



## 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	10 µg/day <sup>[1]</sup>
Cadmium	µg/serving	µg/g	4.1 µg/day <sup>[1]</sup>
Lead	µg/serving	µg/g	6 µg/day <sup>[1]</sup>
Mercury	µg/serving	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*	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		

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:** 23-009657/D003.R000  
**Report Date:** 08/21/2023  
**ORELAP#:** OR100028  
**Purchase Order:** 2595910  
**Received:** 08/14/23 16:11

**Customer:** Etz Hayim Holdings  
**Product identity:** CYCL-GMY.D9.RNA10.V2-FH05  
**Client/Metric ID:** .  
**Laboratory ID:** 23-009657-0001

### Summary

**Potency:**

Analyte per 1g	Result	Limits	Units	Status	
CBC-A per 1g	0.0377		mg/1g		CBD-Total per Serving Size 2.85 mg/1g
CBD per 1g	2.26		mg/1g		
CBD-A per 1g	0.669		mg/1g		THC-Total per Serving Size 2.04 mg/1g
Δ8-THC per 1g	0.0336		mg/1g		(Reported in milligrams per serving)
Δ9-THC per 1g	2.04		mg/1g		

**Residual Solvents:**

All analytes passing and less than LOQ.

**Pesticides:**

All analytes passing and less than LOQ.

**Metals:**

Less than LOQ for all analytes.

**Microbiology:**

Less than LOQ for all analytes.



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

**Product identity:** CYCL-GMY.D9.RNA10.V2-FH05

**Client/Metric ID:** .

**Sample Date:**

**Laboratory ID:** 23-009657-0001

**Evidence of Cooling:** No

**Temp:** 28.2 °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: 2310135	Analyze: 8/18/23 12:31:00 AM	
Analyte	Result	Limits	Units	LOQ	Notes
CBC per 1g	< LOQ		mg/1g	0.0328	
CBC-A per 1g	0.0377		mg/1g	0.0328	
CBC-Total per 1g	< LOQ		mg/1g	0.0615	
CBD per 1g	2.26		mg/1g	0.0328	
CBD-A per 1g	0.669		mg/1g	0.0328	
CBD-Total per 1g	2.85		mg/1g	0.0615	
CBDV per 1g	< LOQ		mg/1g	0.0328	
CBDV-A per 1g	< LOQ		mg/1g	0.0328	
CBDV-Total per 1g	< LOQ		mg/1g	0.0612	
CBE per 1g	< LOQ		mg/1g	0.0328	
CBG per 1g	< LOQ		mg/1g	0.0328	
CBG-A per 1g	< LOQ		mg/1g	0.0328	
CBG-Total per 1g	< LOQ		mg/1g	0.0612	
CBL per 1g	< LOQ		mg/1g	0.0328	
CBL-A per 1g	< LOQ		mg/1g	0.0328	
CBL-Total per 1g	< LOQ		mg/1g	0.0615	
CBN per 1g	< LOQ		mg/1g	0.0328	
CBT per 1g	< LOQ		mg/1g	0.0328	
Δ8-THCV per 1g	< LOQ		mg/1g	0.0328	
Δ10-THC-9R per 1g	< LOQ		mg/1g	0.0328	
Δ10-THC-9S per 1g	< LOQ		mg/1g	0.0328	
Δ10-THC-Total per 1g	< LOQ		mg/1g	0.0656	
Δ8-THC per 1g	0.0336		mg/1g	0.0328	
Δ9-THC per 1g	2.04		mg/1g	0.0328	
delta-9-THCP per 1g	< LOQ		mg/1g	0.0328	
exo-THC per 1g	< LOQ		mg/1g	0.0328	
THC-A per 1g	< LOQ		mg/1g	0.0328	
THC-Total per 1g	2.04		mg/1g	0.0615	
THCV per 1g	< LOQ		mg/1g	0.0328	
THCV-A per 1g	< LOQ		mg/1g	0.0328	



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**Purchase Order:** 2595910  
**Received:** 08/14/23 16:11

