



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



**Report Number:** 23-008138/D003.R000  
**Report Date:** 07/18/2023  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 07/11/23 14:30

**Customer:** The Hemp Collect  
**Product identity:** Live CBD Daytrip Extract - Strawberry Haze  
**Client/Metric ID:** 104323192011  
**Laboratory ID:** 23-008138-0007

### Summary

#### Potency:

Analyte	Result (%)			
CBD	40.6			
CBG	8.83			
CBT	6.12			
CBD-A	4.75			
Δ8-THCV	4.19			
CBE	4.14			
CBDV	1.34			
CBDV-A	1.33			
CBC-A	0.364			
Δ8-THC	0.208			
THC-A	0.190			
CBC	0.168			
				CBD-Total 44.8%
				THC-Total 0.167%
			(Reported in percent of total sample)	

#### Residual Solvents:

Analyte	Result (µg/g)	Limits (µg/g)	Status
n-Butane	223		

#### Pesticides:

All analytes passing and less than LOQ.

#### Metals:

Less than LOQ for all analytes.



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**Customer:** The Hemp Collect  
825 NW 16th Ave  
Portland Oregon 97209  
United States of America (USA)

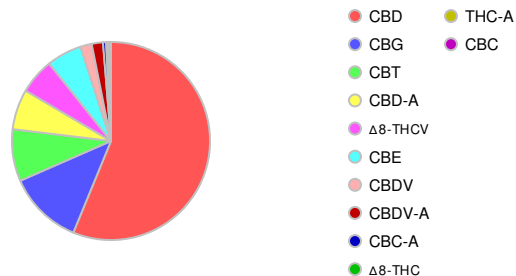
**Product identity:** Live CBD Daytrip Extract - Strawberry Haze  
**Client/Metric ID:** 104323192011  
**Sample Date:**  
**Laboratory ID:** 23-008138-0007  
**Evidence of Cooling:** No  
**Temp:** 22.6 °C  
**Relinquished by:** Hinton



### Sample Results

Potency **Method:** J AOAC 2015 V98-6 (mod)<sup>P</sup> **Units %** **Batch:** 2309035 **Analyze:** 7/12/23 11:06:00 PM

Analyte	As Received	Dry weight	LOQ	Notes
CBC	0.168		0.0676	
CBC-A	0.364		0.0676	
CBC-Total	0.487		0.127	
CBD	40.6		0.676	
CBD-A	4.75		0.0676	
CBD-Total	44.8		0.735	
CBDV	1.34		0.0676	
CBDV-A	1.33		0.0676	
CBDV-Total	2.49		0.126	
CBE	4.14		0.0676	
CBG	8.83		0.0676	
CBG-A	< LOQ		0.0676	
CBG-Total	8.83		0.126	
CBL	< LOQ		0.0676	
CBL-A	< LOQ		0.0676	
CBL-Total	< LOQ		0.127	
CBN	< LOQ		0.0676	
CBT	6.12		0.0676	
Δ10-THC-9R	< LOQ		0.0676	
Δ10-THC-9S	< LOQ		0.0676	
Δ10-THC-Total	< LOQ		0.135	
Δ8-THC	0.208		0.0676	
Δ8-THCV	4.19		0.0676	
Δ9-THC	< LOQ		0.0676	
delta-9-THCP	< LOQ		0.0676	
exo-THC	< LOQ		0.0676	
THC-A	0.190		0.0676	
THC-Total	0.167		0.127	
THCV	< LOQ		0.0676	
THCV-A	< LOQ		0.0676	
THCV-Total	< LOQ		0.126	
<b>Total Cannabinoids</b>	<b>72.2</b>			





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Potency **Method:** J AOAC 2015 V98-6 (mod)<sup>p</sup> **Units %** **Batch:** 2309035 **Analyze:** 7/12/23 11:06:00 PM

