



Customer: The Hemp Collect
2014 SE 9th Ave
Portland Oregon 97214
United States of America (USA)

Product identity: Live D9 Caramel, Knockout, Naturally Derived

Metrc ID: .

Material: Cannabinoid Edible

Laboratory ID: 25-011744-0001

Evidence of Cooling: No

Temp: 20.9 °C

Lot #: 5005SH_100125

Serving Size #1: 18 g



**THE HEMP
COLLECT**

Sample Results

Potency		Method: J AOAC 2015 V98-6 (mod) ^p			Batch: 2507279		Analyze: 10/03/25	
Analyte	Result	Units	LOQ	Notes	Serving Size #1			
					Result	Units	LOQ	
CBD [±]	< LOQ	%	0.0071		< LOQ	mg/18g	1.3	
CBD-A [±]	0.0122	%	0.0071		2.20	mg/18g	1.3	
CBD-Total [±]	< LOQ	%	0.0133		< LOQ	mg/18g	2.40	
CBG	0.00744	%	0.0071		1.34	mg/18g	1.3	
CBG-A	< LOQ	%	0.0071		< LOQ	mg/18g	1.3	
CBG-Total	< LOQ	%	0.0132		< LOQ	mg/18g	2.38	
CBN	0.0623	%	0.0071		11.2	mg/18g	1.3	
Δ10-THC-9R	< LOQ	%	0.0071		< LOQ	mg/18g	1.3	
Δ10-THC-9S	< LOQ	%	0.0071		< LOQ	mg/18g	1.3	
Δ10-THC-Total	< LOQ	%	0.0142		< LOQ	mg/18g	2.55	
Δ8-THC [±]	< LOQ	%	0.0071		< LOQ	mg/18g	1.3	
Δ9-THC [±]	0.228	%	0.0071		41.1	mg/18g	1.3	
Δ9-THC-A [±]	< LOQ	%	0.0071		< LOQ	mg/18g	1.3	
Δ9-THC-Total [±]	0.228	%	0.0133		41.0	mg/18g	2.40	
Total Cannabinoids	0.310	%			55.8	mg/18g		

Microbiology							
Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status Notes
Salmonella spp. [±]	Negative		/25g		2507665	10/22/25 AOAC 2020.02 ^p	
EHEC including STEC [±]	Negative		/25g		2507666	10/22/25 AOAC 2020.06 ^p	

Solvents		Method: Residual Solvents by HS-GC-MS ^p				Units µg/g		Batch 2507764		Analyze: 10/23/25	
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		



Solvents		Method: Residual Solvents by HS-GC-MS ^B				Units µg/g	Batch 2507764	Analyze: 10/23/25			
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
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 [⊥]	< 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)				Units mg/kg	Batch 2507725	Analyze: 10/22/25			
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
Abamectin [⊥]	< LOQ	0.50	0.250	pass		Acephate [⊥]	< LOQ	0.40	0.200	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-ethyl [⊥]	< LOQ	0.20	0.100	pass	
Clofentezine [⊥]	< LOQ	0.20	0.100	pass		Cyfluthrin (sum) [⊥]	< LOQ	1.0	0.500	pass	
Cypermethrin (sum) [⊥]	< 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		Flonicamid [⊥]	< 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		Paclobutrazole [⊥]	< LOQ	0.40	0.200	pass	
Parathion-methyl [⊥]	< LOQ	0.20	0.100	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	0.100	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	



Pesticides		Method: AOAC 2007.01 & EN 15662 (mod)				Units mg/kg	Batch 2507725	Analyze: 10/22/25			
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
Thiacloprid [±]	< LOQ	0.20	0.100	pass		Thiamethoxam [±]	< LOQ	0.20	0.100	pass	
Trifloxystrobin [±]	< LOQ	0.20	0.100	pass							

