

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.



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



Report Number: 23-006209/D005.R000
Report Date: 06/07/2023
ORELAP#: OR100028
Purchase Order:
Received: 05/24/23 16:17

Customer: Etz Hayim Holdings
Product identity: FORM-TN.FS.SAL20-FE06
Client/Metric ID: .
Laboratory ID: 23-006209-0001

Summary

Potency:

Analyte per 1g	Result	Limits	Units	Status	
CBC per 1g	0.914		mg/1g		CBD-Total per Serving Size 21.7 mg/1g
CBD per 1g	21.7		mg/1g		
CBDV per 1g	0.131		mg/1g		THC-Total per Serving Size 0.594 mg/1g
CBE per 1g	0.360		mg/1g		(Reported in milligrams per serving)
CBG per 1g	0.332		mg/1g		
CBN per 1g	0.0727		mg/1g		
CBT per 1g	0.466		mg/1g		
Δ9-THC per 1g	0.594		mg/1g		

Residual Solvents:

All analytes passing and less than LOQ.

Pesticides:

All analytes passing and less than LOQ.

Metals:

Analyte	Result	Units	Limit	Status
Arsenic*	1.57	mg/kg		

Microbiology:

Less than LOQ for all analytes.



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

Product identity: FORM-TN.FS.SAL20-FE06

Client/Metric ID: .

Sample Date:

Laboratory ID: 23-006209-0001

Evidence of Cooling: No

Temp: 25.3

Relinquished by: courier

Serving Size #1: 1 g

Sample Results

Potency per 1g	Method: J AOAC 2015 V98-6 (mod) ^b	Units mg/se	Batch: 2307675	Analyze: 5/26/23 3:10:00 AM	
Analyte	Result	Limits	Units	LOQ	Notes
CBC per 1g	0.914		mg/1g	0.0305	
CBC-A per 1g	< LOQ		mg/1g	0.0305	
CBC-Total per 1g	0.914		mg/1g	0.0572	
CBD per 1g	21.7		mg/1g	0.305	
CBD-A per 1g	< LOQ		mg/1g	0.0305	
CBD-Total per 1g	21.7		mg/1g	0.332	
CBDV per 1g	0.131		mg/1g	0.0305	
CBDV-A per 1g	< LOQ		mg/1g	0.0305	
CBDV-Total per 1g	0.131		mg/1g	0.0569	
CBE per 1g	0.360		mg/1g	0.0305	
CBG per 1g	0.332		mg/1g	0.0305	
CBG-A per 1g	< LOQ		mg/1g	0.0305	
CBG-Total per 1g	0.332		mg/1g	0.0569	
CBL per 1g	< LOQ		mg/1g	0.0305	
CBL-A per 1g	< LOQ		mg/1g	0.0305	
CBL-Total per 1g	< LOQ		mg/1g	0.0572	
CBN per 1g	0.0727		mg/1g	0.0305	
CBT per 1g	0.466		mg/1g	0.0305	
Δ8-THCV per 1g	< LOQ		mg/1g	0.0305	
Δ10-THC-9R per 1g	< LOQ		mg/1g	0.0305	
Δ10-THC-9S per 1g	< LOQ		mg/1g	0.0305	
Δ10-THC-Total per 1g	< LOQ		mg/1g	0.0610	
Δ8-THC per 1g	< LOQ		mg/1g	0.0305	
Δ9-THC per 1g	0.594		mg/1g	0.0305	
delta-9-THCP per 1g	< LOQ		mg/1g	0.0305	
exo-THC per 1g	< LOQ		mg/1g	0.0305	
THC-A per 1g	< LOQ		mg/1g	0.0305	
THC-Total per 1g	0.594		mg/1g	0.0572	
THCV per 1g	< LOQ		mg/1g	0.0305	
THCV-A per 1g	< LOQ		mg/1g	0.0305	



