



Customer: KIK By Kalibloom
Product identity: Mango Kush
Client/Metric ID: Batch # 012422
Laboratory ID: 22-006215-0007

Summary

Potency:

Analyte	Result (%)	 ● Δ8-THC ● Δ8-THCV ● CBT ● CBD ● CBG-A		
Δ8-THC	84.6		CBD-Total	<LOQ
Δ8-THCV	0.325		THC-Total	<LOQ
CBT [†]	0.183		(Reported in percent of total sample)	
CBD	0.137			
CBG-A [†]	0.0772			

Residual Solvents:

All analytes passing and less than LOQ.

Pesticides:

All analytes passing and less than LOQ.

Terpenes:

Analyte	Percent by weight	Percent of Total	Analyte	Percent by weight	Percent of Total
β-Myrcene [†]	0.918	31.76%	α-pinene [†]	0.664	22.98%
β-Caryophyllene [†]	0.432	14.95%	(-)-β-Pinene [†]	0.308	10.66%
(R)-(+)-Limonene [†]	0.182	6.30%	α-Bisabolol [†]	0.130	4.50%
Humulene [†]	0.112	3.88%	Linalool [†]	0.0982	3.40%
Terpinolene [†]	0.0272	0.94%	p-Cymene [†]	0.0224	0.78%
Total Terpenes[†]	2.89	100.00%			

Metals:

Less than LOQ for all analytes.



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503-254-1794

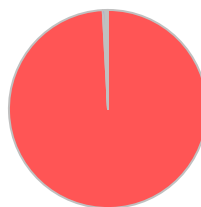


Report Number: 22-006215/D005.R000
Report Date: 06/06/2022
ORELAP#: OR100028
Purchase Order: Delta 8
Received: 05/27/22 15:39

Customer: KIK By Kalibloom
United States of America (USA)
Product identity: Mango Kush
Client/Metric ID: Batch # 012422
Sample Date:
Laboratory ID: 22-006215-0007
Evidence of Cooling: No
Temp: 21.4 °C
Relinquished by: FedEx
Batch Number: 012422

Sample Results

Potency	Method J AOAC 2015 V98-6 (mod)	Units %	Batch: 2204739	Analyze: 6/2/22 10:32:00 PM
Analyte	As Received	Dry weight	LOQ	Notes
CBC	< LOQ		0.0753	
CBC-A [†]	< LOQ		0.0753	
CBC-Total [†]	< LOQ		0.141	
CBD	0.137		0.0753	
CBD-A	< LOQ		0.0753	
CBD-Total	< LOQ		0.141	
CBDV [†]	< LOQ		0.0753	
CBDV-A [†]	< LOQ		0.0753	
CBDV-Total [†]	< LOQ		0.141	
CBE [†]	< LOQ		0.0753	
CBG [†]	< LOQ		0.0753	
CBG-A [†]	0.0772		0.0753	
CBG-Total	< LOQ		0.141	
CBL [†]	< LOQ		0.0753	
CBL-A [†]	< LOQ		0.0753	
CBL-Total [†]	< LOQ		0.141	
CBN	< LOQ		0.0753	
CBT [†]	0.183		0.0753	
Δ8-THC	84.6		0.753	
Δ8-THCV	0.325		0.0753	
Δ9-THC	< LOQ		0.0753	
exo-THC	< LOQ		0.0753	
THC-A	< LOQ		0.0753	
THC-Total	< LOQ		0.141	
THCV [†]	< LOQ		0.0753	
THCV-A [†]	< LOQ		0.0753	
THCV-Total [†]	< LOQ		0.141	
Total Cannabinoids[†]	85.3			



Δ8-THC
Δ8-THCV
CBT
CBD
CBG-A



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Received: 05/27/22 15:39

Solvents		Method Residual Solvents by GC/MS				Units µg/g	Batch 2204738	Analyze 06/06/22 10:09 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		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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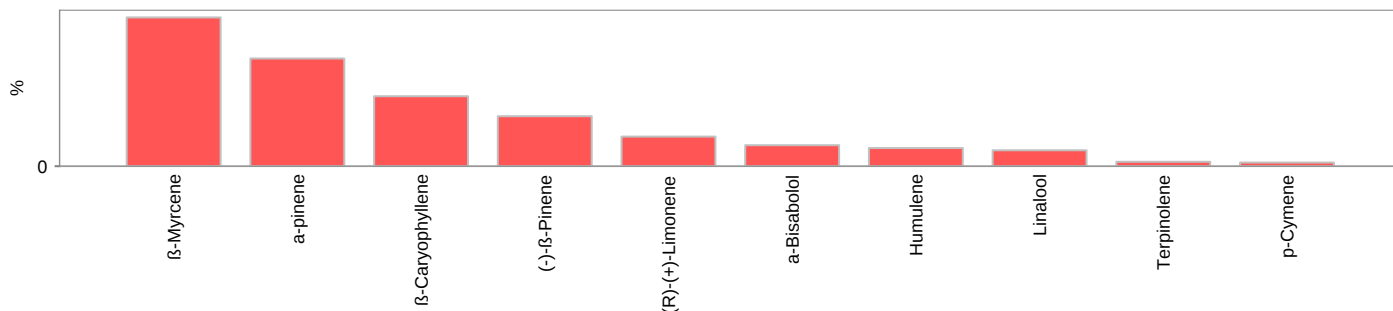


