

CONSOLIDATED TEST RESULTS SUMMARY

Please see the following pages for full test results.

BULK SKU	BATCH #	LOQ: Limit Of Quantitation LOD: Limit Of Detection $1 \text{ g} = 10^{-3} \text{ kg} = 10^3 \text{ mg} = 10^6 \text{ } \mu\text{g}$ $1 \text{ mg/kg} = 1 \text{ ppm} = 1000 \text{ ppb}$
PRODUCT NAME	SERVING SIZE	
LABORATORY :	OREGON ACCREDITATION: OR100028	

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	$\mu\text{g/serving}$	$\mu\text{g/g}$	10 $\mu\text{g/day}$ ^[1]
Cadmium	$\mu\text{g/serving}$	$\mu\text{g/g}$	4.1 $\mu\text{g/day}$ ^[1]
Lead	$\mu\text{g/serving}$	$\mu\text{g/g}$	6 $\mu\text{g/day}$ ^[1]
Mercury	$\mu\text{g/serving}$	$\mu\text{g/g}$	2 $\mu\text{g/day}$ ^[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	$\mu\text{g/g}$	50,000 mg/day
Heptane	$\mu\text{g/g}$	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: 23-000158/D005.R000
Report Date: 01/25/2023
ORELAP#: OR100028
Purchase Order:
Received: 01/04/23 15:05

Customer: Etz Hayim Holdings
Product identity: FORM-TN.FS.SLP50-EK49(A)
Client/Metric ID: .
Laboratory ID: 23-000158-0002

Summary

Potency:

Analyte per 1g	Result	Limits	Units	Status	
CBC per 1g	1.10		mg/1g		CBD-Total per Serving Size 30.7 mg/1g
CBD per 1g	30.7		mg/1g		
CBDV per 1g	0.535		mg/1g		THC-Total per Serving Size 1.25 mg/1g
CBE per 1g	0.471		mg/1g		(Reported in milligrams per serving)
CBG per 1g	12.0		mg/1g		
CBL per 1g	0.0932		mg/1g		
CBN per 1g	12.2		mg/1g		
CBT per 1g	0.582		mg/1g		
Δ9-THC per 1g	1.25		mg/1g		

Residual Solvents:

All analytes passing and less than LOQ.

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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Report Date: 01/25/2023
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Purchase Order:
Received: 01/04/23 15:05

Customer: Etz Hayim Holdings
 16427 NE Airport Way
 PORTLAND 97230
 United States of America (USA)
Product identity: FORM-TN.FS.SLP50-EK49(A)
Client/Metric ID: .
Sample Date:
Laboratory ID: 23-000158-0002
Evidence of Cooling: No
Temp: 18.8 °C
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: 2300233	Analyze: 1/7/23 8:52:00 AM
Analyte	Result	Limits	Units	LOQ	Notes
CBC per 1g	1.10		mg/1g	0.0326	
CBC-A per 1g	< LOQ		mg/1g	0.0326	
CBC-Total per 1g	1.10		mg/1g	0.0612	
CBD per 1g	30.7		mg/1g	0.326	
CBD-A per 1g	< LOQ		mg/1g	0.0326	
CBD-Total per 1g	30.7		mg/1g	0.354	
CBDV per 1g	0.535		mg/1g	0.0326	
CBDV-A per 1g	< LOQ		mg/1g	0.0326	
CBDV-Total per 1g	0.535		mg/1g	0.0608	
CBE per 1g	0.471		mg/1g	0.0326	
CBG per 1g	12.0		mg/1g	0.326	
CBG-A per 1g	< LOQ		mg/1g	0.0326	
CBG-Total per 1g	12.0		mg/1g	0.354	
CBL per 1g	0.0932		mg/1g	0.0326	
CBL-A per 1g	< LOQ		mg/1g	0.0326	
CBL-Total per 1g	0.0932		mg/1g	0.0612	
CBN per 1g	12.2		mg/1g	0.326	
CBT per 1g	0.582		mg/1g	0.0326	
Δ8-THCV per 1g	< LOQ		mg/1g	0.0326	
Δ10-THC-9R per 1g	< LOQ		mg/1g	0.0326	
Δ8-THC per 1g	< LOQ		mg/1g	0.0326	
Δ9-THC per 1g	1.25		mg/1g	0.0326	
exo-THC per 1g	< LOQ		mg/1g	0.0326	
THC-A per 1g	< LOQ		mg/1g	0.0326	
THC-Total per 1g	1.25		mg/1g	0.0612	
THCV per 1g	< LOQ		mg/1g	0.0326	
THCV-A per 1g	< LOQ		mg/1g	0.0326	
THCV-Total per 1g	< LOQ		mg/1g	0.0612	
Total Cannabinoids per 1g	58.9		mg/1g		



