



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

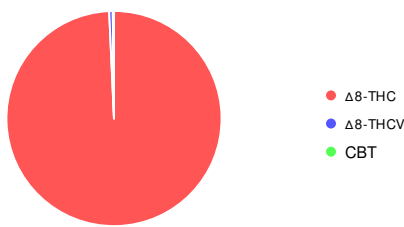


Report Number: 22-014446/D003.R000
Report Date: 12/01/2022
ORELAP#: OR100028
Purchase Order: Delta 8
Received: 11/28/22 10:32

Customer: KIK By Kalibloom
Product identity: Mimosa
Client/Metric ID: Batch # 210038
Laboratory ID: 22-014446-0002

Summary

Potency:

Analyte	Result (%)		(Reported in percent of total sample)		
Δ8-THC	78.0				
Δ8-THCV	0.421				
CBT	0.142				
		CBD-Total			<LOQ
		THC-Total			<LOQ

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	0.468	31.20%	(R)-(+)-Limonene	0.317	21.13%
α-pinene	0.268	17.87%	β-Caryophyllene	0.170	11.33%
(-)-β-Pinene	0.130	8.67%	Linalool	0.0762	5.08%
Humulene	0.0511	3.41%	α-Bisabolol	0.0241	1.61%
Total Terpenes	1.50	100.00%			

Metals:

Less than LOQ for all analytes.



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Received: 11/28/22 10:32

Customer: KIK By Kalibloom
United States of America (USA)

Product identity: Mimosa
Client/Metric ID: Batch # 210038

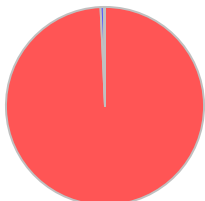
Sample Date:
Laboratory ID: 22-014446-0002

Evidence of Cooling: No

Temp: 14.6

Relinquished by: ups

Sample Results

Potency	Method: J AOAC 2015 V98-6 (mod) ^p			Units %	Batch: 2210176	Analyze: 11/29/22 11:32:00 P
Analyte	As Received	Dry weight	LOQ	Notes		
CBC	< LOQ		0.0738			
CBC-A	< LOQ		0.0738			
CBC-Total	< LOQ		0.139			
CBD	< LOQ		0.0738			
CBD-A	< LOQ		0.0738			
CBD-Total	< LOQ		0.139			
CBDV	< LOQ		0.0738			
CBDV-A	< LOQ		0.0738			
CBDV-Total	< LOQ		0.138			
CBE	< LOQ		0.0738			
CBG	< LOQ		0.0738			
CBG-A	< LOQ		0.0738			
CBG-Total	< LOQ		0.138			
CBL	< LOQ		0.0738			
CBL-A	< LOQ		0.0738			
CBL-Total	< LOQ		0.139			
CBN	< LOQ		0.0738			
CBT	0.142		0.0738			
Δ10-THC	< LOQ		0.0738			
Δ8-THC	78.0		0.738			
Δ8-THCV	0.421		0.0738			
Δ9-THC	< LOQ		0.0738			
exo-THC	< LOQ		0.0738			
THC-A	< LOQ		0.0738			
THC-Total	< LOQ		0.139			
THCV	< LOQ		0.0738			
THCV-A	< LOQ		0.0738			
THCV-Total	< LOQ		0.138			
Total Cannabinoids	78.6					



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Solvents		Method: Residual Solvents by GC/MS ^b				Units µg/g	Batch 2210164	Analyze 11/30/22 09:13 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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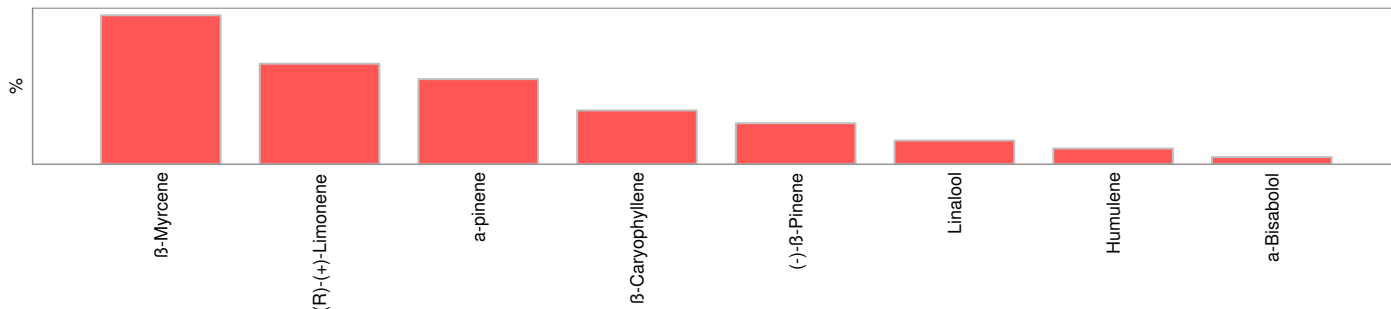