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Pesticides					Method AOAC 2007.01 & EN 15662 (mod)	Units mg/kg	Batch 2204640	Analyze 06/01/22 11:45 AM				
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes	
Abamectin	< LOQ	0.50	0.250	pass		Acephate	< LOQ	0.40	0.250	pass		
Acequinocyl	< LOQ	2.0	1.00	pass		Acetamiprid	< LOQ	0.20	0.100	pass		
Aldicarb	< LOQ	0.40	0.200	pass		Azoxystrobin	< LOQ	0.20	0.100	pass		
Bifenazate	< LOQ	0.20	0.100	pass		Bifenthrin	< LOQ	0.20	0.100	pass		
Boscalid	< LOQ	0.40	0.200	pass		Carbaryl	< LOQ	0.20	0.100	pass		
Carbofuran	< LOQ	0.20	0.100	pass		Chlorantraniliprole	< LOQ	0.20	0.100	pass		
Chlorfenapyr	< LOQ	1.0	0.500	pass		Chlorpyrifos	< LOQ	0.20	0.100	pass		
Clofentezine	< LOQ	0.20	0.100	pass		Cyfluthrin	< LOQ	1.0	0.500	pass		
Cypermethrin	< LOQ	1.0	0.500	pass		Daminozide	< LOQ	1.0	0.500	pass		
Diazinon	< LOQ	0.20	0.100	pass		Dichlorvos	< LOQ	1.0	0.500	pass		
Dimethoate	< LOQ	0.20	0.100	pass		Ethoprophos	< LOQ	0.20	0.100	pass		
Etofenprox	< LOQ	0.40	0.200	pass		Etoxazole	< LOQ	0.20	0.100	pass		
Fenoxycarb	< LOQ	0.20	0.100	pass		Fenpyroximate	< LOQ	0.40	0.200	pass		
Fipronil	< LOQ	0.40	0.200	pass		Flonicamid	< LOQ	1.0	0.400	pass		
Fludioxonil	< LOQ	0.40	0.200	pass		Hexythiazox	< LOQ	1.0	0.400	pass		
Imazalil	< LOQ	0.20	0.100	pass		Imidacloprid	< LOQ	0.40	0.200	pass		
Kresoxim-methyl	< LOQ	0.40	0.200	pass		Malathion	< LOQ	0.20	0.100	pass		
Metalaxyl	< LOQ	0.20	0.100	pass		Methiocarb	< LOQ	0.20	0.100	pass		
Methomyl	< LOQ	0.40	0.200	pass		MGK-264	< LOQ	0.20	0.100	pass		
Myclobutanil	< LOQ	0.20	0.100	pass		Naled	< LOQ	0.50	0.250	pass		
Oxamyl	< LOQ	1.0	0.500	pass		Paclobutrazole	< LOQ	0.40	0.200	pass		
Parathion-Methyl	< LOQ	0.20	0.200	pass		Permethrin	< LOQ	0.20	0.100	pass		
Phosmet	< LOQ	0.20	0.100	pass		Piperonyl butoxide	< LOQ	2.0	1.00	pass		
Prallethrin	< LOQ	0.20	0.200	pass		Propiconazole	< LOQ	0.40	0.200	pass		
Propoxur	< LOQ	0.20	0.100	pass		Pyrethrin I (total)	< LOQ	1.0	0.500	pass		
Pyridaben	< LOQ	0.20	0.100	pass		Spinosad	< LOQ	0.20	0.100	pass		
Spiromesifen	< LOQ	0.20	0.100	pass		Spirotetramat	< LOQ	0.20	0.100	pass		
Spiroxamine	< LOQ	0.40	0.200	pass		Tebuconazole	< LOQ	0.40	0.200	pass		
Thiacloprid	< LOQ	0.20	0.100	pass		Thiamethoxam	< LOQ	0.20	0.100	pass		
Trifloxystrobin	< LOQ	0.20	0.100	pass								



