

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: 23-012700/D007.R000
Report Date: 11/08/2023
ORELAP#: OR100028
Purchase Order: 2690082
Received: 10/26/23 08:56

Customer: Etz Hayim Holdings
Product identity: FORM-TN.ISO.SBH300-FJ09
Client/Metric ID: .
Laboratory ID: 23-012700-0004

Summary

Potency:

Analyte per 1g	Result	Limits	Units	Status	
CBC per 1g	0.116		mg/1g		CBD-Total per Serving Size 311 mg/1g
CBD per 1g	311		mg/1g		
CBDV per 1g	1.94		mg/1g		THC-Total per Serving Size 0.0946 mg/1g
CBT per 1g	0.0423		mg/1g		
Δ9-THC per 1g	0.0946		mg/1g		(Reported in milligrams per serving)

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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Received: 10/26/23 08:56

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

Product identity: FORM-TN.ISO.SBH300-FJ09

Client/Metric ID: .

Sample Date:

Laboratory ID: 23-012700-0004

Evidence of Cooling: No

Temp: 19.6 °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: 2312369	Analyze: 10/31/23 1:42:00 PM	
Analyte	Result	Limits	Units	LOQ	Notes
CBC per 1g	0.116		mg/1g	0.0329	
CBC-A per 1g	< LOQ		mg/1g	0.0329	
CBC-Total per 1g	0.116		mg/1g	0.0618	
CBD per 1g	311		mg/1g	3.29	
CBD-A per 1g	< LOQ		mg/1g	0.0329	
CBD-Total per 1g	311		mg/1g	3.32	
CBDV per 1g	1.94		mg/1g	0.0329	
CBDV-A per 1g	< LOQ		mg/1g	0.0329	
CBDV-Total per 1g	1.94		mg/1g	0.0615	
CBE per 1g	< LOQ		mg/1g	0.0329	
CBG per 1g	< LOQ		mg/1g	0.0329	
CBG-A per 1g	< LOQ		mg/1g	0.0329	
CBG-Total per 1g	< LOQ		mg/1g	0.0615	
CBL per 1g	< LOQ		mg/1g	0.0329	
CBL-A per 1g	< LOQ		mg/1g	0.0329	
CBL-Total per 1g	< LOQ		mg/1g	0.0618	
CBN per 1g	< LOQ		mg/1g	0.0329	
CBT per 1g	0.0423		mg/1g	0.0329	
Δ8-THCV per 1g	< LOQ		mg/1g	0.0329	
Δ10-THC-9R per 1g	< LOQ		mg/1g	0.0329	
Δ10-THC-9S per 1g	< LOQ		mg/1g	0.0329	
Δ10-THC-Total per 1g	< LOQ		mg/1g	0.0659	
Δ8-THC per 1g	< LOQ		mg/1g	0.0329	
Δ9-THC per 1g	0.0946		mg/1g	0.0329	
delta-9-THCP per 1g	< LOQ		mg/1g	0.0329	
exo-THC per 1g	< LOQ		mg/1g	0.0329	
THC-A per 1g	< LOQ		mg/1g	0.0329	
THC-Total per 1g	0.0946		mg/1g	0.0618	
THCV per 1g	< LOQ		mg/1g	0.0329	
THCV-A per 1g	< LOQ		mg/1g	0.0329	



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Potency per 1g						
Analyte	Method: J AOAC 2015 V98-6 (mod) ^P	Result	Limits	Units mg/se	Batch: 2312369	Analyze: 10/31/23 1:42:00 PM
THCV-Total per 1g		< LOQ		mg/1g		0.0619
Total Cannabinoids per 1g		313		mg/1g		

Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2312416	11/04/23 AOAC 991.14 (Petrifilm) ^P		
Total Coliforms	< LOQ		cfu/g	10	2312416	11/04/23 AOAC 991.14 (Petrifilm) ^P		
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2312417	11/05/23 AOAC 2014.05 (RAPID) ^P		
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2312417	11/05/23 AOAC 2014.05 (RAPID) ^P		

