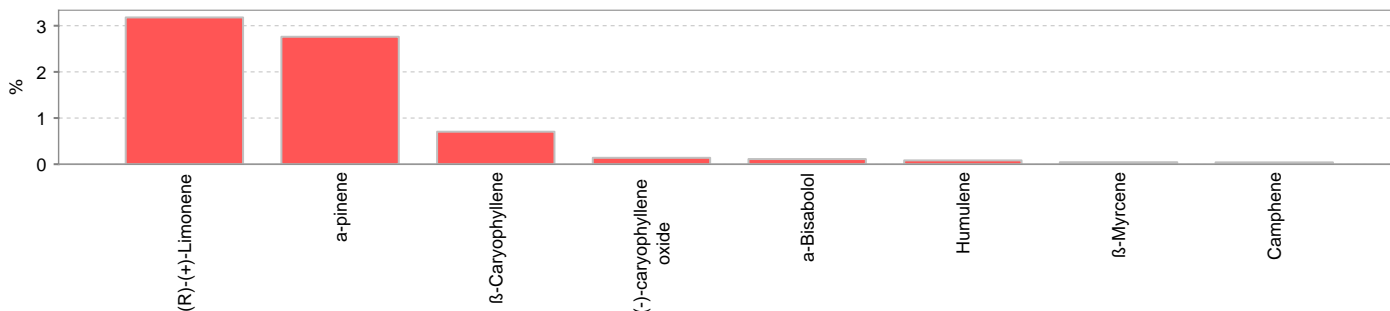


**Report Number:** 23-006471/D012.R000  
**Report Date:** 06/22/2023  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 05/31/23 16:23

Pesticides											
Method: AOAC 2007.01 & EN 15662 (mod) <sup>b</sup>											
Units mg/kg Batch 2308371 Analyze 06/20/23 04:12 PM											
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
Abamectin <sup>‡</sup>	< LOQ	0.50	0.250	pass		Acephate <sup>‡</sup>	< LOQ	0.40	0.200	pass	
Acequinocyl <sup>‡</sup>	< LOQ	2.0	1.00	pass		Acetamiprid <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Aldicarb <sup>‡</sup>	< LOQ	0.40	0.200	pass		Azoxystrobin <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Bifenazate <sup>‡</sup>	< LOQ	0.20	0.100	pass		Bifenthrin <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Boscalid <sup>‡</sup>	< LOQ	0.40	0.200	pass		Carbaryl <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Carbofuran <sup>‡</sup>	< LOQ	0.20	0.100	pass		Chlorantraniliprole <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Chlorfenapyr <sup>‡</sup>	< LOQ	1.0	0.500	pass		Chlorpyrifos <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Clofentezine <sup>‡</sup>	< LOQ	0.20	0.100	pass		Cyfluthrin <sup>‡</sup>	< LOQ	1.0	0.500	pass	
Cypermethrin <sup>‡</sup>	< LOQ	1.0	0.500	pass		Daminozide <sup>‡</sup>	< LOQ	1.0	0.500	pass	
Diazinon <sup>‡</sup>	< LOQ	0.20	0.100	pass		Dichlorvos <sup>‡</sup>	< LOQ	1.0	0.500	pass	
Dimethoate <sup>‡</sup>	< LOQ	0.20	0.100	pass		Ethoprophos <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Etofenprox <sup>‡</sup>	< LOQ	0.40	0.200	pass		Etoxazole <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Fenoxycarb <sup>‡</sup>	< LOQ	0.20	0.100	pass		Fenpyroximate <sup>‡</sup>	< LOQ	0.40	0.200	pass	
Fipronil <sup>‡</sup>	< LOQ	0.40	0.200	pass		Fonicamid <sup>‡</sup>	< LOQ	1.0	0.400	pass	
Fludioxonil <sup>‡</sup>	< LOQ	0.40	0.200	pass		Hexythiazox <sup>‡</sup>	< LOQ	1.0	0.400	pass	
Imazalil <sup>‡</sup>	< LOQ	0.20	0.100	pass		Imidacloprid <sup>‡</sup>	< LOQ	0.40	0.200	pass	
Kresoxim-methyl <sup>‡</sup>	< LOQ	0.40	0.200	pass		Malathion <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Metalaxyl <sup>‡</sup>	< LOQ	0.20	0.100	pass		Methiocarb <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Methomyl <sup>‡</sup>	< LOQ	0.40	0.200	pass		MGK-264 <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Myclobutanil <sup>‡</sup>	< LOQ	0.20	0.100	pass		Naled <sup>‡</sup>	< LOQ	0.50	0.250	pass	
Oxamyl <sup>‡</sup>	< LOQ	1.0	0.500	pass		Paclotbutrazole <sup>‡</sup>	< LOQ	0.40	0.200	pass	
Parathion-Methyl <sup>‡</sup>	< LOQ	0.20	0.100	pass		Permethrin <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Phosmet <sup>‡</sup>	< LOQ	0.20	0.100	pass		Piperonyl butoxide <sup>‡</sup>	< LOQ	2.0	1.00	pass	
Prallethrin <sup>‡</sup>	< LOQ	0.20	0.100	pass		Propiconazole <sup>‡</sup>	< LOQ	0.40	0.200	pass	
Propoxur <sup>‡</sup>	< LOQ	0.20	0.100	pass		Pyrethrin I (total) <sup>‡</sup>	< LOQ	1.0	0.500	pass	
Pyridaben <sup>‡</sup>	< LOQ	0.20	0.100	pass		Spinosad <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Spiromesifen <sup>‡</sup>	< LOQ	0.20	0.100	pass		Spirotetramat <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Spiroxamine <sup>‡</sup>	< LOQ	0.40	0.200	pass		Tebuconazole <sup>‡</sup>	< LOQ	0.40	0.200	pass	
Thiacloprid <sup>‡</sup>	< LOQ	0.20	0.100	pass		Thiamethoxam <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Trifloxystrobin <sup>‡</sup>	< LOQ	0.20	0.100	pass							