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Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2300404	01/15/23 AOAC 991.14 (Petrifilm) ^P		
Total Coliforms	< LOQ		cfu/g	10	2300404	01/15/23 AOAC 991.14 (Petrifilm) ^P		
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2300405	01/16/23 AOAC 2014.05 (RAPID) ^P		
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2300405	01/16/23 AOAC 2014.05 (RAPID) ^P		

Solvents **Method:** Residual Solvents by GC/MS^b **Units** µg/g **Batch** 2300722 **Analyze** 01/24/23 12:13 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							



Pesticides											
Method: AOAC 2007.01 & EN 15662 (mod) ^b						Units mg/kg	Batch 2300525	Analyze 01/18/23 07:53 AM			
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	< LOQ	0.200	mg/kg	0.0903	2300594	01/18/23	AOAC 2013.06 (mod.) ^b	pass		
Cadmium	< LOQ	0.200	mg/kg	0.0903	2300594	01/18/23	AOAC 2013.06 (mod.) ^b	pass		
Lead	< LOQ	0.500	mg/kg	0.0903	2300594	01/18/23	AOAC 2013.06 (mod.) ^b	pass		
Mercury	< LOQ	0.100	mg/kg	0.0452	2300594	01/18/23	AOAC 2013.06 (mod.) ^b	pass		



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Received: 01/04/23 15:05

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: 23-000158/D005.R000
Report Date: 01/25/2023
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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: 2300233

Laboratory Control Sample									
Analyte	LCS	Result	Spike	Units	% Rec	Limits		Evaluation	Notes
CBDVA	2	0.0347	0.034	%	103	80.0	- 120	Acceptable	
CBDV	2	0.0390	0.037	%	106	80.0	- 120	Acceptable	
CBE	2	0.0362	0.035	%	102	80.0	- 120	Acceptable	
CBDA	1	0.0332	0.034	%	96.3	90.0	- 110	Acceptable	
CBGA	1	0.0335	0.034	%	97.1	80.0	- 120	Acceptable	
CBG	1	0.0337	0.035	%	97.4	80.0	- 120	Acceptable	
CBD	1	0.0334	0.035	%	96.3	90.0	- 110	Acceptable	
THCV	2	0.0371	0.035	%	105	80.0	- 120	Acceptable	
d8THCV	2	0.0360	0.036	%	101	80.0	- 120	Acceptable	
THCVA	2	0.0339	0.033	%	103	80.0	- 120	Acceptable	
CBN	1	0.0348	0.036	%	97.4	80.0	- 120	Acceptable	
exo-THC	2	0.0352	0.034	%	103	80.0	- 120	Acceptable	
d9THC	1	0.0363	0.037	%	97.6	90.0	- 110	Acceptable	
d8THC	1	0.0351	0.036	%	97.5	90.0	- 110	Acceptable	
CBL	2	0.0350	0.033	%	105	80.0	- 120	Acceptable	
d10THC	1	NA	0.033	%	NA	80.0	- 120	Acceptable	Q6
CBG	2	0.0372	0.036	%	102	80.0	- 120	Acceptable	
THCA	1	0.0328	0.034	%	96.8	90.0	- 110	Acceptable	
CBGA	2	0.0356	0.034	%	104	80.0	- 120	Acceptable	
CBLA	2	0.0360	0.035	%	103	80.0	- 120	Acceptable	
CBT	2	0.0376	0.036	%	104	80.0	- 120	Acceptable	

