



This is an amended version of report# 24-010742/D007.R000.

Reason: updated client information

Customer: Seventh Hill Releaf LLC
Product identity: Focus CBG Caramel
Client/Metric ID: .
Laboratory ID: 24-010742-0007

Summary

Potency:

Analyte	Result	Limits	Units	Status	
CBD	0.109		%		CBD-Total per Serving Size 18.5 mg/17g
CBG	0.115		%		
Δ9-THC	0.0295		%		Delta-9-THC-Total per 5.02 mg/17g
Analyte per 17g	Result	Limits	Units	Status	(Reported in milligrams per serving)
CBD per 17g	18.5		mg/17g		
CBG per 17g	19.6		mg/17g		
Δ9-THC per 17g	5.02		mg/17g		

Residual Solvents:

All analytes passing and less than LOQ.

Pesticides:

All analytes passing and less than LOQ.

Metals:

Less than LOQ for all analytes.

Microbiology:

Less than LOQ for all analytes.



Customer: Seventh Hill Releaf LLC
 215 S 19th St
 Springfield Oregon 97477
 United States of America (USA)

Product identity: Focus CBG Caramel
Client/Metric ID: .
Sample Date:
Laboratory ID: 24-010742-0007
Evidence of Cooling: No
Temp: 23 °C
Relinquished by: shipping
Lot #: FCCBG24251
Serving Size #1: 17 g

Sample Results

Potency	Method: J AOAC 2015 V98-6 (mod) ^b			Units %	Batch: 2407447	Analyze: 9/26/24 2:40:00 AM
Analyte	Result	Limits	Units	LOQ	Notes	
CBC	< LOQ		%	0.00308		
CBC-A	< LOQ		%	0.00308		
CBC-Total	< LOQ		%	0.00579		
CBD [±]	0.109		%	0.00308		
CBD-A [±]	< LOQ		%	0.00308		
CBD-Total [±]	0.109		%	0.00579		
CBDV	< LOQ		%	0.00308		
CBDV-A	< LOQ		%	0.00308		
CBDV-Total	< LOQ		%	0.00575		
CBE	< LOQ		%	0.00308		
CBG	0.115		%	0.00308		
CBG-A	< LOQ		%	0.00308		
CBG-Total	0.115		%	0.00575		
CBL	< LOQ		%	0.00308		
CBL-A	< LOQ		%	0.00308		
CBL-Total	< LOQ		%	0.00579		
CBN	< LOQ		%	0.00308		
CBT	< LOQ		%	0.00308		
Δ10-THC-9R	< LOQ		%	0.00308		
Δ10-THC-9S	< LOQ		%	0.00308		
Δ10-THC-Total	< LOQ		%	0.00616		
Δ8-THC [±]	< LOQ		%	0.00308		
Δ8-THCV	< LOQ		%	0.00308		
Δ9-THC [±]	0.0295		%	0.00308		
Δ9-THC-A [±]	< LOQ		%	0.00308		
Δ9-THC-Total [±]	0.0295		%	0.00579		
Δ9-THCP	< LOQ		%	0.00308		
Δ9-THCV	< LOQ		%	0.00308		
Δ9-THCV-A	< LOQ		%	0.00308		



Potency	Method: J AOAC 2015 V98-6 (mod) ^b		Units %	Batch: 2407447	Analyze: 9/26/24 2:40:00 AM
Analyte	Result	Limits	Units	LOQ	Notes
Δ9-THCV-Total	< LOQ		%	0.00575	
exo-THC	< LOQ		%	0.00308	
Total Cannabinoids	0.254		%		

Potency per 17g	Method: J AOAC 2015 V98-6 (mod) ^b	Units mg/se	Batch: 2407447	Analyze: 9/26/24 2:40:00 AM	
Analyte	Result	Limits	Units	LOQ	Notes
CBC per 17g	< LOQ		mg/17g	0.524	
CBC-A per 17g	< LOQ		mg/17g	0.524	
CBC-Total per 17g	< LOQ		mg/17g	0.983	
CBD per 17g	18.5		mg/17g	0.524	
CBD-A per 17g [⊥]	< LOQ		mg/17g	0.524	
CBD-Total per 17g [⊥]	18.5		mg/17g	0.983	
CBDV per 17g	< LOQ		mg/17g	0.524	
CBDV-A per 17g	< LOQ		mg/17g	0.524	
CBDV-Total per 17g	< LOQ		mg/17g	0.978	
CBE per 17g	< LOQ		mg/17g	0.524	
CBG per 17g	19.6		mg/17g	0.524	
CBG-A per 17g	< LOQ		mg/17g	0.524	
CBG-Total per 17g	19.6		mg/17g	0.978	
CBL per 17g	< LOQ		mg/17g	0.524	
CBL-A per 17g	< LOQ		mg/17g	0.524	
CBL-Total per 17g	< LOQ		mg/17g	0.983	
CBN per 17g	< LOQ		mg/17g	0.524	
CBT per 17g	< LOQ		mg/17g	0.524	
Δ10-THC-9R per 17g	< LOQ		mg/17g	0.524	
Δ10-THC-9S per 17g	< LOQ		mg/17g	0.524	
Δ10-THC-Total per 17g	< LOQ		mg/17g	1.05	
Δ8-THC per 17g [⊥]	< LOQ		mg/17g	0.524	
Δ8-THCV per 17g	< LOQ		mg/17g	0.524	
Δ9-THC per 17g [⊥]	5.02		mg/17g	0.524	
Δ9-THC-Total per 17g	5.02		mg/17g	0.983	
Δ9-THCP per 17g	< LOQ		mg/17g	0.524	
Δ9-THCV per 17g	< LOQ		mg/17g	0.524	
Δ9-THCV-A per 17g	< LOQ		mg/17g	0.524	
Δ9-THCV-Total per 17g	< LOQ		mg/17g	0.984	
exo-THC per 17g	< LOQ		mg/17g	0.524	
THC-A per 17g [⊥]	< LOQ		mg/17g	0.524	
Total Cannabinoids per 17g	43.1		mg/17g		

Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2407387	09/27/24 AOAC 991.14 (Petrifilm)		
Total Coliforms	< LOQ		cfu/g	10	2407387	09/27/24 AOAC 991.14 (Petrifilm)		
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2407388	09/28/24 AOAC 2014.05 (RAPID)		
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2407388	09/28/24 AOAC 2014.05 (RAPID)		



Solvents						Method: Residual Solvents by HS-GC-MS ^b		Units µg/g	Batch 2407455	Analyze 09/26/24 11:19 AM				
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes			
1,4-Dioxane [⊥]	< LOQ		100			2-Butanol [⊥]	< LOQ		200					
2-Ethoxyethanol [⊥]	< LOQ		30.0			2-Methylbutane (Isopentane) [⊥]	< LOQ		200					
2-Methylpentane [⊥]	< LOQ		30.0			2-Propanol (IPA) [⊥]	< LOQ		200					
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		200			Acetonitrile [⊥]	< LOQ		100					
Benzene [⊥]	< LOQ		1.00			Butanes (sum) [⊥]	< LOQ		400					
Cyclohexane [⊥]	< LOQ		200			Ethyl acetate [⊥]	< LOQ		200					
Ethyl benzene	< LOQ		200			Ethyl ether [⊥]	< LOQ		200					
Ethylene glycol [⊥]	< LOQ		200			Ethylene oxide [⊥]	< LOQ		20.0					
Hexanes (sum) [⊥]	< LOQ		150			Isopropyl acetate [⊥]	< LOQ		200					
Isopropylbenzene (Cumene) [⊥]	< LOQ		30.0			m,p-Xylene [⊥]	< LOQ		200					
Methanol [⊥]	< LOQ		200			Methylene chloride [⊥]	< LOQ		60.0					
Methylpropane (Isobutane) [⊥]	< LOQ		200			n-Butane [⊥]	< LOQ		200					
n-Heptane [⊥]	< LOQ		200			n-Hexane [⊥]	< LOQ		30.0					
n-Pentane [⊥]	< LOQ		200			o-Xylene [⊥]	< LOQ		200					
Pentanes (sum)	< LOQ		600			Propane	< LOQ		200					
Tetrahydrofuran [⊥]	< LOQ		100			Toluene [⊥]	< LOQ		100					
Total Xylenes [⊥]	< LOQ		400			Total Xylenes and Ethyl benzene	< LOQ		600					



Pesticides						Method: AOAC 2007.01 & EN 15662 (mod)						Units mg/kg		Batch 2407457		Analyze 09/26/24 11:38 AM					
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ
Abamectin [±]	< LOQ		0.250			Acephate	< LOQ		0.200												
Acequinocyl [±]	< LOQ		1.00			Acetamiprid	< LOQ		0.100												
Aldicarb [±]	< LOQ		0.200			Azoxystrobin [±]	< LOQ		0.100												
Bifenazate [±]	< LOQ		0.100			Bifenthrin [±]	< LOQ		0.100												
Boscalid [±]	< LOQ		0.200			Carbaryl [±]	< LOQ		0.100												
Carbofuran [±]	< LOQ		0.100			Chlorantraniliprole [±]	< LOQ		0.100												
Chlorfenapyr [±]	< LOQ		0.500			Chlorpyrifos-ethyl [±]	< LOQ		0.100												
Clofentezine [±]	< LOQ		0.100			Cyfluthrin (sum) [±]	< LOQ		0.500												
Cypermethrin and	< LOQ		0.500			Daminozide [±]	< LOQ		0.500												
Diazinon [±]	< LOQ		0.100			Dichlorvos [±]	< LOQ		0.500												
Dimethoate [±]	< LOQ		0.100			Ethoprophos [±]	< LOQ		0.100												
Etofenprox [±]	< LOQ		0.200			Etoxazole [±]	< LOQ		0.100												
Fenoxycarb [±]	< LOQ		0.100			Fenpyroximate [±]	< LOQ		0.200												
Fipronil [±]	< LOQ		0.200			Flonicamid [±]	< LOQ		0.400												
Fludioxonil [±]	< LOQ		0.200			Hexythiazox [±]	< LOQ		0.400												
Imazalil [±]	< LOQ		0.100			Imidacloprid [±]	< LOQ		0.200												
Kresoxim-methyl [±]	< LOQ		0.200			Malathion [±]	< LOQ		0.100												
Metalaxyl [±]	< LOQ		0.100			Methiocarb [±]	< LOQ		0.100												
Methomyl [±]	< LOQ		0.200			MGK-264 [±]	< LOQ		0.100												
Myclobutanil [±]	< LOQ		0.100			Naled [±]	< LOQ		0.250												
Oxamyl [±]	< LOQ		0.500			Paclobutrazole [±]	< LOQ		0.200												
Parathion-methyl [±]	< LOQ		0.100			Permethrin [±]	< LOQ		0.100												
Phosmet [±]	< LOQ		0.100			Piperonyl butoxide [±]	< LOQ		1.00												
Prallethrin [±]	< LOQ		0.100			Propiconazole [±]	< LOQ		0.200												
Propoxur [±]	< LOQ		0.100			Pyrethrin I (total) [±]	< LOQ		0.500												
Pyridaben [±]	< LOQ		0.100			Spinosad [±]	< LOQ		0.100												
Spiromesifen [±]	< LOQ		0.100			Spirotetramat [±]	< LOQ		0.100												
Spiroxamine [±]	< LOQ		0.200			Tebuconazole [±]	< LOQ		0.200												
Thiacloprid [±]	< LOQ		0.100			Thiamethoxam [±]	< LOQ		0.100												
Trifloxystrobin [±]	< LOQ		0.100																		