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Potency per 1g	Method: J AOAC 2015 V98-6 (mod) ^P	Units mg/se	Batch: 2307675	Analyze: 5/26/23 3:10:00 AM	
Analyte	Result	Limits	Units	LOQ	Notes
THCV-Total per 1g	< LOQ		mg/1g	0.0573	
Total Cannabinoids per 1g	24.6		mg/1g		

Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2307783	06/02/23 AOAC 991.14 (Petrifilm) ^P		
Total Coliforms	< LOQ		cfu/g	10	2307783	06/02/23 AOAC 991.14 (Petrifilm) ^P		
Mold (RAPID Petrifilm)	< LOQ		cfu/mL	10	2307784	06/02/23 AOAC 2014.05 (RAPID) ^P		
Yeast (RAPID Petrifilm)	< LOQ		cfu/mL	10	2307784	06/02/23 AOAC 2014.05 (RAPID) ^P		

Solvents

Method: Residual Solvents by GC/MS ^P						Units µg/g	Batch 2307994	Analyze 06/07/23 02:18 PM			
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
1,4-Dioxane	< LOQ	380	100	pass		2-Butanol	< LOQ	5000	200	pass	
2-Ethoxyethanol	< LOQ	160	30.0	pass		2-Methylbutane (Isopentane)	< LOQ		200		
2-Methylpentane	< LOQ		30.0			2-Propanol (IPA)	< LOQ	5000	200	pass	
2,2-Dimethylbutane	< LOQ		30.0			2,2-Dimethylpropane (neo-pentane)	< LOQ		200		
2,3-Dimethylbutane	< LOQ		30.0			3-Methylpentane	< LOQ		30.0		
Acetone	< LOQ	5000	200	pass		Acetonitrile	< LOQ	410	100	pass	
Benzene	< LOQ	2.00	1.00	pass		Butanes (sum)	< LOQ	5000	400	pass	
Cyclohexane	< LOQ	3880	200	pass		Ethanol	< LOQ		200		
Ethyl acetate	< LOQ	5000	200	pass		Ethyl benzene	< LOQ		200		
Ethyl ether	< LOQ	5000	200	pass		Ethylene glycol	< LOQ	620	200	pass	
Ethylene oxide	< LOQ	50.0	20.0	pass		Hexanes (sum)	< LOQ	290	150	pass	
Isopropyl acetate	< LOQ	5000	200	pass		Isopropylbenzene (Cumene)	< LOQ	70.0	30.0	pass	
m,p-Xylene	< LOQ		200			Methanol	< LOQ	3000	200	pass	
Methylene chloride	< LOQ	600	60.0	pass		Methylpropane (Isobutane)	< LOQ		200		
n-Butane	< LOQ		200			n-Heptane	< LOQ	5000	200	pass	
n-Hexane	< LOQ		30.0			n-Pentane	< LOQ		200		
o-Xylene	< LOQ		200			Pentanes (sum)	< LOQ	5000	600	pass	
Propane	< LOQ	5000	200	pass		Tetrahydrofuran	< LOQ	720	100	pass	
Toluene	< LOQ	890	100	pass		Total Xylenes	< LOQ		400		
Total Xylenes and Ethyl benzene	< LOQ	2170	600	pass							



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Pesticides											
Method: AOAC 2007.01 & EN 15662 (mod) ^b											
Units mg/kg Batch 2307861 Analyze 06/02/23 12:29 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		Pacllobutrazole [‡]	< 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 [‡]	1.57		mg/kg	0.0964	2307833	06/01/23 AOAC 2013.06 (mod.) ^b				
Cadmium [‡]	< LOQ		mg/kg	0.0964	2307833	06/01/23 AOAC 2013.06 (mod.) ^b				
Lead [‡]	< LOQ		mg/kg	0.0964	2307833	06/01/23 AOAC 2013.06 (mod.) ^b				
Mercury [‡]	< LOQ		mg/kg	0.0482	2307833	06/01/23 AOAC 2013.06 (mod.) ^b				