Solvents

Method: Residual Solvents by GC/MS ^P											
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 2312435		Analyze 11/02/23 02:19 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 [‡]	< LOQ	0.200	mg/kg	0.0796	2312545	11/06/23	AOAC 2013.06 (mod.) ^b	pass		
Cadmium [‡]	< LOQ	0.200	mg/kg	0.0796	2312545	11/06/23	AOAC 2013.06 (mod.) ^b	pass		
Lead [‡]	< LOQ	0.500	mg/kg	0.0796	2312545	11/06/23	AOAC 2013.06 (mod.) ^b	pass		
Mercury [‡]	< LOQ	0.100	mg/kg	0.0398	2312545	11/06/23	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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Revision: 4 Document ID: 7148
 Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

JAOAC2015 V98-6 Batch ID: 2312369

Laboratory Control Sample									
Analyte	LCS	Result	Spike	Units	% Rec	Limits		Evaluation	Notes
CBDVA	2	0.0309	0.0323	%	95.5	80.0	- 120	Acceptable	
CBDV	2	0.0335	0.0337	%	99.5	80.0	- 120	Acceptable	
CBE	2	0.0339	0.0358	%	94.7	80.0	- 120	Acceptable	
CBD	1	0.0308	0.0322	%	95.6	90.0	- 110	Acceptable	
CBD A	1	0.0312	0.0329	%	94.7	80.0	- 120	Acceptable	
CBD B	1	0.0355	0.0368	%	96.5	80.0	- 120	Acceptable	
CBD C	1	0.0322	0.0313	%	103	90.0	- 110	Acceptable	
THCV	2	0.0323	0.0345	%	93.7	80.0	- 120	Acceptable	
Δ8THCV	2	0.0286	0.0283	%	101	80.0	- 120	Acceptable	
THCV A	2	0.0299	0.0312	%	96.1	80.0	- 120	Acceptable	
CBN	1	0.0313	0.0329	%	95.3	80.0	- 120	Acceptable	
exo-THC	2	0.0309	0.0315	%	98.2	80.0	- 120	Acceptable	
Δ9THC	1	0.0346	0.0365	%	94.8	90.0	- 110	Acceptable	
Δ8THC	1	0.0332	0.0340	%	97.6	90.0	- 110	Acceptable	
9SΔ10THC	1	0.0321	0.0337	%	95.1	80.0	- 120	Acceptable	
CBL	2	0.0327	0.0332	%	98.7	80.0	- 120	Acceptable	
9RΔ10THC	1	0.0321	0.0336	%	95.5	80.0	- 120	Acceptable	
CBC	2	0.0328	0.0342	%	96.0	80.0	- 120	Acceptable	
THCA	1	0.0317	0.0337	%	94.1	90.0	- 110	Acceptable	
CBCA	2	0.0314	0.0338	%	93.1	80.0	- 120	Acceptable	
CBLA	2	0.0339	0.0342	%	99.0	80.0	- 120	Acceptable	
Δ9THCP	2	0.0310	0.0334	%	93.0	80.0	- 120	Acceptable	
CBT	2	0.0367	0.0343	%	107	80.0	- 120	Acceptable	

Method Blank						
Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBDV	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBE	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBD	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBD A	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBD B	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBD C	<LOQ	0.00322	%	< 0.00322	Acceptable	
THCV	<LOQ	0.00322	%	< 0.00322	Acceptable	
Δ8THCV	<LOQ	0.00322	%	< 0.00322	Acceptable	
THCV A	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBN	<LOQ	0.00322	%	< 0.00322	Acceptable	
exo-THC	<LOQ	0.00322	%	< 0.00322	Acceptable	
Δ9THC	<LOQ	0.00322	%	< 0.00322	Acceptable	
Δ8THC	<LOQ	0.00322	%	< 0.00322	Acceptable	
9SΔ10THC	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBL	<LOQ	0.00322	%	< 0.00322	Acceptable	
9RΔ10THC	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBC	<LOQ	0.00322	%	< 0.00322	Acceptable	
THCA	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBCA	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBLA	<LOQ	0.00322	%	< 0.00322	Acceptable	
Δ9THCP	<LOQ	0.00322	%	< 0.00322	Acceptable	
CBT	<LOQ	0.00322	%	< 0.00322	Acceptable	

