



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



Report Number: 24-005230/D001.R002
Report Date: 06/26/2024
ORELAP#: OR100028
Purchase Order:
Received: 05/13/24 11:26

This is an amended version of report# 24-005230/D001.R001.
Reason: updated client information

Customer: Mary & Jane
Product identity: Sunny Melts
Client/Metric ID: Lot: 0099-04-001
Laboratory ID: 24-005230-0001

Summary

Potency:

Analyte	Result	Limits	Units	Status	
Δ9-THC	0.173		%		Delta-9-THC-Total per 0.952 mg/0.55g
Analyte per 0.55g	Result	Limits	Units	Status	CBD-Total per Serving Size <LOQ
Δ9-THC per 0.55g	0.952		mg/0.55g		(Reported in milligrams per serving)

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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Customer: Mary & Jane
2256 Terminal Road
Saint Paul Minnesota 55113
United States of America (USA)

Product identity: Sunny Melts
Client/Metric ID: Lot: 0099-04-001
Sample Date:
Laboratory ID: 24-005230-0001
Evidence of Cooling: No
Temp: 20.3 °C
Relinquished by: client
Serving Size #1: 0.55 g

Sample Results

Potency	Method: J AOAC 2015 V98-6 (mod) ^b	Units %	Batch: 2403711	Analyze: 5/14/24 6:25:00 PM	
Analyte	Result	Limits	Units	LOQ	Notes
CBC	< LOQ		%	0.00313	
CBC-A	< LOQ		%	0.00313	
CBC-Total	< LOQ		%	0.00588	
CBD [±]	< LOQ		%	0.00313	
CBD-A [±]	< LOQ		%	0.00313	
CBD-Total [±]	< LOQ		%	0.00588	
CBDV	< LOQ		%	0.00313	
CBDV-A	< LOQ		%	0.00313	
CBDV-Total	< LOQ		%	0.00585	
CBE	< LOQ		%	0.00313	
CBG	< LOQ		%	0.00313	
CBG-A	< LOQ		%	0.00313	
CBG-Total	< LOQ		%	0.00585	
CBL	< LOQ		%	0.00313	
CBL-A	< LOQ		%	0.00313	
CBL-Total	< LOQ		%	0.00588	
CBN	< LOQ		%	0.00313	
CBT	< LOQ		%	0.00313	
Δ10-THC-9R	< LOQ		%	0.00313	
Δ10-THC-9S	< LOQ		%	0.00313	
Δ10-THC-Total	< LOQ		%	0.00627	
Δ8-THC [±]	< LOQ		%	0.00313	
Δ8-THCV	< LOQ		%	0.00313	
Δ9-THC [±]	0.173		%	0.00313	
Δ9-THC-A [±]	< LOQ		%	0.00313	
Δ9-THC-Total [±]	0.173		%	0.00588	
Δ9-THCP	< LOQ		%	0.00313	
Δ9-THCV	< LOQ		%	0.00313	
Δ9-THCV-A	< LOQ		%	0.00313	
Δ9-THCV-Total	< LOQ		%	0.00585	



Potency	Method: J AOAC 2015 V98-6 (mod) ^b	Units %	Batch: 2403711	Analyze: 5/14/24 6:25:00 PM	
Analyte	Result	Limits	Units	LOQ	Notes
exo-THC	< LOQ		%	0.00313	
Total Cannabinoids	0.173		%		

