

CONSOLIDATED TEST RESULTS SUMMARY

Please see the following pages for full test results.

BULK SKU TN.O.FS.CM50

BATCH # FE29

PRODUCT NAME Organic Chocolate Mint CBD Tincture

SERVING SIZE 1 Dropper (1mL)

LABORATORY: Columbia Laboratories

OREGON ACCREDITATION: OR100028

LOQ: Limit Of Quantitation
LOD: Limit Of Detection

1 g = 10⁻³ kg = 10³ mg = 10⁶ µg
1 mg/kg = 1 ppm = 1000 ppb

POTENCY	PER SERVING	PER GRAM	Percent
Cannabidiol (CBD)	52.45 mg/serving	56.4 mg/g	5.64 %
Total THC (d9-THC, THCA)	1.84 mg/serving	1.98 mg/g	0.20 %
Cannabigerol (CBG)	2.17 mg/serving	2.33 mg/g	0.23 %
Cannabinol (CBN)	0.04 mg/serving	0.05 mg/g	0 %
Cannabichromene (CBC)	4.32 mg/serving	4.65 mg/g	0.47 %
Tetrahydrocannabinolic Acid (THCA)	<LOQ mg/serving	<LOQ mg/g	<LOQ %
Delta-9-THC (d9-THC)	1.84 mg/serving	1.98 mg/g	0.20 %
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: 23-007387/D012.R000
Report Date: 07/05/2023
ORELAP#: OR100028
Purchase Order:
Received: 06/21/23 16:24

Customer: Etz Hayim Holdings
Product identity: FORM-TN.O.FS.CM50-FE29
Client/Metric ID: .
Laboratory ID: 23-007387-0002

Summary

Potency:

Analyte per 1g	Result	Limits	Units	Status	
CBC per 1g	4.65		mg/1g		CBD-Total per Serving Size 56.4 mg/1g
CBD per 1g	56.4		mg/1g		
CBDV per 1g	0.667		mg/1g		THC-Total per Serving Size 1.98 mg/1g
CBE per 1g	1.01		mg/1g		(Reported in milligrams per serving)
CBG per 1g	2.33		mg/1g		
CBL per 1g	0.378		mg/1g		
CBN per 1g	0.0465		mg/1g		
CBT per 1g	1.57		mg/1g		
Δ9-THC per 1g	1.98		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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Received: 06/21/23 16:24

Customer: Etz Hayim Holdings
 16427 NE Airport Way
 PORTLAND 97230
 United States of America (USA)
Product identity: FORM-TN.O.FS.CM50-FE29
Client/Metric ID: .
Sample Date:
Laboratory ID: 23-007387-0002
Evidence of Cooling: No
Temp: 22.6 °C
Relinquished by: Client
Serving Size #1: 1 g

Sample Results

Potency per 1g					
Method: J AOAC 2015 V98-6 (mod) ^p		Units mg/se		Batch: 2308537	
				Analyze: 6/24/23 1:02:00 AM	
Analyte	Result	Limits	Units	LOQ	Notes
CBC per 1g	4.65		mg/1g	0.0328	
CBC-A per 1g	< LOQ		mg/1g	0.0328	
CBC-Total per 1g	4.65		mg/1g	0.0616	
CBD per 1g	56.4		mg/1g	0.328	
CBD-A per 1g	< LOQ		mg/1g	0.0328	
CBD-Total per 1g	56.4		mg/1g	0.357	
CBDV per 1g	0.667		mg/1g	0.0328	
CBDV-A per 1g	< LOQ		mg/1g	0.0328	
CBDV-Total per 1g	0.667		mg/1g	0.0613	
CBE per 1g	1.01		mg/1g	0.0328	
CBG per 1g	2.33		mg/1g	0.0328	
CBG-A per 1g	< LOQ		mg/1g	0.0328	
CBG-Total per 1g	2.33		mg/1g	0.0613	
CBL per 1g	0.378		mg/1g	0.0328	
CBL-A per 1g	< LOQ		mg/1g	0.0328	
CBL-Total per 1g	0.378		mg/1g	0.0616	
CBN per 1g	0.0465		mg/1g	0.0328	
CBT per 1g	1.57		mg/1g	0.0328	
Δ8-THCV per 1g	< LOQ		mg/1g	0.0328	
Δ10-THC-9R per 1g	< LOQ		mg/1g	0.0328	
Δ10-THC-9S per 1g	< LOQ		mg/1g	0.0328	
Δ10-THC-Total per 1g	< LOQ		mg/1g	0.0657	
Δ8-THC per 1g	< LOQ		mg/1g	0.0328	
Δ9-THC per 1g	1.98		mg/1g	0.0328	
delta-9-THCP per 1g	< LOQ		mg/1g	0.0328	
exo-THC per 1g	< LOQ		mg/1g	0.0328	
THC-A per 1g	< LOQ		mg/1g	0.0328	
THC-Total per 1g	1.98		mg/1g	0.0617	
THCV per 1g	< LOQ		mg/1g	0.0328	
THCV-A per 1g	< LOQ		mg/1g	0.0328	



