

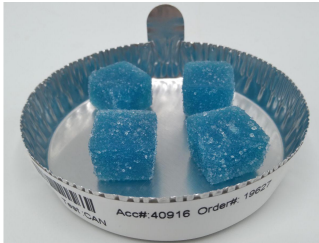
Bayou City Hemp Company

16700 Park Row
Houston, TX 77084
shorton@bayoucityhemp.com
281-928-9181

Sample: 10-31-2023-40916

Sample Received: 10/31/2023;
Report Created: 11/01/2023; Expires: 10/31/2024

G08Wb31023
Ingestible, Soft Chew



0.026 %

Total THC

0.026 %

Δ-9 THC

22.627 mg/unit
Total Cannabinoids

ND mg/unit
Total CBD

Cannabinoids

(Testing Method: HPLC, CON-P-3000)
Date Tested: 10/31/2023

Complete

Analyte	LOD	LOQ	Mass	Mass	Mass	
	mg/unit	mg/unit	mg/unit	mg/g	%	
Δ-8-Tetrahydrocannabinol (Δ-8 THC)	0.408	0.608	21.572	5.397	0.540	<div style="width: 5.397%;"></div>
Δ-9-Tetrahydrocannabinol (Δ-9 THC)	0.408	0.608	1.055	0.264	0.026	<div style="width: 0.264%;"></div>
Δ-9-Tetrahydrocannabinolic Acid (THCA-A)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
Δ-9-Tetrahydrocannabiphorol (Δ-9-THCP)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
Δ-9-Tetrahydrocannabivarin (Δ-9-THCV)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
Δ-9-Tetrahydrocannabivarinic Acid (Δ-9-THCVA)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
R-Δ-10-Tetrahydrocannabinol (R-Δ-10-THC)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
S-Δ-10-Tetrahydrocannabinol (S-Δ-10-THC)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
9R-Hexahydrocannabinol (9R-HHC)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
9S-Hexahydrocannabinol (9S-HHC)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
Tetrahydrocannabinol Acetate (THCO)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
Cannabidivarin (CBDV)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
Cannabidivarinic Acid (CBDVA)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
Cannabidiol (CBD)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
Cannabidiolic Acid (CBDA)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
Cannabigerol (CBG)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
Cannabigerolic Acid (CBGA)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
Cannabinol (CBN)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
Cannabinolic Acid (CBNA)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
Cannabichromene (CBC)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
Cannabichromenic Acid (CBCA)	0.408	0.608	ND	ND	ND	<div style="width: 0%;"></div>
Total			22.627	5.661	0.566	

Total THC = THCa * 0.877 + Δ9-THC; Total CBD = CBDA * 0.877 + CBD; LOQ = Limit of Quantitation; ND = Not Detected.

Total THC Measurement of Uncertainty: ± 0.050%
Total CBD Measurement of Uncertainty: ± 2.000%
THCO potency analysis does not designate quantitative specificity of Δ-8-THCO and Δ-9-THCO isomers

Unit Size: 3.997 g Unit: 1 Gummy



New Bloom Labs
6121 Heritage Park Drive, A500
Chattanooga, TN 37416
(844) 837-8223
TN DEA#: RN0563975
ANAB Testing Laboratory (AT-2868): ISO/IEC
17025:2017

Natalie Siracusa
Laboratory Director

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



Report Number: 23-009673/D002.R000
Report Date: 08/21/2023
ORELAP#: OR100028
Purchase Order:
Received: 08/15/23 11:03

Customer: Bayou City Hemp Company
Product identity: CD26COLA-D8-CH11
Client/Metric ID: .
Laboratory ID: 23-009673-0013

Summary

Potency:

Analyte	Result (%)	<ul style="list-style-type: none"> ● 8-THC ● 8-THCV ● CBT ● CBD 	CBD-Total	<LOQ	
Δ8-THC	80.5		(Reported in percent of total sample)	THC-Total	<LOQ
Δ8-THCV	0.195				
CBT	0.0839				
CBD	0.0701				

Residual Solvents:

All analytes passing and less than LOQ.

Pesticides:

All analytes passing and less than LOQ.

Metals:

Less than LOQ for all analytes.

Microbiology:

Less than LOQ for all analytes.



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 16700 Park Row
 Houston Texas 77084
 United States of America (USA)

Product identity: CD26COLA-D8-CH11

Client/Metric ID: .