**Analyte** **As Received** **Dry weight** **LOQ** **Notes**

Total Xylenes

Solvents **Method:** Residual Solvents by GC/MS<sup>b</sup> **Units µg/g** **Batch** 2309154 **Analyze** 07/17/23 02:28 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		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	223		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) <sup>b</sup>											
Units mg/kg Batch 2309040 Analyze 07/13/23 02:39 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	
Bifentazate <sup>¥</sup>	< LOQ	0.20	0.100	pass		Bifenthrin <sup>¥</sup>	< LOQ	0.20	0.100	pass	
Boscalid <sup>¥</sup>	< LOQ	0.40	0.200	pass		Carbaryl <sup>¥</sup>	< LOQ	0.20	0.100	pass	
Carbofuran <sup>¥</sup>	< LOQ	0.20	0.100	pass		Chlorantraniliprole <sup>¥</sup>	< LOQ	0.20	0.100	pass	
Chlorfenapyr <sup>¥</sup>	< LOQ	1.0	0.500	pass		Chlorpyrifos <sup>¥</sup>	< LOQ	0.20	0.100	pass	
Clofentezine <sup>¥</sup>	< LOQ	0.20	0.100	pass		Cyfluthrin <sup>¥</sup>	< LOQ	1.0	0.500	pass	
Cypermethrin <sup>¥</sup>	< LOQ	1.0	0.500	pass		Daminozide <sup>¥</sup>	< LOQ	1.0	0.500	pass	
Diazinon <sup>¥</sup>	< LOQ	0.20	0.100	pass		Dichlorvos <sup>¥</sup>	< LOQ	1.0	0.500	pass	
Dimethoate <sup>¥</sup>	< LOQ	0.20	0.100	pass		Ethoprophos <sup>¥</sup>	< LOQ	0.20	0.100	pass	
Etofenprox <sup>¥</sup>	< LOQ	0.40	0.200	pass		Etoazole <sup>¥</sup>	< LOQ	0.20	0.100	pass	
Fenoxycarb <sup>¥</sup>	< LOQ	0.20	0.100	pass		Fenpyroximate <sup>¥</sup>	< LOQ	0.40	0.200	pass	
Fipronil <sup>¥</sup>	< LOQ	0.40	0.200	pass		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		Paclobutrazole <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	0.0822	2309029	07/12/23	AOAC 2013.06 (mod.) <sup>b</sup>	pass		
Cadmium <sup>¥</sup>	< LOQ	0.200	mg/kg	0.0822	2309029	07/12/23	AOAC 2013.06 (mod.) <sup>b</sup>	pass		
Lead <sup>¥</sup>	< LOQ	0.500	mg/kg	0.0822	2309029	07/12/23	AOAC 2013.06 (mod.) <sup>b</sup>	pass		
Mercury <sup>¥</sup>	< LOQ	0.100	mg/kg	0.0411	2309029	07/12/23	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**

µg/g = Microgram per gram

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

% = 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: 2309035

Laboratory Control Sample									
Analyte	LCS	Result	Spike	Units	% Rec	Limits		Evaluation	Notes
CBDVA	2	0.0789	0.0729	%	108	80.0	- 120	Acceptable	
CBDV	2	0.0782	0.0727	%	108	80.0	- 120	Acceptable	
CBE	2	0.0852	0.0803	%	106	80.0	- 120	Acceptable	
CBDA	1	0.0762	0.0750	%	102	90.0	- 110	Acceptable	
CBGA	1	0.0768	0.0753	%	102	80.0	- 120	Acceptable	
CBG	1	0.0780	0.0766	%	102	80.0	- 120	Acceptable	
CBD	1	0.0792	0.0779	%	102	90.0	- 110	Acceptable	
THCV	2	0.0574	0.0546	%	105	80.0	- 120	Acceptable	
d8THCV	2	0.0666	0.0644	%	103	80.0	- 120	Acceptable	
THCVA	2	0.0785	0.0711	%	110	80.0	- 120	Acceptable	
CBN	1	0.0787	0.0784	%	100	80.0	- 120	Acceptable	
exo-THC	2	0.0668	0.0653	%	102	80.0	- 120	Acceptable	
d9THC	1	0.0778	0.0759	%	103	90.0	- 110	Acceptable	
d8THC	1	0.0727	0.0738	%	98.5	90.0	- 110	Acceptable	
9S-d10THC	1	0.0782	0.0791	%	98.9	80.0	- 120	Acceptable	
CBL	2	0.0757	0.0718	%	105	80.0	- 120	Acceptable	
9R-d10THC	1	0.0709	0.0722	%	98.2	80.0	- 120	Acceptable	
CBC	2	0.0688	0.0675	%	102	80.0	- 120	Acceptable	
THCA	1	0.0733	0.0744	%	98.6	90.0	- 110	Acceptable	
CBCA	2	0.0797	0.0737	%	108	80.0	- 120	Acceptable	
CBLA	2	0.0748	0.0698	%	107	80.0	- 120	Acceptable	
d9THCP	2	0.0746	0.0752	%	99.3	80.0	- 120	Acceptable	
CBT	2	0.0708	0.0753	%	94.1	80.0	- 120	Acceptable	

