

CONSOLIDATED TEST RESULTS SUMMARY

Please see the following pages for full test results.

BULK SKU	BATCH #	LOQ: Limit Of Quantitation	
PRODUCT NAME	SERVING SIZE	LOD: Limit Of Detection	
LABORATORY :	OREGON ACCREDITATION: OR100028	1 g = 10 ⁻³ kg = 10 ³ mg = 10 ⁶ 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 ^[1]
Cadmium	µg/serving	µg/g	4.1 µg/day ^[1]
Lead	µg/serving	µg/g	6 µg/day ^[1]
Mercury	µg/serving	µg/g	2 µg/day ^[1]
PESTICIDES	REGULATORY ACTION LEVEL		
None of the other 59 pesticides tested found above limit of detection in the sample.			10 ppb ^[1]
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: 23-000397/D003.R000
Report Date: 01/20/2023
ORELAP#: OR100028
Purchase Order:
Received: 01/10/23 15:40

Customer: Etz Hayim Holdings
Product identity: FORM-GMY.SLP25-EL56
Client/Metric ID: .
Laboratory ID: 23-000397-0002

Summary

Potency:

Analyte per 1g	Result	Limits	Units	Status	
CBC per 1g	0.134		mg/1g		CBD-Total per Serving Size 3.20 mg/1g
CBD per 1g	3.20		mg/1g		
CBDV per 1g	0.0465		mg/1g		THC-Total per Serving Size 0.117 mg/1g
CBG per 1g	1.07		mg/1g		(Reported in milligrams per serving)
CBN per 1g	1.06		mg/1g		
CBT per 1g	0.0716		mg/1g		
Δ9-THC per 1g	0.117		mg/1g		

Residual Solvents:

Analyte	Result (µg/g)	Limits (µg/g)	Status
Ethanol	273		

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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Purchase Order:
Received: 01/10/23 15:40

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

Product identity: FORM-GMY.SLP25-EL56

Client/Metric ID: .

Sample Date:

Laboratory ID: 23-000397-0002

Evidence of Cooling: No

Temp: 19.7

Relinquished by: client

Serving Size #1: 1 g

Sample Results

Potency per 1g	Method: J AOAC 2015 V98-6 (mod) ^b	Units mg/se	Batch: 2300463	Analyze: 1/14/23 5:52:00 AM	
Analyte	Result	Limits	Units	LOQ	Notes
CBC per 1g	0.134		mg/1g	0.0309	
CBC-A per 1g	< LOQ		mg/1g	0.0309	
CBC-Total per 1g	0.134		mg/1g	0.0580	
CBD per 1g	3.20		mg/1g	0.0309	
CBD-A per 1g	< LOQ		mg/1g	0.0309	
CBD-Total per 1g	3.20		mg/1g	0.0580	
CBDV per 1g	0.0465		mg/1g	0.0309	
CBDV-A per 1g	< LOQ		mg/1g	0.0309	
CBDV-Total per 1g	< LOQ		mg/1g	0.0577	
CBE per 1g	< LOQ		mg/1g	0.0309	
CBG per 1g	1.07		mg/1g	0.0309	
CBG-A per 1g	< LOQ		mg/1g	0.0309	
CBG-Total per 1g	1.07		mg/1g	0.0577	
CBL per 1g	< LOQ		mg/1g	0.0309	
CBL-A per 1g	< LOQ		mg/1g	0.0309	
CBL-Total per 1g	< LOQ		mg/1g	0.0580	
CBN per 1g	1.06		mg/1g	0.0309	
CBT per 1g	0.0716		mg/1g	0.0309	
Δ8-THCV per 1g	< LOQ		mg/1g	0.0309	
Δ10-THC-9R per 1g	< LOQ		mg/1g	0.0309	
Δ8-THC per 1g	< LOQ		mg/1g	0.0309	
Δ9-THC per 1g	0.117		mg/1g	0.0309	
exo-THC per 1g	< LOQ		mg/1g	0.0309	
THC-A per 1g	< LOQ		mg/1g	0.0309	
THC-Total per 1g	0.117		mg/1g	0.0580	
THCV per 1g	< LOQ		mg/1g	0.0309	
THCV-A per 1g	< LOQ		mg/1g	0.0309	
THCV-Total per 1g	< LOQ		mg/1g	0.0580	
Total Cannabinoids per 1g	5.70		mg/1g		