Terpenes				Method J AOAC 2015 V98-6	Units %	Batch 2204642	Analyze 06/01/22 05:06 AM		
Analyte	Result	LOQ	% of Total	Notes	Analyte	Result	LOQ	% of Total	Notes
β-Myrcene [†]	0.918	0.018	31.765%		α-pinene [†]	0.664	0.018	22.976%	
β-Caryophyllene [†]	0.432	0.018	14.948%		(-)-β-Pinene [†]	0.308	0.018	10.657%	
(R)-(+)-Limonene [†]	0.182	0.018	6.298%		α-Bisabolol [†]	0.130	0.018	4.498%	
Humulene [†]	0.112	0.018	3.875%		Linalool [†]	0.0982	0.018	3.3979%	
Terpinolene [†]	0.0272	0.018	0.9412%		p-Cymene [†]	0.0224	0.018	0.7751%	
(-)-α-Terpineol [†]	< LOQ	0.018	0.00%		Camphene [†]	< LOQ	0.018	0.00%	
(+)-fenchol [†]	< LOQ	0.018	0.00%		(-)-caryophyllene oxide [†]	< LOQ	0.018	0.00%	
farnesene [†]	< LOQ	0.018	0.00%		Geraniol [†]	< LOQ	0.018	0.00%	
Geranyl acetate [†]	< LOQ	0.018	0.00%		(±)-Camphor [†]	< LOQ	0.018	0.00%	
(-)-Isopulegol [†]	< LOQ	0.018	0.00%		nerol [†]	< LOQ	0.018	0.00%	
d-3-Carene [†]	< LOQ	0.018	0.00%		(+)-Pulegone [†]	< LOQ	0.018	0.00%	
Isoborneol [†]	< LOQ	0.018	0.00%		(+)-Borneol [†]	< LOQ	0.018	0.00%	
Menthol [†]	< LOQ	0.018	0.00%		(±)-trans-Nerolidol [†]	< LOQ	0.018	0.00%	
(-)-Guaiol [†]	< LOQ	0.018	0.00%		(+)-Cedrol [†]	< LOQ	0.018	0.00%	
(±)-cis-Nerolidol [†]	< LOQ	0.018	0.00%		(±)-fenchone [†]	< LOQ	0.018	0.00%	
α-cedrene [†]	< LOQ	0.018	0.00%		α-phellandrene [†]	< LOQ	0.018	0.00%	
α-Terpinene [†]	< LOQ	0.018	0.00%		cis-β-Ocimene [†]	< LOQ	0.006	0.00%	
Eucalyptol [†]	< LOQ	0.018	0.00%		gamma-Terpinene [†]	< LOQ	0.018	0.00%	
Sabinene [†]	< LOQ	0.018	0.00%		Sabinene hydrate [†]	< LOQ	0.018	0.00%	
trans-β-Ocimene [†]	< LOQ	0.012	0.00%		valencene [†]	< LOQ	0.018	0.00%	
Total Terpenes	2.89								



Metals

Analyte	Result	Limits	Units	LOQ	Batch	Analyze	Method	Status	Notes
Arsenic	< LOQ	0.200	mg/kg	0.0946	2204705	06/02/22	AOAC 2013.06 (mod.)	pass	
Cadmium	< LOQ	0.200	mg/kg	0.0946	2204705	06/02/22	AOAC 2013.06 (mod.)	pass	
Lead	< LOQ	0.500	mg/kg	0.0823	2204743	06/03/22	AOAC 2013.06 (mod.)	pass	
Mercury	< LOQ	0.100	mg/kg	0.0473	2204705	06/02/22	AOAC 2013.06 (mod.)	pass	



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

Abbreviations

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

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

† = Analyte not NELAP accredited.

Units of Measure

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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Hemp / Cannabis Usable / Extract Chain of Custody Record

Revision: 3.01 Control#: CF023 Rev 02/26/2020 Eff: 02/27/2020
ORELAP ID: OR100028

Company: Kik by Kalibloom Contact: Taylor Street: 3315 E Russell Rd STEAL # 346 City: Las Vegas State: NV Zip: 89120 ☐ Email Results: Kalibloomworldwide@gmail.com Ph: () Fx Results: () Billing (if different):				Analysis Requested				PO Number: Delta 8 Project Number: Project Name: Custom Reporting: Report to State: ☐ METRC or ☐ Other: Turnaround time: ☐ Standard ☐ Rush * ☒ Priority Rush * *Ask for availability Sampled by:			
Lab ID	Client Sample Identification	Date	Time	Potency	Metals	Solvents	Pesticides	Terpene	Sample Type	Weight (Units)	Comments/Metric ID
	Lemonade			X	X	X	X	X			Batch # 012422
	Tropicana Kush			X	X	X	X	X			
	Pineapple Express			X	X	X	X	X			
	Strawberry Cough			X	X	X	X	X			
	Purple			X	X	X	X	X			
	Fire OG			X	X	X	X	X			
	Manap Kush			X	X	X	X	X			
	Watermelon Kush			X	X	X	X	X			
	Kuntz			X	X	X	X	X			
Relinquished By:		Date	Time	Received By:		Date	Time	Lab Use Only:			
				DS		7/27/22	15:39	Shipped Via: <u>Fedex</u> or ☐ Client drop Evidence of cooling: ☐ Yes ☒ No - Temp (°C): <u>21.4° C</u> Sample in good condition: ☒ Yes ☐ No ☐ Cash ☐ Check ☐ CC ☐ Net: Prelog storage:			

† - Sample Type Codes: Vegetation (V) ; Isolates (S) ; Extract/Concentrate (C)

Samples submitted to Columbia Laboratories with testing requirements constitute an agreement for services in accordance with the current terms of service associated with this COC. By signing "Relinquished by" you are agreeing to these terms
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P: (503) 254-1794 | Fax: (503) 254-1452
info@columbialaboratories.com
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Document ID: 3177 Revision: 3
Effective: 04/26/2022
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PACKAGE RECEIVING FORM

Delivery Date: ☒ 5/27/22 ☐ Same as Opened By Date ☐ Unsure
How was the package delivered?
UPS ☒ FEDEX ☐ USPS ☐ DHL ☐ OTHER: _____
Tracking Number: 2735 5580 6608