Metals										
Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method		Status	Notes	
Arsenic [±]	< LOQ	0.200	mg/kg	0.0140	2507791	10/23/25	AOAC 2013.06 (mod.) ^b	pass		
Cadmium [±]	< LOQ	0.200	mg/kg	0.0140	2507791	10/23/25	AOAC 2013.06 (mod.) ^b	pass		
Lead [±]	< LOQ	0.500	mg/kg	0.0140	2507791	10/23/25	AOAC 2013.06 (mod.) ^b	pass		
Mercury [±]	< LOQ	0.100	mg/kg	0.00701	2507791	10/23/25	AOAC 2013.06 (mod.) ^b	pass		

Mycotoxins										
Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method		Status	Notes	
Aflatoxin B1 [±]	< LOQ		µg/kg	5.00	2507801	10/24/25	Mycotoxins by AOAC 2007.01			
Aflatoxin B2 [±]	< LOQ		µg/kg	5.00	2507801	10/24/25	Mycotoxins by AOAC 2007.01			
Aflatoxin G1 [±]	< LOQ		µg/kg	5.00	2507801	10/24/25	Mycotoxins by AOAC 2007.01			
Aflatoxin G2 [±]	< LOQ		µg/kg	5.00	2507801	10/24/25	Mycotoxins by AOAC 2007.01			
Ochratoxin A	< LOQ	20.0	µg/kg	5.00	2507801	10/24/25	Mycotoxins by AOAC 2007.01 ^b	pass		
Total Aflatoxins	< LOQ	20.0	µg/kg	20.0		10/27/25	Mycotoxins by AOAC 2007.01 ^b	pass		



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

Report Number: 25-011744/D005.R000
Report Date: 10/27/2025
ORELAP#: OR100028
Purchase Order:
Received: 10/01/25 09:46



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.

Threshold Note: OAR 333-007-0400

Ⓟ = ISO/IEC 17025:2017 accredited method.

Ⓡ = TNI accredited analyte.

Units of Measure

/25g = Per 25g

µg/g = Microgram per gram

µg/kg = Micrograms per kilogram = parts per billion (ppb)

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

% = Percentage of sample

mg/18g = Milligram per 18g

% wt = µg/g divided by 10,000

Approved Signatory

Derrick Tanner
General Manager



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

Report Number: 25-011744/D005.R000
Report Date: 10/27/2025
ORELAP#: OR100028
Purchase Order:
Received: 10/01/25 09:46



**Hemp & Cannabis
Chain of Custody**

**The-Hemp-Collect-
1759271099**

Company Details Company: <u>The Hemp Collect</u> Contact: <u>Cris Kingsland</u> Street Address: <u>2014 SE 9th</u> City, State, Zip: <u>Portland, OR 97214</u> Email: <u>coas@thehempcollect.com</u> Contact Phone: <u>7707220962</u> Billing Information Billing Email: <u>accounting@thehempcollect.com</u>				Project Details Turnaround Time: <u>4 Business Days Surcharges Apply</u> Relinquishment Sampling, Courier & Shipping Options: <u>By Shipping Service (USPS, UPS, Fedex)</u> Additional Comments for Project: <u>will be requesting further test after reviewing potency.</u> Receipt Information Evidence of Cooling?: No Sample Condition: Satisfactory Prelog Storage: Canna Shelves			Testing ±
#	Sample Name	Lot Additional Sample ID	Material	Amount Provided	Reporting Unit	Serving Size	
1	Live D9 Caramel, Knockout, Naturally Derived	5005SH_100125	Cannabinoid Edible	72 g	mg/g & mg/serving	18 g	✓
2	Live D9 Caramel, Anytime, Naturally Derived	5004SH_100125	Cannabinoid Edible	72 g	mg/g & mg/serving	18 g	✓

Relinquished By	Date	Time	Received By	Date	Time	Received Temp., °C	IR Therm. CL#
<i>Cris Kingsland</i>	<i>09/30/2025</i>	<i>15:24</i>	<i>rls</i>	<i>10/01/2025</i>	<i>09:46</i>	<i>20.90</i>	<i>CL-0530</i>

Samples submitted to Columbia Laboratories with testing requirements constitute an agreement for services in accordance with the [current terms of services](#) associated with this COC. By signing "Relinquished by" you are agreeing to these terms.
 Columbia Laboratories
 12423 NE Whitaker Way
 Portland, OR 97230
 P: (503) 254-1794
info@columbiaboratories.com
 Page 1 of 1
www.columbiaboratories.com