Terpenes				Method: J AOAC 2015 V98-6	Units %	Batch 2308421	Analyze 06/20/23 01:59 PM		
Analyte	Result	LOQ	% of Total	Notes	Analyte	Result	LOQ	% of Total	Notes
(R)-(+)-Limonene	3.18	0.019	45.04%		a-pinene	2.76	0.019	39.09%	
β-Caryophyllene	0.704	0.019	9.972%		(-)-caryophyllene oxide	0.138	0.019	1.955%	
α-Bisabolol	0.112	0.019	1.586%		Humulene	0.0827	0.019	1.1714%	
β-Myrcene	0.0429	0.019	0.6076%		Camphene	0.0395	0.019	0.5595%	
(-)-β-Pinene	< LOQ	0.019	0.00%		Linalool	< LOQ	0.019	0.00%	
(±)-trans-Nerolidol	< LOQ	0.019	0.00%		Geranyl acetate	< LOQ	0.019	0.00%	
(+)-Cedrol	< LOQ	0.019	0.00%		(-)-Guaiol	< LOQ	0.019	0.00%	
(+)-fenchol	< LOQ	0.019	0.00%		Sabinene	< LOQ	0.019	0.00%	
Geraniol	< LOQ	0.019	0.00%		(±)-Camphor	< LOQ	0.019	0.00%	
(+)-Borneol	< LOQ	0.019	0.00%		(±)-fenchone	< LOQ	0.019	0.00%	
valencene	< LOQ	0.019	0.00%		Isoborneol	< LOQ	0.019	0.00%	
(±)-cis-Nerolidol	< LOQ	0.019	0.00%		(+)-Pulegone	< LOQ	0.019	0.00%	
nerol	< LOQ	0.019	0.00%		(-)-Isopulegol	< LOQ	0.019	0.00%	
Menthol	< LOQ	0.019	0.00%		(-)-a-Terpineol	< LOQ	0.019	0.00%	
a-Terpinene	< LOQ	0.019	0.00%		a-cedrene	< LOQ	0.019	0.00%	
a-phellandrene	< LOQ	0.019	0.00%		cis-β-Ocimene	< LOQ	0.006	0.00%	
d-3-Carene	< LOQ	0.019	0.00%		Eucalyptol	< LOQ	0.019	0.00%	
farnesene	< LOQ	0.019	0.00%		gamma-Terpinene	< LOQ	0.019	0.00%	
p-Cymene	< LOQ	0.019	0.00%		Sabinene hydrate	< LOQ	0.019	0.00%	
Terpinolene	< LOQ	0.019	0.00%		trans-β-Ocimene	< LOQ	0.012	0.00%	
<b>Total Terpenes</b>	<b>7.06</b>								



Metals									
Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes	
Arsenic <sup>‡</sup>	< LOQ	0.200	mg/kg	0.0920	2308392	06/20/23 AOAC 2013.06 (mod.) <sup>‡</sup>	pass		
Cadmium <sup>‡</sup>	< LOQ	0.200	mg/kg	0.0920	2308392	06/20/23 AOAC 2013.06 (mod.) <sup>‡</sup>	pass		
Lead <sup>‡</sup>	< LOQ	0.500	mg/kg	0.0920	2308392	06/20/23 AOAC 2013.06 (mod.) <sup>‡</sup>	pass		
Mercury <sup>‡</sup>	< LOQ	0.100	mg/kg	0.0460	2308392	06/20/23 AOAC 2013.06 (mod.) <sup>‡</sup>	pass		



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**Report Number:** 23-006471/D012.R000  
**Report Date:** 06/22/2023  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 05/31/23 16:23

### 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.

<sup>p</sup> = ISO/IEC 17025:2017 accredited method.

<sup>¥</sup> = 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

### Glossary of Qualifiers

I: Insufficient sample received to meet method requirements.

Approved Signatory

Derrick Tanner  
General Manager



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**Report Number:** 23-006471/D012.R000  
**Report Date:** 06/22/2023  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 05/31/23 16:23

Revision: 1 Document ID: 7148  
Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

JAOAC2015 V986 Batch ID: 2307856

Analyte	LCS	Result	Spike	Units	% Rec	Limits	Evaluation	Notes
CBDA	2	0.0283	0.0283	%	99.8	80.0 - 120	Acceptable	
CBV	2	0.0282	0.0291	%	96.8	80.0 - 120	Acceptable	
CEE	2	0.0313	0.0344	%	91.2	80.0 - 120	Acceptable	
CBDA	1	0.0289	0.0311	%	92.9	90.0 - 110	Acceptable	
CBSA	1	0.0305	0.0311	%	98.1	80.0 - 120	Acceptable	
CBS	1	0.0308	0.0322	%	95.7	80.0 - 120	Acceptable	
CB	1	0.0326	0.0323	%	101	90.0 - 110	Acceptable	
THCV	2	0.0193	0.0201	%	95.8	80.0 - 120	Acceptable	
δ8THCV	2	0.0254	0.0268	%	94.4	80.0 - 120	Acceptable	
THCVA	2	0.0295	0.0299	%	98.8	80.0 - 120	Acceptable	
CBN	1	0.0329	0.0329	%	99.9	80.0 - 120	Acceptable	
exo-THC	2	0.0286	0.0292	%	97.8	80.0 - 120	Acceptable	
δ9THC	1	0.0311	0.0341	%	91.0	90.0 - 110	Acceptable	
δ8THC	1	0.0399	0.0420	%	94.9	90.0 - 110	Acceptable	
9SaTHC	1	0.0246	0.0240	%	102	80.0 - 120	Acceptable	
CB	2	0.0318	0.0315	%	101	80.0 - 120	Acceptable	
9RaTHC	1	0.0313	0.0310	%	101	80.0 - 120	Acceptable	
CB	2	0.0315	0.0309	%	102	80.0 - 120	Acceptable	
THCA	1	0.0315	0.0314	%	100	90.0 - 110	Acceptable	
CBA	2	0.0336	0.0326	%	103	80.0 - 120	Acceptable	
CBA	2	0.0334	0.0331	%	101	80.0 - 120	Acceptable	
δ9THCP	2	0.0327	0.0321	%	102	80.0 - 120	Acceptable	
CB	2	0.0332	0.0327	%	102	80.0 - 120	Acceptable	

Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDA	<LOQ	0.00323	%	< 0.00323	Acceptable	
CBV	<LOQ	0.00323	%	< 0.00323	Acceptable	
CEE	<LOQ	0.00323	%	< 0.00323	Acceptable	
CBDA	<LOQ	0.00323	%	< 0.00323	Acceptable	
CBSA	<LOQ	0.00323	%	< 0.00323	Acceptable	
CBS	<LOQ	0.00323	%	< 0.00323	Acceptable	
CB	<LOQ	0.00323	%	< 0.00323	Acceptable	
THCV	<LOQ	0.00323	%	< 0.00323	Acceptable	
δ8THCV	<LOQ	0.00323	%	< 0.00323	Acceptable	
THCVA	<LOQ	0.00323	%	< 0.00323	Acceptable	
CBN	<LOQ	0.00323	%	< 0.00323	Acceptable	
exo-THC	<LOQ	0.00323	%	< 0.00323	Acceptable	
δ9THC	<LOQ	0.00323	%	< 0.00323	Acceptable	
δ8THC	<LOQ	0.00323	%	< 0.00323	Acceptable	
9SaTHC	<LOQ	0.00323	%	< 0.00323	Acceptable	
CB	<LOQ	0.00323	%	< 0.00323	Acceptable	
9RaTHC	<LOQ	0.00323	%	< 0.00323	Acceptable	
CB	<LOQ	0.00323	%	< 0.00323	Acceptable	
THCA	<LOQ	0.00323	%	< 0.00323	Acceptable	
CBA	<LOQ	0.00323	%	< 0.00323	Acceptable	
CBA	<LOQ	0.00323	%	< 0.00323	Acceptable	
δ9THCP	<LOQ	0.00323	%	< 0.00323	Acceptable	
CB	<LOQ	0.00323	%	< 0.00323	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:** 23-006471/D012.R000  
**Report Date:** 06/22/2023  
**ORELAP#:** OR100028  
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Revision: 1 Document ID: 7148  
 Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