		CIRCLE ONE	
1) Was package sealed with no evidence of holes/tampering?	YES	NO	
Further custody seal/tampering notes: _____			
2) Was packing material used?	YES	NO	
If YES: ☐ PEANUTS ☐ BUBBLE ☐ WRAP ☐ FOAM PAPER			
3) Was a Complete Chain of Custody (COC) received?	YES	NO	
Comment (PT?, Email?): not reling.			
4) Sample temperature upon arrival?	21.4	°C	
5) Evidence of cooling?	YES	NO	
If YES, What kind? ☐ ICE ☐ FREEZER PACK ☐ DRY ICE			
Insulation? ☐ PLASTIC COOLER ☐ STYROFOAM ☐ OTHER: _____			
6) Were sample containers sealed in separate plastic bags/secondary containment?	YES	NO	
7) Did sample containers arrive in good condition?	YES	NO	
If NO: ☐ LEAKED ☐ BROKEN ☐ OTHER: _____			
If NO: Suspect contamination of other samples? ☐ YES ☐ NO			
8) Sample labels present?	YES	NO	
9) Do sample labels agree with COC?	YES	NO	
If NO, number of sample containers received: _____			

Sample pre-log location:

R39 R44 F44 R99 ☒ CANNA SHELF ☐ FOOD SHELF Other: _____

Other Notes:

Received By (initials): DS Date: 5/27/22 Time: 5:39



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Revision: 3 Document ID: 3120
Legacy ID: CFL-C21 Worksheet Validated 10/30/2020

Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662			Units: mg/Kg		Batch ID: 2204640				
Method Blank				Laboratory Control Sample					
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes	
Abamectin	0.000	< 0.250		1.039	1.000	103.9	50.0 150		
Acephate	0.000	< 0.250		1.008	1.000	100.8	60.0 120		
Acequinocyl	0.000	< 1.000		4.027	4.000	100.7	40.0 160		
Acetamiprid	0.000	< 0.100		0.388	0.400	96.9	60.0 120		
Aldicarb	0.000	< 0.200		0.765	0.800	95.6	60.0 120		
Azoxystrobin	0.000	< 0.100		0.388	0.400	96.9	60.0 120		
Bifenazate	0.000	< 0.100		0.381	0.400	95.3	60.0 120		
Bifenthrin	0.000	< 0.100		0.403	0.400	100.7	50.0 150		
Boscalid	0.000	< 0.200		0.865	0.800	108.2	60.0 120		
Carbaryl	0.000	< 0.100		0.392	0.400	98.1	60.0 120		
Carbofuran	0.000	< 0.100		0.378	0.400	94.4	60.0 120		
Chlorantraniliprole	0.000	< 0.100		0.385	0.400	96.2	60.0 120		
Chlorfenapyr	0.000	< 0.500		1.708	2.000	85.4	60.0 120		
Chlorpyrifos	0.000	< 0.100		0.396	0.400	99.0	60.0 120		
Clofentezine	0.000	< 0.100		0.396	0.400	99.1	60.0 120		
Cyfluthrin	0.000	< 0.500		2.161	2.000	108.1	50.0 150		
Cypermethrin	0.000	< 0.500		2.036	2.000	101.8	50.0 150		
Daminozide	0.000	< 0.500		1.963	2.000	98.2	60.0 120		
Diazinon	0.000	< 0.100		0.397	0.400	99.2	60.0 120		
Dichlorvos	0.000	< 0.500		1.947	2.000	97.3	60.0 120		
Dimethoate	0.000	< 0.100		0.384	0.400	96.1	60.0 120		
Ethoprophos	0.000	< 0.100		0.384	0.400	96.1	60.0 120		
Etofenprox	0.000	< 0.200		0.782	0.800	97.8	50.0 150		
Etiozazole	0.000	< 0.100		0.399	0.400	99.9	60.0 120		
Fenoxycarb	0.000	< 0.100		0.397	0.400	99.2	60.0 120		
Fenpyroximate	0.000	< 0.200		0.800	0.800	100.0	60.0 120		
Fipronil	0.000	< 0.200		0.783	0.800	97.8	60.0 120		
Flonicamid	0.000	< 0.250		0.974	1.000	97.4	60.0 120		
Fludioxonil	0.000	< 0.200		0.800	0.800	100.0	50.0 150		
Hexythiazox	0.000	< 0.250		0.989	1.000	98.9	60.0 120		
Imazalil	0.000	< 0.100		0.405	0.400	101.4	60.0 120		
Imidacloprid	0.000	< 0.200		0.779	0.800	97.4	60.0 120		
Kresoxin-methyl	0.000	< 0.200		0.822	0.800	102.7	60.0 120		
Malathion	0.000	< 0.100		0.400	0.400	100.1	60.0 120		
Metaxalyl	0.000	< 0.100		0.381	0.400	95.2	60.0 120		
Methiocarb	0.000	< 0.100		0.386	0.400	96.5	60.0 120		
Methomyl	0.000	< 0.200		0.789	0.800	98.6	60.0 120		
MGK-264	0.000	< 0.100		0.401	0.400	100.2	50.0 150		
Myclobutanil	0.000	< 0.100		0.354	0.400	88.4	60.0 120		
Naled	0.000	< 0.250		0.977	1.000	97.7	50.0 150		
Oxamyl	0.000	< 0.500		1.948	2.000	97.4	60.0 120		
Paclobutrazole	0.000	< 0.200		0.774	0.800	96.8	60.0 120		
Parathion-Methyl	0.000	< 0.200		0.889	0.800	111.1	50.0 150		
Permethrin	0.000	< 0.100		0.404	0.400	101.0	50.0 150		
Phosmet	0.000	< 0.100		0.387	0.400	96.8	50.0 150		
Piperonyl butoxide	0.000	< 0.500		1.947	2.000	97.4	60.0 120		
Prallethrin	0.000	< 0.100		0.391	0.400	97.7	60.0 120		
Propiconazole	0.000	< 0.200		0.797	0.800	99.6	60.0 120		
Propoxur	0.000	< 0.100		0.387	0.400	96.8	60.0 120		
Pyrethrin (Summe)	0.000	< 0.100		0.413	0.413	99.9	60.0 120		
Pyridaben	0.000	< 0.100		0.402	0.400	100.6	50.0 150		
Spinosad	0.000	< 0.100		0.393	0.388	101.3	50.0 150		
Spiromesifen	0.000	< 0.100		0.408	0.400	102.0	60.0 120		
Spirotetramat	0.000	< 0.100		0.398	0.400	99.5	60.0 120		
Spiroxamine	0.000	< 0.200		0.781	0.800	97.6	60.0 120		
Tebuconazole	0.000	< 0.200		0.798	0.800	99.7	60.0 120		
Thiacloprid	0.000	< 0.100		0.392	0.400	98.0	60.0 120		
Thiamethoxam	0.000	< 0.100		0.372	0.400	93.0	60.0 120		
Trifloxystrobin	0.000	< 0.100		0.388	0.400	97.0	60.0 120		



