

## CONSOLIDATED TEST RESULTS SUMMARY

Please see the following pages for full test results.

<b>BULK SKU</b>	<b>BATCH #</b>	<b>LOQ: Limit Of Quantitation</b>	
<b>PRODUCT NAME</b>	<b>SERVING SIZE</b>	<b>LOD: Limit Of Detection</b>	
<b>LABORATORY :</b>	<b>OREGON ACCREDITATION: OR100028</b>	1 g = 10 <sup>-3</sup> kg = 10 <sup>3</sup> mg = 10 <sup>6</sup> µg 1 mg/kg = 1 ppm = 1000 ppb	
POTENCY	PER SERVING	PER GRAM	Percent
Cannabidiol (CBD)	mg/serving	mg/g	%
Total THC (d9-THC, THCA)	mg/serving	mg/g	%
Cannabigerol (CBG)	mg/serving	mg/g	%
Cannabinol (CBN)	mg/serving	mg/g	%
Cannabichromene (CBC)	mg/serving	mg/g	%
Tetrahydrocannabinolic Acid (THCA)	mg/serving	mg/g	%
Delta-9-THC (d9-THC)	mg/serving	mg/g	%
Delta-8-THC (d8-THC)	mg/serving	mg/g	%
HEAVY METALS	PER SERVING	PER GRAM	REGULATORY ACTION LEVEL
Arsenic	µg/serving	µg/g	1.5 ppm
Cadmium	µg/serving	µg/g	0.5 ppm
Lead	µg/serving	µg/g	0.5 ppm
Mercury	µg/serving	µg/g	3.0 ppm
PESTICIDES			REGULATORY ACTION LEVEL
None of the other 59 pesticides tested found above limit of detection in the sample.			10 ppb <sup>[1]</sup>
RESIDUAL SOLVENTS	Results	REGULATORY ACTION LEVEL	
Ethanol*	µg/g	5,000 ppm	
Heptane	µg/g	5,000 ppm	
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:** 24-000398/D006.R000  
**Report Date:** 01/22/2024  
**ORELAP#:** OR100028  
**Purchase Order:** 2801495  
**Received:** 01/10/24 16:17

**Customer:** Etz Hayim Holdings  
**Product identity:** FORM-GMY.D9.HB5.V2-FL22  
**Client/Metric ID:** .  
**Laboratory ID:** 24-000398-0004

### Summary

**Potency:**

Analyte per 1g	Result	Limits	Units	Status	
CBD per 1g	5.23		mg/1g		CBD-Total per Serving Size 5.23 mg/1g
CBDV per 1g	0.0331		mg/1g		
Δ9-THC per 1g	1.07		mg/1g		THC-Total per Serving Size 1.07 mg/1g
(Reported in milligrams per serving)					

**Residual Solvents:**

Analyte	Result (µg/g)	Limits (µg/g)	Status
Ethanol	224		

**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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**Purchase Order:** 2801495  
**Received:** 01/10/24 16:17

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

**Product identity:** FORM-GMY.D9.HB5.V2-FL22

**Client/Metric ID:** .

**Sample Date:**

**Laboratory ID:** 24-000398-0004

**Evidence of Cooling:** No

**Temp:** 20.5 °C

**Relinquished by:** client

**Serving Size #1:** 1 g

### Sample Results

Potency per 1g	Method: J AOAC 2015 V98-6 (mod) <sup>b</sup>	Units mg/se	Batch: 2400364	Analyze: 1/12/24 4:19:00 PM	
Analyte	Result	Limits	Units	LOQ	Notes
CBC per 1g	< LOQ		mg/1g	0.0305	
CBC-A per 1g	< LOQ		mg/1g	0.0305	
CBC-Total per 1g	< LOQ		mg/1g	0.0572	
CBD per 1g	5.23		mg/1g	0.0305	
CBD-A per 1g	< LOQ		mg/1g	0.0305	
CBD-Total per 1g	5.23		mg/1g	0.0572	
CBDV per 1g	0.0331		mg/1g	0.0305	
CBDV-A per 1g	< LOQ		mg/1g	0.0305	
CBDV-Total per 1g	< LOQ		mg/1g	0.0569	
CBE per 1g	< LOQ		mg/1g	0.0305	
CBG per 1g	< LOQ		mg/1g	0.0305	
CBG-A per 1g	< LOQ		mg/1g	0.0305	
CBG-Total per 1g	< LOQ		mg/1g	0.0569	
CBL per 1g	< LOQ		mg/1g	0.0305	
CBL-A per 1g	< LOQ		mg/1g	0.0305	
CBL-Total per 1g	< LOQ		mg/1g	0.0572	
CBN per 1g	< LOQ		mg/1g	0.0305	
CBT per 1g	< LOQ		mg/1g	0.0305	
Δ8-THCV per 1g	< LOQ		mg/1g	0.0305	
Δ10-THC-9R per 1g	< LOQ		mg/1g	0.0305	
Δ10-THC-9S per 1g	< LOQ		mg/1g	0.0305	
Δ10-THC-Total per 1g	< LOQ		mg/1g	0.0609	
Δ8-THC per 1g	< LOQ		mg/1g	0.0305	
Δ9-THC per 1g	1.07		mg/1g	0.0305	
delta-9-THCP per 1g	< LOQ		mg/1g	0.0305	
exo-THC per 1g	< LOQ		mg/1g	0.0305	
THC-A per 1g	< LOQ		mg/1g	0.0305	
THC-Total per 1g	1.07		mg/1g	0.0572	
THCV per 1g	< LOQ		mg/1g	0.0305	
THCV-A per 1g	< LOQ		mg/1g	0.0305	



