

## CONSOLIDATED TEST RESULTS SUMMARY

Please see the following pages for full test results.

<b>BULK SKU</b> MO.O.FS8	<b>BATCH #</b> EC19	<b>LOQ:</b> Limit Of Quantitation
<b>PRODUCT NAME</b> CBD Massage Oil	<b>SERVING SIZE</b> 3 tsp	<b>LOD:</b> Limit Of Detection
<b>LABORATORY:</b> Columbia Laboratories	<b>OREGON ACCREDITATION:</b> OR100028	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)	131.19 mg/serving	8.870 mg/g	0.887 %
Total THC (d9-THC, THCA)	6.17 mg/serving	0.417 mg/g	0.042 %
Cannabigerol (CBG)	4.33 mg/serving	0.293 mg/g	0.029 %
Cannabinol (CBN)	<LOQ mg/serving	<LOQ mg/g	<LOQ %
Cannabichromene (CBC)	4.70 mg/serving	0.318 mg/g	0.032 %
Tetrahydrocannabinolic Acid (THCA)	<LOQ mg/serving	<LOQ mg/g	<LOQ %
Delta-9-THC (d9-THC)	6.17 mg/serving	0.417 mg/g	0.042 %
Delta-8-THC (d8-THC)	<LOQ mg/serving	<LOQ mg/g	<LOQ %

  

HEAVY METALS	PER SERVING	PER GRAM	REGULATORY ACTION LEVEL
Arsenic	<LOQ µg/serving	<LOQ µg/g	10 µg/day <sup>[1]</sup>
Cadmium	<LOQ µg/serving	<LOQ µg/g	4.1 µg/day <sup>[1]</sup>
Lead	<LOQ µg/serving	<LOQ µg/g	3.5 µg/day <sup>[2]</sup>
Mercury	<LOQ µg/serving	<LOQ µg/g	2 µg/day <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>[1]</sup>

  

RESIDUAL SOLVENTS	Results	REGULATORY ACTION LEVEL
Ethanol	<LOQ	50,000 mg/day
Heptane	<LOQ	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



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

2. US Food and Drug Administration. (2019). Lead in Food, Foodwares, and Dietary Supplements. Washington DC: FDA. US Food and Drug Administration. (2019). Lead in Food, Foodwares, and Dietary Supplements. Washington DC: FDA.



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



**Report Number:** 22-003105/D002.R000  
**Report Date:** 03/24/2022  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 03/17/22 14:20

**Customer:** Etz Hayim Holdings  
**Product identity:** EC19-MO.O.FS8  
**Client/Metric ID:** .  
**Laboratory ID:** 22-003105-0001

### Summary

**Potency:**

Analyte per 1g	Result	Limits	Units	Status	
CBC per 1g <sup>†</sup>	0.318		mg/1g		CBD Total per 1g 8.87 mg/1g
CBD per 1g	8.87		mg/1g		
CBE per 1g <sup>†</sup>	0.0889		mg/1g		THC-Total per 1g 0.417 mg/1g
CBG per 1g <sup>†</sup>	0.293		mg/1g		
CBT per 1g <sup>†</sup>	0.124		mg/1g		
Δ9 THC per 1g	0.417		mg/1g		(Reported in milligrams per serving)



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**Purchase Order:**  
**Received:** 03/17/22 14:20

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

**Product identity:** EC19-MO.O.FS8

**Client/Metric ID:** .

**Sample Date:**

**Laboratory ID:** 22-003105-0001

**Evidence of Cooling:** No

**Temp:** 20.8 °C

**Relinquished by:** Client

**Serving Size #1:** 1 g

**Density:** 0.8920 g/ml

### Sample Results

Potency per 1g					
Method J AOAC 2015 V98-6 (mod)Units mg/se Batch: 2202413 Analyze: 3/21/22 10:58:00 AM					
Analyte	Result	Limits	Units	LOQ	Notes
CBC per 1g <sup>†</sup>	0.318		mg/1g	0.0325	
CBC-A per 1g <sup>†</sup>	< LOQ		mg/1g	0.0290	
CBC-Total per 1g <sup>†</sup>	0.318		mg/1g	0.0611	
CBD per 1g	8.87		mg/1g	0.0325	
CBD-A per 1g	< LOQ		mg/1g	0.0290	
CBD-Total per 1g	8.87		mg/1g	0.0611	
CBDV per 1g <sup>†</sup>	< LOQ		mg/1g	0.0290	
CBDV-A per 1g <sup>†</sup>	< LOQ		mg/1g	0.0290	
CBDV-Total per 1g <sup>†</sup>	< LOQ		mg/1g	0.0542	
CBE per 1g <sup>†</sup>	0.0889		mg/1g	0.0325	
CBG per 1g <sup>†</sup>	0.293		mg/1g	0.0325	
CBG-A per 1g <sup>†</sup>	< LOQ		mg/1g	0.0290	
CBG-Total per 1g <sup>†</sup>	0.293		mg/1g	0.0607	
CBL per 1g <sup>†</sup>	< LOQ		mg/1g	0.0290	
CBL-A per 1g <sup>†</sup>	< LOQ		mg/1g	0.0290	
CBL-Total per 1g <sup>†</sup>	< LOQ		mg/1g	0.0545	
CBN per 1g	< LOQ		mg/1g	0.0290	
CBT per 1g <sup>†</sup>	0.124		mg/1g	0.0325	
Δ8-THCV per 1g <sup>†</sup>	< LOQ		mg/1g	0.0290	
Δ8-THC per 1g <sup>†</sup>	< LOQ		mg/1g	0.0290	
Δ9-THC per 1g	0.417		mg/1g	0.0325	
exo-THC per 1g <sup>†</sup>	< LOQ		mg/1g	0.0290	
THC-A per 1g	< LOQ		mg/1g	0.0290	
THC-Total per 1g	0.417		mg/1g	0.0611	
THCV per 1g <sup>†</sup>	< LOQ		mg/1g	0.0290	
THCV-A per 1g <sup>†</sup>	< LOQ		mg/1g	0.0290	
THCV-Total per 1g <sup>†</sup>	< LOQ		mg/1g	0.0545	
Total Cannabinoids per 1g	10.1		mg/1g		



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These test results are representative of the individual sample selected and submitted by the client.