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

Units of Measure:
 %- Percent



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Laboratory Quality Control Results

JAOAC2015 V98-6		Batch ID: 2312369						
Sample Duplicate		Sample ID: 23-012700-0004						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBDV	0.194	0.194	0.00330	%	0.149	< 20	Acceptable	
CBE	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBD	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBD ^A	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBD ^B	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBD	31.6	31.1	0.00330	%	1.74	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
Δ ⁸ THCV	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
THCV/A	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBN	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
Δ ⁹ THC	0.00946	0.00946	0.00330	%	0.0718	< 20	Acceptable	
Δ ⁸ THC	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
9S-Δ ¹⁰ THC	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
9R-Δ ¹⁰ THC	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBC	0.0116	0.0116	0.00330	%	0.110	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
Δ ⁹ THCP	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBT	0.00429	0.00423	0.00330	%	1.35	< 20	Acceptable	

Abbreviations

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

Units of Measure:

%- Percent



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Laboratory Pesticide Quality Control Results

AOAC2007.1 &EN 15662		Units: mg/Kg			Batch ID 2312435			
Method Blank	Blank Result	Blank Limits	Notes	LCS Result	LCS Spk	LCS % Re	Limits	Notes
Abamectin	0.000	< 0.250		0.922	1.000	92.2	50.0	150
Acetate	0.069	< 0.200		0.626	0.800	78.3	60.0	120
Acetaminophen	0.033	< 1.000		2.797	4.000	69.9	40.0	160
Acetamiprid	0.005	< 0.100		0.345	0.400	86.2	60.0	120
Aldicarb	0.000	< 0.200		0.677	0.800	84.6	60.0	120
Azoxystrobin	0.008	< 0.100		0.337	0.400	84.3	60.0	120
Bifenazate	0.000	< 0.100		0.360	0.400	90.0	60.0	120
Bifenthrin	0.011	< 0.100		0.327	0.400	81.6	50.0	150
Boscalid	0.000	< 0.200		0.676	0.800	84.6	60.0	120
Carbaryl	0.003	< 0.100		0.333	0.400	83.3	60.0	120
Carbendazim	0.000	< 0.100		0.350	0.400	87.5	60.0	120
Chlorantraniliprole	0.000	< 0.100		0.351	0.400	87.7	60.0	120
Chlorfenapyr	0.000	< 0.500		1.667	2.000	84.3	60.0	120
Chlorpyrifos	0.000	< 0.100		0.342	0.400	85.6	60.0	120