Metals											
Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method		Status Notes			
Arsenic [±]	< LOQ		mg/kg	0.0169	2407535	09/27/24	AOAC 2013.06 (mod.) ^b				
Cadmium [±]	< LOQ		mg/kg	0.0169	2407535	09/27/24	AOAC 2013.06 (mod.) ^b				
Lead [±]	< LOQ		mg/kg	0.0169	2407535	09/27/24	AOAC 2013.06 (mod.) ^b				
Mercury [±]	< LOQ		mg/kg	0.00844	2407535	09/27/24	AOAC 2013.06 (mod.) ^b				



Mycotoxins

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
Aflatoxin B1 [⊥]	< LOQ		µg/kg	5.00	2407532	09/30/24 AOAC 2007.01 & EN 15662 (mod)		
Aflatoxin B2 [⊥]	< LOQ		µg/kg	5.00	2407532	09/30/24 AOAC 2007.01 & EN 15662 (mod)		
Aflatoxin G1 [⊥]	< LOQ		µg/kg	5.00	2407532	09/30/24 AOAC 2007.01 & EN 15662 (mod)		
Aflatoxin G2 [⊥]	< LOQ		µg/kg	5.00	2407532	09/30/24 AOAC 2007.01 & EN 15662 (mod)		
Ochratoxin A [⊥]	< LOQ		µg/kg	5.00	2407532	09/30/24 AOAC 2007.01 & EN 15662 (mod)		
Total Aflatoxins	< LOQ		µg/kg	20.0		10/01/24 AOAC 2007.01 & EN 15662 (mod) ^p		



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



Report Number: 24-010742/D007.R001
Report Date: 10/02/2024
ORELAP#: OR100028
Purchase Order:
Received: 09/24/24 11:17

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 = Gram

µg/g = Microgram per gram

µg/kg = Micrograms per kilogram = parts per billion (ppb)

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

mg/17g = Milligram per 17g

% = Percentage of sample

% wt = µg/g divided by 10,000

Approved Signatory

Derrick Tanner
General Manager



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



Report Number: 24-010742/D007.R001
Report Date: 10/02/2024
ORELAP#: OR100028
Purchase Order:
Received: 09/24/24 11:17



Hemp & Cannabis
Chain of Custody

Seventh-Hill-CBD-
1727115653

							Testing	
Company Details Company: <u>Seventh Hill CBD</u> Contact: <u>Jordan Dunn</u> Street Address: <u>215 South 19th Street</u> City, State, Zip: <u>Springfield, OR 97477</u> Email: <u>jordandunn@seventhhillcbd.com</u> Contact Phone: <u>5415912620</u> Company Phone: <u>5415912620</u> Billing Information Billing Phone: <u>5415912620</u> Billing Email: <u>jordandunn@seventhhillcbd.com</u>				Project Details Turnaround Time: <u>3 Business Days</u> <u>Surcharges Apply</u> Relinquishment Sampling, Courier & Shipping Options: <u>By Shipping Service (USPS, UPS, Fedex)</u> Receipt Information Prelog Storage: <u>Canna Shelves</u> Sample Condition: <u>Satisfactory</u>			CH001 - Standard Potency + Safety Package H0014 - Potency Cannabis (Basic)	
#	Sample Name	Material	METRC ID	Amount Provided	Reporting Unit	Specifications		
1	3-1 Caramel	Cannabinoid Edible	CC3124251	30 g	% & mg/serving	17 Grams per Serving	✓	
2	6-1 Caramel	Cannabinoid Edible	CC6124251	30 g	% & mg/serving	17 grams per serving	✓	
3	3-1 THCV Caramel	Cannabinoid Edible	CC31TV2425	30 g	% & mg/serving	17 Grams per serving	✓	
4	6-1-1 Gummy Caramel	Cannabinoid Edible	GC6112425	30 g	% & mg/serving	17 Grams per serving	✓	
5	1-1 Pumpkin Caramel	Cannabinoid Edible	CC11PS2425	30 g	% & mg/serving	17 Grams per serving		✓
6	3-1 Fudge	Cannabinoid Edible	FF3124251	30 g	% & mg/serving	12.5 Grams per Serving	✓	
7	Focus CBG Caramel	Cannabinoid Edible	FCBG624251	30 g	% & mg/serving	17 grams	✓	
8	1-1 Caramel	Cannabinoid Edible	CC1124251	30 g	% & mg/serving	17 Grams per Serving	✓	

