

CONSOLIDATED TEST RESULTS SUMMARY

Please see the following pages for full test results.

BULK SKU PET.DIG10	BATCH # EK45	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 Digestive Health CBD Dog Chews - Pumpkin	SERVING SIZE 1 treat (~5.6g)		
LABORATORY: Columbia Laboratories	OREGON ACCREDITATION: OR100028		
POTENCY	PER SERVING	PER GRAM	Percent
Cannabidiol (CBD)	11.20 mg/serving	2.00 mg/g	0.20 %
Total THC (d9-THC, THCA)	0.40 mg/serving	0.07 mg/g	0.01 %
Cannabigerol (CBG)	0.20 mg/serving	0.04 mg/g	0.00 %
Cannabinol (CBN)	<LOQ mg/serving	<LOQ mg/g	<LOQ %
Cannabichromene (CBC)	0.33 mg/serving	0.06 mg/g	0.01 %
Tetrahydrocannabinolic Acid (THCA)	<LOQ mg/serving	<LOQ mg/g	<LOQ %
Delta-9-THC (d9-THC)	0.40 mg/serving	0.07 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	2.23 µg/serving	0.40 µg/g	10 µg/day ^[1]
Cadmium	<LOQ µg/serving	<LOQ µg/g	4.1 µg/day ^[1]
Lead	0.21 µg/serving	0.04 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.



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



Report Number: 22-014843/D006.R000
Report Date: 12/12/2022
ORELAP#: OR100028
Purchase Order:
Received: 12/05/22 15:05

Customer: Etz Hayim Holdings
Product identity: FORM-PET.DIG10-EK45
Client/Metric ID: .
Laboratory ID: 22-014843-0001

Summary

Potency:

Analyte per 1g	Result	Limits	Units	Status	
CBC per 1g	0.0581		mg/1g		CBD-Total per Serving Size 2.00 mg/1g
CBD per 1g	1.98		mg/1g		
CBDV per 1g	0.0320		mg/1g		THC-Total per Serving Size 0.0721 mg/1g
CBG per 1g	0.0363		mg/1g		
Δ9-THC per 1g	0.0721		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:

Analyte	Result	Units	Limit	Status	Analyte	Result	Units	Limit	Status
Arsenic	0.398	mg/kg	1.500	pass	Lead	0.0374	mg/kg	0.500	pass

Microbiology:

Less than LOQ for all analytes.



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Purchase Order:
Received: 12/05/22 15:05

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

Product identity: FORM-PET.DIG10-EK45

Client/Metric ID: .

Sample Date:

Laboratory ID: 22-014843-0001

Evidence of Cooling: No

Temp: 18.4 °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: 2210437	Analyze: 12/7/22 9:08:00 PM	
Analyte	Result	Limits	Units	LOQ	Notes
CBC per 1g	0.0581		mg/1g	0.0318	
CBC-A per 1g	< LOQ		mg/1g	0.0318	
CBC-Total per 1g	< LOQ		mg/1g	0.0598	
CBD per 1g	1.98		mg/1g	0.0318	
CBD-A per 1g	< LOQ		mg/1g	0.0318	
CBD-Total per 1g	2.00		mg/1g	0.0598	
CBDV per 1g	0.0320		mg/1g	0.0318	
CBDV-A per 1g	< LOQ		mg/1g	0.0318	
CBDV-Total per 1g	< LOQ		mg/1g	0.0594	
CBE per 1g	< LOQ		mg/1g	0.0318	
CBG per 1g	0.0363		mg/1g	0.0318	
CBG-A per 1g	< LOQ		mg/1g	0.0318	
CBG-Total per 1g	< LOQ		mg/1g	0.0594	
CBL per 1g	< LOQ		mg/1g	0.0318	
CBL-A per 1g	< LOQ		mg/1g	0.0318	
CBL-Total per 1g	< LOQ		mg/1g	0.0598	
CBN per 1g	< LOQ		mg/1g	0.0318	
CBT per 1g	< LOQ		mg/1g	0.0318	
Δ8-THCV per 1g	< LOQ		mg/1g	0.0318	
Δ10-THC per 1g	< LOQ		mg/1g	0.0318	
Δ8-THC per 1g	< LOQ		mg/1g	0.0318	
Δ9-THC per 1g	0.0721		mg/1g	0.0318	
exo-THC per 1g	< LOQ		mg/1g	0.0318	
THC-A per 1g	< LOQ		mg/1g	0.0318	
THC-Total per 1g	0.0721		mg/1g	0.0598	
THCV per 1g	< LOQ		mg/1g	0.0318	
THCV-A per 1g	< LOQ		mg/1g	0.0318	
THCV-Total per 1g	< LOQ		mg/1g	0.0598	
Total Cannabinoids per 1g	2.25		mg/1g		