Laboratory Quality Control Results

J AOAC 2015 V98-6 Batch ID: 2507279

Laboratory Control Sample

Analyte	LCS	Result	Spike	Units	% Rec	Limits	Evaluation	Notes
CBDVA	2	0.0281	0.0274	%	102	80.0 - 120	Acceptable	
CBDV	2	0.0296	0.0288	%	103	80.0 - 120	Acceptable	
CBE	2	0.0316	0.0312	%	101	80.0 - 120	Acceptable	
CBDA	1	0.0282	0.0257	%	110	90.0 - 110	Acceptable	
CBGA	1	0.0299	0.0292	%	102	80.0 - 120	Acceptable	
CBG	1	0.0273	0.0272	%	100	80.0 - 120	Acceptable	
CBD	1	0.0287	0.0270	%	106	90.0 - 110	Acceptable	
THCV	2	0.0302	0.0296	%	102	80.0 - 120	Acceptable	
d8THCV	2	0.0304	0.0298	%	102	80.0 - 120	Acceptable	
THCVA	2	0.0269	0.0260	%	103	80.0 - 120	Acceptable	
CBN	1	0.0288	0.0274	%	105	80.0 - 120	Acceptable	
exo-THC	2	0.0274	0.0272	%	101	80.0 - 120	Acceptable	
d9THC	1	0.0315	0.0295	%	107	90.0 - 110	Acceptable	
d8THC	1	0.0280	0.0271	%	103	90.0 - 110	Acceptable	
9S-d10THC	1	0.0319	0.0303	%	105	80.0 - 120	Acceptable	
CBL	2	0.0294	0.0282	%	104	80.0 - 120	Acceptable	
9R-d10THC	1	0.0318	0.0301	%	105	80.0 - 120	Acceptable	
CBC	2	0.0302	0.0298	%	102	80.0 - 120	Acceptable	
THCA	1	0.0334	0.0317	%	105	90.0 - 110	Acceptable	
CBCA	2	0.0288	0.0279	%	103	80.0 - 120	Acceptable	
CBLA	2	0.0282	0.0277	%	102	80.0 - 120	Acceptable	
d9THCP	2	0.0278	0.0271	%	103	80.0 - 120	Acceptable	
CBT	2	0.0301	0.0291	%	103	80.0 - 120	Acceptable	

Method Blank

Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.00709	%	< 0.00709	Acceptable	
CBDV	<LOQ	0.00709	%	< 0.00709	Acceptable	
CBE	<LOQ	0.00709	%	< 0.00709	Acceptable	
CBDA	<LOQ	0.00709	%	< 0.00709	Acceptable	
CBGA	<LOQ	0.00709	%	< 0.00709	Acceptable	
CBG	<LOQ	0.00709	%	< 0.00709	Acceptable	
CBD	<LOQ	0.00709	%	< 0.00709	Acceptable	
THCV	<LOQ	0.00709	%	< 0.00709	Acceptable	
d8THCV	<LOQ	0.00709	%	< 0.00709	Acceptable	
THCVA	<LOQ	0.00709	%	< 0.00709	Acceptable	
CBN	<LOQ	0.00709	%	< 0.00709	Acceptable	
exo-THC	<LOQ	0.00709	%	< 0.00709	Acceptable	
d9THC	<LOQ	0.00709	%	< 0.00709	Acceptable	
d8THC	<LOQ	0.00709	%	< 0.00709	Acceptable	
9S-d10THC	<LOQ	0.00709	%	< 0.00709	Acceptable	
CBL	<LOQ	0.00709	%	< 0.00709	Acceptable	
9R-d10THC	<LOQ	0.00709	%	< 0.00709	Acceptable	
CBC	<LOQ	0.00709	%	< 0.00709	Acceptable	
THCA	<LOQ	0.00709	%	< 0.00709	Acceptable	
CBCA	<LOQ	0.00709	%	< 0.00709	Acceptable	
CBLA	<LOQ	0.00709	%	< 0.00709	Acceptable	
d9THCP	<LOQ	0.00709	%	< 0.00709	Acceptable	
CBT	<LOQ	0.00709	%	< 0.00709	Acceptable	