**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.

† = Analyte not NELAP accredited.

**Units of Measure**

- g = g
- g/ml = Gram per milliliter
- mg/1g = Milligram per 1g
- % = Percentage of sample
- % wt =  $\mu\text{g/g}$  divided by 10,000

Approved Signatory

Derrick Tanner  
General Manager



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12423 NE Whitaker Way Portland OR 97230 p.503-254-1794

**Cannabis Chain of Custody Record**

ORELAP ID: OR100028

Form Dept Field ID		Date/Time Collected	Pesticides - OR 59 compounds	Pesticide Multi-Residue - 379 compounds	Potency	Residual Solvents	Water Activity	Moisture	Terpenes	Micro: Yeast and Mold	Micro: E. Coli and Total Coliform	Heavy Metals	Mycotoxins	Other	Matrix	Weight	Serving size for edibles	Comments/Metric ID
EC19-MD.O.FS8		3/14/22			X										Liquid	Typical	mg/g	La 2. Met. Discant
EC19-MD.O.FS8		3/14/22	X			X			X	X	X	X				7TN		potency 1st

Purchase Order Number:  
Project Number:  
Project Name:  
 Report Instructions:  
 Send to State - METRC  
 Email Final Results:  
 Fax Final Results  
 Cash/Check/CC/Net 30  
Other:

Collected By:	Relinquished By:	Date	Time	Received by:	Date	Time	Lab Use Only:
<input checked="" type="checkbox"/> Standard (5 day)							Client Alias:
<input type="checkbox"/> Rush (3-4 day) (1.5x Standard)							Order Number:
<input type="checkbox"/> Priority Rush (2 day) (2x Standard)							Proper Container
							Sample Condition
							Temperature: 20.8°C
							Shipped Via: client
							Evidence of cooling: <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No

SUBMISSION OF SAMPLES WITH TESTING REQUIREMENTS TO PIXIS WILL BE UNDERSTOOD TO BE AN AGREEMENT FOR SERVICES IN ACCORDANCE WITH THE CONDITIONS LISTED ON THE BACK OF THIS FORM  
Revision: 1.02 Control#: CF023 Effective 01/31/2019 Revised 01/31/2019  
www.pixislabs.com  
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**Purchase Order:**  
**Received:** 03/17/22 14:20



Revision 1 Documen D 7148  
Legacy D Workshee Valida ed 04/20/2021

**Laboratory Quality Control Results**

**J AOAC 2015 V98-6** **Batch ID: 2202413**

Laboratory Control Sample								
Analyte	Result	Spike	Units	% Rec	Limits		Evaluation	Notes
CBDVA	0.00951	0.01	%	95.1	80.0	- 120	Acceptable	
CBDV	0.0127	0.01	%	127	80.0	- 120	Acceptable	Q6
CBE	0.00955	0.01	%	95.5	80.0	- 120	Acceptable	
CBDA	0.0102	0.01	%	102	80.0	- 120	Acceptable	
CBGA	0.00956	0.01	%	95.6	80.0	- 120	Acceptable	
CBG	0.00975	0.01	%	97.5	80.0	- 120	Acceptable	
CBD	0.00993	0.01	%	99.3	80.0	- 120	Acceptable	
THCV	0.00971	0.01	%	97.1	80.0	- 120	Acceptable	
d8THCV	0.0100	0.01	%	100	80.0	- 120	Acceptable	
THCVA	0.00919	0.01	%	91.9	80.0	- 120	Acceptable	
CBN	0.0104	0.01	%	104	80.0	- 120	Acceptable	
exo-THC	0.00914	0.01	%	91.4	80.0	- 120	Acceptable	
d9THC	0.0100	0.01	%	100	80.0	- 120	Acceptable	
d8THC	0.0100	0.01	%	100	80.0	- 120	Acceptable	
CBL	0.00908	0.01	%	90.8	80.0	- 120	Acceptable	
CBC	0.0101	0.01	%	101	80.0	- 120	Acceptable	
THCA	0.00959	0.01	%	95.9	80.0	- 120	Acceptable	
CBCA	0.00988	0.01	%	98.8	80.0	- 120	Acceptable	
CBLA	0.00958	0.01	%	95.8	80.0	- 120	Acceptable	
CBT	0.0101	0.01	%	101	80.0	- 120	Acceptable	

**Method Blank**

Analyte	Result	LOQ	Units	Limits		Evaluation	Notes
CBDVA	< LOQ	0.003	%	< 0.003		Acceptable	
CBDV	< LOQ	0.003	%	< 0.003		Acceptable	
CBE	< LOQ	0.003	%	< 0.003		Acceptable	
CBDA	< LOQ	0.003	%	< 0.003		Acceptable	
CBGA	< LOQ	0.003	%	< 0.003		Acceptable	
CBG	< LOQ	0.003	%	< 0.003		Acceptable	
CBD	< LOQ	0.003	%	< 0.003		Acceptable	
THCV	< LOQ	0.003	%	< 0.003		Acceptable	
d8THCV	< LOQ	0.003	%	< 0.003		Acceptable	
THCVA	< LOQ	0.003	%	< 0.003		Acceptable	
CBN	< LOQ	0.003	%	< 0.003		Acceptable	
exo-THC	< LOQ	0.003	%	< 0.003		Acceptable	
d9THC	< LOQ	0.003	%	< 0.003		Acceptable	
d8THC	< LOQ	0.003	%	< 0.003		Acceptable	
CBL	< LOQ	0.003	%	< 0.003		Acceptable	
CBC	< LOQ	0.003	%	< 0.003		Acceptable	
THCA	< LOQ	0.003	%	< 0.003		Acceptable	
CBCA	< LOQ	0.003	%	< 0.003		Acceptable	
CBLA	< LOQ	0.003	%	< 0.003		Acceptable	
CBT	< LOQ	0.003	%	< 0.003		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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**Report Number:** 22-003105/D002.R000  
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Revision 1 Document D 7148  
Legacy D Workshee Validated 04/20/2021