Package Details

Standard Potency + Safety Package: Cannabis Heavy Metals Profile OR • E. coli/Coliform Count (EC) Petrifilm • Pesticides (OR - Cannabis) • Potency Cannabis (Basic+Expanded) • RAPID Yeast and Mold Count (RYM) Petrifilm • Residual Solvents (Cannabis - Oregon)

Relinquished By	Date	Time	Received By	Date	Time	Received Temp., °C	Evidence of Cooling?
<i>Jordan Taylor Dunn</i>	<i>09/23/2024</i>	<i>11:20</i>	<i>edm</i>	<i>09/24/2024</i>	<i>11:17</i>	<i>23.00</i>	<i>No</i>

Samples submitted to Columbia Laboratories with testing requirements constitute an agreement for services in accordance with the [current terms of services](#) associated with this COC. By signing "Relinquished by" you are agreeing to these terms.

Columbia Laboratories
12423 NE Whitaker Way
Portland, OR 97230

P: (503) 254-1794
info@columbialaboratories.com

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www.columbialaboratories.com


Laboratory Quality Control Results
J AOAC 2015 V98-6 **Batch ID: 2407447**
Laboratory Control Sample

Analyte	LCS	Result	Spike	Units	% Rec	Limits	Evaluation	Notes
CBDVA	2	0.0313	0.0309	%	101	80.0 - 120	Acceptable	
CBDV	2	0.0280	0.0289	%	96.8	80.0 - 120	Acceptable	
CBE	2	0.0350	0.0352	%	99.6	80.0 - 120	Acceptable	
CBDA	1	0.0361	0.0358	%	101	90.0 - 110	Acceptable	
CBGA	1	0.0355	0.0351	%	101	80.0 - 120	Acceptable	
CBG	1	0.0318	0.0314	%	101	80.0 - 120	Acceptable	
CBD	1	0.0350	0.0351	%	99.8	90.0 - 110	Acceptable	
THCV	2	0.0335	0.0338	%	99.1	80.0 - 120	Acceptable	
d8THCV	2	0.0349	0.0348	%	100	80.0 - 120	Acceptable	
THCVA	2	0.0306	0.0308	%	99.3	80.0 - 120	Acceptable	
CBN	1	0.0354	0.0352	%	101	80.0 - 120	Acceptable	
exo-THC	2	0.0317	0.0326	%	97.3	80.0 - 120	Acceptable	
d9THC	1	0.0333	0.0337	%	98.8	90.0 - 110	Acceptable	
d8THC	1	0.0278	0.0284	%	97.9	90.0 - 110	Acceptable	
9S-d10THC	1	0.0344	0.0354	%	97.2	80.0 - 120	Acceptable	
CBL	2	0.0300	0.0307	%	97.5	80.0 - 120	Acceptable	
9R-d10THC	1	0.0352	0.0365	%	96.5	80.0 - 120	Acceptable	
CBC	2	0.0336	0.0347	%	96.9	80.0 - 120	Acceptable	
THCA	1	0.0366	0.0359	%	102	90.0 - 110	Acceptable	
CBCA	2	0.0311	0.0321	%	96.8	80.0 - 120	Acceptable	
CBLA	2	0.0311	0.0325	%	95.7	80.0 - 120	Acceptable	
d9THCP	2	0.0305	0.0322	%	94.7	80.0 - 120	Acceptable	
CBT	2	0.0307	0.0334	%	91.7	80.0 - 120	Acceptable	

Method Blank

Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBDV	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBE	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBDA	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBGA	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBG	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBD	<LOQ	0.00327	%	< 0.00327	Acceptable	
THCV	<LOQ	0.00327	%	< 0.00327	Acceptable	
d8THCV	<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	
d9THC	<LOQ	0.00327	%	< 0.00327	Acceptable	
d8THC	<LOQ	0.00327	%	< 0.00327	Acceptable	
9S-d10THC	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBL	<LOQ	0.00327	%	< 0.00327	Acceptable	
9R-d10THC	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBC	<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	
d9THCP	<LOQ	0.00327	%	< 0.00327	Acceptable	
CBT	<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