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



Report Number: 22-006215/D005.R000
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Received: 05/27/22 15:39

Revision: 3 Document ID: 3120
Legacy ID: CFL-C21 Worksheet Validated 10/30/2020

Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662										
Units: mg/Kg										
Batch ID: 2204640										
Matrix Spike/Matrix Spike Duplicate Recoveries										
Sample ID: 22-005900-0001										
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes
Abamectin	0.000	0.805	0.725	1.000	10.5%	< 30	80.5%	72.5%	50 - 150	
Acephate	0.000	0.930	0.909	1.000	2.2%	< 30	93.0%	90.9%	50 - 150	
Acequinocyl	0.000	2.621	3.638	4.000	32.5%	< 30	65.5%	90.9%	50 - 150	R
Acetamiprid	0.000	0.341	0.342	0.400	0.2%	< 30	85.3%	85.4%	50 - 150	
Aldicarb	0.000	0.713	0.714	0.800	0.1%	< 30	89.1%	89.3%	50 - 150	
Azoxystrobin	0.000	0.248	0.251	0.400	1.0%	< 30	62.1%	62.7%	50 - 150	
Bifenazate	0.000	0.395	0.371	0.400	6.4%	< 30	98.8%	92.7%	50 - 150	
Bifenthrin	0.033	0.349	0.379	0.400	9.0%	< 30	79.0%	86.5%	50 - 150	
Boscalid	0.000	0.624	0.583	0.800	6.8%	< 30	78.0%	72.8%	50 - 150	
Carbaryl	0.000	0.315	0.300	0.400	4.6%	< 30	78.6%	75.1%	50 - 150	
Carbofuran	0.000	0.305	0.305	0.400	0.2%	< 30	76.2%	76.4%	50 - 150	
Chlorantraniliprole	0.000	0.413	0.392	0.400	5.1%	< 30	103.2%	98.0%	50 - 150	
Chlorfenvapir	0.000	0.679	0.736	2.000	8.2%	< 30	33.9%	36.8%	50 - 150	Q
Chlorpyrifos	0.000	0.389	0.362	0.400	7.2%	< 30	97.2%	90.5%	50 - 150	
Clofentezine	0.000	0.379	0.356	0.400	6.3%	< 30	94.7%	89.0%	50 - 150	
Cyfluthrin	0.000	0.986	1.056	2.000	6.9%	< 30	49.3%	52.8%	30 - 150	
Cypermethrin	0.000	1.013	1.008	2.000	0.5%	< 30	50.7%	50.4%	50 - 150	
Daminozide	0.000	1.906	1.821	2.000	4.6%	< 30	95.3%	91.0%	30 - 150	
Diazinon	0.000	0.343	0.346	0.400	0.7%	< 30	85.8%	86.5%	50 - 150	
Dichlorvos	0.000	2.030	2.058	2.000	1.4%	< 30	101.5%	102.9%	50 - 150	
Dimethoate	0.000	0.368	0.363	0.400	1.5%	< 30	92.0%	90.7%	50 - 150	
Ethoprophos	0.000	0.316	0.321	0.400	1.6%	< 30	79.0%	80.2%	50 - 150	
Etofenprox	0.000	0.526	0.513	0.800	2.6%	< 30	65.8%	64.1%	50 - 150	
Etoxazole	0.000	0.342	0.336	0.400	1.9%	< 30	85.6%	84.0%	50 - 150	
Fenoxycarb	0.000	0.343	0.317	0.400	7.7%	< 30	85.7%	79.3%	50 - 150	
Fenpyroximate	0.000	0.451	0.414	0.800	8.4%	< 30	56.4%	51.8%	50 - 150	
Fipronil	0.000	0.536	0.529	0.800	1.3%	< 30	66.9%	66.1%	50 - 150	
