

CONSOLIDATED TEST RESULTS SUMMARY

Please see the following pages for full test results.

BULK SKU TN.ISO.SBH300		BATCH # FJ09		LOQ: Limit Of Quantitation LOD: Limit Of Detection 1 g = 10 ⁻³ kg = 10 ³ mg = 10 ⁶ µg 1 mg/kg = 1 ppm = 1000 ppb			
PRODUCT NAME Strawberry Hibiscus CBD Tincture		SERVING SIZE 1 mL (~0.949 g)					
LABORATORY: Columbia Laboratories		OREGON ACCREDITATION: OR100028					
POTENCY		PER SERVING		PER GRAM		Percent	
Cannabidiol (CBD)		295.39	mg/serving	311	mg/g	31.10	%
Total THC (d9-THC, THCA)		0.10	mg/serving	0.095	mg/g	0.01	%
Cannabigerol (CBG)		<LOQ	mg/serving	<LOQ	mg/g	<LOQ	%
Cannabinol (CBN)		<LOQ	mg/serving	<LOQ	mg/g	<LOQ	%
Cannabichromene (CBC)		0.12	mg/serving	0.116	mg/g	0.01	%
Tetrahydrocannabinolic Acid (THCA)		<LOQ	mg/serving	<LOQ	mg/g	<LOQ	%
Delta-9-THC (d9-THC)		0.10	mg/serving	0.095	mg/g	0.01	%
Delta-8-THC (d8-THC)		<LOQ	mg/serving	<LOQ	mg/g	<LOQ	%
HEAVY METALS		PER SERVING		PER GRAM		REGULATORY ACTION LEVEL	
Arsenic		<LOQ	µg/serving	<LOQ	µg/g	10	µg/day ^[1]
Cadmium		<LOQ	µg/serving	<LOQ	µg/g	4.1	µg/day ^[1]
Lead		<LOQ	µg/serving	<LOQ	µg/g	6	µg/day ^[1]
Mercury		<LOQ	µg/serving	<LOQ	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*		<LOQ	µg/g	50,000 mg/day			
Heptane		<LOQ	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		
Δ ⁹ -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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ORELAP#: OR100028
Purchase Order: 2690082
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)	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	
Δ ⁸ -THCV per 1g	< LOQ		mg/1g	0.0329	
Δ ¹⁰ -THC-9R per 1g	< LOQ		mg/1g	0.0329	
Δ ¹⁰ -THC-9S per 1g	< LOQ		mg/1g	0.0329	
Δ ¹⁰ -THC-Total per 1g	< LOQ		mg/1g	0.0659	
Δ ⁸ -THC per 1g	< LOQ		mg/1g	0.0329	
Δ ⁹ -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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Received: 10/26/23 08:56

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
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) ^b		
Total Coliforms	< LOQ		cfu/g	10	2312416	11/04/23 AOAC 991.14 (Petrifilm) ^b		
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2312417	11/05/23 AOAC 2014.05 (RAPID) ^b		
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2312417	11/05/23 AOAC 2014.05 (RAPID) ^b		

Solvents		Method: Residual Solvents by GC/MS ^a				Units µg/g	Batch 2312529	Analyze 11/06/23 03:22 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 2312435				Analyze 11/02/23 02:19 PM			
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result
Abamectin [‡]	< LOQ	0.50	0.250	pass		Acephate [‡]	< LOQ	0.40	0.200	pass		Acetamiprid [‡]	< LOQ	0.20	0.100	pass		Azoxystrobin [‡]	< LOQ