Laboratory Quality Control Results

J AOAC 2015 V98-6			Batch ID: 2407447					
Sample Duplicate			Sample ID: 24-010717-0001					
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
CBDV	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
CBE	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
CBDA	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
CBG	0.227	0.226	0.00309	%	0.337	< 20	Acceptable	
CBD	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
CBN	0.00320	0.00315	0.00309	%	1.65	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
d9THC	0.257	0.256	0.00309	%	0.498	< 20	Acceptable	
d8THC	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
9S-d10THC	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
9R-d10THC	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
CBC	0.257	0.256	0.00309	%	0.243	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
CBCA	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
d9THCP	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	
CBT	<LOQ	<LOQ	0.00309	%	NA	< 20	Acceptable	

Abbreviations

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

Units of Measure:

% - Percent


Laboratory Quality Control Results

Residual Solvents					Batch ID: 2407455				
Method Blank					Laboratory Control Sample				
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		435	585	µg/g	74.4	60 - 120	
Isobutane	ND	< 200		559	770	µg/g	72.6	60 - 120	
Butane	ND	< 200		570	769	µg/g	74.1	60 - 120	
2,2-Dimethylpropane	ND	< 200		706	956	µg/g	73.8	60 - 120	
Methanol	ND	< 200		1320	1630	µg/g	81.0	60 - 120	
Ethylene Oxide	ND	< 30		43.9	57.7	µg/g	76.1	60 - 120	
2-Methylbutane	ND	< 200		1270	1620	µg/g	78.4	60 - 120	
Pentane	ND	< 200		1250	1620	µg/g	77.2	60 - 120	
Ethanol	ND	< 200		1350	1620	µg/g	83.3	70 - 130	
Ethyl Ether	ND	< 200		1480	1620	µg/g	91.4	60 - 120	
2,2-Dimethylbutane	ND	< 30		144	179	µg/g	80.4	60 - 120	
Acetone	ND	< 200		1300	1620	µg/g	80.2	60 - 120	
2-Propanol	ND	< 200		1310	1620	µg/g	80.9	60 - 120	
Ethyl Formate	ND	< 500		1170	1610	µg/g	72.7	70 - 130	
Acetonitrile	ND	< 100		391	502	µg/g	77.9	60 - 120	
Methyl Acetate	ND	< 500		1390	1610	µg/g	86.3	70 - 130	
2,3-Dimethylbutane	ND	< 30		147	180	µg/g	81.7	60 - 120	
Dichloromethane	ND	< 60		434	533	µg/g	81.4	60 - 120	
2-Methylpentane	ND	< 30		151	181	µg/g	83.4	60 - 120	
MTBE	ND	< 500		1330	1600	µg/g	83.1	70 - 130	
3-Methylpentane	ND	< 30		148	177	µg/g	83.6	60 - 120	
Hexane	ND	< 30		147	182	µg/g	80.8	60 - 120	
1-Propanol	ND	< 500		1310	1610	µg/g	81.4	70 - 130	
Methylethylketone	ND	< 500		1350	1600	µg/g	84.4	70 - 130	
Ethyl acetate	ND	< 200		1330	1620	µg/g	82.1	60 - 120	
2-Butanol	ND	< 200		1320	1630	µg/g	81.0	60 - 120	
Tetrahydrofuran	ND	< 100		411	499	µg/g	82.4	60 - 120	
Cyclohexane	ND	< 200		1280	1610	µg/g	79.5	60 - 120	
2-methyl-1-propanol	ND	< 500		1330	1600	µg/g	83.1	70 - 130	
Benzene	ND	< 1		3.96	5.01	µg/g	79.0	60 - 120	
Isopropyl Acetate	ND	< 200		1310	1620	µg/g	80.9	60 - 120	
Heptane	ND	< 200		1270	1610	µg/g	78.9	60 - 120	
1-Butanol	ND	< 500		1390	1600	µg/g	86.9	70 - 130	
Propyl Acetate	ND	< 500		1410	1600	µg/g	88.1	70 - 130	
1,4-Dioxane	ND	< 100		398	493	µg/g	80.7	60 - 120	
2-Ethoxyethanol	ND	< 30		142	182	µg/g	78.0	60 - 120	
Methylisobutylketone	ND	< 500		1390	1610	µg/g	86.3	70 - 130	
3-Methyl-1-butanol	ND	< 500		1370	1600	µg/g	85.6	70 - 130	
Ethylene Glycol	ND	< 200		322	501	µg/g	64.3	60 - 120	
Toluene	ND	< 100		400	501	µg/g	79.8	60 - 120	
Isobutyl Acetate	ND	< 500		1320	1600	µg/g	82.5	70 - 130	
1-Pentanol	ND	< 500		1320	1600	µg/g	82.5	70 - 130	
Butyl Acetate	ND	< 500		1340	1600	µg/g	83.8	70 - 130	
Ethylbenzene	ND	< 200		756	981	µg/g	77.1	60 - 120	
m,p-Xylene	ND	< 200		728	1000	µg/g	72.8	60 - 120	
o-Xylene	ND	< 200		834	981	µg/g	85.0	60 - 120	
Cumene	ND	< 30		153	177	µg/g	86.4	60 - 120	
Anisole	ND	< 500		1550	1610	µg/g	96.3	70 - 130	
DMSO	ND	< 500		1360	1600	µg/g	85.0	70 - 130	
1,2-dimethoxyethane	ND	< 50		127	161	µg/g	78.9	70 - 130	
Triethylamine	ND	< 500		1310	1600	µg/g	81.9	70 - 130	
N,N-dimethylformamide	ND	< 150		423	484	µg/g	87.4	70 - 130	
N,N-dimethylacetamide	ND	< 150		507	497	µg/g	102.0	70 - 130	
Pyridine	ND	< 50		124	162	µg/g	76.5	70 - 130	
Sulfolane	ND	< 50		128	166	µg/g	77.1	70 - 130	
1,2-Dichloroethane	ND	< 1		0.864	1	µg/g	86.4	70 - 130	
Chloroform	ND	< 1		0.84	1	µg/g	84.0	70 - 130	
Trichloroethylene	ND	< 1		0.841	1	µg/g	84.1	70 - 130	
1,1-Dichloroethane	ND	< 1		0.858	1	µg/g	85.8	70 - 130	


