

## CONSOLIDATED TEST RESULTS SUMMARY

Please see the following pages for full test results.

<b>BULK SKU</b> BEV.D9.IT50.4PK	<b>BATCH #</b> FK37-1C	<b>LOQ:</b> Limit Of Quantitation <b>LOD:</b> Limit Of Detection  $1\text{ g} = 10^{-3}\text{ kg} = 10^3\text{ mg} = 10^6\text{ }\mu\text{g}$ $1\text{ mg/kg} = 1\text{ ppm} = 1000\text{ ppb}$
<b>PRODUCT NAME</b> Sparkling Iced Tea THC Lemonade	<b>SERVING SIZE</b> 1 Can (473mL)	
<b>LABORATORY:</b> Columbia Laboratories	<b>OREGON ACCREDITATION:</b> OR100028	

POTENCY	PER SERVING	PER GRAM	Percent
Cannabidiol (CBD)	43.4 mg/serving	0.09 mg/g	0.01 %
Total THC (d9-THC, THCA)	<b>44.3 mg/serving</b>	0.094 mg/g	0.01 %
Cannabigerol (CBG)	<LO mg/serving	<LO mg/g	<LO %
Cannabinol (CBN)	<LO mg/serving	<LO mg/g	<LO %
Cannabichromene (CBC)	<LO mg/serving	<LO mg/g	<LO %
Tetrahydrocannabinolic Acid (THCA)	<LO mg/serving	<LO mg/g	<LO %
Delta-9-THC (d9-THC)	<b>44.3 mg/serving</b>	0.094 mg/g	0.01 %
Delta-8-THC (d8-THC)	<b>14.1 mg/serving</b>	0.030 mg/g	0.00 %

HEAVY METALS	PER SERVING	PER GRAM	REGULATORY ACTION LEVEL
Arsenic	<b>2.48</b> $\mu\text{g/serving}$	0.005 $\mu\text{g/g}$	1.5 ppm
Cadmium	<LO $\mu\text{g/serving}$	<LO $\mu\text{g/g}$	0.5 ppm
Lead	<LO $\mu\text{g/serving}$	<LO $\mu\text{g/g}$	0.5 ppm
Mercury	<LO $\mu\text{g/serving}$	<LO $\mu\text{g/g}$	3.0 ppm

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*	375 $\mu\text{g/g}$	5,000 ppm
Heptane	<LO $\mu\text{g/g}$	5,000 ppm

None of the 34 residual solvents tested found above limit of quantitation in the sample.

MICROBIAL	PASS/FAIL
Yeast & Mold	Pass
Coliform	Pass



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:** 24-000281/D023.R000  
**Report Date:** 01/19/2024  
**ORELAP#:** OR100028  
**Purchase Order:** 2797832  
**Received:** 01/08/24 16:15

**Customer:** Etz Hayim Holdings  
**Product identity:** CYCL-BEV.D9.IT50.4PK-FK37-1C  
**Client/Metric ID:** .  
**Laboratory ID:** 24-000281-0003

### Summary

**Potency:**

Analyte per 1g	Result	Limits	Units	Status	
CBD per 1g	0.0918		mg/1g		THC-Total per Serving Size 0.0938 mg/1g
Δ8-THC per 1g	0.0300		mg/1g		
Δ9-THC per 1g	0.0938		mg/1g		CBD-Total per Serving Size 0.0918 mg/1g
(Reported in milligrams per serving)					

**Residual Solvents:**

Analyte	Result (µg/g)	Limits (µg/g)	Status
Ethanol	375		

**Pesticides:**

*All analytes passing and less than LOQ.*

**Metals:**

Analyte	Result	Units	Limit	Status
Arsenic*	0.00525	mg/kg	0.200	pass

**Microbiology:**

*Less than LOQ for all analytes.*



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**Purchase Order:** 2797832  
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**Customer:** Etz Hayim Holdings  
 16427 NE Airport Way  
 PORTLAND 97230  
 United States of America (USA)

