

CONSOLIDATED TEST RESULTS SUMMARY

Please see the following pages for full test results.

BULK SKU SLZ.D9.LM2.6PK		BATCH # EH22V			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 Lemon Light Seltzer		SERVING SIZE 1 can (255 mL)					
LABORATORY: Columbia Laboratories		OREGON ACCREDITATION: OR100028					
POTENCY		PER SERVING		PER GRAM		Percent	
Cannabidiol (CBD)		3.62	mg/serving	0.01	mg/g	0.00	%
Total THC (d9-THC, THCA)		1.58	mg/serving	0.00	mg/g	0.00	%
Cannabigerol (CBG)		<LOQ	mg/serving	<LOQ	mg/g	<LOQ	%
Cannabinol (CBN)		<LOQ	mg/serving	<LOQ	mg/g	<LOQ	%
Cannabichromene (CBC)		<LOQ	mg/serving	<LOQ	mg/g	<LOQ	%
Tetrahydrocannabinolic Acid (THCA)		<LOQ	mg/serving	<LOQ	mg/g	<LOQ	%
Delta-9-THC (d9-THC)		1.58	mg/serving	0.00	mg/g	0.00	%
Delta-8-THC (d8-THC)		0.53	mg/serving	0.00	mg/g	0.00	%
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*		780	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: 22-013057/D004.R000
Report Date: 11/01/2022
ORELAP#: OR100028
Purchase Order:
Received: 10/25/22 15:42

Customer: Etz Hayim Holdings
Product identity: CYCL-SLZ.DP.LM2.6PK-EH22V
Client/Metric ID: .
Laboratory ID: 22-013057-0003

Summary

Potency:

Analyte per 1g	Result	Limits	Units	Status	
CBD per 1g	0.0102		mg/1g		CBD-Total per Serving Size 0.0102 mg/1g
Δ10-THC per 1g	0.000		mg/1g		
Δ8-THC per 1g	0.00148		mg/1g		THC-Total per Serving Size 0.00446 mg/1g
Δ9-THC per 1g	0.00446		mg/1g		
(Reported in milligrams per serving)					

Residual Solvents:

Analyte	Result (μg/g)	Limits (μg/g)	Status
Ethanol	780		

Pesticides:

All analytes passing and less than LOQ.

Metals:

Less than LOQ for all analytes.

Microbiology:

Less than LOQ for all analytes.



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Report Date: 11/01/2022
ORELAP#: OR100028
Purchase Order:
Received: 10/25/22 15:42

Customer: Etz Hayim Holdings
16427 NE Airport Way
PORTLAND 97230
United States of America (USA)
Product identity: CYCL-SLZ.DP.LM2.6PK-EH22V
Client/Metric ID: .
Sample Date:
Laboratory ID: 22-013057-0003
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: 2209202	Analyze: 10/26/22 9:27:00 PM
Analyte	Result	Limits	Units	LOQ	Notes
CBC per 1g	< LOQ		mg/1g	0.000999	
CBC-A per 1g	< LOQ		mg/1g	0.000999	
CBC-Total per 1g	< LOQ		mg/1g	0.00188	
CBD per 1g	0.0102		mg/1g	0.000999	
CBD-A per 1g	< LOQ		mg/1g	0.000999	
CBD-Total per 1g	0.0102		mg/1g	0.00188	
CBDV per 1g	< LOQ		mg/1g	0.000999	
CBDV-A per 1g	< LOQ		mg/1g	0.000999	
CBDV-Total per 1g	< LOQ		mg/1g	0.00186	
CBE per 1g	< LOQ		mg/1g	0.000999	
CBG per 1g	< LOQ		mg/1g	0.000999	
CBG-A per 1g	< LOQ		mg/1g	0.000999	
CBG-Total per 1g	< LOQ		mg/1g	0.00186	
CBL per 1g	< LOQ		mg/1g	0.000999	
CBL-A per 1g	< LOQ		mg/1g	0.000999	
CBL-Total per 1g	< LOQ		mg/1g	0.00188	
CBN per 1g	< LOQ		mg/1g	0.000999	
CBT per 1g	< LOQ		mg/1g	0.000999	
Δ8-THCV per 1g	< LOQ		mg/1g	0.000999	
Δ10-THC per 1g	0.000		mg/1g	0.000	
Δ8-THC per 1g	0.00148		mg/1g	0.000999	
Δ9-THC per 1g	0.00446		mg/1g	0.000999	
exo-THC per 1g	< LOQ		mg/1g	0.000999	
THC-A per 1g	< LOQ		mg/1g	0.000999	
THC-Total per 1g	0.00446		mg/1g	0.00188	
THCV per 1g	< LOQ		mg/1g	0.000999	
THCV-A per 1g	< LOQ		mg/1g	0.000999	
THCV-Total per 1g	< LOQ		mg/1g	0.00188	
Total Cannabinoids per 1g	0.0168		mg/1g		



Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2209204	10/29/22 AOAC 991.14 (Petrifilm) ^P		X
Total Coliforms	< LOQ		cfu/g	10	2209204	10/29/22 AOAC 991.14 (Petrifilm) ^P		X
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2209207	10/30/22 AOAC 2014.05 (RAPID) ^P		X
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2209207	10/30/22 AOAC 2014.05 (RAPID) ^P		X

Solvents

Method: Residual Solvents by GC/MS^b **Units** µg/g **Batch** 2209296 **Analyze** 10/31/22 11:43 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	780		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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Received: 10/25/22 15:42

Pesticides						Method: AOAC 2007.01 & EN 15662 (mod) ^b		Units mg/kg		Batch 2209305		Analyze 10/31/22 01:33 PM	
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes		
Abamectin [¥]	< LOQ	0.50	0.250	pass		Acephate [¥]	< LOQ	0.40	0.250	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		Etioazole [¥]	< 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	0.200	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	0.200	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	0.00395	2209286	10/28/22	AOAC 2013.06 (mod.) ^b	pass	X	
Cadmium	< LOQ	0.200	mg/kg	0.00395	2209286	10/28/22	AOAC 2013.06 (mod.) ^b	pass	X	
Lead	< LOQ	0.500	mg/kg	0.00395	2209286	10/28/22	AOAC 2013.06 (mod.) ^b	pass	X	
Mercury	< LOQ	0.100	mg/kg	0.00198	2209286	10/28/22	AOAC 2013.06 (mod.) ^b	pass	X	

Nutrition										
Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method		Status	Notes	
Water Activity	0.988		Aw	0.030	2209211	10/27/22	AOAC 978.18 ^b		X	



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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

Aw = Water Activity

% wt = µg/g divided by 10,000

Glossary of Qualifiers

X: Not ORELAP accredited.

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

13057



ORELAP ID: OR100028

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
			X	X	X					X	X	X			Beverage		mg/g	Passionfruit flavor
			X		X	X				X	X	X			Beverage		mg/g	watermelon flavor
			X		X	X				X	X	X			Beverage		mg/g	lemon flavor
			X		X	X				X	X	X			Beverage		mg/g	wild cherry flavor
			X		X	X				X	X	X			Beverage		mg/g	ruby grapefruit flavor
																		Parallel path - all
																		LazNat Discount

Collected By:	Relinquished By:	Date	Time	Received by:	Date	Time	Lab Use Only:
☒ Standard (5 day)							Client Alias:
☐ Rush (3-4 day) (1.5x Standard)							Order Number:
☐ Priority Rush (2 day) (2x Standard)							Proper Container
							Sample Condition
							Temperature: 18.4
							Shipped Via: client drop-off
							Evidence of cooling: ☐ Yes ☒ 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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Portland, OR 97230
503-254-1794



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Revision: 1 Document ID: 7148
Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