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Potency per 1g	Method: J AOAC 2015 V98-6 (mod) <sup>P</sup>	Units mg/se	Batch: 2400364	Analyze: 1/12/24 4:19:00 PM	
Analyte	Result	Limits	Units	LOQ	Notes
THCV-Total per 1g	< LOQ		mg/1g	0.0572	
Total Cannabinoids per 1g	6.33		mg/1g		

Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2400283	01/13/24 AOAC 991.14 (Petrifilm) <sup>P</sup>		
Total Coliforms	< LOQ		cfu/g	10	2400283	01/13/24 AOAC 991.14 (Petrifilm) <sup>P</sup>		
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2400284	01/14/24 AOAC 2014.05 (RAPID) <sup>P</sup>		
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2400284	01/14/24 AOAC 2014.05 (RAPID) <sup>P</sup>		

Solvents	Method: Residual Solvents by GC/MS <sup>P</sup>	Units µg/g	Batch 2400463	Analyze 01/22/24 10:36 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	224		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) <sup>b</sup>											
Units mg/kg Batch 2400371 Analyze 01/15/24 02:49 PM											
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
Abamectin <sup>‡</sup>	< LOQ	0.50	0.250	pass		Acephate <sup>‡</sup>	< LOQ	0.40	0.200	pass	
Acequinocyl <sup>‡</sup>	< LOQ	2.0	1.00	pass		Acetamiprid <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Aldicarb <sup>‡</sup>	< LOQ	0.40	0.200	pass		Azoxystrobin <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Bifenazate <sup>‡</sup>	< LOQ	0.20	0.100	pass		Bifenthrin <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Boscalid <sup>‡</sup>	< LOQ	0.40	0.200	pass		Carbaryl <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Carbofuran <sup>‡</sup>	< LOQ	0.20	0.100	pass		Chlorantraniliprole <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Chlorfenapyr <sup>‡</sup>	< LOQ	1.0	0.500	pass		Chlorpyrifos <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Clofentezine <sup>‡</sup>	< LOQ	0.20	0.100	pass		Cyfluthrin <sup>‡</sup>	< LOQ	1.0	0.500	pass	
Cypermethrin <sup>‡</sup>	< LOQ	1.0	0.500	pass		Daminozide <sup>‡</sup>	< LOQ	1.0	0.500	pass	
Diazinon <sup>‡</sup>	< LOQ	0.20	0.100	pass		Dichlorvos <sup>‡</sup>	< LOQ	1.0	0.500	pass	
Dimethoate <sup>‡</sup>	< LOQ	0.20	0.100	pass		Ethoprophos <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Etofenprox <sup>‡</sup>	< LOQ	0.40	0.200	pass		Etoxazole <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Fenoxycarb <sup>‡</sup>	< LOQ	0.20	0.100	pass		Fenpyroximate <sup>‡</sup>	< LOQ	0.40	0.200	pass	
Fipronil <sup>‡</sup>	< LOQ	0.40	0.200	pass		Flonicamid <sup>‡</sup>	< LOQ	1.0	0.400	pass	
Fludioxonil <sup>‡</sup>	< LOQ	0.40	0.200	pass		Hexythiazox <sup>‡</sup>	< LOQ	1.0	0.400	pass	
Imazalil <sup>‡</sup>	< LOQ	0.20	0.100	pass		Imidacloprid <sup>‡</sup>	< LOQ	0.40	0.200	pass	
Kresoxim-methyl <sup>‡</sup>	< LOQ	0.40	0.200	pass		Malathion <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Metalaxyl <sup>‡</sup>	< LOQ	0.20	0.100	pass		Methiocarb <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Methomyl <sup>‡</sup>	< LOQ	0.40	0.200	pass		MGK-264 <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Myclobutanil <sup>‡</sup>	< LOQ	0.20	0.100	pass		Naled <sup>‡</sup>	< LOQ	0.50	0.250	pass	
Oxamyl <sup>‡</sup>	< LOQ	1.0	0.500	pass		Pacllobutrazole <sup>‡</sup>	< LOQ	0.40	0.200	pass	
Parathion-Methyl <sup>‡</sup>	< LOQ	0.20	0.100	pass		Permethrin <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Phosmet <sup>‡</sup>	< LOQ	0.20	0.100	pass		Piperonyl butoxide <sup>‡</sup>	< LOQ	2.0	1.00	pass	
Prallethrin <sup>‡</sup>	< LOQ	0.20	0.100	pass		Propiconazole <sup>‡</sup>	< LOQ	0.40	0.200	pass	
Propoxur <sup>‡</sup>	< LOQ	0.20	0.100	pass		Pyrethrin I (total) <sup>‡</sup>	< LOQ	1.0	0.500	pass	
Pyridaben <sup>‡</sup>	< LOQ	0.20	0.100	pass		Spinosad <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Spiromesifen <sup>‡</sup>	< LOQ	0.20	0.100	pass		Spirotetramat <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Spiroxamine <sup>‡</sup>	< LOQ	0.40	0.200	pass		Tebuconazole <sup>‡</sup>	< LOQ	0.40	0.200	pass	
Thiacloprid <sup>‡</sup>	< LOQ	0.20	0.100	pass		Thiamethoxam <sup>‡</sup>	< LOQ	0.20	0.100	pass	
Trifloxystrobin <sup>‡</sup>	< LOQ	0.20	0.100	pass							