Abbreviations

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

Units of Measure:

% - Percent



Laboratory Quality Control Results

AOAC 2015 V98-6		Batch ID: 2507279						
Sample Duplicate		Sample ID: 25-011637-0001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	
CBDV	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	
CBE	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	
CBDA	<LOQ	<LOQ	0.00707	%	NA	< 10	Acceptable	
CBGA	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	
CBG	0.00866	0.00814	0.00707	%	6.19	< 20	Acceptable	
CBD	<LOQ	<LOQ	0.00707	%	NA	< 10	Acceptable	
THCV	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	
CBN	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	
d9THC	0.269	0.268	0.00707	%	0.400	< 10	Acceptable	
d8THC	<LOQ	<LOQ	0.00707	%	NA	< 10	Acceptable	
9S-d10THC	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	
9R-d10THC	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	
CBC	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.00707	%	NA	< 10	Acceptable	
CBCA	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	
d9THCP	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	
CBT	<LOQ	<LOQ	0.00707	%	NA	< 20	Acceptable	

Abbreviations

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

Units of Measure:

% - Percent



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: 2507725			
Method Blank			Laboratory Control Sample					
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes
Abamectin	0.000	< 0.070		0.229	0.280	81.7	50.0	150
Acephate	0.005	< 0.020		0.073	0.080	91.6	60.0	120
Acequinocyl	0.003	< 0.020		0.063	0.080	78.5	40.0	160
Acetamiprid	0.000	< 0.020		0.071	0.080	89.3	60.0	120
Aldicarb	0.000	< 0.100		0.363	0.400	90.8	60.0	120
Azoxystrobin	0.001	< 0.010		0.036	0.040	89.0	60.0	120
Bifenazate	0.000	< 0.010		0.038	0.040	93.9	60.0	120
Bifenthrin	0.003	< 0.100		0.338	0.400	84.4	50.0	150
Boscalid	0.001	< 0.010		0.037	0.040	93.3	60.0	120
Carbaryl	0.003	< 0.020		0.073	0.080	90.7	60.0	120
Carbofuran	0.000	< 0.010		0.035	0.040	88.3	60.0	120
Chlorantraniliprole	0.000	< 0.010		0.034	0.040	85.1	60.0	120
Chlorfenapyr	0.024	< 0.050		0.145	0.200	72.4	60.0	120
Chlorpyrifos	0.000	< 0.010		0.036	0.040	89.6	60.0	120
Clofentezine	0.001	< 0.010		0.034	0.040	84.8	60.0	120
Cyfluthrin	0.018	< 0.200		0.651	0.800	81.4	50.0	150
Cypermethrin	0.012	< 0.200		0.689	0.800	86.1	50.0	150
Daminozide	0.000	< 0.050		0.065	0.200	32.7	60.0	120
Diazinon	0.000	< 0.010		0.034	0.040	85.2	60.0	120
Dichlorvos	0.004	< 0.050		0.184	0.200	91.9	60.0	120
Dimethoate	0.001	< 0.010		0.034	0.040	85.8	60.0	120
Ethoprophos	0.000	< 0.010		0.033	0.040	82.8	60.0	120
Etofenprox	0.000	< 0.010		0.033	0.040	83.2	50.0	150
Etoazole	0.001	< 0.010		0.034	0.040	84.1	60.0	120
Fenoxycarb	0.000	< 0.010		0.033	0.040	83.4	60.0	120
Fenpyroximate	0.004	< 0.020		0.067	0.080	84.3	60.0	120
Fipronil	0.001	< 0.010		0.032	0.040	80.9	60.0	120
Fonicamid	0.001	< 0.020		0.066	0.080	81.9	60.0	120
Fludioxonil	0.000	< 0.010		0.034	0.040	83.8	50.0	150
Hexythiazox	0.000	< 0.010		0.031	0.040	76.7	60.0	120
Imazalil	0.000	< 0.010		0.036	0.040	90.2	60.0	120
Imidacloprid	0.000	< 0.010		0.036	0.040	89.4	60.0	120
Kresoxim-methyl	0.000	< 0.010		0.033	0.040	83.3	60.0	120
Malathion	0.000	< 0.010		0.033	0.040	81.7	60.0	120
Metalaxyl	0.000	< 0.010		0.040	0.040	98.9	60.0	120
Methiocarb	0.001	< 0.010		0.036	0.040	89.8	60.0	120
Methomyl	0.000	< 0.020		0.076	0.080	95.1	60.0	120
MGK-264	0.004	< 0.050		0.172	0.200	86.0	50.0	150
Myclobutanil	0.003	< 0.010		0.036	0.040	90.5	60.0	120
Naled	0.006	< 0.100		0.348	0.400	87.1	50.0	150
Oxamyl	0.006	< 0.200		0.750	0.800	93.7	60.0	120
Paclobutrazole	0.000	< 0.010		0.036	0.040	89.6	60.0	120
Parathion-Methyl	0.000	< 0.030		0.106	0.120	88.4	50.0	150
Permethrin	0.000	< 0.040		0.130	0.160	81.1	50.0	150
Phosmet	0.000	< 0.010		0.036	0.040	89.6	50.0	150
Piperonyl butoxide	0.013	< 0.200		0.692	0.800	86.5	60.0	120
Prallethrin	0.007	< 0.050		0.176	0.200	88.1	60.0	120
Propiconazole	0.002	< 0.010		0.032	0.040	80.7	60.0	120
Propoxur	0.000	< 0.010		0.035	0.040	88.4	60.0	120
Pyrethrin (Summe)	0.002	< 0.100		0.089	0.100	88.9	60.0	120
Pyridaben	0.003	< 0.020		0.066	0.080	82.8	50.0	150
Spinosad	0.000	< 0.100		0.034	0.040	85.4	50.0	150
Spiromesifen	0.000	< 0.030		0.096	0.120	80.3	60.0	120
Spirotetramat	0.000	< 0.010		0.035	0.040	88.7	60.0	120
Spiroxamine	0.001	< 0.010		0.034	0.040	85.5	60.0	120