**Laboratory Quality Control Results**

J AOAC 2015 V98-6								
Batch ID: 2202413								
Sample Duplicate								
Sample D 22-003006-0001								
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
CBDV	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
CBE	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
CBDA	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
CBGA	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
CBG	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
CBD	14.7	15.3	0.003	%	4.11	< 20	Acceptable	
THCV	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
d8THCV	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
THCVA	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
CBN	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
exo-THC	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
d9THC	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
d8THC	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
CBL	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
CBC	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
THCA	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
CBCA	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
CBLA	< LOQ	< LOQ	0.003	%	NA	< 20	Acceptable	
CBT	< LOQ	< LOQ	0.003	%	NA	< 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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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:** 22-003105/D004.R000  
**Report Date:** 04/07/2022  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 03/17/22 14:20

**Customer:** Etz Hayim Holdings  
**Product identity:** EC19-MO.O.FS8  
**Client/Metric ID:** .  
**Laboratory ID:** 22-003105-0002

### Summary

-----  
**Residual Solvents:**

All analytes passing and less than LOQ.

**Pesticides:**

All analytes passing and less than LOQ.

**Terpenes:**

Less than LOQ for all analytes.

**Metals:**

Less than LOQ for all analytes.

**Microbiology:**

Less than LOQ for all analytes.  
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**Purchase Order:**  
**Received:** 03/17/22 14:20

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

**Product identity:** EC19-MO.O.FS8

**Client/Metric ID:** .

**Sample Date:**

**Laboratory ID:** 22-003105-0002

**Evidence of Cooling:** No

**Temp:** 20.8 °C

**Relinquished by:** Client

### Sample Results

#### Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyze	Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2202804	04/03/22	AOAC 991.14 (Petrifilm)	X, I	
Total Coliforms	< LOQ		cfu/g	10	2202804	04/03/22	AOAC 991.14 (Petrifilm)	X, I	
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2202806	04/04/22	AOAC 2014.05 (RAPID)	X, I	
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2202806	04/04/22	AOAC 2014.05 (RAPID)	X, I	

Solvents		Method Residual Solvents by GC/MS				Units µg/g	Batch 2202942	Analyze 04/06/22 08:55 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-Dimethyl butane	< LOQ		30.0			2,2-Dimethylpropane (neo-pentane)	< LOQ		200		
2,3-Dimethyl butane	< 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 <sup>l</sup>	< 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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Pesticides											
Method AOAC 2007.01 & EN 15662 (mod) Units mg/kg Batch 2202924 Analyze 04/05/22 02:43 PM											
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
Abamectin	< LOQ	0.50	0.250	pass		Acephate	< LOQ	0.40	0.250	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	
Bifenazate	< 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		Etoxazole	< 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	
Imazalil	< 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		Paclotrazole	< LOQ	0.40	0.200	pass	
Parathion-Methyl	< LOQ	0.20	0.200	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	0.200	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							



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**Report Number:** 22-003105/D004.R000  
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**Purchase Order:**  
**Received:** 03/17/22 14:20

Terpenes				Method J AOAC 2015 V98-6	Units %	Batch 2202887	Analyze 04/04/22 06:26 PM		
Analyte	Result	LOQ	% of Total	Notes	Analyte	Result	LOQ	% of Total	Notes
Geranyl acetate <sup>†</sup>	< LOQ	0.018	0.00%		Geraniol <sup>†</sup>	< LOQ	0.018	0.00%	
nerol <sup>†</sup>	< LOQ	0.018	0.00%		farnesene <sup>†</sup>	< LOQ	0.018	0.00%	
Menthol <sup>†</sup>	< LOQ	0.018	0.00%		(-)-a-Terpineol <sup>†</sup>	< LOQ	0.018	0.00%	
(+)-Pulegone <sup>†</sup>	< LOQ	0.018	0.00%		β-Caryophyllene <sup>†</sup>	< LOQ	0.018	0.00%	
a-cedrene <sup>†</sup>	< LOQ	0.018	0.00%		(+)-Cedrol <sup>†</sup>	< LOQ	0.018	0.00%	
(-)-caryophyllene oxide <sup>†</sup>	< LOQ	0.018	0.00%		(-)-Guaiol <sup>†</sup>	< LOQ	0.018	0.00%	
valencene <sup>†</sup>	< LOQ	0.018	0.00%		(-)-Isopulegol <sup>†</sup>	< LOQ	0.018	0.00%	
(-)-β-Pinene <sup>†</sup>	< LOQ	0.018	0.00%		(+)-Borneol <sup>†</sup>	< LOQ	0.018	0.00%	
(+)-fenchol <sup>†</sup>	< LOQ	0.018	0.00%		(±)-Camphor <sup>†</sup>	< LOQ	0.018	0.00%	
(±)-cis-Nerolidol <sup>†</sup>	< LOQ	0.018	0.00%		(±)-fenchone <sup>†</sup>	< LOQ	0.018	0.00%	
(±)-trans-Nerolidol <sup>†</sup>	< LOQ	0.018	0.00%		(R)-(+)-Limonene <sup>†</sup>	< LOQ	0.018	0.00%	
a-Bisabolol <sup>†</sup>	< LOQ	0.018	0.00%		a-phellandrene <sup>†</sup>	< LOQ	0.018	0.00%	
a-pinene <sup>†</sup>	< LOQ	0.018	0.00%		a-Terpinene <sup>†</sup>	< LOQ	0.018	0.00%	
Camphene <sup>†</sup>	< LOQ	0.018	0.00%		cis-β-Ocimene <sup>†</sup>	< LOQ	0.006	0.00%	
d-3-Carene <sup>†</sup>	< LOQ	0.018	0.00%		Eucalyptol <sup>†</sup>	< LOQ	0.018	0.00%	
gamma-Terpinene <sup>†</sup>	< LOQ	0.018	0.00%		Humulene <sup>†</sup>	< LOQ	0.018	0.00%	
Isoborneol <sup>†</sup>	< LOQ	0.018	0.00%		Linalool <sup>†</sup>	< LOQ	0.018	0.00%	
p-Cymene <sup>†</sup>	< LOQ	0.018	0.00%		Sabinene <sup>†</sup>	< LOQ	0.018	0.00%	
Sabinene hydrate <sup>†</sup>	< LOQ	0.018	0.00%		β-Myrcene <sup>†</sup>	< LOQ	0.018	0.00%	
Terpinolene <sup>†</sup>	< LOQ	0.018	0.00%		trans-β-Ocimene <sup>†</sup>	< LOQ	0.012	0.00%	
<b>Total Terpenes</b>	<b>&lt; LOQ</b>								