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Potency per 1g Method: J AOAC 2015 V98-6 (mod)^b Units mg/se Batch: 2308537 Analyze: 6/24/23 1:02:00 AM

Analyte	Result	Limits	Units	LOQ	Notes
THCV-Total per 1g	< LOQ		mg/1g	0.0617	
Total Cannabinoids per 1g	69.0		mg/1g		

Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2308648	07/01/23 AOAC 991.14 (Petrifilm) ^b		I
Total Coliforms	< LOQ		cfu/g	10	2308648	07/01/23 AOAC 991.14 (Petrifilm) ^b		I
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2308649	07/01/23 AOAC 2014.05 (RAPID) ^b		I
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2308649	07/01/23 AOAC 2014.05 (RAPID) ^b		I

Solvents Method: Residual Solvents by GC/MS^b Units µg/g Batch 2308701 Analyze 06/30/23 09:23 AM

Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
1,4-Dioxane	< LOQ	380	100	pass		2-Butanol	< LOQ	5000	200	pass	
2-Ethoxyethanol	< LOQ	160	30.0	pass		2-Methylbutane (Isopentane)	< LOQ		200		
2-Methylpentane	< LOQ		30.0			2-Propanol (IPA)	< LOQ	5000	200	pass	
2,2-Dimethylbutane	< LOQ		30.0			2,2-Dimethylpropane (neo-pentane)	< LOQ		200		
2,3-Dimethylbutane	< LOQ		30.0			3-Methylpentane	< LOQ		30.0		
Acetone	< LOQ	5000	200	pass		Acetonitrile	< LOQ	410	100	pass	
Benzene	< LOQ	2.00	1.00	pass		Butanes (sum)	< LOQ	5000	400	pass	
Cyclohexane	< LOQ	3880	200	pass		Ethanol	< 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) ^p						Units mg/kg		Batch 2308728		Analyze 07/03/23 06:48 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	
Bifentazate [‡]	< 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		Etoazole [‡]	< LOQ	0.20	0.100	pass	
Fenoxycarb [‡]	< LOQ	0.20	0.100	pass		Fenproximate [‡]	< LOQ	0.40	0.200	pass	
Fipronil [‡]	< LOQ	0.40	0.200	pass		Fonicamid [‡]	< 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		Paclobutrazole [‡]	< 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	0.0986	2308688	06/29/23	AOAC 2013.06 (mod.) ^p		pass		
Cadmium [‡]	< LOQ	0.200	mg/kg	0.0986	2308688	06/29/23	AOAC 2013.06 (mod.) ^p		pass		
Lead [‡]	< LOQ	0.500	mg/kg	0.0986	2308688	06/29/23	AOAC 2013.06 (mod.) ^p		pass		
Mercury [‡]	< LOQ	0.100	mg/kg	0.0493	2308688	06/29/23	AOAC 2013.06 (mod.) ^p		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

Glossary of Qualifiers

I: Insufficient sample received to meet method requirements.

