

## 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 KG	REGULATORY ACTION LEVEL
Arsenic	mg/serving	mg/kg	12.5 mg/kg <sup>[1]</sup>
Cadmium	mg/serving	mg/kg	10 mg/kg <sup>[1]</sup>
Lead	mg/serving	mg/kg	10 mg/kg <sup>[1]</sup>
Mercury	mg/serving	mg/kg	0.267 mg/kg <sup>[1]</sup>
PESTICIDES	REGULATORY ACTION LEVEL		
None of the other 59 pesticides tested found above limit of detection in the sample.			10 ppb <sup>[2]</sup>
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		
TERPENES	% OF SAMPLE		
Farnesene	%		
β-Caryophyllene	%		
α-Bisabolol	%		
Guaiol	%		
Humulene	%		
Caryophyllene Oxide	%		



1. Deemy, M. Benjamin, L. FDA. (2019). *CVM CY15-17 Report on Heavy Metals in Animal Food*. Retrieved from: <https://www.fda.gov/media/132046/download>

2. 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:** 24-000030/D008.R000  
**Report Date:** 01/10/2024  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 01/02/24 16:18

**Customer:** Etz Hayim Holdings  
**Product identity:** FORM-TN.FS.SAL20-FL62  
**Client/Metric ID:** .  
**Laboratory ID:** 24-000030-0004

### Summary

#### Potency:

Analyte per 1g	Result	Limits	Units	Status	
CBC per 1g	0.851		mg/1g		CBD-Total per Serving Size 23.0 mg/1g
CBD per 1g	23.0		mg/1g		
CBDV per 1g	0.127		mg/1g		THC-Total per Serving Size 0.596 mg/1g
CBE per 1g	0.434		mg/1g		(Reported in milligrams per serving)
CBG per 1g	0.398		mg/1g		
CBN per 1g	0.0768		mg/1g		
CBT per 1g	0.707		mg/1g		
Δ9-THC per 1g	0.596		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	0.0314	38.01%	a-Bisabolol	0.0258	31.23%
(-)-caryophyllene oxide	0.0254	30.75%	<b>Total Terpenes</b>	<b>0.0826</b>	<b>100.00%</b>

#### Metals:

Analyte	Result	Units	Limit	Status
Arsenic*	0.656	mg/kg		

#### Microbiology:

Less than LOQ for all analytes.



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

**Product identity:** FORM-TN.FS.SAL20-FL62

**Client/Metric ID:** .

**Sample Date:**

**Laboratory ID:** 24-000030-0004

**Evidence of Cooling:** No

**Temp:** 18.8 °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: 2400093	Analyze: 1/3/24 11:13:00 PM	
Analyte	Result	Limits	Units	LOQ	Notes
CBC per 1g	0.851		mg/1g	0.0329	
CBC-A per 1g	< LOQ		mg/1g	0.0329	
CBC-Total per 1g	0.851		mg/1g	0.0617	
CBD per 1g	23.0		mg/1g	0.329	
CBD-A per 1g	< LOQ		mg/1g	0.0329	
CBD-Total per 1g	23.0		mg/1g	0.358	
CBDV per 1g	0.127		mg/1g	0.0329	
CBDV-A per 1g	< LOQ		mg/1g	0.0329	
CBDV-Total per 1g	0.127		mg/1g	0.0614	
CBE per 1g	0.434		mg/1g	0.0329	
CBG per 1g	0.398		mg/1g	0.0329	
CBG-A per 1g	< LOQ		mg/1g	0.0329	
CBG-Total per 1g	0.398		mg/1g	0.0614	
CBL per 1g	< LOQ		mg/1g	0.0329	
CBL-A per 1g	< LOQ		mg/1g	0.0329	
CBL-Total per 1g	< LOQ		mg/1g	0.0617	
CBN per 1g	0.0768		mg/1g	0.0329	
CBT per 1g	0.707		mg/1g	0.0329	
Δ8-THCV per 1g	< LOQ		mg/1g	0.0329	
Δ10-THC-9R per 1g	< LOQ		mg/1g	0.0329	
Δ10-THC-9S per 1g	< LOQ		mg/1g	0.0329	
Δ10-THC-Total per 1g	< LOQ		mg/1g	0.0658	
Δ8-THC per 1g	< LOQ		mg/1g	0.0329	
Δ9-THC per 1g	0.596		mg/1g	0.0329	
delta-9-THCP per 1g	< LOQ		mg/1g	0.0329	
exo-THC per 1g	< LOQ		mg/1g	0.0329	
THC-A per 1g	< LOQ		mg/1g	0.0329	
THC-Total per 1g	0.596		mg/1g	0.0617	
THCV per 1g	< LOQ		mg/1g	0.0329	
THCV-A per 1g	< LOQ		mg/1g	0.0329	



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Potency per 1g	Method: J AOAC 2015 V98-6 (mod) <sup>P</sup>	Units mg/se	Batch: 2400093	Analyze: 1/3/24 11:13:00 PM	
Analyte	Result	Limits	Units	LOQ	Notes
THCV-Total per 1g	< LOQ		mg/1g	0.0618	
Total Cannabinoids per 1g	26.2		mg/1g		

### Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2400134	01/07/24 AOAC 991.14 (Petrifilm) <sup>P</sup>		
Total Coliforms	< LOQ		cfu/g	10	2400134	01/07/24 AOAC 991.14 (Petrifilm) <sup>P</sup>		
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2400135	01/08/24 AOAC 2014.05 (RAPID) <sup>P</sup>		
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2400135	01/08/24 AOAC 2014.05 (RAPID) <sup>P</sup>		