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Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2300342	01/14/23 AOAC 991.14 (Petrifilm) ^P		
Total Coliforms	< LOQ		cfu/g	10	2300342	01/14/23 AOAC 991.14 (Petrifilm) ^P		
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2300343	01/15/23 AOAC 2014.05 (RAPID) ^P		
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2300343	01/15/23 AOAC 2014.05 (RAPID) ^P		

Solvents **Method:** Residual Solvents by GC/MS^b **Units** µg/g **Batch** 2300638 **Analyze** 01/20/23 10:57 AM

Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
1,4-Dioxane	< LOQ	380	100	pass		2-Butanol	< LOQ	5000	200	pass	
2-Ethoxyethanol	< LOQ	160	30.0	pass		2-Methylbutane (Isopentane)	< LOQ		200		
2-Methylpentane	< LOQ		30.0			2-Propanol (IPA)	< LOQ	5000	200	pass	
2,2-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	273		200		
Ethyl acetate	< LOQ	5000	200	pass		Ethyl benzene	< LOQ		200		
Ethyl ether	< LOQ	5000	200	pass		Ethylene glycol	< LOQ	620	200	pass	
Ethylene oxide	< LOQ	50.0	20.0	pass		Hexanes (sum)	< LOQ	290	150	pass	
Isopropyl acetate	< LOQ	5000	200	pass		Isopropylbenzene (Cumene)	< LOQ	70.0	30.0	pass	
m,p-Xylene	< LOQ		200			Methanol	< LOQ	3000	200	pass	
Methylene chloride	< LOQ	600	60.0	pass		Methylpropane (Isobutane)	< LOQ		200		
n-Butane	< LOQ		200			n-Heptane	< LOQ	5000	200	pass	
n-Hexane	< LOQ		30.0			n-Pentane	< LOQ		200		
o-Xylene	< LOQ		200			Pentanes (sum)	< LOQ	5000	600	pass	
Propane	< LOQ	5000	200	pass		Tetrahydrofuran	< LOQ	720	100	pass	
Toluene	< LOQ	890	100	pass		Total Xylenes	< LOQ		400		
Total Xylenes and Ethyl benzene	< LOQ	2170	600	pass							



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Pesticides											
Method: AOAC 2007.01 & EN 15662 (mod) ^b											
Units mg/kg Batch 2300380 Analyze 01/12/23 02:26 PM											
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
Abamectin [‡]	< LOQ	0.50	0.250	pass		Acephate [‡]	< LOQ	0.40	0.200	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 [‡]	< LOQ	0.20	0.100	pass	
Clofentezine [‡]	< LOQ	0.20	0.100	pass		Cyfluthrin [‡]	< LOQ	1.0	0.500	pass	
Cypermethrin [‡]	< LOQ	1.0	0.500	pass		Daminozide [‡]	< LOQ	1.0	0.500	pass	
Diazinon [‡]	< LOQ	0.20	0.100	pass		Dichlorvos [‡]	< LOQ	1.0	0.500	pass	
Dimethoate [‡]	< LOQ	0.20	0.100	pass		Ethoprophos [‡]	< LOQ	0.20	0.100	pass	
Etofenprox [‡]	< LOQ	0.40	0.200	pass		Etoxazole [‡]	< LOQ	0.20	0.100	pass	
Fenoxycarb [‡]	< LOQ	0.20	0.100	pass		Fenpyroximate [‡]	< LOQ	0.40	0.200	pass	
Fipronil [‡]	< LOQ	0.40	0.200	pass		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		Paclobotrazole [‡]	< 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	
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	0.0176	2300511	01/17/23	AOAC 2013.06 (mod.) ^b	pass		
Cadmium	< LOQ	0.200	mg/kg	0.0176	2300511	01/17/23	AOAC 2013.06 (mod.) ^b	pass		
Lead	< LOQ	0.500	mg/kg	0.0176	2300511	01/17/23	AOAC 2013.06 (mod.) ^b	pass		
Mercury	< LOQ	0.100	mg/kg	0.00878	2300511	01/17/23	AOAC 2013.06 (mod.) ^b	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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Report Number: 23-000397/D003.R000
Report Date: 01/20/2023
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Purchase Order:
Received: 01/10/23 15:40