Metals									
Analyte	Result	Limits	Units	LOQ	Batch	Analyze	Method	Status	Notes
Arsenic	< LOQ	0.200	mg/kg	0.0875	2202970	04/06/22	AOAC 2013.06 (mod.)	pass	X
Cadmium	< LOQ	0.200	mg/kg	0.0875	2202970	04/06/22	AOAC 2013.06 (mod.)	pass	X
Lead	< LOQ	0.500	mg/kg	0.0875	2202970	04/06/22	AOAC 2013.06 (mod.)	pass	X
Mercury	< LOQ	0.100	mg/kg	0.0437	2202970	04/06/22	AOAC 2013.06 (mod.)	pass	X



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**Report Number:** 22-003105/D004.R000  
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**Purchase Order:**  
**Received:** 03/17/22 14:20

These test results are representative of the individual sample selected and submitted by the client.

**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.

† = Analyte not NELAP accredited.

**Units of Measure**

cfu/g = Colony forming units per gram

µg/g = Microgram per gram

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

% = Percentage of sample

% wt = µg/g divided by 10,000

**Glossary of Qualifiers**

I: Insufficient sample received to meet method requirements.

X: Not ORELAP accredited.

Approved Signatory

Derrick Tanner  
General Manager



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Report Number: 22-003105/D004.R000  
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12423 NE Whitaker Way Portland OR 97230 p.503-254-1794

**Cannabis Chain of Custody Record**

ORELAP ID: OR100028

Form Dept Field ID		Date/Time Collected	Pesticides - OR 59 compounds	Pesticide Multi-Residue - 379 compounds	Potency	Residual Solvents	Water Activity	Moisture	Terpenes	Micro: Yeast and Mold	Micro: E. Coli and Total Coliform	Heavy Metals	Mycotoxins	Other:	Matrix	Weight	Serving size for edibles	Comments/Metric ID
EC19-MD.O.FS8		3/14 Sp			X										Liquid	Typical of TN	mg/g	La 2. Ndt. Discourt
EC19-MD.O.FS8		3/14 Sp	X			X			X	X	X	X						potency   SE

Purchase Order Number:  
Project Number:  
Project Name:  
 Report Instructions:  
 Send to State - METRC  
 Email Final Results:  
 Fax Final Results  
 Cash/Check/CC/Net 30  
Other:

Collected By:	Relinquished By:	Date	Time	Received by:	Date	Time	Lab Use Only:
<input checked="" type="checkbox"/> Standard (5 day)							Client Alias:
<input type="checkbox"/> Rush (3-4 day) (1.5x Standard)							Order Number:
<input type="checkbox"/> Priority Rush (2 day) (2x Standard)							Proper Container
							Sample Condition
							Temperature: 20.8°C
							Shipped Via: Client
							Evidence of cooling: <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No

SUBMISSION OF SAMPLES WITH TESTING REQUIREMENTS TO PIXIS WILL BE UNDERSTOOD TO BE AN AGREEMENT FOR SERVICES IN ACCORDANCE WITH THE CONDITIONS LISTED ON THE BACK OF THIS FORM  
Revision: 1.02 Control#: CF023 Effective 01/31/2019 Revised 01/31/2019 www.pixislabs.com Page 1 of 2



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

**Terpenes Quality Control Results**

Method Reference EPA 5035				Batch ID 2202887					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	LCS	Units	LCS % Rec	Limits	Notes
a-pinene	< OQ	< 200		497	500	µg/g	99%	70 - 30	
Camphene	< OQ	< 200		49	500	µg/g	98%	70 - 30	
Sabinene	< OQ	< 200		496	500	µg/g	99%	70 - 30	
b-Pinene	< OQ	< 200		482	500	µg/g	96%	70 - 30	
b-Myrcene	< OQ	< 200		492	500	µg/g	98%	70 - 30	
a-phellandrene	< OQ	< 200		40	500	µg/g	80%	70 - 30	
d-3-Carene	< OQ	< 200		547	500	µg/g	109%	70 - 30	
a-Terpinene	< OQ	< 200		449	500	µg/g	90%	70 - 30	
p-Cymene	< OQ	< 200		487	500	µg/g	97%	70 - 30	
D-imonene	< OQ	< 200		45	500	µg/g	90%	70 - 30	
α-caryophyllol	< OQ	< 200		466	500	µg/g	93%	70 - 30	
b-cis-Ocimene	< OQ	< 67		49	67	µg/g	90%	70 - 30	
b-trans-Ocimene	< OQ	< 33		279	333	µg/g	84%	70 - 30	
γ-Terpinene	< OQ	< 200		4	500	µg/g	82%	70 - 30	
Sabinene hydrate	< OQ	< 200		454	500	µg/g	91%	70 - 30	
Terpinolene	< OQ	< 200		443	500	µg/g	89%	70 - 30	
D-enchone	< OQ	< 200		433	500	µg/g	87%	70 - 30	
linalool	< OQ	< 200		446	500	µg/g	89%	70 - 30	
α-enchol	< OQ	< 200		467	500	µg/g	93%	70 - 30	
Camphor	< OQ	< 200		482	500	µg/g	96%	70 - 30	
sopulego	< OQ	< 200		485	500	µg/g	97%	70 - 30	
soborneol	< OQ	< 200		48	500	µg/g	96%	70 - 30	
Borneol	< OQ	< 200		474	500	µg/g	95%	70 - 30	
D-Menthol	< OQ	< 200		489	500	µg/g	98%	70 - 30	
Terpineol	< OQ	< 200		404	500	µg/g	81%	70 - 30	
Nerol	< OQ	< 200		429	500	µg/g	86%	70 - 30	
Pulegone	< OQ	< 200		476	500	µg/g	95%	70 - 30	
Geraniol	< OQ	< 200		377	500	µg/g	75%	70 - 30	
Geranyl Acetate	< OQ	< 200		496	500	µg/g	99%	70 - 30	
α-Cedrene	< OQ	< 200		49	500	µg/g	98%	70 - 30	
b-Caryophyllene	< OQ	< 200		500	500	µg/g	100%	70 - 30	
α-Humulene	< OQ	< 200		483	500	µg/g	97%	70 - 30	
Valenene	< OQ	< 200		495	500	µg/g	99%	70 - 30	
cis-Nerolidol	< OQ	< 200		503	500	µg/g	101%	70 - 30	
α-arnesene	< OQ	< 200		564	500	µg/g	113%	70 - 30	
trans-Nerolidol	< OQ	< 200		55	500	µg/g	103%	70 - 30	
Caryophyllene Oxide	< OQ	< 200		498	500	µg/g	100%	70 - 30	
Guaiol	< OQ	< 200		529	500	µg/g	106%	70 - 30	
Cedrol	< OQ	< 200		53	500	µg/g	103%	70 - 30	
α-Bisabolol	< OQ	< 200		55	500	µg/g	103%	70 - 30	