Approved Signatory

Derrick Tanner
General Manager



Revision: 4 Document ID: 7148
Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

J AOAC 2015 V98-6							
Laboratory Control Sample							
Batch ID: 2308537							
Analyte	LCS	Result	Spike	Units	% Rec	Limits	Evaluation
CBDVA	2	0.0307	0.0316	%	97.3	80.0 - 120	Acceptable
CBDV	2	0.0308	0.0315	%	97.6	80.0 - 120	Acceptable
CBE	2	0.0334	0.0348	%	95.8	80.0 - 120	Acceptable
CBDA	1	0.0314	0.0333	%	94.4	90.0 - 110	Acceptable
CBGA	1	0.0311	0.0330	%	94.3	80.0 - 120	Acceptable
CBG	1	0.0356	0.0380	%	93.7	80.0 - 120	Acceptable
CBD	1	0.0348	0.0370	%	94.2	90.0 - 110	Acceptable
THCV	2	0.0231	0.0236	%	97.7	80.0 - 120	Acceptable
Δ8THCV	2	0.0269	0.0279	%	96.5	80.0 - 120	Acceptable
THCVA	2	0.0299	0.0308	%	97.0	80.0 - 120	Acceptable
CBN	1	0.0333	0.0350	%	94.9	80.0 - 120	Acceptable
exo-THC	2	0.0271	0.0283	%	95.8	80.0 - 120	Acceptable
Δ9THC	1	0.0341	0.0361	%	94.4	90.0 - 110	Acceptable
Δ8THC	1	0.0429	0.0450	%	95.4	90.0 - 110	Acceptable
9S-Δ10THC	1	0.0242	0.0255	%	94.7	80.0 - 120	Acceptable
CBL	2	0.0304	0.0311	%	97.7	80.0 - 120	Acceptable
9R-Δ10THC	1	0.0310	0.0329	%	94.3	80.0 - 120	Acceptable
CBC	2	0.0286	0.0293	%	97.7	80.0 - 120	Acceptable
THCA	1	0.0312	0.0331	%	94.2	90.0 - 110	Acceptable
CBCA	2	0.0309	0.0320	%	96.9	80.0 - 120	Acceptable
CBLA	2	0.0292	0.0302	%	96.5	80.0 - 120	Acceptable
Δ9THCP	2	0.0310	0.0326	%	95.1	80.0 - 120	Acceptable
CBT	2	0.0315	0.0326	%	96.7	80.0 - 120	Acceptable
Method Blank							
Analyte	Result	LOQ	Units	Limits	Evaluation	Notes	
CBDVA	<LOQ	0.00317	%	< 0.00317	Acceptable		
CBDV	<LOQ	0.00317	%	< 0.00317	Acceptable		
CBE	<LOQ	0.00317	%	< 0.00317	Acceptable		
CBDA	<LOQ	0.00317	%	< 0.00317	Acceptable		
CBGA	<LOQ	0.00317	%	< 0.00317	Acceptable		
CBG	<LOQ	0.00317	%	< 0.00317	Acceptable		
CBD	<LOQ	0.00317	%	< 0.00317	Acceptable		
THCV	<LOQ	0.00317	%	< 0.00317	Acceptable		
Δ8THCV	<LOQ	0.00317	%	< 0.00317	Acceptable		
THCVA	<LOQ	0.00317	%	< 0.00317	Acceptable		
CBN	<LOQ	0.00317	%	< 0.00317	Acceptable		
exo-THC	<LOQ	0.00317	%	< 0.00317	Acceptable		
Δ9THC	<LOQ	0.00317	%	< 0.00317	Acceptable		
Δ8THC	<LOQ	0.00317	%	< 0.00317	Acceptable		
9S-Δ10THC	<LOQ	0.00317	%	< 0.00317	Acceptable		
CBL	<LOQ	0.00317	%	< 0.00317	Acceptable		
9R-Δ10THC	<LOQ	0.00317	%	< 0.00317	Acceptable		
CBC	<LOQ	0.00317	%	< 0.00317	Acceptable		
THCA	<LOQ	0.00317	%	< 0.00317	Acceptable		
CBCA	<LOQ	0.00317	%	< 0.00317	Acceptable		
CBLA	<LOQ	0.00317	%	< 0.00317	Acceptable		
Δ9THCP	<LOQ	0.00317	%	< 0.00317	Acceptable		
CBT	<LOQ	0.00317	%	< 0.00317	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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Revision: 4 Document ID: 7148
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Laboratory Quality Control Results