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: 2300380			
Method Blank		Laboratory Control Sample						
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes
Abamectin	0.000	< 0.250		0.894	1.000	89.4	50.0	150
Acephate	0.000	< 0.200		0.648	0.800	81.0	60.0	120
Acequinocyl	0.000	< 1.000		3.361	4.000	84.0	40.0	160
Acetamiprid	0.000	< 0.100		0.350	0.400	87.4	60.0	120
Aldicarb	0.000	< 0.200		0.678	0.800	84.7	60.0	120
Azoxystrobin	0.000	< 0.100		0.353	0.400	88.2	60.0	120
Bifenazate	0.000	< 0.100		0.381	0.400	95.3	60.0	120
Bifenthrin	0.000	< 0.100		0.346	0.400	86.5	50.0	150
Boscalid	0.000	< 0.200		0.705	0.800	88.2	60.0	120
Carbaryl	0.000	< 0.100		0.353	0.400	88.2	60.0	120
Carbofuran	0.000	< 0.100		0.369	0.400	92.3	60.0	120
Chlorantraniliprole	0.000	< 0.100		0.337	0.400	84.4	60.0	120
Chlorfenapyr	0.000	< 0.500		1.761	2.000	88.1	60.0	120
Chlorpyrifos	0.000	< 0.100		0.376	0.400	94.0	60.0	120
Clofentazine	0.000	< 0.100		0.292	0.400	73.1	60.0	120
Cyfluthrin	0.000	< 0.500		1.663	2.000	83.1	50.0	150
Cypermethrin	0.000	< 0.500		1.753	2.000	87.6	50.0	150
Daminozide	0.000	< 0.500		0.718	2.000	35.9	60.0	120
Diazinon	0.000	< 0.100		0.349	0.400	87.4	60.0	120
Dichlorvos	0.000	< 0.500		1.590	2.000	79.5	60.0	120
Dimethoate	0.000	< 0.100		0.357	0.400	89.2	60.0	120
Ethoprophos	0.000	< 0.100		0.353	0.400	88.3	60.0	120
Etofenprox	0.000	< 0.200		0.688	0.800	86.0	50.0	150
Etoxazole	0.000	< 0.100		0.345	0.400	86.4	60.0	120
Fenoxycarb	0.000	< 0.100		0.341	0.400	85.4	60.0	120
Fenpyroximate	0.000	< 0.200		0.700	0.800	87.5	60.0	120
Fipronil	0.000	< 0.200		0.702	0.800	87.8	60.0	120
Fonicamid	0.000	< 0.250		0.936	1.000	93.6	60.0	120
Fludioxonil	0.000	< 0.200		0.766	0.800	95.8	50.0	150
Hexythiazox	0.000	< 0.250		0.873	1.000	87.3	60.0	120
Imazalil	0.000	< 0.100		0.357	0.400	89.2	60.0	120
Imidacloprid	0.000	< 0.200		0.697	0.800	87.1	60.0	120
Kresoxim-methyl	0.000	< 0.200		0.700	0.800	87.5	60.0	120
Malathion	0.000	< 0.100		0.339	0.400	84.8	60.0	120
Metaxalyl	0.000	< 0.100		0.353	0.400	88.1	60.0	120
Methiocarb	0.000	< 0.100		0.348	0.400	87.1	60.0	120
Methomyl	0.000	< 0.200		0.746	0.800	93.2	60.0	120
MGK-264	0.000	< 0.100		0.338	0.400	84.5	50.0	150
Myclobutanil	0.000	< 0.100		0.347	0.400	86.7	60.0	120
Naled	0.000	< 0.250		0.840	1.000	84.0	50.0	150
Oxamyl	0.000	< 0.500		1.759	2.000	88.0	60.0	120
Paclbutrazole	0.000	< 0.200		0.660	0.800	82.5	60.0	120
Parathion-Methyl	0.000	< 0.100		0.373	0.400	93.2	50.0	150
Permethrin	0.000	< 0.100		0.347	0.400	86.8	50.0	150
Phosmet	0.000	< 0.100		0.344	0.400	85.9	50.0	150
Piperonyl butoxide	0.000	< 0.500		1.672	2.000	83.6	60.0	120
Prallethrin	0.000	< 0.100		0.337	0.400	84.3	60.0	120
Propiconazole	0.001	< 0.200		0.674	0.800	84.3	60.0	120
Propoxur	0.000	< 0.100		0.352	0.400	88.1	60.0	120
Pyrethrin (Summe)	0.000	< 0.100		0.406	0.488	83.3	60.0	120
Pyridaben	0.000	< 0.100		0.338	0.400	84.4	50.0	150
Spinosad	0.000	< 0.100		0.334	0.388	86.2	50.0	150
Spiromesifen	0.000	< 0.100		0.360	0.400	90.1	60.0	120
Spirotetramat	0.000	< 0.100		0.354	0.400	88.5	60.0	120
Spiroxamine	0.000	< 0.200		0.670	0.800	83.8	60.0	120
Tebuconazole	0.000	< 0.200		0.664	0.800	83.0	60.0	120
Thiacloprid	0.000	< 0.100		0.347	0.400	86.8	60.0	120
Thiamethoxam	0.000	< 0.100		0.368	0.400	91.9	60.0	120
Trifloxystrobin	0.000	< 0.100		0.348	0.400	87.0	60.0	120