Method Blank						
Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.0735	%	< 0.0735	Acceptable	
CBDV	<LOQ	0.0735	%	< 0.0735	Acceptable	
CBE	<LOQ	0.0735	%	< 0.0735	Acceptable	
CBDA	<LOQ	0.0735	%	< 0.0735	Acceptable	
CBGA	<LOQ	0.0735	%	< 0.0735	Acceptable	
CBG	<LOQ	0.0735	%	< 0.0735	Acceptable	
CBD	<LOQ	0.0735	%	< 0.0735	Acceptable	
THCV	<LOQ	0.0735	%	< 0.0735	Acceptable	
d8THCV	<LOQ	0.0735	%	< 0.0735	Acceptable	
THCVA	<LOQ	0.0735	%	< 0.0735	Acceptable	
CBN	<LOQ	0.0735	%	< 0.0735	Acceptable	
exo-THC	<LOQ	0.0735	%	< 0.0735	Acceptable	
d9THC	<LOQ	0.0735	%	< 0.0735	Acceptable	
d8THC	<LOQ	0.0735	%	< 0.0735	Acceptable	
9S-d10THC	<LOQ	0.0735	%	< 0.0735	Acceptable	
CBL	<LOQ	0.0735	%	< 0.0735	Acceptable	
9R-d10THC	<LOQ	0.0735	%	< 0.0735	Acceptable	
CBC	<LOQ	0.0735	%	< 0.0735	Acceptable	
THCA	<LOQ	0.0735	%	< 0.0735	Acceptable	
CBCA	<LOQ	0.0735	%	< 0.0735	Acceptable	
CBLA	<LOQ	0.0735	%	< 0.0735	Acceptable	
d9THCP	<LOQ	0.0735	%	< 0.0735	Acceptable	
CBT	<LOQ	0.0735	%	< 0.0735	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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Revision: 4 Document ID: 7148  
Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

J AOAC 2015 V98-6		Batch ID: 2309035						
Sample Duplicate		Sample ID: 23-008095-0001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.0747	%	NA	< 20	Acceptable	
CBDV	<LOQ	<LOQ	0.0747	%	NA	< 20	Acceptable	
CBE	<LOQ	<LOQ	0.0747	%	NA	< 20	Acceptable	
CBDA	0.138	0.131	0.0747	%	5.21	< 20	Acceptable	
CBGA	1.36	1.34	0.0747	%	1.29	< 20	Acceptable	
CBG	0.159	0.158	0.0747	%	0.891	< 20	Acceptable	
CBD	<LOQ	<LOQ	0.0747	%	NA	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.0747	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.0747	%	NA	< 20	Acceptable	
THCVA	0.369	0.365	0.0747	%	0.965	< 20	Acceptable	
CBN	<LOQ	<LOQ	0.0747	%	NA	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.0747	%	NA	< 20	Acceptable	
d9THC	2.08	1.99	0.0747	%	4.11	< 20	Acceptable	
d8THC	0.562	0.553	0.0747	%	1.47	< 20	Acceptable	
9S-d10THC	<LOQ	<LOQ	0.0747	%	NA	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.0747	%	NA	< 20	Acceptable	
9R-d10THC	<LOQ	<LOQ	0.0747	%	NA	< 20	Acceptable	
CBC	<LOQ	<LOQ	0.0747	%	NA	< 20	Acceptable	
THCA	73.5	72.9	0.0747	%	0.815	< 20	Acceptable	
CBCA	1.14	1.07	0.0747	%	6.66	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.0747	%	NA	< 20	Acceptable	
d9THCP	<LOQ	<LOQ	0.0747	%	NA	< 20	Acceptable	
CBT	<LOQ	<LOQ	0.0747	%	NA	< 20	Acceptable	