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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

cfu/mL = Colony forming units per milliliter

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

Laboratory Quality Control Results

J AOAC 2015 V98-6 Batch ID: 2307675

Laboratory Control Sample									
Analyte	LCS	Result	Spike	Units	% Rec	Limits		Evaluation	Notes
CBDVA	2	0.0285	0.0283	%	101	80.0	- 120	Acceptable	
CBDV	2	0.0299	0.0291	%	102	80.0	- 120	Acceptable	
CBE	2	0.0339	0.0344	%	98.6	80.0	- 120	Acceptable	
CBDA	1	0.0297	0.0311	%	95.3	90.0	- 110	Acceptable	
CBGA	1	0.0311	0.0311	%	100	80.0	- 120	Acceptable	
CBG	1	0.0322	0.0322	%	100	80.0	- 120	Acceptable	
CBD	1	0.0319	0.0323	%	98.8	90.0	- 110	Acceptable	
THCV	2	0.0207	0.0201	%	103	80.0	- 120	Acceptable	
d8THCV	2	0.0265	0.0268	%	98.7	80.0	- 120	Acceptable	
THCVA	2	0.0303	0.0299	%	101	80.0	- 120	Acceptable	
CBN	1	0.0334	0.0329	%	102	80.0	- 120	Acceptable	
exo-THC	2	0.0293	0.0292	%	100	80.0	- 120	Acceptable	
d9THC	1	0.0341	0.0341	%	99.9	90.0	- 110	Acceptable	
d8THC	1	0.0427	0.0420	%	101	90.0	- 110	Acceptable	
9S-d10THC	1	0.0241	0.0240	%	100	80.0	- 120	Acceptable	
CBL	2	0.0323	0.0315	%	103	80.0	- 120	Acceptable	
9R-d10THC	1	0.0309	0.0310	%	99.5	80.0	- 120	Acceptable	
CBC	2	0.0310	0.0309	%	100	80.0	- 120	Acceptable	
THCA	1	0.0312	0.0314	%	99.4	90.0	- 110	Acceptable	
CBCA	2	0.0331	0.0326	%	101	80.0	- 120	Acceptable	
CBLA	2	0.0331	0.0331	%	100.0	80.0	- 120	Acceptable	
d9THCP	2	0.0323	0.0321	%	101	80.0	- 120	Acceptable	
CBT	2	0.0329	0.0327	%	101	80.0	- 120	Acceptable	

Method Blank							
Analyte	Result	LOQ	Units	Limits		Evaluation	Notes
CBDVA	<LOQ	0.00314	%	< 0.00314		Acceptable	
CBDV	<LOQ	0.00314	%	< 0.00314		Acceptable	
CBE	<LOQ	0.00314	%	< 0.00314		Acceptable	
CBDA	<LOQ	0.00314	%	< 0.00314		Acceptable	
CBGA	<LOQ	0.00314	%	< 0.00314		Acceptable	
CBG	<LOQ	0.00314	%	< 0.00314		Acceptable	
CBD	<LOQ	0.00314	%	< 0.00314		Acceptable	
THCV	<LOQ	0.00314	%	< 0.00314		Acceptable	
d8THCV	<LOQ	0.00314	%	< 0.00314		Acceptable	
THCVA	<LOQ	0.00314	%	< 0.00314		Acceptable	
CBN	<LOQ	0.00314	%	< 0.00314		Acceptable	
exo-THC	<LOQ	0.00314	%	< 0.00314		Acceptable	
d9THC	<LOQ	0.00314	%	< 0.00314		Acceptable	
d8THC	<LOQ	0.00314	%	< 0.00314		Acceptable	
9S-d10THC	<LOQ	0.00314	%	< 0.00314		Acceptable	
CBL	<LOQ	0.00314	%	< 0.00314		Acceptable	
9R-d10THC	<LOQ	0.00314	%	< 0.00314		Acceptable	
CBC	<LOQ	0.00314	%	< 0.00314		Acceptable	
THCA	<LOQ	0.00314	%	< 0.00314		Acceptable	
CBCA	<LOQ	0.00314	%	< 0.00314		Acceptable	
CBLA	<LOQ	0.00314	%	< 0.00314		Acceptable	
d9THCP	<LOQ	0.00314	%	< 0.00314		Acceptable	
CBT	<LOQ	0.00314	%	< 0.00314		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: 2307675						
Sample Duplicate		Sample ID: 23-006168-0001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
CBDV	0.0389	0.0399	0.00327	%	2.58	< 20	Acceptable	
CBE	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
CBD	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
CBDA	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
CBG	0.470	0.472	0.00327	%	0.505	< 20	Acceptable	
CBD	16.9	16.6	0.00327	%	1.90	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
CBN	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
d9THC	0.00743	0.00724	0.00327	%	2.58	< 20	Acceptable	
d8THC	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
9S-d10THC	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
9R-d10THC	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
CBC	0.139	0.140	0.00327	%	0.454	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
CBCA	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
d9THCP	<LOQ	<LOQ	0.00327	%	NA	< 20	Acceptable	
CBT	0.00882	0.00879	0.00327	%	0.319	< 20	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 Pesticide Quality Control Results