J AOAC 2015 V98-6 Batch ID: 2209202

Laboratory Control Sample

Analyte	LCS	Result	Spike	Units	% Rec	Limits	Evaluation	Notes
CBDVA	2	0.0009	0.001	%	91.3	80.0 - 120	Acceptable	
CBDV	2	0.0011	0.001	%	102	80.0 - 120	Acceptable	
CBE	2	0.0010	0.001	%	99.9	80.0 - 120	Acceptable	
CBDA	1	0.0010	0.001	%	97.5	90.0 - 110	Acceptable	
CBGA	1	0.0010	0.001	%	98.4	80.0 - 120	Acceptable	
CBG	1	0.0011	0.001	%	106	80.0 - 120	Acceptable	
CBD	1	0.0011	0.001	%	110	90.0 - 110	Acceptable	
THCV	2	0.0011	0.001	%	101	80.0 - 120	Acceptable	
d8THCV	2	0.0011	0.001	%	99.9	80.0 - 120	Acceptable	
THCVA	2	0.0009	0.001	%	92.2	80.0 - 120	Acceptable	
CBN	1	0.0011	0.001	%	108	90.0 - 110	Acceptable	
exo-THC	2	0.0010	0.001	%	101	80.0 - 120	Acceptable	
d9THC	1	0.0011	0.001	%	108	90.0 - 110	Acceptable	
d8THC	1	0.0011	0.001	%	106	90.0 - 110	Acceptable	
CBL	2	0.0010	0.001	%	106	80.0 - 120	Acceptable	
d10THC	1	0.0010	0.001	%	108	80.0 - 120	Acceptable	
CBG	2	0.0011	0.001	%	103	80.0 - 120	Acceptable	
THCA	1	0.0010	0.001	%	97.7	90.0 - 110	Acceptable	
CBGA	2	0.0009	0.001	%	89.5	80.0 - 120	Acceptable	
CBLA	2	0.0010	0.001	%	93.4	80.0 - 120	Acceptable	
CBT	2	0.0011	0.001	%	101	80.0 - 120	Acceptable	

Method Blank

Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBDV	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBE	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBDA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBGA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBG	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBD	<LOQ	0.0001	%	< 0.0001	Acceptable	
THCV	<LOQ	0.0001	%	< 0.0001	Acceptable	
d8THCV	<LOQ	0.0001	%	< 0.0001	Acceptable	
THCVA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBN	<LOQ	0.0001	%	< 0.0001	Acceptable	
exo-THC	<LOQ	0.0001	%	< 0.0001	Acceptable	
d9THC	<LOQ	0.0001	%	< 0.0001	Acceptable	
d8THC	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBL	<LOQ	0.0001	%	< 0.0001	Acceptable	
d10THC	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBG	<LOQ	0.0001	%	< 0.0001	Acceptable	
THCA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBGA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBLA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBT	<LOQ	0.0001	%	< 0.0001	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: 1 Document ID: 7148
Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

J AOAC 2015 V98-6							
Batch ID: 2209202							
Sample Duplicate							
Sample ID: 22-012919-0001							
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation
CBDVA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
CBDV	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
CBE	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
CBD	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
CBGA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
CBG	0.0002	0.0002	0.0001	%	0.155	< 20	Acceptable
CBDA	0.0071	0.0071	0.0001	%	0.539	< 20	Acceptable
THCV	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
d8THCV	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
THCVA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
CBN	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
exo-THC	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
d9THC	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
d8THC	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
CBL	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
d10THC	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
CBC	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
THCA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
CBCA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
CBLA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable
CBT	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable

Abbreviations

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

Units of Measure:



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Laboratory Quality Control Results