Metals										
Analyte	Result	Limits	Units	LOQ	Batch	Analyzed	Method	Status	Notes	
Arsenic <sup>‡</sup>	< LOQ	0.200	mg/kg	0.0141	2400340	01/12/24	AOAC 2013.06 (mod.) <sup>b</sup>	pass		
Cadmium <sup>‡</sup>	< LOQ	0.200	mg/kg	0.0141	2400340	01/12/24	AOAC 2013.06 (mod.) <sup>b</sup>	pass		
Lead <sup>‡</sup>	< LOQ	0.500	mg/kg	0.0141	2400340	01/12/24	AOAC 2013.06 (mod.) <sup>b</sup>	pass		
Mercury <sup>‡</sup>	< LOQ	0.100	mg/kg	0.00705	2400340	01/12/24	AOAC 2013.06 (mod.) <sup>b</sup>	pass		



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**Received:** 01/10/24 16:17

### Abbreviations

**Limits:** Action Levels per OAR-333-007-0400, OAR-333-007-0210, OAR-333-007-0220, CCR title 16-division 42. BCC-section 5723

**Limit(s) of Quantitation (LOQ):** The minimum levels, concentrations, or quantities of a target variable (e.g., target analyte) that can be reported with a specified degree of confidence.

Ⓐ = ISO/IEC 17025:2017 accredited method.

Ⓜ = TNI accredited analyte.

### Units of Measure

cfu/g = Colony forming units per gram

g = g

µg/g = Microgram per gram

mg/kg = Milligram per kilogram = parts per million (ppm)