### Solvents

Method: Residual Solvents by GC/MS <sup>P</sup>						Units µg/g	Batch 2400188	Analyze 01/08/24 02:07 PM			
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							



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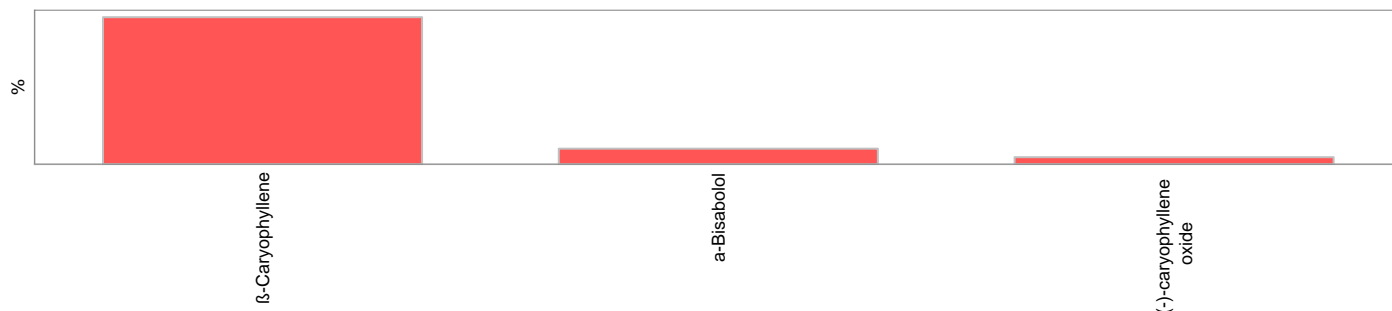


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Pesticides											
Method: AOAC 2007.01 & EN 15662 (mod) <sup>b</sup>											
Units mg/kg Batch 2400180 Analyze 01/08/24 10:57 AM											
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
Abamectin <sup>‡</sup>	< 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	
Bifentazate <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		Etoazole <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	
Imazali <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 2400206	Analyze 01/08/24 07:24 PM		
Analyte	Result	LOQ	% of Total	Notes	Analyte	Result	LOQ	% of Total	Notes
β-Caryophyllene	0.0314	0.019	38.0145%		α-Bisabolol	0.0258	0.019	31.2349%	
(-)-caryophyllene oxide	0.0254	0.019	30.7506%		Humulene	< LOQ	0.019	0.00%	
(-)-Guaiol	< LOQ	0.019	0.00%		Geraniol	< LOQ	0.019	0.00%	
Linalool	< LOQ	0.019	0.00%		Terpinolene	< LOQ	0.019	0.00%	
(±)-fenchone	< LOQ	0.019	0.00%		(+)-fenchol	< LOQ	0.019	0.00%	
(±)-cis-Nerolidol	< LOQ	0.019	0.00%		(±)-trans-Nerolidol	< LOQ	0.019	0.00%	
α-Terpinene	< LOQ	0.019	0.00%		(-)-β-Pinene	< LOQ	0.019	0.00%	
Menthol	< LOQ	0.019	0.00%		valencene	< LOQ	0.019	0.00%	
(+)-Borneol	< LOQ	0.019	0.00%		β-Myrcene	< LOQ	0.019	0.00%	
Geranyl acetate	< LOQ	0.019	0.00%		(+)-Cedrol	< LOQ	0.019	0.00%	
(-)-α-Terpineol	< LOQ	0.019	0.00%		(-)-Isopulegol	< LOQ	0.019	0.00%	
(+)-Pulegone	< LOQ	0.019	0.00%		(±)-Camphor	< LOQ	0.019	0.00%	
(R)-(+)-Limonene	< 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%		Eucalyptol	< LOQ	0.019	0.00%	
farnesene	< LOQ	0.019	0.00%		γ-Terpinene	< LOQ	0.019	0.00%	
Isoborneol	< LOQ	0.019	0.00%		nerol	< LOQ	0.019	0.00%	
p-Cymene	< LOQ	0.019	0.00%		Sabinene	< LOQ	0.019	0.00%	
Sabinene hydrate	< LOQ	0.019	0.00%		trans-β-Ocimene	< LOQ	0.012	0.00%	
<b>Total Terpenes</b>	<b>0.0826</b>								



Metals							
Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status Notes
Arsenic <sup>‡</sup>	0.656		mg/kg	0.0898	2400223	01/08/24 AOAC 2013.06 (mod.) <sup>‡</sup>	
Cadmium <sup>‡</sup>	< LOQ		mg/kg	0.0898	2400223	01/08/24 AOAC 2013.06 (mod.) <sup>‡</sup>	
Lead <sup>‡</sup>	< LOQ		mg/kg	0.0898	2400223	01/08/24 AOAC 2013.06 (mod.) <sup>‡</sup>	
Mercury <sup>‡</sup>	< LOQ		mg/kg	0.0449	2400223	01/08/24 AOAC 2013.06 (mod.) <sup>‡</sup>	