JAOAC2015 V986		Batch ID: 2307856						
Sample Duplicate		Sample ID: 23-006038000101						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDA	<LOQ	<LOQ	0.00313	%	NA	< 20	Acceptable	
CBDV	0.0139	0.0142	0.00313	%	2.49	< 20	Acceptable	
CBF	<LOQ	<LOQ	0.00313	%	NA	< 20	Acceptable	
CBDA	<LOQ	<LOQ	0.00313	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.00313	%	NA	< 20	Acceptable	
CBG	0.00782	0.00810	0.00313	%	3.59	< 20	Acceptable	
CB	2.71	2.84	0.00313	%	4.64	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.00313	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.00313	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.00313	%	NA	< 20	Acceptable	
CBN	<LOQ	<LOQ	0.00313	%	NA	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.00313	%	NA	< 20	Acceptable	
d9THC	0.132	0.137	0.00313	%	3.98	< 20	Acceptable	
d8THC	0.0199	0.0206	0.00313	%	3.44	< 20	Acceptable	
9Sa10THC	<LOQ	<LOQ	0.00313	%	NA	< 20	Acceptable	
CB	<LOQ	<LOQ	0.00313	%	NA	< 20	Acceptable	
9Rd10THC	<LOQ	<LOQ	0.00313	%	NA	< 20	Acceptable	
CB	0.0103	0.0108	0.00313	%	4.45	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.00313	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.00313	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.00313	%	NA	< 20	Acceptable	
d9THCP	<LOQ	<LOQ	0.00313	%	NA	< 20	Acceptable	
CB	0.0216	0.0224	0.00313	%	3.46	< 20	Acceptable	

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Laboratory Quality Control Results

JAOAC2015 V986 Batch ID: 2308195

Analyte	LCS	Result	Spike	Units	% Rec	Limits	Evaluation	Notes
CBDA	2	0.0290	0.0283	%	102	80.0 - 120	Acceptable	
CBV	2	0.0301	0.0291	%	103	80.0 - 120	Acceptable	
CBE	2	0.0350	0.0344	%	102	80.0 - 120	Acceptable	
CBDA	1	0.0318	0.0324	%	98.0	90.0 - 110	Acceptable	
CBSA	1	0.0316	0.0328	%	96.3	80.0 - 120	Acceptable	
CBS	1	0.0329	0.0340	%	96.7	80.0 - 120	Acceptable	
CB	1	0.0329	0.0343	%	95.9	90.0 - 110	Acceptable	
THCV	2	0.0210	0.0201	%	104	80.0 - 120	Acceptable	
δ8THCV	2	0.0267	0.0268	%	99.5	80.0 - 120	Acceptable	
THCVA	2	0.0309	0.0299	%	103	80.0 - 120	Acceptable	
CBN	1	0.0337	0.0347	%	96.9	80.0 - 120	Acceptable	
exo-THC	2	0.0293	0.0292	%	100	80.0 - 120	Acceptable	
δ9THC	1	0.0346	0.0351	%	98.5	90.0 - 110	Acceptable	
δ8THC	1	0.0434	0.0428	%	101	90.0 - 110	Acceptable	
9SaTHC	1	0.0245	0.0246	%	99.7	80.0 - 120	Acceptable	
CB	2	0.0324	0.0315	%	103	80.0 - 120	Acceptable	
9RaTHC	1	0.0314	0.0330	%	95.2	80.0 - 120	Acceptable	
CB	2	0.0315	0.0309	%	102	80.0 - 120	Acceptable	
THCA	1	0.0316	0.0332	%	95.3	90.0 - 110	Acceptable	
CBA	2	0.0334	0.0326	%	102	80.0 - 120	Acceptable	
CBA	2	0.0337	0.0331	%	102	80.0 - 120	Acceptable	
δ9THCP	2	0.0326	0.0321	%	102	80.0 - 120	Acceptable	
CB	2	0.0330	0.0327	%	101	80.0 - 120	Acceptable	

Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDA	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBV	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBE	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBDA	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBSA	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBS	<LOQ	0.00327	%	< 0.00327	Acceptable	
CB	<LOQ	0.00327	%	< 0.00327	Acceptable	
THCV	<LOQ	0.00327	%	< 0.00327	Acceptable	
δ8THCV	<LOQ	0.00327	%	< 0.00327	Acceptable	
THCVA	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBN	<LOQ	0.00327	%	< 0.00327	Acceptable	
exo-THC	<LOQ	0.00327	%	< 0.00327	Acceptable	
δ9THC	<LOQ	0.00327	%	< 0.00327	Acceptable	
δ8THC	<LOQ	0.00327	%	< 0.00327	Acceptable	
9SaTHC	<LOQ	0.00327	%	< 0.00327	Acceptable	
CB	<LOQ	0.00327	%	< 0.00327	Acceptable	
9RaTHC	<LOQ	0.00327	%	< 0.00327	Acceptable	
CB	<LOQ	0.00327	%	< 0.00327	Acceptable	
THCA	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBA	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBA	<LOQ	0.00327	%	< 0.00327	Acceptable	
δ9THCP	<LOQ	0.00327	%	< 0.00327	Acceptable	
CB	<LOQ	0.00327	%	< 0.00327	Acceptable	

**Abbreviations**  
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12423 NE Whitaker Way  
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Laboratory Quality Control Results

JAOAC2015 V986		Batch ID: 2308195						
Sample Duplicate		Sample ID: 23-006471000302						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDA	<LOQ	<LOQ	0.00311	%	NA	< 20	Acceptable	
CBDV	0.0896	0.0897	0.00311	%	0.112	< 20	Acceptable	
CBF	1.09	1.10	0.00311	%	0.467	< 20	Acceptable	
CBDA	<LOQ	<LOQ	0.00311	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.00311	%	NA	< 20	Acceptable	
CBG	0.0772	0.0749	0.00311	%	3.03	< 20	Acceptable	
CB	11.7	11.8	0.00311	%	0.471	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.00311	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.00311	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.00311	%	NA	< 20	Acceptable	
CBN	0.0935	0.0938	0.00311	%	0.254	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.00311	%	NA	< 20	Acceptable	
d9THC	0.126	0.128	0.00311	%	1.75	< 20	Acceptable	
d8THC	<LOQ	<LOQ	0.00311	%	NA	< 20	Acceptable	
9Sa10THC	<LOQ	<LOQ	0.00311	%	NA	< 20	Acceptable	
CB	0.0601	0.0635	0.00311	%	5.45	< 20	Acceptable	
9Rd10THC	<LOQ	<LOQ	0.00311	%	NA	< 20	Acceptable	
CB	0.494	0.493	0.00311	%	0.219	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.00311	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.00311	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.00311	%	NA	< 20	Acceptable	
d9THCP	<LOQ	<LOQ	0.00311	%	NA	< 20	Acceptable	
CB	0.379	0.379	0.00311	%	0.0001	< 20	Acceptable	