Q6



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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: 2300380				
Matrix Spike/Matrix Spike Duplicate Recoveries						Sample ID: 23-000399-0001					
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes	
Abamectin	0.000	0.833	0.935	1.000	11.6%	< 30	83.3%	93.5%	50 - 150		
Acephate	0.000	0.573	0.629	0.800	9.3%	< 30	71.7%	78.7%	50 - 150		
Acetaminocyl	0.000	3.145	3.285	4.000	4.3%	< 30	78.6%	82.1%	50 - 150		
Acetamiprid	0.000	0.330	0.346	0.400	4.7%	< 30	82.4%	86.4%	50 - 150		
Aldicarb	0.000	0.643	0.672	0.800	4.3%	< 30	80.4%	84.0%	50 - 150		
Azoxystrobin	0.000	0.316	0.322	0.400	1.8%	< 30	79.1%	80.5%	50 - 150		
Bifenazate	0.000	0.335	0.348	0.400	3.7%	< 30	83.7%	86.9%	50 - 150		
Bifenthrin	0.000	0.324	0.331	0.400	2.3%	< 30	80.9%	82.8%	50 - 150		
Boscalid	0.000	0.619	0.637	0.800	2.9%	< 30	77.4%	79.6%	50 - 150		
Carbaryl	0.000	0.329	0.339	0.400	3.0%	< 30	82.3%	84.8%	50 - 150		
Carbofuran	0.000	0.343	0.354	0.400	3.1%	< 30	85.8%	88.5%	50 - 150		
Chlorantraniliprole	0.000	0.324	0.338	0.400	4.3%	< 30	81.0%	84.6%	50 - 150		
Chlorfenapyr	0.000	1.726	1.845	2.000	6.7%	< 30	86.3%	92.2%	50 - 150		
Chlorpyrifos	0.000	0.328	0.342	0.400	4.1%	< 30	82.0%	85.4%	50 - 150		
Clofentezine	0.000	0.049	0.052	0.400	4.2%	< 30	12.4%	12.9%	50 - 150	Q	
Cyfluthrin	0.000	1.551	1.647	2.000	6.0%	< 30	77.5%	82.4%	30 - 150		
Cypermethrin	0.000	1.689	1.720	2.000	1.8%	< 30	84.4%	86.0%	50 - 150		
Daminozide	0.000	0.629	0.665	2.000	5.6%	< 30	31.4%	33.2%	30 - 150		
Diazinon	0.000	0.308	0.311	0.400	0.8%	< 30	77.1%	77.6%	50 - 150		
Dichlorvos	0.000	1.526	1.539	2.000	0.9%	< 30	76.3%	76.9%	50 - 150		
Dimethoate	0.000	0.335	0.349	0.400	4.1%	< 30	83.7%	87.2%	50 - 150		
Ethoprophos	0.000	0.319	0.336	0.400	5.3%	< 30	79.8%	84.1%	50 - 150		
Etofenprox	0.000	0.641	0.663	0.800	3.5%	< 30	80.1%	82.9%	50 - 150		
Etoxazole	0.000	0.315	0.337	0.400	6.5%	< 30	78.9%	84.1%	50 - 150		
Fenoxycarb	0.000	0.307	0.322	0.400	4.6%	< 30	76.9%	80.5%	50 - 150		
Fenpyroximate	0.000	0.730	0.773	0.800	5.7%	< 30	91.2%	96.6%	50 - 150		
Fipronil	0.000	0.658	0.683	0.800	3.7%	< 30	82.3%	85.3%	50 - 150		
Flonicamid	0.000	0.919	0.961	1.000	4.5%	< 30	91.9%	96.1%	50 - 150		