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

Laboratory Quality Control Results

JAOAC2015 V986 Batch ID: 2400093

Laboratory Control Sample									
Analyte	LCS	Result	Spike	Units	% Rec	Limits		Evaluation	Notes
CBDA	2	0.0318	0.0329	%	96.7	80.0	- 120	Acceptable	
CBV	2	0.0314	0.0322	%	97.5	80.0	- 120	Acceptable	
CBE	2	0.0335	0.0350	%	95.6	80.0	- 120	Acceptable	
CBDA	1	0.0302	0.0317	%	95.2	90.0	- 110	Acceptable	
CBSA	1	0.0301	0.0315	%	95.5	80.0	- 120	Acceptable	
CBS	1	0.0294	0.0309	%	95.2	80.0	- 120	Acceptable	
CB	1	0.0317	0.0330	%	95.9	90.0	- 110	Acceptable	
THCV	2	0.0338	0.0352	%	96.1	80.0	- 120	Acceptable	
deltaTHCV	2	0.0296	0.0307	%	96.2	80.0	- 120	Acceptable	
THCVA	2	0.0308	0.0320	%	96.3	80.0	- 120	Acceptable	
CBN	1	0.0321	0.0330	%	97.1	80.0	- 120	Acceptable	
exo-THC	2	0.0299	0.0313	%	95.5	80.0	- 120	Acceptable	
deltaTHC	1	0.0326	0.0337	%	96.8	90.0	- 110	Acceptable	
deltaTHC	1	0.0328	0.0336	%	97.6	90.0	- 110	Acceptable	
9SdeltaTHC	1	0.0315	0.0326	%	96.9	80.0	- 120	Acceptable	
CB	2	0.0321	0.0326	%	98.5	80.0	- 120	Acceptable	
9SHHC	3	0.0307	0.0316	%	97.2	80.0	- 120	Acceptable	
9RdeltaTHC	1	0.0305	0.0318	%	95.9	80.0	- 120	Acceptable	
CB	2	0.0317	0.0328	%	96.5	80.0	- 120	Acceptable	
9RHHC	3	0.0283	0.0293	%	96.7	80.0	- 120	Acceptable	
THCA	1	0.0308	0.0322	%	95.6	90.0	- 110	Acceptable	
CBSA	2	0.0322	0.0334	%	96.5	80.0	- 120	Acceptable	
CLA	2	0.0327	0.0343	%	95.1	80.0	- 120	Acceptable	
deltaTHCP	2	0.0316	0.0328	%	96.1	80.0	- 120	Acceptable	
deltaTHCO	3	0.0335	0.0336	%	99.7	80.0	- 120	Acceptable	
CB	2	0.0326	0.0341	%	95.6	80.0	- 120	Acceptable	
deltaTHCO	3	0.0342	0.0329	%	104	80.0	- 120	Acceptable	

Method Blank

Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDA	<LOQ	0.00310	%	< 0.00310	Acceptable	
CBV	<LOQ	0.00310	%	< 0.00310	Acceptable	
CBE	<LOQ	0.00310	%	< 0.00310	Acceptable	
CBDA	<LOQ	0.00310	%	< 0.00310	Acceptable	
CBSA	<LOQ	0.00310	%	< 0.00310	Acceptable	
CBS	<LOQ	0.00310	%	< 0.00310	Acceptable	
CB	<LOQ	0.00310	%	< 0.00310	Acceptable	
THCV	<LOQ	0.00310	%	< 0.00310	Acceptable	
deltaTHCV	<LOQ	0.00310	%	< 0.00310	Acceptable	
THCVA	<LOQ	0.00310	%	< 0.00310	Acceptable	
CBN	<LOQ	0.00310	%	< 0.00310	Acceptable	
exo-THC	<LOQ	0.00310	%	< 0.00310	Acceptable	
deltaTHC	<LOQ	0.00310	%	< 0.00310	Acceptable	
deltaTHC	<LOQ	0.00310	%	< 0.00310	Acceptable	
9SdeltaTHC	<LOQ	0.00310	%	< 0.00310	Acceptable	
CB	<LOQ	0.00310	%	< 0.00310	Acceptable	
9SHHC	<LOQ	0.00310	%	< 0.00310	Acceptable	
9RdeltaTHC	<LOQ	0.00310	%	< 0.00310	Acceptable	
CB	<LOQ	0.00310	%	< 0.00310	Acceptable	
9RHHC	<LOQ	0.00310	%	< 0.00310	Acceptable	
THCA	<LOQ	0.00310	%	< 0.00310	Acceptable	
CBSA	<LOQ	0.00310	%	< 0.00310	Acceptable	
CLA	<LOQ	0.00310	%	< 0.00310	Acceptable	
deltaTHCP	<LOQ	0.00310	%	< 0.00310	Acceptable	
deltaTHCO	<LOQ	0.00310	%	< 0.00310	Acceptable	
CB	<LOQ	0.00310	%	< 0.00310	Acceptable	
deltaTHCO	<LOQ	0.00310	%	< 0.00310	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:** 24-000030/D008.R000  
**Report Date:** 01/10/2024  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 01/02/24 16:18