Potency per 1g	Method: J AOAC 2015 V98-6 (mod) <sup>P</sup>	Units mg/se	Batch: 2310135	Analyze: 8/18/23	12:31:00 AM
Analyte	Result	Limits	Units	LOQ	Notes
THCV-Total per 1g	< LOQ		mg/1g	0.0616	
Total Cannabinoids per 1g	5.04		mg/1g		

**Microbiology**

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2310050	08/18/23 AOAC 991.14 (Petrifilm) <sup>P</sup>		
Total Coliforms	< LOQ		cfu/g	10	2310050	08/18/23 AOAC 991.14 (Petrifilm) <sup>P</sup>		
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2310051	08/18/23 AOAC 2014.05 (RAPID) <sup>P</sup>		
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2310051	08/18/23 AOAC 2014.05 (RAPID) <sup>P</sup>		

**Solvents** Method: Residual Solvents by GC/MS<sup>P</sup> Units µg/g Batch 2310111 Analyze 08/17/23 12:54 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							


**Pesticides**      **Method:** AOAC 2007.01 & EN 15662 (mod)<sup>b</sup>      **Units** mg/kg      **Batch** 2310121      **Analyze** 08/17/23 03:24 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		Flonicamid <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		Pacllobutrazole <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							

**Metals**

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
Arsenic <sup>‡</sup>	< LOQ	0.200	mg/kg	0.0175	2310198	08/21/23 AOAC 2013.06 (mod.) <sup>b</sup>	pass	
Cadmium <sup>‡</sup>	< LOQ	0.200	mg/kg	0.0175	2310198	08/21/23 AOAC 2013.06 (mod.) <sup>b</sup>	pass	
Lead <sup>‡</sup>	< LOQ	0.500	mg/kg	0.0175	2310198	08/21/23 AOAC 2013.06 (mod.) <sup>b</sup>	pass	
Mercury <sup>‡</sup>	< LOQ	0.100	mg/kg	0.00874	2310198	08/21/23 AOAC 2013.06 (mod.) <sup>b</sup>	pass	



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**Abbreviations**

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

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

Ⓐ = ISO/IEC 17025:2017 accredited method.

Ⓜ = TNI accredited analyte.