Flonicamid	0.000	0.789	0.892	1.000	12.3%	< 30	78.9%	89.2%	50 - 150	
Fludioxonil	0.000	0.981	0.994	0.800	1.4%	< 30	122.6%	124.3%	50 - 150	
Hexythiazox	0.000	0.594	0.527	1.000	11.9%	< 30	59.4%	52.7%	50 - 150	
Imazalil	0.000	0.384	0.324	0.400	17.0%	< 30	96.0%	80.9%	50 - 150	
Imidacloprid	0.000	0.783	0.755	0.800	3.6%	< 30	97.8%	94.3%	50 - 150	
Kresoxim-methyl	0.000	0.640	0.615	0.800	4.0%	< 30	80.0%	76.8%	50 - 150	
Malathion	0.000	0.307	0.286	0.400	7.2%	< 30	76.8%	71.4%	50 - 150	
Metaxyl	0.002	0.368	0.341	0.400	7.5%	< 30	91.5%	84.9%	50 - 150	
Methiocarb	0.000	0.332	0.322	0.400	3.0%	< 30	83.0%	80.6%	50 - 150	
Methomyl	0.000	0.659	0.646	0.800	2.0%	< 30	82.4%	80.8%	50 - 150	
MGK-264	0.000	0.253	0.237	0.400	6.3%	< 30	63.3%	59.4%	50 - 150	
Myclobutanil	0.000	0.299	0.261	0.400	13.6%	< 30	74.6%	65.1%	50 - 150	
Naled	0.000	0.035	FALSE	1.000	200.0%	< 30	3.5%	0.0%	50 - 150	R,Q
Oxamyl	0.000	1.706	1.973	2.000	14.5%	< 30	85.3%	98.6%	50 - 150	
Paclobutrazole	0.000	0.668	0.627	0.800	6.2%	< 30	83.4%	78.4%	50 - 150	
Parathion-Methyl	0.000	0.334	0.345	0.800	3.1%	< 30	41.8%	43.1%	30 - 150	
Permethrin	0.000	0.230	0.216	0.400	6.2%	< 30	57.6%	54.1%	50 - 150	
Phosmet	0.000	0.313	0.300	0.400	4.3%	< 30	78.4%	75.1%	50 - 150	
Piperonyl butoxide	0.000	1.532	1.464	2.000	4.6%	< 30	76.6%	73.2%	50 - 150	
Prallethrin	0.000	0.781	0.823	0.400	5.3%	< 30	195.2%	205.7%	50 - 150	Q
Propiconazole	0.011	0.819	0.833	0.800	1.7%	< 30	101.0%	102.7%	50 - 150	
Propoxur	0.000	0.335	0.324	0.400	3.2%	< 30	83.7%	81.0%	50 - 150	
Pyrethrin (Summe)	0.002	0.688	0.737	0.413	6.8%	< 30	166.2%	178.0%	50 - 150	Q
Pyridaben	0.000	0.313	0.277	0.400	12.2%	< 30	78.2%	69.3%	50 - 150	
Spinosad	0.000	0.377	0.391	0.388	3.6%	< 30	97.1%	100.7%	50 - 150	
Spiromesifen	0.000	0.402	0.385	0.400	4.3%	< 30	100.6%	96.3%	50 - 150	
Spirotetramat	0.000	0.480	0.459	0.400	4.6%	< 30	120.1%	114.7%	50 - 150	
Spiroxamine	0.000	0.785	0.732	0.800	7.0%	< 30	98.2%	91.6%	50 - 150	
Tebuconazole	0.000	0.710	0.670	0.800	5.9%	< 30	88.8%	83.7%	50 - 150	
Thiacloprid	0.000	0.356	0.345	0.400	3.2%	< 30	89.0%	86.1%	50 - 150	
Thiamethoxam	0.000	0.318	0.351	0.400	9.9%	< 30	79.4%	87.7%	50 - 150	
Trifloxystrobin	0.000	0.291	0.280	0.400	3.7%	< 30	72.7%	70.0%	50 - 150	



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Laboratory Quality Control Results