Sample Date:

Laboratory ID: 23-009673-0013

Evidence of Cooling: No

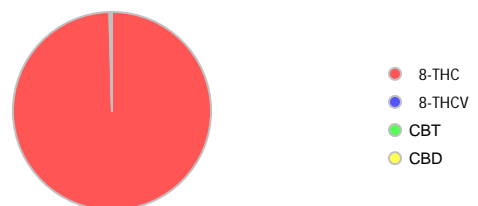
Temp: 27.2 °C

Relinquished by: client

Sample Results

Potency **Method:** J AOAC 2015 V98-6 (mod)^P **Units %** **Batch:** 2310117 **Analyze:** 8/16/23 11:34:00 PM

Analyte	As Received	Dry weight	LOQ	Notes
CBC	< LOQ		0.0676	
CBC	< LOQ		0.0676	
CBC-A	< LOQ		0.0676	
CBC-A	< LOQ		0.0676	
CBC-Total	< LOQ		0.127	
CBC-Total	< LOQ		0.127	
CBD	0.0701		0.0676	
CBD	0.0701		0.0676	
CBD-A	< LOQ		0.0676	
CBD-A	< LOQ		0.0676	
CBD-Total	< LOQ		0.127	
CBD-Total	< LOQ		0.127	
CBDV	< LOQ		0.0676	
CBDV	< LOQ		0.0676	
CBDV-A	< LOQ		0.0676	
CBDV-A	< LOQ		0.0676	
CBDV-Total	< LOQ		0.126	
CBDV-Total	< LOQ		0.126	
CBE	< LOQ		0.0676	
CBE	< LOQ		0.0676	
CBG	< LOQ		0.0676	
CBG	< LOQ		0.0676	
CBG-A	< LOQ		0.0676	
CBG-A	< LOQ		0.0676	
CBG-Total	< LOQ		0.126	
CBG-Total	< LOQ		0.126	
CBL	< LOQ		0.0676	
CBL	< LOQ		0.0676	
CBL-A	< LOQ		0.0676	
CBL-A	< LOQ		0.0676	
CBL-Total	< LOQ		0.127	





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Potency **Method:** J AOAC 2015 V98-6 (mod)^P **Units %** **Batch:** 2310117 **Analyze:** 8/16/23 11:34:00 PM

Analyte	As Received	Dry weight	LOQ	Notes
CBL-Total	< LOQ		0.127	
CBN	< LOQ		0.0676	
CBN	< LOQ		0.0676	
CBT	0.0839		0.0676	
CBT	0.0839		0.0676	
Δ10-THC-9R	< LOQ		0.0676	
Δ10-THC-9R	< LOQ		0.0676	
Δ10-THC-9S	< LOQ		0.0676	
Δ10-THC-9S	< LOQ		0.0676	
Δ10-THC-Total	< LOQ		0.135	
Δ10-THC-Total	< LOQ		0.135	
Δ8-THC	80.5		0.676	
Δ8-THC	80.5		0.676	
Δ8-THCV	0.195		0.0676	
Δ8-THCV	0.195		0.0676	
Δ9-THC	< LOQ		0.0676	
Δ9-THC	< LOQ		0.0676	
delta-9-THCP	< LOQ		0.0676	
delta-9-THCP	< LOQ		0.0676	
exo-THC	< LOQ		0.0676	
exo-THC	< LOQ		0.0676	
THC-A	< LOQ		0.0676	
THC-A	< LOQ		0.0676	
THC-Total	< LOQ		0.127	
THC-Total	< LOQ		0.127	
THCV	< LOQ		0.0676	
THCV	< LOQ		0.0676	
THCV-A	< LOQ		0.0676	
THCV-A	< LOQ		0.0676	
THCV-Total	< LOQ		0.126	
THCV-Total	< LOQ		0.126	
Total Cannabinoids	80.8			
Total Cannabinoids	80.8			

Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2310050	08/18/23 AOAC 991.14 (Petrifilm) ^P		
Total Coliforms	< LOQ		cfu/g	10	2310050	08/18/23 AOAC 991.14 (Petrifilm) ^P		
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2310051	08/18/23 AOAC 2014.05 (RAPID) ^P		
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2310051	08/18/23 AOAC 2014.05 (RAPID) ^P		