**Units of Measure**

cfu/g = Colony forming units per gram

g = g

µg/g = Microgram per gram

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

mg/1g = Milligram per 1g

% = Percentage of sample

% wt = µg/g divided by 10,000

Approved Signatory

Derrick Tanner  
General Manager



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

Laboratory Quality Control Results

Residual Solvents				Batch ID: 2310111					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		486	584	µg/g	83.2	60 - 120	
Isobutane	ND	< 200		695	767	µg/g	90.6	60 - 120	
Butane	ND	< 200		682	782	µg/g	87.2	60 - 120	
2,2-Dimethylpropane	ND	< 200		794	939	µg/g	84.6	60 - 120	
Methanol	ND	< 200		1550	1670	µg/g	92.8	60 - 120	
Ethylene Oxide	ND	< 30		50	57.1	µg/g	87.6	60 - 120	
2-Methylbutane	ND	< 200		1430	1680	µg/g	85.1	60 - 120	
Pentane	ND	< 200		1400	1670	µg/g	83.8	60 - 120	
Ethanol	ND	< 200		1500	1660	µg/g	90.4	70 - 130	
Ethyl Ether	ND	< 200		1430	1670	µg/g	85.6	60 - 120	
2,2-Dimethylbutane	ND	< 30		169	189	µg/g	89.4	60 - 120	
Acetone	ND	< 200		1450	1670	µg/g	86.8	60 - 120	
2-Propanol	ND	< 200		1430	1630	µg/g	87.7	60 - 120	
Ethyl Formate	ND	< 500		4530	1600	µg/g	283.1	70 - 130	Q6
Acetonitrile	ND	< 100		412	492	µg/g	83.7	60 - 120	
Methyl Acetate	ND	< 500		1560	1600	µg/g	97.5	70 - 130	
2,3-Dimethylbutane	ND	< 30		152	180	µg/g	84.4	60 - 120	
Dichloromethane	ND	< 60		427	488	µg/g	87.5	60 - 120	
2-Methylpentane	ND	< 30		155	182	µg/g	85.2	60 - 120	
MTBE	ND	< 500		1590	1610	µg/g	98.8	70 - 130	
3-Methylpentane	ND	< 30		159	177	µg/g	89.8	60 - 120	
Hexane	ND	< 30		150	177	µg/g	84.7	60 - 120	
1-Propanol	ND	< 500		1640	1600	µg/g	102.5	70 - 130	
Methylethylketone	ND	< 500		1530	1610	µg/g	95.0	70 - 130	
Ethyl acetate	ND	< 200		1370	1630	µg/g	84.0	60 - 120	
2-Butanol	ND	< 200		1400	1630	µg/g	85.9	60 - 120	
Tetrahydrofuran	ND	< 100		417	488	µg/g	85.5	60 - 120	
Cyclohexane	ND	< 200		1380	1610	µg/g	85.7	60 - 120	
2-methyl-1-propanol	ND	< 500		1710	1610	µg/g	106.2	70 - 130	
Benzene	ND	< 1		3.58	4.79	µg/g	74.7	60 - 120	
Isopropyl Acetate	ND	< 200		1380	1650	µg/g	83.6	60 - 120	
Heptane	ND	< 200		1350	1630	µg/g	82.8	60 - 120	
1-Butanol	ND	< 500		1770	1600	µg/g	110.6	70 - 130	
Propyl Acetate	ND	< 500		1580	1600	µg/g	98.8	70 - 130	
1,4-Dioxane	ND	< 100		439	523	µg/g	83.9	60 - 120	
2-Ethoxyethanol	ND	< 30		146	179	µg/g	81.6	60 - 120	
Methylisobutylketone	ND	< 500		1610	1600	µg/g	100.6	70 - 130	
3-Methyl-1-butanol	ND	< 500		1710	1600	µg/g	106.9	70 - 130	
Ethylene Glycol	ND	< 200		319	506	µg/g	63.0	60 - 120	
Toluene	ND	< 100		428	496	µg/g	86.3	60 - 120	
Isobutyl Acetate	ND	< 500		1610	1610	µg/g	100.0	70 - 130	
1-Pentanol	ND	< 500		1860	1600	µg/g	116.3	70 - 130	
Butyl Acetate	ND	< 500		1610	1610	µg/g	100.0	70 - 130	
Ethylbenzene	ND	< 200		794	978	µg/g	81.2	60 - 120	
m,p-Xylene	ND	< 200		801	994	µg/g	80.6	60 - 120	
o-Xylene	ND	< 200		797	982	µg/g	81.2	60 - 120	
Cumene	ND	< 30		131	171	µg/g	76.6	60 - 120	
Anisole	ND	< 500		1760	1600	µg/g	110.0	70 - 130	
DMSO	ND	< 500		1400	1620	µg/g	86.4	70 - 130	
1,2-dimethoxyethane	ND	< 50		173	185	µg/g	93.0	70 - 130	
Triethylamine	ND	< 500		1430	1600	µg/g	89.4	70 - 130	
N,N-dimethylformamide	ND	< 150		483	480	µg/g	100.6	70 - 130	
N,N-dimethylacetamide	ND	< 150		478	483	µg/g	99.0	70 - 130	
Pyridine	ND	< 50		161	168	µg/g	95.8	70 - 130	
Sulfolane	ND	< 50		164	161	µg/g	101.9	70 - 130	
1,2-Dichloroethane	ND	< 1		0.837	1	µg/g	83.7	70 - 130	
Chloroform	ND	< 1		0.951	1	µg/g	95.1	70 - 130	
Trichloroethylene	ND	< 1		0.926	1	µg/g	92.6	70 - 130	
1,1-Dichloroethane	ND	< 1		0.758	1	µg/g	75.8	70 - 130	