**Product identity:** CYCL-BEV.D9.IT50.4PK-FK37-1C

**Client/Metric ID:** .

**Sample Date:**

**Laboratory ID:** 24-000281-0003

**Evidence of Cooling:** No

**Temp:** 17.8 °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: 2400305	Analyze: 1/10/24 8:17:00 PM	
Analyte	Result	Limits	Units	LOQ	Notes
CBC per 1g	< LOQ		mg/1g	0.000998	
CBC-A per 1g	< LOQ		mg/1g	0.000998	
CBC-Total per 1g	< LOQ		mg/1g	0.00187	
CBD per 1g	0.0918		mg/1g	0.000998	
CBD-A per 1g	< LOQ		mg/1g	0.000998	
CBD-Total per 1g	0.0918		mg/1g	0.00187	
CBDV per 1g	< LOQ		mg/1g	0.000998	
CBDV-A per 1g	< LOQ		mg/1g	0.000998	
CBDV-Total per 1g	< LOQ		mg/1g	0.00186	
CBE per 1g	< LOQ		mg/1g	0.000998	
CBG per 1g	< LOQ		mg/1g	0.000998	
CBG-A per 1g	< LOQ		mg/1g	0.000998	
CBG-Total per 1g	< LOQ		mg/1g	0.00186	
CBL per 1g	< LOQ		mg/1g	0.000998	
CBL-A per 1g	< LOQ		mg/1g	0.000998	
CBL-Total per 1g	< LOQ		mg/1g	0.00187	
CBN per 1g	< LOQ		mg/1g	0.000998	
CBT per 1g	< LOQ		mg/1g	0.000998	
Δ8-THCV per 1g	< LOQ		mg/1g	0.000998	
Δ10-THC-9R per 1g	< LOQ		mg/1g	0.000998	
Δ10-THC-9S per 1g	< LOQ		mg/1g	0.000998	
Δ10-THC-Total per 1g	< LOQ		mg/1g	0.00200	
Δ8-THC per 1g	0.0300		mg/1g	0.000998	
Δ9-THC per 1g	0.0938		mg/1g	0.000998	
delta-9-THCP per 1g	< LOQ		mg/1g	0.000998	
exo-THC per 1g	< LOQ		mg/1g	0.000998	
THC-A per 1g	< LOQ		mg/1g	0.000998	
THC-Total per 1g	0.0938		mg/1g	0.00187	
THCV per 1g	< LOQ		mg/1g	0.000998	
THCV-A per 1g	< LOQ		mg/1g	0.000998	



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Potency per 1g						
Analyte	Method: J AOAC 2015 V98-6 (mod) <sup>P</sup>	Result	Limits	Units mg/se	Batch: 2400305	Analyze: 1/10/24 8:17:00 PM
THCV-Total per 1g		< LOQ		mg/1g		0.00187
Total Cannabinoids per 1g		0.216		mg/1g		

Microbiology							
Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status Notes
E.coli	< LOQ		cfu/g	10	2400247	01/12/24 AOAC 991.14 (Petrifilm) <sup>P</sup>	
Total Coliforms	< LOQ		cfu/g	10	2400247	01/12/24 AOAC 991.14 (Petrifilm) <sup>P</sup>	
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2400248	01/13/24 AOAC 2014.05 (RAPID) <sup>P</sup>	
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2400248	01/13/24 AOAC 2014.05 (RAPID) <sup>P</sup>	

Solvents											
Method: Residual Solvents by GC/MS <sup>P</sup>						Units µg/g		Batch 2400422		Analyze 01/19/24 01:50 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	375		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 2400371		Analyze 01/15/24 02:49 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		Acetamidrid <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		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		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>	0.00525	0.200	mg/kg	0.00389	2400340	01/12/24	AOAC 2013.06 (mod.) <sup>b</sup>	pass		
Cadmium <sup>¥</sup>	< LOQ	0.200	mg/kg	0.00389	2400340	01/12/24	AOAC 2013.06 (mod.) <sup>b</sup>	pass		
Lead <sup>¥</sup>	< LOQ	0.500	mg/kg	0.00389	2400340	01/12/24	AOAC 2013.06 (mod.) <sup>b</sup>	pass		
Mercury <sup>¥</sup>	< LOQ	0.100	mg/kg	0.00195	2400340	01/12/24	AOAC 2013.06 (mod.) <sup>b</sup>	pass		



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**Received:** 01/08/24 16:15