Residual Solvents				Batch ID: 2209296					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		604	572	µg/g	105.6	60 - 120	
Isobutane	ND	< 200		778	731	µg/g	106.4	60 - 120	
Butane	ND	< 200		755	731	µg/g	103.3	60 - 120	
2,2-Dimethylpropane	ND	< 200		1180	936	µg/g	126.1	60 - 120	Q1
Methanol	ND	< 200		1990	1650	µg/g	120.6	60 - 120	Q1
Ethylene Oxide	ND	< 30		58.7	56.2	µg/g	104.4	60 - 120	
2-Methylbutane	ND	< 200		1900	1650	µg/g	115.2	60 - 120	
Pentane	ND	< 200		1920	1650	µg/g	116.4	60 - 120	
Ethanol	ND	< 200		1890	1660	µg/g	113.9	70 - 130	
Ethyl Ether	ND	< 200		1830	1630	µg/g	112.3	60 - 120	
2,2-Dimethylbutane	ND	< 30		205	189	µg/g	108.5	60 - 120	
Acetone	ND	< 200		1890	1650	µg/g	114.5	60 - 120	
2-Propanol	ND	< 200		1890	1650	µg/g	114.5	60 - 120	
Ethyl Formate	ND	< 500		1350	1610	µg/g	83.9	70 - 130	
Acetonitrile	ND	< 100		578	504	µg/g	114.7	60 - 120	
Methyl Acetate	ND	< 500		1570	1630	µg/g	96.3	70 - 130	
2,3-Dimethylbutane	ND	< 30		188	174	µg/g	108.0	60 - 120	
Dichloromethane	ND	< 60		530	521	µg/g	101.7	60 - 120	
2-Methylpentane	ND	< 30		192	187	µg/g	102.7	60 - 120	
MTBE	ND	< 500		1520	1600	µg/g	95.0	70 - 130	
3-Methylpentane	ND	< 30		203	188	µg/g	108.0	60 - 120	
Hexane	ND	< 30		203	182	µg/g	111.5	60 - 120	
1-Propanol	ND	< 500		1600	1610	µg/g	99.4	70 - 130	
Methylethylketone	ND	< 500		1510	1600	µg/g	94.4	70 - 130	
Ethyl acetate	ND	< 200		1870	1630	µg/g	114.7	60 - 120	
2-Butanol	ND	< 200		1860	1630	µg/g	114.1	60 - 120	
Tetrahydrofuran	ND	< 100		518	506	µg/g	102.4	60 - 120	
Cyclohexane	ND	< 200		1640	1640	µg/g	100.0	60 - 120	
2-methyl-1-propanol	ND	< 500		1400	1620	µg/g	86.4	70 - 130	
Benzene	ND	< 1		4.68	4.93	µg/g	94.9	60 - 120	
Isopropyl Acetate	ND	< 200		1820	1640	µg/g	111.0	60 - 120	
Heptane	ND	< 200		1680	1630	µg/g	103.1	60 - 120	
1-Butanol	ND	< 500		1420	1600	µg/g	88.8	70 - 130	
Propyl Acetate	ND	< 500		1490	1620	µg/g	92.0	70 - 130	
1,4-Dioxane	ND	< 100		460	493	µg/g	93.3	60 - 120	
2-Ethoxyethanol	ND	< 30		183	171	µg/g	107.0	60 - 120	
Methylisobutylketone	ND	< 500		1310	1620	µg/g	80.9	70 - 130	
3-Methyl-1-butanol	ND	< 500		1350	1610	µg/g	83.9	70 - 130	
Ethylene Glycol	ND	< 200		429	494	µg/g	86.8	60 - 120	
Toluene	ND	< 100		442	506	µg/g	87.4	60 - 120	
Isobutyl Acetate	ND	< 500		1410	1620	µg/g	87.0	70 - 130	
1-Pentanol	ND	< 500		1330	1610	µg/g	82.6	70 - 130	
Butyl Acetate	ND	< 500		1360	1610	µg/g	84.5	70 - 130	
Ethylbenzene	ND	< 200		815	996	µg/g	81.8	60 - 120	
m,p-Xylene	ND	< 200		813	1010	µg/g	80.5	60 - 120	
o-Xylene	ND	< 200		758	979	µg/g	77.4	60 - 120	
Cumene	ND	< 30		139	188	µg/g	73.9	60 - 120	
Anisole	ND	< 500		1110	1610	µg/g	68.9	70 - 130	Q6
DMSO	ND	< 500		1160	1600	µg/g	72.5	70 - 130	
1,2-dimethoxyethane	ND	< 50		174	190	µg/g	91.6	70 - 130	
Triethylamine	ND	< 500		1360	1610	µg/g	84.5	70 - 130	
N,N-dimethylformamide	ND	< 150		373	496	µg/g	75.2	70 - 130	
N,N-dimethylacetamide	ND	< 150		368	483	µg/g	76.2	70 - 130	
Pyridine	ND	< 50		130	167	µg/g	77.8	70 - 130	
Sulfolane	ND	< 50		97.6	161	µg/g	60.6	70 - 130	Q6
1,2-Dichloroethane	ND	< 1		1	1	µg/g	100.0	70 - 130	
Chloroform	ND	< 1		0.983	1	µg/g	98.3	70 - 130	
Trichloroethylene	ND	< 1		0.934	1	µg/g	93.4	70 - 130	
1,1-Dichloroethane	ND	< 1		1.04	1	µg/g	104.0	70 - 130	