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

QC - Sample Duplicate
Sample ID: 24-010616-0001

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	370	404	200	µg/g	8.8	< 20	Acceptable	
2-Butanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Tetrahydrofuran	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Cyclohexane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2-methyl-1-propanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Benzene	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Isopropyl Acetate	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Heptane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
1-Butanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Propyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
1,4-Dioxane	ND	ND	100	µg/g	0.0	< 20	Acceptable	
2-Ethoxyethanol	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Methylisobutylketone	ND	ND	500	µg/g	0.0	< 20	Acceptable	
3-Methyl-1-butanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Ethylene Glycol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Toluene	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Isobutyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
1-Pentanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Butyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Ethylbenzene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
m,p-Xylene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
o-Xylene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Cumene	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Anisole	ND	ND	500	µg/g	0.0	< 20	Acceptable	
DMSO	ND	ND	500	µg/g	0.0	< 20	Acceptable	
1,2-dimethoxyethane	ND	ND	50	µg/g	0.0	< 20	Acceptable	
Triethylamine	ND	ND	500	µg/g	0.0	< 20	Acceptable	
N,N-dimethylformamide	ND	ND	150	µg/g	0.0	< 20	Acceptable	
N,N-dimethylacetamide	ND	ND	150	µg/g	0.0	< 20	Acceptable	
Pyridine	ND	ND	50	µg/g	0.0	< 20	Acceptable	
Sulfolane	ND	ND	50	µg/g	0.0	< 20	Acceptable	
1,2-Dichloroethane	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Chloroform	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Trichloroethylene	ND	ND	1	µg/g	0.0	< 20	Acceptable	
1,1-Dichloroethane	ND	ND	1	µg/g	0.0	< 20	Acceptable	

Abbreviations

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

Units of Measure:

µg/g- Microgram per gram or ppm



Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662			Units: mg/Kg		Batch ID: 2407457			
Method Blank			Laboratory Control Sample					
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes
Abamectin	0.000	< 0.250		0.981	1.000	98.1	50.0 150	
Acephate	0.011	< 0.200		0.685	0.800	85.6	60.0 120	
Acequinocyl	0.000	< 1.000		3.338	4.000	83.4	40.0 160	
Acetamiprid	0.011	< 0.100		0.370	0.400	92.6	60.0 120	
Aldicarb	0.000	< 0.200		0.716	0.800	89.5	60.0 120	
Azoxystrobin	0.015	< 0.100		0.352	0.400	87.9	60.0 120	
Bifenazate	0.000	< 0.100		0.380	0.400	95.0	60.0 120	
Bifenthrin	0.019	< 0.100		0.387	0.400	96.7	50.0 150	
Boscalid	0.020	< 0.200		0.734	0.800	91.8	60.0 120	
Carbaryl	0.011	< 0.100		0.356	0.400	88.9	60.0 120	
Carbofuran	0.010	< 0.100		0.359	0.400	89.9	60.0 120	
Chlorantraniliprole	0.009	< 0.100		0.347	0.400	86.7	60.0 120	
Chlorfenapyr	0.000	< 0.500		1.856	2.000	92.8	60.0 120	
Chlorpyrifos	0.003	< 0.100		0.368	0.400	92.0	60.0 120	
Clofentezine	0.014	< 0.100		0.287	0.400	71.8	60.0 120	
Cyfluthrin	0.047	< 0.500		1.904	2.000	95.2	50.0 150	
Cypermethrin	0.074	< 0.500		1.869	2.000	93.4	50.0 150	
Daminozide	0.072	< 0.500		0.760	2.000	38.0	60.0 120	Q7
Diazinon	0.007	< 0.100		0.369	0.400	92.3	60.0 120	
Dichlorvos	0.060	< 0.500		1.695	2.000	84.7	60.0 120	
Dimethoate	0.006	< 0.100		0.364	0.400	90.9	60.0 120	
Ethoprophos	0.005	< 0.100		0.363	0.400	90.6	60.0 120	
Etofenprox	0.052	< 0.200		0.805	0.800	100.6	50.0 150	
Etoxazole	0.004	< 0.100		0.403	0.400	100.7	60.0 120	
Fenoxycarb	0.009	< 0.100		0.353	0.400	88.2	60.0 120	
Fenpyroximate	0.016	< 0.200		0.744	0.800	93.0	60.0 120	
Fipronil	0.034	< 0.200		0.714	0.800	89.2	60.0 120	
Flonicamid	0.000	< 0.250		0.871	1.000	87.1	60.0 120	
Fludioxonil	0.001	< 0.200		0.745	0.800	93.1	50.0 150	
Hexythiazox	0.008	< 0.250		0.941	1.000	94.1	60.0 120	
Imazalil	0.008	< 0.100		0.371	0.400	92.8	60.0 120	
Imidacloprid	0.045	< 0.200		0.663	0.800	82.8	60.0 120	
Kresoxim-methyl	0.021	< 0.200		0.687	0.800	85.9	60.0 120	
Malathion	0.000	< 0.100		0.375	0.400	93.9	60.0 120	
Metalaxyl	0.000	< 0.100		0.344	0.400	86.0	60.0 120	
Methiocarb	0.009	< 0.100		0.355	0.400	88.9	60.0 120	
Methomyl	0.000	< 0.200		0.763	0.800	95.3	60.0 120	
MGK-264	0.000	< 0.100		0.364	0.400	91.1	50.0 150	
Myclobutanil	0.022	< 0.100		0.350	0.400	87.4	60.0 120	
Naled	0.000	< 0.250		0.843	1.000	84.3	50.0 150	
Oxamyl	0.000	< 0.500		1.862	2.000	93.1	60.0 120	
Paclobutrazole	0.021	< 0.200		0.690	0.800	86.3	60.0 120	
Parathion-Methyl	0.000	< 0.100		0.350	0.400	87.6	50.0 150	
Permethrin	0.013	< 0.100		0.405	0.400	101.2	50.0 150	
Phosmet	0.016	< 0.100		0.345	0.400	86.3	50.0 150	
Piperonyl butoxide	0.021	< 0.500		2.030	2.000	101.5	60.0 120	
Prallethrin	0.000	< 0.100		0.351	0.400	87.7	60.0 120	
Propiconazole	0.011	< 0.200		0.724	0.800	90.5	60.0 120	
Propoxur	0.006	< 0.100		0.356	0.400	89.1	60.0 120	
Pyrethrin (Summe)	0.000	< 0.100		0.231	0.488	47.2	60.0 120	84.3
Pyridaben	0.007	< 0.100		0.402	0.400	100.6	50.0 150	
Spinosad	0.000	< 0.100		0.350	0.388	90.1	50.0 150	
Spiromesifen	0.000	< 0.100		0.386	0.400	96.6	60.0 120	
Spirotetramat	0.006	< 0.100		0.374	0.400	93.6	60.0 120	



Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662				Units: mg/Kg			Batch ID: 2407457				
Matrix Spike/Matrix Spike Duplicate Recoveries				Sample ID: 24-010641-0001							
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes	
Abamectin	0.000	0.849	0.906	1.000	6.5%	< 30	84.9%	90.6%	50 - 150	R,Q	
Acephate	0.000	0.656	0.645	0.800	1.8%	< 30	82.1%	80.6%	50 - 150		
Acequinocyl	0.000	3.375	0.393	4.000	158.3%	< 30	84.4%	9.8%	50 - 150		
Acetamiprid	0.000	0.356	0.359	0.400	0.7%	< 30	89.0%	89.7%	50 - 150		
Aldicarb	0.000	0.640	0.649	0.800	1.4%	< 30	80.0%	81.1%	50 - 150	R,Q	
Azoxystrobin	0.000	0.329	0.319	0.400	3.4%	< 30	82.4%	79.6%	50 - 150		
Bifenazate	0.000	0.346	0.336	0.400	2.9%	< 30	86.5%	84.0%	50 - 150		
Bifenthrin	0.000	0.292	0.190	0.400	42.3%	< 30	73.1%	47.6%	50 - 150		
Boscalid	0.000	0.643	0.662	0.800	2.9%	< 30	80.4%	82.7%	50 - 150	R,Q	
Carbaryl	0.000	0.320	0.324	0.400	1.4%	< 30	79.9%	81.0%	50 - 150		
Carbofuran	0.000	0.327	0.334	0.400	2.0%	< 30	81.8%	83.5%	50 - 150		
Chlorantraniliprole	0.000	0.324	0.335	0.400	3.3%	< 30	81.0%	83.7%	50 - 150		
Chlorfenapyr	0.000	1.340	1.555	2.000	14.9%	< 30	67.0%	77.8%	50 - 150	R,Q	
Chlorpyrifos	0.000	0.323	0.357	0.400	9.8%	< 30	80.8%	89.2%	50 - 150		
Clofentezine	0.000	0.321	0.304	0.400	5.5%	< 30	80.3%	76.0%	50 - 150		
Cyfluthrin	0.000	1.595	1.611	2.000	1.0%	< 30	79.7%	80.6%	30 - 150		
Cypermethrin	0.000	1.146	1.066	2.000	7.2%	< 30	57.3%	53.3%	50 - 150	R,Q	
Daminozide	0.000	0.728	0.777	2.000	6.5%	< 30	36.4%	38.9%	30 - 150		
Diazinon	0.000	0.341	0.343	0.400	0.7%	< 30	85.1%	85.7%	50 - 150		
Dichlorvos	0.000	1.675	1.694	2.000	1.1%	< 30	83.7%	84.7%	50 - 150		
Dimethoate	0.000	0.347	0.360	0.400	3.7%	< 30	86.8%	90.0%	50 - 150	R,Q	
Ethoprophos	0.000	0.328	0.343	0.400	4.5%	< 30	81.9%	85.7%	50 - 150		
Etofenprox	0.000	0.519	0.372	0.800	33.0%	< 30	64.9%	46.5%	50 - 150		
Etoxazole	0.000	0.354	0.345	0.400	2.4%	< 30	88.4%	86.2%	50 - 150		
Fenoxycarb	0.000	0.327	0.322	0.400	1.6%	< 30	81.8%	80.5%	50 - 150	R,Q	
Fenpyroximate	0.000	0.604	0.520	0.800	15.0%	< 30	75.5%	65.0%	50 - 150		
Fipronil	0.000	0.593	0.582	0.800	1.8%	< 30	74.1%	72.8%	50 - 150		
Flonicamid	0.000	0.797	0.877	1.000	9.5%	< 30	79.7%	87.7%	50 - 150		
Fludioxonil	0.000	0.711	0.742	0.800	4.3%	< 30	88.8%	92.7%	50 - 150	R,Q	
Hexythiazox	0.000	0.643	0.635	1.000	1.2%	< 30	64.3%	63.5%	50 - 150		
Imazalil	0.000	0.345	0.344	0.400	0.1%	< 30	86.2%	86.1%	50 - 150		
Imidacloprid	0.000	0.611	0.636	0.800	3.9%	< 30	76.4%	79.4%	50 - 150		
Kresoxim-methyl	0.000	0.628	0.662	0.800	5.3%	< 30	78.5%	82.7%	50 - 150	R,Q	
Malathion	0.000	0.347	0.368	0.400	5.8%	< 30	86.9%	92.1%	50 - 150		
Metalaxyl	0.000	0.329	0.312	0.400	5.2%	< 30	82.2%	78.0%	50 - 150		
Methiocarb	0.000	0.324	0.329	0.400	1.7%	< 30	80.9%	82.3%	50 - 150		
Methomyl	0.000	0.689	0.707	0.800	2.5%	< 30	86.1%	88.3%	50 - 150	R,Q	
MGK-264	0.000	0.306	0.297	0.400	2.7%	< 30	76.4%	74.4%	50 - 150		
Myclobutanil	0.000	0.325	0.305	0.400	6.5%	< 30	81.3%	76.2%	50 - 150		
Naled	0.000	0.746	0.726	1.000	2.7%	< 30	74.6%	72.6%	50 - 150		
Oxamyl	0.000	1.764	1.925	2.000	8.7%	< 30	88.2%	96.2%	50 - 150	R,Q	
Paclobutrazole	0.000	0.620	0.605	0.800	2.5%	< 30	77.5%	75.6%	50 - 150		
Parathion-Methyl	0.000	0.217	0.235	0.400	7.9%	< 30	54.3%	58.8%	30 - 150		
Permethrin	0.000	0.281	0.261	0.400	7.3%	< 30	70.1%	65.2%	50 - 150		
Phosmet	0.000	0.318	0.305	0.400	4.3%	< 30	79.5%	76.2%	50 - 150	R,Q	
Piperonyl butoxide	0.000	1.713	1.713	2.000	0.0%	< 30	85.7%	85.6%	50 - 150		
Prallethrin	0.000	0.283	0.275	0.400	2.9%	< 30	70.7%	68.7%	50 - 150		
Propiconazole	0.000	0.618	0.637	0.800	3.0%	< 30	77.3%	79.6%	50 - 150		
Propoxur	0.000	0.323	0.331	0.400	2.6%	< 30	80.6%	82.8%	50 - 150	R,Q	
Pyrethrin (Summe)	0.000	0.196	0.226	0.488	14.2%	< 30	71.5%	82.5%	50 - 150		
Pyridaben	0.000	0.252	0.242	0.400	3.9%	< 30	63.0%	60.6%	50 - 150		
Spinosad	0.000	0.312	0.302	0.388	3.4%	< 30	80.4%	77.8%	50 - 150		
Spiromesifen	0.000	0.232	0.218	0.400	6.4%	< 30	58.1%	54.5%	50 - 150	R,Q	
Spiratetramat	0.000	0.337	0.328	0.400	2.8%	< 30	84.2%	81.9%	50 - 150		
Spiratetramat	0.000	0.337	0.328	0.400	2.8%	< 30	84.2%	81.9%	50 - 150		
Spiratetramat	0.000	0.337	0.328	0.400	2.8%	< 30	84.2%	81.9%	50 - 150		



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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.