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

QC - Sample Duplicate		Sample ID: 23-009604-0001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Accept/Fail	Notes
Propane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Isobutane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Butane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2,2-Dimethylpropane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Methanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethylene Oxide	ND	ND	30	µg/g	0.0	< 20	Acceptable	
2-Methylbutane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Pentane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethyl Ether	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2,2-Dimethylbutane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Acetone	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2-Propanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethyl Formate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Acetonitrile	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Methyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
2,3-Dimethylbutane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Dichloromethane	ND	ND	60	µg/g	0.0	< 20	Acceptable	
2-Methylpentane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
MTBE	ND	ND	500	µg/g	0.0	< 20	Acceptable	
3-Methylpentane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Hexane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
1-Propanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Methyl ethyl ketone	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Ethyl acetate	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2-Butanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Tetrahydrofuran	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Cyclohexane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2-methyl-1-propanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Benzene	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Isopropyl Acetate	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Heptane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
1-Butanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Propyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
1,4-Dioxane	ND	ND	100	µg/g	0.0	< 20	Acceptable	
2-Ethoxyethanol	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Methylisobutylketone	ND	ND	500	µg/g	0.0	< 20	Acceptable	
3-Methyl-1-butanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Ethylene Glycol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Toluene	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Isobutyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
1-Pentanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Butyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Ethylbenzene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
m,p-Xylene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
o-Xylene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Cumene	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Anisole	ND	ND	500	µg/g	0.0	< 20	Acceptable	
DMSO	ND	ND	500	µg/g	0.0	< 20	Acceptable	
1,2-dimethoxyethane	ND	ND	50	µg/g	0.0	< 20	Acceptable	
Triethylamine	ND	ND	500	µg/g	0.0	< 20	Acceptable	
N,N-dimethylformamide	ND	ND	150	µg/g	0.0	< 20	Acceptable	
N,N-dimethylacetamide	ND	ND	150	µg/g	0.0	< 20	Acceptable	
Pyridine	ND	ND	50	µg/g	0.0	< 20	Acceptable	
Sulfolane	ND	ND	50	µg/g	0.0	< 20	Acceptable	
1,2-Dichloroethane	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Chloroform	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Trichloroethylene	ND	ND	1	µg/g	0.0	< 20	Acceptable	
1,1-Dichloroethane	ND	ND	1	µg/g	0.0	< 20	Acceptable	

**Abbreviations**

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

**Units of Measure:**

µg/g- Microgram per gram or ppm





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



**Report Number:** 23-009657/D003.R000  
**Report Date:** 08/21/2023  
**ORELAP#:** OR100028  
**Purchase Order:** 2595910  
**Received:** 08/14/23 16:11

Revision: 3 Document ID: 3120  
Legacy ID: CFL-C21 Worksheet Validated 10/30/2020