Potency per 0.55g	Method: J AOAC 2015 V98-6 (mod) ^b	Units mg/se	Batch: 2403711	Analyze: 5/14/24 6:25:00 PM	
Analyte	Result	Limits	Units	LOQ	Notes
CBC per 0.55g	< LOQ		mg/0.55g	0.0172	
CBC-A per 0.55g	< LOQ		mg/0.55g	0.0172	
CBC-Total per 0.55g	< LOQ		mg/0.55g	0.0323	
CBD per 0.55g	< LOQ		mg/0.55g	0.0172	
CBD-A per 0.55g [±]	< LOQ		mg/0.55g	0.0172	
CBD-Total per 0.55g [±]	< LOQ		mg/0.55g	0.0323	
CBDV per 0.55g	< LOQ		mg/0.55g	0.0172	
CBDV-A per 0.55g	< LOQ		mg/0.55g	0.0172	
CBDV-Total per 0.55g	< LOQ		mg/0.55g	0.0322	
CBE per 0.55g	< LOQ		mg/0.55g	0.0172	
CBG per 0.55g	< LOQ		mg/0.55g	0.0172	
CBG-A per 0.55g	< LOQ		mg/0.55g	0.0172	
CBG-Total per 0.55g	< LOQ		mg/0.55g	0.0322	
CBL per 0.55g	< LOQ		mg/0.55g	0.0172	
CBL-A per 0.55g	< LOQ		mg/0.55g	0.0172	
CBL-Total per 0.55g	< LOQ		mg/0.55g	0.0323	
CBN per 0.55g	< LOQ		mg/0.55g	0.0172	
CBT per 0.55g	< LOQ		mg/0.55g	0.0172	
Δ10-THC-9R per 0.55g	< LOQ		mg/0.55g	0.0172	
Δ10-THC-9S per 0.55g	< LOQ		mg/0.55g	0.0172	
Δ10-THC-Total per 0.55g	< LOQ		mg/0.55g	0.0345	
Δ8-THC per 0.55g [±]	< LOQ		mg/0.55g	0.0172	
Δ8-THCV per 0.55g	< LOQ		mg/0.55g	0.0172	
Δ9-THC per 0.55g [±]	0.952		mg/0.55g	0.0172	
Δ9-THC-Total per 0.55g	0.952		mg/0.55g	0.0323	
Δ9-THCP per 0.55g	< LOQ		mg/0.55g	0.0172	
Δ9-THCV per 0.55g	< LOQ		mg/0.55g	0.0172	
Δ9-THCV-A per 0.55g	< LOQ		mg/0.55g	0.0172	
Δ9-THCV-Total per 0.55g	< LOQ		mg/0.55g	0.0324	
exo-THC per 0.55g	< LOQ		mg/0.55g	0.0172	
THC-A per 0.55g [±]	< LOQ		mg/0.55g	0.0172	
Total Cannabinoids per 0.55g	0.952		mg/0.55g		

Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2403651	05/16/24 AOAC 991.14 (Petrifilm)		
Total Coliforms	< LOQ		cfu/g	10	2403651	05/16/24 AOAC 991.14 (Petrifilm)		
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2403652	05/17/24 AOAC 2014.05 (RAPID)		
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2403652	05/17/24 AOAC 2014.05 (RAPID)		



Solvents		Method: Residual Solvents by HS-GC-MS ^b				Units µg/g	Batch 2403762	Analyze 05/16/24 12:37 PM			
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
1,4-Dioxane [⊥]	< LOQ		100			2-Butanol [⊥]	< LOQ		200		
2-Ethoxyethanol [⊥]	< LOQ		30.0			2-Methylbutane (Isopentane) [⊥]	< LOQ		200		
2-Methylpentane [⊥]	< LOQ		30.0			2-Propanol (IPA) [⊥]	< LOQ		200		
2,2-Dimethylbutane [⊥]	< LOQ		30.0			2,2-Dimethylpropane (neo-pentane) [⊥]	< LOQ		200		
2,3-Dimethylbutane [⊥]	< LOQ		30.0			3-Methylpentane [⊥]	< LOQ		30.0		
Acetone [⊥]	< LOQ		200			Acetonitrile [⊥]	< LOQ		100		
Benzene [⊥]	< LOQ		1.00			Butanes (sum) [⊥]	< LOQ		400		
Cyclohexane [⊥]	< LOQ		200			Ethyl acetate [⊥]	< LOQ		200		
Ethyl benzene	< LOQ		200			Ethyl ether [⊥]	< LOQ		200		
Ethylene glycol [⊥]	< LOQ		200			Ethylene oxide [⊥]	< LOQ		20.0		
Hexanes (sum) [⊥]	< LOQ		150			Isopropyl acetate [⊥]	< LOQ		200		
Isopropylbenzene (Cumene) [⊥]	< LOQ		30.0			m,p-Xylene [⊥]	< LOQ		200		
Methanol [⊥]	< LOQ		200			Methylene chloride [⊥]	< LOQ		60.0		
Methylpropane (Isobutane) [⊥]	< LOQ		200			n-Butane [⊥]	< LOQ		200		
n-Heptane [⊥]	< LOQ		200			n-Hexane [⊥]	< LOQ		30.0		
n-Pentane [⊥]	< LOQ		200			o-Xylene [⊥]	< LOQ		200		
Pentanes (sum)	< LOQ		600			Propane	< LOQ		200		
Tetrahydrofuran [⊥]	< LOQ		100			Toluene [⊥]	< LOQ		100		
Total Xylenes [⊥]	< LOQ		400			Total Xylenes and Ethyl benzene	< LOQ		600		