mg/1g = Milligram per 1g

% = Percentage of sample

% wt = µg/g divided by 10,000

Approved Signatory

Derrick Tanner  
General Manager



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

Laboratory Quality Control Results

JAOAC2015 V986 Batch ID: 2400364

Analyte	LCS	Result	Spike	Units	% Rec	Limits	Evaluation	Notes
CBVA	2	0.0324	0.0324	%	100	80.0 - 120	Acceptable	
CBV	2	0.0343	0.0347	%	98.8	80.0 - 120	Acceptable	
CEE	2	0.0348	0.0350	%	99.3	80.0 - 120	Acceptable	
CBDA	1	0.0321	0.0317	%	101	90.0 - 110	Acceptable	
CBSA	1	0.0319	0.0315	%	101	80.0 - 120	Acceptable	
CBG	1	0.0310	0.0309	%	101	80.0 - 120	Acceptable	
CB	1	0.0334	0.0330	%	101	90.0 - 110	Acceptable	
THCV	2	0.0344	0.0345	%	99.8	80.0 - 120	Acceptable	
δ8THCV	2	0.0301	0.0303	%	99.2	80.0 - 120	Acceptable	
THCVA	2	0.0317	0.0318	%	99.7	80.0 - 120	Acceptable	
CBN	1	0.0339	0.0330	%	103	80.0 - 120	Acceptable	
exo-THC	2	0.0326	0.0330	%	98.6	80.0 - 120	Acceptable	
δ9THC	1	0.0343	0.0337	%	102	90.0 - 110	Acceptable	
δ8THC	1	0.0344	0.0336	%	102	90.0 - 110	Acceptable	
9SaTHC	1	0.0331	0.0326	%	102	80.0 - 120	Acceptable	
CB	2	0.0319	0.0326	%	97.9	80.0 - 120	Acceptable	
9RaTHC	1	0.0319	0.0318	%	100	80.0 - 120	Acceptable	
CB	2	0.0343	0.0349	%	98.4	80.0 - 120	Acceptable	
THCA	1	0.0330	0.0322	%	102	90.0 - 110	Acceptable	
CBCA	2	0.0332	0.0334	%	99.4	80.0 - 120	Acceptable	
CBLA	2	0.0336	0.0338	%	99.2	80.0 - 120	Acceptable	
δ9THCP	2	0.0317	0.0328	%	96.8	80.0 - 120	Acceptable	
CB	2	0.0327	0.0347	%	94.4	80.0 - 120	Acceptable	

Analyte	Result	LOG	Units	Limits	Evaluation	Notes
CBVA	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBV	<LOQ	0.00327	%	< 0.00327	Acceptable	
CEE	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBDA	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBSA	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBG	<LOQ	0.00327	%	< 0.00327	Acceptable	
CB	<LOQ	0.00327	%	< 0.00327	Acceptable	
THCV	<LOQ	0.00327	%	< 0.00327	Acceptable	
δ8THCV	<LOQ	0.00327	%	< 0.00327	Acceptable	
THCVA	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBN	<LOQ	0.00327	%	< 0.00327	Acceptable	
exo-THC	<LOQ	0.00327	%	< 0.00327	Acceptable	
δ9THC	<LOQ	0.00327	%	< 0.00327	Acceptable	
δ8THC	<LOQ	0.00327	%	< 0.00327	Acceptable	
9SaTHC	<LOQ	0.00327	%	< 0.00327	Acceptable	
CB	<LOQ	0.00327	%	< 0.00327	Acceptable	
9RaTHC	<LOQ	0.00327	%	< 0.00327	Acceptable	
CB	<LOQ	0.00327	%	< 0.00327	Acceptable	
THCA	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBCA	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBLA	<LOQ	0.00327	%	< 0.00327	Acceptable	
δ9THCP	<LOQ	0.00327	%	< 0.00327	Acceptable	
CB	<LOQ	0.00327	%	< 0.00327	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  
 Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

JAOAC2015 V986		Batch ID: 2400364						
Sample Duplicate		Sample ID: 24-0003320001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
CBDV	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
CBE	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
CBDVA	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
CBSA	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
CBS	0.00398	0.00375	0.00316	%	5.94	< 20	Acceptable	
CBD	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
δ8THCV	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
CBN	0.00822	0.00783	0.00316	%	4.91	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
δ9THC	0.259	0.246	0.00316	%	5.08	< 20	Acceptable	
δ8THC	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
9Sα10THC	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
CB	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
9Rα10THC	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
CB	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
CBSA	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
δ9THCP	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	
CB	<LOQ	<LOQ	0.00316	%	NA	< 20	Acceptable	

Abbreviations

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

Units of Measure:

% - Percent





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



Report Number: 24-000398/D006.R000  
Report Date: 01/22/2024  
ORELAP#: OR100028  
Purchase Order: 2801495  
Received: 01/10/24 16:17

Revision: 3 Document ID: 3120  
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Laboratory Pesticide Quality Control Results