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Received: 12/05/22 15:05

Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2210345	12/08/22 AOAC 991.14 (Petrifilm) ^P		
Total Coliforms	< LOQ		cfu/g	10	2210345	12/08/22 AOAC 991.14 (Petrifilm) ^P		
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2210347	12/09/22 AOAC 2014.05 (RAPID) ^P		
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2210347	12/09/22 AOAC 2014.05 (RAPID) ^P		

Solvents **Method:** Residual Solvents by GC/MS^b **Units** µg/g **Batch** 2210461 **Analyze** 12/09/22 10:40 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-Dimethyl butane	< LOQ		30.0			2,2-Dimethylpropane (neo-pentane)	< LOQ		200		
2,3-Dimethyl butane	< 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		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	
Isopropyl benzene (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 2210400		Analyze 12/07/22 01:29 PM	
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
Abamectin [‡]	< LOQ	0.50	0.250	pass		Acephate [‡]	< LOQ	0.40	0.200	pass	
Acequinocyl [‡]	< LOQ	2.0	1.00	pass		Acetamiprid [‡]	< LOQ	0.20	0.100	pass	
Aldicarb [‡]	< LOQ	0.40	0.200	pass		Azoxystrobin [‡]	< LOQ	0.20	0.100	pass	
Bifenazate [‡]	< LOQ	0.20	0.100	pass		Bifenthrin [‡]	< LOQ	0.20	0.100	pass	
Boscalid [‡]	< LOQ	0.40	0.200	pass		Carbaryl [‡]	< LOQ	0.20	0.100	pass	
Carbofuran [‡]	< LOQ	0.20	0.100	pass		Chlorantraniliprole [‡]	< LOQ	0.20	0.100	pass	
Chlorfenapyr [‡]	< LOQ	1.0	0.500	pass		Chlorpyrifos [‡]	< LOQ	0.20	0.100	pass	
Clofentezine [‡]	< LOQ	0.20	0.100	pass		Cyfluthrin [‡]	< LOQ	1.0	0.500	pass	
Cypermethrin [‡]	< LOQ	1.0	0.500	pass		Daminozide [‡]	< LOQ	1.0	0.500	pass	
Diazinon [‡]	< LOQ	0.20	0.100	pass		Dichlorvos [‡]	< LOQ	1.0	0.500	pass	
Dimethoate [‡]	< LOQ	0.20	0.100	pass		Ethoprophos [‡]	< LOQ	0.20	0.100	pass	
Etofenprox [‡]	< LOQ	0.40	0.200	pass		Etoxazole [‡]	< LOQ	0.20	0.100	pass	
Fenoxycarb [‡]	< LOQ	0.20	0.100	pass		Fenpyroximate [‡]	< LOQ	0.40	0.200	pass	
Fipronil [‡]	< LOQ	0.40	0.200	pass		Flonicamid [‡]	< LOQ	1.0	0.400	pass	
Fludioxonil [‡]	< LOQ	0.40	0.200	pass		Hexythiazox [‡]	< LOQ	1.0	0.400	pass	
Imazalil [‡]	< LOQ	0.20	0.100	pass		Imidacloprid [‡]	< LOQ	0.40	0.200	pass	
Kresoxim-methyl [‡]	< LOQ	0.40	0.200	pass		Malathion [‡]	< LOQ	0.20	0.100	pass	
Metalaxyl [‡]	< LOQ	0.20	0.100	pass		Methiocarb [‡]	< LOQ	0.20	0.100	pass	
Methomyl [‡]	< LOQ	0.40	0.200	pass		MGK-264 [‡]	< LOQ	0.20	0.100	pass	
Myclobutanil [‡]	< LOQ	0.20	0.100	pass		Naled [‡]	< LOQ	0.50	0.250	pass	
Oxamyl [‡]	< LOQ	1.0	0.500	pass		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	0.398	1.500	mg/kg	0.0146	2210479	12/09/22	AOAC 2013.06 (mod.) ^b	pass		
Cadmium	< LOQ	0.500	mg/kg	0.0146	2210479	12/09/22	AOAC 2013.06 (mod.) ^b	pass		
Lead	0.0374	0.500	mg/kg	0.0146	2210479	12/09/22	AOAC 2013.06 (mod.) ^b	pass		
Mercury	< LOQ	3.000	mg/kg	0.00731	2210479	12/09/22	AOAC 2013.06 (mod.) ^b	pass		