AOAC 2007.1 & EN 15662		Units: mg/Kg			Batch ID: 2307861			
Method Blank		Laboratory Control Sample						
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes
Abamectin	0.000	< 0.250		1.059	1.000	105.9	50.0	150
Acephate	0.039	< 0.200		0.657	0.800	82.1	60.0	120
Acequinocyl	0.000	< 1.000		4.176	4.000	104.4	40.0	160
Acetamiprid	0.002	< 0.100		0.380	0.400	95.0	60.0	120
Aldicarb	0.000	< 0.200		0.869	0.800	108.6	60.0	120
Azoxystrobin	0.008	< 0.100		0.380	0.400	95.0	60.0	120
Bifenazate	0.000	< 0.100		0.401	0.400	100.3	60.0	120
Bifenthrin	0.000	< 0.100		0.399	0.400	99.7	50.0	150
Boscalid	0.000	< 0.200		0.785	0.800	98.1	60.0	120
Carbaryl	0.000	< 0.100		0.379	0.400	94.7	60.0	120
Carbofuran	0.000	< 0.100		0.387	0.400	96.7	60.0	120
Chlorantraniliprole	0.000	< 0.100		0.380	0.400	95.1	60.0	120
Chlorfenapyr	0.000	< 0.500		2.083	2.000	104.1	60.0	120
Chlorpyrifos	0.009	< 0.100		0.345	0.400	86.3	60.0	120
Clofentazine	0.000	< 0.100		0.393	0.400	98.3	60.0	120
Cyfluthrin	0.000	< 0.500		1.909	2.000	95.4	50.0	150
Cypermethrin	0.000	< 0.500		1.927	2.000	96.4	50.0	150
Daminozide	0.000	< 0.500		0.734	2.000	36.7	60.0	120
Diazinon	0.000	< 0.100		0.418	0.400	104.5	60.0	120
Dichlorvos	0.011	< 0.500		1.883	2.000	94.2	60.0	120
Dimethoate	0.002	< 0.100		0.374	0.400	93.5	60.0	120
Ethoprophos	0.003	< 0.100		0.376	0.400	94.1	60.0	120
Etofenprox	0.005	< 0.200		0.795	0.800	99.4	50.0	150
Etoxazole	0.000	< 0.100		0.408	0.400	102.0	60.0	120
Fenoxycarb	0.000	< 0.100		0.408	0.400	102.1	60.0	120
Fenpyroximate	0.003	< 0.200		0.807	0.800	100.9	60.0	120
Fipronil	0.000	< 0.200		0.814	0.800	101.7	60.0	120
Fonicamid	0.000	< 0.250		0.921	1.000	92.1	60.0	120
Fludioxonil	0.000	< 0.200		0.754	0.800	94.2	50.0	150
Hexythiazox	0.000	< 0.250		0.960	1.000	96.0	60.0	120
Imazalil	0.000	< 0.100		0.391	0.400	97.8	60.0	120
Imidacloprid	0.022	< 0.200		0.781	0.800	97.6	60.0	120
Kresoxim-methyl	0.000	< 0.200		0.829	0.800	103.7	60.0	120
Malathion	0.000	< 0.100		0.394	0.400	98.4	60.0	120
Metaxalyl	0.000	< 0.100		0.410	0.400	102.5	60.0	120
Methiocarb	0.000	< 0.100		0.405	0.400	101.3	60.0	120
Methomyl	0.000	< 0.200		0.769	0.800	96.2	60.0	120
MGK-264	0.000	< 0.100		0.427	0.400	106.7	50.0	150
Myclobutanil	0.000	< 0.100		0.370	0.400	92.6	60.0	120
Naled	0.000	< 0.250		0.946	1.000	94.6	50.0	150
Oxamyl	0.000	< 0.500		1.785	2.000	89.2	60.0	120
Paclotrazole	0.000	< 0.200		0.810	0.800	101.3	60.0	120
Parathion-Methyl	0.000	< 0.100		0.433	0.400	108.3	50.0	150
Permethrin	0.000	< 0.100		0.378	0.400	94.4	50.0	150
Phosmet	0.000	< 0.100		0.424	0.400	106.0	50.0	150
Piperonyl butoxide	0.000	< 0.500		2.156	2.000	107.8	60.0	120
Prallethrin	0.000	< 0.100		0.429	0.400	107.2	60.0	120
Propiconazole	0.005	< 0.200		0.787	0.800	98.4	60.0	120
Propoxur	0.003	< 0.100		0.369	0.400	92.3	60.0	120
Pyrethrin (Summe)	0.000	< 0.100		0.508	0.488	104.1	60.0	120
Pyridaben	0.004	< 0.100		0.409	0.400	102.3	50.0	150
Spirosad	0.000	< 0.100		0.394	0.388	101.6	50.0	150
Spiromesifen	0.000	< 0.100		0.384	0.400	96.0	60.0	120
Spirotetramat	0.000	< 0.100		0.376	0.400	94.0	60.0	120
Spiroxamine	0.000	< 0.200		0.825	0.800	103.1	60.0	120
Tebuconazole	0.000	< 0.200		0.798	0.800	99.7	60.0	120
Thiacloprid	0.000	< 0.100		0.386	0.400	96.4	60.0	120
Thiamethoxam	0.000	< 0.100		0.386	0.400	96.6	60.0	120
Trifloxystrobin	0.001	< 0.100		0.397	0.400	99.3	60.0	120