AOAC2007.1 & EN 15662		Units: mg/Kg			Batch ID 2400371				
Method	Blank	Laboratory Control Sample							
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spk	LCS % Re	Limits	Notes	
Abamectin	0.00	< 0.250		0.920	1.000	92.0	50.0 150		
Acophate	0.014	< 0.200		0.655	0.800	81.9	60.0 120		
Acquinocyl	0.000	< 1.000		4.015	4.000	100.4	40.0 160		
Acetamiprid	0.000	< 0.100		0.351	0.400	87.8	60.0 120		
Aldicarb	0.000	< 0.200		0.749	0.800	93.6	60.0 120		
Azoxystrobin	0.005	< 0.100		0.369	0.400	92.2	60.0 120		
Bifenazate	0.000	< 0.100		0.370	0.400	92.5	60.0 120		
Bifenthrin	0.000	< 0.100		0.355	0.400	89.1	50.0 150		
Boscalid	0.000	< 0.200		0.719	0.800	89.9	60.0 120		
Carbaryl	0.000	< 0.100		0.357	0.400	89.4	60.0 120		
Carbofuran	0.000	< 0.100		0.353	0.400	88.2	60.0 120		
Chlorantraniliprole	0.000	< 0.100		0.347	0.400	86.7	60.0 120		
Chlorfenapyr	0.029	< 0.500		2.013	2.000	100.6	60.0 120		
Chlorpyrifos	0.000	< 0.100		0.338	0.400	84.5	60.0 120		
Clofentezine	0.000	< 0.100		0.325	0.400	81.3	60.0 120		
Cyfluthrin	0.000	< 0.500		1.837	2.000	91.9	50.0 150		
Cypermethrin	0.000	< 0.500		1.816	2.000	90.8	50.0 150		
Daminozide	0.000	< 0.500		0.631	2.000	31.6	60.0 120	Q6	
Diazinon	0.000	< 0.100		0.375	0.400	93.7	60.0 120		
Dichlorvos	0.000	< 0.500		1.585	2.000	79.3	60.0 120		
Dimethoate	0.000	< 0.100		0.351	0.400	87.7	60.0 120		
Ethoprophos	0.000	< 0.100		0.341	0.400	85.4	60.0 120		
Etofenprox	0.000	< 0.200		0.753	0.800	94.1	50.0 150		
Etoxazole	0.000	< 0.100		0.389	0.400	97.2	60.0 120		
Fenoxycarb	0.000	< 0.100		0.367	0.400	91.8	60.0 120		
Fenpyroximate	0.000	< 0.200		0.737	0.800	92.1	60.0 120		
Fipronil	0.000	< 0.200		0.730	0.800	91.2	60.0 120		
Fonicamid	0.000	< 0.250		0.678	1.000	67.8	60.0 120		
Fludioxonil	0.000	< 0.200		0.681	0.800	85.1	50.0 150		
Hexythiazox	0.000	< 0.250		0.927	1.000	92.7	60.0 120		
Imazalil	0.000	< 0.100		0.339	0.400	84.7	60.0 120		
Imidacloprid	0.000	< 0.200		0.597	0.800	74.6	60.0 120		
Kresoxim-methyl	0.000	< 0.200		0.713	0.800	89.1	60.0 120		
Malathion	0.000	< 0.100		0.362	0.400	90.5	60.0 120		
Metaxalyl	0.000	< 0.100		0.351	0.400	87.7	60.0 120		
Methiocarb	0.000	< 0.100		0.366	0.400	91.4	60.0 120		
Methomyl	0.000	< 0.200		0.555	0.800	69.3	60.0 120		
MCK-264	0.000	< 0.100		0.362	0.400	90.6	50.0 150		
Mydobutanol	0.000	< 0.100		0.333	0.400	83.3	60.0 120		
Naled	0.000	< 0.250		0.848	1.000	84.8	50.0 150		
Oxamyl	0.000	< 0.500		1.490	2.000	74.5	60.0 120		
Padobutrazole	0.000	< 0.200		0.711	0.800	88.9	60.0 120		
Parathion-Methyl	0.024	< 0.100		0.359	0.400	89.7	50.0 150		
Permethrin	0.000	< 0.100		0.365	0.400	91.5	50.0 150		
Phosmet	0.000	< 0.100		0.374	0.400	93.4	50.0 150		
Piperonyl butoxide	0.000	< 0.500		1.821	2.000	91.1	60.0 120		
Prallethrin	0.000	< 0.100		0.349	0.400	87.3	60.0 120		
Propiconazole	0.000	< 0.200		0.730	0.800	91.2	60.0 120		
Propoxur	0.000	< 0.100		0.345	0.400	86.6	60.0 120		
Pyrethrin (Summe)	0.000	< 0.100		0.455	0.488	93.2	60.0 120		
Pyridaben	0.000	< 0.100		0.357	0.400	89.3	50.0 150		
Spinosad	0.000	< 0.100		0.353	0.388	90.9	50.0 150		
Spiromesifen	0.000	< 0.100		0.360	0.400	90.0	60.0 120		
Spirotetramat	0.000	< 0.100		0.339	0.400	84.6	60.0 120		
Spiroxamine	0.000	< 0.200		0.716	0.800	89.5	60.0 120		
Tebuconazole	0.000	< 0.200		0.712	0.800	89.0	60.0 120		
Thiadoprid	0.000	< 0.100		0.359	0.400	89.7	60.0 120		
Thiamethoxam	0.000	< 0.100		0.304	0.400	75.9	60.0 120		
Trifloxystrobin	0.000	< 0.100		0.345	0.400	86.2	60.0 120		