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Solvents											Method: Residual Solvents by GC/MS ^b					Units µg/g		Batch 2310111		Analyze 08/17/23 12:54 PM				
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		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) ^b											
Units mg/kg Batch 2310069 Analyze 08/16/23 01:11 PM											
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
Abamectin [‡]	< LOQ	0.50	0.250	pass		Acephate [‡]	< LOQ	0.40	0.200	pass	
Acequinocyl [‡]	< LOQ	2.0	1.00	pass		Acetamiprid [‡]	< LOQ	0.20	0.100	pass	
Aldicarb [‡]	< LOQ	0.40	0.200	pass		Azoxystrobin [‡]	< LOQ	0.20	0.100	pass	
Bifenazate [‡]	< LOQ	0.20	0.100	pass		Bifenthrin [‡]	< LOQ	0.20	0.100	pass	
Boscalid [‡]	< LOQ	0.40	0.200	pass		Carbaryl [‡]	< LOQ	0.20	0.100	pass	
Carbofuran [‡]	< LOQ	0.20	0.100	pass		Chlorantraniliprole [‡]	< LOQ	0.20	0.100	pass	
Chlorfenapyr [‡]	< LOQ	1.0	0.500	pass		Chlorpyrifos [‡]	< LOQ	0.20	0.100	pass	
Clofentezine [‡]	< LOQ	0.20	0.100	pass		Cyfluthrin [‡]	< LOQ	1.0	0.500	pass	
Cypermethrin [‡]	< LOQ	1.0	0.500	pass		Daminozide [‡]	< LOQ	1.0	0.500	pass	
Diazinon [‡]	< LOQ	0.20	0.100	pass		Dichlorvos [‡]	< LOQ	1.0	0.500	pass	
Dimethoate [‡]	< LOQ	0.20	0.100	pass		Ethoprophos [‡]	< LOQ	0.20	0.100	pass	
Etofenprox [‡]	< LOQ	0.40	0.200	pass		Etoxazole [‡]	< LOQ	0.20	0.100	pass	
Fenoxycarb [‡]	< LOQ	0.20	0.100	pass		Fenpyroximate [‡]	< LOQ	0.40	0.200	pass	
Fipronil [‡]	< LOQ	0.40	0.200	pass		Flonicamid [‡]	< LOQ	1.0	0.400	pass	
Fludioxonil [‡]	< LOQ	0.40	0.200	pass		Hexythiazox [‡]	< LOQ	1.0	0.400	pass	
Imazalil [‡]	< LOQ	0.20	0.100	pass		Imidacloprid [‡]	< LOQ	0.40	0.200	pass	
Kresoxim-methyl [‡]	< LOQ	0.40	0.200	pass		Malathion [‡]	< LOQ	0.20	0.100	pass	
Metalaxyl [‡]	< LOQ	0.20	0.100	pass		Methiocarb [‡]	< LOQ	0.20	0.100	pass	
Methomyl [‡]	< LOQ	0.40	0.200	pass		MGK-264 [‡]	< LOQ	0.20	0.100	pass	
Myclobutanil [‡]	< LOQ	0.20	0.100	pass		Naled [‡]	< LOQ	0.50	0.250	pass	
Oxamyl [‡]	< LOQ	1.0	0.500	pass		Pacllobutrazole [‡]	< LOQ	0.40	0.200	pass	
Parathion-Methyl [‡]	< LOQ	0.20	0.100	pass		Permethrin [‡]	< LOQ	0.20	0.100	pass	
Phosmet [‡]	< LOQ	0.20	0.100	pass		Piperonyl butoxide [‡]	< LOQ	2.0	1.00	pass	
Prallethrin [‡]	< LOQ	0.20	0.100	pass		Propiconazole [‡]	< LOQ	0.40	0.200	pass	
Propoxur [‡]	< LOQ	0.20	0.100	pass		Pyrethrin I (total) [‡]	< LOQ	1.0	0.500	pass	
Pyridaben [‡]	< LOQ	0.20	0.100	pass		Spinosad [‡]	< LOQ	0.20	0.100	pass	
Spiromesifen [‡]	< LOQ	0.20	0.100	pass		Spirotetramat [‡]	< LOQ	0.20	0.100	pass	
Spiroxamine [‡]	< LOQ	0.40	0.200	pass		Tebuconazole [‡]	< LOQ	0.40	0.200	pass	
Thiacloprid [‡]	< LOQ	0.20	0.100	pass		Thiamethoxam [‡]	< LOQ	0.20	0.100	pass	
Trifloxystrobin [‡]	< LOQ	0.20	0.100	pass							

Metals										
Analyte	Result	Limits	Units	LOQ	Batch	Analyzed	Method	Status	Notes	
Arsenic [‡]	< LOQ	0.200	mg/kg	0.0770	2310146	08/17/23	AOAC 2013.06 (mod.) ^b	pass		
Cadmium [‡]	< LOQ	0.200	mg/kg	0.0770	2310146	08/17/23	AOAC 2013.06 (mod.) ^b	pass		
Lead [‡]	< LOQ	0.500	mg/kg	0.0770	2310146	08/17/23	AOAC 2013.06 (mod.) ^b	pass		
Mercury [‡]	< LOQ	0.100	mg/kg	0.0385	2310146	08/17/23	AOAC 2013.06 (mod.) ^b	pass		



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Abbreviations

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

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

Ⓟ = ISO/IEC 17025:2017 accredited method.

Ⓜ = TNI accredited analyte.