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

Laboratory Quality Control Results

JAOAC2015 V986		Batch ID: 2400093						
Sample Duplicate		Sample ID: 23-0151260001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDA	<LOQ	<LOQ	0.00314	%	NA	< 20	Acceptable	
CBDV	0.132	0.131	0.00314	%	0.314	< 20	Acceptable	
CBE	0.0640	0.0631	0.00314	%	1.31	< 20	Acceptable	
CBDA	0.393	0.391	0.00314	%	0.569	< 20	Acceptable	
CBDA	0.0351	0.0352	0.00314	%	0.174	< 20	Acceptable	
CBG	0.177	0.177	0.00314	%	0.0429	< 20	Acceptable	
CBG	9.36	9.27	0.00314	%	0.924	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.00314	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.00314	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.00314	%	NA	< 20	Acceptable	
CBN	0.0306	0.0306	0.00314	%	0.241	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.00314	%	NA	< 20	Acceptable	
d9THC	0.754	0.752	0.00314	%	0.303	< 20	Acceptable	
d8THC	0.00564	0.00583	0.00314	%	3.27	< 20	Acceptable	
9Sd10THC	<LOQ	<LOQ	0.00314	%	NA	< 20	Acceptable	
CB	<LOQ	<LOQ	0.00314	%	NA	< 20	Acceptable	
9SHHC	<LOQ	<LOQ	0.00314	%	NA	< 20	Acceptable	
9Rd10THC	<LOQ	<LOQ	0.00314	%	NA	< 20	Acceptable	
CB	0.305	0.304	0.00314	%	0.317	< 20	Acceptable	
9RHHC	<LOQ	<LOQ	0.00314	%	NA	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.00314	%	NA	< 20	Acceptable	
CBDA	0.0366	0.0363	0.00314	%	0.716	< 20	Acceptable	
CHA	<LOQ	<LOQ	0.00314	%	NA	< 20	Acceptable	
d9THCP	<LOQ	<LOQ	0.00314	%	NA	< 20	Acceptable	
d8THCO	<LOQ	<LOQ	0.00314	%	NA	< 20	Acceptable	
CB	0.148	0.148	0.00314	%	0.0812	< 20	Acceptable	
d9THCO	<LOQ	<LOQ	0.00314	%	NA	< 20	Acceptable	

Abbreviations

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**Purchase Order:**  
**Received:** 01/02/24 16:18



Revision: 3 Document ID: 3120  
 LegacyID: CFLC21WorksheetValidated 10/30/2020

Laboratory Pesticide Quality Control Results

AOAC2007.1 &EN 15662		Units: mg/Kg			Batch ID 2400180			
Method Blank	Blank Result	Blank Limits	Notes	LCS Result	LCS Spk	LCS % Re	Limits	Notes
Abamectin	0.000	< 0.250		1.002	1.000	100.2	50.0	150
Acaphate	0.115	< 0.200		0.782	0.800	97.8	60.0	120
Acetaminophen	0.454	< 1.000		4.289	4.000	107.2	40.0	160
Acetamiprid	0.000	< 0.100		0.386	0.400	91.4	60.0	120
Aldicarb	0.000	< 0.200		0.754	0.800	94.3	60.0	120
Azoxystrobin	0.000	< 0.100		0.337	0.400	84.1	60.0	120
Bifenazate	0.000	< 0.100		0.372	0.400	92.9	60.0	120
Bifenthrin	0.000	< 0.100		0.389	0.400	97.4	50.0	150
Boscalid	0.000	< 0.200		0.772	0.800	96.5	60.0	120
Carbaryl	0.000	< 0.100		0.375	0.400	93.7	60.0	120
Carbendazim	0.000	< 0.100		0.390	0.400	97.4	60.0	120
Chlorantraniliprole	0.000	< 0.100		0.378	0.400	94.5	60.0	120
Chlorfenapyr	0.000	< 0.500		1.668	2.000	83.4	60.0	120
Chlorpyrifos	0.000	< 0.100		0.367	0.400	91.8	60.0	120
Clofentezine	0.000	< 0.100		0.358	0.400	89.6	60.0	120
Cyfluthrin	0.000	< 0.500		1.907	2.000	95.4	50.0	150
Cypermethrin	0.000	< 0.500		1.904	2.000	95.2	50.0	150
Daminozide	0.000	< 0.500		0.660	2.000	33.0	60.0	120
Diazinon	0.000	< 0.100		0.364	0.400	91.0	60.0	120
Dichlorvos	0.000	< 0.500		1.753	2.000	87.7	60.0	120
Dimethoate	0.000	< 0.100		0.375	0.400	93.8	60.0	120
Ethionphos	0.000	< 0.100		0.353	0.400	88.3	60.0	120
Etofenprox	0.015	< 0.200		0.795	0.800	99.4	50.0	150
Etoxazole	0.000	< 0.100		0.376	0.400	94.0	60.0	120
Fenoxycarb	0.000	< 0.100		0.368	0.400	91.9	60.0	120
Fenpyroximate	0.000	< 0.200		0.758	0.800	94.7	60.0	120
Fipronil	0.000	< 0.200		0.768	0.800	96.0	60.0	120
Fonicamid	0.000	< 0.250		1.012	1.000	101.2	60.0	120
Fludioxonil	0.000	< 0.200		0.702	0.800	87.7	50.0	150
Hexythiazox	0.000	< 0.250		0.944	1.000	94.4	60.0	120
Imazalil	0.000	< 0.100		0.359	0.400	89.8	60.0	120
Imidacloprid	0.000	< 0.200		0.723	0.800	90.3	60.0	120
Kiesoxim-methyl	0.000	< 0.200		0.733	0.800	91.7	60.0	120
Malathion	0.000	< 0.100		0.398	0.400	99.4	60.0	120
Metaxyl	0.000	< 0.100		0.377	0.400	94.2	60.0	120
Methiocarb	0.000	< 0.100		0.377	0.400	94.2	60.0	120
Methomyl	0.000	< 0.200		0.839	0.800	104.9	60.0	120
MCK-264	0.000	< 0.100		0.349	0.400	87.2	50.0	150
Mydobutanol	0.000	< 0.100		0.362	0.400	90.5	60.0	120
Naled	0.000	< 0.250		0.912	1.000	91.2	50.0	150
Oxaryl	0.000	< 0.500		2.011	2.000	100.6	60.0	120
Padobutrazole	0.000	< 0.200		0.782	0.800	97.8	60.0	120
Parathion-Methyl	0.000	< 0.100		0.393	0.400	98.3	50.0	150
Permethrin	0.015	< 0.100		0.397	0.400	99.2	50.0	150
Phosmet	0.000	< 0.100		0.362	0.400	90.6	50.0	150
Piperonyl butoxide	0.000	< 0.500		1.932	2.000	96.6	60.0	120
Prallethrin	0.000	< 0.100		0.343	0.400	85.8	60.0	120
Propiconazole	0.000	< 0.200		0.724	0.800	90.5	60.0	120
Propoxur	0.000	< 0.100		0.388	0.400	97.0	60.0	120
Pyrethrin (Summe)	0.000	< 0.100		0.448	0.488	91.9	60.0	120
Pyridaben	0.000	< 0.100		0.377	0.400	94.3	50.0	150
Spirosad	0.000	< 0.100		0.363	0.388	93.5	50.0	150
Spiromesfen	0.000	< 0.100		0.357	0.400	89.2	60.0	120
Spirotetramat	0.000	< 0.100		0.359	0.400	89.9	60.0	120
Spiroxamine	0.000	< 0.200		0.746	0.800	93.2	60.0	120
Tebuconazole	0.000	< 0.200		0.759	0.800	94.9	60.0	120
Thiadoprid	0.000	< 0.100		0.372	0.400	92.9	60.0	120
Thiamethoxam	0.000	< 0.100		0.399	0.400	99.7	60.0	120
Trifloxystrobin	0.000	< 0.100		0.351	0.400	87.9	60.0	120