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Pesticides					Method: AOAC 2007.01 & EN 15662 (mod) ^b	Units mg/kg	Batch 2210196	Analyze 11/30/22 03:36 PM				
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes	
Abamectin [¥]	< LOQ	0.50	0.250	pass		Acephate [¥]	< LOQ	0.40	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		Paclobutrazole [¥]	< 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								



Terpenes					Method: J AOAC 2015 V98-6	Units %	Batch 2210177	Analyze 11/29/22 07:56 PM	
Analyte	Result	LOQ	% of Total	Notes	Analyte	Result	LOQ	% of Total	Notes
β-Myrcene	0.468	0.019	31.200%		(R)-(+)-Limonene	0.317	0.019	21.133%	
α-pinene	0.268	0.019	17.867%		β-Caryophyllene	0.170	0.019	11.333%	
(-)-β-Pinene	0.130	0.019	8.667%		Linalool	0.0762	0.019	5.0800%	
Humulene	0.0511	0.019	3.4067%		α-Bisabolol	0.0241	0.019	1.6067%	
Geraniol	< LOQ	0.019	0.00%		(+)-Pulegone	< LOQ	0.019	0.00%	
farnesene	< LOQ	0.019	0.00%		Isoborneol	< LOQ	0.019	0.00%	
valencene	< LOQ	0.019	0.00%		p-Cymene	< LOQ	0.019	0.00%	
(+)-fenchol	< LOQ	0.019	0.00%		Camphene	< LOQ	0.019	0.00%	
(±)-Camphor	< LOQ	0.019	0.00%		(-)-Isopulegol	< LOQ	0.019	0.00%	
Geranyl acetate	< LOQ	0.019	0.00%		(±)-trans-Nerolidol	< LOQ	0.019	0.00%	
(±)-fenchone	< LOQ	0.019	0.00%		Terpinolene	< LOQ	0.019	0.00%	
Menthol	< LOQ	0.019	0.00%		(-)-Guaiol	< LOQ	0.019	0.00%	
(-)-caryophyllene oxide	< LOQ	0.019	0.00%		Eucalyptol	< LOQ	0.019	0.00%	
(+)-Cedrol	< LOQ	0.019	0.00%		(±)-cis-Nerolidol	< LOQ	0.019	0.00%	
cis-β-Ocimene	< LOQ	0.006	0.00%		α-Terpinene	< LOQ	0.019	0.00%	
Sabinene	< LOQ	0.019	0.00%		α-phellandrene	< LOQ	0.019	0.00%	
(-)-α-Terpineol	< LOQ	0.019	0.00%		(+)-Borneol	< LOQ	0.019	0.00%	
α-cedrene	< LOQ	0.019	0.00%		d-3-Carene	< LOQ	0.019	0.00%	
gamma-Terpinene	< LOQ	0.019	0.00%		nerol	< LOQ	0.019	0.00%	
Sabinene hydrate	< LOQ	0.019	0.00%		trans-β-Ocimene	< LOQ	0.012	0.00%	
Total Terpenes	1.50								



Metals

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
Arsenic	< LOQ	0.200	mg/kg	0.0985	2210243	12/01/22 AOAC 2013.06 (mod.) ^p	pass	
Cadmium	< LOQ	0.200	mg/kg	0.0985	2210243	12/01/22 AOAC 2013.06 (mod.) ^p	pass	
Lead	< LOQ	0.500	mg/kg	0.0985	2210243	12/01/22 AOAC 2013.06 (mod.) ^p	pass	
Mercury	< LOQ	0.100	mg/kg	0.0493	2210243	12/01/22 AOAC 2013.06 (mod.) ^p	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.

[®] = ISO/IEC 17025:2017 accredited method.

[™] = TNI accredited analyte.

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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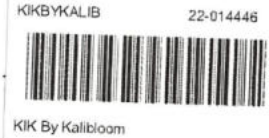


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Hemp / Cannabis Usable / Extract
Chain of Custody Record
Revision: 3.01 Control#: CFO23 Rev 02/26/2020 Eff: 02/27/2020
ORELAP ID: OR100028

11-16



Company: <u>Kik by Kalibloom</u>				Analysis Requested				PO Number: <u>Delta 8</u>			
Contact: <u>Taylor</u>								Project Number:			
Street: <u>3315 E Russell Rd STEALL #346</u>								Project Name:			
City: <u>Las Vegas</u> State: <u>NV</u> Zip: <u>89120</u>								Custom Reporting:			
☐ Email Results: <u>Taylor@kalibloom.com</u>								Report to State - ☐ METRC or ☐ Other:			
Ph: () Fx Results: ()								Turnaround time: ☐ Standard ☐ Rush * ☒ Priority Rush *			
Billing (if different):								*Ask for availability			
Lab ID	Client Sample Identification	Date	Time	Potency	Metals	Solvents	Pesticides	Terpene	Sample Type	Weight (Units)	Comments/Metric ID
	<u>Blue Dream</u>			X	X	X	X	X			<u>Batch #</u> <u>210038</u>
	<u>Mimosa</u>			X	X	X	X	X			
	<u>Fire OG</u>			X	X	X	X	X			
Relinquished By:				Date	Time	Received By:		Date	Time	Lab Use Only:	
						<u>JSF</u>		<u>11/28</u>	<u>10:32</u>	☒ Shipped Via: <u>VPS</u> or ☐ Client drop	
										Evidence of cooling: ☐ Yes ☒ No - Temp (°C): <u>14.4</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
12423 NE Whitaker Way
Portland, OR 97230