**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

J AOAC 2015 V98-6 Batch ID: 2400305

Laboratory Control Sample									
Analyte	LCS	Result	Spike	Units	% Rec	Limits		Evaluation	Notes
CBDVA	2	0.0009	0.0010	%	94.4	80.0	- 120	Acceptable	
CBDV	2	0.00108	0.00104	%	104	80.0	- 120	Acceptable	
CBE	2	0.00106	0.00105	%	100	80.0	- 120	Acceptable	
CBDA	1	0.0009	0.0009	%	90.1	90.0	- 110	Acceptable	
CBGA	1	0.0009	0.0009	%	92.2	80.0	- 120	Acceptable	
CBG	1	0.0009	0.0009	%	94.7	80.0	- 120	Acceptable	
CBD	1	0.00100	0.0010	%	101	90.0	- 110	Acceptable	
THCV	2	0.00104	0.00104	%	101	80.0	- 120	Acceptable	
d8THCV	2	0.0009	0.0009	%	101	80.0	- 120	Acceptable	
THCVA	2	0.0009	0.0010	%	93.1	80.0	- 120	Acceptable	
CBN	1	0.00100	0.0010	%	101	80.0	- 120	Acceptable	
exo-THC	2	0.0010	0.0010	%	99.8	80.0	- 120	Acceptable	
d9THC	1	0.00104	0.00101	%	103	90.0	- 110	Acceptable	
d8THC	1	0.00103	0.00101	%	102	90.0	- 110	Acceptable	
9S-d10THC	1	0.0010	0.0010	%	101	80.0	- 120	Acceptable	
CBL	2	0.0010	0.0010	%	101	80.0	- 120	Acceptable	
9R-d10THC	1	0.0010	0.0010	%	100	80.0	- 120	Acceptable	
CB	2	0.00105	0.00105	%	99.9	80.0	- 120	Acceptable	
THCA	1	0.0009	0.0010	%	88.8	90.0	- 110	Acceptable	Q6
CBCA	2	0.0009	0.00100	%	90.0	80.0	- 120	Acceptable	
CBLA	2	0.0010	0.00102	%	93.7	80.0	- 120	Acceptable	
d9THCP	2	0.0010	0.0010	%	100	80.0	- 120	Acceptable	
CBT	2	0.00103	0.00104	%	99.4	80.0	- 120	Acceptable	

Method Blank						
Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBDV	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBE	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBDA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBGA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBG	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBD	<LOQ	0.0001	%	< 0.0001	Acceptable	
THCV	<LOQ	0.0001	%	< 0.0001	Acceptable	
d8THCV	<LOQ	0.0001	%	< 0.0001	Acceptable	
THCVA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBN	<LOQ	0.0001	%	< 0.0001	Acceptable	
exo-THC	<LOQ	0.0001	%	< 0.0001	Acceptable	
d9THC	<LOQ	0.0001	%	< 0.0001	Acceptable	
d8THC	<LOQ	0.0001	%	< 0.0001	Acceptable	
9S-d10THC	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBL	<LOQ	0.0001	%	< 0.0001	Acceptable	
9R-d10THC	<LOQ	0.0001	%	< 0.0001	Acceptable	
CB	<LOQ	0.0001	%	< 0.0001	Acceptable	
THCA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBCA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBLA	<LOQ	0.0001	%	< 0.0001	Acceptable	
d9THCP	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBT	<LOQ	0.0001	%	< 0.0001	Acceptable	

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

**Units of Measure:**  
 % - Percent



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Laboratory Quality Control Results

J AOAC 2015 V98-6		Batch ID: 2400305						
Sample Duplicate		Sample ID: 24-000278-0001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBDV	0.0002	0.0002	0.0001	%	0.0216	< 20	Acceptable	
CBE	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBDA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBG	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBD	0.0153	0.0152	0.0001	%	0.417	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBN	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
d9THC	0.00809	0.00805	0.0001	%	0.465	< 20	Acceptable	
d8THC	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
9S-d10THC	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
9R-d10THC	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBC	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBCA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
d9THCP	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBT	<LOQ	<LOQ	0.0001	%	NA	< 20	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:** 24-000281/D023.R000  
**Report Date:** 01/19/2024  
**ORELAP#:** OR100028  
**Purchase Order:** 2797832  
**Received:** 01/08/24 16:15