Units of Measure

cfu/g = Colony forming units per gram

µg/g = Microgram per gram

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

% = Percentage of sample

% wt = µg/g divided by 10,000

Approved Signatory

Derrick Tanner
General Manager



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Revision: 3 Document ID: 3120
LegacyID: CFLC21WorksheetValidated 10/30/2020

Laboratory Pesticide Quality Control Results

AOAC2007.1 &EN 15662		Units: mg/Kg			Batch ID 2310069			
Method Blank	Laboratory Control Sample							
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spk	LCS % Re	Limits	Notes
Abamectin	0.00	< 0.250		0.922	1.00	92.2	50.0	150
Acaphate	0.031	< 0.200		0.702	0.80	87.8	60.0	120
Acquinocyl	0.00	< 1.000		3.482	4.00	87.1	40.0	160
Acetamiprid	0.00	< 0.100		0.372	0.40	93.0	60.0	120
Aldicarb	0.00	< 0.200		0.746	0.80	93.3	60.0	120
Azoxystrobin	0.01	< 0.100		0.374	0.40	93.6	60.0	120
Bifenazate	0.00	< 0.100		0.383	0.40	95.7	60.0	120
Bifenthrin	0.00	< 0.100		0.334	0.40	83.5	50.0	150
Boscalid	0.00	< 0.200		0.710	0.80	88.8	60.0	120
Carbaryl	0.00	< 0.100		0.378	0.40	94.5	60.0	120
Carbifuran	0.011	< 0.100		0.371	0.40	92.7	60.0	120
Chlorantraniliprole	0.005	< 0.100		0.355	0.40	89.1	60.0	120
Chlorfenapyr	0.00	< 0.500		1.889	2.00	94.5	60.0	120
Chlorpyrifos	0.005	< 0.100		0.411	0.40	102.8	60.0	120
Clofentezane	0.016	< 0.100		0.364	0.40	90.9	60.0	120
Cyfluthrin	0.00	< 0.500		2.233	2.00	111.7	50.0	150
Cypermethrin	0.00	< 0.500		1.799	2.00	89.9	50.0	150
Daminozide	0.249	< 0.500		2.016	2.00	100.8	60.0	120
Diazinon	0.00	< 0.100		0.370	0.40	92.4	60.0	120
Dichlorvos	0.093	< 0.500		1.999	2.00	99.9	60.0	120
Dimethoate	0.00	< 0.100		0.367	0.40	91.8	60.0	120
Ethiofoprofos	0.00	< 0.100		0.378	0.40	94.6	60.0	120
Etofenprox	0.00	< 0.200		0.734	0.80	91.7	50.0	150
Etoxazole	0.013	< 0.100		0.371	0.40	92.7	60.0	120
Fenoxycarb	0.001	< 0.100		0.380	0.40	95.0	60.0	120
Fenpyroximate	0.001	< 0.200		0.735	0.80	91.8	60.0	120
Fipronil	0.00	< 0.200		0.727	0.80	90.9	60.0	120
Fonicamid	0.00	< 0.250		0.893	1.00	89.3	60.0	120
Fludioxonil	0.00	< 0.200		0.795	0.80	99.4	50.0	150
Hexythiazox	0.032	< 0.250		0.929	1.00	92.9	60.0	120
Imazalil	0.019	< 0.100		0.371	0.40	92.8	60.0	120
Imidacloprid	0.037	< 0.200		0.692	0.80	86.5	60.0	120
Kiesoxim-methyl	0.00	< 0.200		0.789	0.80	98.7	60.0	120
Malathion	0.00	< 0.100		0.380	0.40	94.9	60.0	120
Metaxyl	0.00	< 0.100		0.374	0.40	93.4	60.0	120
Methiocarb	0.025	< 0.100		0.374	0.40	93.5	60.0	120
Methomyl	0.00	< 0.200		0.698	0.80	87.2	60.0	120
MCK-264	0.00	< 0.100		0.325	0.40	81.4	50.0	150
Mydobutani	0.028	< 0.100		0.370	0.40	92.6	60.0	120
Naled	0.00	< 0.250		0.920	1.00	92.0	50.0	150
Oxaryl	0.00	< 0.500		1.666	2.00	83.3	60.0	120
Padobutrazole	0.00	< 0.200		0.735	0.80	91.9	60.0	120
Parathion-Methyl	0.00	< 0.100		0.379	0.40	94.8	50.0	150
Permethrin	0.00	< 0.100		0.349	0.40	87.3	50.0	150
Phosmet	0.00	< 0.100		0.385	0.40	96.2	50.0	150
Piperonyl butoxide	0.008	< 0.500		1.882	2.00	94.1	60.0	120
Prallethrin	0.017	< 0.100		0.410	0.40	102.4	60.0	120
Propiconazole	0.00	< 0.200		0.711	0.80	88.9	60.0	120
Propoxur	0.010	< 0.100		0.372	0.40	93.1	60.0	120
Pyrethrin (Summe)	0.027	< 0.100		0.459	0.48	94.1	60.0	120
Pyridaben	0.00	< 0.100		0.372	0.40	93.0	50.0	150
Spinosad	0.00	< 0.100		0.357	0.38	92.0	50.0	150
Spiromesfen	0.00	< 0.100		0.379	0.40	94.6	60.0	120
Spirotetramat	0.00	< 0.100		0.386	0.40	96.5	60.0	120
Spiroxamine	0.00	< 0.200		0.759	0.80	94.9	60.0	120
Tebuconazole	0.00	< 0.200		0.754	0.80	94.2	60.0	120
Thiadoprid	0.00	< 0.100		0.383	0.40	90.7	60.0	120
Thiamethoxam	0.00	< 0.100		0.384	0.40	96.1	60.0	120
Trifloxystrobin	0.00	< 0.100		0.378	0.40	94.4	60.0	120