J AOAC 2015 V98-6		Batch ID: 2308537						
Sample Duplicate		Sample ID: 23-007387-0002						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.00323	%	NA	< 20	Acceptable	
CBDV	0.0664	0.0667	0.00323	%	0.482	< 20	Acceptable	
CBE	0.101	0.101	0.00323	%	0.310	< 20	Acceptable	
CBDA	<LOQ	<LOQ	0.00323	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.00323	%	NA	< 20	Acceptable	
CBG	0.234	0.233	0.00323	%	0.654	< 20	Acceptable	
CBD	5.69	5.64	0.00323	%	0.925	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.00323	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.00323	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.00323	%	NA	< 20	Acceptable	
CBN	0.00467	0.00465	0.00323	%	0.392	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.00323	%	NA	< 20	Acceptable	
d9THC	0.198	0.198	0.00323	%	0.132	< 20	Acceptable	
d8THC	<LOQ	<LOQ	0.00323	%	NA	< 20	Acceptable	
9S-d10THC	<LOQ	<LOQ	0.00323	%	NA	< 20	Acceptable	
CBL	0.0376	0.0378	0.00323	%	0.589	< 20	Acceptable	
9R-d10THC	<LOQ	<LOQ	0.00323	%	NA	< 20	Acceptable	
CBC	0.465	0.465	0.00323	%	0.0596	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.00323	%	NA	< 20	Acceptable	
CBCA	<LOQ	<LOQ	0.00323	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.00323	%	NA	< 20	Acceptable	
d9THCP	<LOQ	<LOQ	0.00323	%	NA	< 20	Acceptable	
CBT	0.157	0.157	0.00323	%	0.301	< 20	Acceptable	

Abbreviations

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

Units of Measure:

% - Percent



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

Laboratory Quality Control Results

Residual Solvents				Batch ID: 2308701						
Method Blank				Laboratory Control Sample						
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes	
Propane	ND	< 200		509	584	µg/g	87.2	60	- 120	
Isobutane	ND	< 200		704	767	µg/g	91.8	60	- 120	
Butane	ND	< 200		701	782	µg/g	89.6	60	- 120	
2,2-Dimethylpropane	ND	< 200		756	939	µg/g	80.5	60	- 120	
Methanol	ND	< 200		1520	1640	µg/g	92.7	60	- 120	
Ethylene Oxide	ND	< 30		50.9	57.1	µg/g	89.1	60	- 120	
2-Methylbutane	ND	< 200		1400	1600	µg/g	87.5	60	- 120	
Pentane	ND	< 200		1430	1620	µg/g	88.3	60	- 120	
Ethanol	ND	< 200		1560	1610	µg/g	96.9	70	- 130	
Ethyl Ether	ND	< 200		1450	1610	µg/g	90.1	60	- 120	
2,2-Dimethylbutane	ND	< 30		152	168	µg/g	90.5	60	- 120	
Acetone	ND	< 200		1480	1620	µg/g	91.4	60	- 120	
2-Propanol	ND	< 200		1620	1600	µg/g	101.3	60	- 120	
Ethyl Formate	ND	< 500		1370	1600	µg/g	85.6	70	- 130	
Acetonitrile	ND	< 100		436	484	µg/g	90.1	60	- 120	
Methyl Acetate	ND	< 500		1470	1610	µg/g	91.3	70	- 130	
2,3-Dimethylbutane	ND	< 30		140	162	µg/g	86.4	60	- 120	
Dichloromethane	ND	< 60		444	483	µg/g	91.9	60	- 120	
2-Methylpentane	ND	< 30		162	174	µg/g	93.1	60	- 120	
MTBE	ND	< 500		1490	1610	µg/g	92.5	70	- 130	
3-Methylpentane	ND	< 30		160	168	µg/g	95.2	60	- 120	
Hexane	ND	< 30		151	168	µg/g	89.9	60	- 120	
1-Propanol	ND	< 500		1500	1600	µg/g	93.8	70	- 130	
Methylethylketone	ND	< 500		1480	1620	µg/g	91.4	70	- 130	
Ethyl acetate	ND	< 200		1520	1600	µg/g	95.0	60	- 120	
2-Butanol	ND	< 200		1650	1600	µg/g	103.1	60	- 120	
Tetrahydrofuran	ND	< 100		467	514	µg/g	90.9	60	- 120	
Cyclohexane	ND	< 200		1500	1600	µg/g	93.8	60	- 120	
2-methyl-1-propanol	ND	< 500		1470	1610	µg/g	91.3	70	- 130	
Benzene	ND	< 1		3.93	5.12	µg/g	76.8	60	- 120	
Isopropyl Acetate	ND	< 200		1530	1620	µg/g	94.4	60	- 120	
Heptane	ND	< 200		1490	1610	µg/g	92.5	60	- 120	
1-Butanol	ND	< 500		1510	1600	µg/g	94.4	70	- 130	
Propyl Acetate	ND	< 500		1450	1600	µg/g	90.6	70	- 130	
1,4-Dioxane	ND	< 100		466	493	µg/g	94.5	60	- 120	
2-Ethoxyethanol	ND	< 30		175	163	µg/g	107.4	60	- 120	
Methylisobutylketone	ND	< 500		1450	1600	µg/g	90.6	70	- 130	
3-Methyl-1-butanol	ND	< 500		1530	1610	µg/g	95.0	70	- 130	
Ethylene Glycol	ND	< 200		288	483	µg/g	59.6	60	- 120	
Toluene	ND	< 100		454	493	µg/g	92.1	60	- 120	
Isobutyl Acetate	ND	< 500		1430	1600	µg/g	89.4	70	- 130	
1-Pentanol	ND	< 500		1570	1600	µg/g	98.1	70	- 130	
Butyl Acetate	ND	< 500		1420	1600	µg/g	88.8	70	- 130	
Ethylbenzene	ND	< 200		901	959	µg/g	93.0	60	- 120	
m,p-Xylene	ND	< 200		892	968	µg/g	92.1	60	- 120	
o-Xylene	ND	< 200		908	976	µg/g	93.0	60	- 120	
Cumene	ND	< 30		148	162	µg/g	91.4	60	- 120	
Anisole	ND	< 500		1390	1610	µg/g	86.3	70	- 130	
DMSO	ND	< 500		1080	1610	µg/g	67.1	70	- 130 Q6	
1,2-dimethoxyethane	ND	< 50		149	164	µg/g	90.9	70	- 130	
Triethylamine	ND	< 500		1280	1600	µg/g	80.0	70	- 130	
N,N-dimethylformamide	ND	< 150		440	484	µg/g	90.9	70	- 130	
N,N-dimethylacetamide	ND	< 150		413	489	µg/g	84.5	70	- 130	
Pyridine	ND	< 50		123	172	µg/g	71.5	70	- 130	
Sulfolane	ND	< 50		108	163	µg/g	66.3	70	- 130 Q6	
1,2-Dichloroethane	ND	< 1		0.959	1	µg/g	95.9	70	- 130	
Chloroform	ND	< 1		1.03	1	µg/g	103.0	70	- 130	
Trichloroethylene	ND	< 1		1.21	1	µg/g	121.0	70	- 130	
1,1-Dichloroethane	ND	< 1		0.994	1	µg/g	99.4	70	- 130	