Q7


 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: 2507725			
Matrix Spike/Matrix Spike Duplicate Recoveries							Sample ID: 25-011744-0001			
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes
Abamectin	0.019	0.244	0.243	0.280	0.3%	< 30	80.2%	79.9%	50 - 150	
Acephate	0.000	0.075	0.072	0.080	4.4%	< 30	94.1%	90.1%	50 - 150	
Acequinocyl	0.000	0.077	0.077	0.080	0.1%	< 30	96.1%	96.0%	50 - 150	
Acetamiprid	0.000	0.073	0.068	0.080	6.4%	< 30	91.0%	85.3%	50 - 150	
Aldicarb	0.000	0.352	0.349	0.400	0.9%	< 30	88.0%	87.2%	50 - 150	
Azoxystrobin	0.001	0.035	0.035	0.040	0.9%	< 30	85.2%	84.4%	50 - 150	
Bifenazate	0.000	0.032	0.033	0.040	2.1%	< 30	78.8%	80.5%	50 - 150	
Bifenthrin	0.002	0.332	0.314	0.400	5.6%	< 30	82.5%	78.0%	50 - 150	
Boscalid	0.001	0.036	0.037	0.040	2.6%	< 30	87.4%	89.7%	50 - 150	
Carbaryl	0.003	0.078	0.070	0.080	12.1%	< 30	94.6%	83.8%	50 - 150	
Carbofuran	0.000	0.036	0.033	0.040	7.3%	< 30	88.8%	82.5%	50 - 150	
Chlorantraniliprole	0.000	0.035	0.031	0.040	10.3%	< 30	86.9%	78.4%	50 - 150	
Chlorfenapyr	0.000	0.115	0.089	0.200	25.0%	< 30	57.3%	44.6%	50 - 150	Q
Chlorpyrifos	0.000	0.032	0.031	0.040	1.3%	< 30	79.4%	78.4%	50 - 150	
Clofentezine	0.001	0.032	0.031	0.040	3.4%	< 30	77.4%	74.8%	50 - 150	
Cyfluthrin	0.002	0.640	0.498	0.800	25.1%	< 30	79.7%	62.0%	30 - 150	
Cypermethrin	0.000	0.677	0.632	0.800	7.0%	< 30	84.7%	79.0%	50 - 150	
Daminozide	0.000	0.078	0.069	0.200	12.8%	< 30	39.0%	34.3%	30 - 150	
Diazinon	0.000	0.036	0.033	0.040	8.3%	< 30	90.2%	83.0%	50 - 150	
Dichlorvos	0.006	0.170	0.175	0.200	2.8%	< 30	82.3%	84.6%	50 - 150	
Dimethoate	0.000	0.036	0.034	0.040	5.3%	< 30	89.1%	84.5%	50 - 150	
Ethoprophos	0.001	0.033	0.033	0.040	0.9%	< 30	80.0%	80.7%	50 - 150	
Etofenprox	0.000	0.035	0.033	0.040	6.4%	< 30	87.1%	81.7%	50 - 150	
Etoxazole	0.001	0.036	0.033	0.040	8.4%	< 30	87.4%	80.3%	50 - 150	
Fenoxycarb	0.000	0.036	0.035	0.040	1.6%	< 30	89.2%	87.9%	50 - 150	
Fenpyroximate	0.004	0.064	0.059	0.080	8.1%	< 30	74.8%	68.9%	50 - 150	