Acequinocyl [‡]	< LOQ	2.0	1.00	pass		Acetamiprid [‡]	< LOQ	0.20	0.100	pass		Azoxystrobin [‡]	< LOQ	0.20	0.100	pass		Bifenthrin [‡]	< LOQ
Aldicarb [‡]	< LOQ	0.40	0.200	pass		Azoxystrobin [‡]	< LOQ	0.20	0.100	pass		Bifenthrin [‡]	< LOQ	0.20	0.100	pass		Carbaryl [‡]	< LOQ
Bifenazate [‡]	< LOQ	0.20	0.100	pass		Bifenthrin [‡]	< LOQ	0.20	0.100	pass		Carbaryl [‡]	< LOQ	0.20	0.100	pass		Chlorantraniliprole [‡]	< LOQ
Boscalid [‡]	< LOQ	0.40	0.200	pass		Carbaryl [‡]	< LOQ	0.20	0.100	pass		Chlorantraniliprole [‡]	< LOQ	0.20	0.100	pass		Chlorpyrifos [‡]	< LOQ
Carbofuran [‡]	< LOQ	0.20	0.100	pass		Chlorantraniliprole [‡]	< LOQ	0.20	0.100	pass		Chlorpyrifos [‡]	< LOQ	0.20	0.100	pass		Cyfluthrin [‡]	< LOQ
Chlorfenapyr [‡]	< LOQ	1.0	0.500	pass		Chlorpyrifos [‡]	< LOQ	1.0	0.500	pass		Cyfluthrin [‡]	< LOQ	1.0	0.500	pass		Daminozide [‡]	< LOQ
Clofentezine [‡]	< LOQ	0.20	0.100	pass		Cyfluthrin [‡]	< LOQ	1.0	0.500	pass		Daminozide [‡]	< LOQ	1.0	0.500	pass		Dichlorvos [‡]	< LOQ
Cypermethrin [‡]	< LOQ	1.0	0.500	pass		Daminozide [‡]	< LOQ	1.0	0.500	pass		Dichlorvos [‡]	< LOQ	1.0	0.500	pass		Ethoprophos [‡]	< LOQ
Diazinon [‡]	< LOQ	0.20	0.100	pass		Dichlorvos [‡]	< LOQ	0.20	0.100	pass		Ethoprophos [‡]	< LOQ	0.20	0.100	pass		Etoxazole [‡]	< LOQ
Dimethoate [‡]	< LOQ	0.20	0.100	pass		Ethoprophos [‡]	< LOQ	0.20	0.100	pass		Etoxazole [‡]	< LOQ	0.20	0.100	pass		Fenpyroximate [‡]	< LOQ
Etofenprox [‡]	< LOQ	0.40	0.200	pass		Etoxazole [‡]	< LOQ	0.40	0.200	pass		Fenpyroximate [‡]	< LOQ	0.40	0.200	pass		Flonicamid [‡]	< LOQ
Fenoxycarb [‡]	< LOQ	0.20	0.100	pass		Fenpyroximate [‡]	< LOQ	1.0	0.400	pass		Flonicamid [‡]	< LOQ	1.0	0.400	pass		Hexythiazox [‡]	< LOQ
Fipronil [‡]	< LOQ	0.40	0.200	pass		Flonicamid [‡]	< LOQ	1.0	0.400	pass		Hexythiazox [‡]	< LOQ	0.40	0.200	pass		Imidacloprid [‡]	< LOQ
Fludioxonil [‡]	< LOQ	0.40	0.200	pass		Hexythiazox [‡]	< LOQ	0.20	0.100	pass		Imidacloprid [‡]	< LOQ	0.20	0.100	pass		Malathion [‡]	< LOQ
Imazalil [‡]	< LOQ	0.20	0.100	pass		Imidacloprid [‡]	< LOQ	0.20	0.100	pass		Malathion [‡]	< LOQ	0.20	0.100	pass		Methiocarb [‡]	< LOQ
Kresoxim-methyl [‡]	< LOQ	0.40	0.200	pass		Malathion [‡]	< LOQ	0.20	0.100	pass		Methiocarb [‡]	< LOQ	0.20	0.100	pass		MGK-264 [‡]	< LOQ
Metalaxy [‡]	< LOQ	0.20	0.100	pass		Methiocarb [‡]	< LOQ	0.50	0.250	pass		MGK-264 [‡]	< LOQ	0.50	0.250	pass		Naled [‡]	< LOQ
Methomyl [‡]	< LOQ	0.40	0.200	pass		Naled [‡]	< LOQ	0.40	0.200	pass		Naled [‡]	< LOQ	0.40	0.200	pass		Paclobutrazole [‡]	< LOQ
Myclobutanil [‡]	< LOQ	1.0	0.500	pass		Paclobutrazole [‡]	< LOQ	0.20	0.100	pass		Paclobutrazole [‡]	< LOQ	0.20	0.100	pass		Permethrin [‡]	< LOQ
Oxamy [‡]	< LOQ	0.20	0.100	pass		Permethrin [‡]	< LOQ	2.0	1.00	pass		Permethrin [‡]	< LOQ	2.0	1.00	pass		Piperonyl butoxide [‡]	< LOQ