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Units of Measure:

% - Percent



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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: 2309040				
Method Blank		Laboratory Control Sample						
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes
Abamectin	0.000	< 0.250		0.996	1.000	99.6	50.0	150
Acephate	0.000	< 0.200		0.671	0.800	83.8	60.0	120
Acetamiprid	0.000	< 1.000		3.881	4.000	97.0	40.0	160
Acetamiprid	0.000	< 0.100		0.379	0.400	94.8	60.0	120
Aldicarb	0.000	< 0.200		0.791	0.800	98.9	60.0	120
Azoxystrobin	0.000	< 0.100		0.395	0.400	98.8	60.0	120
Bifenazate	0.000	< 0.100		0.377	0.400	94.2	60.0	120
Bifenthrin	0.000	< 0.100		0.389	0.400	97.3	50.0	150
Boscalid	0.000	< 0.200		0.718	0.800	89.8	60.0	120
Carbaryl	0.000	< 0.100		0.382	0.400	95.4	60.0	120
Carbofuran	0.000	< 0.100		0.379	0.400	94.9	60.0	120
Chlorantraniliprole	0.000	< 0.100		0.395	0.400	98.8	60.0	120
Chlorfenapyr	0.000	< 0.500		1.619	2.000	80.9	60.0	120
Chlorpyrifos	0.000	< 0.100		0.347	0.400	86.7	60.0	120
Clofentazine	0.000	< 0.100		0.395	0.400	98.7	60.0	120
Cyfluthrin	0.000	< 0.500		2.064	2.000	103.2	50.0	150
Cypermethrin	0.000	< 0.500		1.916	2.000	95.8	50.0	150
Daminozide	0.383	< 0.500		2.187	2.000	109.4	60.0	120
Diazinon	0.000	< 0.100		0.371	0.400	92.8	60.0	120
Dichlorvos	0.000	< 0.500		1.856	2.000	92.8	60.0	120
Dimethoate	0.000	< 0.100		0.401	0.400	100.3	60.0	120
Ethoprophos	0.000	< 0.100		0.375	0.400	93.7	60.0	120
Etofenprox	0.000	< 0.200		0.770	0.800	96.3	50.0	150
Etoxazole	0.000	< 0.100		0.369	0.400	92.2	60.0	120
Fenoxycarb	0.000	< 0.100		0.395	0.400	98.7	60.0	120
Fenpyroximate	0.000	< 0.200		0.761	0.800	95.2	60.0	120
Fipronil	0.000	< 0.200		0.800	0.800	100.1	60.0	120
Fonicamid	0.000	< 0.250		1.118	1.000	111.8	60.0	120
Fludioxonil	0.000	< 0.200		0.747	0.800	93.3	50.0	150
Hexythiazox	0.000	< 0.250		0.983	1.000	98.3	60.0	120
Imazalil	0.000	< 0.100		0.392	0.400	98.1	60.0	120
Imidacloprid	0.000	< 0.200		0.776	0.800	97.0	60.0	120
Kresoxim-methyl	0.000	< 0.200		0.800	0.800	100.0	60.0	120
Malathion	0.000	< 0.100		0.383	0.400	95.8	60.0	120
Metlaxyl	0.000	< 0.100		0.391	0.400	97.6	60.0	120
Methiocarb	0.000	< 0.100		0.384	0.400	96.1	60.0	120
Methomyl	0.000	< 0.200		0.845	0.800	105.7	60.0	120
MGK-264	0.000	< 0.100		0.423	0.400	105.6	50.0	150
Myclobutanil	0.000	< 0.100		0.369	0.400	92.2	60.0	120
Naled	0.000	< 0.250		0.965	1.000	96.5	50.0	150
Oxamyl	0.000	< 0.500		2.031	2.000	101.5	60.0	120
Pacllobutrazole	0.000	< 0.200		0.762	0.800	95.2	60.0	120
Parathion-Methyl	0.000	< 0.100		0.376	0.400	94.0	50.0	150
Permethrin	0.000	< 0.100		0.371	0.400	92.9	50.0	150
Phosmet	0.000	< 0.100		0.406	0.400	101.4	50.0	150
Piperonyl butoxide	0.000	< 0.500		1.940	2.000	97.0	60.0	120
Prallethrin	0.000	< 0.100		0.378	0.400	94.6	60.0	120
Propiconazole	0.000	< 0.200		0.751	0.800	93.9	60.0	120
Propoxur	0.000	< 0.100		0.376	0.400	94.1	60.0	120
Pyrethrin (Summe)	0.001	< 0.100		0.493	0.488	101.0	60.0	120
Pyridaben	0.000	< 0.100		0.389	0.400	97.2	50.0	150
Spirosad	0.000	< 0.100		0.382	0.388	98.5	50.0	150
Spiromesifen	0.000	< 0.100		0.395	0.400	98.8	60.0	120
Spirotetramat	0.000	< 0.100		0.395	0.400	98.7	60.0	120
Spiroxamine	0.000	< 0.200		0.759	0.800	94.9	60.0	120
Tebuconazole	0.000	< 0.200		0.778	0.800	97.3	60.0	120
Thiacloprid	0.000	< 0.100		0.378	0.400	94.5	60.0	120
Thiamethoxam	0.000	< 0.100		0.421	0.400	105.1	60.0	120
Trifloxystrobin	0.000	< 0.100		0.388	0.400	97.1	60.0	120