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: 2400371			
Method Blank			Laboratory Control Sample					
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes
Abamectin	0.000	< 0.250		0.920	1.000	92.0	50.0 150	
Acephate	0.014	< 0.200		0.655	0.800	81.9	60.0 120	
Acequinocyl	0.000	< 1.000		4.015	4.000	100.4	40.0 160	
Acetamiprid	0.000	< 0.100		0.351	0.400	87.8	60.0 120	
Aldicarb	0.000	< 0.200		0.749	0.800	93.6	60.0 120	
Azoxystrobin	0.005	< 0.100		0.369	0.400	92.2	60.0 120	
Bifenazate	0.000	< 0.100		0.370	0.400	92.5	60.0 120	
Bifenthrin	0.000	< 0.100		0.356	0.400	89.1	50.0 150	
Boscalid	0.000	< 0.200		0.719	0.800	89.9	60.0 120	
Carbaryl	0.000	< 0.100		0.357	0.400	89.4	60.0 120	
Carbofuran	0.000	< 0.100		0.353	0.400	88.2	60.0 120	
Chlorantraniliprole	0.000	< 0.100		0.347	0.400	86.7	60.0 120	
Chlorfenapyr	0.029	< 0.500		2.013	2.000	100.6	60.0 120	
Chlorpyrifos	0.000	< 0.100		0.338	0.400	84.5	60.0 120	
Clofentezine	0.000	< 0.100		0.325	0.400	81.3	60.0 120	
Cyfluthrin	0.000	< 0.500		1.837	2.000	91.9	50.0 150	
Cypermethrin	0.000	< 0.500		1.816	2.000	90.8	50.0 150	
Daminozide	0.000	< 0.500		0.631	2.000	31.6	60.0 120	Q6
Diazinon	0.000	< 0.100		0.375	0.400	93.7	60.0 120	
Dichlorvos	0.000	< 0.500		1.586	2.000	79.3	60.0 120	
Dimethoate	0.000	< 0.100		0.351	0.400	87.7	60.0 120	
Ethoprophos	0.000	< 0.100		0.341	0.400	85.4	60.0 120	
Etofenprox	0.000	< 0.200		0.753	0.800	94.1	50.0 150	
Etoxazole	0.000	< 0.100		0.389	0.400	97.2	60.0 120	
Fenoxycarb	0.000	< 0.100		0.367	0.400	91.8	60.0 120	
Fenpyroximate	0.000	< 0.200		0.737	0.800	92.1	60.0 120	
Fipronil	0.000	< 0.200		0.730	0.800	91.2	60.0 120	
Flonicamid	0.000	< 0.250		0.678	1.000	67.8	60.0 120	
Fludioxonil	0.000	< 0.200		0.681	0.800	85.1	50.0 150	
Hexythiazox	0.000	< 0.250		0.927	1.000	92.7	60.0 120	
Imazalil	0.000	< 0.100		0.339	0.400	84.7	60.0 120	
Imidacloprid	0.000	< 0.200		0.597	0.800	74.6	60.0 120	
Kresoxim-methyl	0.000	< 0.200		0.713	0.800	89.1	60.0 120	
Malathion	0.000	< 0.100		0.362	0.400	90.5	60.0 120	
Metalaxyl	0.000	< 0.100		0.351	0.400	87.7	60.0 120	
Methiocarb	0.000	< 0.100		0.366	0.400	91.4	60.0 120	
Methomyl	0.000	< 0.200		0.555	0.800	69.3	60.0 120	
MGK-264	0.000	< 0.100		0.362	0.400	90.6	50.0 150	
Myclobutanil	0.000	< 0.100		0.333	0.400	83.3	60.0 120	
Naled	0.000	< 0.250		0.848	1.000	84.8	50.0 150	
Oxamyl	0.000	< 0.500		1.490	2.000	74.5	60.0 120	
Paclobutrazole	0.000	< 0.200		0.711	0.800	88.9	60.0 120	
Parathion-Methyl	0.024	< 0.100		0.359	0.400	89.7	50.0 150	
Permethrin	0.000	< 0.100		0.366	0.400	91.5	50.0 150	
Phosmet	0.000	< 0.100		0.374	0.400	93.4	50.0 150	
Piperonyl butoxide	0.000	< 0.500		1.821	2.000	91.1	60.0 120	
Prallethrin	0.000	< 0.100		0.349	0.400	87.3	60.0 120	
Propiconazole	0.000	< 0.200		0.730	0.800	91.2	60.0 120	
Propoxur	0.000	< 0.100		0.346	0.400	86.6	60.0 120	
Pyrethrin (Summe)	0.000	< 0.100		0.455	0.488	93.2	60.0 120	
Pyridaben	0.000	< 0.100		0.357	0.400	89.3	50.0 150	
Spinosad	0.000	< 0.100		0.353	0.388	90.9	50.0 150	
Spiromesifen	0.000	< 0.100		0.360	0.400	90.0	60.0 120	
Spirotetramat	0.000	< 0.100		0.339	0.400	84.6	60.0 120	
Spiroxamine	0.000	< 0.200		0.716	0.800	89.5	60.0 120	
Tebuconazole	0.000	< 0.200		0.712	0.800	89.0	60.0 120	
Thiacloprid	0.000	< 0.100		0.359	0.400	89.7	60.0 120	
Thiamethoxam	0.000	< 0.100		0.304	0.400	75.9	60.0 120	
Trifloxystrobin	0.000	< 0.100		0.345	0.400	86.2	60.0 120	