Parathion-Methyl [‡]	< LOQ	0.20	0.100	pass		Piperonyl butoxide [‡]	< LOQ	0.40	0.200	pass		Piperonyl butoxide [‡]	< LOQ	0.40	0.200	pass		Pyrethrin I (total) [‡]	< LOQ
Phosmet [‡]	< LOQ	0.20	0.100	pass		Pyrethrin I (total) [‡]	< LOQ	0.20	0.100	pass		Pyrethrin I (total) [‡]	< LOQ	0.20	0.100	pass		Spinosad [‡]	< LOQ
Prallethrin [‡]	< LOQ	0.20	0.100	pass		Spinosad [‡]	< LOQ	0.20	0.100	pass		Spinosad [‡]	< LOQ	0.20	0.100	pass		Spirotetramat [‡]	< LOQ
Propoxur [‡]	< LOQ	0.20	0.100	pass		Spirotetramat [‡]	< LOQ	0.40	0.200	pass		Spirotetramat [‡]	< LOQ	0.40	0.200	pass		Tebuconazole [‡]	< LOQ
Pyridaben [‡]	< LOQ	0.20	0.100	pass		Tebuconazole [‡]	< LOQ	0.20	0.100	pass		Tebuconazole [‡]	< LOQ	0.20	0.100	pass		Thiamethoxam [‡]	< LOQ
Spiromesifen [‡]	< LOQ	0.40	0.200	pass		Thiamethoxam [‡]	< LOQ					Thiamethoxam [‡]	< LOQ						
Spiroxamine [‡]	< LOQ	0.20	0.100	pass															
Thiacloprid [‡]	< 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							
JAOAC 2015 V98-6 Batch ID: 2312369							
Laboratory Control Sample							
Analyte	LCS	Result	Spike	Units	% Rec	Limits	Evaluation
CBDA	2	0.0309	0.0323	%	95.5	80.0 - 120	Acceptable
CBDA	2	0.0335	0.0337	%	99.5	80.0 - 120	Acceptable
CBE	2	0.0339	0.0358	%	94.7	80.0 - 120	Acceptable
CBDA	1	0.0308	0.0322	%	95.6	90.0 - 110	Acceptable
CBGA	1	0.0312	0.0329	%	94.7	80.0 - 120	Acceptable
CBG	1	0.0355	0.0368	%	96.5	80.0 - 120	Acceptable
CBG	1	0.0322	0.0313	%	103	90.0 - 110	Acceptable
THCV	2	0.0323	0.0345	%	93.7	80.0 - 120	Acceptable
g8THCV	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
g9THC	1	0.0346	0.0365	%	94.8	90.0 - 110	Acceptable
g8THC	1	0.0332	0.0340	%	97.6	90.0 - 110	Acceptable
9S-d10THC	1	0.0321	0.0337	%	95.1	80.0 - 120	Acceptable
CBL	2	0.0327	0.0332	%	98.7	80.0 - 120	Acceptable
9R-d10THC	1	0.0321	0.0336	%	95.5	80.0 - 120	Acceptable
CBG	2	0.0326	0.0342	%	96.0	80.0 - 120	Acceptable
THCA	1	0.0317	0.0337	%	94.1	90.0 - 110	Acceptable
CBGA	2	0.0314	0.0338	%	93.1	80.0 - 120	Acceptable
CBLA	2	0.0339	0.0342	%	99.0	80.0 - 120	Acceptable
g9THCF	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	
CBDA	<LOQ	0.00322	%	< 0.00322	Acceptable		
CBDA	<LOQ	0.00322	%	< 0.00322	Acceptable		
CBE	<LOQ	0.00322	%	< 0.00322	Acceptable		
CBDA	<LOQ	0.00322	%	< 0.00322	Acceptable		
CBGA	<LOQ	0.00322	%	< 0.00322	Acceptable		
CBG	<LOQ	0.00322	%	< 0.00322	Acceptable		
CBG	<LOQ	0.00322	%	< 0.00322	Acceptable		
THCV	<LOQ	0.00322	%	< 0.00322	Acceptable		
g8THCV	<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		
g9THC	<LOQ	0.00322	%	< 0.00322	Acceptable		
g8THC	<LOQ	0.00322	%	< 0.00322	Acceptable		
9S-d10THC	<LOQ	0.00322	%	< 0.00322	Acceptable		
CBL	<LOQ	0.00322	%	< 0.00322	Acceptable		
9R-d10THC	<LOQ	0.00322	%	< 0.00322	Acceptable		
CBG	<LOQ	0.00322	%	< 0.00322	Acceptable		
THCA	<LOQ	0.00322	%	< 0.00322	Acceptable		