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



**Report Number:** 23-008138/D003.R000  
**Report Date:** 07/18/2023  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 07/11/23 14:30

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: 2309040				
Matrix Spike/Matrix Spike Duplicate Recoveries		Sample ID: 23-008095-0001									
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes	
Abamectin	0.000	0.794	0.906	1.000	13.3%	< 30	79.4%	90.6%	50 - 150		
Acephate	0.000	0.842	0.667	0.800	23.2%	< 30	105.3%	83.4%	50 - 150		
Acequinocyl	0.000	2.429	3.099	4.000	24.2%	< 30	60.7%	77.5%	50 - 150		
Acetamiprid	0.000	0.364	0.373	0.400	2.4%	< 30	91.0%	93.2%	50 - 150		
Aldicarb	0.000	0.760	0.756	0.800	0.5%	< 30	95.0%	94.6%	50 - 150		
Azoxystrobin	0.000	0.360	0.342	0.400	5.4%	< 30	90.1%	85.4%	50 - 150		
Bifenazate	0.000	0.361	0.384	0.400	6.2%	< 30	90.2%	95.9%	50 - 150		
Bifenthrin	0.000	0.219	0.278	0.400	23.9%	< 30	54.6%	69.4%	50 - 150		
Boscalid	0.000	0.742	0.869	0.800	15.7%	< 30	92.8%	108.6%	50 - 150		
Carbaryl	0.000	0.354	0.365	0.400	2.9%	< 30	88.5%	91.2%	50 - 150		
Carbofuran	0.000	0.363	0.375	0.400	3.3%	< 30	90.8%	93.8%	50 - 150		
Chlorantraniliprole	0.000	0.365	0.374	0.400	2.4%	< 30	91.2%	93.4%	50 - 150		
Chlorfenapyr	0.000	1.575	1.602	2.000	1.7%	< 30	78.7%	80.1%	50 - 150		
Chlorpyrifos	0.000	0.339	0.366	0.400	7.6%	< 30	84.8%	91.5%	50 - 150		
Clofentazine	0.000	0.419	0.483	0.400	14.3%	< 30	104.7%	120.8%	50 - 150		
Cyfluthrin	0.000	1.331	1.382	2.000	3.8%	< 30	66.5%	69.1%	30 - 150		
Cypermethrin	0.000	1.246	1.345	2.000	7.6%	< 30	62.3%	67.2%	50 - 150		
Daminozide	0.000	3.429	2.241	2.000	41.9%	< 30	171.5%	112.1%	30 - 150	R, Q	
Diazinon	0.000	0.368	0.396	0.400	7.2%	< 30	92.1%	99.0%	50 - 150		
Dichlorvos	0.000	1.741	1.734	2.000	0.4%	< 30	87.1%	86.7%	50 - 150		
Dimethoate	0.000	0.337	0.371	0.400	9.7%	< 30	84.1%	92.8%	50 - 150		
Ethoprophos	0.000	0.346	0.361	0.400	4.3%	< 30	86.4%	90.2%	50 - 150		
Etofenprox	0.000	0.035	0.035	0.800	0.9%	< 30	4.4%	4.3%	50 - 150	Q	
Etoxazole	0.000	0.295	0.315	0.400	6.4%	< 30	73.8%	78.7%	50 - 150		
Fenoxycarb	0.000	0.368	0.385	0.400	4.6%	< 30	91.9%	96.2%	50 - 150		
Fenpyroximate	0.000	0.598	0.639	0.800	6.6%	< 30	74.8%	79.9%	50 - 150		
Fipronil	0.000	0.834	0.897	0.800	7.3%	< 30	104.3%	112.2%	50 - 150		