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QC - Sample Duplicate		Sample ID: 23-007062-0002						
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	
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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Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662		Units: mg/Kg			Batch ID: 2308728			
Method Blank		Laboratory Control Sample						
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes
Abamectin	0.000	< 0.250		0.912	1.000	91.2	50.0	150
Acephate	0.035	< 0.200		0.798	0.800	99.8	60.0	120
Acequinocyl	0.000	< 1.000		3.817	4.000	95.4	40.0	160
Acetamiprid	0.000	< 0.100		0.385	0.400	96.1	60.0	120
Aldicarb	0.000	< 0.200		0.793	0.800	99.1	60.0	120
Azoxystrobin	0.000	< 0.100		0.374	0.400	93.6	60.0	120
Bifenazate	0.000	< 0.100		0.385	0.400	96.3	60.0	120
Bifenthrin	0.000	< 0.100		0.370	0.400	92.5	50.0	150
Boscalid	0.000	< 0.200		0.738	0.800	92.3	60.0	120
Carbaryl	0.000	< 0.100		0.388	0.400	97.1	60.0	120
Carbofuran	0.000	< 0.100		0.386	0.400	96.5	60.0	120
Chlorantraniliprole	0.000	< 0.100		0.383	0.400	95.8	60.0	120
Chlorfenapyr	0.000	< 0.500		2.149	2.000	107.4	60.0	120
Chlorpyrifos	0.002	< 0.100		0.373	0.400	93.2	60.0	120
Clofentezine	0.000	< 0.100		0.366	0.400	91.5	60.0	120
Cyfluthrin	0.010	< 0.500		1.796	2.000	89.8	50.0	150
Cypermethrin	0.000	< 0.500		1.928	2.000	96.4	50.0	150
Daminozide	0.000	< 0.500		0.816	2.000	40.8	60.0	120
Diazinon	0.000	< 0.100		0.409	0.400	102.4	60.0	120
Dichlorvos	0.000	< 0.500		1.935	2.000	96.8	60.0	120
Dimethoate	0.000	< 0.100		0.371	0.400	92.9	60.0	120
Ethoprophos	0.000	< 0.100		0.395	0.400	98.7	60.0	120
Etofenprox	0.000	< 0.200		0.773	0.800	96.6	50.0	150
Etoxazole	0.001	< 0.100		0.389	0.400	97.3	60.0	120
Fenoxycarb	0.000	< 0.100		0.376	0.400	93.9	60.0	120
Fenpyroximate	0.000	< 0.200		0.758	0.800	94.8	60.0	120
Fipronil	0.000	< 0.200		0.778	0.800	97.3	60.0	120
Fonicamid	0.000	< 0.250		1.011	1.000	101.1	60.0	120
Fludioxonil	0.000	< 0.200		0.791	0.800	98.9	50.0	150
Hexythiazox	0.004	< 0.250		0.952	1.000	95.2	60.0	120
Imazalil	0.000	< 0.100		0.395	0.400	98.7	60.0	120
Imidacloprid	0.000	< 0.200		0.723	0.800	90.3	60.0	120
Kresoxim-methyl	0.000	< 0.200		0.816	0.800	102.0	60.0	120
Malathion	0.000	< 0.100		0.386	0.400	96.6	60.0	120
Metalaxyl	0.000	< 0.100		0.389	0.400	97.2	60.0	120
Methiocarb	0.000	< 0.100		0.391	0.400	97.6	60.0	120
Methomyl	0.000	< 0.200		0.737	0.800	92.1	60.0	120
MGK-264	0.000	< 0.100		0.392	0.400	98.0	50.0	150
Myclobutanil	0.000	< 0.100		0.383	0.400	95.7	60.0	120
Naled	0.000	< 0.250		0.932	1.000	93.2	50.0	150
Oxamyl	0.000	< 0.500		1.799	2.000	90.0	60.0	120
Pacllobutrazole	0.000	< 0.200		0.783	0.800	97.9	60.0	120
Parathion-Methyl	0.000	< 0.100		0.340	0.400	84.9	50.0	150
Permethrin	0.000	< 0.100		0.382	0.400	95.6	50.0	150
Phosmet	0.000	< 0.100		0.389	0.400	97.2	50.0	150
Piperonyl butoxide	0.000	< 0.500		1.968	2.000	98.4	60.0	120
Prallethrin	0.000	< 0.100		0.379	0.400	94.6	60.0	120
Propiconazole	0.000	< 0.200		0.761	0.800	95.1	60.0	120
Propoxur	0.000	< 0.100		0.390	0.400	97.6	60.0	120
Pyrethrin (Summe)	0.002	< 0.100		0.471	0.488	96.4	60.0	120
Pyridaben	0.000	< 0.100		0.387	0.400	96.7	50.0	150
Spinosad	0.000	< 0.100		0.375	0.388	96.6	50.0	150
Spiromesifen	0.000	< 0.100		0.382	0.400	95.5	60.0	120
Spirotetramat	0.000	< 0.100		0.377	0.400	94.3	60.0	120
Spiroxamine	0.004	< 0.200		0.760	0.800	95.0	60.0	120
Tebuconazole	0.000	< 0.200		0.781	0.800	97.6	60.0	120
Thiacloprid	0.000	< 0.100		0.384	0.400	96.0	60.0	120
Thiamethoxam	0.000	< 0.100		0.371	0.400	92.8	60.0	120
Trifloxystrobin	0.003	< 0.100		0.375	0.400	93.8	60.0	120