Pesticides		Method: AOAC 2007.01 & EN 15662 (mod)				Units mg/kg	Batch 2403785	Analyze 05/17/24 09:01 AM			
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
Abamectin [±]	< LOQ		0.250			Acephate	< LOQ		0.200		
Acequinocyl [±]	< LOQ		1.00			Acetamiprid	< LOQ		0.100		
Aldicarb [±]	< LOQ		0.200			Azoxystrobin [±]	< LOQ		0.100		
Bifenazate [±]	< LOQ		0.100			Bifenthrin [±]	< LOQ		0.100		
Boscalid [±]	< LOQ		0.200			Carbaryl [±]	< LOQ		0.100		
Carbofuran [±]	< LOQ		0.100			Chlorantraniliprole [±]	< LOQ		0.100		
Chlorfenapyr [±]	< LOQ		0.500			Chlorpyrifos-ethyl [±]	< LOQ		0.100		
Clofentezine [±]	< LOQ		0.100			Cyfluthrin (sum) [±]	< LOQ		0.500		
Cypermethrin and	< LOQ		0.500			Daminozide [±]	< LOQ		0.500		
Diazinon [±]	< LOQ		0.100			Dichlorvos [±]	< LOQ		0.500		
Dimethoate [±]	< LOQ		0.100			Ethoprophos [±]	< LOQ		0.100		
Etofenprox [±]	< LOQ		0.200			Etoxazole [±]	< LOQ		0.100		
Fenoxycarb [±]	< LOQ		0.100			Fenpyroximate [±]	< LOQ		0.200		
Fipronil [±]	< LOQ		0.200			Flonicamid [±]	< LOQ		0.400		
Fludioxonil [±]	< LOQ		0.200			Hexythiazox [±]	< LOQ		0.400		
Imazalil [±]	< LOQ		0.100			Imidacloprid [±]	< LOQ		0.200		
Kresoxim-methyl [±]	< LOQ		0.200			Malathion [±]	< LOQ		0.100		
Metalaxyl [±]	< LOQ		0.100			Methiocarb [±]	< LOQ		0.100		
Methomyl [±]	< LOQ		0.200			MGK-264 [±]	< LOQ		0.100		
Myclobutanil [±]	< LOQ		0.100			Naled [±]	< LOQ		0.250		
Oxamyl [±]	< LOQ		0.500			Paclobutrazole [±]	< LOQ		0.200		
Parathion-Methyl [±]	< LOQ		0.100			Permethrin and isomers	< LOQ		0.100		
Phosmet [±]	< LOQ		0.100			Piperonyl butoxide [±]	< LOQ		1.00		
Prallethrin [±]	< LOQ		0.100			Propiconazole [±]	< LOQ		0.200		
Propoxur [±]	< LOQ		0.100			Pyrethrin I (total) [±]	< LOQ		0.500		
Pyridaben [±]	< LOQ		0.100			Spinosad [±]	< LOQ		0.100		
Spiromesifen [±]	< LOQ		0.100			Spirotetramat [±]	< LOQ		0.100		
Spiroxamine [±]	< LOQ		0.200			Tebuconazole [±]	< LOQ		0.200		
Thiacloprid [±]	< LOQ		0.100			Thiamethoxam [±]	< LOQ		0.100		
Trifloxystrobin [±]	< LOQ		0.100								

Metals											
Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method		Status	Notes		
Arsenic [±]	< LOQ		mg/kg	0.0200	2403796	05/17/24	AOAC 2013.06 (mod.) ^p				
Cadmium [±]	< LOQ		mg/kg	0.0200	2403796	05/17/24	AOAC 2013.06 (mod.) ^p				
Lead [±]	< LOQ		mg/kg	0.0200	2403796	05/17/24	AOAC 2013.06 (mod.) ^p				
Mercury [±]	< LOQ		mg/kg	0.0100	2403796	05/17/24	AOAC 2013.06 (mod.) ^p				



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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.

^p = ISO/IEC 17025:2017 accredited method.

[⊥] = TNI accredited analyte.

Units of Measure

cfu/g = Colony forming units per gram

g = Gram

µg/g = Microgram per gram

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

mg/0.55g = Milligram per 0.55g

% = Percentage of sample

% wt = µg/g divided by 10,000

Approved Signatory

Derrick Tanner
General Manager



12423 NE Whitaker Way
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Hemp & Cannabis
 Chain of Custody

Minsky-&-Paul-1715619361

ORELAP ID: OR1000028 ANAB ISO 17025 ID: AT1508

Company Details Company: <u>Minsky & Paul</u> Contact: <u>Laura Roos</u> Street Address: <u>2256 Terminal Road</u> City, State, Zip: <u>St. Paul, under ned 55113</u> Email: <u>laura@minnyandpaul.com</u> Contact Phone: <u>9522123349</u> Company Phone: <u>9522123349</u> Billing Information Billing Phone: <u>9522123349</u> Billing Email: <u>laura@minnyandpaul.com</u>			Project Details Turnaround Time: <u>3 Business Days Surcharges Apply</u> Relinquishment Sampling, Courier & Shipping Options: <u>By Shipping Service (USPS, UPS, FedEx)</u> Compliance: <u>Compliance</u> Receipt Information			Testing H10010 - Potency Cannabis (Basic+Expanded)
#	Sample Name	Material	Amount Provided	Reporting Unit	Speci cations	
1	Sunny Melts	Cannabinoid Edible	22 each	mg/g	Tablets are .495g	