Fipronil	0.000	0.032	0.033	0.040	2.1%	< 30	80.1%	81.9%	50 - 150	
Fonicamid	0.000	0.067	0.061	0.080	9.5%	< 30	83.6%	76.0%	50 - 150	
Fludioxonil	0.000	0.035	0.033	0.040	6.3%	< 30	88.6%	83.1%	50 - 150	
Hexythiazox	0.000	0.036	0.034	0.040	3.6%	< 30	88.7%	85.5%	50 - 150	
Imazalil	0.000	0.038	0.030	0.040	24.8%	< 30	95.3%	74.3%	50 - 150	
Imidacloprid	0.000	0.039	0.033	0.040	15.3%	< 30	96.5%	82.8%	50 - 150	
Kresoxim-methyl	0.000	0.033	0.031	0.040	6.6%	< 30	81.8%	76.6%	50 - 150	
Malathion	0.000	0.034	0.030	0.040	11.2%	< 30	84.0%	75.1%	50 - 150	
Metalaxyl	0.000	0.037	0.035	0.040	4.5%	< 30	91.9%	87.9%	50 - 150	
Methiocarb	0.001	0.036	0.034	0.040	5.8%	< 30	87.1%	82.2%	50 - 150	
Methomyl	0.002	0.072	0.070	0.080	2.7%	< 30	87.9%	85.6%	50 - 150	
MGK-264	0.003	0.174	0.161	0.200	8.1%	< 30	85.3%	78.6%	50 - 150	
Myclobutanil	0.002	0.036	0.036	0.040	0.1%	< 30	84.8%	84.9%	50 - 150	
Naled	0.000	0.349	0.330	0.400	5.5%	< 30	87.2%	82.5%	50 - 150	
Oxamyl	0.000	0.690	0.700	0.800	1.4%	< 30	86.3%	87.5%	50 - 150	
Paclobutrazole	0.000	0.037	0.037	0.040	1.2%	< 30	93.0%	91.9%	50 - 150	
Parathion-Methyl	0.012	0.122	0.108	0.120	13.5%	< 30	91.5%	79.9%	30 - 150	
Permethrin	0.000	0.138	0.123	0.160	11.5%	< 30	86.3%	76.9%	50 - 150	
Phosmet	0.000	0.035	0.033	0.040	6.6%	< 30	88.0%	82.4%	50 - 150	
Piperonyl butoxide	0.014	0.662	0.667	0.800	0.8%	< 30	80.9%	81.6%	50 - 150	
Prallethrin	0.006	0.174	0.165	0.200	5.2%	< 30	83.9%	79.6%	50 - 150	
Propiconazole	0.001	0.035	0.032	0.040	6.7%	< 30	84.6%	79.1%	50 - 150	
Propoxur	0.000	0.036	0.035	0.040	2.7%	< 30	88.7%	86.3%	50 - 150	
Pyrethrin (Summe)	0.000	0.087	0.064	0.100	31.1%	< 30	87.5%	63.9%	50 - 150	R
Pyridaben	0.003	0.061	0.065	0.080	6.2%	< 30	73.1%	77.7%	50 - 150	
Spinosad	0.000	0.033	0.031	0.040	5.1%	< 30	82.5%	78.4%	50 - 150	
Spiromesifen	0.000	0.104	0.098	0.120	6.5%	< 30	87.1%	81.6%	50 - 150	
Spirotetramat	0.000	0.035	0.033	0.040	6.1%	< 30	88.1%	82.9%	50 - 150	
Spiroxamine	0.001	0.035	0.033	0.040	5.0%	< 30	84.9%	80.8%	50 - 150	