Fludioxonil	0.000	0.729	0.693	0.800	5.0%	< 30	91.1%	86.6%	50 - 150		
Hexythiazox	0.000	0.794	0.812	1.000	2.2%	< 30	79.4%	81.2%	50 - 150		
Imazalil	0.000	0.330	0.348	0.400	5.4%	< 30	82.4%	87.0%	50 - 150		
Imidacloprid	0.000	0.648	0.680	0.800	4.8%	< 30	81.0%	84.9%	50 - 150		
Kresoxim-methyl	0.000	0.634	0.661	0.800	4.1%	< 30	79.3%	82.6%	50 - 150		
Malathion	0.000	0.317	0.327	0.400	3.0%	< 30	79.3%	81.8%	50 - 150		
Metaxalyl	0.000	0.323	0.330	0.400	2.2%	< 30	80.7%	82.6%	50 - 150		
Methiocarb	0.000	0.326	0.340	0.400	4.3%	< 30	81.4%	85.0%	50 - 150		
Methomyl	0.000	0.707	0.747	0.800	5.5%	< 30	88.4%	93.4%	50 - 150		
MGK-264	0.000	0.308	0.318	0.400	3.0%	< 30	77.0%	79.4%	50 - 150		
Myclobutanil	0.000	0.317	0.338	0.400	6.6%	< 30	79.2%	84.6%	50 - 150		
Naled	0.000	0.795	0.820	1.000	3.1%	< 30	79.5%	82.0%	50 - 150		
Oxamyl	0.000	1.734	1.813	2.000	4.4%	< 30	86.7%	90.6%	50 - 150		
Pacllobutrazole	0.000	0.607	0.639	0.800	5.1%	< 30	75.9%	79.9%	50 - 150		
Parathion-Methyl	0.000	0.317	0.306	0.400	3.7%	< 30	79.4%	76.4%	30 - 150		
Permethrin	0.000	0.319	0.330	0.400	3.5%	< 30	79.6%	82.5%	50 - 150		
Phosmet	0.000	0.322	0.326	0.400	1.4%	< 30	80.4%	81.6%	50 - 150		
Piperonyl butoxide	0.000	1.580	1.611	2.000	1.9%	< 30	79.0%	80.5%	50 - 150		
Prallethrin	0.000	0.307	0.320	0.400	4.1%	< 30	76.8%	80.0%	50 - 150		
Propiconazole	0.000	0.619	0.644	0.800	3.9%	< 30	77.4%	80.5%	50 - 150		
Propoxur	0.000	0.328	0.336	0.400	2.5%	< 30	82.0%	84.1%	50 - 150		
Pyrethrin (Summe)	0.000	0.337	0.353	0.488	4.5%	< 30	69.1%	72.4%	50 - 150		
Pyridaben	0.000	0.300	0.314	0.400	4.5%	< 30	74.9%	78.4%	50 - 150		
Spirosad	0.000	0.318	0.336	0.388	5.5%	< 30	82.0%	86.7%	50 - 150		
Spiromesifen	0.000	0.341	0.349	0.400	2.2%	< 30	85.3%	87.2%	50 - 150		
Spirotetramat	0.000	0.326	0.335	0.400	2.5%	< 30	81.6%	83.7%	50 - 150		
Spiroxamine	0.000	0.638	0.640	0.800	0.3%	< 30	79.8%	80.1%	50 - 150		
Tebuconazole	0.000	0.610	0.639	0.800	4.7%	< 30	76.2%	79.8%	50 - 150		
Thiacloprid	0.000	0.341	0.348	0.400	2.3%	< 30	85.2%	87.1%	50 - 150		
Thiamethoxam	0.000	0.344	0.352	0.400	2.1%	< 30	86.1%	87.9%	50 - 150		
Trifloxystrobin	0.000	0.320	0.333	0.400	4.0%	< 30	80.1%	83.4%	50 - 150		