P: (503) 254-1794 | Fax: (503) 254-1452
info@columbiaboratories.com

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www.columbiaboratories.com



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

Laboratory Quality Control Results

Residual Solvents				Batch ID: 2210164					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		601	572	µg/g	105.1	60 - 120	
Isobutane	ND	< 200		751	731	µg/g	102.7	60 - 120	
Butane	ND	< 200		744	731	µg/g	101.8	60 - 120	
2,2-Dimethylpropane	ND	< 200		977	936	µg/g	104.4	60 - 120	
Methanol	ND	< 200		1730	1650	µg/g	104.8	60 - 120	
Ethylene Oxide	ND	< 30		58.5	56.2	µg/g	104.1	60 - 120	
2-Methylbutane	ND	< 200		1660	1650	µg/g	100.6	60 - 120	
Pentane	ND	< 200		1660	1650	µg/g	100.6	60 - 120	
Ethanol	ND	< 200		1650	1660	µg/g	99.4	70 - 130	
Ethyl Ether	ND	< 200		1580	1630	µg/g	96.9	60 - 120	
2,2-Dimethylbutane	ND	< 30		202	189	µg/g	106.9	60 - 120	
Acetone	ND	< 200		1700	1650	µg/g	103.0	60 - 120	
2-Propanol	ND	< 200		1710	1650	µg/g	103.6	60 - 120	
Ethyl Formate	ND	< 500		1510	1610	µg/g	93.8	70 - 130	
Acetonitrile	ND	< 100		551	504	µg/g	109.3	60 - 120	
Methyl Acetate	ND	< 500		1740	1630	µg/g	106.7	70 - 130	
2,3-Dimethylbutane	ND	< 30		165	174	µg/g	94.8	60 - 120	
Dichloromethane	ND	< 60		514	521	µg/g	98.7	60 - 120	
2-Methylpentane	ND	< 30		187	187	µg/g	100.0	60 - 120	
MTBE	ND	< 500		1640	1600	µg/g	102.5	70 - 130	
3-Methylpentane	ND	< 30		198	188	µg/g	105.3	60 - 120	
Hexane	ND	< 30		199	182	µg/g	109.3	60 - 120	
1-Propanol	ND	< 500		1720	1610	µg/g	106.8	70 - 130	
Methylethylketone	ND	< 500		1690	1600	µg/g	105.6	70 - 130	
Ethyl acetate	ND	< 200		1660	1630	µg/g	101.8	60 - 120	
2-Butanol	ND	< 200		1650	1630	µg/g	101.2	60 - 120	
Tetrahydrofuran	ND	< 100		532	506	µg/g	105.1	60 - 120	
Cyclohexane	ND	< 200		1580	1640	µg/g	96.3	60 - 120	
2-methyl-1-propanol	ND	< 500		1610	1620	µg/g	99.4	70 - 130	
Benzene	ND	< 1		4.51	4.93	µg/g	91.5	60 - 120	
Isopropyl Acetate	ND	< 200		1650	1640	µg/g	100.6	60 - 120	
Heptane	ND	< 200		1650	1630	µg/g	101.2	60 - 120	
1-Butanol	ND	< 500		1630	1600	µg/g	101.9	70 - 130	
Propyl Acetate	ND	< 500		1670	1620	µg/g	103.1	70 - 130	
1,4-Dioxane	ND	< 100		471	493	µg/g	95.5	60 - 120	
2-Ethoxyethanol	ND	< 30		169	171	µg/g	98.8	60 - 120	
Methylisobutylketone	ND	< 500		1660	1620	µg/g	102.5	70 - 130	
3-Methyl-1-butanol	ND	< 500		1680	1610	µg/g	104.3	70 - 130	
Ethylene Glycol	ND	< 200		495	494	µg/g	100.2	60 - 120	
Toluene	ND	< 100		482	506	µg/g	95.3	60 - 120	
Isobutyl Acetate	ND	< 500		1700	1620	µg/g	104.9	70 - 130	
1-Pentanol	ND	< 500		1680	1610	µg/g	104.3	70 - 130	
Butyl Acetate	ND	< 500		1610	1610	µg/g	100.0	70 - 130	
Ethylbenzene	ND	< 200		938	996	µg/g	94.2	60 - 120	
m,p-Xylene	ND	< 200		913	1010	µg/g	90.4	60 - 120	
o-Xylene	ND	< 200		921	979	µg/g	94.1	60 - 120	
Cumene	ND	< 30		157	188	µg/g	83.5	60 - 120	
Anisole	ND	< 500		1620	1610	µg/g	100.6	70 - 130	
DMSO	ND	< 500		1560	1600	µg/g	97.5	70 - 130	
1,2-dimethoxyethane	ND	< 50		205	190	µg/g	107.9	70 - 130	
Triethylamine	ND	< 500		1540	1610	µg/g	95.7	70 - 130	
N,N-dimethylformamide	ND	< 150		434	496	µg/g	87.5	70 - 130	
N,N-dimethylacetamide	ND	< 150		515	483	µg/g	106.6	70 - 130	
Pyridine	ND	< 50		157	167	µg/g	94.0	70 - 130	
Sulfolane	ND	< 50		115	161	µg/g	71.4	70 - 130	