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These test results are representative of the individual sample selected and submitted by the client.

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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12423 NE Whitaker Way Portland OR 97230 p.503-254-1794

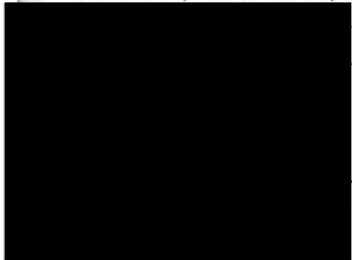
Cannabis Chain of Custody Record

ETZHAYIM 22-014843



Analysis Requested

Etz Hayim Holdings



Project Name:

- Report Instructions:
- Send to State - METRC
- Email Final Results:
- Fax Final Results
- Cash/Check/CC/Net 30

Other:

Field ID	Date/Time Collected	Pesticides - OR 59 compounds	Pesticide Multi-Residue - 379 compounds	Potency	Residual Solvents	Water Activity	Moisture	Terpenes	Micro: Yeast and Mold	Micro: E.Coli and Total Coliform	Heavy Metals	Mycotoxins	Other	Matrix	Weight	Serving size for edibles	Comments/Metric ID
FORM-PET.DIG10-EK45		X		X	X				X	X	X					mg/g	Parallel Path
FORM-GMY.RLX25-220217-D		X		X	X				X	X	X			Edible		mg/g	Parallel Path
FORM-GMY.RLX25-220311-A		X		X	X				X	X	X			Edible		mg/g	Parallel Path
CYCL-GMY.D9.HB5-220811-A		X		X	X				X	X	X			Edible		mg/g	Parallel Path
CYCL-GMY.D9.HB5-220711-B		X		X	X				X	X	X			Edible		mg/g	Parallel Path
																	LazNat Discount

Collected By:	Relinquished By:	Date:	Time:	Received by:	Date:	Time:	Lab Use Only:
<input checked="" type="checkbox"/> Standard (5 day)							Client Alias:
<input type="checkbox"/> Rush (3-4 day) (1.5x Standard)							Order Number:
<input type="checkbox"/> Priority Rush (2 day) (2x Standard)							Proper Container
							Sample Condition
							Temperature: 18.9
							Shipped Via: Courier
							Evidence of cooling: <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No

SUBMISSION OF SAMPLES WITH TESTING REQUIREMENTS TO PIXIS WILL BE UNDERSTOOD TO BE AN AGREEMENT FOR SERVICES IN ACCORDANCE WITH THE CONDITIONS LISTED ON THE BACK OF THIS FORM