Q6



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Revision: 3 Document ID: 3120
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Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662		Units: mg/Kg				Batch ID: 2307861				
Matrix Spike/Matrix Spike Duplicate Recoveries					Sample ID: 23-006209-0001					
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes
Abamectin	0.000	0.508	0.569	1.000	11.3%	< 30	50.8%	56.9%	50 - 150	
Acephate	0.042	0.707	0.686	0.800	3.2%	< 30	83.1%	80.6%	50 - 150	
Acetaminophen	0.000	0.650	0.669	4.000	2.8%	< 30	16.3%	16.7%	50 - 150	Q
Acetamiprid	0.000	0.361	0.364	0.400	0.8%	< 30	90.2%	91.0%	50 - 150	
Aldicarb	0.000	0.831	0.860	0.800	3.5%	< 30	103.9%	107.6%	50 - 150	
Azoxystrobin	0.007	0.315	0.316	0.400	0.5%	< 30	76.9%	77.4%	50 - 150	
Bifenazate	0.010	0.328	0.344	0.400	4.9%	< 30	79.5%	83.4%	50 - 150	
Bifenthrin	0.000	0.000	0.000	0.400	NA	< 30	0.0%	0.0%	50 - 150	Q, R1
Boscalid	0.000	0.674	0.697	0.800	3.4%	< 30	84.3%	87.1%	50 - 150	
Carbaryl	0.000	0.332	0.324	0.400	2.4%	< 30	83.0%	81.1%	50 - 150	
Carbofuran	0.000	0.347	0.347	0.400	0.1%	< 30	86.6%	86.8%	50 - 150	
Chlorantraniliprole	0.000	0.355	0.372	0.400	4.7%	< 30	88.8%	93.0%	50 - 150	
Chlorfenapyr	0.000	1.364	1.038	2.000	27.1%	< 30	68.2%	51.9%	50 - 150	
Chlorpyrifos	0.010	0.328	0.316	0.400	3.8%	< 30	79.5%	76.5%	50 - 150	
Clofentezine	0.001	0.305	0.321	0.400	4.9%	< 30	76.0%	79.8%	50 - 150	
Cyfluthrin	0.000	1.104	1.094	2.000	1.0%	< 30	55.2%	54.7%	30 - 150	
Cypermethrin	0.000	1.028	1.084	2.000	5.3%	< 30	51.4%	54.2%	50 - 150	
Daminozide	0.000	0.687	0.716	2.000	4.0%	< 30	34.4%	35.8%	30 - 150	
Diazinon	0.000	0.303	0.315	0.400	3.8%	< 30	75.8%	78.7%	50 - 150	
Dichlorvos	0.008	1.742	1.749	2.000	0.4%	< 30	86.7%	87.1%	50 - 150	
Dimethoate	0.002	0.370	0.374	0.400	1.3%	< 30	91.8%	93.0%	50 - 150	
Ethoprophos	0.003	0.307	0.314	0.400	2.3%	< 30	75.9%	77.6%	50 - 150	
Etofenprox	0.000	0.298	0.299	0.800	0.5%	< 30	37.2%	37.4%	50 - 150	Q
Etoxazole	0.001	0.298	0.305	0.400	2.2%	< 30	74.4%	76.1%	50 - 150	
Fenoxycarb	0.000	0.328	0.347	0.400	5.9%	< 30	81.9%	86.9%	50 - 150	
Fenpyroximate	0.000	0.334	0.337	0.800	0.8%	< 30	41.7%	42.1%	50 - 150	Q
Fipronil	0.017	0.554	0.562	0.800	1.6%	< 30	67.1%	68.2%	50 - 150	