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

QC - Sample Duplicate

Sample ID: 22-014208-0001

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

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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Portland, OR 97230
503-254-1794



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Report Date: 12/01/2022
ORELAP#: OR100028
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Received: 11/28/22 10:32

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

Laboratory Quality Control Results

J AOAC 2015 V98-6 Batch ID: 2210176

Laboratory Control Sample										
Analyte	LCS	Result	Spike	Units	% Rec	Limits			Evaluation	Notes
CBDVA	2	0.105	0.102	%	102	80.0	-	120	Acceptable	
CBDV	2	0.108	0.106	%	102	80.0	-	120	Acceptable	
CBE	2	0.106	0.106	%	100	80.0	-	120	Acceptable	
CBD	1	0.102	0.096	%	107	90.0	-	110	Acceptable	
CBGA	1	0.102	0.097	%	106	80.0	-	120	Acceptable	
CBG	1	0.103	0.095	%	108	80.0	-	120	Acceptable	
CBD	1	0.103	0.096	%	108	90.0	-	110	Acceptable	
THCV	2	0.107	0.102	%	105	80.0	-	120	Acceptable	
d8THCV	2	0.110	0.109	%	101	80.0	-	120	Acceptable	
THCVA	2	0.102	0.100	%	102	80.0	-	120	Acceptable	
CBN	1	0.104	0.099	%	105	80.0	-	120	Acceptable	
exo-THC	2	0.103	0.098	%	104	80.0	-	120	Acceptable	
d9THC	1	0.108	0.102	%	106	90.0	-	110	Acceptable	
d8THC	1	0.0992	0.100	%	99.1	90.0	-	110	Acceptable	
CBL	2	0.103	0.100	%	103	80.0	-	120	Acceptable	
9S-HHC	3	0.100	0.100	%	100	80.0	-	120	Acceptable	
d10THC	1	0.0981	0.092	%	106	80.0	-	120	Acceptable	
CBG	2	0.108	0.105	%	103	80.0	-	120	Acceptable	
9R-HHC	3	0.0920	0.100	%	92.0	80.0	-	120	Acceptable	
THCA	1	0.104	0.096	%	108	90.0	-	110	Acceptable	
CBGA	2	0.101	0.103	%	98.0	80.0	-	120	Acceptable	
CBLA	2	0.107	0.106	%	101	80.0	-	120	Acceptable	
d8THCO	3	0.107	0.100	%	107	80.0	-	120	Acceptable	
CBT	2	0.101	0.110	%	92.0	80.0	-	120	Acceptable	
d9THCO	3	0.107	0.100	%	107	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	
CBD	<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	
9S-HHC	<LOQ	0.077	%	< 0.077	Acceptable	
d10THC	<LOQ	0.077	%	< 0.077	Acceptable	
CBG	<LOQ	0.077	%	< 0.077	Acceptable	
9R-HHC	<LOQ	0.077	%	< 0.077	Acceptable	
THCA	<LOQ	0.077	%	< 0.077	Acceptable	
CBGA	<LOQ	0.077	%	< 0.077	Acceptable	
CBLA	<LOQ	0.077	%	< 0.077	Acceptable	
d8THCO	<LOQ	0.077	%	< 0.077	Acceptable	
CBT	<LOQ	0.077	%	< 0.077	Acceptable	
d9THCO	<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



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Portland, OR 97230
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Laboratory Quality Control Results

J AOAC 2015 V98-6 Batch ID: 2210176								
Sample Duplicate Sample ID: 22-014446-0001								
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	
CBD	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBG	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
d8THCV	0.415	0.406	0.077	%	2.23	< 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	75.3	76.5	0.077	%	1.50	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
9S-HHC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
d10THC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
9R-HHC	<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	
d8THCO	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBT	0.138	0.139	0.077	%	0.438	< 20	Acceptable	
d9THCO	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	

Abbreviations

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

Units of Measure:



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Terpenes Quality Control Results