Clofentezine	0.000	< 0.100		0.322	0.400	80.5	60.0	120
Cyfluthrin	0.000	< 0.500		1.663	2.000	83.1	50.0	150
Cypermethrin	0.000	< 0.500		1.654	2.000	82.7	50.0	150
Daminozide	0.000	< 0.500		0.628	2.000	31.4	60.0	120
Diazonon	0.000	< 0.100		0.323	0.400	80.9	60.0	120
Dichlorvos	0.000	< 0.500		1.720	2.000	86.0	60.0	120
Dimethoate	0.007	< 0.100		0.325	0.400	81.3	60.0	120
Ethionphos	0.001	< 0.100		0.350	0.400	87.5	60.0	120
Etofenprox	0.001	< 0.200		0.654	0.800	81.7	50.0	150
Etoxazole	0.000	< 0.100		0.339	0.400	84.8	60.0	120
Fenoxycarb	0.000	< 0.100		0.349	0.400	87.2	60.0	120
Fenprophate	0.000	< 0.200		0.664	0.800	83.0	60.0	120
Fipronil	0.000	< 0.200		0.662	0.800	82.8	60.0	120
Fonicamid	0.000	< 0.250		0.810	1.000	81.0	60.0	120
Fludioxonil	0.000	< 0.200		0.676	0.800	84.5	50.0	150
Hexythiazox	0.000	< 0.250		0.860	1.000	86.0	60.0	120
Imazalil	0.000	< 0.100		0.374	0.400	93.6	60.0	120
Imidacloprid	0.000	< 0.200		0.658	0.800	82.2	60.0	120
Kiesoxim-methyl	0.000	< 0.200		0.691	0.800	86.3	60.0	120
Malathion	0.000	< 0.100		0.360	0.400	89.9	60.0	120
Metolaxyl	0.007	< 0.100		0.342	0.400	85.5	60.0	120
Methiocarb	0.002	< 0.100		0.327	0.400	81.7	60.0	120
Methomyl	0.000	< 0.200		0.634	0.800	79.2	60.0	120
MCK-264	0.000	< 0.100		0.334	0.400	83.4	50.0	150
Mydobutanol	0.000	< 0.100		0.350	0.400	87.5	60.0	120
Naled	0.000	< 0.250		0.843	1.000	84.3	50.0	150
Oxamyl	0.000	< 0.500		1.577	2.000	78.9	60.0	120
Padobutrazole	0.000	< 0.200		0.697	0.800	87.1	60.0	120
Parathion-Methyl	0.000	< 0.100		0.392	0.400	97.9	50.0	150
Permethrin	0.000	< 0.100		0.361	0.400	90.3	50.0	150
Phosmet	0.000	< 0.100		0.332	0.400	82.9	50.0	150
Piperonyl butoxide	0.000	< 0.500		1.702	2.000	85.1	60.0	120
Prallethrin	0.000	< 0.100		0.347	0.400	86.9	60.0	120
Propiconazole	0.000	< 0.200		0.734	0.800	91.8	60.0	120
Propoxur	0.002	< 0.100		0.350	0.400	87.5	60.0	120
Pyrethrin (Summe)	0.002	< 0.100		0.413	0.488	84.5	60.0	120
Pyridaben	0.003	< 0.100		0.347	0.400	86.8	50.0	150
Spirosad	0.000	< 0.100		0.346	0.388	89.1	50.0	150
Spiromesfen	0.000	< 0.100		0.341	0.400	85.1	60.0	120
Spirotetramat	0.000	< 0.100		0.374	0.400	93.6	60.0	120
Spiroxamine	0.000	< 0.200		0.694	0.800	86.8	60.0	120
Tebuconazole	0.000	< 0.200		0.720	0.800	90.0	60.0	120
Thiadoprid	0.000	< 0.100		0.338	0.400	84.6	60.0	120
Thiamethoxam	0.000	< 0.100		0.324	0.400	80.9	60.0	120
Trifloxystrobin	0.000	< 0.100		0.336	0.400	83.9	60.0	120