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 LegacyID: CFLC21WorksheetValidated 10/30/2020

Laboratory Pesticide Quality Control Results

AOAC2007.1 & EN 15662		Units: mg/Kg				Batch ID 2310069				
Matrix Spke/Matrix Spke Duplicate Recoveries	Result	MS Res	MSD Res	Spike	RFD%	Limit	MS% Re	MSD % Re	Limits	Notes
Abamectin	0.000	0.893	0.959	1.000	7.0%	< 30	89.3%	95.9%	50 - 150	
Acephate	0.000	0.526	0.643	0.800	20.0%	< 30	65.8%	80.4%	50 - 150	
Acetaminophyl	0.000	3.776	4.157	4.000	9.6%	< 30	94.4%	103.9%	50 - 150	
Acetamiprid	0.000	0.367	0.370	0.400	0.6%	< 30	91.9%	92.4%	50 - 150	
Aldicarb	0.000	0.754	0.762	0.800	1.1%	< 30	94.2%	95.2%	50 - 150	
Azoxystrobin	0.009	0.368	0.377	0.400	2.3%	< 30	89.9%	92.0%	50 - 150	
Bifenazate	0.000	0.377	0.390	0.400	3.5%	< 30	94.1%	97.9%	50 - 150	
Bifenthrin	0.014	0.383	0.392	0.400	2.4%	< 30	92.2%	94.9%	50 - 150	
Boscalid	0.073	0.783	0.699	0.800	12.9%	< 30	88.8%	78.3%	50 - 150	
Carbaryl	0.000	0.364	0.373	0.400	2.4%	< 30	91.0%	93.2%	50 - 150	
Carbofuran	0.009	0.354	0.359	0.400	1.6%	< 30	86.1%	87.5%	50 - 150	
Chlorantraniliprole	0.008	0.331	0.374	0.400	12.8%	< 30	80.7%	91.8%	50 - 150	
Chlorfenapyr	0.000	0.684	0.803	2.000	16.0%	< 30	34.2%	40.1%	50 - 150	Q
Chlorpyrifos	0.008	0.391	0.380	0.400	3.0%	< 30	95.8%	93.0%	50 - 150	
Clofentezane	0.014	0.325	0.351	0.400	7.8%	< 30	78.0%	84.3%	50 - 150	
Cyfluthrin	0.000	1.438	1.569	2.000	8.7%	< 30	71.9%	78.4%	30 - 150	
Cypermethrin	0.000	1.483	1.649	2.000	10.8%	< 30	74.1%	82.5%	50 - 150	
Daminozide	0.198	1.981	1.939	2.000	2.3%	< 30	89.1%	87.1%	30 - 150	
Diazinon	0.026	0.378	0.355	0.400	6.8%	< 30	88.1%	82.4%	50 - 150	
Dichlorvos	0.081	1.863	1.904	2.000	2.3%	< 30	89.1%	91.2%	50 - 150	
Dimethoate	0.019	0.394	0.353	0.400	8.9%	< 30	91.3%	83.9%	50 - 150	
Ethiofoprophos	0.001	0.357	0.382	0.400	6.9%	< 30	88.9%	95.3%	50 - 150	
Etofenprox	0.000	0.761	0.762	0.800	0.1%	< 30	95.1%	95.2%	50 - 150	
Etoxazole	0.008	0.327	0.332	0.400	1.5%	< 30	79.8%	81.0%	50 - 150	
Fenoxycarb	0.000	0.365	0.369	0.400	1.0%	< 30	91.3%	92.2%	50 - 150	
Fenpyroximate	0.000	0.564	0.616	0.800	8.7%	< 30	70.8%	77.0%	50 - 150	
Fipronil	0.000	0.655	0.662	0.800	1.0%	< 30	81.9%	82.7%	50 - 150	