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QC - Sample Duplicate		Sample ID: 22-012856-0001						
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	296	306	200	µg/g	3.3	< 20	Acceptable	
Ethyl Ether	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2,2-Dimethylbutane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Acetone	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2-Propanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethyl Formate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Acetonitrile	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Methyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
2,3-Dimethylbutane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Dichloromethane	ND	ND	60	µg/g	0.0	< 20	Acceptable	
2-Methylpentane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
MTBE	ND	ND	500	µg/g	0.0	< 20	Acceptable	
3-Methylpentane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Hexane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
1-Propanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Methyl ethyl ketone	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Ethyl acetate	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2-Butanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Tetrahydrofuran	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Cyclohexane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2-methyl-1-propanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Benzene	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Isopropyl Acetate	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Heptane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
1-Butanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Propyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
1,4-Dioxane	ND	ND	100	µg/g	0.0	< 20	Acceptable	
2-Ethoxyethanol	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Methylisobutylketone	ND	ND	500	µg/g	0.0	< 20	Acceptable	
3-Methyl-1-butanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Ethylene Glycol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Toluene	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Isobutyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
1-Pentanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Butyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Ethylbenzene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
m,p-Xylene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
o-Xylene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Cumene	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Anisole	ND	ND	500	µg/g	0.0	< 20	Acceptable	
DMSO	ND	ND	500	µg/g	0.0	< 20	Acceptable	
1,2-dimethoxyethane	ND	ND	50	µg/g	0.0	< 20	Acceptable	
Triethylamine	ND	ND	500	µg/g	0.0	< 20	Acceptable	
N,N-dimethylformamide	ND	ND	150	µg/g	0.0	< 20	Acceptable	
N,N-dimethylacetamide	ND	ND	150	µg/g	0.0	< 20	Acceptable	
Pyridine	ND	ND	50	µg/g	0.0	< 20	Acceptable	
Sulfolane	ND	ND	50	µg/g	0.0	< 20	Acceptable	
1,2-Dichloroethane	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Chloroform	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Trichloroethylene	ND	ND	1	µg/g	0.0	< 20	Acceptable	
1,1-Dichloroethane	ND	ND	1	µg/g	0.0	< 20	Acceptable	

Abbreviations

ND - None Detected at or above MRL
RPD - Relative Percent Difference
LOQ - Limit of Quantitation
Q1 - Quality control result biased high. Only non-detect samples reported.
Q6 - Quality control outside QC limits. Data acceptable based on remaining QC.

Units of Measure:

µg/g- Microgram per gram or ppm



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Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662									
Units: mg/Kg									
Batch ID: 2209305									
Method Blank			Laboratory Control Sample						
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes	
Abamectin	0.000	< 0.250		0.836	1.000	83.6	50.0 150		
Acephate	0.000	< 0.250		0.904	1.000	90.4	60.0 120		
Acequinocyl	0.000	< 1.000		3.382	4.000	84.6	40.0 160		
Acetamiprid	0.000	< 0.100		0.357	0.400	89.3	60.0 120		
Aldicarb	0.000	< 0.200		0.721	0.800	90.1	60.0 120		
Azoxystrobin	0.000	< 0.100		0.325	0.400	81.2	60.0 120		
Bifenazate	0.000	< 0.100		0.343	0.400	85.7	60.0 120		
Bifenthrin	0.000	< 0.100		0.359	0.400	89.8	50.0 150		
Boscalid	0.000	< 0.200		0.648	0.800	81.0	60.0 120		
Carbaryl	0.000	< 0.100		0.356	0.400	88.9	60.0 120		
Carbofuran	0.000	< 0.100		0.365	0.400	91.1	60.0 120		
Chlorantraniliprole	0.000	< 0.100		0.330	0.400	82.6	60.0 120		
Chlorfenvapir	0.000	< 0.500		1.859	2.000	93.0	60.0 120		
Chlorpyrifos	0.000	< 0.100		0.323	0.400	80.8	60.0 120		
Clofentezine	0.000	< 0.100		0.048	0.400	12.0	60.0 120		Q6
Cyfluthrin	0.000	< 0.500		1.780	2.000	89.0	50.0 150		
Cypermethrin	0.000	< 0.500		1.834	2.000	91.7	50.0 150		
Daminozide	0.000	< 0.500		0.562	2.000	28.1	60.0 120		Q6
Diazinon	0.000	< 0.100		0.322	0.400	80.4	60.0 120		
Dichlorvos	0.000	< 0.500		1.771	2.000	88.5	60.0 120		
Dimethoate	0.000	< 0.100		0.363	0.400	90.7	60.0 120		
Ethoprophos	0.000	< 0.100		0.350	0.400	87.6	60.0 120		
Etofenprox	0.000	< 0.200		0.696	0.800	87.0	50.0 150		
Etoxazole	0.000	< 0.100		0.347	0.400	86.8	60.0 120		
Fenoxycarb	0.000	< 0.100		0.330	0.400	82.4	60.0 120		
Fenpyroximate	0.000	< 0.200		0.696	0.800	87.1	60.0 120		
Fipronil	0.000	< 0.200		0.672	0.800	84.1	60.0 120		
Fonicamid	0.000	< 0.250		0.909	1.000	90.9	60.0 120		
Fludioxonil	0.000	< 0.200		0.699	0.800	87.3	50.0 150		
Hexythiazox	0.000	< 0.250		0.845	1.000	84.5	60.0 120		
Imazalil	0.000	< 0.100		0.340	0.400	85.1	60.0 120		
Imidacloprid	0.000	< 0.200		0.722	0.800	90.2	60.0 120		
Kresoxim-methyl	0.000	< 0.200		0.650	0.800	81.3	60.0 120		
Malathion	0.000	< 0.100		0.341	0.400	85.1	60.0 120		
Metaxyl	0.000	< 0.100		0.358	0.400	89.5	60.0 120		
Methiocarb	0.000	< 0.100		0.342	0.400	85.6	60.0 120		
Methomyl	0.000	< 0.200		0.718	0.800	89.7	60.0 120		
MGK-264	0.000	< 0.100		0.329	0.400	82.3	50.0 150		
Myclobutanil	0.000	< 0.100		0.340	0.400	84.9	60.0 120		
Naled	0.000	< 0.250		0.842	1.000	84.2	50.0 150		
Oxamyl	0.000	< 0.500		1.802	2.000	90.1	60.0 120		
Paclobutrazole	0.000	< 0.200		0.691	0.800	86.4	60.0 120		
Parathion-Methyl	0.000	< 0.200		0.655	0.800	81.9	50.0 150		
Permethrin	0.000	< 0.100		0.357	0.400	89.1	50.0 150		
Phosmet	0.000	< 0.100		0.343	0.400	85.7	50.0 150		
Piperonyl butoxide	0.000	< 0.500		1.609	2.000	80.4	60.0 120		
Prallethrin	0.000	< 0.100		0.331	0.400	82.8	60.0 120		
Propiconazole	0.000	< 0.200		0.682	0.800	85.3	60.0 120		
Propoxur	0.000	< 0.100		0.376	0.400	94.0	60.0 120		
Pyrethrin (Summe)	0.000	< 0.100		0.346	0.413	83.7	60.0 120		
Pyridaben	0.000	< 0.100		0.339	0.400	84.6	50.0 150		
Spirosad	0.000	< 0.100		0.319	0.388	82.2	50.0 150		
Spiromesifen	0.000	< 0.100		0.352	0.400	88.1	60.0 120		
Spirotetramat	0.000	< 0.100		0.337	0.400	84.2	60.0 120		
Spiroxamine	0.000	< 0.200		0.655	0.800	81.9	60.0 120		
Tebuconazole	0.000	< 0.200		0.686	0.800	85.7	60.0 120		
Thiacloprid	0.000	< 0.100		0.353	0.400	88.4	60.0 120		
Thiamethoxam	0.000	< 0.100		0.358	0.400	89.4	60.0 120		
Trifloxystrobin	0.000	< 0.100		0.322	0.400	80.6	60.0 120		