Method Blank						
Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.003	%	< 0.003	Acceptable	
CBDV	<LOQ	0.003	%	< 0.003	Acceptable	
CBE	<LOQ	0.003	%	< 0.003	Acceptable	
CBDA	<LOQ	0.003	%	< 0.003	Acceptable	
CBGA	<LOQ	0.003	%	< 0.003	Acceptable	
CBG	<LOQ	0.003	%	< 0.003	Acceptable	
CBD	<LOQ	0.003	%	< 0.003	Acceptable	
THCV	<LOQ	0.003	%	< 0.003	Acceptable	
d8THCV	<LOQ	0.003	%	< 0.003	Acceptable	
THCVA	<LOQ	0.003	%	< 0.003	Acceptable	
CBN	<LOQ	0.003	%	< 0.003	Acceptable	
exo-THC	<LOQ	0.003	%	< 0.003	Acceptable	
d9THC	<LOQ	0.003	%	< 0.003	Acceptable	
d8THC	<LOQ	0.003	%	< 0.003	Acceptable	
CBL	<LOQ	0.003	%	< 0.003	Acceptable	
d10THC	<LOQ	0.003	%	< 0.003	Acceptable	
CBG	<LOQ	0.003	%	< 0.003	Acceptable	
THCA	<LOQ	0.003	%	< 0.003	Acceptable	
CBGA	<LOQ	0.003	%	< 0.003	Acceptable	
CBLA	<LOQ	0.003	%	< 0.003	Acceptable	
CBT	<LOQ	0.003	%	< 0.003	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: 2300233						
Sample Duplicate		Sample ID: 22-011971-0002						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBDV	0.0084	0.0083	0.003	%	0.528	< 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	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBD	1.86	1.84	0.003	%	1.18	< 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	<LOQ	<LOQ	0.003	%	NA	< 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:



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Laboratory Pesticide Quality Control Results