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**Report Number:** 24-000281/D023.R000  
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**Purchase Order:** 2797832  
**Received:** 01/08/24 16:15

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: 2400371				
Matrix Spike/Matrix Spike Duplicate Recoveries					Sample ID: 24-000281-0003					
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes
Abamectin	0.000	0.877	0.900	1.000	2.6%	< 30	87.7%	90.0%	50 - 150	
Acephate	0.000	0.659	0.652	0.800	1.0%	< 30	82.3%	81.5%	50 - 150	
Acequinocyl	0.000	4.924	3.673	4.000	29.1%	< 30	123.1%	91.8%	50 - 150	
Acetamiprid	0.000	0.378	0.349	0.400	8.1%	< 30	94.6%	87.3%	50 - 150	
Aldicarb	0.000	0.748	0.746	0.800	0.2%	< 30	93.5%	93.3%	50 - 150	
Azoxystrobin	0.005	0.351	0.319	0.400	9.8%	< 30	86.5%	78.5%	50 - 150	
Bifenazate	0.000	0.357	0.344	0.400	3.7%	< 30	89.3%	86.1%	50 - 150	
Bifenthrin	0.000	0.310	0.339	0.400	8.8%	< 30	77.6%	84.7%	50 - 150	
Boscalid	0.000	0.763	0.697	0.800	9.1%	< 30	95.4%	87.1%	50 - 150	
Carbaryl	0.000	0.363	0.320	0.400	12.6%	< 30	90.9%	80.1%	50 - 150	
Carbofuran	0.000	0.328	0.311	0.400	5.3%	< 30	82.1%	77.8%	50 - 150	
Chlorantraniliprole	0.000	0.330	0.332	0.400	0.7%	< 30	82.5%	83.1%	50 - 150	
Chlorfenapyr	0.028	1.727	1.810	2.000	4.8%	< 30	85.0%	89.1%	50 - 150	
Chlorpyrifos	0.000	0.325	0.310	0.400	4.5%	< 30	81.2%	77.6%	50 - 150	
Clofentezine	0.000	0.298	0.281	0.400	6.1%	< 30	74.6%	70.1%	50 - 150	
Cyfluthrin	0.000	1.953	1.960	2.000	0.4%	< 30	97.7%	98.0%	30 - 150	
Cypermethrin	0.000	1.920	1.963	2.000	2.2%	< 30	96.0%	98.2%	50 - 150	
Daminozide	0.000	0.616	0.688	2.000	11.0%	< 30	30.8%	34.4%	30 - 150	
Diazinon	0.000	0.346	0.362	0.400	4.7%	< 30	86.4%	90.6%	50 - 150	
Dichlorvos	0.000	1.694	1.617	2.000	4.6%	< 30	84.7%	80.9%	50 - 150	
Dimethoate	0.000	0.353	0.342	0.400	3.1%	< 30	88.2%	85.5%	50 - 150	
Ethoprofos	0.000	0.338	0.338	0.400	0.1%	< 30	84.5%	84.6%	50 - 150	
Etofenprox	0.000	0.631	0.677	0.800	7.0%	< 30	78.8%	84.6%	50 - 150	
Etoxazole	0.000	0.369	0.370	0.400	0.5%	< 30	92.2%	92.6%	50 - 150	
Fenoxycarb	0.000	0.339	0.340	0.400	0.3%	< 30	84.8%	85.0%	50 - 150	
Fenpyroximate	0.000	0.706	0.727	0.800	2.9%	< 30	88.2%	90.9%	50 - 150	
Fipronil	0.000	0.767	0.712	0.800	7.4%	< 30	95.8%	89.0%	50 - 150	