Abbreviations

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Laboratory Pesticide Quality Control Results

AOAC2007.1 &EN 15662		Units: mg/Kg			Batch ID 2308371			
Method Blank	Blank Result	Blank Limits	Notes	LCS Result	LCS Spke	LCS% Re	Limits	Notes
Abamectin	0.00	< 0.250		0.942	1.000	94.2	50.0	150
Acaphate	0.035	< 0.200		0.666	0.800	83.3	60.0	120
Acquinocyl	0.00	< 1.000		3.539	4.000	90.0	40.0	160
Acetamiprid	0.002	< 0.100		0.371	0.400	92.7	60.0	120
Aldicarb	0.00	< 0.200		0.751	0.800	93.8	60.0	120
Azoxystrobin	0.001	< 0.100		0.367	0.400	91.7	60.0	120
Bifenazate	0.00	< 0.100		0.367	0.400	91.7	60.0	120
Bifenthrin	0.00	< 0.100		0.369	0.400	92.2	50.0	150
Boscalid	0.00	< 0.200		0.769	0.800	96.1	60.0	120
Carbaryl	0.00	< 0.100		0.361	0.400	90.2	60.0	120
Carbifuran	0.00	< 0.100		0.369	0.400	92.3	60.0	120
Chlorantraniliprole	0.00	< 0.100		0.374	0.400	93.4	60.0	120
Chlorfenapyr	0.00	< 0.500		1.787	2.000	89.4	60.0	120
Chlorpyrifos	0.003	< 0.100		0.357	0.400	89.4	60.0	120
Clofentazane	0.00	< 0.100		0.370	0.400	92.6	60.0	120
Cyfluthrin	0.00	< 0.500		1.794	2.000	89.7	50.0	150
Cypermethrin	0.00	< 0.500		1.831	2.000	91.5	50.0	150
Daminozide	0.018	< 0.500		0.640	2.000	32.0	60.0	120
Diazonon	0.00	< 0.100		0.383	0.400	95.6	60.0	120
Dichlorvos	0.00	< 0.500		1.798	2.000	89.9	60.0	120
Dimethoate	0.00	< 0.100		0.389	0.400	97.3	60.0	120
Ethiofoprofos	0.001	< 0.100		0.370	0.400	92.4	60.0	120
Etofenprox	0.00	< 0.200		0.743	0.800	92.9	50.0	150
Etoxazole	0.005	< 0.100		0.382	0.400	95.4	60.0	120
Fenoxycarb	0.00	< 0.100		0.375	0.400	93.8	60.0	120
Fenpyroximate	0.00	< 0.200		0.748	0.800	93.5	60.0	120
Fipronil	0.00	< 0.200		0.747	0.800	93.4	60.0	120
Fonicamid	0.00	< 0.250		0.982	1.000	98.2	60.0	120
Fludioxonil	0.00	< 0.200		0.746	0.800	93.3	50.0	150
Hexythiazox	0.00	< 0.250		0.912	1.000	91.2	60.0	120
Imazalil	0.004	< 0.100		0.370	0.400	92.5	60.0	120
Imidacloprid	0.00	< 0.200		0.701	0.800	87.6	60.0	120
Kiesoxim-methyl	0.00	< 0.200		0.712	0.800	89.0	60.0	120
Malathion	0.001	< 0.100		0.368	0.400	92.0	60.0	120
Metaxyl	0.00	< 0.100		0.371	0.400	92.6	60.0	120
Methiocarb	0.002	< 0.100		0.367	0.400	91.7	60.0	120
Methomyl	0.00	< 0.200		0.769	0.800	96.6	60.0	120
MCK-264	0.00	< 0.100		0.363	0.400	90.8	50.0	150
Mydobutani	0.00	< 0.100		0.359	0.400	89.7	60.0	120
Naled	0.00	< 0.250		0.919	1.000	91.9	50.0	150
Oxamyl	0.00	< 0.500		1.866	2.000	93.3	60.0	120
Padobutrazole	0.002	< 0.200		0.728	0.800	91.0	60.0	120
Parathion-Methyl	0.00	< 0.100		0.330	0.400	82.5	50.0	150
Permethrin	0.00	< 0.100		0.362	0.400	90.4	50.0	150
Phosmet	0.00	< 0.100		0.369	0.400	92.4	50.0	150
Piperonyl butoxide	0.00	< 0.500		1.874	2.000	93.7	60.0	120
Prallethrin	0.00	< 0.100		0.372	0.400	93.1	60.0	120
Propiconazole	0.00	< 0.200		0.740	0.800	92.4	60.0	120
Propoxur	0.004	< 0.100		0.373	0.400	93.3	60.0	120
Pyrethrin (Summe)	0.00	< 0.100		0.453	0.488	92.8	60.0	120
Pyridaben	0.005	< 0.100		0.377	0.400	94.4	50.0	150
Spinosad	0.00	< 0.100		0.347	0.388	89.6	50.0	150
Spiromesfen	0.00	< 0.100		0.371	0.400	92.7	60.0	120
Spirotetramat	0.00	< 0.100		0.373	0.400	93.3	60.0	120
Spiroxamine	0.008	< 0.200		0.755	0.800	94.4	60.0	120
Tebuconazole	0.00	< 0.200		0.728	0.800	91.1	60.0	120
Thiadoprid	0.00	< 0.100		0.373	0.400	93.3	60.0	120
Thiamethoxam	0.00	< 0.100		0.391	0.400	97.7	60.0	120
Trifloxystrobin	0.002	< 0.100		0.369	0.400	92.2	60.0	120

Q6



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Laboratory Pesticide Quality Control Results