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

Laboratory Quality Control Results

Residual Solvents				Batch ID: 2507764						
Method Blank				Laboratory Control Sample						
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes	
1,1,1-Trichloroethane	ND	< 5		6.33	5	µg/g	126.6	50-150		
1,1-Dichloroethane	ND	< 1		1.28	1	µg/g	128.0	50-150		
1,2-Dichloroethane,cis-	ND	< 1		1.22	1	µg/g	122.0	50-150		
1,4-Dioxane	ND	< 100		573	496	µg/g	115.5	60-120		
1-Propanol	ND	< 500		2110	1690	µg/g	124.9	50-150		
2,2-Dimethylbutane	ND	< 30		201	172	µg/g	116.9	60-120		
2,2-Dimethylpropane	ND	< 200		967	956	µg/g	101.2	60-120		
2,3-Dimethylbutane	ND	< 30		212	173	µg/g	122.5	60-120	Q1	
2-Butanol	ND	< 200		1810	1610	µg/g	112.4	60-120		
2-Ethoxyethanol	ND	< 30		160	177	µg/g	90.4	60-120		
2-Methylbutane	ND	< 200		1990	1630	µg/g	122.1	60-120	Q1	
2-Methylpentane	ND	< 30		184	164	µg/g	112.2	60-120		
2-Propanol	ND	< 200		1780	1610	µg/g	110.6	60-120		
3-Methylpentane	ND	< 30		209	183	µg/g	114.2	60-120		
Acetone	ND	< 200		1800	1620	µg/g	111.1	60-120		
Acetonitrile	ND	< 100		602	493	µg/g	122.1	60-120	Q1	
Anisole	ND	< 500		1420	1680	µg/g	84.5	50-150		
Benzene	ND	< 1		1.21	1	µg/g	121.0	50-150		
Butane	ND	< 200		828	769	µg/g	107.7	60-120		
Chlorobenzene	ND	< 1		1.2	1	µg/g	120.0	50-150		
Chloroform	ND	< 1		1.25	1	µg/g	125.0	50-150		
Cumene	ND	< 30		215	174	µg/g	123.6	60-120	Q1	
Cyclohexane	ND	< 200		1810	1630	µg/g	111.0	60-120		
Dichloromethane	ND	< 1		1.23	1	µg/g	123.0	50-150		
Ethanol	ND	< 200		1970	1630	µg/g	120.9	60-120	Q1	
Ethyl acetate	ND	< 200		1810	1630	µg/g	111.0	60-120		
Ethyl Ether	ND	< 200		1810	1620	µg/g	111.7	60-120		
Ethyl Formate	ND	< 500		1320	1680	µg/g	78.6	50-150		
Ethylbenzene	ND	< 200		1080	976	µg/g	110.7	60-120		
Ethylene Glycol	ND	< 200		361	484	µg/g	74.6	60-120		
Ethylene Oxide	ND	< 1		1.35	1	µg/g	135.0	50-150		
Heptane	ND	< 200		1490	1600	µg/g	93.1	60-120		
Hexane	ND	< 30		193	172	µg/g	112.2	60-120		
Isobutane	ND	< 200		815	770	µg/g	105.8	60-120		
Isopropyl Acetate	ND	< 200		1450	1610	µg/g	90.1	60-120		
m,p-Xylene	ND	< 200		1090	988	µg/g	110.3	60-120		
Methanol	ND	< 200		1980	1650	µg/g	120.0	60-120	Q1	
Methyl Acetate	ND	< 500		1820	1650	µg/g	110.3	50-150		
o-Xylene	ND	< 200		1060	975	µg/g	108.7	60-120		
Pentane	ND	< 200		1950	1610	µg/g	121.1	60-120	Q1	
Propane	ND	< 200		575	585	µg/g	98.3	60-120		
Propyl Acetate	ND	< 500		1360	1660	µg/g	81.9	50-150		
Sulfolane	ND	< 50		60.8	220	µg/g	27.6	50-150	Q7	
Tetrahydrofuran	ND	< 100		529	486	µg/g	108.8	60-120		
Toluene	ND	< 100		543	485	µg/g	112.0	60-120		