Q7



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Laboratory Pesticide Quality Control Results

AOAC2007.1 & EN 15662		Units: mg/Kg					Batch ID 2312435			
Matrix Spke/Matrix Spke Duplicate Recoveries	Result	MS Res	MSD Res	Spike	RFD%	Limit	MS % Re	MSD % Re	Limits	Notes
Abamectin	0.00	0.646	0.588	1.00	9.2%	< 30	64.9%	58.8%	50 - 150	
Acephate	0.052	0.657	0.641	0.80	2.6%	< 30	75.8%	73.6%	50 - 150	
Acetaminophyl	0.00	1.737	1.492	4.00	15.2%	< 30	43.4%	37.3%	50 - 150	Q
Acetamiprid	0.00	0.364	0.348	0.40	4.5%	< 30	91.1%	87.1%	50 - 150	
Aldicarb	0.00	0.649	0.646	0.80	0.6%	< 30	81.2%	80.7%	50 - 150	
Azoxystrobin	0.009	0.310	0.299	0.40	3.7%	< 30	75.2%	72.9%	50 - 150	
Bifenazate	0.00	0.310	0.317	0.40	2.3%	< 30	77.9%	79.3%	50 - 150	
Bifenthrin	0.011	0.157	0.155	0.40	1.7%	< 30	36.8%	36.0%	50 - 150	Q
Boscalid	0.00	0.602	0.612	0.80	1.7%	< 30	75.2%	76.9%	50 - 150	
Carbaryl	0.002	0.308	0.306	0.40	0.7%	< 30	76.8%	75.9%	50 - 150	
Carbendazim	0.005	0.324	0.315	0.40	2.7%	< 30	79.8%	77.8%	50 - 150	
Chlorantraniliprole	0.00	0.308	0.276	0.40	10.8%	< 30	77.0%	69.1%	50 - 150	
Chlorfenapyr	0.00	1.470	1.449	2.00	1.4%	< 30	73.9%	72.9%	50 - 150	
Chlorpyrifos	0.00	0.251	0.259	0.40	3.1%	< 30	62.7%	64.7%	50 - 150	
Clofentezine	0.00	0.245	0.256	0.40	4.4%	< 30	61.2%	63.9%	50 - 150	
Cyfluthrin	0.00	0.907	0.904	2.00	0.4%	< 30	45.4%	45.2%	30 - 150	
Cypermethrin	0.00	1.023	0.966	2.00	5.7%	< 30	51.2%	48.3%	50 - 150	Q
Daminozide	0.00	0.623	0.573	2.00	8.4%	< 30	31.2%	28.6%	30 - 150	Q
Diazinon	0.00	0.283	0.284	0.40	0.1%	< 30	70.9%	70.9%	50 - 150	
Dichlorvos	0.032	1.567	1.542	2.00	1.6%	< 30	76.8%	75.9%	50 - 150	
Dimethoate	0.007	0.323	0.308	0.40	4.7%	< 30	79.0%	75.4%	50 - 150	
Ethioniazole	0.001	0.314	0.317	0.40	1.0%	< 30	78.1%	78.9%	50 - 150	
Etofenprox	0.001	0.375	0.363	0.80	3.4%	< 30	46.8%	45.2%	50 - 150	Q
Etoxazole	0.00	0.234	0.242	0.40	3.5%	< 30	58.9%	60.8%	50 - 150	
Fenoxycarb	0.00	0.292	0.291	0.40	0.5%	< 30	73.1%	72.7%	50 - 150	
Fenpyroximate	0.016	0.542	0.531	0.80	2.1%	< 30	65.7%	64.3%	50 - 150	
Fipronil	0.00	0.414	0.373	0.80	10.4%	< 30	51.8%	46.8%	50 - 150	Q