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Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662 Units: mg/Kg Batch ID: 2209305										
Matrix Spike/Matrix Spike Duplicate Recoveries Sample ID: 22-013054-0001										
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes
Abamectin	0.000	0.985	0.948	1.000	3.8%	< 30	98.5%	94.8%	50 - 150	
Acephate	0.000	1.279	1.231	1.000	3.8%	< 30	127.9%	123.1%	50 - 150	
Acequinocyl	0.000	3.853	3.826	4.000	0.7%	< 30	96.3%	95.7%	50 - 150	
Acetamiprid	0.000	0.524	0.508	0.400	3.0%	< 30	131.0%	127.1%	50 - 150	
Aldicarb	0.000	1.042	0.998	0.800	4.2%	< 30	130.2%	124.8%	50 - 150	
Azoxystrobin	0.000	0.414	0.426	0.400	2.9%	< 30	103.5%	106.5%	50 - 150	
Bifenazate	0.000	0.386	0.389	0.400	0.6%	< 30	96.6%	97.2%	50 - 150	
Bifenthrin	0.000	0.391	0.390	0.400	0.2%	< 30	97.7%	97.5%	50 - 150	
Boscalid	0.000	0.789	0.801	0.800	1.5%	< 30	98.6%	100.1%	50 - 150	
Carbaryl	0.000	0.508	0.500	0.400	1.6%	< 30	126.9%	124.9%	50 - 150	
Carbofuran	0.000	0.583	0.579	0.400	0.7%	< 30	145.7%	144.6%	50 - 150	
Chlorantraniliprole	0.000	0.424	0.424	0.400	0.1%	< 30	106.0%	105.9%	50 - 150	
Chlorfenvapir	0.000	2.228	2.025	2.000	9.5%	< 30	111.4%	101.3%	50 - 150	
Chlorpyrifos	0.000	0.345	0.345	0.400	0.1%	< 30	86.1%	86.2%	50 - 150	
Clofentazine	0.000	0.257	0.304	0.400	17.0%	< 30	64.1%	76.0%	50 - 150	
Cyfluthrin	0.000	1.705	1.650	2.000	3.3%	< 30	85.3%	82.5%	30 - 150	
Cypermethrin	0.000	1.641	1.638	2.000	0.2%	< 30	82.0%	81.9%	50 - 150	
Daminozide	0.000	1.615	1.564	2.000	3.2%	< 30	80.7%	78.2%	30 - 150	
Diazinon	0.000	0.367	0.383	0.400	4.3%	< 30	91.7%	95.8%	50 - 150	
Dichlorvos	0.000	2.371	2.183	2.000	8.3%	< 30	118.6%	109.1%	50 - 150	
Dimethoate	0.000	0.449	0.438	0.400	2.5%	< 30	112.3%	109.6%	50 - 150	
Ethoprophos	0.000	0.389	0.379	0.400	2.7%	< 30	97.2%	94.7%	50 - 150	
Etofenprox	0.000	0.797	0.790	0.800	0.9%	< 30	99.6%	98.7%	50 - 150	
Etoxazole	0.000	0.379	0.380	0.400	0.3%	< 30	94.6%	95.0%	50 - 150	
Fenoxycarb	0.000	0.403	0.392	0.400	2.7%	< 30	100.7%	98.0%	50 - 150	
Fenpyroximate	0.000	0.695	0.688	0.800	1.0%	< 30	86.8%	86.0%	50 - 150	
Fipronil	0.000	1.002	0.978	0.800	2.4%	< 30	125.3%	122.3%	50 - 150	
Flonicamid	0.000	0.857	0.826	1.000	3.7%	< 30	85.7%	82.6%	50 - 150	
Fludioxonil	0.000	0.656	0.627	0.800	4.6%	< 30	82.0%	78.3%	50 - 150	