Q6



12423 NE Whitaker Way
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Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662		Units: mg/Kg				Batch ID: 2308728				
Matrix Spike/Matrix Spike Duplicate Recoveries		Sample ID: 23-007725-0001								
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes
Abamectin	0.000	0.803	0.758	1.000	5.7%	< 30	80.3%	75.8%	50 - 150	
Acephate	0.087	0.721	0.747	0.800	4.0%	< 30	79.2%	82.4%	50 - 150	
Acequinocyl	0.000	1.881	1.792	4.000	4.8%	< 30	47.0%	44.8%	50 - 150	Q
Acetamidrid	0.000	0.357	0.368	0.400	3.2%	< 30	89.2%	92.0%	50 - 150	
Aldicarb	0.000	0.739	0.772	0.800	4.4%	< 30	92.4%	96.5%	50 - 150	
Azoxystrobin	0.000	0.321	0.340	0.400	5.6%	< 30	80.3%	84.9%	50 - 150	
Bifenazate	0.000	0.341	0.356	0.400	4.4%	< 30	85.3%	89.1%	50 - 150	
Bifenthrin	0.000	0.204	0.204	0.400	0.2%	< 30	50.9%	51.0%	50 - 150	
Boscalid	0.062	0.733	0.716	0.800	2.7%	< 30	83.9%	81.6%	50 - 150	
Carbaryl	0.000	0.318	0.308	0.400	3.2%	< 30	79.4%	76.9%	50 - 150	
Carbofuran	0.000	0.347	0.339	0.400	2.3%	< 30	86.7%	84.7%	50 - 150	
Chlorantraniliprole	0.000	0.373	0.385	0.400	3.1%	< 30	93.3%	96.2%	50 - 150	
Chlorfenvinpyr	0.000	1.420	1.421	2.000	0.0%	< 30	71.0%	71.0%	50 - 150	
Chlorpyrifos	0.003	0.334	0.329	0.400	1.6%	< 30	82.9%	81.6%	50 - 150	
Clofentezine	0.003	0.303	0.308	0.400	1.8%	< 30	74.8%	76.2%	50 - 150	
Cyfluthrin	0.000	0.752	0.749	2.000	0.4%	< 30	37.6%	37.4%	30 - 150	
Cypermethrin	0.000	1.636	1.587	2.000	3.0%	< 30	81.8%	79.4%	50 - 150	
Daminozide	0.004	0.807	0.869	2.000	7.4%	< 30	40.2%	43.2%	30 - 150	
Diazinon	0.002	0.346	0.378	0.400	8.8%	< 30	86.0%	93.9%	50 - 150	
Dichlorvos	0.000	1.723	1.743	2.000	1.1%	< 30	86.2%	87.1%	50 - 150	
Dimethoate	0.000	0.396	0.410	0.400	3.4%	< 30	99.1%	102.4%	50 - 150	
Ethoprophos	0.000	0.353	0.360	0.400	1.7%	< 30	88.3%	89.9%	50 - 150	
Etofenprox	0.000	0.550	0.499	0.800	9.8%	< 30	68.8%	62.3%	50 - 150	
Etoxazole	0.000	0.318	0.312	0.400	2.1%	< 30	79.6%	77.9%	50 - 150	
Fenoxycarb	0.000	0.329	0.349	0.400	6.0%	< 30	82.2%	87.3%	50 - 150	
Fenpyroximate	0.000	0.445	0.441	0.800	0.8%	< 30	55.6%	55.2%	50 - 150	
Fipronil	0.000	0.604	0.606	0.800	0.3%	< 30	75.5%	75.7%	50 - 150	
Flonicamid	0.000	1.061	1.135	1.000	6.8%	< 30	106.1%	113.5%	50 - 150	
Fludioxonil	0.000	0.795	0.797	0.800	0.2%	< 30	99.4%	99.6%	50 - 150	
Hexythiazox	0.004	0.596	0.629	1.000	5.5%	< 30	59.2%	62.5%	50 - 150	