Definitions

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





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**Report Number:** 22-003105/D004.R000  
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**Purchase Order:**  
**Received:** 03/17/22 14:20

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

**Terpenes Quality Control Results**

Method Reference EPA 5035		Batch ID 2202887					
Sample/Sample Duplicate		Sample ID 22-003397-0002					
Analyte	Result	Org. Result	LOQ	Units	% RPD	LIMIT	Notes
a-pinene	< OQ	< OQ	83	µg/g	0%	< 20	
Camphene	< OQ	< OQ	83	µg/g	0%	< 20	
Sabinene	< OQ	< OQ	83	µg/g	0%	< 20	
b-Pinene	< OQ	< OQ	83	µg/g	0%	< 20	
b-Myrcene	46600	44 00	83	µg/g	6%	< 20	
a-phellandrene	< OQ	< OQ	83	µg/g	0%	< 20	
d-3-Carene	523	524	83	µg/g	0%	< 20	
a-Terpinene	< OQ	< OQ	83	µg/g	0%	< 20	
p-Cymene	< OQ	< OQ	83	µg/g	0%	< 20	
D-imonene	0000	08000	83	µg/g	2%	< 20	
α-caryophyllol	< OQ	< OQ	83	µg/g	0%	< 20	
b-cis-Ocimene	< OQ	< OQ	60.9	µg/g	0%	< 20	
b-trans-Ocimene	< OQ	< OQ	22	µg/g	0%	< 20	
g-Terpinene	< OQ	< OQ	83	µg/g	0%	< 20	
Sabinene hydrate	< OQ	< OQ	83	µg/g	0%	< 20	
Terpinolene	< OQ	< OQ	83	µg/g	0%	< 20	
D-enchone	< OQ	< OQ	83	µg/g	0%	< 20	
α-inalool	< OQ	< OQ	83	µg/g	0%	< 20	
α-enchol	< OQ	< OQ	83	µg/g	0%	< 20	
Camphor	< OQ	< OQ	83	µg/g	0%	< 20	
α-sopulego	< OQ	< OQ	83	µg/g	0%	< 20	
α-soborneol	< OQ	< OQ	83	µg/g	0%	< 20	
Borneol	< OQ	< OQ	83	µg/g	0%	< 20	
D-Menthol	< OQ	< OQ	83	µg/g	0%	< 20	
Terpineol	< OQ	< OQ	83	µg/g	0%	< 20	
Nerol	< OQ	< OQ	83	µg/g	0%	< 20	
Pulegone	< OQ	< OQ	83	µg/g	0%	< 20	
Geraniol	< OQ	< OQ	83	µg/g	0%	< 20	
Geranyl acetate	< OQ	< OQ	83	µg/g	0%	< 20	
α-Cedrene	< OQ	< OQ	83	µg/g	0%	< 20	
b-Caryophyllene	9900	9900	83	µg/g	0%	< 20	
α-Humulene	420	4 0	83	µg/g	1%	< 20	
Valenene	< OQ	< OQ	83	µg/g	0%	< 20	
cis-Nerolidol	< OQ	< OQ	83	µg/g	0%	< 20	
α-arnesene	< OQ	< OQ	83	µg/g	0%	< 20	
trans-Nerolidol	< OQ	< OQ	83	µg/g	0%	< 20	
Caryophyllene Oxide	228	235	83	µg/g	3%	< 20	
Guaiol	< OQ	< OQ	83	µg/g	0%	< 20	
Cedrol	< OQ	< OQ	83	µg/g	0%	< 20	
α-Bisabolol	< OQ	< OQ	83	µg/g	0%	< 20	

Definitions

RPD Relative Percent Difference



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**Laboratory Pesticide Quality Control Results**