CBGA	<LOQ	0.00322	%	< 0.00322	Acceptable		
CBLA	<LOQ	0.00322	%	< 0.00322	Acceptable		
g9THCF	<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								
JAOAC 2015 V98-6								
Batch ID: 2312369								
Sample ID: 23-012700-0004								
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDV/A	<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/A	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBD/A	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBD	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
THC/V	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
THC/V	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
THC/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	
9THC	0.00946	0.00946	0.00330	%	0.0718	< 20	Acceptable	
9THC	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
9S-10THC	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
9R-10THC	<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	
CBGA	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
9THCP	<LOQ	<LOQ	0.00330	%	NA	< 20	Acceptable	
CB1	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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Report Number: 23-012700/D007.R000
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Laboratory Pesticide Quality Control Results

AQAC 2007.1 & EN 15662		Units: mg/kg		Batch ID: 2312435			
Method Blank		Laboratory Control Sample					
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS % Re	Limits	Notes
Abamectin	0.00	< 0.25		0.92	1.00	92.2	50.0 150
Acophate	0.00	< 0.20		0.63	0.80	78.3	60.0 120
Acequinool	0.03	< 1.00		2.79	4.00	69.9	40.0 160
Acetamiprid	0.05	< 0.10		0.34	0.40	86.2	60.0 120
Aldicarb	0.00	< 0.20		0.67	0.80	84.6	60.0 120
Azoxystrobin	0.00	< 0.10		0.33	0.40	84.3	60.0 120
Bifenazale	0.00	< 0.10		0.30	0.40	90.0	60.0 120
Bifenthrin	0.01	< 0.10		0.32	0.40	81.6	50.0 150
Boscalid	0.00	< 0.20		0.66	0.80	84.6	60.0 120
Carbaryl	0.00	< 0.10		0.33	0.40	83.3	60.0 120
Carbendazim	0.00	< 0.10		0.30	0.40	87.5	60.0 120
Chlorantraniliprole	0.00	< 0.10		0.35	0.40	87.7	60.0 120
Chlorfenapyr	0.00	< 0.50		1.68	2.00	84.3	60.0 120
Chlorpyrifos	0.00	< 0.10		0.34	0.40	85.6	60.0 120
Clofentezine	0.00	< 0.10		0.32	0.40	80.5	60.0 120
Cyfluthrin	0.00	< 0.50		1.66	2.00	83.1	50.0 150
Cypermethrin	0.00	< 0.50		1.64	2.00	82.7	50.0 150
Daminozide	0.00	< 0.50		0.63	2.00	31.4	60.0 120
Diazinon	0.00	< 0.10		0.33	0.40	80.9	60.0 120
Dichlorvos	0.00	< 0.50		1.70	2.00	86.0	60.0 120
Dimethoate	0.00	< 0.10		0.35	0.40	81.3	60.0 120
Ethionphos	0.00	< 0.10		0.30	0.40	87.5	60.0 120
Efenprox	0.00	< 0.20		0.64	0.80	81.7	50.0 150
Etoxazole	0.00	< 0.10		0.33	0.40	84.8	60.0 120
Fenoxycarb	0.00	< 0.10		0.34	0.40	87.2	60.0 120
Fenpyroximate	0.00	< 0.20		0.64	0.80	83.0	60.0 120