AOAC2007.1 & EN 15662		Units: mg/Kg				Batch ID 2308371				
Matrix Spke/Matrix Spke Duplicate Recoveries	Result	MS Res	MSD Res	Spike	RFD%	Limit	MS% Re	MSD % Re	Limits	Notes
Abamectin	0.000	0.987	0.987	1.000	0.0%	< 30	98.7%	98.7%	50 - 150	
Acephate	0.031	0.707	0.708	0.800	0.3%	< 30	84.8%	84.7%	50 - 150	
Acetamiprid	0.096	3.769	3.766	4.000	0.1%	< 30	91.8%	91.7%	50 - 150	
Acetamiprid	0.002	0.385	0.380	0.400	1.4%	< 30	95.7%	94.3%	50 - 150	
Aldicarb	0.000	0.778	0.772	0.800	0.8%	< 30	97.3%	96.5%	50 - 150	
Azoxystrobin	0.001	0.377	0.378	0.400	0.4%	< 30	94.0%	94.4%	50 - 150	
Bifenazate	0.000	0.385	0.379	0.400	1.4%	< 30	96.2%	94.8%	50 - 150	
Bifenthrin	0.000	0.383	0.386	0.400	0.8%	< 30	95.7%	96.4%	50 - 150	
Boscalid	0.000	0.757	0.745	0.800	1.6%	< 30	94.8%	93.1%	50 - 150	
Carbaryl	0.000	0.383	0.380	0.400	0.8%	< 30	95.7%	94.9%	50 - 150	
Carbofuran	0.000	0.383	0.382	0.400	0.2%	< 30	95.8%	95.8%	50 - 150	
Chlorantraniliprole	0.000	0.402	0.382	0.400	4.9%	< 30	100.4%	95.8%	50 - 150	
Chlorfenapyr	0.000	1.824	1.871	2.000	2.6%	< 30	91.2%	93.8%	50 - 150	
Chlorpyrifos	0.003	0.378	0.374	0.400	1.3%	< 30	93.9%	92.7%	50 - 150	
Clofentezine	0.000	0.352	0.344	0.400	2.4%	< 30	88.1%	86.0%	50 - 150	
Cyfluthrin	0.000	1.886	1.808	2.000	4.2%	< 30	94.3%	90.4%	30 - 150	
Cypermethrin	0.000	1.963	1.941	2.000	1.1%	< 30	98.1%	97.1%	50 - 150	
Daminozide	0.005	0.648	0.660	2.000	1.8%	< 30	32.2%	32.8%	30 - 150	
Diazinon	0.000	0.405	0.399	0.400	1.5%	< 30	101.2%	99.7%	50 - 150	
Dichlorvos	0.022	1.978	1.977	2.000	0.1%	< 30	97.8%	97.8%	50 - 150	
Dimethoate	0.000	0.389	0.393	0.400	1.2%	< 30	97.2%	98.4%	50 - 150	
Ethionphos	0.001	0.383	0.377	0.400	1.4%	< 30	95.8%	94.1%	50 - 150	
Etofenprox	0.000	0.771	0.759	0.800	1.5%	< 30	96.4%	94.9%	50 - 150	
Etoxazole	0.004	0.411	0.410	0.400	0.1%	< 30	101.6%	101.4%	50 - 150	
Fenoxycarb	0.000	0.384	0.387	0.400	0.7%	< 30	96.0%	96.8%	50 - 150	
Fenpyroximate	0.000	0.804	0.794	0.800	1.3%	< 30	100.5%	99.3%	50 - 150	
Fipronil	0.000	0.778	0.778	0.800	0.0%	< 30	97.3%	97.3%	50 - 150	
Fonicamid	0.000	1.005	0.949	1.000	5.8%	< 30	100.5%	94.9%	50 - 150	
Fludioxonil	0.000	0.751	0.762	0.800	1.4%	< 30	93.9%	95.2%	50 - 150	
Hexythiazox	0.010	0.943	0.930	1.000	1.4%	< 30	93.3%	92.0%	50 - 150	
Imazalil	0.004	0.390	0.376	0.400	3.6%	< 30	96.8%	93.1%	50 - 150	
Imidacloprid	0.000	0.732	0.724	0.800	1.1%	< 30	91.5%	90.5%	50 - 150	
Kiesoxim-methyl	0.000	0.776	0.741	0.800	4.6%	< 30	97.1%	92.7%	50 - 150	
Malathion	0.001	0.384	0.378	0.400	1.4%	< 30	95.8%	94.4%	50 - 150	
Metolaxyl	0.000	0.389	0.388	0.400	0.2%	< 30	97.3%	97.1%	50 - 150	
Methiocarb	0.002	0.387	0.381	0.400	1.6%	< 30	96.3%	94.8%	50 - 150	
Methomyl	0.000	0.786	0.795	0.800	1.1%	< 30	98.3%	99.3%	50 - 150	
MCK-264	0.000	0.375	0.383	0.400	2.0%	< 30	93.9%	95.8%	50 - 150	
Mydobutani	0.000	0.392	0.388	0.400	1.2%	< 30	98.1%	96.9%	50 - 150	
Naled	0.000	0.970	0.950	1.000	2.1%	< 30	97.0%	95.0%	50 - 150	
Oxaryl	0.000	1.939	1.767	2.000	9.3%	< 30	97.0%	88.3%	50 - 150	
Padobutrazole	0.000	0.790	0.762	0.800	3.6%	< 30	98.8%	95.2%	50 - 150	
Parathion-Methyl	0.000	0.400	0.364	0.400	9.3%	< 30	100.0%	91.1%	30 - 150	
Permethrin	0.000	0.386	0.367	0.400	4.8%	< 30	96.4%	91.9%	50 - 150	
Phosmet	0.000	0.388	0.373	0.400	3.8%	< 30	96.9%	93.3%	50 - 150	
Piperonyl butoxide	0.000	1.929	1.918	2.000	0.6%	< 30	96.8%	95.9%	50 - 150	
Prallethrin	0.000	0.380	0.387	0.400	1.8%	< 30	95.0%	96.7%	50 - 150	
Propiconazole	0.000	0.776	0.765	0.800	1.5%	< 30	97.0%	95.8%	50 - 150	
Propoxur	0.004	0.388	0.385	0.400	0.8%	< 30	96.0%	95.3%	50 - 150	
Pyrethrin (Summe)	0.009	0.376	0.369	0.488	2.1%	< 30	75.2%	73.8%	50 - 150	
Pyridaben	0.005	0.372	0.373	0.400	0.1%	< 30	91.9%	92.0%	50 - 150	
Spirosad	0.000	0.359	0.362	0.388	0.8%	< 30	92.8%	93.4%	50 - 150	
Spiromesfen	0.000	0.390	0.384	0.400	1.6%	< 30	97.8%	96.0%	50 - 150	
Spirotetramat	0.000	0.382	0.378	0.400	1.1%	< 30	95.8%	94.8%	50 - 150	
Spiroxamine	0.000	0.791	0.781	0.800	1.3%	< 30	98.8%	97.8%	50 - 150	
Tebuconazole	0.000	0.779	0.769	0.800	1.2%	< 30	97.4%	96.2%	50 - 150	
Thiadoprid	0.000	0.392	0.383	0.400	2.1%	< 30	97.9%	95.8%	50 - 150	
Thiamethoxam	0.000	0.392	0.373	0.400	4.9%	< 30	98.0%	93.3%	50 - 150	
Trifloxystrobin	0.002	0.387	0.387	0.400	0.2%	< 30	96.4%	96.2%	50 - 150	