Fonicamid	0.000	0.973	1.012	1.000	4.0%	< 30	97.3%	101.2%	50 - 150		
Fludioxonil	0.000	0.751	0.730	0.800	2.8%	< 30	93.8%	91.3%	50 - 150		
Hexythiazox	0.000	0.863	0.931	1.000	7.6%	< 30	86.3%	93.1%	50 - 150		
Imazalil	0.000	0.344	0.374	0.400	8.4%	< 30	85.9%	93.4%	50 - 150		
Imidacloprid	0.000	0.678	0.715	0.800	5.4%	< 30	84.7%	89.4%	50 - 150		
Kresoxim-methyl	0.000	0.741	0.720	0.800	2.9%	< 30	92.6%	90.0%	50 - 150		
Malathion	0.000	0.386	0.375	0.400	2.8%	< 30	96.5%	93.8%	50 - 150		
Metaxalyl	0.000	0.378	0.378	0.400	0.1%	< 30	94.5%	94.6%	50 - 150		
Methiocarb	0.000	0.378	0.384	0.400	1.4%	< 30	94.6%	95.9%	50 - 150		
Methomyl	0.000	0.648	0.799	0.800	20.8%	< 30	81.1%	99.9%	50 - 150		
MGK-264	0.000	0.388	0.414	0.400	6.6%	< 30	96.9%	103.5%	50 - 150		
Myclobutanil	0.000	0.351	0.352	0.400	0.4%	< 30	87.7%	88.1%	50 - 150		
Naled	0.000	0.902	0.908	1.000	0.7%	< 30	90.2%	90.8%	50 - 150		
Oxamyl	0.000	1.515	1.760	2.000	15.0%	< 30	75.7%	88.0%	50 - 150		
Pacllobutrazole	0.000	0.718	0.738	0.800	2.8%	< 30	89.8%	92.3%	50 - 150		
Parathion-Methyl	0.067	0.386	0.401	0.400	4.5%	< 30	79.9%	83.6%	30 - 150		
Permethrin	0.000	0.237	0.259	0.400	8.8%	< 30	59.2%	64.6%	50 - 150		
Phosmet	0.000	0.375	0.376	0.400	0.1%	< 30	93.8%	93.9%	50 - 150		
Piperonyl butoxide	0.000	1.534	1.787	2.000	15.2%	< 30	76.7%	89.3%	50 - 150		
Prallethrin	0.000	0.409	0.413	0.400	0.8%	< 30	102.4%	103.2%	50 - 150		
Propiconazole	0.000	0.676	0.731	0.800	7.9%	< 30	84.5%	91.4%	50 - 150		
Propoxur	0.000	0.356	0.364	0.400	2.3%	< 30	89.0%	91.0%	50 - 150		
Pyrethrin (Summe)	0.033	0.702	0.728	0.488	3.8%	< 30	137.1%	142.4%	50 - 150		
Pyridaben	0.000	0.220	0.247	0.400	11.5%	< 30	55.1%	61.9%	50 - 150		
Spirosad	0.000	0.291	0.319	0.388	9.3%	< 30	74.9%	82.2%	50 - 150		
Spiromesifen	0.000	0.312	0.325	0.400	4.0%	< 30	78.1%	81.2%	50 - 150		
Spirotetramat	0.000	0.384	0.398	0.400	3.6%	< 30	96.0%	99.5%	50 - 150		
Spiroxamine	0.000	0.714	0.787	0.800	9.8%	< 30	89.2%	98.4%	50 - 150		
Tebuconazole	0.000	0.703	0.755	0.800	7.1%	< 30	87.8%	94.3%	50 - 150		
Thiacloprid	0.000	0.370	0.382	0.400	3.3%	< 30	92.5%	95.5%	50 - 150		
Thiamethoxam	0.000	0.349	0.383	0.400	9.3%	< 30	87.2%	95.7%	50 - 150		
Trifloxystrobin	0.000	0.309	0.341	0.400	10.0%	< 30	77.2%	85.3%	50 - 150		