Method Reference: EPA 5035				Batch ID: 2210177					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	LCS	Units	LCS % Rec	Limits	Notes
a-pinene	<LOQ	< 200		475	500	µg/g	95%	70 - 130	
Camphene	<LOQ	< 200		406	500	µg/g	81%	70 - 130	
Sabinene	<LOQ	< 200		471	500	µg/g	94%	70 - 130	
b-Pinene	<LOQ	< 200		466	500	µg/g	93%	70 - 130	
b-Myrcene	<LOQ	< 200		400	500	µg/g	80%	70 - 130	
a-phellandrene	<LOQ	< 200		411	500	µg/g	82%	70 - 130	
d-3-Carene	<LOQ	< 200		403	500	µg/g	81%	70 - 130	
a-Terpinene	<LOQ	< 200		492	500	µg/g	98%	70 - 130	
p-Cymene	<LOQ	< 200		403	500	µg/g	81%	70 - 130	
D-Limonene	<LOQ	< 200		481	500	µg/g	96%	70 - 130	
Eucalyptol	<LOQ	< 200		397	500	µg/g	79%	70 - 130	
b-cis-Ocimene	<LOQ	< 67		145	167	µg/g	87%	70 - 130	
b-trans-Ocimene	<LOQ	< 133		280	333	µg/g	84%	70 - 130	
g-Terpinene	<LOQ	< 200		480	500	µg/g	96%	70 - 130	
Sabinene_Hydrate	<LOQ	< 200		478	500	µg/g	96%	70 - 130	
Terpinolene	<LOQ	< 200		493	500	µg/g	99%	70 - 130	
D-Fenchone	<LOQ	< 200		473	500	µg/g	95%	70 - 130	
Linalool	<LOQ	< 200		474	500	µg/g	95%	70 - 130	
Fenchol	<LOQ	< 200		498	500	µg/g	100%	70 - 130	
Camphor	<LOQ	< 200		410	500	µg/g	82%	70 - 130	
Isopulego	<LOQ	< 200		439	500	µg/g	88%	70 - 130	
Isoborneol	<LOQ	< 200		420	500	µg/g	84%	70 - 130	
Borneol	<LOQ	< 200		490	500	µg/g	98%	70 - 130	
DL-Menthol	<LOQ	< 200		408	500	µg/g	82%	70 - 130	
Terpineol	<LOQ	< 200		506	500	µg/g	101%	70 - 130	
Nerol	<LOQ	< 200		463	500	µg/g	93%	70 - 130	
Pulegone	<LOQ	< 200		524	500	µg/g	105%	70 - 130	
Geraniol	<LOQ	< 200		586	500	µg/g	117%	70 - 130	
Geranyl_Acetate	<LOQ	< 200		410	500	µg/g	82%	70 - 130	
a-Cedrene	<LOQ	< 200		478	500	µg/g	96%	70 - 130	
b-Caryophyllene	<LOQ	< 200		392	500	µg/g	78%	70 - 130	
a-Humulene	<LOQ	< 200		499	500	µg/g	100%	70 - 130	
Valenene	<LOQ	< 200		400	500	µg/g	80%	70 - 130	
cis-Nerolidol	<LOQ	< 200		459	500	µg/g	92%	70 - 130	
a-Farnesene	<LOQ	< 200		478	500	µg/g	96%	70 - 130	
trans-Nerolidol	<LOQ	< 200		515	500	µg/g	103%	70 - 130	
Caryophyllene_Oxide	<LOQ	< 200		408	500	µg/g	82%	70 - 130	
Guaiol	<LOQ	< 200		518	500	µg/g	104%	70 - 130	
Cedrol	<LOQ	< 200		417	500	µg/g	83%	70 - 130	
a-Bisabolol	<LOQ	< 200		430	500	µg/g	86%	70 - 130	

Definitions

LOQ	Limit of Quantitation
LCS	Laboratory Control Sample
% REC	Percent Recovery



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503-254-1794



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Terpenes Quality Control Results