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Laboratory Pesticide Quality Control Results

AOAC2007.1 & EN 15662		Units: mg/Kg				Batch ID 2400371					
Matrix Spke/Matrix Spke Duplicate Recoveries		Sample ID: 24-0002810003									
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Re	MSD % Re	Limits	Notes	
Abamectin	0.00	0.877	0.900	1.000	2.6%	< 30	87.7%	90.0%	50 - 150		
Acophate	0.00	0.659	0.652	0.800	1.0%	< 30	82.3%	81.5%	50 - 150		
Acquinocyl	0.00	4.924	3.673	4.000	29.1%	< 30	123.1%	91.8%	50 - 150		
Acetamiprid	0.00	0.378	0.349	0.400	8.1%	< 30	94.6%	87.3%	50 - 150		
Aldicarb	0.00	0.748	0.746	0.800	0.2%	< 30	93.5%	93.3%	50 - 150		
Azoxystrobin	0.00	0.351	0.319	0.400	9.8%	< 30	86.5%	78.5%	50 - 150		
Bifenazate	0.00	0.357	0.344	0.400	3.7%	< 30	89.3%	86.1%	50 - 150		
Bifenthrin	0.00	0.310	0.339	0.400	8.8%	< 30	77.6%	84.7%	50 - 150		
Boscalid	0.00	0.763	0.697	0.800	9.1%	< 30	95.4%	87.1%	50 - 150		
Carbaryl	0.00	0.363	0.320	0.400	12.6%	< 30	90.9%	80.1%	50 - 150		
Carbofuran	0.00	0.328	0.311	0.400	5.3%	< 30	82.1%	77.8%	50 - 150		
Chlorantraniliprole	0.00	0.330	0.332	0.400	0.7%	< 30	82.5%	83.1%	50 - 150		
Chlorfenapyr	0.028	1.727	1.810	2.000	4.8%	< 30	85.0%	89.1%	50 - 150		
Chlorpyrifos	0.00	0.325	0.310	0.400	4.5%	< 30	81.2%	77.6%	50 - 150		
Cofentazine	0.00	0.298	0.281	0.400	6.1%	< 30	74.6%	70.1%	50 - 150		
Cyfluthrin	0.00	1.953	1.960	2.000	0.4%	< 30	97.7%	98.0%	30 - 150		
Cypermethrin	0.00	1.920	1.963	2.000	2.2%	< 30	96.0%	98.2%	50 - 150		
Daminozide	0.00	0.616	0.688	2.000	11.0%	< 30	30.8%	34.4%	30 - 150		
Diazinon	0.00	0.346	0.362	0.400	4.7%	< 30	86.4%	90.6%	50 - 150		
Dichlorvos	0.00	1.694	1.617	2.000	4.6%	< 30	84.7%	80.9%	50 - 150		
Dimethoate	0.00	0.353	0.342	0.400	3.1%	< 30	88.2%	85.5%	50 - 150		
Ethoprophos	0.00	0.338	0.338	0.400	0.1%	< 30	84.5%	84.6%	50 - 150		
Etofenprox	0.00	0.631	0.677	0.800	7.0%	< 30	78.8%	84.6%	50 - 150		
Etoxazole	0.00	0.369	0.370	0.400	0.5%	< 30	92.2%	92.6%	50 - 150		
Fenoxycarb	0.00	0.339	0.340	0.400	0.3%	< 30	84.8%	85.0%	50 - 150		
Fenpyroximate	0.00	0.706	0.727	0.800	2.9%	< 30	88.2%	90.9%	50 - 150		
Fipronil	0.00	0.767	0.712	0.800	7.4%	< 30	95.8%	89.0%	50 - 150		