Q7



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**Purchase Order:**  
**Received:** 01/02/24 16:18

Revision: 3 Document ID: 3120  
 LegacyID: CFLC21WorksheetValidated 10/30/2020

Laboratory Pesticide Quality Control Results

AOAC2007.1 & EN 15662		Units: mg/Kg					Batch ID 2400180			
Matrix Spke/Matrix Spke Duplicate Recoveries	Result	MS Res	MSD Res	Spike	RFD%	Limit	MS% Re	MSD% Re	Limits	Notes
Abamectin	0.00	0.909	0.947	1.00	4.0%	< 30	90.9%	94.7%	50 - 150	
Acephate	0.134	0.791	0.746	0.800	7.0%	< 30	82.1%	76.5%	50 - 150	
Acetaminophen	0.00	3.405	4.455	4.00	26.7%	< 30	85.1%	111.4%	50 - 150	
Acetamiprid	0.00	0.376	0.372	0.400	1.0%	< 30	94.4%	93.1%	50 - 150	
Aldicarb	0.00	0.716	0.721	0.800	0.7%	< 30	89.5%	90.2%	50 - 150	
Azoxystrobin	0.00	0.371	0.348	0.400	6.6%	< 30	92.8%	86.9%	50 - 150	
Bifenazate	0.00	0.355	0.346	0.400	2.4%	< 30	88.8%	86.0%	50 - 150	
Bifenthrin	0.00	0.348	0.402	0.400	14.5%	< 30	87.0%	100.6%	50 - 150	
Boscalid	0.00	0.720	0.730	0.800	1.3%	< 30	90.0%	91.2%	50 - 150	
Carbaryl	0.00	0.357	0.357	0.400	0.1%	< 30	89.3%	89.3%	50 - 150	
Carbofuran	0.00	0.357	0.363	0.400	1.7%	< 30	89.2%	90.7%	50 - 150	
Chlorantraniliprole	0.00	0.335	0.341	0.400	1.7%	< 30	83.9%	85.3%	50 - 150	
Chlorfenapyr	0.00	1.818	1.567	2.00	14.8%	< 30	90.9%	78.4%	50 - 150	
Chlorpyrifos	0.00	0.322	0.329	0.400	2.1%	< 30	80.6%	82.4%	50 - 150	
Clofentezine	0.00	0.297	0.290	0.400	2.4%	< 30	74.3%	72.3%	50 - 150	
Cyfluthrin	0.00	1.988	2.084	2.00	5.7%	< 30	98.4%	104.2%	30 - 150	
Cypermethrin	0.00	2.105	2.132	2.00	1.3%	< 30	105.3%	106.6%	50 - 150	
Daminozide	0.00	0.646	0.635	2.00	1.7%	< 30	32.3%	31.8%	30 - 150	
Diazinon	0.00	0.332	0.335	0.400	0.9%	< 30	82.9%	83.0%	50 - 150	
Dichlorvos	0.00	1.805	1.858	2.00	2.9%	< 30	90.2%	92.9%	50 - 150	
Dimethoate	0.00	0.362	0.359	0.400	0.8%	< 30	90.5%	89.7%	50 - 150	
Ethionphos	0.00	0.353	0.346	0.400	2.1%	< 30	88.3%	86.5%	50 - 150	
Etofenprox	0.00	0.680	0.779	0.800	13.8%	< 30	85.0%	97.4%	50 - 150	
Etoxazole	0.00	0.351	0.316	0.400	10.7%	< 30	87.9%	79.0%	50 - 150	
Fenoxycarb	0.00	0.331	0.342	0.400	3.5%	< 30	82.7%	85.6%	50 - 150	
Fenpyroximate	0.00	0.717	0.741	0.800	3.2%	< 30	89.7%	92.8%	50 - 150	
Fipronil	0.00	0.695	0.741	0.800	6.3%	< 30	86.9%	92.8%	50 - 150	
Fonicamid	0.00	1.005	1.004	1.00	0.3%	< 30	100.6%	100.4%	50 - 150	
Fludioxonil	0.00	0.570	0.605	0.800	6.1%	< 30	71.2%	75.7%	50 - 150	
Hexythiazox	0.00	0.874	0.878	1.00	0.4%	< 30	87.4%	87.8%	50 - 150	
Imazalil	0.00	0.343	0.344	0.400	0.3%	< 30	85.7%	86.0%	50 - 150	
Imidacloprid	0.00	0.684	0.648	0.800	5.5%	< 30	85.5%	80.9%	50 - 150	
Kiesoxim-methyl	0.00	0.684	0.657	0.800	4.0%	< 30	85.5%	82.2%	50 - 150	
Malathion	0.00	0.344	0.347	0.400	0.8%	< 30	86.0%	86.7%	50 - 150	
Metolaxyl	0.00	0.360	0.350	0.400	2.6%	< 30	89.9%	87.8%	50 - 150	
Methiocarb	0.00	0.352	0.337	0.400	4.4%	< 30	87.9%	84.2%	50 - 150	
Methomyl	0.00	0.833	0.802	0.800	3.8%	< 30	104.1%	100.2%	50 - 150	
MCK-264	0.00	0.323	0.335	0.400	3.5%	< 30	80.9%	83.7%	50 - 150	
Mydobutani	0.00	0.354	0.338	0.400	4.7%	< 30	88.4%	84.4%	50 - 150	
Naled	0.00	0.872	0.851	1.00	2.5%	< 30	87.2%	85.1%	50 - 150	
Oxaryl	0.00	1.969	1.893	2.00	3.9%	< 30	98.5%	94.7%	50 - 150	
Padobutrazole	0.00	0.771	0.772	0.800	0.1%	< 30	96.4%	96.5%	50 - 150	
Parathion-Methyl	0.02	0.359	0.418	0.400	16.0%	< 30	84.8%	99.5%	30 - 150	
Permethrin	0.03	0.346	0.384	0.400	11.0%	< 30	83.2%	92.9%	50 - 150	
Phosmet	0.00	0.333	0.345	0.400	3.4%	< 30	83.3%	86.1%	50 - 150	
Piperonyl butoxide	0.00	1.765	1.767	2.00	0.1%	< 30	88.3%	88.4%	50 - 150	
Prallethrin	0.00	0.326	0.327	0.400	0.4%	< 30	81.5%	81.8%	50 - 150	
Propiconazole	0.00	0.667	0.655	0.800	4.1%	< 30	83.4%	86.9%	50 - 150	
Propoxur	0.00	0.346	0.361	0.400	4.2%	< 30	86.5%	90.2%	50 - 150	
Pyrethrin (Summe)	0.00	0.433	0.437	0.488	0.8%	< 30	88.8%	89.5%	50 - 150	
Pyridaben	0.00	0.400	0.404	0.400	0.9%	< 30	100.1%	101.0%	50 - 150	
Spirosad	0.00	0.329	0.335	0.388	1.8%	< 30	84.7%	86.3%	50 - 150	
Spiromesfen	0.00	0.343	0.347	0.400	1.0%	< 30	85.9%	86.7%	50 - 150	
Spirotetramat	0.00	0.327	0.328	0.400	0.2%	< 30	81.8%	81.9%	50 - 150	
Spiroxamine	0.00	0.704	0.705	0.800	0.2%	< 30	88.0%	88.2%	50 - 150	
Tebuconazole	0.00	0.707	0.727	0.800	2.9%	< 30	88.3%	90.9%	50 - 150	
Thiadoprid	0.00	0.357	0.352	0.400	1.5%	< 30	89.3%	88.0%	50 - 150	
Thiamethoxam	0.00	0.395	0.354	0.400	11.1%	< 30	98.8%	88.4%	50 - 150	
Trifloxystrobin	0.00	0.328	0.336	0.400	2.4%	< 30	82.0%	84.0%	50 - 150	