Revision: 1.02 Control#: CF023
Effective 01/31/2019 Revised 01/31/2019

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Page 1 of 2



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Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662		Units: mg/Kg		Batch ID: 2210400				
Method Blank		Laboratory Control Sample						
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes
Abamectin	0.000	< 0.250		1.131	1.000	113.1	50.0	150
Acephate	0.000	< 0.200		0.834	0.800	104.2	60.0	120
Acetamiprid	0.000	< 1.000		4.224	4.000	105.6	40.0	160
Acetamiprid	0.000	< 0.100		0.419	0.400	104.8	60.0	120
Aldicarb	0.000	< 0.200		0.885	0.800	110.6	60.0	120
Azoxystrobin	0.000	< 0.100		0.425	0.400	106.4	60.0	120
Bifenazate	0.000	< 0.100		0.431	0.400	107.9	60.0	120
Bifenthrin	0.000	< 0.100		0.419	0.400	104.8	50.0	150
Boscalid	0.000	< 0.200		0.794	0.800	99.3	60.0	120
Carbaryl	0.000	< 0.100		0.419	0.400	104.7	60.0	120
Carbofuran	0.000	< 0.100		0.404	0.400	100.9	60.0	120
Chlorantraniliprole	0.000	< 0.100		0.418	0.400	104.4	60.0	120
Chlorfenapyr	0.000	< 0.500		2.013	2.000	100.7	60.0	120
Chlorpyrifos	0.000	< 0.100		0.395	0.400	98.8	60.0	120
Clofentazine	0.000	< 0.100		0.354	0.400	88.5	60.0	120
Cyfluthrin	0.000	< 0.500		2.127	2.000	106.3	50.0	150
Cypermethrin	0.000	< 0.500		2.070	2.000	103.5	50.0	150
Daminozide	0.000	< 0.500		0.698	2.000	34.9	60.0	120
Diazinon	0.000	< 0.100		0.463	0.400	115.7	60.0	120
Dichlorvos	0.000	< 0.500		1.965	2.000	98.2	60.0	120
Dimethoate	0.000	< 0.100		0.414	0.400	103.6	60.0	120
Ethoprophos	0.000	< 0.100		0.399	0.400	99.7	60.0	120
Etofenprox	0.000	< 0.200		0.828	0.800	103.5	50.0	150
Etoxazole	0.000	< 0.100		0.478	0.400	119.6	60.0	120
Fenoxycarb	0.000	< 0.100		0.428	0.400	107.0	60.0	120
Fenpyroximate	0.000	< 0.200		0.827	0.800	103.3	60.0	120
Fipronil	0.000	< 0.200		0.882	0.800	110.3	60.0	120
Fonicamid	0.000	< 0.250		1.055	1.000	105.5	60.0	120
Fludioxonil	0.000	< 0.200		0.844	0.800	105.5	50.0	150
Hexythiazox	0.000	< 0.250		1.023	1.000	102.3	60.0	120
Imazalil	0.000	< 0.100		0.418	0.400	104.6	60.0	120
Imidacloprid	0.000	< 0.200		0.833	0.800	104.1	60.0	120
Kresoxim methyl	0.000	< 0.200		0.896	0.800	112.0	60.0	120
Malathion	0.000	< 0.100		0.420	0.400	105.0	60.0	120
Metaxalyl	0.000	< 0.100		0.418	0.400	104.5	60.0	120
Methiocarb	0.000	< 0.100		0.422	0.400	105.6	60.0	120
Methomyl	0.000	< 0.200		0.924	0.800	115.5	60.0	120
MGK 264	0.000	< 0.100		0.422	0.400	105.5	50.0	150
Myclobutanil	0.000	< 0.100		0.414	0.400	103.4	60.0	120
Naled	0.000	< 0.250		1.042	1.000	104.2	50.0	150
Oxamyl	0.000	< 0.500		2.277	2.000	113.9	60.0	120
Paclobotrazole	0.000	< 0.200		0.858	0.800	107.2	60.0	120
Parathion Methyl	0.000	< 0.100		0.398	0.400	99.5	50.0	150
Permethrin	0.000	< 0.100		0.405	0.400	101.3	50.0	150
Phosmet	0.000	< 0.100		0.422	0.400	105.6	50.0	150
Piperonyl butoxide	0.000	< 0.500		2.128	2.000	106.4	60.0	120
Prallethrin	0.000	< 0.100		0.415	0.400	103.8	60.0	120
Propiconazole	0.000	< 0.200		0.864	0.800	108.0	60.0	120
Propoxur	0.000	< 0.100		0.415	0.400	103.7	60.0	120
Pyrethrin (Summe)	0.000	< 0.100		0.529	0.488	108.4	60.0	120
Pyridaben	0.000	< 0.100		0.416	0.400	104.1	50.0	150
Spinosad	0.000	< 0.100		0.414	0.388	106.6	50.0	150
Spiromesifen	0.000	< 0.100		0.413	0.400	103.2	60.0	120
Spirotetramat	0.000	< 0.100		0.394	0.400	98.6	60.0	120
Spiroxamine	0.000	< 0.200		0.840	0.800	105.0	60.0	120
ebuconazole	0.000	< 0.200		0.858	0.800	107.3	60.0	120
hiacloprid	0.000	< 0.100		0.425	0.400	106.1	60.0	120
hiamethoxam	0.000	< 0.100		0.429	0.400	107.1	60.0	120
rifloxystrobin	0.000	< 0.100		0.413	0.400	103.3	60.0	120