AOAC 2007.1 & EN 15662		Units: mg/Kg			Batch ID: 2202924			
Method Blank		Laboratory Control Sample						
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes
Abamectin	0.000	< 0.250		1.889	1.000	188.9	50.0	150
Acephate	0.000	< 0.250		1.017	1.000	101.7	60.0	120
Acetaminocyl	0.000	< 1.000		4.263	4.000	106.6	40.0	160
Acetamiprid	0.000	< 0.100		0.395	0.400	98.8	60.0	120
Aldicarb	0.000	< 0.200		0.779	0.800	97.4	60.0	120
Azoxystrobin	0.000	< 0.100		0.385	0.400	96.3	60.0	120
Bifenazate	0.000	< 0.100		0.387	0.400	96.7	60.0	120
Bifenthrin	0.000	< 0.100		0.414	0.400	103.5	50.0	150
Boscalid	0.000	< 0.200		0.841	0.800	105.2	60.0	120
Carbaryl	0.000	< 0.100		0.397	0.400	99.4	60.0	120
Carbofuran	0.000	< 0.100		0.391	0.400	97.6	60.0	120
Chlorantraniliprole	0.000	< 0.100		0.384	0.400	95.9	60.0	120
Chlorfenapyr	0.000	< 0.500		2.217	2.000	110.8	60.0	120
Chlorpyrifos	0.000	< 0.100		0.404	0.400	101.1	60.0	120
Clofentazine	0.000	< 0.100		0.400	0.400	100.0	60.0	120
Cyfluthrin	0.000	< 0.500		1.642	2.000	82.1	50.0	150
Cypermethrin	0.000	< 0.500		1.997	2.000	99.8	50.0	150
Daminozide	0.000	< 0.500		2.435	2.000	121.8	60.0	120
Diazinon	0.000	< 0.100		0.404	0.400	100.9	60.0	120
Dichlorvos	0.000	< 0.500		2.057	2.000	102.8	60.0	120
Dimethoate	0.000	< 0.100		0.404	0.400	101.0	60.0	120
Ethoprophos	0.000	< 0.100		0.403	0.400	100.8	60.0	120
Etofenprox	0.000	< 0.200		0.819	0.800	102.4	50.0	150
Etoxazole	0.000	< 0.100		0.412	0.400	102.9	60.0	120
Fenoxycarb	0.000	< 0.100		0.393	0.400	98.2	60.0	120
Fenpyroximate	0.000	< 0.200		0.796	0.800	99.6	60.0	120
Fipronil	0.000	< 0.200		0.787	0.800	98.4	60.0	120
Fonicamid	0.000	< 0.250		1.065	1.000	106.5	60.0	120
Fludioxonil	0.000	< 0.200		0.794	0.800	99.3	50.0	150
Hexythiazox	0.000	< 0.250		1.027	1.000	102.7	60.0	120
Imazalil	0.000	< 0.100		0.395	0.400	98.8	60.0	120
Imidacloprid	0.000	< 0.200		0.784	0.800	98.0	60.0	120
Kresoxim methyl	0.000	< 0.200		0.840	0.800	105.1	60.0	120
Malathion	0.000	< 0.100		0.403	0.400	100.6	60.0	120
Metlaxyl	0.000	< 0.100		0.395	0.400	98.8	60.0	120
Methiocarb	0.000	< 0.100		0.385	0.400	96.3	60.0	120
Methomyl	0.000	< 0.200		0.728	0.800	90.9	60.0	120
MGK 264	0.000	< 0.100		0.401	0.400	100.2	50.0	150
Myclobutanil	0.000	< 0.100		0.394	0.400	98.4	60.0	120
Naled	0.000	< 0.250		0.976	1.000	97.6	50.0	150
Oxamyl	0.000	< 0.500		1.994	2.000	99.7	60.0	120
Paclobotrazole	0.000	< 0.200		0.782	0.800	97.8	60.0	120
Parathion Methyl	0.000	< 0.200		0.893	0.800	111.6	50.0	150
Permethrin	0.000	< 0.100		0.401	0.400	100.2	50.0	150
Phosmet	0.000	< 0.100		0.396	0.400	99.0	50.0	150
Piperonyl butoxide	0.000	< 0.500		2.376	2.000	118.8	60.0	120
Prallethrin	0.000	< 0.100		0.411	0.400	102.8	60.0	120
Propiconazole	0.000	< 0.200		0.819	0.800	102.3	60.0	120
Propoxur	0.000	< 0.100		0.397	0.400	99.3	60.0	120
Pyrethrin (Summe)	0.000	< 0.100		0.418	0.413	101.1	60.0	120
Pyridaben	0.000	< 0.100		0.409	0.400	102.3	50.0	150
Spirosad	0.000	< 0.100		0.409	0.388	105.4	50.0	150
Spiromesifen	0.000	< 0.100		0.422	0.400	105.4	60.0	120
Spirotetramat	0.000	< 0.100		0.404	0.400	101.1	60.0	120
Spiroxamine	0.000	< 0.200		0.806	0.800	100.8	60.0	120
ebuconazole	0.000	< 0.200		0.801	0.800	100.1	60.0	120
hiacloprid	0.000	< 0.100		0.395	0.400	98.8	60.0	120
hiamethoxam	0.000	< 0.100		0.424	0.400	105.9	60.0	120
rifloxystrobin	0.000	< 0.100		0.402	0.400	100.4	60.0	120