Fonicamid	0.000	0.934	0.905	1.000	3.1%	< 30	93.4%	90.5%	50 - 150	
Fludioxonil	0.000	0.757	0.770	0.800	1.8%	< 30	94.8%	96.2%	50 - 150	
Hexythiazox	0.000	0.327	0.357	1.000	8.7%	< 30	32.7%	35.7%	50 - 150	Q
Imazalil	0.016	0.369	0.356	0.400	3.6%	< 30	88.1%	85.0%	50 - 150	
Imidacloprid	0.031	0.714	0.705	0.800	1.2%	< 30	85.4%	84.4%	50 - 150	
Kiesoxim-methyl	0.000	0.720	0.703	0.800	2.5%	< 30	90.0%	87.9%	50 - 150	
Malathion	0.000	0.361	0.359	0.400	0.6%	< 30	90.3%	89.8%	50 - 150	
Metaxyl	0.004	0.379	0.376	0.400	0.6%	< 30	93.7%	93.1%	50 - 150	
Methiocarb	0.028	0.363	0.381	0.400	0.4%	< 30	88.8%	88.3%	50 - 150	
Methomyl	0.000	0.730	0.725	0.800	0.6%	< 30	91.3%	90.7%	50 - 150	
MCK-264	0.000	0.299	0.338	0.400	12.5%	< 30	74.8%	84.8%	50 - 150	
Mydobutani	0.019	0.364	0.372	0.400	2.5%	< 30	86.2%	88.4%	50 - 150	
Naled	0.000	0.875	0.918	1.000	4.8%	< 30	87.5%	91.8%	50 - 150	
Oxaryl	0.000	1.897	1.860	2.000	2.0%	< 30	94.9%	93.0%	50 - 150	
Padobutrazole	0.000	0.718	0.755	0.800	5.2%	< 30	89.7%	94.5%	50 - 150	
Parathion-Methyl	0.000	0.312	0.461	0.400	38.7%	< 30	78.0%	115.3%	30 - 150	R
Permethrin	0.000	0.396	0.385	0.400	2.7%	< 30	98.9%	96.3%	50 - 150	
Phosmet	0.000	0.360	0.378	0.400	4.9%	< 30	90.1%	94.8%	50 - 150	
Piperonyl butoxide	0.005	1.600	1.720	2.000	7.2%	< 30	79.8%	85.7%	50 - 150	
Prallethrin	0.000	0.290	0.296	0.400	1.9%	< 30	72.8%	74.0%	50 - 150	
Propiconazole	0.000	0.596	0.657	0.800	9.6%	< 30	74.5%	82.1%	50 - 150	
Propoxur	0.009	0.368	0.368	0.400	0.2%	< 30	89.8%	89.9%	50 - 150	
Pyrethrin (Summe)	0.021	0.527	0.546	0.488	3.7%	< 30	103.7%	107.6%	50 - 150	
Pyridaben	0.000	0.463	0.449	0.400	2.9%	< 30	115.6%	112.3%	50 - 150	
Spinosad	0.000	0.379	0.395	0.388	4.4%	< 30	97.7%	102.1%	50 - 150	
Spiromesfen	0.000	0.443	0.432	0.400	2.5%	< 30	110.7%	108.0%	50 - 150	
Spirotetramat	0.000	0.369	0.376	0.400	1.8%	< 30	92.2%	93.9%	50 - 150	
Spiroxamine	0.000	0.760	0.744	0.800	2.2%	< 30	95.0%	92.9%	50 - 150	
Tebuconazole	0.000	0.716	0.652	0.800	9.4%	< 30	89.9%	81.5%	50 - 150	
Thiadoprid	0.000	0.354	0.355	0.400	0.2%	< 30	88.5%	88.7%	50 - 150	
Thiamethoxam	0.000	0.394	0.355	0.400	10.1%	< 30	98.4%	88.9%	50 - 150	
Trifloxystrobin	0.000	0.206	0.223	0.400	8.1%	< 30	51.4%	55.8%	50 - 150	