Q6

Test results relate only to the parameters tested and to the samples as received by the laboratory. Test results meet all requirements of NELAP and the Columbia Laboratories quality assurance plan unless otherwise noted. This report shall not be reproduced, except in full, without the written consent of this laboratory. Samples will be retained for a maximum of 30 days from the receipt date unless prior arrangements have been made.

Testing in accordance with: OAR 333-007-0390 OAR 333-007-0400 OAR 333-007-0410 OAR 333-007-0430



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



Report Number: 22-014843/D006.R000
 Report Date: 12/12/2022
 ORELAP#: OR100028
 Purchase Order:
 Received: 12/05/22 15:05

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Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662		Units: mg/Kg					Batch ID: 2210400				
Matrix Spike/Matrix Spike Duplicate Recoveries						Sample ID: 22-014843-0001					
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes	
Abamectin	0.000	1.219	1.288	1.000	5.6%	< 30	121.9%	128.8%	50 150		
Acephate	0.000	0.710	0.721	0.800	1.4%	< 30	88.8%	90.1%	50 150		
Acetaminopyr	0.000	3.929	3.853	4.000	1.9%	< 30	98.2%	96.3%	50 150		
Acetamiprid	0.000	0.422	0.432	0.400	2.4%	< 30	105.5%	108.0%	50 150		
Aldicarb	0.000	0.864	0.886	0.800	2.4%	< 30	108.0%	110.7%	50 150		
Azoxystrobin	0.000	0.410	0.422	0.400	2.9%	< 30	102.5%	105.5%	50 150		
Bifenazate	0.000	0.437	0.443	0.400	1.4%	< 30	109.2%	110.7%	50 150		
Bifenthrin	0.000	0.387	0.392	0.400	1.2%	< 30	96.8%	97.9%	50 150		
Boscalid	0.000	0.773	0.824	0.800	6.4%	< 30	96.6%	103.0%	50 150		
Carbaryl	0.000	0.406	0.413	0.400	1.7%	< 30	101.6%	103.3%	50 150		
Carbofuran	0.000	0.400	0.402	0.400	0.3%	< 30	100.1%	100.4%	50 150		
Chlorantraniliprole	0.000	0.410	0.421	0.400	2.7%	< 30	102.6%	105.3%	50 150		
Chlorfenapyr	0.000	1.862	2.081	2.000	11.1%	< 30	93.1%	104.1%	50 150		
Chlorpyrifos	0.000	0.412	0.419	0.400	1.6%	< 30	103.1%	104.8%	50 150		
Clofentezine	0.000	0.105	0.107	0.400	1.4%	< 30	26.3%	26.7%	50 150	Q	
Cyfluthrin	0.000	2.305	2.330	2.000	1.1%	< 30	115.3%	116.5%	30 150		
Cypermethrin	0.000	2.569	2.563	2.000	0.2%	< 30	128.4%	128.2%	50 150		
Daminozide	0.000	0.471	0.484	2.000	2.9%	< 30	23.5%	24.2%	30 150	Q	
Diazinon	0.000	0.416	0.422	0.400	1.3%	< 30	104.1%	105.4%	50 150		
Dichlorvos	0.000	1.980	2.022	2.000	2.1%	< 30	99.0%	101.1%	50 150		
Dimethoate	0.000	0.410	0.423	0.400	3.1%	< 30	102.5%	105.7%	50 150		
Ethoprophos	0.000	0.385	0.412	0.400	6.8%	< 30	96.3%	103.1%	50 150		
Etofenprox	0.000	0.671	0.688	0.800	2.5%	< 30	83.9%	86.0%	50 150		