Flonicamid	0.000	0.756	0.846	1.000	11.2%	< 30	75.6%	84.6%	50 - 150	
Fludioxonil	0.000	0.682	0.601	0.800	12.7%	< 30	85.3%	75.1%	50 - 150	
Hexythiazox	0.000	1.485	1.453	1.000	2.2%	< 30	148.5%	145.3%	50 - 150	
Imazalil	0.000	0.345	0.338	0.400	2.1%	< 30	86.2%	84.4%	50 - 150	
Imidacloprid	0.000	0.649	0.651	0.800	0.3%	< 30	81.1%	81.3%	50 - 150	
Kresoxim-methyl	0.000	0.700	0.645	0.800	8.2%	< 30	87.5%	80.6%	50 - 150	
Malathion	0.000	0.321	0.331	0.400	2.9%	< 30	80.3%	82.7%	50 - 150	
Metalaxyl	0.000	0.347	0.330	0.400	5.0%	< 30	86.7%	82.5%	50 - 150	
Methiocarb	0.000	0.370	0.352	0.400	4.9%	< 30	92.4%	88.0%	50 - 150	
Methomyl	0.000	0.605	0.694	0.800	13.8%	< 30	75.6%	86.8%	50 - 150	
MGK-264	0.000	0.333	0.311	0.400	7.0%	< 30	83.3%	77.6%	50 - 150	
Myclobutanil	0.000	0.319	0.347	0.400	8.5%	< 30	79.7%	86.8%	50 - 150	
Naled	0.000	0.796	0.800	1.000	0.5%	< 30	79.6%	80.0%	50 - 150	
Oxamyl	0.000	1.584	1.651	2.000	4.2%	< 30	79.2%	82.6%	50 - 150	
Paclobutrazole	0.000	0.707	0.626	0.800	12.2%	< 30	88.4%	78.2%	50 - 150	
Parathion-Methyl	0.024	0.350	0.366	0.400	4.9%	< 30	81.5%	85.6%	30 - 150	
Permethrin	0.000	0.312	0.334	0.400	6.8%	< 30	78.0%	83.5%	50 - 150	
Phosmet	0.000	0.367	0.350	0.400	4.8%	< 30	91.8%	87.5%	50 - 150	
Piperonyl butoxide	0.000	1.734	1.643	2.000	5.4%	< 30	86.7%	82.2%	50 - 150	
Prallethrin	0.000	0.344	0.342	0.400	0.7%	< 30	86.0%	85.4%	50 - 150	
Propiconazole	0.000	0.710	0.699	0.800	1.6%	< 30	88.8%	87.4%	50 - 150	
Propoxur	0.000	0.332	0.332	0.400	0.0%	< 30	83.0%	83.1%	50 - 150	
Pyrethrin (Summe)	0.000	0.431	0.433	0.488	0.5%	< 30	88.2%	88.7%	50 - 150	
Pyridaben	0.000	0.374	0.396	0.400	5.8%	< 30	93.4%	99.0%	50 - 150	
Spinosad	0.000	0.348	0.324	0.388	7.2%	< 30	89.7%	83.4%	50 - 150	
Spiromesfen	0.000	0.358	0.359	0.400	0.4%	< 30	89.5%	89.8%	50 - 150	
Spirotetramat	0.000	0.348	0.331	0.400	5.2%	< 30	87.1%	82.6%	50 - 150	
Spiroxamine	0.000	0.681	0.694	0.800	1.9%	< 30	85.1%	86.7%	50 - 150	
Tebuconazole	0.000	0.728	0.735	0.800	1.0%	< 30	91.0%	91.9%	50 - 150	
Thiacloprid	0.000	0.358	0.350	0.400	2.3%	< 30	89.6%	87.5%	50 - 150	
Thiamethoxam	0.000	0.338	0.376	0.400	10.7%	< 30	84.5%	94.0%	50 - 150	
Trifloxystrobin	0.000	0.329	0.307	0.400	7.0%	< 30	82.3%	76.7%	50 - 150	