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Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662		Units: mg/Kg					Batch ID: 2202924				
Matrix Spike/Matrix Spike Duplicate Recoveries						Sample ID: 22-003658-0001					
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes	
Abamectin	0.000	2.116	2.090	1.000	1.2%	< 30	211.6%	209.0%	50 150	Q	
Acephate	0.000	1.001	0.986	1.000	1.5%	< 30	100.1%	98.6%	50 150		
Acetaminocyl	0.000	5.131	4.659	4.000	9.6%	< 30	128.3%	116.5%	50 150		
Acetamiprid	0.000	0.388	0.391	0.400	0.7%	< 30	97.0%	97.7%	50 150		
Aldicarb	0.000	0.764	0.757	0.800	0.9%	< 30	95.5%	94.7%	50 150		
Azoxystrobin	0.000	0.388	0.401	0.400	3.5%	< 30	96.9%	100.4%	50 150		
Bifenazate	0.000	0.422	0.412	0.400	2.5%	< 30	105.6%	103.0%	50 150		
Bifenthrin	0.000	0.455	0.459	0.400	0.9%	< 30	113.8%	114.8%	50 150		
Boscalid	0.000	0.781	0.747	0.800	4.4%	< 30	97.6%	93.4%	50 150		
Carbaryl	0.000	0.389	0.389	0.400	0.0%	< 30	97.2%	97.2%	50 150		
Carbofuran	0.000	0.382	0.390	0.400	2.3%	< 30	95.4%	97.6%	50 150		
Chlorantraniliprole	0.000	0.369	0.379	0.400	2.7%	< 30	92.2%	94.7%	50 150		
Chlorfenapyr	0.000	1.985	1.761	2.000	12.0%	< 30	99.3%	88.1%	50 150		
Chlorpyrifos	0.000	0.414	0.360	0.400	13.9%	< 30	103.6%	90.1%	50 150		
Clofentazine	0.000	0.453	0.453	0.400	0.1%	< 30	113.2%	113.4%	50 150		
Cyfluthrin	0.000	2.018	1.862	2.000	8.1%	< 30	100.9%	93.1%	30 150		
Cypermethrin	0.000	2.057	1.600	2.000	25.0%	< 30	102.9%	80.0%	50 150		
Daminozide	0.173	2.300	2.274	2.000	1.2%	< 30	106.3%	105.0%	30 150		
Diazinon	0.000	0.392	0.382	0.400	2.7%	< 30	98.0%	95.4%	50 150		
Dichlorvos	0.000	1.982	1.908	2.000	3.8%	< 30	99.1%	95.4%	50 150		
Dimethoate	0.000	0.390	0.391	0.400	0.4%	< 30	97.4%	97.8%	50 150		
Ethoprophos	0.000	0.384	0.411	0.400	6.8%	< 30	95.9%	102.6%	50 150		
Etofenprox	0.000	0.962	0.921	0.800	4.3%	< 30	120.2%	115.2%	50 150		
Etoxazole	0.000	0.519	0.530	0.400	2.1%	< 30	129.7%	132.4%	50 150		
Fenoxycarb	0.000	0.388	0.384	0.400	1.0%	< 30	97.0%	96.1%	50 150		
Fenpyroximate	0.000	0.861	0.888	0.800	3.1%	< 30	107.6%	111.0%	50 150		
Fipronil	0.000	0.872	0.901	0.800	3.3%	< 30	108.9%	112.6%	50 150		
Fonicamid	0.000	1.032	1.012	1.000	2.0%	< 30	103.2%	101.2%	50 150		
Fludioxonil	0.000	0.739	0.764	0.800	3.4%	< 30	92.3%	95.6%	50 150		
Hexythiazox	0.000	1.054	1.097	1.000	4.0%	< 30	105.4%	109.7%	50 150		
Imazalil	0.000	0.314	0.310	0.400	1.3%	< 30	78.4%	77.4%	50 150		
Imidacloprid	0.000	0.770	0.765	0.800	0.7%	< 30	96.3%	95.6%	50 150		
Kresoxim methyl	0.000	0.823	0.796	0.800	3.4%	< 30	102.9%	99.5%	50 150		
Malathion	0.000	0.396	0.389	0.400	1.9%	< 30	99.0%	97.2%	50 150		
Metaxalyl	0.000	0.373	0.372	0.400	0.1%	< 30	93.2%	93.1%	50 150		
Methiocarb	0.000	0.381	0.381	0.400	0.0%	< 30	95.2%	95.3%	50 150		
Methomyl	0.000	0.650	0.753	0.800	14.7%	< 30	81.2%	94.1%	50 150		
MGK 264	0.000	0.418	0.412	0.400	1.4%	< 30	104.5%	103.1%	50 150		
Myclobutanil	0.000	0.371	0.378	0.400	1.7%	< 30	92.8%	94.5%	50 150		
Naled	0.000	0.882	0.919	1.000	4.1%	< 30	88.2%	91.9%	50 150		
Oxamyl	0.000	1.914	1.812	2.000	5.5%	< 30	95.7%	90.6%	50 150		
Paclobotrazole	0.000	0.746	0.736	0.800	1.4%	< 30	93.3%	92.0%	50 150		
Parathion Methyl	0.000	0.758	0.770	0.800	1.6%	< 30	94.7%	96.2%	30 150		
Permethrin	0.000	0.538	0.566	0.400	5.1%	< 30	134.6%	141.6%	50 150		
Phosmet	0.000	0.373	0.378	0.400	1.3%	< 30	93.2%	94.4%	50 150		
Piperonyl butoxide	0.000	2.345	2.435	2.000	3.8%	< 30	117.2%	121.7%	50 150		
Prallethrin	0.000	0.477	0.473	0.400	0.8%	< 30	119.2%	118.2%	50 150		
Propiconazole	0.000	0.797	0.791	0.800	0.6%	< 30	99.6%	98.9%	50 150		
Propoxur	0.000	0.388	0.392	0.400	1.1%	< 30	96.9%	98.0%	50 150		
Pyrethrin (Summe)	4.058	5.235	5.381	0.413	11.7%	< 30	284.9%	320.2%	50 150	Q	
Pyridaben	0.000	0.544	0.509	0.400	6.7%	< 30	135.9%	127.1%	50 150		
Spinosad	0.000	0.440	0.400	0.388	9.4%	< 30	113.3%	103.1%	50 150		
Spiromesifen	0.000	0.412	0.414	0.400	0.4%	< 30	103.0%	103.5%	50 150		
Spirotetramat	0.000	0.398	0.406	0.400	2.0%	< 30	99.5%	101.5%	50 150		
Spiroxamine	0.000	0.785	0.761	0.800	3.1%	< 30	98.1%	95.1%	50 150		
ebuconazole	0.000	0.786	0.768	0.800	2.4%	< 30	98.3%	96.0%	50 150		
hiacloprid	0.000	0.377	0.385	0.400	2.1%	< 30	94.1%	96.1%	50 150		
hiamethoxam	0.000	0.397	0.419	0.400	5.3%	< 30	99.3%	104.7%	50 150		
rifloxystrobin	0.000	0.419	0.415	0.400	1.1%	< 30	104.8%	103.7%	50 150		