Etoxazole	0.000	0.455	0.471	0.400	3.5%	< 30	113.8%	117.8%	50 150		
Fenoxycarb	0.000	0.386	0.391	0.400	1.4%	< 30	96.4%	97.8%	50 150		
Fenpyroximate	0.000	0.838	0.843	0.800	0.6%	< 30	104.7%	105.4%	50 150		
Fipronil	0.000	0.514	0.534	0.800	3.9%	< 30	64.2%	66.7%	50 150		
Flonicamid	0.000	1.075	1.098	1.000	2.1%	< 30	107.5%	109.8%	50 150		
Fludioxonil	0.000	0.891	0.911	0.800	2.1%	< 30	111.4%	113.8%	50 150		
Hexythiazox	0.000	1.019	1.003	1.000	1.6%	< 30	101.9%	100.3%	50 150		
Imazalil	0.000	0.409	0.420	0.400	2.6%	< 30	102.2%	104.9%	50 150		
Imidacloprid	0.000	0.997	1.042	0.800	4.4%	< 30	124.6%	130.2%	50 150		
Kresoxim methyl	0.000	0.779	0.789	0.800	1.4%	< 30	97.3%	98.6%	50 150		
Malathion	0.000	0.392	0.395	0.400	0.8%	< 30	98.0%	98.8%	50 150		
Metaxalyl	0.000	0.408	0.413	0.400	1.3%	< 30	102.0%	103.4%	50 150		
Methiocarb	0.000	0.405	0.416	0.400	2.7%	< 30	101.3%	104.1%	50 150		
Methomyl	0.000	0.912	0.907	0.800	0.5%	< 30	114.0%	113.4%	50 150		
MGK 264	0.000	0.391	0.410	0.400	4.8%	< 30	97.7%	102.5%	50 150		
Myclobutanil	0.000	0.402	0.420	0.400	4.4%	< 30	100.5%	104.9%	50 150		
Naled	0.000	0.943	0.944	1.000	0.1%	< 30	94.3%	94.4%	50 150		
Oxamyl	0.000	2.261	2.257	2.000	0.2%	< 30	113.1%	112.9%	50 150		
Paclobotrazole	0.000	0.778	0.787	0.800	1.2%	< 30	97.2%	98.4%	50 150		
Parathion Methyl	0.000	0.374	0.408	0.400	8.8%	< 30	93.5%	102.1%	30 150		
Permethrin	0.000	0.444	0.448	0.400	0.9%	< 30	111.0%	112.0%	50 150		
Phosmet	0.000	0.392	0.389	0.400	0.8%	< 30	98.1%	97.3%	50 150		
Piperonyl butoxide	0.099	2.120	2.129	2.000	0.4%	< 30	101.0%	101.5%	50 150		
Prallethrin	0.029	0.431	0.446	0.400	3.6%	< 30	100.5%	104.2%	50 150		
Propiconazole	0.000	0.728	0.737	0.800	1.2%	< 30	91.0%	92.2%	50 150		
Propoxur	0.000	0.402	0.420	0.400	4.5%	< 30	100.5%	105.1%	50 150		
Pyrethrin (Summe)	0.000	0.514	0.522	0.488	1.6%	< 30	105.2%	106.9%	50 150		
Pyridaben	0.000	0.340	0.342	0.400	0.8%	< 30	84.9%	85.6%	50 150		
Spinosad	0.000	0.385	0.393	0.388	2.2%	< 30	99.1%	101.4%	50 150		
Spiromesifen	0.000	0.380	0.383	0.400	0.8%	< 30	94.9%	95.7%	50 150		
Spirotetramat	0.000	0.437	0.448	0.400	2.5%	< 30	109.2%	112.0%	50 150		
Spiroxamine	0.000	0.836	0.851	0.800	1.8%	< 30	104.5%	106.4%	50 150		
ebuconazole	0.000	0.793	0.805	0.800	1.6%	< 30	99.1%	100.7%	50 150		
hiacloprid	0.000	0.422	0.438	0.400	3.8%	< 30	105.5%	109.6%	50 150		
hiamethoxam	0.000	0.464	0.442	0.400	4.9%	< 30	116.0%	110.5%	50 150		
rifloxystrobin	0.000	0.388	0.395	0.400	1.6%	< 30	97.1%	98.6%	50 150		



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Laboratory Quality Control Results