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**Report Number:** 24-000281/D023.R000  
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**Purchase Order:** 2797832  
**Received:** 01/08/24 16:15



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

Laboratory Quality Control Results

Residual Solvents				Batch ID: 2400422					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		603	584	µg/g	103.3	60 - 120	
Isobutane	ND	< 200		795	767	µg/g	103.7	60 - 120	
Butane	ND	< 200		794	782	µg/g	101.5	60 - 120	
2,2-Dimethylpropane	ND	< 200		993	939	µg/g	105.8	60 - 120	
Methanol	ND	< 200		1400	1600	µg/g	87.5	60 - 120	
Ethylene Oxide	ND	< 30		58.7	57.1	µg/g	102.8	60 - 120	
2-Methylbutane	ND	< 200		1400	1600	µg/g	87.5	60 - 120	
Pentane	ND	< 200		1360	1600	µg/g	85.0	60 - 120	
Ethanol	ND	< 200		1330	1600	µg/g	83.1	70 - 130	
Ethyl Ether	ND	< 200		1420	1600	µg/g	88.8	60 - 120	
2,2-Dimethylbutane	ND	< 30		142	161	µg/g	88.2	60 - 120	
Acetone	ND	< 200		1450	1600	µg/g	90.6	60 - 120	
2-Propanol	ND	< 200		1460	1600	µg/g	91.3	60 - 120	
Acetonitrile	ND	< 100		431	488	µg/g	88.3	60 - 120	
2,3-Dimethylbutane	ND	< 30		135	163	µg/g	82.8	60 - 120	
Dichloromethane	ND	< 60		427	488	µg/g	87.5	60 - 120	
2-Methylpentane	ND	< 30		77.2	161	µg/g	48.0	60 - 120	Q6
3-Methylpentane	ND	< 30		140	162	µg/g	86.4	60 - 120	
Hexane	ND	< 30		127	161	µg/g	78.9	60 - 120	
Ethyl acetate	ND	< 200		1420	1610	µg/g	88.2	60 - 120	
2-Butanol	ND	< 200		1430	1610	µg/g	88.8	60 - 120	
Tetrahydrofuran	ND	< 100		432	483	µg/g	89.4	60 - 120	
Cyclohexane	ND	< 200		1410	1600	µg/g	88.1	60 - 120	
Benzene	ND	< 1		2.73	4.99	µg/g	54.7	60 - 120	Q6
Isopropyl Acetate	ND	< 200		1430	1600	µg/g	89.4	60 - 120	
Heptane	ND	< 200		1370	1600	µg/g	85.6	60 - 120	
1,4-Dioxane	ND	< 100		402	480	µg/g	83.8	60 - 120	
2-Ethoxyethanol	ND	< 30		120	161	µg/g	74.5	60 - 120	
Ethylene Glycol	ND	< 200		241	481	µg/g	50.1	60 - 120	Q6
Toluene	ND	< 100		419	483	µg/g	86.7	60 - 120	
Ethylbenzene	ND	< 200		849	962	µg/g	88.3	60 - 120	
m,p-Xylene	ND	< 200		859	972	µg/g	88.4	60 - 120	
o-Xylene	ND	< 200		847	965	µg/g	87.8	60 - 120	
Cumene	ND	< 30		153	169	µg/g	90.5	60 - 120	



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

QC - Sample Duplicate		Sample ID: 24-000281-0001						
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	
Acetonitrile	ND	ND	100	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	
3-Methylpentane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Hexane	ND	ND	30	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	
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	
1,4-Dioxane	ND	ND	100	µg/g	0.0	< 20	Acceptable	
2-Ethoxyethanol	ND	ND	30	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	
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	

**Abbreviations**

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

**Units of Measure:**

µg/g - Microgram per gram or ppm



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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.