Fipronil	0.00	< 0.20		0.62	0.80	82.8	60.0 120
Flonicamid	0.00	< 0.25		0.80	1.00	81.0	60.0 120
Fludioxonil	0.00	< 0.20		0.66	0.80	84.5	50.0 150
Hexythiazox	0.00	< 0.25		0.80	1.00	86.0	60.0 120
Imazalil	0.00	< 0.10		0.34	0.40	93.6	60.0 120
Imidacloprid	0.00	< 0.20		0.68	0.80	82.2	60.0 120
Kiesoxim-methyl	0.00	< 0.20		0.69	0.80	86.3	60.0 120
Malathion	0.00	< 0.10		0.30	0.40	89.9	60.0 120
Metaboxy	0.00	< 0.10		0.34	0.40	85.5	60.0 120
Methiocarb	0.00	< 0.10		0.32	0.40	81.7	60.0 120
Methomyl	0.00	< 0.20		0.64	0.80	79.2	60.0 120
MKG-264	0.00	< 0.10		0.34	0.40	83.4	50.0 150
Myndobutani	0.00	< 0.10		0.30	0.40	87.5	60.0 120
Naled	0.00	< 0.25		0.84	1.00	84.3	50.0 150
Oxamyl	0.00	< 0.50		1.57	2.00	78.9	60.0 120
Pedacutrazole	0.00	< 0.20		0.69	0.80	87.1	60.0 120
Permethrin-Methyl	0.00	< 0.10		0.32	0.40	97.9	50.0 150
Permethrin	0.00	< 0.10		0.36	0.40	90.3	50.0 150
Phosmet	0.00	< 0.10		0.32	0.40	82.9	50.0 150
Piperonyl butoxide	0.00	< 0.50		1.70	2.00	85.1	60.0 120
Prallethrin	0.00	< 0.10		0.34	0.40	86.9	60.0 120
Propiconazole	0.00	< 0.20		0.74	0.80	91.8	60.0 120
Propoxur	0.00	< 0.10		0.30	0.40	87.5	60.0 120
Pyrethrin (Summe)	0.00	< 0.10		0.43	0.40	84.5	60.0 120
Pyridaben	0.00	< 0.10		0.34	0.40	86.8	50.0 150
Spinosad	0.00	< 0.10		0.34	0.38	89.1	50.0 150
Spiromesifen	0.00	< 0.10		0.34	0.40	85.1	60.0 120
Spirotetramat	0.00	< 0.10		0.34	0.40	93.6	60.0 120
Spiromame	0.00	< 0.20		0.64	0.80	86.8	60.0 120
Tebuconazole	0.00	< 0.20		0.70	0.80	90.0	60.0 120
Thiadoprid	0.00	< 0.10		0.33	0.40	84.6	60.0 120
Thiamethoxam	0.00	< 0.10		0.34	0.40	80.9	60.0 120
Trifloxystrobin	0.00	< 0.10		0.33	0.40	83.9	60.0 120

Q7



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Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662		Units: mg/Kg				Batch ID 2312435				
Matrix Spike/Matrix Spike Duplicate Recoveries		Sample ID: 23-0126354001								
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Re	MSD % Re	Limits	Notes
Abamectin	0.000	0.645	0.588	1.000	9.2%	< 30	64.9%	58.8%	50 - 150	
Acopiate	0.052	0.657	0.641	0.800	2.6%	< 30	75.6%	73.6%	50 - 150	
Acetamiprid	0.000	1.737	1.492	4.000	15.2%	< 30	43.4%	37.3%	50 - 150	Q
Acetamiprid	0.000	0.364	0.348	0.400	4.5%	< 30	91.1%	87.1%	50 - 150	
Aldicarb	0.000	0.648	0.645	0.800	0.6%	< 30	81.2%	80.7%	50 - 150	
Azoxystrobin	0.009	0.310	0.299	0.400	3.7%	< 30	75.2%	72.5%	50 - 150	
Bifenazate	0.000	0.310	0.317	0.400	2.3%	< 30	77.9%	79.3%	50 - 150	
Bifenthrin	0.011	0.157	0.155	0.400	1.7%	< 30	36.6%	36.0%	50 - 150	Q
Boscalid	0.000	0.602	0.612	0.800	1.7%	< 30	75.2%	76.5%	50 - 150	
Carbaryl	0.002	0.308	0.306	0.400	0.7%	< 30	76.9%	75.9%	50 - 150	
Carbendazim	0.005	0.324	0.315	0.400	2.7%	< 30	79.8%	77.6%	50 - 150	
Chlorantraniliprole	0.000	0.308	0.276	0.400	10.8%	< 30	77.0%	69.1%	50 - 150	