AOAC2007.1 & EN 15662		Units: mg/Kg			Batch ID 2300525			
Method Bank	Laboratory Control Sample							
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spk	LCS % Re	Limits	Notes
Abamectin	0.000	< 0.250		0.895	1.000	89.5	50.0	150
Acephate	0.000	< 0.200		0.691	0.800	86.4	60.0	120
Acetaminocyl	0.000	< 1.000		3.509	4.000	87.7	40.0	160
Acetamidiprid	0.000	< 0.100		0.368	0.400	91.9	60.0	120
Aldicarb	0.000	< 0.200		0.732	0.800	91.5	60.0	120
Azoxystrobin	0.000	< 0.100		0.354	0.400	88.5	60.0	120
Bifenazate	0.000	< 0.100		0.395	0.400	98.7	60.0	120
Bifenthrin	0.000	< 0.100		0.351	0.400	87.7	50.0	150
Boscalid	0.000	< 0.200		0.664	0.800	83.0	60.0	120
Carbaryl	0.000	< 0.100		0.362	0.400	90.5	60.0	120
Carbofuran	0.000	< 0.100		0.374	0.400	93.6	60.0	120
Chlorantraniliprole	0.000	< 0.100		0.363	0.400	90.8	60.0	120
Chlorfenapyr	0.000	< 0.500		1.868	2.000	93.4	60.0	120
Chlorpyrifos	0.000	< 0.100		0.372	0.400	93.0	60.0	120
Clofentazine	0.000	< 0.100		0.366	0.400	91.5	60.0	120
Cyfluthrin	0.000	< 0.500		1.826	2.000	91.3	50.0	150
Cypermethrin	0.000	< 0.500		1.821	2.000	91.0	50.0	150
Daminozide	0.000	< 0.500		1.804	2.000	90.2	60.0	120
Diazinon	0.000	< 0.100		0.387	0.400	96.8	60.0	120
Dichlorvos	0.000	< 0.500		1.900	2.000	95.0	60.0	120
Dimethoate	0.000	< 0.100		0.359	0.400	89.8	60.0	120
Ethoprophos	0.000	< 0.100		0.369	0.400	92.2	60.0	120
Etofenprox	0.000	< 0.200		0.712	0.800	89.0	50.0	150
Etoxazole	0.000	< 0.100		0.366	0.400	91.6	60.0	120
Fenoxycarb	0.000	< 0.100		0.363	0.400	90.8	60.0	120
Fenpyroximate	0.000	< 0.200		0.742	0.800	92.7	60.0	120
Fipronil	0.000	< 0.200		0.713	0.800	89.1	60.0	120
Fonicamid	0.000	< 0.250		0.839	1.000	83.9	60.0	120
Fludioxonil	0.000	< 0.200		0.777	0.800	97.2	50.0	150
Hexythiazox	0.000	< 0.250		0.885	1.000	88.5	60.0	120
Imazalil	0.000	< 0.100		0.377	0.400	94.3	60.0	120
Imidacloprid	0.000	< 0.200		0.721	0.800	90.2	60.0	120
Kresoxim-methyl	0.000	< 0.200		0.733	0.800	91.6	60.0	120
Malathion	0.000	< 0.100		0.366	0.400	91.5	60.0	120
Metaxalaxyl	0.000	< 0.100		0.373	0.400	93.3	60.0	120
Methiocarb	0.000	< 0.100		0.372	0.400	92.9	60.0	120
Methomyl	0.000	< 0.200		0.685	0.800	85.6	60.0	120
MGK-264	0.000	< 0.100		0.362	0.400	90.4	50.0	150
Myclobutanil	0.000	< 0.100		0.366	0.400	91.6	60.0	120
Naled	0.000	< 0.250		0.929	1.000	92.9	50.0	150
Oxamyl	0.000	< 0.500		1.752	2.000	87.6	60.0	120
Pacllobutrazole	0.000	< 0.200		0.728	0.800	91.0	60.0	120
Parathion-Methyl	0.000	< 0.100		0.304	0.400	76.1	50.0	150
Permethrin	0.000	< 0.100		0.364	0.400	91.0	50.0	150
Phosmet	0.000	< 0.100		0.349	0.400	87.3	50.0	150
Piperonyl butoxide	0.000	< 0.500		1.796	2.000	89.8	60.0	120
Prallethrin	0.000	< 0.100		0.364	0.400	91.1	60.0	120
Propiconazole	0.000	< 0.200		0.730	0.800	91.3	60.0	120
Propoxur	0.000	< 0.100		0.372	0.400	93.0	60.0	120
Pyrethrin (Summe)	0.000	< 0.100		0.466	0.488	95.5	60.0	120
Pyridaben	0.000	< 0.100		0.371	0.400	92.6	50.0	150
Spirosad	0.000	< 0.100		0.344	0.388	88.6	50.0	150
Spiromesifen	0.000	< 0.100		0.363	0.400	90.8	60.0	120
Spirotetramat	0.000	< 0.100		0.370	0.400	92.5	60.0	120
Spiroxamine	0.000	< 0.200		0.691	0.800	86.4	60.0	120
Tebuconazole	0.000	< 0.200		0.730	0.800	91.2	60.0	120
Thiacloprid	0.000	< 0.100		0.363	0.400	90.7	60.0	120
Thiamethoxam	0.000	< 0.100		0.381	0.400	95.3	60.0	120
Trifloxystrobin	0.000	< 0.100		0.361	0.400	90.2	60.0	120



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Laboratory Pesticide Quality Control Results