J AOAC 2015 V98-6 Batch ID: 2210437

Laboratory Control Sample									
Analyte	LCS	Result	Spike	Units	% Rec	Limits		Evaluation	Notes
CBDVA	2	0.0353	0.034	%	104	80.0	- 120	Acceptable	
CBDV	2	0.0363	0.035	%	103	80.0	- 120	Acceptable	
CBE	2	0.0365	0.035	%	103	80.0	- 120	Acceptable	
CBDA	1	0.0319	0.032	%	100	90.0	- 110	Acceptable	
CBGA	1	0.0322	0.032	%	100	80.0	- 120	Acceptable	
CBG	1	0.0320	0.032	%	101	80.0	- 120	Acceptable	
CBD	1	0.0315	0.032	%	98.8	90.0	- 110	Acceptable	
THCV	2	0.0350	0.034	%	103	80.0	- 120	Acceptable	
d8THCV	2	0.0373	0.036	%	103	80.0	- 120	Acceptable	
THCVA	2	0.0342	0.033	%	103	80.0	- 120	Acceptable	
CBN	1	0.0331	0.033	%	100	80.0	- 120	Acceptable	
exo-THC	2	0.0336	0.033	%	103	80.0	- 120	Acceptable	
d9THC	1	0.0346	0.034	%	102	90.0	- 110	Acceptable	
d8THC	1	0.0332	0.033	%	99.5	90.0	- 110	Acceptable	
CBL	2	0.0349	0.033	%	105	80.0	- 120	Acceptable	
d10THC	1	0.0304	0.031	%	99.1	80.0	- 120	Acceptable	
CBG	2	0.0352	0.035	%	101	80.0	- 120	Acceptable	
THCA	1	0.0323	0.032	%	101	90.0	- 110	Acceptable	
CBCA	2	0.0353	0.034	%	103	80.0	- 120	Acceptable	
CBLA	2	0.0360	0.035	%	102	80.0	- 120	Acceptable	
CBT	2	0.0367	0.037	%	100	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	
CBCA	<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: 2210437						
Sample Duplicate		Sample ID: 22-014843-0001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBDV	0.00317	0.00320	0.003	%	1.15	< 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	0.00356	0.00363	0.003	%	1.83	< 20	Acceptable	
CBD	0.197	0.198	0.003	%	0.581	< 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	0.00709	0.00721	0.003	%	1.61	< 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	0.00570	0.00581	0.003	%	1.92	< 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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Revision 2 Document D 7087
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Laboratory Quality Control Results