Chlorfenvinphos	0.000	1.470	1.448	2.000	1.4%	< 30	73.9%	72.5%	50 - 150	
Chlorpyrifos	0.000	0.251	0.259	0.400	3.1%	< 30	62.7%	64.7%	50 - 150	
Clofentezine	0.000	0.245	0.256	0.400	4.4%	< 30	61.2%	63.9%	50 - 150	
Cyfluthrin	0.000	0.907	0.904	2.000	0.4%	< 30	45.4%	45.2%	30 - 150	
Cypermethrin	0.000	1.023	0.986	2.000	5.7%	< 30	51.2%	48.3%	50 - 150	Q
Daminozide	0.000	0.623	0.573	2.000	8.4%	< 30	31.2%	28.6%	30 - 150	Q
Diazinon	0.000	0.283	0.284	0.400	0.1%	< 30	70.9%	70.9%	50 - 150	
Disulfoton	0.032	1.567	1.542	2.000	1.6%	< 30	76.8%	75.5%	50 - 150	
Dimethoate	0.007	0.323	0.308	0.400	4.7%	< 30	79.0%	75.4%	50 - 150	
Ethionphos	0.001	0.314	0.317	0.400	1.0%	< 30	78.1%	78.9%	50 - 150	
Etofenprox	0.001	0.375	0.363	0.800	3.4%	< 30	46.8%	45.2%	50 - 150	Q
Etoxazole	0.000	0.234	0.242	0.400	3.5%	< 30	58.9%	60.6%	50 - 150	
Fenoxycarb	0.000	0.292	0.291	0.400	0.5%	< 30	73.1%	72.7%	50 - 150	
Fenprophos	0.016	0.542	0.531	0.800	2.1%	< 30	65.7%	64.3%	50 - 150	
Fipronil	0.000	0.414	0.373	0.800	10.4%	< 30	51.8%	46.6%	50 - 150	Q
Flonicamid	0.000	0.870	0.803	1.000	8.0%	< 30	87.0%	80.3%	50 - 150	
Fludioxonil	0.000	0.778	0.807	0.800	3.6%	< 30	97.3%	100.8%	50 - 150	
Hexythiazox	0.000	0.324	0.314	1.000	3.3%	< 30	32.4%	31.4%	50 - 150	Q
Imazalil	0.005	0.329	0.335	0.400	1.7%	< 30	80.9%	82.3%	50 - 150	
Imidacloprid	0.021	0.682	0.635	0.800	7.4%	< 30	82.9%	76.6%	50 - 150	
Kiesoxim-methyl	0.012	0.584	0.548	0.800	6.4%	< 30	71.9%	67.1%	50 - 150	
Malathion	0.015	0.304	0.314	0.400	3.5%	< 30	72.2%	74.8%	50 - 150	
Metolachlor	0.007	0.318	0.307	0.400	3.7%	< 30	77.6%	74.8%	50 - 150	
Methidathion	0.003	0.290	0.289	0.400	7.6%	< 30	71.8%	68.6%	50 - 150	
Methomyl	0.000	0.580	0.653	0.800	11.9%	< 30	72.9%	81.6%	50 - 150	
MK264	0.000	0.251	0.248	0.400	1.2%	< 30	62.8%	61.9%	50 - 150	
Myndobutanol	0.000	0.329	0.306	0.400	7.1%	< 30	82.1%	76.5%	50 - 150	
Naled	0.000	0.775	0.748	1.000	4.3%	< 30	77.5%	74.3%	50 - 150	
Oxamyl	0.000	1.763	1.695	2.000	3.9%	< 30	88.2%	84.8%	50 - 150	
Pedacoutazole	0.000	0.614	0.577	0.800	6.2%	< 30	76.9%	72.2%	50 - 150	
Permethrin-Methyl	0.000	0.289	0.287	0.400	0.6%	< 30	72.2%	71.7%	30 - 150	
Permethrin	0.000	0.250	0.251	0.400	0.3%	< 30	62.9%	62.7%	50 - 150	
Phosmet	0.000	0.308	0.305	0.400	0.9%	< 30	77.0%	76.3%	50 - 150	
Piperonyl butoxide	0.009	1.233	1.236	2.000	0.3%	< 30	61.2%	61.4%	50 - 150	
Prallethrin	0.001	0.204	0.205	0.400	1.3%	< 30	50.8%	51.3%	50 - 150	
Propiconazole	0.000	0.539	0.539	0.800	0.0%	< 30	67.4%	67.4%	50 - 150	
Propoxur	0.003	0.328	0.307	0.400	6.5%	< 30	81.1%	76.0%	50 - 150	
Pyrethrin (Summe)	0.000	0.323	0.326	0.488	1.2%	< 30	66.0%	66.8%	50 - 150	
Pyridaben	0.000	0.179	0.174	0.400	2.6%	< 30	44.7%	43.5%	50 - 150	Q
Spinosad	0.000	0.248	0.243	0.388	1.9%	< 30	63.9%	62.7%	50 - 150	