AOAC2007.1 & EN 15662		Units: mg/Kg					Batch ID 2300525			
Matrix Spke/Matrix Spke Duplicate Recoveries	Result	MS Res	MSD Res	Spike	RFD%	Limit	MS % Re	MSD % Re	Limits	Notes
Abamectin	0.000	0.836	0.901	1.000	7.4%	< 30	83.6%	90.1%	50 - 150	
Acephate	0.000	0.728	0.678	0.800	7.1%	< 30	91.0%	84.7%	50 - 150	
Acetaminocyl	0.000	3.197	3.409	4.000	6.4%	< 30	79.9%	85.2%	50 - 150	
Acetamiprid	0.000	0.373	0.367	0.400	1.6%	< 30	93.3%	91.8%	50 - 150	
Aldicarb	0.000	0.720	0.716	0.800	0.5%	< 30	90.0%	89.6%	50 - 150	
Azoxystrobin	0.000	0.347	0.347	0.400	0.1%	< 30	86.7%	86.8%	50 - 150	
Bifenazate	0.000	0.410	0.420	0.400	2.2%	< 30	102.6%	104.9%	50 - 150	
Bifenthrin	0.000	0.307	0.312	0.400	1.6%	< 30	76.7%	78.0%	50 - 150	
Boscalid	0.000	0.674	0.714	0.800	5.8%	< 30	84.2%	89.3%	50 - 150	
Carbaryl	0.000	0.352	0.349	0.400	0.9%	< 30	88.0%	87.2%	50 - 150	
Carbofuran	0.000	0.360	0.356	0.400	1.3%	< 30	90.1%	88.9%	50 - 150	
Chlorantraniliprole	0.000	0.381	0.374	0.400	1.8%	< 30	95.2%	93.4%	50 - 150	
Chlorfenapyr	0.000	1.384	1.484	2.000	6.9%	< 30	69.2%	74.2%	50 - 150	
Chlorpyrifos	0.000	0.389	0.386	0.400	0.7%	< 30	97.2%	96.5%	50 - 150	
Clofentezine	0.000	0.353	0.352	0.400	0.4%	< 30	88.3%	88.0%	50 - 150	
Cyfluthrin	0.000	1.058	1.128	2.000	6.4%	< 30	52.9%	56.4%	30 - 150	
Cypermethrin	0.000	1.113	1.056	2.000	5.2%	< 30	55.6%	52.8%	50 - 150	
Daminozide	0.000	1.780	1.797	2.000	1.0%	< 30	89.0%	89.8%	30 - 150	
Diazinon	0.000	0.335	0.335	0.400	0.2%	< 30	83.8%	83.6%	50 - 150	
Dichlorvos	0.000	1.884	1.814	2.000	3.8%	< 30	94.2%	90.7%	50 - 150	
Dimethoate	0.000	0.365	0.361	0.400	0.9%	< 30	91.2%	90.3%	50 - 150	
Ethoprophos	0.000	0.347	0.354	0.400	2.1%	< 30	86.7%	88.5%	50 - 150	
Etofenprox	0.000	0.592	0.622	0.800	4.8%	< 30	74.0%	77.7%	50 - 150	
Etoxazole	0.000	0.329	0.334	0.400	1.6%	< 30	82.2%	83.5%	50 - 150	
Fenoxycarb	0.000	0.353	0.358	0.400	1.4%	< 30	88.2%	89.4%	50 - 150	
Fenpyroximate	0.000	0.407	0.409	0.800	0.6%	< 30	50.9%	51.2%	50 - 150	
Fipronil	0.000	0.644	0.574	0.800	11.4%	< 30	80.5%	71.8%	50 - 150	
Fonicamid	0.000	0.929	0.917	1.000	1.3%	< 30	92.9%	91.7%	50 - 150	