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 503-254-1794

Report Number: 23-006471/D012.R000  
 Report Date: 06/22/2023  
 ORELAP#: OR100028  
 Purchase Order:  
 Received: 05/31/23 16:23



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

Laboratory Quality Control Results

Residual Solvents				Batch ID: 2308405					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		538	584	µg/g	92.1	60 - 120	
Isobutane	ND	< 200		656	767	µg/g	85.5	60 - 120	
Butane	ND	< 200		677	782	µg/g	86.6	60 - 120	
2,2-Dimethylpropane	ND	< 200		807	939	µg/g	85.9	60 - 120	
Methanol	ND	< 200		1390	1640	µg/g	84.8	60 - 120	
Ethylene Oxide	ND	< 30		51.8	57.1	µg/g	90.7	60 - 120	
2-Methylbutane	ND	< 200		1360	1600	µg/g	85.0	60 - 120	
Pentane	ND	< 200		1400	1620	µg/g	86.4	60 - 120	
Ethanol	ND	< 200		1410	1610	µg/g	87.6	70 - 130	
Ethyl Ether	ND	< 200		1440	1610	µg/g	89.4	60 - 120	
2,2-Dimethylbutane	ND	< 30		144	168	µg/g	85.7	60 - 120	
Acetone	ND	< 200		1410	1620	µg/g	87.0	60 - 120	
2-Propanol	ND	< 200		1420	1600	µg/g	88.8	60 - 120	
Ethyl Formate	ND	< 500		1200	1600	µg/g	75.0	70 - 130	
Acetonitrile	ND	< 100		414	484	µg/g	85.5	60 - 120	
Methyl Acetate	ND	< 500		1440	1610	µg/g	89.4	70 - 130	
2,3-Dimethylbutane	ND	< 30		150	162	µg/g	92.6	60 - 120	
Dichloromethane	ND	< 60		449	483	µg/g	93.0	60 - 120	
2-Methylpentane	ND	< 30		136	174	µg/g	78.2	60 - 120	
MTBE	ND	< 500		1490	1610	µg/g	92.5	70 - 130	
3-Methylpentane	ND	< 30		151	168	µg/g	89.9	60 - 120	
Hexane	ND	< 30		146	168	µg/g	86.9	60 - 120	
1-Propanol	ND	< 500		1410	1600	µg/g	88.1	70 - 130	
Methylethylketone	ND	< 500		1480	1620	µg/g	91.4	70 - 130	
Ethyl acetate	ND	< 200		1430	1600	µg/g	89.4	60 - 120	
2-Butanol	ND	< 200		1400	1600	µg/g	87.5	60 - 120	
Tetrahydrofuran	ND	< 100		453	514	µg/g	88.1	60 - 120	
Cyclohexane	ND	< 200		1460	1600	µg/g	91.3	60 - 120	
2-methyl-1-propanol	ND	< 500		1460	1610	µg/g	90.7	70 - 130	
Benzene	ND	< 1		3.85	5.12	µg/g	75.2	60 - 120	
Isopropyl Acetate	ND	< 200		1440	1620	µg/g	88.9	60 - 120	
Heptane	ND	< 200		1450	1610	µg/g	90.1	60 - 120	
1-Butanol	ND	< 500		1390	1600	µg/g	86.9	70 - 130	
Propyl Acetate	ND	< 500		1200	1600	µg/g	75.0	70 - 130	
1,4-Dioxane	ND	< 100		400	493	µg/g	81.1	60 - 120	
2-Ethoxyethanol	ND	< 30		137	163	µg/g	84.0	60 - 120	
Methylisobutylketone	ND	< 500		1450	1600	µg/g	90.6	70 - 130	
3-Methyl-1-butanol	ND	< 500		1490	1610	µg/g	92.5	70 - 130	
Ethylene Glycol	ND	< 200		527	483	µg/g	109.1	60 - 120	
Toluene	ND	< 100		431	493	µg/g	87.4	60 - 120	
Isobutyl Acetate	ND	< 500		1400	1600	µg/g	87.5	70 - 130	
1-Pentanol	ND	< 500		1380	1600	µg/g	86.3	70 - 130	
Butyl Acetate	ND	< 500		1360	1600	µg/g	85.0	70 - 130	
Ethylbenzene	ND	< 200		803	969	µg/g	82.9	60 - 120	
m,p-Xylene	ND	< 200		748	968	µg/g	77.3	60 - 120	
o-Xylene	ND	< 200		815	976	µg/g	83.5	60 - 120	
Cumene	ND	< 30		129	162	µg/g	79.6	60 - 120	
Anisole	ND	< 500		1340	1610	µg/g	83.2	70 - 130	
DMSO	ND	< 500		1270	1610	µg/g	78.9	70 - 130	
1,2-dimethoxyethane	ND	< 50		144	164	µg/g	87.8	70 - 130	
Triethylamine	ND	< 500		1480	1600	µg/g	92.5	70 - 130	
N,N-dimethylformamide	ND	< 150		430	484	µg/g	88.8	70 - 130	
N,N-dimethylacetamide	ND	< 150		459	489	µg/g	93.9	70 - 130	
Pyridine	ND	< 50		143	172	µg/g	83.1	70 - 130	
Sulfolane	ND	< 50		75.2	163	µg/g	46.1	70 - 130	Q6
1,2-Dichloroethane	ND	< 1		0.883	1	µg/g	88.3	70 - 130	
Chloroform	ND	< 1		0.916	1	µg/g	91.6	70 - 130	
Trichloroethylene	ND	< 1		0.893	1	µg/g	89.3	70 - 130	
1,1-Dichloroethane	ND	< 1		0.933	1	µg/g	93.3	70 - 130	



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**Report Number:** 23-006471/D012.R000  
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**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 05/31/23 16:23

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

QC- Sample Duplicate Sample ID: 23-006471-0004

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



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**Report Number:** 23-006471/D012.R000  
**Report Date:** 06/22/2023  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 05/31/23 16:23