Test results relate only to the parameters tested and to the samples as received by the laboratory. Test results meet all requirements of NELAP and the Columbia Laboratories quality assurance plan unless otherwise noted. This report shall not be reproduced, except in full, without the written consent of this laboratory. Samples will be retained for a maximum of 30 days from the receipt date unless prior arrangements have been made.

Testing in accordance with: OAR 333-007-0390 OAR 333-007-0400 OAR 333-007-0410 OAR 333-007-0430



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



Report Number: 23-000397/D003.R000
Report Date: 01/20/2023
ORELAP#: OR100028
Purchase Order:
Received: 01/10/23 15:40

Revision: 1 Document ID: 7148
Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

J AOAC 2015 V98-6 Batch ID: 2300463

Laboratory Control Sample									
Analyte	LCS	Result	Spike	Units	% Rec	Limits		Evaluation	Notes
CBDVA	2	0.0334	0.033	%	99.9	80.0	- 120	Acceptable	
CBDV	2	0.0354	0.035	%	100	80.0	- 120	Acceptable	
CBE	2	0.0347	0.035	%	99.1	80.0	- 120	Acceptable	
CBDA	1	0.0312	0.032	%	97.4	90.0	- 110	Acceptable	
CBGA	1	0.0315	0.032	%	98.2	80.0	- 120	Acceptable	
CBG	1	0.0324	0.033	%	98.4	80.0	- 120	Acceptable	
CBD	1	0.0313	0.032	%	96.5	90.0	- 110	Acceptable	
THCV	2	0.0351	0.035	%	99.1	80.0	- 120	Acceptable	
d8THCV	2	0.0348	0.034	%	101	80.0	- 120	Acceptable	
THCVA	2	0.0328	0.033	%	99.5	80.0	- 120	Acceptable	
CBN	1	0.0335	0.034	%	98.7	80.0	- 120	Acceptable	
exo-THC	2	0.0324	0.032	%	100	80.0	- 120	Acceptable	
d9THC	1	0.0352	0.035	%	101	90.0	- 110	Acceptable	
d8THC	1	0.0319	0.033	%	95.3	90.0	- 110	Acceptable	
CBL	2	0.0346	0.035	%	100	80.0	- 120	Acceptable	
d10THC	1	0.0154	0.016	%	97.6	80.0	- 120	Acceptable	
CBC	2	0.0342	0.035	%	98.8	80.0	- 120	Acceptable	
THCA	1	0.0317	0.032	%	100	90.0	- 110	Acceptable	
CBCA	2	0.0342	0.034	%	99.4	80.0	- 120	Acceptable	
CBLA	2	0.0346	0.035	%	99.1	80.0	- 120	Acceptable	
CBT	2	0.0344	0.035	%	97.9	80.0	- 120	Acceptable	