Method Reference: EPA 5035			Batch ID: 2210177				
Sample/Sample Duplicate			Sample ID: 22-014396-0001				
Analyte	Result	Org. Result	LOQ	Units	% RPD	LIMIT	Notes
a-pinene	<LOQ	<LOQ	191	µg/g	0%	< 20	
Camphene	<LOQ	<LOQ	191	µg/g	0%	< 20	
Sabinene	<LOQ	<LOQ	191	µg/g	0%	< 20	
b-Pinene	<LOQ	<LOQ	191	µg/g	0%	< 20	
b-Myrcene	<LOQ	<LOQ	191	µg/g	0%	< 20	
a-phellandrene	<LOQ	<LOQ	191	µg/g	0%	< 20	
d-3-Carene	<LOQ	<LOQ	191	µg/g	0%	< 20	
a-Terpinene	<LOQ	<LOQ	191	µg/g	0%	< 20	
p-Cymene	<LOQ	<LOQ	191	µg/g	0%	< 20	
D-Limonene	1760	1740	191	µg/g	1%	< 20	
Eucalyptol	<LOQ	<LOQ	191	µg/g	0%	< 20	
b-cis-Ocimene	<LOQ	<LOQ	63.5	µg/g	0%	< 20	
b-trans-Ocimene	<LOQ	<LOQ	127	µg/g	0%	< 20	
g-Terpinene	<LOQ	<LOQ	191	µg/g	0%	< 20	
Sabinene_Hydrate	<LOQ	<LOQ	191	µg/g	0%	< 20	
Terpinolene	<LOQ	<LOQ	191	µg/g	0%	< 20	
D-Fenchone	<LOQ	<LOQ	191	µg/g	0%	< 20	
Linalool	384	383	191	µg/g	0%	< 20	
Fenchol	<LOQ	<LOQ	191	µg/g	0%	< 20	
Camphor	<LOQ	<LOQ	191	µg/g	0%	< 20	
Isopulego	<LOQ	<LOQ	191	µg/g	0%	< 20	
Isoborneol	<LOQ	<LOQ	191	µg/g	0%	< 20	
Borneol	<LOQ	<LOQ	191	µg/g	0%	< 20	
DL-Menthol	<LOQ	<LOQ	191	µg/g	0%	< 20	
Terpineol	<LOQ	<LOQ	191	µg/g	0%	< 20	
Nerol	<LOQ	<LOQ	191	µg/g	0%	< 20	
Pulegone	<LOQ	<LOQ	191	µg/g	0%	< 20	
Gereniol	<LOQ	<LOQ	191	µg/g	0%	< 20	
Geranyl_Acetate	<LOQ	<LOQ	191	µg/g	0%	< 20	
a-Cedrene	306	302	191	µg/g	1%	< 20	
b-Caryophyllene	<LOQ	<LOQ	191	µg/g	0%	< 20	
a-Humulene	<LOQ	<LOQ	191	µg/g	0%	< 20	
Valenene	<LOQ	<LOQ	191	µg/g	0%	< 20	
cis-Nerolidol	<LOQ	<LOQ	191	µg/g	0%	< 20	
a-Farnesene	<LOQ	<LOQ	191	µg/g	0%	< 20	
trans-Nerolidol	<LOQ	<LOQ	191	µg/g	0%	< 20	
Caryophyllene_Oxide	<LOQ	<LOQ	191	µg/g	0%	< 20	
Guaiol	<LOQ	<LOQ	191	µg/g	0%	< 20	
Cedrol	193	192	191	µg/g	1%	< 20	
a-Bisabolol	<LOQ	<LOQ	191	µg/g	0%	< 20	

Definitions

RPD Relative Percent Difference



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Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662		Units: mg/Kg		Batch ID: 2210196				
Method Blank				Laboratory Control Sample				
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes
Abamectin	0.000	< 0.250		1.067	1.000	106.7	50.0	150
Acephate	0.000	< 0.200		0.765	0.800	95.6	60.0	120
Acequinocyl	0.000	< 1.000		4.080	4.000	102.0	40.0	160
Acetamiprid	0.000	< 0.100		0.412	0.400	103.1	60.0	120
Aldicarb	0.000	< 0.200		0.870	0.800	108.7	60.0	120
Azoxystrobin	0.000	< 0.100		0.404	0.400	100.9	60.0	120
Bifenazate	0.000	< 0.100		0.398	0.400	99.5	60.0	120
Bifenthrin	0.000	< 0.100		0.408	0.400	102.1	50.0	150
Boscalid	0.000	< 0.200		0.878	0.800	109.7	60.0	120
Carbaryl	0.000	< 0.100		0.435	0.400	108.7	60.0	120
Carbofuran	0.000	< 0.100		0.449	0.400	112.3	60.0	120