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Revision Document ID  
 Legacy ID Effective

Laboratory Quality Control Results									
Residual Solvents									
Method Blank									
Laboratory Control Sample									
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	200		535	5 2	µg/g	93 5	60	20
Isobutane	ND	200		9	3	µg/g	02 5	60	20
Butane	ND	200		66	3	µg/g	0 8	60	20
2,2-Dimethylpropane	ND	200		89	596	µg/g	95 8	60	20
Methanol	ND	200		5 0	620	µg/g	95	60	20
Ethylene Oxide	ND	0		59 5	56 2	µg/g	05 9	60	20
2-Methylbutane	ND	200		590	620	µg/g	98	60	20
Pentane	ND	200		5 0	6 0	µg/g	95	60	20
Ethanol	ND	200		590	630	µg/g	9 5	0	30
Ethyl Ether	ND	200		80	620	µg/g	9	60	20
2,2-Dimethylbutane	ND	0		5		µg/g	90 2	60	20
Acetone	ND	200		360	650	µg/g	82	60	20
2-Propanol	ND	200		0	6 0	µg/g	28 8	60	20
Ethyl Formate	ND	500		0	600	µg/g	28	0	30
Acetonitrile	ND	0		2	98	µg/g	9 8	60	20
Methyl Acetate	ND	500		90	6 0	µg/g	92 5	0	30
2,3-Dimethylbutane	ND	0		6	6	µg/g	9 5	60	20
Dichloromethane	ND	60		85	5 0	µg/g	95 3	60	20
2-Methylpentane	ND	0		50	6	µg/g	85 2	60	20
m-NE	ND	500		0	600	µg/g	90 0	0	30
3-Methylpentane	ND	0		55	5	µg/g	88 6	60	20
Hexane	ND	0		3		µg/g	9	60	20
Propanol	ND	500		5 0	6 0	µg/g	93 8	0	30
Methyl ethyl ketone	ND	500		5 0	600	µg/g	9	0	30
Ethyl acetate	ND	200		500	630	µg/g	92 0	60	20
2-Butanol	ND	200		0	620	µg/g	88 3	60	20
Tetrahydrofuran	ND	0		3	500	µg/g	8	60	20
Cyclohexane	ND	200		390	620	µg/g	85 8	60	20
2-methyl propanol	ND	500		3 0	620	µg/g	80 9	0	30
Benzene	ND			5	5 32	µg/g	85 3	60	20
Isopropyl Acetate	ND	200		3 0	620	µg/g	8 6	60	20
Heptane	ND	200		0		µg/g	9	60	20
Butanol	ND	500		250	600	µg/g	8	0	30
Propyl Acetate	ND	500		50	600	µg/g	90 6	0	30
Dioxane	ND	0		23	50	µg/g	83 9	60	20
2-Ethoxyethanol	ND	0		56	8	µg/g	86 2	60	20
Methylisobutylketone	ND	500		360	6 0	µg/g	8 5	0	30
3-Methyl butanol	ND	500		200	6 0	µg/g	5	0	30
Ethylene Glycol	ND	200		36	9	µg/g	3	60	20
Octane	ND	200		0	9	µg/g	8 6	60	20
Isobutyl Acetate	ND	500		500	600	µg/g	93 8	0	30
Pentanol	ND	500		5 0	6 0	µg/g	95 0	0	30
Butyl Acetate	ND	500		0	6 0	µg/g	9 3	0	30
Ethyl benzene	ND	200		86	9 3	µg/g	88 5	60	20
m-p Xylene	ND	200		85	596	µg/g	86 0	60	20
o-Xylene	ND	200		890	9 3	µg/g	9 5	60	20
Cumene	ND	0		0		µg/g	82 9	60	20
Anisole	ND	500		350	6 0	µg/g	83 9	0	30
DMSO	ND	500		560	630	µg/g	95	0	30
2-dimethoxyethane	ND	50		8	6	µg/g	90 2	0	30
Diethylamine	ND	500		360	600	µg/g	85 0	0	30
N,N-dimethylformamide	ND	50		0	9	µg/g	86 5	0	30
N,N-dimethylacetamide	ND	50		88	98	µg/g	98 0	0	30
Pyridine	ND	50		5	80	µg/g	9 2	0	30
1,2-Dichloroethane	ND			0		µg/g	0	0	30
Chloroform	ND			03		µg/g	03 0	0	30
Trichloroethylene	ND					µg/g	0	0	30



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QC Sample Duplicate		Sample ID: 22 003693 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	60	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	
Mt SE	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	
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	60	µg/g	0.0	20	Acceptable	
Cyclohexane	ND	ND	200	µg/g	0.0	20	Acceptable	
2 methyl propanol	ND	ND	500	µg/g	0.0	20	Acceptable	
Benzene	ND	ND	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	
Butanol	ND	ND	500	µg/g	0.0	20	Acceptable	
Propyl Acetate	ND	ND	500	µg/g	0.0	20	Acceptable	
Dioxane	ND	ND	60	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 butanol	ND	ND	500	µg/g	0.0	20	Acceptable	
Ethylene Glycol	ND	ND	200	µg/g	0.0	20	Acceptable	
okane	ND	ND	200	µg/g	0.0	20	Acceptable	
Isobutyl Acetate	ND	ND	500	µg/g	0.0	20	Acceptable	
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	
2 dimethoxyethane	ND	ND	50	µg/g	0.0	20	Acceptable	
riethylamine	ND	ND	500	µg/g	0.0	20	Acceptable	
N N d methylformam de	ND	ND	50	µg/g	0.0	20	Acceptable	
N N d methylacetamide	ND	ND	50	µg/g	0.0	20	Acceptable	
Pyridine	ND	ND	50	µg/g	0.0	20	Acceptable	
2 Dichloroethane	ND	ND	µg/g	0.0	20	Acceptable		
Chloroform	ND	ND	µg/g	0.0	20	Acceptable		
dichloroethylene	ND	ND	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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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.