Method Blank						
Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.0006	%	< 0.0006	Acceptable	
CBDV	<LOQ	0.0006	%	< 0.0006	Acceptable	
CBE	<LOQ	0.0006	%	< 0.0006	Acceptable	
CBDA	<LOQ	0.0006	%	< 0.0006	Acceptable	
CBGA	<LOQ	0.0006	%	< 0.0006	Acceptable	
CBG	<LOQ	0.0006	%	< 0.0006	Acceptable	
CBD	<LOQ	0.0006	%	< 0.0006	Acceptable	
THCV	<LOQ	0.0006	%	< 0.0006	Acceptable	
d8THCV	<LOQ	0.0006	%	< 0.0006	Acceptable	
THCVA	<LOQ	0.0006	%	< 0.0006	Acceptable	
CBN	<LOQ	0.0006	%	< 0.0006	Acceptable	
exo-THC	<LOQ	0.0006	%	< 0.0006	Acceptable	
d9THC	<LOQ	0.0006	%	< 0.0006	Acceptable	
d8THC	<LOQ	0.0006	%	< 0.0006	Acceptable	
CBL	<LOQ	0.0006	%	< 0.0006	Acceptable	
d10THC	<LOQ	0.0006	%	< 0.0006	Acceptable	
CBC	<LOQ	0.0006	%	< 0.0006	Acceptable	
THCA	<LOQ	0.0006	%	< 0.0006	Acceptable	
CBCA	<LOQ	0.0006	%	< 0.0006	Acceptable	
CBLA	<LOQ	0.0006	%	< 0.0006	Acceptable	
CBT	<LOQ	0.0006	%	< 0.0006	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
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503-254-1794



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Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

J AOAC 2015 V98-6		Batch ID: 2300463						
Sample Duplicate		Sample ID: 23-000249-0001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBDV	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBE	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBDA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBG	0.0172	0.0172	0.003	%	0.159	< 20	Acceptable	
CBD	0.587	0.590	0.003	%	0.512	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBN	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
d9THC	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
d8THC	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
d10THC	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBC	0.00543	0.00546	0.003	%	0.548	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBCA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBT	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	

Abbreviations

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

Units of Measure:



12423 NE Whitaker Way
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 503-254-1794

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

Laboratory Quality Control Results

Residual Solvents				Batch ID: 2300638					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		545	572	µg/g	95.3	60 - 120	
Isobutane	ND	< 200		640	731	µg/g	87.6	60 - 120	
Butane	ND	< 200		625	731	µg/g	85.5	60 - 120	
2,2-Dimethylpropane	ND	< 200		839	936	µg/g	89.6	60 - 120	
Methanol	ND	< 200		1680	1620	µg/g	103.7	60 - 120	
Ethylene Oxide	ND	< 30		49.8	56.2	µg/g	88.6	60 - 120	
2-Methylbutane	ND	< 200		1400	1610	µg/g	87.0	60 - 120	
Pentane	ND	< 200		1400	1600	µg/g	87.5	60 - 120	
Ethanol	ND	< 200		1180	1610	µg/g	73.3	70 - 130	
Ethyl Ether	ND	< 200		1410	1630	µg/g	86.5	60 - 120	
2,2-Dimethylbutane	ND	< 30		161	171	µg/g	94.2	60 - 120	
Acetone	ND	< 200		1510	1630	µg/g	92.6	60 - 120	
2-Propanol	ND	< 200		1520	1620	µg/g	93.8	60 - 120	
Ethyl Formate	ND	< 500		1600	1670	µg/g	95.8	70 - 130	
Acetonitrile	ND	< 100		435	498	µg/g	87.3	60 - 120	
Methyl Acetate	ND	< 500		1640	1730	µg/g	94.8	70 - 130	
2,3-Dimethylbutane	ND	< 30		151	171	µg/g	88.3	60 - 120	
Dichloromethane	ND	< 60		433	483	µg/g	89.6	60 - 120	
2-Methylpentane	ND	< 30		143	168	µg/g	85.1	60 - 120	
MTBE	ND	< 500		1500	1650	µg/g	90.9	70 - 130	
3-Methylpentane	ND	< 30		133	167	µg/g	79.6	60 - 120	
Hexane	ND	< 30		202	182	µg/g	111.0	60 - 120	
1-Propanol	ND	< 500		1400	1620	µg/g	86.4	70 - 130	
Methylethylketone	ND	< 500		1610	1620	µg/g	99.4	70 - 130	
Ethyl acetate	ND	< 200		1520	1610	µg/g	94.4	60 - 120	
2-Butanol	ND	< 200		1450	1600	µg/g	90.6	60 - 120	
Tetrahydrofuran	ND	< 100		386	483	µg/g	79.9	60 - 120	
Cyclohexane	ND	< 200		1520	1610	µg/g	94.4	60 - 120	
2-methyl-1-propanol	ND	< 500		1250	1620	µg/g	77.2	70 - 130	
Benzene	ND	< 1		4.05	5.02	µg/g	80.7	60 - 120	
Isopropyl Acetate	ND	< 200		1430	1620	µg/g	88.3	60 - 120	
Heptane	ND	< 200		1580	1610	µg/g	98.1	60 - 120	
1-Butanol	ND	< 500		1010	1630	µg/g	62.0	70 - 130	Q6
Propyl Acetate	ND	< 500		1290	1610	µg/g	80.1	70 - 130	
1,4-Dioxane	ND	< 100		415	491	µg/g	84.5	60 - 120	
2-Ethoxyethanol	ND	< 30		112	181	µg/g	61.9	60 - 120	
Methylisobutylketone	ND	< 500		1420	1620	µg/g	87.7	70 - 130	
3-Methyl-1-butanol	ND	< 500		1220	1630	µg/g	74.8	70 - 130	
Ethylene Glycol	ND	< 200		513	484	µg/g	106.0	60 - 120	
Toluene	ND	< 100		400	485	µg/g	82.5	60 - 120	
Isobutyl Acetate	ND	< 500		1550	1630	µg/g	95.1	70 - 130	
1-Pentanol	ND	< 500		1230	1620	µg/g	75.9	70 - 130	
Butyl Acetate	ND	< 500		1390	1620	µg/g	85.8	70 - 130	
Ethylbenzene	ND	< 200		838	969	µg/g	86.5	60 - 120	
m,p-Xylene	ND	< 200		853	994	µg/g	85.8	60 - 120	
o-Xylene	ND	< 200		770	967	µg/g	79.6	60 - 120	
Cumene	ND	< 30		129	171	µg/g	75.4	60 - 120	
Anisole	ND	< 500		1010	1630	µg/g	62.0	70 - 130	Q6
DMSO	ND	< 500		1410	1680	µg/g	83.9	70 - 130	
1,2-dimethoxyethane	ND	< 50		159	169	µg/g	94.1	70 - 130	
Triethylamine	ND	< 500		1540	1630	µg/g	94.5	70 - 130	
N,N-dimethylformamide	ND	< 150		459	482	µg/g	95.2	70 - 130	
N,N-dimethylacetamide	ND	< 150		333	510	µg/g	65.3	70 - 130	Q6
Pyridine	ND	< 50		174	203	µg/g	85.7	70 - 130	
Sulfone	ND	< 50		133	172	µg/g	77.3	70 - 130	
1,2-Dichloroethane	ND	< 1		0.801	1	µg/g	80.1	70 - 130	
Chloroform	ND	< 1		0.83	1	µg/g	83.0	70 - 130	
Trichloroethylene	ND	< 1		0.825	1	µg/g	82.5	70 - 130	
1,1-Dichloroethane	ND	< 1		0.779	1	µg/g	77.9	70 - 130	



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Received: 01/10/23 15:40

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QC - Sample Duplicate		Sample ID: 23-000249-0001						
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
 Q6 - Quality control outside QC limits. Data acceptable based on remaining QC.

Units of Measure:

µg/g- Microgram per gram or ppm



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



Report Number: 23-000397/D003.R000
Report Date: 01/20/2023
ORELAP#: OR100028
Purchase Order:
Received: 01/10/23 15:40





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.