Chlorantraniliprole	0.000	< 0.100		0.388	0.400	97.0	60.0	120
Chlorfenvapir	0.000	< 0.500		1.954	2.000	97.7	60.0	120
Chlorpyrifos	0.000	< 0.100		0.389	0.400	97.2	60.0	120
Clofentezine	0.000	< 0.100		0.430	0.400	107.6	60.0	120
Cyfluthrin	0.000	< 0.500		2.332	2.000	116.6	50.0	150
Cypermethrin	0.000	< 0.500		2.009	2.000	100.4	50.0	150
Daminozide	0.225	< 0.500		1.589	2.000	79.5	60.0	120
Diazinon	0.000	< 0.100		0.416	0.400	104.0	60.0	120
Dichlorvos	0.000	< 0.500		2.064	2.000	103.2	60.0	120
Dimethoate	0.000	< 0.100		0.427	0.400	106.8	60.0	120
Ethoprophos	0.000	< 0.100		0.432	0.400	108.1	60.0	120
Etofenprox	0.000	< 0.200		0.817	0.800	102.2	50.0	150
Etiozazole	0.000	< 0.100		0.409	0.400	102.2	60.0	120
Fenoxycarb	0.000	< 0.100		0.418	0.400	104.6	60.0	120
Fenpyroximate	0.000	< 0.200		0.809	0.800	101.1	60.0	120
Fipronil	0.000	< 0.200		0.900	0.800	112.5	60.0	120
Flonicamid	0.000	< 0.250		1.180	1.000	118.0	60.0	120
Fludioxonil	0.000	< 0.200		0.852	0.800	106.5	50.0	150
Hexythiazox	0.000	< 0.250		1.027	1.000	102.7	60.0	120
Imazalil	0.000	< 0.100		0.400	0.400	100.1	60.0	120
Imidacloprid	0.000	< 0.200		0.724	0.800	90.6	60.0	120
Kresoxim-methyl	0.000	< 0.200		0.849	0.800	106.1	60.0	120
Malathion	0.000	< 0.100		0.413	0.400	103.4	60.0	120
Metaxaryl	0.000	< 0.100		0.421	0.400	105.3	60.0	120
Methiocarb	0.000	< 0.100		0.425	0.400	106.4	60.0	120
Methomyl	0.000	< 0.200		0.919	0.800	114.9	60.0	120
MGK-264	0.000	< 0.100		0.483	0.400	120.9	50.0	150
Myclobutanil	0.000	< 0.100		0.397	0.400	99.2	60.0	120
Naled	0.000	< 0.250		1.050	1.000	105.0	50.0	150
Oxamyl	0.000	< 0.500		2.050	2.000	102.5	60.0	120
Paclobutrazole	0.000	< 0.200		0.853	0.800	106.6	60.0	120
Parathion-Methyl	0.000	< 0.100		0.504	0.400	126.0	50.0	150
Permethrin	0.000	< 0.100		0.428	0.400	106.9	50.0	150
Phosmet	0.000	< 0.100		0.423	0.400	105.8	50.0	150
Piperonyl butoxide	0.000	< 0.500		2.045	2.000	102.2	60.0	120
Prallethrin	0.000	< 0.100		0.426	0.400	106.4	60.0	120
Propiconazole	0.000	< 0.200		0.836	0.800	104.5	60.0	120
Propoxur	0.000	< 0.100		0.438	0.400	109.5	60.0	120
Pyrethrin (Summe)	0.000	< 0.100		0.512	0.488	104.9	60.0	120
Pyridaben	0.000	< 0.100		0.406	0.400	101.6	50.0	150
Spinosad	0.000	< 0.100		0.394	0.388	101.5	50.0	150
Spiromesifen	0.000	< 0.100		0.416	0.400	103.9	60.0	120
Spirotetramat	0.000	< 0.100		0.401	0.400	100.1	60.0	120
Spiroxamine	0.000	< 0.200		0.787	0.800	98.4	60.0	120
Tebuconazole	0.000	< 0.200		0.849	0.800	106.2	60.0	120
Thiacloprid	0.000	< 0.100		0.400	0.400	100.1	60.0	120
Thiamethoxam	0.000	< 0.100		0.439	0.400	109.8	60.0	120
Trifloxystrobin	0.000	< 0.100		0.410	0.400	102.4	60.0	120