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

Laboratory Quality Control Results

Residual Solvents				Batch ID: 2400188					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		531	584	µg/g	90.9	60 - 120	
Isobutane	ND	< 200		593	767	µg/g	77.3	60 - 120	
Butane	ND	< 200		587	782	µg/g	75.1	60 - 120	
2,2-Dimethylpropane	ND	< 200		842	939	µg/g	89.7	60 - 120	
Methanol	ND	< 200		1420	1600	µg/g	88.8	60 - 120	
Ethylene Oxide	ND	< 30		46.6	57.1	µg/g	81.6	60 - 120	
2-Methylbutane	ND	< 200		1390	1600	µg/g	86.9	60 - 120	
Pentane	ND	< 200		1360	1600	µg/g	85.0	60 - 120	
Ethanol	ND	< 200		1300	1600	µg/g	81.3	70 - 130	
Ethyl Ether	ND	< 200		1370	1600	µg/g	85.6	60 - 120	
2,2-Dimethylbutane	ND	< 30		138	161	µg/g	85.7	60 - 120	
Acetone	ND	< 200		1440	1600	µg/g	90.0	60 - 120	
2-Propanol	ND	< 200		1410	1600	µg/g	88.1	60 - 120	
Ethyl Formate	ND	< 500		1140	1600	µg/g	71.3	70 - 130	
Acetonitrile	ND	< 100		430	488	µg/g	88.1	60 - 120	
Methyl Acetate	ND	< 500		1410	1610	µg/g	87.6	70 - 130	
2,3-Dimethylbutane	ND	< 30		140	163	µg/g	85.9	60 - 120	
Dichloromethane	ND	< 60		407	488	µg/g	83.4	60 - 120	
2-Methylpentane	ND	< 30		132	161	µg/g	82.0	60 - 120	
MTBE	ND	< 500		1530	1650	µg/g	92.7	70 - 130	
3-Methylpentane	ND	< 30		134	162	µg/g	82.7	60 - 120	
Hexane	ND	< 30		118	161	µg/g	73.3	60 - 120	
1-Propanol	ND	< 500		1450	1620	µg/g	89.5	70 - 130	
Methyl ethyl ketone	ND	< 500		1450	1610	µg/g	90.1	70 - 130	
Ethyl acetate	ND	< 200		1420	1610	µg/g	88.2	60 - 120	
2-Butanol	ND	< 200		1400	1610	µg/g	87.0	60 - 120	
Tetrahydrofuran	ND	< 100		409	483	µg/g	84.7	60 - 120	
Cyclohexane	ND	< 200		1340	1600	µg/g	83.8	60 - 120	
2-methyl-1-propanol	ND	< 500		1290	1600	µg/g	80.6	70 - 130	
Benzene	ND	< 1		2.66	4.99	µg/g	53.3	60 - 120	Q6
Isopropyl Acetate	ND	< 200		1440	1600	µg/g	90.0	60 - 120	
Heptane	ND	< 200		1410	1600	µg/g	88.1	60 - 120	
1-Butanol	ND	< 500		1340	1610	µg/g	83.2	70 - 130	
Propyl Acetate	ND	< 500		1490	1610	µg/g	92.5	70 - 130	
1,4-Dioxane	ND	< 100		379	480	µg/g	79.0	60 - 120	
2-Ethoxyethanol	ND	< 30		113	161	µg/g	70.2	60 - 120	
Methylisobutylketone	ND	< 500		1440	1610	µg/g	89.4	70 - 130	
3-Methyl-1-butanol	ND	< 500		1330	1610	µg/g	82.6	70 - 130	
Ethylene Glycol	ND	< 200		288	481	µg/g	59.9	60 - 120	Q6
Toluene	ND	< 100		410	483	µg/g	84.9	60 - 120	
Isobutyl Acetate	ND	< 500		1490	1610	µg/g	92.5	70 - 130	
1-Pentanol	ND	< 500		1320	1610	µg/g	82.0	70 - 130	
Butyl Acetate	ND	< 500		1420	1600	µg/g	88.8	70 - 130	
Ethylbenzene	ND	< 200		850	962	µg/g	88.4	60 - 120	
m,p-Xylene	ND	< 200		850	972	µg/g	87.4	60 - 120	
o-Xylene	ND	< 200		834	965	µg/g	86.4	60 - 120	
Cumene	ND	< 30		133	169	µg/g	78.7	60 - 120	
Anisole	ND	< 500		1310	1600	µg/g	81.9	70 - 130	
DMSO	ND	< 500		973	1600	µg/g	60.8	70 - 130	Q6
1,2-dimethoxyethane	ND	< 50		151	163	µg/g	92.6	70 - 130	
Triethylamine	ND	< 500		1140	1600	µg/g	71.3	70 - 130	
N,N-dimethylformamide	ND	< 150		368	482	µg/g	76.3	70 - 130	
N,N-dimethylacetamide	ND	< 150		397	483	µg/g	82.2	70 - 130	