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Laboratory Quality Control Results

Residual Solvents				Batch ID: 2310111					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		486	584	µg/g	83.2	60 - 120	
Isobutane	ND	< 200		695	767	µg/g	90.6	60 - 120	
Butane	ND	< 200		682	782	µg/g	87.2	60 - 120	
2,2-Dimethylpropane	ND	< 200		794	939	µg/g	84.6	60 - 120	
Methanol	ND	< 200		1550	1670	µg/g	92.8	60 - 120	
Ethylene Oxide	ND	< 30		50	57.1	µg/g	87.6	60 - 120	
2-Methylbutane	ND	< 200		1430	1680	µg/g	85.1	60 - 120	
Pentane	ND	< 200		1400	1670	µg/g	83.8	60 - 120	
Ethanol	ND	< 200		1500	1660	µg/g	90.4	70 - 130	
Ethyl Ether	ND	< 200		1430	1670	µg/g	85.6	60 - 120	
2,2-Dimethylbutane	ND	< 30		169	189	µg/g	89.4	60 - 120	
Acetone	ND	< 200		1450	1670	µg/g	86.8	60 - 120	
2-Propanol	ND	< 200		1430	1630	µg/g	87.7	60 - 120	
Ethyl Formate	ND	< 500		4530	1600	µg/g	283.1	70 - 130	Q6
Acetonitrile	ND	< 100		412	492	µg/g	83.7	60 - 120	
Methyl Acetate	ND	< 500		1560	1600	µg/g	97.5	70 - 130	
2,3-Dimethylbutane	ND	< 30		152	180	µg/g	84.4	60 - 120	
Dichloromethane	ND	< 60		427	488	µg/g	87.5	60 - 120	
2-Methylpentane	ND	< 30		155	182	µg/g	85.2	60 - 120	
MTBE	ND	< 500		1590	1610	µg/g	98.8	70 - 130	
3-Methylpentane	ND	< 30		159	177	µg/g	89.8	60 - 120	
Hexane	ND	< 30		150	177	µg/g	84.7	60 - 120	
1-Propanol	ND	< 500		1640	1600	µg/g	102.5	70 - 130	
Methylethylketone	ND	< 500		1530	1610	µg/g	95.0	70 - 130	
Ethyl acetate	ND	< 200		1370	1630	µg/g	84.0	60 - 120	
2-Butanol	ND	< 200		1400	1630	µg/g	85.9	60 - 120	
Tetrahydrofuran	ND	< 100		417	488	µg/g	85.5	60 - 120	
Cyclohexane	ND	< 200		1380	1610	µg/g	85.7	60 - 120	
2-methyl-1-propanol	ND	< 500		1710	1610	µg/g	106.2	70 - 130	
Benzene	ND	< 1		3.58	4.79	µg/g	74.7	60 - 120	
Isopropyl Acetate	ND	< 200		1380	1650	µg/g	83.6	60 - 120	
Heptane	ND	< 200		1350	1630	µg/g	82.8	60 - 120	
1-Butanol	ND	< 500		1770	1600	µg/g	110.6	70 - 130	
Propyl Acetate	ND	< 500		1580	1600	µg/g	98.8	70 - 130	
1,4-Dioxane	ND	< 100		439	523	µg/g	83.9	60 - 120	
2-Ethoxyethanol	ND	< 30		146	179	µg/g	81.6	60 - 120	
Methylisobutylketone	ND	< 500		1610	1600	µg/g	100.6	70 - 130	
3-Methyl-1-butanol	ND	< 500		1710	1600	µg/g	106.9	70 - 130	
Ethylene Glycol	ND	< 200		319	508	µg/g	63.0	60 - 120	
Toluene	ND	< 100		428	496	µg/g	86.3	60 - 120	
Isobutyl Acetate	ND	< 500		1610	1610	µg/g	100.0	70 - 130	
1-Pentanol	ND	< 500		1860	1600	µg/g	116.3	70 - 130	
Butyl Acetate	ND	< 500		1610	1610	µg/g	100.0	70 - 130	
Ethylbenzene	ND	< 200		794	978	µg/g	81.2	60 - 120	
m,p-Xylene	ND	< 200		801	994	µg/g	80.6	60 - 120	
o-Xylene	ND	< 200		797	982	µg/g	81.2	60 - 120	
Cumene	ND	< 30		131	171	µg/g	76.6	60 - 120	
Anisole	ND	< 500		1760	1600	µg/g	110.0	70 - 130	
DMSO	ND	< 500		1400	1620	µg/g	86.4	70 - 130	
1,2-dimethoxyethane	ND	< 50		173	186	µg/g	93.0	70 - 130	
Triethylamine	ND	< 500		1430	1600	µg/g	89.4	70 - 130	
N,N-dimethylformamide	ND	< 150		483	480	µg/g	100.6	70 - 130	
N,N-dimethylacetamide	ND	< 150		478	483	µg/g	99.0	70 - 130	
Pyridine	ND	< 50		161	168	µg/g	95.8	70 - 130	
Silolane	ND	< 50		164	161	µg/g	101.9	70 - 130	
1,2-Dichloroethane	ND	< 1		0.837	1	µg/g	83.7	70 - 130	
Chloroform	ND	< 1		0.951	1	µg/g	95.1	70 - 130	
Trichloroethylene	ND	< 1		0.926	1	µg/g	92.6	70 - 130	
1,1,1-Trichloroethane	ND	< 1		0.758	1	µg/g	75.8	70 - 130	



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QC- Sample Duplicate		Sample ID: 23-009604-0001						
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

Units of Measure:

µg/g - Microgram per gram or ppm



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Laboratory Quality Control Results