Fonicamid	0.00	0.870	0.803	1.00	8.0%	< 30	87.0%	80.3%	50 - 150	
Fludioxonil	0.00	0.778	0.807	0.80	3.6%	< 30	97.3%	100.8%	50 - 150	
Hexythiazox	0.00	0.324	0.314	1.00	3.3%	< 30	32.4%	31.4%	50 - 150	Q
Imazalil	0.005	0.329	0.335	0.40	1.7%	< 30	80.9%	82.3%	50 - 150	
Imidacloprid	0.021	0.682	0.635	0.80	7.4%	< 30	82.9%	76.8%	50 - 150	
Kiesoxim-methyl	0.012	0.584	0.549	0.80	6.4%	< 30	71.9%	67.1%	50 - 150	
Malathion	0.015	0.304	0.314	0.40	3.5%	< 30	72.2%	74.8%	50 - 150	
Metaxyl	0.007	0.318	0.307	0.40	3.7%	< 30	77.8%	74.8%	50 - 150	
Methiocarb	0.003	0.230	0.269	0.40	7.6%	< 30	71.8%	66.8%	50 - 150	
Methomyl	0.00	0.580	0.653	0.80	11.9%	< 30	72.9%	81.6%	50 - 150	
MCK-264	0.00	0.251	0.248	0.40	1.2%	< 30	62.8%	61.9%	50 - 150	
Mydobutani	0.00	0.329	0.306	0.40	7.1%	< 30	82.1%	76.9%	50 - 150	
Naled	0.00	0.775	0.743	1.00	4.3%	< 30	77.9%	74.3%	50 - 150	
Oxamyl	0.00	1.763	1.695	2.00	3.9%	< 30	88.2%	84.8%	50 - 150	
Padobutrazole	0.00	0.614	0.577	0.80	6.2%	< 30	76.8%	72.2%	50 - 150	
Parathion-Methyl	0.00	0.289	0.287	0.40	0.6%	< 30	72.2%	71.7%	30 - 150	
Permethrin	0.00	0.250	0.251	0.40	0.3%	< 30	62.9%	62.7%	50 - 150	
Phosmet	0.00	0.308	0.305	0.40	0.9%	< 30	77.0%	76.3%	50 - 150	
Piperonyl butoxide	0.009	1.233	1.235	2.00	0.3%	< 30	61.2%	61.4%	50 - 150	
Prallethrin	0.001	0.204	0.206	0.40	1.3%	< 30	50.8%	51.3%	50 - 150	
Propiconazole	0.00	0.539	0.539	0.80	0.0%	< 30	67.4%	67.4%	50 - 150	
Propoxur	0.003	0.328	0.307	0.40	6.5%	< 30	81.1%	76.0%	50 - 150	
Pyrethrin (Summe)	0.00	0.323	0.326	0.488	1.2%	< 30	66.0%	66.8%	50 - 150	
Pyridaben	0.00	0.179	0.174	0.40	2.6%	< 30	44.7%	43.9%	50 - 150	Q
Spirosad	0.00	0.248	0.243	0.388	1.9%	< 30	63.9%	62.7%	50 - 150	
Spiromesfen	0.003	0.228	0.220	0.40	3.5%	< 30	56.1%	54.1%	50 - 150	
Spirotetramat	0.00	0.369	0.353	0.40	4.3%	< 30	92.1%	88.3%	50 - 150	
Spiroxamine	0.00	0.657	0.639	0.80	2.8%	< 30	82.1%	79.8%	50 - 150	
Tebuconazole	0.00	0.607	0.602	0.80	0.9%	< 30	75.9%	75.3%	50 - 150	
Thiadoprid	0.00	0.337	0.320	0.40	5.1%	< 30	84.3%	80.1%	50 - 150	
Thiamethoxam	0.00	0.333	0.313	0.40	6.1%	< 30	83.2%	78.3%	50 - 150	
Trifloxystrobin	0.00	0.254	0.260	0.40	2.5%	< 30	63.4%	65.0%	50 - 150	