Pyridine	ND	< 50		125	161	µg/g	77.6	70 - 130	
Silfolane	ND	< 50		77.8	163	µg/g	47.7	70 - 130	Q6
1,2-Dichloroethane	ND	< 1		0.835	1	µg/g	83.5	70 - 130	
Chloroform	ND	< 1		0.728	1	µg/g	72.8	70 - 130	
Trichloroethylene	ND	< 1		0.77	1	µg/g	77.0	70 - 130	
1,1,1-Trichloroethane	ND	< 1		0.855	1	µg/g	85.5	70 - 130	



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

QC- Sample Duplicate		Sample ID: 23-015123-0001						
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		
Methylethylketone	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

**Units of Measure:**

µg/g - Microgram per gram or ppm



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**Report Number:** 24-000030/D008.R000  
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 Legacy ID: CFL-E57Worksheet Validated 11/04/2020

Terpenes Quality Control Results

Method Reference: EPA5035				Batch ID: 2400206					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	LCS	Units	LCS% Rec	Limits	Notes
a-pinene	<LOQ	< 200		429	500	µg/g	86%	70 - 130	
Camphene	<LOQ	< 200		440	500	µg/g	88%	70 - 130	
Sabinene	<LOQ	< 200		429	500	µg/g	86%	70 - 130	
b-Pinene	<LOQ	< 200		429	500	µg/g	86%	70 - 130	
b-Myrcene	<LOQ	< 200		436	500	µg/g	87%	70 - 130	
a-phellandrene	<LOQ	< 200		448	500	µg/g	90%	70 - 130	
d-3-Carene	<LOQ	< 200		450	500	µg/g	90%	70 - 130	
a-Terpinene	<LOQ	< 200		443	500	µg/g	89%	70 - 130	
p-Cymene	<LOQ	< 200		439	500	µg/g	88%	70 - 130	
D-Limonene	<LOQ	< 200		433	500	µg/g	87%	70 - 130	
Eucalyptol	<LOQ	< 200		430	500	µg/g	86%	70 - 130	
b-cis-Cimene	<LOQ	< 67		149	167	µg/g	89%	70 - 130	
b-trans-Cimene	<LOQ	< 133		294	333	µg/g	88%	70 - 130	
g-Terpinene	<LOQ	< 200		447	500	µg/g	89%	70 - 130	
Sabinene Hydrate	<LOQ	< 200		447	500	µg/g	89%	70 - 130	
Terpinolene	<LOQ	< 200		439	500	µg/g	88%	70 - 130	
D-Fenchone	<LOQ	< 200		432	500	µg/g	86%	70 - 130	
Linalool	<LOQ	< 200		458	500	µg/g	92%	70 - 130	
Fenchol	<LOQ	< 200		450	500	µg/g	90%	70 - 130	
Camphor	<LOQ	< 200		439	500	µg/g	88%	70 - 130	
Isopulego	<LOQ	< 200		483	500	µg/g	97%	70 - 130	
Isoborneol	<LOQ	< 200		465	500	µg/g	93%	70 - 130	
Borneol	<LOQ	< 200		452	500	µg/g	90%	70 - 130	
DL-Menthol	<LOQ	< 200		455	500	µg/g	91%	70 - 130	
Terpineol	<LOQ	< 200		464	500	µg/g	93%	70 - 130	
Nerd	<LOQ	< 200		449	500	µg/g	90%	70 - 130	
Pulegone	<LOQ	< 200		468	500	µg/g	94%	70 - 130	
Geraniol	<LOQ	< 200		453	500	µg/g	91%	70 - 130	
Geranyl_Acdate	<LOQ	< 200		462	500	µg/g	92%	70 - 130	
a-Cedrene	<LOQ	< 200		470	500	µg/g	94%	70 - 130	
b-Caryophyllene	<LOQ	< 200		484	500	µg/g	97%	70 - 130	
a-Humulene	<LOQ	< 200		468	500	µg/g	94%	70 - 130	
Valene	<LOQ	< 200		471	500	µg/g	94%	70 - 130	
cis-Nerolidol	<LOQ	< 200		483	500	µg/g	97%	70 - 130	
a-Farnesene	<LOQ	< 200		510	500	µg/g	102%	70 - 130	
trans-Nerolidol	<LOQ	< 200		477	500	µg/g	95%	70 - 130	
Caryophyllene_Oxide	<LOQ	< 200		471	500	µg/g	94%	70 - 130	
Guaiol	<LOQ	< 200		484	500	µg/g	97%	70 - 130	
Cedrol	<LOQ	< 200		473	500	µg/g	95%	70 - 130	
a-Bisabolol	<LOQ	< 200		474	500	µg/g	95%	70 - 130	