Imazalil	0.003	0.348	0.368	0.400	5.5%	< 30	86.4%	91.2%	50 - 150	
Imidacloprid	0.000	0.838	0.916	0.800	8.8%	< 30	104.8%	114.5%	50 - 150	
Kresoxim-methyl	0.000	0.669	0.674	0.800	0.7%	< 30	83.6%	84.2%	50 - 150	
Malathion	0.000	0.346	0.350	0.400	1.0%	< 30	86.6%	87.5%	50 - 150	
Metaxalyl	0.000	0.351	0.372	0.400	6.0%	< 30	87.7%	93.1%	50 - 150	
Methiocarb	0.000	0.328	0.351	0.400	7.0%	< 30	81.9%	87.8%	50 - 150	
Methomyl	0.000	0.733	0.794	0.800	7.9%	< 30	91.7%	99.2%	50 - 150	
MGK-264	0.000	0.326	0.350	0.400	7.1%	< 30	81.5%	87.5%	50 - 150	
Myclobutanil	0.000	0.344	0.346	0.400	0.5%	< 30	86.0%	86.5%	50 - 150	
Naled	0.000	0.806	0.779	1.000	3.4%	< 30	80.6%	77.9%	50 - 150	
Oxamyl	0.000	1.967	2.077	2.000	5.4%	< 30	98.3%	103.8%	50 - 150	
Paclobutrazole	0.000	0.660	0.699	0.800	5.7%	< 30	82.5%	87.3%	50 - 150	
Parathion-Methyl	0.000	0.274	0.288	0.400	5.3%	< 30	68.4%	72.1%	30 - 150	
Permethrin	0.000	0.229	0.230	0.400	0.4%	< 30	57.3%	57.6%	50 - 150	
Phosmet	0.000	0.316	0.348	0.400	9.7%	< 30	79.1%	87.1%	50 - 150	
Piperonyl butoxide	0.000	1.546	1.630	2.000	5.2%	< 30	77.3%	81.5%	50 - 150	
Prallethrin	0.000	0.411	0.431	0.400	4.7%	< 30	102.7%	107.6%	50 - 150	
Propiconazole	0.001	0.703	0.714	0.800	1.5%	< 30	87.7%	89.1%	50 - 150	
Propoxur	0.000	0.358	0.341	0.400	4.9%	< 30	89.4%	85.1%	50 - 150	
Pyrethrin (Summe)	0.000	0.310	0.321	0.488	3.5%	< 30	63.4%	65.7%	50 - 150	
Pyridaben	0.000	0.342	0.326	0.400	4.7%	< 30	85.5%	81.6%	50 - 150	
Spinosad	0.000	0.307	0.313	0.388	1.7%	< 30	79.2%	80.6%	50 - 150	
Spiromesifen	0.000	0.337	0.339	0.400	0.5%	< 30	84.3%	84.7%	50 - 150	
Spirotetramat	0.000	0.508	0.533	0.400	4.8%	< 30	127.0%	133.3%	50 - 150	
Spiroxamine	0.004	0.682	0.756	0.800	10.3%	< 30	84.8%	94.0%	50 - 150	
Tebuconazole	0.000	0.658	0.711	0.800	7.6%	< 30	82.3%	88.8%	50 - 150	
Thiacloprid	0.000	0.354	0.379	0.400	6.8%	< 30	88.5%	94.7%	50 - 150	
Thiamethoxam	0.000	0.399	0.427	0.400	6.7%	< 30	99.8%	106.8%	50 - 150	
Trifloxystrobin	0.000	0.254	0.272	0.400	6.8%	< 30	63.5%	68.0%	50 - 150	



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Report Number: 23-007387/D012.R000
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ORELAP#: OR100028
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Received: 06/21/23 16:24





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.