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Laboratory Quality Control Results

Residual Solvents				Batch ID: 2312529					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		556	584	µg/g	95.2	60 - 120	
Isobutane	ND	< 200		749	767	µg/g	97.7	60 - 120	
Butane	ND	< 200		747	782	µg/g	95.5	60 - 120	
2,2-Dimethylpropane	ND	< 200		959	939	µg/g	102.1	60 - 120	
Methanol	ND	< 200		1820	1670	µg/g	109.0	60 - 120	
Ethylene Oxide	ND	< 30		59.4	57.1	µg/g	104.0	60 - 120	
2-Methylbutane	ND	< 200		1660	1680	µg/g	98.8	60 - 120	
Pentane	ND	< 200		1650	1670	µg/g	98.8	60 - 120	
Ethanol	ND	< 200		1760	1660	µg/g	106.0	70 - 130	
Ethyl Ether	ND	< 200		1700	1670	µg/g	101.8	60 - 120	
2,2-Dimethylbutane	ND	< 30		192	189	µg/g	101.6	60 - 120	
Acetone	ND	< 200		1740	1670	µg/g	104.2	60 - 120	
2-Propanol	ND	< 200		1720	1630	µg/g	105.5	60 - 120	
Ethyl Formate	ND	< 500		1180	1600	µg/g	73.8	70 - 130	
Acetonitrile	ND	< 100		498	492	µg/g	101.2	60 - 120	
Methyl Acetate	ND	< 500		1510	1600	µg/g	94.4	70 - 130	
2,3-Dimethylbutane	ND	< 30		194	180	µg/g	107.8	60 - 120	
Dichloromethane	ND	< 60		506	488	µg/g	103.7	60 - 120	
2-Methylpentane	ND	< 30		168	182	µg/g	92.3	60 - 120	
MTBE	ND	< 500		1530	1610	µg/g	95.0	70 - 130	
3-Methylpentane	ND	< 30		188	177	µg/g	106.2	60 - 120	
Hexane	ND	< 30		180	177	µg/g	101.7	60 - 120	
1-Propanol	ND	< 500		1530	1600	µg/g	95.6	70 - 130	
Methylethylketone	ND	< 500		1480	1610	µg/g	91.9	70 - 130	
Ethyl acetate	ND	< 200		1690	1630	µg/g	103.7	60 - 120	
2-Butanol	ND	< 200		1670	1630	µg/g	102.5	60 - 120	
Tetrahydrofuran	ND	< 100		507	488	µg/g	103.9	60 - 120	
Cyclohexane	ND	< 200		1590	1610	µg/g	98.8	60 - 120	
2-methyl-1-propanol	ND	< 500		1500	1610	µg/g	93.2	70 - 130	
Benzene	ND	< 1		5.13	4.79	µg/g	107.1	60 - 120	
Isopropyl Acetate	ND	< 200		1720	1650	µg/g	104.2	60 - 120	
Heptane	ND	< 200		1580	1630	µg/g	96.9	60 - 120	
1-Butanol	ND	< 500		1550	1600	µg/g	96.9	70 - 130	
Propyl Acetate	ND	< 500		1440	1600	µg/g	90.0	70 - 130	
1,4-Dioxane	ND	< 100		563	523	µg/g	107.6	60 - 120	
2-Ethoxyethanol	ND	< 30		173	179	µg/g	96.6	60 - 120	
Methylisobutylketone	ND	< 500		1450	1600	µg/g	90.6	70 - 130	
3-Methyl-1-butanol	ND	< 500		1310	1600	µg/g	81.9	70 - 130	
Ethylene Glycol	ND	< 200		380	508	µg/g	75.1	60 - 120	
Toluene	ND	< 100		507	496	µg/g	102.2	60 - 120	
Isobutyl Acetate	ND	< 500		1430	1610	µg/g	88.8	70 - 130	
1-Pentanol	ND	< 500		1360	1600	µg/g	85.0	70 - 130	
Butyl Acetate	ND	< 500		1410	1610	µg/g	87.6	70 - 130	
Ethylbenzene	ND	< 200		924	978	µg/g	94.5	60 - 120	
m,p-Xylene	ND	< 200		925	994	µg/g	93.1	60 - 120	
o-Xylene	ND	< 200		930	982	µg/g	94.7	60 - 120	
Cumene	ND	< 30		164	171	µg/g	95.9	60 - 120	
Anisole	ND	< 500		1450	1600	µg/g	90.6	70 - 130	
DMSO	ND	< 500		1380	1620	µg/g	85.2	70 - 130	
1,2-dimethoxyethane	ND	< 50		163	186	µg/g	87.6	70 - 130	
Triethylamine	ND	< 500		1430	1600	µg/g	89.4	70 - 130	
N,N-dimethylformamide	ND	< 150		420	480	µg/g	87.5	70 - 130	
N,N-dimethylacetamide	ND	< 150		479	483	µg/g	99.2	70 - 130	
Pyridine	ND	< 50		133	168	µg/g	79.2	70 - 130	
Silfolane	ND	< 50		101	161	µg/g	62.7	70 - 130	Q6
1,2-Dichloroethane	ND	< 1		0.94	1	µg/g	94.0	70 - 130	
Chloroform	ND	< 1		0.976	1	µg/g	97.6	70 - 130	
Trichloroethylene	ND	< 1		0.968	1	µg/g	96.8	70 - 130	
1,1-Dichloroethane	ND	< 1		0.997	1	µg/g	99.7	70 - 130	



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QC- Sample Duplicate		Sample ID: 23-012700-0004						
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		
Methylethylketone	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: 23-012700/D007.R000
Report Date: 11/08/2023
ORELAP#: OR100028
Purchase Order: 2690082
Received: 10/26/23 08:56





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.