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**Report Date:** 07/18/2023  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 07/11/23 14:30

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

Laboratory Quality Control Results

Residual Solvents				Batch ID: 2309154					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		626	584	µg/g	107.2	60 - 120	
Isobutane	ND	< 200		837	767	µg/g	109.1	60 - 120	
Butane	ND	< 200		815	782	µg/g	104.2	60 - 120	
2,2-Dimethylpropane	ND	< 200		1030	939	µg/g	109.7	60 - 120	
Methanol	ND	< 200		1700	1640	µg/g	103.7	60 - 120	
Ethylene Oxide	ND	< 30		70.2	57.1	µg/g	122.9	60 - 120	
2-Methylbutane	ND	< 200		1540	1600	µg/g	96.3	60 - 120	
Pentane	ND	< 200		1560	1620	µg/g	96.3	60 - 120	
Ethanol	ND	< 200		1730	1610	µg/g	107.5	70 - 130	
Ethyl Ether	ND	< 200		1590	1610	µg/g	98.8	60 - 120	
2,2-Dimethylbutane	ND	< 30		167	168	µg/g	99.4	60 - 120	
Acetone	ND	< 200		1620	1620	µg/g	100.0	60 - 120	
2-Propanol	ND	< 200		1720	1600	µg/g	107.5	60 - 120	
Ethyl Formate	ND	< 500		1530	1600	µg/g	95.6	70 - 130	
Acetonitrile	ND	< 100		474	484	µg/g	97.9	60 - 120	
Methyl Acetate	ND	< 500		1630	1610	µg/g	101.2	70 - 130	
2,3-Dimethylbutane	ND	< 30		152	162	µg/g	93.8	60 - 120	
Dichloromethane	ND	< 60		474	483	µg/g	98.1	60 - 120	
2-Methylpentane	ND	< 30		176	174	µg/g	101.1	60 - 120	
MTBE	ND	< 500		1600	1610	µg/g	99.4	70 - 130	
3-Methylpentane	ND	< 30		171	168	µg/g	101.8	60 - 120	
Hexane	ND	< 30		160	168	µg/g	95.2	60 - 120	
1-Propanol	ND	< 500		1750	1600	µg/g	109.4	70 - 130	
Methylethylketone	ND	< 500		1620	1620	µg/g	100.0	70 - 130	
Ethyl acetate	ND	< 200		1600	1600	µg/g	100.0	60 - 120	
2-Butanol	ND	< 200		1710	1600	µg/g	106.9	60 - 120	
Tetrahydrofuran	ND	< 100		496	514	µg/g	96.5	60 - 120	
Cyclohexane	ND	< 200		1550	1600	µg/g	96.9	60 - 120	
2-methyl-1-propanol	ND	< 500		1830	1610	µg/g	113.7	70 - 130	
Benzene	ND	< 1		4.07	5.12	µg/g	79.5	60 - 120	
Isopropyl Acetate	ND	< 200		1600	1620	µg/g	98.8	60 - 120	