Residual Solvents				Batch ID: 2204738					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		477	572	µg/g	83.4	60 - 120	
Isobutane	ND	< 200		704	731	µg/g	96.3	60 - 120	
Butane	ND	< 200		691	731	µg/g	94.5	60 - 120	
2,2-Dimethylpropane	ND	< 200		821	936	µg/g	87.7	60 - 120	
Methanol	ND	< 200		1700	1620	µg/g	104.9	60 - 120	
Ethylene Oxide	ND	< 30		58.8	56.2	µg/g	104.6	60 - 120	
2-Methylbutane	ND	< 200		1800	1620	µg/g	111.1	60 - 120	
Pentane	ND	< 200		1770	1610	µg/g	109.9	60 - 120	
Ethanol	ND	< 200		1750	1630	µg/g	107.4	70 - 130	
Ethyl Ether	ND	< 200		1740	1620	µg/g	107.4	60 - 120	
2,2-Dimethylbutane	ND	< 30		182	174	µg/g	104.6	60 - 120	
Acetone	ND	< 200		1730	1650	µg/g	104.8	60 - 120	
2-Propanol	ND	< 200		1610	1610	µg/g	100.0	60 - 120	
Ethyl Formate	ND	< 500		1510	1600	µg/g	94.4	70 - 130	
Acetonitrile	ND	< 100		518	498	µg/g	104.0	60 - 120	
Methyl Acetate	ND	< 500		1550	1610	µg/g	96.3	70 - 130	
2,3-Dimethylbutane	ND	< 30		167	176	µg/g	94.9	60 - 120	
Dichloromethane	ND	< 60		556	510	µg/g	109.0	60 - 120	
2-Methylpentane	ND	< 30		173	176	µg/g	98.3	60 - 120	
MTBE	ND	< 500		1570	1600	µg/g	98.1	70 - 130	
3-Methylpentane	ND	< 30		179	175	µg/g	102.3	60 - 120	
Hexane	ND	< 30		183	177	µg/g	103.4	60 - 120	
1-Propanol	ND	< 500		1510	1610	µg/g	93.8	70 - 130	
Methyl ethyl ketone	ND	< 500		1550	1600	µg/g	96.9	70 - 130	
Ethyl acetate	ND	< 200		1650	1630	µg/g	101.2	60 - 120	
2-Butanol	ND	< 200		1530	1620	µg/g	94.4	60 - 120	
Tetrahydrofuran	ND	< 100		492	500	µg/g	98.4	60 - 120	
Cyclohexane	ND	< 200		1610	1620	µg/g	99.4	60 - 120	
2-methyl-1-propanol	ND	< 500		1450	1620	µg/g	89.5	70 - 130	
Benzene	ND	< 1		4.91	5.32	µg/g	92.3	60 - 120	
Isopropyl Acetate	ND	< 200		1600	1620	µg/g	98.8	60 - 120	
Heptane	ND	< 200		1640	1770	µg/g	92.7	60 - 120	
1-Butanol	ND	< 500		1510	1600	µg/g	94.4	70 - 130	
Propyl Acetate	ND	< 500		1540	1600	µg/g	96.3	70 - 130	
1,4-Dioxane	ND	< 100		457	504	µg/g	90.7	60 - 120	
2-Ethoxyethanol	ND	< 30		137	161	µg/g	75.7	60 - 120	
Methylisobutylketone	ND	< 500		1390	1610	µg/g	86.3	70 - 130	
3-Methyl-1-butanol	ND	< 500		1390	1610	µg/g	86.3	70 - 130	
Ethylene Glycol	ND	< 200		337	494	µg/g	68.2	60 - 120	
Toluene	ND	< 100		425	491	µg/g	86.6	60 - 120	
Isobutyl Acetate	ND	< 500		1480	1600	µg/g	92.5	70 - 130	
1-Pentanol	ND	< 500		1340	1610	µg/g	83.2	70 - 130	
Butyl Acetate	ND	< 500		1260	1610	µg/g	78.3	70 - 130	
Ethylbenzene	ND	< 200		748	973	µg/g	76.9	60 - 120	
m,p-Xylene	ND	< 200		754	996	µg/g	75.7	60 - 120	
o-Xylene	ND	< 200		720	973	µg/g	74.0	60 - 120	
Cumene	ND	< 30		121	170	µg/g	71.2	60 - 120	
Anisole	ND	< 500		1230	1610	µg/g	76.4	70 - 130	
DMSO	ND	< 500		1080	1630	µg/g	66.3	70 - 130	Q6
1,2-dimethoxyethane	ND	< 50		156	164	µg/g	95.1	70 - 130	
Triethylamine	ND	< 500		1220	1600	µg/g	76.3	70 - 130	
N,N-dimethylformamide	ND	< 150		388	497	µg/g	78.1	70 - 130	
N,N-dimethylacetamide	ND	< 150		354	498	µg/g	71.1	70 - 130	
Pyridine	ND	< 50		151	180	µg/g	83.9	70 - 130	
1,2-Dichloroethane	ND	< 1		0.848	1	µg/g	84.8	70 - 130	
Chloroform	ND	< 1		0.863	1	µg/g	86.3	70 - 130	
Trichloroethylene	ND	< 1		0.812	1	µg/g	81.2	70 - 130	