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

Terpenes Quality Control Results

Method Reference: EPA5035				Batch ID: 2308421					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	LCS	Units	LCS% Rec	Limits	Notes
a-pinene	<LOQ	< 200		407	500	µg/g	81%	70 - 130	
Camphene	<LOQ	< 200		416	500	µg/g	83%	70 - 130	
Sabinene	<LOQ	< 200		404	500	µg/g	81%	70 - 130	
b-Pinene	<LOQ	< 200		403	500	µg/g	81%	70 - 130	
b-Myrcene	<LOQ	< 200		398	500	µg/g	80%	70 - 130	
a-phellandrene	<LOQ	< 200		419	500	µg/g	84%	70 - 130	
d-3-Carene	<LOQ	< 200		407	500	µg/g	81%	70 - 130	
a-Terpinene	<LOQ	< 200		412	500	µg/g	82%	70 - 130	
p-Cymene	<LOQ	< 200		401	500	µg/g	80%	70 - 130	
D-Limonene	<LOQ	< 200		409	500	µg/g	82%	70 - 130	
Eucalyptol	<LOQ	< 200		409	500	µg/g	82%	70 - 130	
b-cis-Cimene	<LOQ	< 67		132	167	µg/g	79%	70 - 130	
b-trans-Cimene	<LOQ	< 133		270	333	µg/g	81%	70 - 130	
g-Terpinene	<LOQ	< 200		409	500	µg/g	82%	70 - 130	
Sabinene Hydrate	<LOQ	< 200		415	500	µg/g	83%	70 - 130	
Terpinolene	<LOQ	< 200		402	500	µg/g	80%	70 - 130	
D-Fenchone	<LOQ	< 200		409	500	µg/g	82%	70 - 130	
Linalool	<LOQ	< 200		412	500	µg/g	82%	70 - 130	
Fenchol	<LOQ	< 200		409	500	µg/g	82%	70 - 130	
Camphor	<LOQ	< 200		406	500	µg/g	81%	70 - 130	
Isopulego	<LOQ	< 200		416	500	µg/g	83%	70 - 130	
Isoborneol	<LOQ	< 200		420	500	µg/g	84%	70 - 130	
Borneol	<LOQ	< 200		423	500	µg/g	85%	70 - 130	
DL-Menthol	<LOQ	< 200		393	500	µg/g	79%	70 - 130	
Terpineol	<LOQ	< 200		402	500	µg/g	80%	70 - 130	
Nerd	<LOQ	< 200		350	500	µg/g	70%	70 - 130	
Pulegone	<LOQ	< 200		384	500	µg/g	77%	70 - 130	
Geraniol	<LOQ	< 200		384	500	µg/g	77%	70 - 130	
Geranyl Acetate	<LOQ	< 200		365	500	µg/g	73%	70 - 130	
a-Cedrene	<LOQ	< 200		402	500	µg/g	80%	70 - 130	
b-Caryophyllene	<LOQ	< 200		415	500	µg/g	83%	70 - 130	
a-Humulene	<LOQ	< 200		419	500	µg/g	84%	70 - 130	
Valene	<LOQ	< 200		399	500	µg/g	80%	70 - 130	
cis-Nerolidol	<LOQ	< 200		405	500	µg/g	81%	70 - 130	
a-Farnesene	<LOQ	< 200		419	500	µg/g	84%	70 - 130	
trans-Nerolidol	<LOQ	< 200		426	500	µg/g	85%	70 - 130	
Caryophyllene Oxide	<LOQ	< 200		423	500	µg/g	85%	70 - 130	
Guaiol	<LOQ	< 200		419	500	µg/g	84%	70 - 130	
Cedrol	<LOQ	< 200		415	500	µg/g	83%	70 - 130	
a-Bisabolol	<LOQ	< 200		406	500	µg/g	81%	70 - 130	

Definitions

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





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**Report Number:** 23-006471/D012.R000  
**Report Date:** 06/22/2023  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 05/31/23 16:23

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

Terpenes Quality Control Results

Method Reference: EPA5035		Batch ID: 2308421					
Sample/ Sample Duplicate		Sample ID: 23-006471-0004					
Analyte	Result	Org. Result	LOQ	Units	% RPD	LIMIT	Notes
a-pinene	27800	27600	193	µg/g	1%	< 20	
Camphene	397	395	193	µg/g	1%	< 20	
Sabinene	<LOQ	<LOQ	193	µg/g	0%	< 20	
b-Pinene	<LOQ	<LOQ	193	µg/g	0%	< 20	
b-Myrcene	431	429	193	µg/g	0%	< 20	
a-phellandrene	<LOQ	<LOQ	193	µg/g	0%	< 20	
d-3-Carene	<LOQ	<LOQ	193	µg/g	0%	< 20	
a-Terpinene	<LOQ	<LOQ	193	µg/g	0%	< 20	
p-Cymene	<LOQ	<LOQ	193	µg/g	0%	< 20	
D-Limonene	32000	31800	193	µg/g	1%	< 20	
Eucalyptol	<LOQ	<LOQ	193	µg/g	0%	< 20	
b-cis-Cimene	<LOQ	<LOQ	642	µg/g	0%	< 20	
b-trans-Cimene	<LOQ	<LOQ	128	µg/g	0%	< 20	
g-Terpinene	<LOQ	<LOQ	193	µg/g	0%	< 20	
Sabinene Hydrate	<LOQ	<LOQ	193	µg/g	0%	< 20	
Terpinolene	<LOQ	<LOQ	193	µg/g	0%	< 20	
D-Fenchone	<LOQ	<LOQ	193	µg/g	0%	< 20	
Linalool	<LOQ	<LOQ	193	µg/g	0%	< 20	
Fenchol	<LOQ	<LOQ	193	µg/g	0%	< 20	
Camphor	<LOQ	<LOQ	193	µg/g	0%	< 20	
Isopulego	<LOQ	<LOQ	193	µg/g	0%	< 20	
Isoborneol	<LOQ	<LOQ	193	µg/g	0%	< 20	
Borneol	<LOQ	<LOQ	193	µg/g	0%	< 20	
DL-Menthhol	<LOQ	<LOQ	193	µg/g	0%	< 20	
Terpineol	<LOQ	<LOQ	193	µg/g	0%	< 20	
Nerd	<LOQ	<LOQ	193	µg/g	0%	< 20	
Pulegone	<LOQ	<LOQ	193	µg/g	0%	< 20	
Geraniol	<LOQ	<LOQ	193	µg/g	0%	< 20	
Geranyl_Acetate	<LOQ	<LOQ	193	µg/g	0%	< 20	
a-Cedrene	<LOQ	<LOQ	193	µg/g	0%	< 20	
b-Caryophyllene	7090	7040	193	µg/g	1%	< 20	
a-Humulene	829	827	193	µg/g	0%	< 20	
Valnene	<LOQ	<LOQ	193	µg/g	0%	< 20	
cis-Nerolidol	<LOQ	<LOQ	193	µg/g	0%	< 20	
a-Farnesene	<LOQ	<LOQ	193	µg/g	0%	< 20	
trans-Nerolidol	<LOQ	<LOQ	193	µg/g	0%	< 20	
Caryophyllene_Oxide	1390	1380	193	µg/g	1%	< 20	
Guaiol	<LOQ	<LOQ	193	µg/g	0%	< 20	
Cedrol	<LOQ	<LOQ	193	µg/g	0%	< 20	
a-Bisabolol	1170	1120	193	µg/g	4%	< 20	

Definitions  
 RPD Relative Percent Difference



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**Report Number:** 23-006471/D012.R000  
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**Received:** 05/31/23 16:23





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.