Heptane	ND	< 200		1550	1610	µg/g	96.3	60 - 120	
1-Butanol	ND	< 500		1810	1600	µg/g	113.1	70 - 130	
Propyl Acetate	ND	< 500		1580	1600	µg/g	98.8	70 - 130	
1,4-Dioxane	ND	< 100		460	493	µg/g	93.3	60 - 120	
2-Ethoxyethanol	ND	< 30		165	163	µg/g	101.2	60 - 120	
Methylisobutylketone	ND	< 500		1600	1600	µg/g	100.0	70 - 130	
3-Methyl-1-butanol	ND	< 500		1770	1610	µg/g	109.9	70 - 130	
Ethylene Glycol	ND	< 200		271	483	µg/g	56.1	60 - 120	Q6
Toluene	ND	< 100		452	493	µg/g	91.7	60 - 120	
Isobutyl Acetate	ND	< 500		1560	1600	µg/g	97.5	70 - 130	
1-Pentanol	ND	< 500		1860	1600	µg/g	116.3	70 - 130	
Butyl Acetate	ND	< 500		1530	1600	µg/g	95.6	70 - 130	
Ethylbenzene	ND	< 200		876	969	µg/g	90.4	60 - 120	
m,p-Xylene	ND	< 200		865	968	µg/g	89.4	60 - 120	
o-Xylene	ND	< 200		872	976	µg/g	89.3	60 - 120	
Cumene	ND	< 30		141	162	µg/g	87.0	60 - 120	
Anisole	ND	< 500		1390	1610	µg/g	86.3	70 - 130	
DMSO	ND	< 500		1460	1610	µg/g	90.7	70 - 130	
1,2-dimethoxyethane	ND	< 50		158	164	µg/g	96.3	70 - 130	
Triethylamine	ND	< 500		1410	1600	µg/g	88.1	70 - 130	
N,N-dimethylformamide	ND	< 150		463	484	µg/g	95.7	70 - 130	
N,N-dimethylacetamide	ND	< 150		448	489	µg/g	91.6	70 - 130	
Pyridine	ND	< 50		121	172	µg/g	70.3	70 - 130	
Sulfolane	ND	< 50		104	163	µg/g	63.8	70 - 130	Q6
1,2-Dichloroethane	ND	< 1		1.02	1	µg/g	102.0	70 - 130	
Chloroform	ND	< 1		1.08	1	µg/g	108.0	70 - 130	
Trichloroethylene	ND	< 1		1.09	1	µg/g	109.0	70 - 130	
1,1-Dichloroethane	ND	< 1		1.08	1	µg/g	108.0	70 - 130	



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**Purchase Order:**  
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Revision: 2 Document ID: 7087  
 Legacy ID: CFL-E33Effective:

QC - Sample Duplicate		Sample ID: 23-008095-0001						
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



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**Purchase Order:**  
**Received:** 07/11/23 14:30

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.