JAOAC2015 V98-6 Batch ID: 2310117

Laboratory Control Sample									
Analyte	LCS	Result	Spike	Units	% Rec	Limits		Evaluation	Notes
CBDVA	2	0.0761	0.0718	%	106	80.0	- 120	Acceptable	
CBDV	2	0.0742	0.0708	%	105	80.0	- 120	Acceptable	
CBE	2	0.0840	0.0805	%	104	80.0	- 120	Acceptable	
CEDA	1	0.0790	0.0776	%	102	90.0	- 110	Acceptable	
CBGA	1	0.0781	0.0774	%	101	80.0	- 120	Acceptable	
CBG	1	0.0811	0.0794	%	102	80.0	- 120	Acceptable	
CBD	1	0.0829	0.0812	%	102	90.0	- 110	Acceptable	
THCV	2	0.0520	0.0513	%	101	80.0	- 120	Acceptable	
Δ8THCV	2	0.0622	0.0627	%	99.1	80.0	- 120	Acceptable	
THCV/A	2	0.0815	0.0715	%	114	80.0	- 120	Acceptable	
CBN	1	0.0789	0.0810	%	97.4	80.0	- 120	Acceptable	
exo-THC	2	0.0706	0.0718	%	98.3	80.0	- 120	Acceptable	
Δ9THC	1	0.0769	0.0796	%	96.7	90.0	- 110	Acceptable	
Δ8THC	1	0.0713	0.0750	%	95.0	90.0	- 110	Acceptable	
9SΔ10THC	1	0.0772	0.0816	%	94.5	80.0	- 120	Acceptable	
CBL	2	0.0719	0.0718	%	100	80.0	- 120	Acceptable	
9RΔ10THC	1	0.0701	0.0745	%	94.1	80.0	- 120	Acceptable	
CBG	2	0.0711	0.0736	%	96.7	80.0	- 120	Acceptable	
THCA	1	0.0720	0.0763	%	94.3	90.0	- 110	Acceptable	
CBGA	2	0.0794	0.0750	%	106	80.0	- 120	Acceptable	
CBLA	2	0.121	0.115	%	105	80.0	- 120	Acceptable	
Δ9THCP	2	0.0688	0.0746	%	92.2	80.0	- 120	Acceptable	
CBT	2	0.0613	0.0725	%	84.5	80.0	- 120	Acceptable	

Method Blank						
Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.0698	%	< 0.0698	Acceptable	
CBDV	<LOQ	0.0698	%	< 0.0698	Acceptable	
CBE	<LOQ	0.0698	%	< 0.0698	Acceptable	
CEDA	<LOQ	0.0698	%	< 0.0698	Acceptable	
CBGA	<LOQ	0.0698	%	< 0.0698	Acceptable	
CBG	<LOQ	0.0698	%	< 0.0698	Acceptable	
CBD	<LOQ	0.0698	%	< 0.0698	Acceptable	
THCV	<LOQ	0.0698	%	< 0.0698	Acceptable	
Δ8THCV	<LOQ	0.0698	%	< 0.0698	Acceptable	
THCV/A	<LOQ	0.0698	%	< 0.0698	Acceptable	
CBN	<LOQ	0.0698	%	< 0.0698	Acceptable	
exo-THC	<LOQ	0.0698	%	< 0.0698	Acceptable	
Δ9THC	<LOQ	0.0698	%	< 0.0698	Acceptable	
Δ8THC	<LOQ	0.0698	%	< 0.0698	Acceptable	
9SΔ10THC	<LOQ	0.0698	%	< 0.0698	Acceptable	
CBL	<LOQ	0.0698	%	< 0.0698	Acceptable	
9RΔ10THC	<LOQ	0.0698	%	< 0.0698	Acceptable	
CBG	<LOQ	0.0698	%	< 0.0698	Acceptable	
THCA	<LOQ	0.0698	%	< 0.0698	Acceptable	
CBGA	<LOQ	0.0698	%	< 0.0698	Acceptable	
CBLA	<LOQ	0.0698	%	< 0.0698	Acceptable	
Δ9THCP	<LOQ	0.0698	%	< 0.0698	Acceptable	
CBT	<LOQ	0.0698	%	< 0.0698	Acceptable	

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

Units of Measure:
 %- Percent



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 Portland, OR 97230
 503-254-1794

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

Laboratory Quality Control Results

JAOAC2015 V98-6		Batch ID: 2310117						
Sample Duplicate		Sample ID: 23-009667-0001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
CBDV	0.345	0.344	0.0755	%	0.194	< 20	Acceptable	
CBE	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
CBD	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
CBD ^A	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
CBD ^B	0.162	0.161	0.0755	%	0.577	< 20	Acceptable	
CBD	93.6	95.2	0.0755	%	1.63	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
Δ8THCV	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
THCV/A	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
CBN	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
Δ9THC	0.181	0.179	0.0755	%	1.28	< 20	Acceptable	
Δ8THC	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
9S-Δ10THC	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
9R-Δ10THC	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
CBC	0.160	0.156	0.0755	%	2.67	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
Δ9THCP	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	
CBT	<LOQ	<LOQ	0.0755	%	NA	< 20	Acceptable	

Abbreviations

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

Units of Measure:

%- Percent



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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.