**Laboratory Pesticide Quality Control Results**

AOAC 2007.1 & EN 15662		Units: mg/Kg			Batch ID: 2310121			
Method Blank		Laboratory Control Sample						
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes
Abamectin	0.000	< 0.250		0.842	1.000	84.2	50.0	150
Acephate	0.000	< 0.200		0.651	0.800	81.4	60.0	120
Acetaminocyl	0.000	< 1.000		3.416	4.000	85.4	40.0	160
Acetamiprid	0.000	< 0.100		0.335	0.400	83.7	60.0	120
Aldicarb	0.000	< 0.200		0.728	0.800	91.1	60.0	120
Azoxystrobin	0.004	< 0.100		0.339	0.400	84.8	60.0	120
Bifenazate	0.000	< 0.100		0.349	0.400	87.2	60.0	120
Bifenthrin	0.000	< 0.100		0.321	0.400	80.1	50.0	150
Boscalid	0.028	< 0.200		0.671	0.800	83.9	60.0	120
Carbaryl	0.000	< 0.100		0.330	0.400	82.6	60.0	120
Carbofuran	0.000	< 0.100		0.331	0.400	82.8	60.0	120
Chlorantraniliprole	0.000	< 0.100		0.327	0.400	81.9	60.0	120
Chlorfenapyr	0.000	< 0.500		1.875	2.000	93.8	60.0	120
Chlorpyrifos	0.000	< 0.100		0.359	0.400	89.8	60.0	120
Clofentazine	0.000	< 0.100		0.308	0.400	76.9	60.0	120
Cyfluthrin	0.000	< 0.500		1.794	2.000	89.7	50.0	150
Cypermethrin	0.000	< 0.500		1.580	2.000	79.0	50.0	150
Daminozide	0.000	< 0.500		0.674	2.000	33.7	60.0	120
Diazinon	0.000	< 0.100		0.342	0.400	85.6	60.0	120
Dichlorvos	0.000	< 0.500		1.847	2.000	92.4	60.0	120
Dimethoate	0.000	< 0.100		0.339	0.400	84.7	60.0	120
Ethoprophos	0.000	< 0.100		0.337	0.400	84.1	60.0	120
Etofenprox	0.000	< 0.200		0.656	0.800	82.0	50.0	150
Etoxazole	0.003	< 0.100		0.334	0.400	83.6	60.0	120
Fenoxycarb	0.000	< 0.100		0.319	0.400	79.8	60.0	120
Fenpyroximate	0.005	< 0.200		0.651	0.800	81.4	60.0	120
Fipronil	0.000	< 0.200		0.643	0.800	80.4	60.0	120
Fonicamid	0.000	< 0.250		0.907	1.000	90.7	60.0	120
Fludioxonil	0.000	< 0.200		0.710	0.800	88.7	50.0	150
Hexythiazox	0.000	< 0.250		0.834	1.000	83.4	60.0	120
Imazalil	0.000	< 0.100		0.334	0.400	83.5	60.0	120
Imidacloprid	0.005	< 0.200		0.685	0.800	85.6	60.0	120
Kresoxim-methyl	0.000	< 0.200		0.656	0.800	81.9	60.0	120
Malathion	0.000	< 0.100		0.326	0.400	81.5	60.0	120
Metaxalyl	0.000	< 0.100		0.337	0.400	84.3	60.0	120
Methiocarb	0.000	< 0.100		0.331	0.400	82.8	60.0	120
Methomyl	0.000	< 0.200		0.723	0.800	90.4	60.0	120
MGK-264	0.000	< 0.100		0.339	0.400	84.6	50.0	150
Myclobutanil	0.000	< 0.100		0.355	0.400	88.9	60.0	120
Naled	0.000	< 0.250		0.856	1.000	85.6	50.0	150
Oxamyl	0.000	< 0.500		1.678	2.000	83.9	60.0	120
Pacllobutrazole	0.000	< 0.200		0.670	0.800	83.8	60.0	120
Parathion-Methyl	0.000	< 0.100		0.362	0.400	90.5	50.0	150
Permethrin	0.000	< 0.100		0.329	0.400	82.2	50.0	150
Phosmet	0.000	< 0.100		0.345	0.400	86.3	50.0	150
Piperonyl butoxide	0.000	< 0.500		1.750	2.000	87.5	60.0	120
Prallethrin	0.000	< 0.100		0.358	0.400	89.4	60.0	120
Propiconazole	0.000	< 0.200		0.612	0.800	76.5	60.0	120
Propoxur	0.000	< 0.100		0.332	0.400	82.9	60.0	120
Pyrethrin (Summe)	0.009	< 0.100		0.420	0.488	86.0	60.0	120
Pyridaben	0.000	< 0.100		0.328	0.400	81.9	50.0	150
Spinosad	0.000	< 0.100		0.320	0.388	82.6	50.0	150
Spiromesifen	0.000	< 0.100		0.350	0.400	87.4	60.0	120
Spirotetramat	0.000	< 0.100		0.334	0.400	83.5	60.0	120
Spiroxamine	0.000	< 0.200		0.684	0.800	85.5	60.0	120
Tebuconazole	0.000	< 0.200		0.613	0.800	76.7	60.0	120
Thiacloprid	0.000	< 0.100		0.318	0.400	79.4	60.0	120
Thiamethoxam	0.000	< 0.100		0.363	0.400	90.6	60.0	120
Trifloxystrobin	0.000	< 0.100		0.335	0.400	83.8	60.0	120

Q6



12423 NE Whitaker Way  
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Revision: 3 Document ID: 3120  
Legacy ID: CFL-C21 Worksheet Validated 10/30/2020

Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662		Units: mg/Kg					Batch ID: 2310121			
Matrix Spike/Matrix Spike Duplicate Recoveries		Sample ID: 23-009707-0001								
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes
Abamectin	0.000	0.805	0.729	1.000	9.8%	< 30	80.5%	72.9%	50 - 150	
Acephate	0.019	0.659	0.703	0.800	6.7%	< 30	80.0%	85.5%	50 - 150	
Acetaminophen	0.000	3.355	3.339	4.000	0.5%	< 30	83.9%	83.5%	50 - 150	
Acetamiprid	0.000	0.339	0.325	0.400	4.1%	< 30	84.7%	81.3%	50 - 150	
Aldicarb	0.000	0.732	0.700	0.800	4.6%	< 30	91.6%	87.4%	50 - 150	
Azoxystrobin	0.004	0.343	0.320	0.400	6.8%	< 30	84.8%	79.2%	50 - 150	
Bifenazate	0.000	0.342	0.338	0.400	1.3%	< 30	85.6%	84.5%	50 - 150	
Bifenthrin	0.000	0.311	0.309	0.400	0.7%	< 30	77.8%	77.2%	50 - 150	
Boscalid	0.024	0.620	0.636	0.800	2.7%	< 30	74.4%	76.4%	50 - 150	
Carbaryl	0.000	0.342	0.326	0.400	4.8%	< 30	85.4%	81.4%	50 - 150	
Carbofuran	0.000	0.331	0.313	0.400	5.6%	< 30	82.8%	78.3%	50 - 150	
Chlorantraniliprole	0.000	0.313	0.293	0.400	6.5%	< 30	78.3%	73.4%	50 - 150	
Chlorfenapyr	0.000	1.360	1.537	2.000	12.3%	< 30	68.0%	76.9%	50 - 150	
Chlorpyrifos	0.000	0.353	0.356	0.400	0.9%	< 30	88.2%	89.0%	50 - 150	
Clofentezine	0.000	0.090	0.103	0.400	13.6%	< 30	22.5%	25.8%	50 - 150	Q
Cyfluthrin	0.000	1.771	1.702	2.000	4.0%	< 30	88.6%	85.1%	30 - 150	
Cypermethrin	0.000	1.702	1.556	2.000	9.0%	< 30	85.1%	77.8%	50 - 150	
Daminozide	0.000	0.684	0.652	2.000	4.8%	< 30	34.2%	32.6%	30 - 150	
Diazinon	0.000	0.328	0.306	0.400	6.9%	< 30	82.0%	76.5%	50 - 150	
Dichlorvos	0.000	1.794	1.739	2.000	3.1%	< 30	89.7%	87.0%	50 - 150	
Dimethoate	0.000	0.342	0.315	0.400	8.2%	< 30	85.5%	78.8%	50 - 150	
Ethoprophos	0.001	0.335	0.318	0.400	5.3%	< 30	83.5%	79.1%	50 - 150	
Etofenprox	0.000	0.628	0.605	0.800	3.6%	< 30	78.5%	75.7%	50 - 150	
Etoxazole	0.004	0.318	0.306	0.400	3.7%	< 30	78.4%	75.6%	50 - 150	
Fenoxycarb	0.000	0.336	0.311	0.400	7.7%	< 30	83.9%	77.6%	50 - 150	
Fenpyroximate	0.005	0.688	0.649	0.800	5.9%	< 30	85.4%	80.5%	50 - 150	
Fipronil	0.000	0.693	0.636	0.800	8.5%	< 30	86.6%	79.5%	50 - 150	
Fonicamid	0.000	0.884	0.780	1.000	12.4%	< 30	88.4%	78.0%	50 - 150	