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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: 2210196										
Matrix Spike/Matrix Spike Duplicate Recoveries					Sample ID: 22-014208-0001					
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes
Abamectin	0.000	1.020	1.117	1.000	9.1%	< 30	102.0%	111.7%	50 - 150	
Acephate	0.010	0.702	0.644	0.800	8.8%	< 30	86.5%	79.2%	50 - 150	
Acequinocyl	0.000	3.851	4.092	4.000	6.1%	< 30	96.3%	102.3%	50 - 150	
Acetamiprid	0.000	0.421	0.430	0.400	2.1%	< 30	105.3%	107.5%	50 - 150	
Aldicarb	0.000	0.893	0.910	0.800	2.0%	< 30	111.6%	113.8%	50 - 150	
Azoxystrobin	0.000	0.408	0.401	0.400	1.7%	< 30	102.0%	100.3%	50 - 150	
Bifenazate	0.000	0.546	0.571	0.400	4.4%	< 30	136.6%	142.7%	50 - 150	
Bifenthrin	0.000	0.378	0.367	0.400	2.9%	< 30	94.5%	91.9%	50 - 150	
Boscalid	0.000	0.889	0.832	0.800	6.6%	< 30	111.2%	104.1%	50 - 150	
Carbaryl	0.000	0.435	0.420	0.400	3.3%	< 30	108.7%	105.1%	50 - 150	
Carbofuran	0.000	0.449	0.451	0.400	0.4%	< 30	112.3%	112.8%	50 - 150	
Chlorantraniliprole	0.000	0.403	0.403	0.400	0.0%	< 30	100.7%	100.7%	50 - 150	
Chlorfenvapir	0.000	1.874	1.805	2.000	3.7%	< 30	93.7%	90.3%	50 - 150	
Chlorpyrifos	0.000	0.388	0.361	0.400	7.1%	< 30	96.9%	90.2%	50 - 150	
Clofentazine	0.000	0.381	0.387	0.400	1.4%	< 30	95.3%	96.7%	50 - 150	
Cyfluthrin	0.109	1.384	1.246	2.000	11.5%	< 30	63.7%	56.8%	30 - 150	
Cypermethrin	0.000	0.897	0.804	2.000	11.0%	< 30	44.9%	40.2%	50 - 150	Q
Daminozide	0.299	1.880	1.847	2.000	2.1%	< 30	79.1%	77.4%	30 - 150	
Diazinon	0.000	0.373	0.386	0.400	3.3%	< 30	93.3%	96.5%	50 - 150	
Dichlorvos	0.000	2.165	2.074	2.000	4.3%	< 30	108.2%	103.7%	50 - 150	
Dimethoate	0.000	0.447	0.447	0.400	0.1%	< 30	111.7%	111.8%	50 - 150	
Ethoprophos	0.000	0.406	0.435	0.400	6.9%	< 30	101.4%	108.7%	50 - 150	
Etofenprox	0.000	0.799	0.770	0.800	3.6%	< 30	99.9%	96.3%	50 - 150	
Etoxazole	0.000	0.395	0.410	0.400	3.7%	< 30	98.8%	102.6%	50 - 150	
Fenoxycarb	0.000	0.424	0.444	0.400	4.6%	< 30	106.1%	111.0%	50 - 150	
Fenpyroximate	0.000	0.477	0.488	0.800	2.3%	< 30	59.6%	61.0%	50 - 150	
Fipronil	0.000	0.835	0.798	0.800	4.5%	< 30	104.3%	99.7%	50 - 150	
Fonicamid	0.000	1.168	1.228	1.000	5.0%	< 30	116.8%	122.8%	50 - 150	
Fludioxonil	0.000	0.954	0.989	0.800	3.6%	< 30	119.3%	123.6%	50 - 150	
Hexythiazox	0.000	1.177	1.218	1.000	3.4%	< 30	117.7%	121.8%	50 - 150	
Imazalil	0.000	0.429	0.430	0.400	0.3%	< 30	107.2%	107.5%	50 - 150	
Imidacloprid	0.000	0.734	0.759	0.800	3.4%	< 30	91.8%	94.9%	50 - 150	
Kresoxim-methyl	0.000	0.855	0.893	0.800	4.4%	< 30	106.8%	111.6%	50 - 150	
Malathion	0.000	0.423	0.440	0.400	4.1%	< 30	105.7%	110.1%	50 - 150	
Metaxyl	0.000	0.427	0.451	0.400	5.6%	< 30	106.7%	112.8%	50 - 150	
Methiocarb	0.000	0.431	0.435	0.400	1.0%	< 30	107.6%	108.7%	50 - 150	
Methomyl	0.000	0.874	0.920	0.800	5.2%	< 30	109.2%	115.1%	50 - 150	
MGK-264	0.000	0.407	0.430	0.400	5.6%	< 30	101.7%	107.5%	50 - 150	
Myclobutanil	0.000	0.365	0.377	0.400	3.3%	< 30	91.3%	94.3%	50 - 150	
Naled	0.000	1.051	1.034	1.000	1.7%	< 30	105.1%	103.4%	50 - 150	
Oxamyl	0.000	2.325	2.539	2.000	8.8%	< 30	116.2%	127.0%	50 - 150	
Paclobutrazole	0.000	0.826	0.838	0.800	1.5%	< 30	103.3%	104.8%	50 - 150	
Parathion-Methyl	0.000	0.494	0.515	0.400	4.1%	< 30	123.6%	128.7%	30 - 150	
Permethrin	0.000	0.376	0.364	0.400	3.2%	< 30	94.0%	91.0%	50 - 150	
Phosmet	0.000	0.451	0.460	0.400	1.8%	< 30	112.9%	114.9%	50 - 150	
Piperonyl butoxide	0.000	2.171	2.285	2.000	5.1%	< 30	108.5%	114.2%	50 - 150	
Prallethrin	0.025	0.523	0.533	0.400	1.9%	< 30	124.6%	127.1%	50 - 150	
Propiconazole	0.000	1.136	1.181	0.800	3.9%	< 30	142.1%	147.7%	50 - 150	
Propoxur	0.000	0.440	0.433	0.400	1.7%	< 30	110.1%	108.2%	50 - 150	
Pyrethrin (Summe)	0.044	0.674	0.725	0.488	7.8%	< 30	129.0%	139.5%	50 - 150	
Pyridaben	0.000	0.484	0.486	0.400	0.5%	< 30	120.9%	121.6%	50 - 150	
Spirosad	0.000	0.379	0.349	0.388	8.2%	< 30	97.6%	89.9%	50 - 150	
Spiromesifen	0.000	0.445	0.434	0.400	2.6%	< 30	111.3%	108.4%	50 - 150	
Spirotetramat	0.000	0.451	0.453	0.400	0.4%	< 30	112.7%	113.2%	50 - 150	
Spiroxamine	0.000	0.835	0.870	0.800	4.1%	< 30	104.4%	108.7%	50 - 150	
Tebuconazole	0.000	0.863	0.875	0.800	1.4%	< 30	107.8%	109.3%	50 - 150	
Thiacloprid	0.000	0.426	0.433	0.400	1.6%	< 30	106.4%	108.2%	50 - 150	
Thiamethoxam	0.000	0.454	0.488	0.400	7.2%	< 30	113.5%	121.9%	50 - 150	
Trifloxystrobin	0.000	0.401	0.387	0.400	3.6%	< 30	100.2%	96.6%	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.