Definitions

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



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Legacy ID: CFL-E57Worksheet Validated 11/04/2020

Terpenes Quality Control Results

Method Reference: EPA5035		Batch ID: 2400206					
Sample/ Sample Duplicate		Sample ID: 24-000030-004					
Analyte	Result	Org. Result	LOQ	Units	% RPD	LIMIT	Notes
a-pinene	<LOQ	<LOQ	191	µg/g	0%	< 20	
Camphene	<LOQ	<LOQ	191	µg/g	0%	< 20	
Sabinene	<LOQ	<LOQ	191	µg/g	0%	< 20	
b-Pinene	<LOQ	<LOQ	191	µg/g	0%	< 20	
b-Myrcene	<LOQ	<LOQ	191	µg/g	0%	< 20	
a-phellandrene	<LOQ	<LOQ	191	µg/g	0%	< 20	
d-3-Carene	<LOQ	<LOQ	191	µg/g	0%	< 20	
a-Terpinene	<LOQ	<LOQ	191	µg/g	0%	< 20	
p-Cymene	<LOQ	<LOQ	191	µg/g	0%	< 20	
D-Limonene	<LOQ	<LOQ	191	µg/g	0%	< 20	
Eucalyptol	<LOQ	<LOQ	191	µg/g	0%	< 20	
b-cis-Cimene	<LOQ	<LOQ	63.6	µg/g	0%	< 20	
b-trans-Cimene	<LOQ	<LOQ	127	µg/g	0%	< 20	
g-Terpinene	<LOQ	<LOQ	191	µg/g	0%	< 20	
Sabinene Hydrate	<LOQ	<LOQ	191	µg/g	0%	< 20	
Terpinolene	<LOQ	<LOQ	191	µg/g	0%	< 20	
D-Fenchone	<LOQ	<LOQ	191	µg/g	0%	< 20	
Linalool	<LOQ	<LOQ	191	µg/g	0%	< 20	
Fenchol	<LOQ	<LOQ	191	µg/g	0%	< 20	
Camphor	<LOQ	<LOQ	191	µg/g	0%	< 20	
Isopulego	<LOQ	<LOQ	191	µg/g	0%	< 20	
Isoborneol	<LOQ	<LOQ	191	µg/g	0%	< 20	
Borneol	<LOQ	<LOQ	191	µg/g	0%	< 20	
DL-Menthhol	<LOQ	<LOQ	191	µg/g	0%	< 20	
Terpineol	<LOQ	<LOQ	191	µg/g	0%	< 20	
Nerd	<LOQ	<LOQ	191	µg/g	0%	< 20	
Pulegone	<LOQ	<LOQ	191	µg/g	0%	< 20	
Geraniol	<LOQ	<LOQ	191	µg/g	0%	< 20	
Geranyl_Acetate	<LOQ	<LOQ	191	µg/g	0%	< 20	
a-Cedrene	<LOQ	<LOQ	191	µg/g	0%	< 20	
b-Caryophyllene	319	314	191	µg/g	2%	< 20	
a-Humulene	<LOQ	<LOQ	191	µg/g	0%	< 20	
Valnene	<LOQ	<LOQ	191	µg/g	0%	< 20	
cis-Nerolidol	<LOQ	<LOQ	191	µg/g	0%	< 20	
a-Farnesene	<LOQ	<LOQ	191	µg/g	0%	< 20	
trans-Nerolidol	<LOQ	<LOQ	191	µg/g	0%	< 20	
Caryophyllene_Oxide	256	254	191	µg/g	1%	< 20	
Guaiol	<LOQ	<LOQ	191	µg/g	0%	< 20	
Cedrol	<LOQ	<LOQ	191	µg/g	0%	< 20	
a-Bisabolol	259	258	191	µg/g	0%	< 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.