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

QC - Sample Duplicate

Sample ID: 25-012461-0003

Analyte	SR Result	SD Result	LOQ	Units	RPD	Limits	Accept/Fail	Notes
1,1,1-Trichloroethane	ND	ND	5	µg/g	0.0	< 20	Acceptable	
1,1-Dichloroethane	ND	ND	1	µg/g	0.0	< 20	Acceptable	
1,2-Dichloroethene,cis-	ND	ND	1	µg/g	0.0	< 20	Acceptable	
1,4-Dioxane	ND	ND	100	µg/g	0.0	< 20	Acceptable	
1-Propanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
2,2-Dimethylbutane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
2,2-Dimethylpropane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2,3-Dimethylbutane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
2-Butanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2-Ethoxyethanol	ND	ND	30	µg/g	0.0	< 20	Acceptable	
2-Methylbutane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2-Methylpentane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
2-Propanol	593	718	200	µg/g	19.1	< 20	Acceptable	
3-Methylpentane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Acetone	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Acetonitrile	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Anisole	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Benzene	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Butane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Chlorobenzene	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Chloroform	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Cumene	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Cyclohexane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Dichloromethane	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Ethanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethyl acetate	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethyl Ether	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethyl Formate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Ethylbenzene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethylene Glycol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethylene Oxide	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Heptane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Hexane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Isobutane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Isopropyl Acetate	ND	ND	200	µg/g	0.0	< 20	Acceptable	
m,p-Xylene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Methanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Methyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
o-Xylene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Pentane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Propane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Propyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Sulfolane	ND	ND	50	µg/g	0.0	< 20	Acceptable	
Tetrahydrofuran	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Toluene	ND	ND	100	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: 25-011744/D005.R000
Report Date: 10/27/2025
ORELAP#: OR100028
Purchase Order:
Received: 10/01/25 09:46





Explanation of QC Flag Comments:

Code	Explanation
A	This analysis was performed on a VOA sample containing headspace.
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.
B3	Dilution water blank of BOD was above the recommended limit; associated samples could be high biased.
CP	Client provided value.
CV	Calculated value.
E	Analyte concentration exceeds the calibration range, results are estimated.
E1	Estimated value.
E2	Estimated value. Matrix interference observed.
H	Holding time was exceeded.
J	Estimated value, above the detection limit and below the LOQ
I	Insufficient sample received to meet method requirements.
LOQ1	Quantitation level raised due to low sample volume and/or dilution.
LOQ2	Quantitation level raised due to matrix interference.
LOQ3	< LOQ could be due to potential inhibition.
N1	See case narrative
P	Not preserved to the proper pH
P1	Storage temperature out of control
P2	Incubator temperature out of control
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.
Q7	Quality control outside QC limits.
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.
RE	Re-extracted and/or re-analyzed.
REH	The original analysis was within holding time; re-analysis past holding time.
S	Surrogate recovery outside control limit.
T	Tentatively Identified Compound (TIC) by library search.
T1	Confirmed by secondary ion
W	Results are reported on dry weight basis.