Fludioxonil	0.000	0.865	0.878	0.800	1.5%	< 30	108.1%	109.7%	50 - 150	
Hexythiazox	0.000	0.887	0.866	1.000	2.5%	< 30	88.7%	86.6%	50 - 150	
Imazalil	0.000	0.376	0.371	0.400	1.3%	< 30	94.0%	92.8%	50 - 150	
Imidacloprid	0.000	0.720	0.713	0.800	1.0%	< 30	90.0%	89.1%	50 - 150	
Kresoxim-methyl	0.000	0.691	0.698	0.800	1.0%	< 30	86.4%	87.3%	50 - 150	
Malathion	0.036	0.321	0.314	0.400	2.7%	< 30	71.4%	69.5%	50 - 150	
Metaxalyl	0.000	0.361	0.368	0.400	2.0%	< 30	90.2%	92.0%	50 - 150	
Methiocarb	0.000	0.351	0.348	0.400	1.0%	< 30	87.9%	87.0%	50 - 150	
Methomyl	0.000	0.749	0.780	0.800	4.0%	< 30	93.7%	97.5%	50 - 150	
MGK-264	0.000	0.289	0.297	0.400	2.8%	< 30	72.1%	74.2%	50 - 150	
Myclobutanil	0.000	0.372	0.371	0.400	0.1%	< 30	93.0%	92.9%	50 - 150	
Naled	0.000	0.825	0.820	1.000	0.5%	< 30	82.5%	82.0%	50 - 150	
Oxamyl	0.000	1.851	1.715	2.000	7.7%	< 30	92.6%	85.7%	50 - 150	
Paclobotrazole	0.000	0.714	0.705	0.800	1.2%	< 30	89.2%	88.2%	50 - 150	
Parathion-Methyl	0.000	0.347	0.267	0.400	26.1%	< 30	86.8%	66.8%	30 - 150	
Permethrin	0.000	0.347	0.338	0.400	2.7%	< 30	86.7%	84.4%	50 - 150	
Phosmet	0.000	0.351	0.350	0.400	0.3%	< 30	87.7%	87.5%	50 - 150	
Piperonyl butoxide	0.000	1.595	1.669	2.000	4.5%	< 30	79.7%	83.4%	50 - 150	
Prallethrin	0.000	0.333	0.334	0.400	0.4%	< 30	83.2%	83.5%	50 - 150	
Propiconazole	0.000	0.778	0.780	0.800	0.2%	< 30	97.3%	97.5%	50 - 150	
Propoxur	0.000	0.367	0.360	0.400	1.9%	< 30	91.7%	90.0%	50 - 150	
Pyrethrin (Summe)	0.018	0.507	0.517	0.488	2.0%	< 30	100.4%	102.4%	50 - 150	
Pyridaben	0.000	0.340	0.336	0.400	1.3%	< 30	85.1%	84.0%	50 - 150	
Spinosad	0.000	0.314	0.314	0.388	0.1%	< 30	81.0%	81.0%	50 - 150	
Spiromesifen	0.000	0.378	0.374	0.400	1.2%	< 30	94.6%	93.5%	50 - 150	
Spirotetramat	0.000	0.428	0.438	0.400	2.3%	< 30	107.1%	109.6%	50 - 150	
Spiroxamine	0.000	0.691	0.705	0.800	2.0%	< 30	86.4%	88.1%	50 - 150	
Tebuconazole	0.000	0.700	0.701	0.800	0.1%	< 30	87.5%	87.6%	50 - 150	
Thiacloprid	0.000	0.363	0.353	0.400	2.8%	< 30	90.9%	88.3%	50 - 150	
Thiamethoxam	0.000	0.423	0.371	0.400	13.1%	< 30	105.7%	92.7%	50 - 150	
Trifloxystrobin	0.000	0.320	0.317	0.400	1.1%	< 30	80.1%	79.2%	50 - 150	