Fludioxonil	0.000	0.696	0.668	0.800	4.0%	< 30	87.0%	83.6%	50 - 150	
Hexythiazox	0.000	0.782	0.746	1.000	4.7%	< 30	78.2%	74.6%	50 - 150	
Imazalil	0.000	0.341	0.324	0.400	5.0%	< 30	85.2%	81.0%	50 - 150	
Imidacloprid	0.005	0.698	0.609	0.800	13.7%	< 30	86.6%	75.5%	50 - 150	
Kresoxim-methyl	0.000	0.667	0.640	0.800	4.2%	< 30	83.4%	80.0%	50 - 150	
Malathion	0.000	0.326	0.308	0.400	5.5%	< 30	81.4%	77.0%	50 - 150	
Metaxalyl	0.000	0.348	0.330	0.400	5.3%	< 30	87.0%	82.5%	50 - 150	
Methiocarb	0.000	0.340	0.332	0.400	2.2%	< 30	84.9%	83.1%	50 - 150	
Methomyl	0.000	0.680	0.642	0.800	5.9%	< 30	85.0%	80.2%	50 - 150	
MGK-264	0.000	0.322	0.320	0.400	0.5%	< 30	80.5%	80.1%	50 - 150	
Myclobutanil	0.000	0.336	0.294	0.400	13.5%	< 30	84.1%	73.5%	50 - 150	
Naled	0.000	0.880	0.847	1.000	3.8%	< 30	88.0%	84.7%	50 - 150	
Oxamyl	0.000	1.828	1.494	2.000	20.1%	< 30	91.4%	74.7%	50 - 150	
Pacllobutrazole	0.001	0.685	0.638	0.800	7.1%	< 30	85.5%	79.7%	50 - 150	
Parathion-Methyl	0.000	0.366	0.296	0.400	21.3%	< 30	91.5%	73.9%	30 - 150	
Permethrin	0.000	0.302	0.302	0.400	0.2%	< 30	75.6%	75.4%	50 - 150	
Phosmet	0.000	0.358	0.342	0.400	4.5%	< 30	89.6%	85.6%	50 - 150	
Piperonyl butoxide	0.000	1.658	1.620	2.000	2.3%	< 30	82.9%	81.0%	50 - 150	
Prallethrin	0.011	0.353	0.358	0.400	1.7%	< 30	85.3%	86.8%	50 - 150	
Propiconazole	0.000	0.647	0.593	0.800	8.7%	< 30	80.8%	74.1%	50 - 150	
Propoxur	0.000	0.343	0.327	0.400	5.1%	< 30	85.9%	81.6%	50 - 150	
Pyrethrin (Summe)	0.001	0.315	0.308	0.488	2.4%	< 30	64.4%	62.9%	50 - 150	
Pyridaben	0.000	0.293	0.279	0.400	5.0%	< 30	73.2%	69.7%	50 - 150	
Spinosad	0.000	0.314	0.300	0.388	4.5%	< 30	81.0%	77.4%	50 - 150	
Spiromesifen	0.000	0.348	0.322	0.400	7.6%	< 30	87.0%	80.6%	50 - 150	
Spirotetramat	0.000	0.335	0.318	0.400	5.1%	< 30	83.7%	79.5%	50 - 150	
Spiroxamine	0.000	0.700	0.678	0.800	3.1%	< 30	87.4%	84.7%	50 - 150	
Tebuconazole	0.000	0.630	0.573	0.800	9.4%	< 30	78.7%	71.7%	50 - 150	
Thiacloprid	0.000	0.325	0.305	0.400	6.5%	< 30	81.4%	76.2%	50 - 150	
Thiamethoxam	0.000	0.357	0.329	0.400	8.0%	< 30	89.2%	82.3%	50 - 150	
Trifloxystrobin	0.000	0.324	0.305	0.400	6.0%	< 30	81.1%	76.4%	50 - 150	