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QC - Sample Duplicate				Sample ID: 22-006204-0002			
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Accept/Fail
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	313	317	200	µg/g	1.3	< 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
Methyl ethyl ketone	ND	ND	500	µg/g	0.0	< 20	Acceptable
Ethyl acetate	ND	ND	200	µg/g	0.0	< 20	Acceptable
2-Butanol	ND	ND	200	µg/g	0.0	< 20	Acceptable
Tetrahydrofuran	ND	ND	100	µg/g	0.0	< 20	Acceptable
Cyclohexane	ND	ND	200	µg/g	0.0	< 20	Acceptable
2-methyl-1-propanol	ND	ND	500	µg/g	0.0	< 20	Acceptable
Benzene	ND	ND	1	µg/g	0.0	< 20	Acceptable
Isopropyl Acetate	ND	ND	200	µg/g	0.0	< 20	Acceptable
Heptane	ND	ND	200	µg/g	0.0	< 20	Acceptable
1-Butanol	ND	ND	500	µg/g	0.0	< 20	Acceptable
Propyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable
1,4-Dioxane	ND	ND	100	µg/g	0.0	< 20	Acceptable
2-Ethoxyethanol	ND	ND	30	µg/g	0.0	< 20	Acceptable
Methyl isobutyl ketone	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
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

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
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Laboratory Quality Control Results

J AOAC 2015 V98-6 Batch ID: 2204739										
Laboratory Control Sample										
Analyte	LCS	Result	Spike	Units	% Rec	Limits			Evaluation	Notes
CBDVA	1	0.104	0.100	%	104	80.0	-	120	Acceptable	
CBDV	1	0.117	0.100	%	117	80.0	-	120	Acceptable	
CBE	1	0.0813	0.100	%	81.3	80.0	-	120	Acceptable	
CBDA	1	0.0911	0.100	%	91.1	90.0	-	110	Acceptable	
CBGA	1	0.0957	0.100	%	95.7	80.0	-	120	Acceptable	
CBG	1	0.0917	0.100	%	91.7	80.0	-	120	Acceptable	
CBD	1	0.1000	0.100	%	100	90.0	-	110	Acceptable	
THCV	1	0.0953	0.100	%	95.3	80.0	-	120	Acceptable	
d8THCV	1	0.0947	0.100	%	94.7	80.0	-	120	Acceptable	
THCVA	1	0.101	0.100	%	101	80.0	-	120	Acceptable	
CBN	1	0.0971	0.100	%	97.1	90.0	-	110	Acceptable	
exo-THC	1	0.0918	0.100	%	91.8	80.0	-	120	Acceptable	
d9THC	1	0.0914	0.100	%	91.4	90.0	-	110	Acceptable	
d8THC	1	0.0887	0.100	%	88.7	80.0	-	120	Acceptable	
CBL	1	0.0907	0.100	%	90.7	80.0	-	120	Acceptable	
CBC	1	0.0972	0.100	%	97.2	80.0	-	120	Acceptable	
THCA	1	0.0958	0.100	%	95.8	90.0	-	110	Acceptable	
CBCA	1	0.105	0.100	%	105	80.0	-	120	Acceptable	
CBLA	1	0.105	0.100	%	105	80.0	-	120	Acceptable	
CBT	1	0.0810	0.100	%	81.0	80.0	-	120	Acceptable	

Method Blank

Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.077	%	< 0.077	Acceptable	
CBDV	<LOQ	0.077	%	< 0.077	Acceptable	
CBE	<LOQ	0.077	%	< 0.077	Acceptable	
CBDA	<LOQ	0.077	%	< 0.077	Acceptable	
CBGA	<LOQ	0.077	%	< 0.077	Acceptable	
CBG	<LOQ	0.077	%	< 0.077	Acceptable	
CBD	<LOQ	0.077	%	< 0.077	Acceptable	
THCV	<LOQ	0.077	%	< 0.077	Acceptable	
d8THCV	<LOQ	0.077	%	< 0.077	Acceptable	
THCVA	<LOQ	0.077	%	< 0.077	Acceptable	
CBN	<LOQ	0.077	%	< 0.077	Acceptable	
exo-THC	<LOQ	0.077	%	< 0.077	Acceptable	
d9THC	<LOQ	0.077	%	< 0.077	Acceptable	
d8THC	<LOQ	0.077	%	< 0.077	Acceptable	
CBL	<LOQ	0.077	%	< 0.077	Acceptable	
CBC	<LOQ	0.077	%	< 0.077	Acceptable	
THCA	<LOQ	0.077	%	< 0.077	Acceptable	
CBCA	<LOQ	0.077	%	< 0.077	Acceptable	
CBLA	<LOQ	0.077	%	< 0.077	Acceptable	
CBT	<LOQ	0.077	%	< 0.077	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: 22-006215/D005.R000
Report Date: 06/06/2022
ORELAP#: OR100028
Purchase Order: Delta 8
Received: 05/27/22 15:39

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

Laboratory Quality Control Results

J AOAC 2015 V98-6				Batch ID: 2204739				
Sample Duplicate				Sample ID: 22-006187-0001				
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBDV	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBE	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBDA	0.0892	0.127	0.077	%	35.0	< 20	Outlier	R
CBGA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBG	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBD	0.380	0.459	0.077	%	18.8	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
d8THCV	0.289	0.240	0.077	%	18.6	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBN	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
d9THC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
d8THC	63.6	54.5	0.077	%	15.4	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBCA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBT	0.0974	<LOQ	0.077	%	NA	< 20	Acceptable	R2

Abbreviations

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

Units of Measure:



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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.