Fonicamid	0.000	0.874	0.916	1.000	4.7%	< 30	87.4%	91.6%	50 - 150	
Fludioxonil	0.000	0.832	0.817	0.800	1.9%	< 30	104.1%	102.1%	50 - 150	
Hexythiazox	0.015	0.583	0.592	1.000	1.5%	< 30	56.8%	57.6%	50 - 150	
Imazalil	0.007	0.367	0.367	0.400	0.1%	< 30	89.9%	90.0%	50 - 150	
Imidacloprid	0.000	0.769	0.819	0.800	6.4%	< 30	96.1%	102.4%	50 - 150	
Kresoxim-methyl	0.000	0.655	0.658	0.800	0.5%	< 30	81.9%	82.3%	50 - 150	
Malathion	0.007	0.331	0.326	0.400	1.5%	< 30	81.0%	79.8%	50 - 150	
Metaxalyl	0.000	0.367	0.379	0.400	3.1%	< 30	91.8%	94.7%	50 - 150	
Methiocarb	0.000	0.329	0.339	0.400	3.0%	< 30	82.3%	84.8%	50 - 150	
Methomyl	0.000	0.736	0.771	0.800	4.7%	< 30	92.0%	96.3%	50 - 150	
MGK-264	0.002	0.259	0.270	0.400	4.4%	< 30	64.3%	67.1%	50 - 150	
Myclobutanil	0.000	0.358	0.338	0.400	5.9%	< 30	89.6%	84.4%	50 - 150	
Naled	0.000	0.822	0.830	1.000	0.9%	< 30	82.2%	83.0%	50 - 150	
Oxamyl	0.000	1.705	1.802	2.000	5.5%	< 30	85.3%	90.1%	50 - 150	
Pacllobutrazole	0.000	0.671	0.723	0.800	7.4%	< 30	83.9%	90.4%	50 - 150	
Parathion-Methyl	0.000	0.304	0.295	0.400	3.1%	< 30	76.0%	73.7%	30 - 150	
Permethrin	0.000	0.127	0.128	0.400	0.5%	< 30	31.7%	31.9%	50 - 150	Q
Phosmet	0.000	0.357	0.364	0.400	1.7%	< 30	89.3%	90.9%	50 - 150	
Piperonyl butoxide	0.000	1.669	1.650	2.000	1.1%	< 30	83.4%	82.5%	50 - 150	
Prallethrin	0.139	0.460	0.493	0.400	9.9%	< 30	80.0%	88.4%	50 - 150	
Propiconazole	0.006	0.652	0.694	0.800	6.4%	< 30	80.7%	86.1%	50 - 150	
Propoxur	0.003	0.350	0.343	0.400	2.1%	< 30	86.6%	84.8%	50 - 150	
Pyrethrin (Summe)	0.000	0.374	0.387	0.488	3.5%	< 30	76.6%	79.3%	50 - 150	
Pyridaben	0.004	0.209	0.222	0.400	6.0%	< 30	51.3%	54.5%	50 - 150	
Spinosad	0.000	0.286	0.304	0.388	5.9%	< 30	73.8%	78.3%	50 - 150	
Spiromesifen	0.008	0.325	0.334	0.400	3.0%	< 30	79.2%	81.7%	50 - 150	
Spirotetramat	0.000	0.429	0.435	0.400	1.4%	< 30	107.2%	108.7%	50 - 150	
Spiroxamine	0.000	0.744	0.771	0.800	3.5%	< 30	93.0%	96.3%	50 - 150	
Tebuconazole	0.000	0.656	0.663	0.800	1.0%	< 30	82.1%	82.9%	50 - 150	
Thiacloprid	0.000	0.376	0.387	0.400	2.8%	< 30	94.1%	96.7%	50 - 150	
Thiamethoxam	0.000	0.380	0.385	0.400	1.4%	< 30	94.9%	96.2%	50 - 150	
Trifloxystrobin	0.001	0.283	0.292	0.400	3.1%	< 30	70.5%	72.7%	50 - 150	