Fonicamid	0.00	0.756	0.846	1.000	11.2%	< 30	75.6%	84.6%	50 - 150		
Fludioxonil	0.00	0.682	0.601	0.800	12.7%	< 30	85.3%	75.1%	50 - 150		
Hexythiazox	0.00	1.486	1.453	1.000	2.2%	< 30	148.5%	145.3%	50 - 150		
Imazalil	0.00	0.345	0.338	0.400	2.1%	< 30	86.2%	84.4%	50 - 150		
Imidacloprid	0.00	0.649	0.651	0.800	0.3%	< 30	81.1%	81.3%	50 - 150		
Kresoxim-methyl	0.00	0.700	0.646	0.800	8.2%	< 30	87.5%	80.6%	50 - 150		
Malathion	0.00	0.321	0.331	0.400	2.9%	< 30	80.3%	82.7%	50 - 150		
Metaxalyl	0.00	0.347	0.330	0.400	5.0%	< 30	86.7%	82.5%	50 - 150		
Methiocarb	0.00	0.370	0.352	0.400	4.9%	< 30	92.4%	88.0%	50 - 150		
Methomyl	0.00	0.605	0.694	0.800	13.8%	< 30	75.6%	86.6%	50 - 150		
MGK-264	0.00	0.333	0.311	0.400	7.0%	< 30	83.3%	77.6%	50 - 150		
Mydobutanol	0.00	0.319	0.347	0.400	8.5%	< 30	79.7%	86.6%	50 - 150		
Naled	0.00	0.796	0.800	1.000	0.5%	< 30	79.6%	80.0%	50 - 150		
Oxaryl	0.00	1.594	1.651	2.000	4.2%	< 30	79.2%	82.6%	50 - 150		
Padobutrazole	0.00	0.707	0.626	0.800	12.2%	< 30	88.4%	78.2%	50 - 150		
Parathion-Methyl	0.024	0.350	0.366	0.400	4.9%	< 30	81.5%	85.6%	30 - 150		
Permethrin	0.00	0.312	0.334	0.400	6.8%	< 30	78.0%	83.3%	50 - 150		
Phosmet	0.00	0.367	0.350	0.400	4.8%	< 30	91.8%	87.9%	50 - 150		
Piperonyl butoxide	0.00	1.734	1.643	2.000	5.4%	< 30	86.7%	82.2%	50 - 150		
Prallethrin	0.00	0.344	0.342	0.400	0.7%	< 30	86.0%	85.4%	50 - 150		
Propiconazole	0.00	0.710	0.699	0.800	1.6%	< 30	88.8%	87.4%	50 - 150		
Propoxur	0.00	0.332	0.332	0.400	0.0%	< 30	83.0%	83.1%	50 - 150		
Pyrethrin (Summe)	0.00	0.431	0.433	0.488	0.5%	< 30	88.2%	88.7%	50 - 150		
Pyridaben	0.00	0.374	0.366	0.400	5.8%	< 30	93.4%	99.0%	50 - 150		
Spinosad	0.00	0.348	0.324	0.388	7.2%	< 30	89.7%	83.4%	50 - 150		
Spiromesfen	0.00	0.358	0.359	0.400	0.4%	< 30	89.9%	89.8%	50 - 150		
Spirotetramat	0.00	0.348	0.331	0.400	5.2%	< 30	87.1%	82.6%	50 - 150		
Spiroxamine	0.00	0.681	0.694	0.800	1.9%	< 30	85.1%	86.7%	50 - 150		
Tebuconazole	0.00	0.728	0.735	0.800	1.0%	< 30	91.0%	91.9%	50 - 150		
Thiadoprid	0.00	0.358	0.350	0.400	2.3%	< 30	89.6%	87.5%	50 - 150		
Thiamethoxam	0.00	0.338	0.376	0.400	10.7%	< 30	84.5%	94.0%	50 - 150		
Trifloxystrobin	0.00	0.329	0.307	0.400	7.0%	< 30	82.3%	76.7%	50 - 150		