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**Purchase Order:** 2595910  
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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: 2310135

Laboratory Control Sample									
Analyte	LCS	Result	Spike	Units	% Rec	Limits		Evaluation	Notes
CBDVA	2	0.0306	0.0309	%	99.1	80.0	- 120	Acceptable	
CBDV	2	0.0310	0.0313	%	99.0	80.0	- 120	Acceptable	
CBE	2	0.0322	0.0329	%	97.8	80.0	- 120	Acceptable	
CBDA	1	0.0335	0.0338	%	99.0	90.0	- 110	Acceptable	
CBGA	1	0.0343	0.0343	%	100	80.0	- 120	Acceptable	
CBG	1	0.0361	0.0363	%	99.3	80.0	- 120	Acceptable	
CBD	1	0.0346	0.0351	%	98.4	90.0	- 110	Acceptable	
THCV	2	0.0198	0.0200	%	99.0	80.0	- 120	Acceptable	
d8THCV	2	0.0273	0.0276	%	99.0	80.0	- 120	Acceptable	
THCVA	2	0.0302	0.0307	%	98.4	80.0	- 120	Acceptable	
CBN	1	0.0343	0.0343	%	99.8	80.0	- 120	Acceptable	
exo-THC	2	0.0297	0.0302	%	98.3	80.0	- 120	Acceptable	
d9THC	1	0.0352	0.0355	%	99.0	90.0	- 110	Acceptable	
d8THC	1	0.0368	0.0364	%	101	90.0	- 110	Acceptable	
9S-d10THC	1	0.0354	0.0354	%	99.9	80.0	- 120	Acceptable	
CBL	2	0.0314	0.0311	%	101	80.0	- 120	Acceptable	
9R-d10THC	1	0.0112	0.0115	%	97.4	80.0	- 120	Acceptable	
CBC	2	0.0327	0.0335	%	97.7	80.0	- 120	Acceptable	
THCA	1	0.0342	0.0344	%	99.5	90.0	- 110	Acceptable	
CBCA	2	0.0318	0.0319	%	99.7	80.0	- 120	Acceptable	
CBLA	2	0.0641	0.0647	%	99.1	80.0	- 120	Acceptable	
d9THCP	2	0.0308	0.0316	%	97.5	80.0	- 120	Acceptable	
CBT	2	0.0300	0.0308	%	97.4	80.0	- 120	Acceptable	

Method Blank						
Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBDV	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBE	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBDA	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBGA	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBG	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBD	<LOQ	0.00322	%	< 0.00322	Acceptable	
THCV	<LOQ	0.00322	%	< 0.00322	Acceptable	
d8THCV	<LOQ	0.00322	%	< 0.00322	Acceptable	
THCVA	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBN	<LOQ	0.00322	%	< 0.00322	Acceptable	
exo-THC	<LOQ	0.00322	%	< 0.00322	Acceptable	
d9THC	<LOQ	0.00322	%	< 0.00322	Acceptable	
d8THC	<LOQ	0.00322	%	< 0.00322	Acceptable	
9S-d10THC	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBL	<LOQ	0.00322	%	< 0.00322	Acceptable	
9R-d10THC	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBC	<LOQ	0.00322	%	< 0.00322	Acceptable	
THCA	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBCA	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBLA	<LOQ	0.00322	%	< 0.00322	Acceptable	
d9THCP	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBT	<LOQ	0.00322	%	< 0.00322	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:** 23-009657/D003.R000  
**Report Date:** 08/21/2023  
**ORELAP#:** OR100028  
**Purchase Order:** 2595910  
**Received:** 08/14/23 16:11

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

Laboratory Quality Control Results

J AOAC 2015 V98-6		Batch ID: 2310135						
Sample Duplicate		Sample ID: 23-009657-0001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
CBDV	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
CBE	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
CBDA	0.0676	0.0669	0.00326	%	1.13	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
CBG	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
CBD	0.227	0.226	0.00326	%	0.285	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
CBN	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
d9THC	0.204	0.204	0.00326	%	0.0561	< 20	Acceptable	
d8THC	<LOQ	0.00336	0.00326	%	NA	< 20	Acceptable	R2
9S-d10THC	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
9R-d10THC	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
CBC	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
CBCA	0.00369	0.00377	0.00326	%	2.12	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
d9THCP	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	
CBT	<LOQ	<LOQ	0.00326	%	NA	< 20	Acceptable	

Abbreviations

- ND - None Detected at or above MRL
- RPD - Relative Percent Difference
- LOQ - Limit of Quantitation
- R2 - Sample replicates RPD non-calculable, as only one replicate is within analytical range.

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.