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Laboratory Quality Control Results

Residual Solvents				Batch ID: 2300722					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		480	572	µg/g	83.9	60 - 120	
Isobutane	ND	< 200		623	731	µg/g	85.2	60 - 120	
Butane	ND	< 200		592	731	µg/g	81.0	60 - 120	
2,2-Dimethylpropane	ND	< 200		812	936	µg/g	86.8	60 - 120	
Methanol	ND	< 200		1410	1620	µg/g	87.0	60 - 120	
Ethylene Oxide	ND	< 30		49	56.2	µg/g	87.2	60 - 120	
2-Methylbutane	ND	< 200		1330	1610	µg/g	82.6	60 - 120	
Pentane	ND	< 200		1330	1600	µg/g	83.1	60 - 120	
Ethanol	ND	< 200		1400	1610	µg/g	87.0	70 - 130	
Ethyl Ether	ND	< 200		1340	1630	µg/g	82.2	60 - 120	
2,2-Dimethylbutane	ND	< 30		138	171	µg/g	80.7	60 - 120	
Acetone	ND	< 200		1340	1630	µg/g	82.2	60 - 120	
2-Propanol	ND	< 200		1440	1620	µg/g	88.9	60 - 120	
Ethyl Formate	ND	< 500		1380	1670	µg/g	82.6	70 - 130	
Acetonitrile	ND	< 100		409	498	µg/g	82.1	60 - 120	
Methyl Acetate	ND	< 500		1460	1730	µg/g	84.4	70 - 130	
2,3-Dimethylbutane	ND	< 30		135	171	µg/g	78.9	60 - 120	
Dichloromethane	ND	< 60		406	483	µg/g	84.1	60 - 120	
2-Methylpentane	ND	< 30		146	168	µg/g	86.9	60 - 120	
MTBE	ND	< 500		1520	1650	µg/g	92.1	70 - 130	
3-Methylpentane	ND	< 30		125	167	µg/g	74.9	60 - 120	
Hexane	ND	< 30		178	182	µg/g	97.8	60 - 120	
1-Propanol	ND	< 500		1420	1620	µg/g	87.7	70 - 130	
Methylethylketone	ND	< 500		1330	1620	µg/g	82.1	70 - 130	
Ethyl acetate	ND	< 200		1360	1610	µg/g	84.5	60 - 120	
2-Butanol	ND	< 200		1430	1600	µg/g	89.4	60 - 120	
Tetrahydrofuran	ND	< 100		397	483	µg/g	82.2	60 - 120	
Cyclohexane	ND	< 200		1300	1610	µg/g	80.7	60 - 120	
2-methyl-1-propanol	ND	< 500		1360	1620	µg/g	84.0	70 - 130	
Benzene	ND	< 1		4.42	5.02	µg/g	88.0	60 - 120	
Isopropyl Acetate	ND	< 200		1450	1620	µg/g	89.5	60 - 120	
Heptane	ND	< 200		1280	1610	µg/g	79.5	60 - 120	
1-Butanol	ND	< 500		1450	1630	µg/g	89.0	70 - 130	
Propyl Acetate	ND	< 500		1310	1610	µg/g	81.4	70 - 130	
1,4-Dioxane	ND	< 100		390	491	µg/g	79.4	60 - 120	
2-Ethoxyethanol	ND	< 30		296	181	µg/g	163.5	60 - 120	Q1
Methylisobutylketone	ND	< 500		1260	1620	µg/g	77.8	70 - 130	
3-Methyl-1-butanol	ND	< 500		1380	1630	µg/g	84.7	70 - 130	
Ethylene Glycol	ND	< 200		652	484	µg/g	134.7	60 - 120	Q1
Toluene	ND	< 100		373	485	µg/g	76.9	60 - 120	
Isobutyl Acetate	ND	< 500		1320	1630	µg/g	81.0	70 - 130	
1-Pentanol	ND	< 500		1330	1620	µg/g	82.1	70 - 130	
Butyl Acetate	ND	< 500		1280	1620	µg/g	79.0	70 - 130	
Ethylbenzene	ND	< 200		712	969	µg/g	73.5	60 - 120	
m,p-Xylene	ND	< 200		720	994	µg/g	72.4	60 - 120	
o-Xylene	ND	< 200		694	967	µg/g	71.8	60 - 120	
Cumene	ND	< 30		126	171	µg/g	73.7	60 - 120	
Anisole	ND	< 500		1120	1630	µg/g	68.7	70 - 130	Q6
DMSO	ND	< 500		2220	1680	µg/g	132.1	70 - 130	Q1
1,2-dimethoxyethane	ND	< 50		147	169	µg/g	87.0	70 - 130	
Triethylamine	ND	< 500		1340	1630	µg/g	82.2	70 - 130	
N,N-dimethylformamide	ND	< 150		573	482	µg/g	118.9	70 - 130	
N,N-dimethylacetamide	ND	< 150		533	510	µg/g	104.5	70 - 130	
Pyridine	ND	< 50		194	203	µg/g	95.6	70 - 130	
Sulfolane	ND	< 50		198	172	µg/g	115.1	70 - 130	
1,2-Dichloroethane	ND	< 1		0.857	1	µg/g	85.7	70 - 130	
Chloroform	ND	< 1		0.892	1	µg/g	89.2	70 - 130	
Trichloroethylene	ND	< 1		0.93	1	µg/g	93.0	70 - 130	
1,1-Dichloroethane	ND	< 1		0.899	1	µg/g	89.9	70 - 130	



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QC - Sample Duplicate		Sample ID: 23-000158-0002						
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
 Q1 - Quality control result biased high. Only non-detect samples reported.
 Q6 - Quality control outside QC limits. Data acceptable based on remaining QC.

Units of Measure:

µg/g- Microgram per gram or ppm



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ORELAP#: OR100028
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Received: 01/04/23 15:05





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.