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Laboratory Quality Control Results

Residual Solvents				Batch ID: 2400463					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		657	584	µg/g	112.5	60 - 120	
Isobutane	ND	< 200		806	767	µg/g	105.1	60 - 120	
Butane	ND	< 200		832	782	µg/g	106.4	60 - 120	
2,2-Dimethylpropane	ND	< 200		1080	939	µg/g	115.0	60 - 120	
Methanol	ND	< 200		1800	1600	µg/g	112.5	60 - 120	
Ethylene Oxide	ND	< 30		63.6	57.1	µg/g	111.4	60 - 120	
2-Methylbutane	ND	< 200		1680	1600	µg/g	105.0	60 - 120	
Pentane	ND	< 200		1640	1600	µg/g	102.5	60 - 120	
Ethanol	ND	< 200		1610	1600	µg/g	100.6	70 - 130	
Ethyl Ether	ND	< 200		1630	1600	µg/g	101.9	60 - 120	
2,2-Dimethylbutane	ND	< 30		165	161	µg/g	102.5	60 - 120	
Acetone	ND	< 200		1700	1600	µg/g	106.3	60 - 120	
2-Propanol	ND	< 200		1750	1600	µg/g	109.4	60 - 120	
Ethyl Formate	ND	< 500		1150	1600	µg/g	71.9	70 - 130	
Acetonitrile	ND	< 100		496	488	µg/g	101.6	60 - 120	
Methyl Acetate	ND	< 500		1640	1610	µg/g	101.9	70 - 130	
2,3-Dimethylbutane	ND	< 30		187	163	µg/g	114.7	60 - 120	
Dichloromethane	ND	< 60		482	488	µg/g	98.8	60 - 120	
2-Methylpentane	ND	< 30		98.4	161	µg/g	61.1	60 - 120	
MTBE	ND	< 500		1790	1650	µg/g	108.5	70 - 130	
3-Methylpentane	ND	< 30		165	162	µg/g	101.9	60 - 120	
Hexane	ND	< 30		151	161	µg/g	93.8	60 - 120	
1-Propanol	ND	< 500		1790	1620	µg/g	110.5	70 - 130	
Methylethylketone	ND	< 500		1660	1610	µg/g	103.1	70 - 130	
Ethyl acetate	ND	< 200		1700	1610	µg/g	105.6	60 - 120	
2-Butanol	ND	< 200		1720	1610	µg/g	106.8	60 - 120	
Tetrahydrofuran	ND	< 100		484	483	µg/g	100.2	60 - 120	
Cyclohexane	ND	< 200		1590	1600	µg/g	99.4	60 - 120	
2-methyl-1-propanol	ND	< 500		1510	1600	µg/g	94.4	70 - 130	
Benzene	ND	< 1		3.07	4.99	µg/g	61.5	60 - 120	
Isopropyl Acetate	ND	< 200		1690	1600	µg/g	105.6	60 - 120	
Heptane	ND	< 200		1670	1600	µg/g	104.4	60 - 120	
1-Butanol	ND	< 500		1540	1610	µg/g	95.7	70 - 130	
Propyl Acetate	ND	< 500		1660	1610	µg/g	103.1	70 - 130	
1,4-Dioxane	ND	< 100		463	480	µg/g	96.5	60 - 120	
2-Ethoxyethanol	ND	< 30		158	161	µg/g	98.1	60 - 120	
Methylisobutylketone	ND	< 500		1610	1610	µg/g	100.0	70 - 130	
3-Methyl-1-butanol	ND	< 500		1520	1610	µg/g	94.4	70 - 130	
Ethylene Glycol	ND	< 200		341	481	µg/g	70.9	60 - 120	
Toluene	ND	< 100		467	483	µg/g	96.7	60 - 120	
Isobutyl Acetate	ND	< 500		1640	1610	µg/g	101.9	70 - 130	
1-Pentanol	ND	< 500		1570	1610	µg/g	97.5	70 - 130	
Butyl Acetate	ND	< 500		1570	1600	µg/g	98.1	70 - 130	
Ethylbenzene	ND	< 200		938	962	µg/g	97.5	60 - 120	
m,p-Xylene	ND	< 200		928	972	µg/g	95.5	60 - 120	
o-Xylene	ND	< 200		914	965	µg/g	94.7	60 - 120	
Cumene	ND	< 30		163	169	µg/g	96.4	60 - 120	
Anisole	ND	< 500		1390	1600	µg/g	86.9	70 - 130	
DMSO	ND	< 500		1000	1600	µg/g	62.5	70 - 130	Q6
1,2-dimethoxyethane	ND	< 50		166	163	µg/g	101.8	70 - 130	
Triethylamine	ND	< 500		1510	1600	µg/g	94.4	70 - 130	
N,N-dimethylformamide	ND	< 150		424	482	µg/g	88.0	70 - 130	
N,N-dimethylacetamide	ND	< 150		480	483	µg/g	99.4	70 - 130	
Pyridine	ND	< 50		165	161	µg/g	102.5	70 - 130	
Silfolane	ND	< 50		112	163	µg/g	68.7	70 - 130	Q6
1,2-Dichloroethane	ND	< 1		0.83	1	µg/g	83.0	70 - 130	
Chloroform	ND	< 1		0.903	1	µg/g	90.3	70 - 130	
Trichloroethylene	ND	< 1		0.703	1	µg/g	70.3	70 - 130	
1,1-Dichloroethane	ND	< 1		0.895	1	µg/g	89.5	70 - 130	



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QC- Sample Duplicate		Sample ID: 24-000384-0001						
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		
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  
 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:** 24-000398/D006.R000  
**Report Date:** 01/22/2024  
**ORELAP#:** OR100028  
**Purchase Order:** 2801495  
**Received:** 01/10/24 16:17





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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.