Residual Solvents				Batch ID: 2210461					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		416	572	µg/g	72.7	60	120
Isobutane	ND	< 200		498	731	µg/g	68.1	60	120
Butane	ND	< 200		474	731	µg/g	64.8	60	120
2,2 Dimethylpropane	ND	< 200		654	936	µg/g	69.9	60	120
Methanol	ND	< 200		1130	1620	µg/g	69.8	60	120
Ethylene Oxide	ND	< 30		39	56.2	µg/g	69.4	60	120
2 Methylbutane	ND	< 200		1160	1610	µg/g	72.0	60	120
Pentane	ND	< 200		1150	1600	µg/g	71.9	60	120
Ethanol	ND	< 200		1060	1610	µg/g	65.8	70	130 Q6
Ethyl Ether	ND	< 200		1170	1630	µg/g	71.8	60	120
2,2 Dimethylbutane	ND	< 30		120	171	µg/g	70.2	60	120
Acetone	ND	< 200		1220	1630	µg/g	74.8	60	120
2 Propanol	ND	< 200		1090	1620	µg/g	67.3	60	120
Ethyl Formate	ND	< 500		1430	1670	µg/g	85.6	70	130
Acetonitrile	ND	< 100		346	498	µg/g	69.5	60	120
Methyl Acetate	ND	< 500		1450	1730	µg/g	83.8	70	130
2,3 Dimethylbutane	ND	< 30		128	171	µg/g	74.9	60	120
Dichloromethane	ND	< 60		349	483	µg/g	72.3	60	120
2 Methylpentane	ND	< 30		116	168	µg/g	69.0	60	120
M BE	ND	< 500		1460	1650	µg/g	88.5	70	130
3 Methylpentane	ND	< 30		103	167	µg/g	61.7	60	120
Hexane	ND	< 30		153	182	µg/g	84.1	60	120
1 Propanol	ND	< 500		1360	1620	µg/g	84.0	70	130
Methylethylketone	ND	< 500		1370	1620	µg/g	84.6	70	130
Ethyl acetate	ND	< 200		1110	1610	µg/g	68.9	60	120
2 Butanol	ND	< 200		1010	1600	µg/g	63.1	60	120
tetrahydrofuran	ND	< 100		330	483	µg/g	68.3	60	120
Cyclohexane	ND	< 200		1140	1610	µg/g	70.8	60	120
2 methyl 1 propanol	ND	< 500		1340	1620	µg/g	82.7	70	130
Benzene	ND	< 1		3.97	5.02	µg/g	79.1	60	120
Isopropyl Acetate	ND	< 200		1140	1620	µg/g	70.4	60	120
Heptane	ND	< 200		1030	1610	µg/g	64.0	60	120
1 Butanol	ND	< 500		1390	1630	µg/g	85.3	70	130
Propyl Acetate	ND	< 500		1350	1610	µg/g	83.9	70	130
1,4 Dioxane	ND	< 100		325	491	µg/g	66.2	60	120
2 Ethoxyethanol	ND	< 30		136	181	µg/g	75.1	60	120
Methylisobutylketone	ND	< 500		1340	1620	µg/g	82.7	70	130
3 Methyl 1 butanol	ND	< 500		1330	1630	µg/g	81.6	70	130
Ethylene Glycol	ND	< 200		305	484	µg/g	63.0	60	120
oluene	ND	< 100		329	485	µg/g	67.8	60	120
Isobutyl Acetate	ND	< 500		1350	1630	µg/g	82.8	70	130
1 Pentanol	ND	< 500		1300	1620	µg/g	80.2	70	130
Butyl Acetate	ND	< 500		1330	1620	µg/g	82.1	70	130
Ethylbenzene	ND	< 200		648	969	µg/g	66.9	60	120
m,p Xylene	ND	< 200		662	994	µg/g	66.6	60	120
o Xylene	ND	< 200		646	967	µg/g	66.8	60	120
Cumene	ND	< 30		111	171	µg/g	64.9	60	120
Anisole	ND	< 500		1310	1630	µg/g	80.4	70	130
DMSO	ND	< 500		1430	1680	µg/g	85.1	70	130
1,2 dimethoxyethane	ND	< 50		146	169	µg/g	86.4	70	130
riethylamine	ND	< 500		1420	1630	µg/g	87.1	70	130
N,N dimethylformamide	ND	< 150		417	482	µg/g	86.5	70	130
N,N dimethylacetamide	ND	< 150		411	510	µg/g	80.6	70	130
Pyridine	ND	< 50		165	203	µg/g	81.3	70	130
Sulfolane	ND	< 50		127	172	µg/g	73.8	70	130
1,2 Dichloroethane	ND	< 1		0.872	1	µg/g	87.2	70	130
Chloroform	ND	< 1		0.818	1	µg/g	81.8	70	130
richloroethylene	ND	< 1		0.802	1	µg/g	80.2	70	130
1,1 Dichloroethane	ND	< 1		0.736	1	µg/g	73.6	70	130



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QC - Sample Duplicate		Sample ID: 22-014435-0001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Accept/Fail	Notes
Propane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Isobutane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Butane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2,2 Dimethylpropane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Methanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethylene Oxide	ND	ND	30	µg/g	0.0	< 20	Acceptable	
2 Methylbutane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Pentane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethyl Ether	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2,2 Dimethylbutane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Acetone	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2 Propanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethyl Formate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Acetonitrile	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Methyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
2,3 Dimethylbutane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Dichloromethane	ND	ND	60	µg/g	0.0	< 20	Acceptable	
2 Methylpentane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
M BE	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	
oluene	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	
richloroethylene	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
 Q6 Quality control outside QC limits. Data acceptable based on remaining QC.

Units of Measure:

µg/g Microgram per gram or ppm



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



Report Number: 22-014843/D006.R000
Report Date: 12/12/2022
ORELAP#: OR100028
Purchase Order:
Received: 12/05/22 15:05





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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.