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

Laboratory Quality Control Results

Residual Solvents				Batch ID: 2307994					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		501	584	µg/g	85.8	60 - 120	
Isobutane	ND	< 200		642	767	µg/g	83.7	60 - 120	
Butane	ND	< 200		630	782	µg/g	80.6	60 - 120	
2,2-Dimethylpropane	ND	< 200		886	939	µg/g	94.4	60 - 120	
Methanol	ND	< 200		1340	1640	µg/g	81.7	60 - 120	
Ethylene Oxide	ND	< 30		43.8	57.1	µg/g	76.7	60 - 120	
2-Methylbutane	ND	< 200		1330	1600	µg/g	83.1	60 - 120	
Pentane	ND	< 200		1340	1620	µg/g	82.7	60 - 120	
Ethanol	ND	< 200		1380	1610	µg/g	85.7	70 - 130	
Ethyl Ether	ND	< 200		1380	1610	µg/g	85.7	60 - 120	
2,2-Dimethylbutane	ND	< 30		145	168	µg/g	86.3	60 - 120	
Acetone	ND	< 200		1380	1620	µg/g	85.2	60 - 120	
2-Propanol	ND	< 200		1370	1600	µg/g	85.6	60 - 120	
Ethyl Formate	ND	< 500		1370	1600	µg/g	85.6	70 - 130	
Acetonitrile	ND	< 100		401	484	µg/g	82.9	60 - 120	
Methyl Acetate	ND	< 500		1510	1610	µg/g	93.8	70 - 130	
2,3-Dimethylbutane	ND	< 30		131	162	µg/g	80.9	60 - 120	
Dichloromethane	ND	< 60		417	483	µg/g	86.3	60 - 120	
2-Methylpentane	ND	< 30		153	174	µg/g	87.9	60 - 120	
MTBE	ND	< 500		1550	1610	µg/g	96.3	70 - 130	
3-Methylpentane	ND	< 30		144	168	µg/g	85.7	60 - 120	
Hexane	ND	< 30		140	168	µg/g	83.3	60 - 120	
1-Propanol	ND	< 500		1460	1600	µg/g	91.3	70 - 130	
Methylethylketone	ND	< 500		1520	1620	µg/g	93.8	70 - 130	
Ethyl acetate	ND	< 200		1330	1600	µg/g	83.1	60 - 120	
2-Butanol	ND	< 200		1360	1600	µg/g	85.0	60 - 120	
Tetrahydrofuran	ND	< 100		443	514	µg/g	86.2	60 - 120	
Cyclohexane	ND	< 200		1400	1600	µg/g	87.5	60 - 120	
2-methyl-1-propanol	ND	< 500		1500	1610	µg/g	93.2	70 - 130	
Benzene	ND	< 1		4.05	5.12	µg/g	79.1	60 - 120	
Isopropyl Acetate	ND	< 200		1350	1620	µg/g	83.3	60 - 120	
Heptane	ND	< 200		1330	1610	µg/g	82.6	60 - 120	