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**Report Number:** 23-011802/D002.R001  
**Report Date:** 10/19/2023  
**ORELAP#:** OR100028  
**Purchase Order:** 2663180  
**Received:** 10/04/23 16:16

Potency per 1g					
Method: J AOAC 2015 V98-6 (mod) <sup>b</sup>					
Analyte	Result	Limits	Units	Batch: 2311963	Analyze: 10/18/23 12:12:00 A
THCV-Total per 1g	< LOQ		mg/1g		0.0574
Total Cannabinoids per 1g	311		mg/1g		

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



**Report Number:** 23-011802/D002.R001  
**Report Date:** 10/19/2023  
**ORELAP#:** OR100028  
**Purchase Order:** 2663180  
**Received:** 10/04/23 16:16

**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.

<sup>p</sup> = ISO/IEC 17025:2017 accredited method.

**Units of Measure**

g = g

mg/1g = Milligram per 1g

% = Percentage of sample

% wt =  $\mu\text{g/g}$  divided by 10,000

Approved Signatory

Derrick Tanner  
General Manager



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



**Report Number:** 23-011802/D002.R001  
**Report Date:** 10/19/2023  
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**Purchase Order:** 2663180  
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Revision: 4 Document ID: 7148  
 Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

J AOAC 2015 V98-6 Batch ID: 2311963

Laboratory Control Sample									
Analyte	LCS	Result	Spike	Units	% Rec	Limits		Evaluation	Notes
CBDVA	2	0.0307	0.0333	%	92.2	80.0	- 120	Acceptable	
CBDV	2	0.0309	0.0324	%	95.4	80.0	- 120	Acceptable	
CBE	2	0.0329	0.0355	%	92.7	80.0	- 120	Acceptable	
CBDA	1	0.0310	0.0322	%	96.0	90.0	- 110	Acceptable	
CBGA	1	0.0314	0.0329	%	95.4	80.0	- 120	Acceptable	
CBG	1	0.0355	0.0368	%	96.7	80.0	- 120	Acceptable	
CBD	1	0.0308	0.0313	%	98.2	90.0	- 110	Acceptable	
THCV	2	0.0280	0.0304	%	91.9	80.0	- 120	Acceptable	
d8THCV	2	0.0291	0.0305	%	95.4	80.0	- 120	Acceptable	
THCVA	2	0.0305	0.0327	%	93.3	80.0	- 120	Acceptable	
CBN	1	0.0316	0.0329	%	96.2	80.0	- 120	Acceptable	
exo-THC	2	0.0298	0.0327	%	91.2	80.0	- 120	Acceptable	
d9THC	1	0.0345	0.0365	%	94.6	90.0	- 110	Acceptable	
d8THC	1	0.0331	0.0340	%	97.4	90.0	- 110	Acceptable	
9S-d10THC	1	0.0326	0.0337	%	96.6	80.0	- 120	Acceptable	
CBL	2	0.0319	0.0337	%	94.6	80.0	- 120	Acceptable	
9R-d10THC	1	0.0318	0.0336	%	94.6	80.0	- 120	Acceptable	
CBC	2	0.0311	0.0338	%	91.8	80.0	- 120	Acceptable	
THCA	1	0.0308	0.0337	%	91.5	90.0	- 110	Acceptable	
CBCA	2	0.0295	0.0333	%	88.6	80.0	- 120	Acceptable	
CBLA	2	0.0315	0.0349	%	90.4	80.0	- 120	Acceptable	
d9THCP	2	0.0293	0.0333	%	88.0	80.0	- 120	Acceptable	
CBT	2	0.0288	0.0322	%	89.4	80.0	- 120	Acceptable	

Method Blank						
Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.00331	%	< 0.00331	Acceptable	
CBDV	<LOQ	0.00331	%	< 0.00331	Acceptable	
CBE	<LOQ	0.00331	%	< 0.00331	Acceptable	
CBDA	<LOQ	0.00331	%	< 0.00331	Acceptable	
CBGA	<LOQ	0.00331	%	< 0.00331	Acceptable	
CBG	<LOQ	0.00331	%	< 0.00331	Acceptable	
CBD	<LOQ	0.00331	%	< 0.00331	Acceptable	
THCV	<LOQ	0.00331	%	< 0.00331	Acceptable	
d8THCV	<LOQ	0.00331	%	< 0.00331	Acceptable	
THCVA	<LOQ	0.00331	%	< 0.00331	Acceptable	
CBN	<LOQ	0.00331	%	< 0.00331	Acceptable	
exo-THC	<LOQ	0.00331	%	< 0.00331	Acceptable	
d9THC	<LOQ	0.00331	%	< 0.00331	Acceptable	
d8THC	<LOQ	0.00331	%	< 0.00331	Acceptable	
9S-d10THC	<LOQ	0.00331	%	< 0.00331	Acceptable	
CBL	<LOQ	0.00331	%	< 0.00331	Acceptable	
9R-d10THC	<LOQ	0.00331	%	< 0.00331	Acceptable	
CBC	<LOQ	0.00331	%	< 0.00331	Acceptable	
THCA	<LOQ	0.00331	%	< 0.00331	Acceptable	
CBCA	<LOQ	0.00331	%	< 0.00331	Acceptable	
CBLA	<LOQ	0.00331	%	< 0.00331	Acceptable	
d9THCP	<LOQ	0.00331	%	< 0.00331	Acceptable	
CBT	<LOQ	0.00331	%	< 0.00331	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 Quality Control Results

J AOAC 2015 V98-6		Batch ID: 2311963						
Sample Duplicate		Sample ID: 23-012215-0001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBDV	0.0360	0.0367	0.00330	%	1.83	< 20	Acceptable	
CBE	0.0584	0.0581	0.00330	%	0.501	< 20	Acceptable	
CBD	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBDA	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBG	0.0429	0.0422	0.00330	%	1.74	< 20	Acceptable	
CBD	3.78	3.84	0.00330	%	1.54	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBN	0.0241	0.0242	0.00330	%	0.349	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
d9THC	0.135	0.135	0.00330	%	0.235	< 20	Acceptable	
d8THC	0.0311	0.0309	0.00330	%	0.505	< 20	Acceptable	
9S-d10THC	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
9R-d10THC	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBC	0.245	0.245	0.00330	%	0.0285	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBCA	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
d9THCP	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBT	0.0673	0.0474	0.00330	%	34.6	< 20	Outlier	R, Q6

Abbreviations

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

Units of Measure:

% - Percent



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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.