Spiromesifen	0.003	0.228	0.220	0.400	3.5%	< 30	56.1%	54.1%	50 - 150	
Spirotetramat	0.000	0.389	0.353	0.400	4.3%	< 30	92.1%	88.3%	50 - 150	
Spirotetramat	0.000	0.657	0.639	0.800	2.8%	< 30	82.1%	79.8%	50 - 150	
Tebuconazole	0.000	0.607	0.602	0.800	0.9%	< 30	75.9%	75.3%	50 - 150	
Thioproprid	0.000	0.337	0.320	0.400	5.1%	< 30	84.3%	80.1%	50 - 150	
Thiamethoxam	0.000	0.333	0.313	0.400	6.1%	< 30	83.2%	79.3%	50 - 150	
Trifloxystrobin	0.000	0.254	0.260	0.400	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		558	584	µg/g	65.2 60 - 120
Isobutane	ND	< 200		749	787	µg/g	67.7 60 - 120
Butane	ND	< 200		747	782	µg/g	65.5 60 - 120
2,2-Dimethylpropane	ND	< 200		656	698	µg/g	102.1 60 - 120
Methanol	ND	< 200		1820	1870	µg/g	109.0 60 - 120
Ethylene Oxide	ND	< 30		59.4	57.1	µg/g	104.0 60 - 120
2-Methylbutane	ND	< 200		1600	1680	µg/g	88.8 60 - 120
Pentane	ND	< 200		1650	1670	µg/g	88.8 60 - 120
Ethanol	ND	< 200		1760	1680	µg/g	106.0 70 - 130
Ethyl Ether	ND	< 200		1704	1670	µg/g	101.8 60 - 120
2,2-Dimethylbutane	ND	< 30		162	168	µg/g	101.6 60 - 120
Acetone	ND	< 200		1740	1670	µg/g	104.2 60 - 120
2-Propanol	ND	< 200		1724	1630	µg/g	105.3 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		508	488	µg/g	103.7 60 - 120
2-Methylpentane	ND	< 30		168	182	µg/g	62.3 60 - 120
MTBE	ND	< 500		1530	1610	µg/g	65.0 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	61.9 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	83.2 70 - 130
Benzene	ND	< 1		5.15	4.70	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		1560	1600	µg/g	98.0 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	178	µg/g	90.0 60 - 120
Methyl isobutyl ketone	ND	< 500		1450	1600	µg/g	90.8 70 - 130
3-Methyl-1-butanol	ND	< 500		1310	1600	µg/g	61.9 70 - 130
Ethylene Glycol	ND	< 200		380	308	µg/g	75.1 60 - 120
Toluene	ND	< 100		507	498	µg/g	102.2 60 - 120
Isobutyl Acetate	ND	< 500		1430	1610	µg/g	88.8 70 - 130
1-Pentanol	ND	< 500		1380	1600	µg/g	85.0 70 - 130
Butyl Acetate	ND	< 500		1410	1610	µg/g	87.8 70 - 130
Ethylbenzene	ND	< 200		924	978	µg/g	64.5 60 - 120
m,p-Xylene	ND	< 200		925	994	µg/g	63.1 60 - 120
o-Xylene	ND	< 200		930	982	µg/g	64.7 60 - 120
Cumene	ND	< 30		164	171	µg/g	65.9 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	188	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
Sulfone	ND	< 50		107	161	µg/g	62.7 70 - 130 Q6
1,2-Dichloroethane	ND	< 1		0.94	1	µg/g	64.0 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		0.907	1	µg/g	60.7 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	
Tetraethylamine	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



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





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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.