Hexythiazox	0.000	1.077	1.098	1.000	2.0%	< 30	107.7%	109.8%	50 - 150	
Imazalil	0.000	0.429	0.424	0.400	1.2%	< 30	107.2%	105.9%	50 - 150	
Imidacloprid	0.000	0.631	0.619	0.800	1.8%	< 30	78.8%	77.4%	50 - 150	
Kresoxim-methyl	0.000	0.800	0.789	0.800	1.3%	< 30	99.9%	98.6%	50 - 150	
Malathion	0.000	0.436	0.414	0.400	5.0%	< 30	108.9%	103.5%	50 - 150	
Metaxyl	0.000	0.441	0.432	0.400	2.0%	< 30	110.2%	108.0%	50 - 150	
Methiocarb	0.000	0.415	0.410	0.400	1.2%	< 30	103.8%	102.5%	50 - 150	
Methomyl	0.000	0.832	0.831	0.800	0.2%	< 30	104.0%	103.9%	50 - 150	
MGK-264	0.000	0.394	0.392	0.400	0.5%	< 30	98.5%	98.0%	50 - 150	
Myclobutanil	0.000	0.405	0.396	0.400	2.2%	< 30	101.1%	98.9%	50 - 150	
Naled	0.000	1.177	1.153	1.000	2.1%	< 30	117.7%	115.3%	50 - 150	
Oxamyl	0.000	1.977	1.942	2.000	1.8%	< 30	98.9%	97.1%	50 - 150	
Paclobutrazole	0.000	1.025	1.011	0.800	1.4%	< 30	128.1%	126.4%	50 - 150	
Parathion-Methyl	0.000	1.192	1.150	0.800	3.6%	< 30	149.0%	143.8%	30 - 150	
Permethrin	0.000	0.355	0.354	0.400	0.2%	< 30	88.7%	88.5%	50 - 150	
Phosmet	0.000	0.424	0.407	0.400	4.2%	< 30	106.1%	101.7%	50 - 150	
Piperonyl butoxide	0.000	1.891	1.861	2.000	1.6%	< 30	94.6%	93.1%	50 - 150	
Prallethrin	0.000	0.416	0.405	0.400	2.7%	< 30	104.0%	101.2%	50 - 150	
Propiconazole	0.000	0.887	0.868	0.800	2.1%	< 30	110.8%	108.5%	50 - 150	
Propoxur	0.000	0.557	0.536	0.400	3.8%	< 30	139.2%	134.0%	50 - 150	
Pyrethrin (Summe)	0.000	0.389	0.387	0.413	0.4%	< 30	94.1%	93.8%	50 - 150	
Pyridaben	0.000	0.385	0.387	0.400	0.5%	< 30	96.3%	96.8%	50 - 150	
Spirosad	0.000	0.381	0.364	0.388	4.6%	< 30	98.1%	93.8%	50 - 150	
Spiromesifen	0.000	0.381	0.387	0.400	1.5%	< 30	95.3%	96.7%	50 - 150	
Spirotetramat	0.000	0.395	0.388	0.400	1.5%	< 30	98.6%	97.1%	50 - 150	
Spiroxamine	0.000	0.781	0.779	0.800	0.2%	< 30	97.7%	97.4%	50 - 150	
Tebuconazole	0.000	0.935	0.933	0.800	0.2%	< 30	116.8%	116.6%	50 - 150	
Thiacloprid	0.000	0.554	0.544	0.400	1.8%	< 30	138.6%	136.1%	50 - 150	
Thiamethoxam	0.000	0.351	0.337	0.400	4.2%	< 30	87.9%	84.2%	50 - 150	
Trifloxystrobin	0.000	0.369	0.369	0.400	0.0%	< 30	92.2%	92.3%	50 - 150	



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Report Number: 22-013057/D004.R000
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Purchase Order:
Received: 10/25/22 15:42





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.