1-Butanol	ND	< 500		1420	1600	µg/g	88.8	70 - 130	
Propyl Acetate	ND	< 500		1470	1600	µg/g	91.9	70 - 130	
1,4-Dioxane	ND	< 100		438	493	µg/g	84.8	60 - 120	
2-Ethoxyethanol	ND	< 30		138	163	µg/g	84.7	60 - 120	
Methylisobutylketone	ND	< 500		1640	1600	µg/g	102.5	70 - 130	
3-Methyl-1-butanol	ND	< 500		1650	1610	µg/g	102.5	70 - 130	
Ethylene Glycol	ND	< 200		419	483	µg/g	86.7	60 - 120	
Toluene	ND	< 100		402	493	µg/g	81.5	60 - 120	
Isobutyl Acetate	ND	< 500		1430	1600	µg/g	89.4	70 - 130	
1-Pentanol	ND	< 500		1330	1600	µg/g	83.1	70 - 130	
Butyl Acetate	ND	< 500		1420	1600	µg/g	88.8	70 - 130	
Ethylbenzene	ND	< 200		800	969	µg/g	82.6	60 - 120	
m,p-Xylene	ND	< 200		798	968	µg/g	82.4	60 - 120	
o-Xylene	ND	< 200		800	976	µg/g	82.0	60 - 120	
Cumene	ND	< 30		132	162	µg/g	81.5	60 - 120	
Anisole	ND	< 500		1350	1610	µg/g	83.9	70 - 130	
DMSO	ND	< 500		1270	1610	µg/g	78.9	70 - 130	
1,2-dimethoxyethane	ND	< 50		157	164	µg/g	95.7	70 - 130	
Triethylamine	ND	< 500		1520	1600	µg/g	95.0	70 - 130	
N,N-dimethylformamide	ND	< 150		437	484	µg/g	90.3	70 - 130	
N,N-dimethylacetamide	ND	< 150		371	489	µg/g	75.9	70 - 130	
Pyridine	ND	< 50		153	172	µg/g	89.0	70 - 130	
Sulfolane	ND	< 50		100	163	µg/g	61.3	70 - 130	Q6
1,2-Dichloroethane	ND	< 1		0.788	1	µg/g	78.8	70 - 130	
Chloroform	ND	< 1		0.782	1	µg/g	78.2	70 - 130	
Trichloroethylene	ND	< 1		0.752	1	µg/g	75.2	70 - 130	
1,1-Dichloroethane	ND	< 1		0.753	1	µg/g	75.3	70 - 130	



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QC - Sample Duplicate		Sample ID: 23-006209-0001						
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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Report Number: 23-006209/D005.R000
Report Date: 06/07/2023
ORELAP#: OR100028
Purchase Order:
Received: 05/24/23 16:17





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.