Relinquished By	Date	Time	Received By	Date	Time	Received Temp., °C	Evidence of Cooling?
<i>Laura Roos</i>	<i>05/13/2024</i>	<i>09:56</i>	<i>rls</i>	<i>05/13/2024</i>	<i>11:26</i>	<i>20.30</i>	<i>No</i>

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

Columbia Laboratories
 12423 NE Whitaker Way
 Portland, OR 97230

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

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www.columbiaboratories.com



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Purchase Order:
Received: 05/13/24 11:26

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

Laboratory Quality Control Results

J AOAC 2015 V98-6 Batch ID: 2403711

Laboratory Control Sample

Analyte	LCS	Result	Spike	Units	% Rec	Limits	Evaluation	Notes
CBDVA	2	0.0366	0.0358	%	102	80.0 - 120	Acceptable	
CBDV	2	0.0345	0.0340	%	101	80.0 - 120	Acceptable	
CBE	2	0.0316	0.0315	%	100	80.0 - 120	Acceptable	
CBDA	1	0.0310	0.0309	%	100	90.0 - 110	Acceptable	
CBGA	1	0.0307	0.0304	%	101	80.0 - 120	Acceptable	
CBG	1	0.0345	0.0344	%	100	80.0 - 120	Acceptable	
CBD	1	0.0339	0.0335	%	101	90.0 - 110	Acceptable	
THCV	2	0.0342	0.0336	%	102	80.0 - 120	Acceptable	
d8THCV	2	0.0361	0.0360	%	100	80.0 - 120	Acceptable	
THCVA	2	0.0324	0.0318	%	102	80.0 - 120	Acceptable	
CBN	1	0.0323	0.0320	%	101	80.0 - 120	Acceptable	
exo-THC	2	0.0331	0.0331	%	100	80.0 - 120	Acceptable	
d9THC	1	0.0330	0.0324	%	102	90.0 - 110	Acceptable	
d8THC	1	0.0321	0.0323	%	99.4	90.0 - 110	Acceptable	
9S-d10THC	1	0.0323	0.0322	%	100	80.0 - 120	Acceptable	
CBL	2	0.0328	0.0331	%	99.1	80.0 - 120	Acceptable	
9R-d10THC	1	0.0312	0.0314	%	99.5	80.0 - 120	Acceptable	
CBC	2	0.0347	0.0349	%	99.6	80.0 - 120	Acceptable	
THCA	1	0.0312	0.0304	%	103	90.0 - 110	Acceptable	
CBCA	2	0.0339	0.0336	%	101	80.0 - 120	Acceptable	
CBLA	2	0.0341	0.0339	%	101	80.0 - 120	Acceptable	
d9THCP	2	0.0322	0.0324	%	99.3	80.0 - 120	Acceptable	
CBT	2	0.0337	0.0351	%	96.0	80.0 - 120	Acceptable	

Method Blank

Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.00328	%	< 0.00328	Acceptable	
CBDV	<LOQ	0.00328	%	< 0.00328	Acceptable	
CBE	<LOQ	0.00328	%	< 0.00328	Acceptable	
CBDA	<LOQ	0.00328	%	< 0.00328	Acceptable	
CBGA	<LOQ	0.00328	%	< 0.00328	Acceptable	
CBG	<LOQ	0.00328	%	< 0.00328	Acceptable	
CBD	<LOQ	0.00328	%	< 0.00328	Acceptable	
THCV	<LOQ	0.00328	%	< 0.00328	Acceptable	
d8THCV	<LOQ	0.00328	%	< 0.00328	Acceptable	
THCVA	<LOQ	0.00328	%	< 0.00328	Acceptable	
CBN	<LOQ	0.00328	%	< 0.00328	Acceptable	
exo-THC	<LOQ	0.00328	%	< 0.00328	Acceptable	
d9THC	<LOQ	0.00328	%	< 0.00328	Acceptable	
d8THC	<LOQ	0.00328	%	< 0.00328	Acceptable	
9S-d10THC	<LOQ	0.00328	%	< 0.00328	Acceptable	
CBL	<LOQ	0.00328	%	< 0.00328	Acceptable	
9R-d10THC	<LOQ	0.00328	%	< 0.00328	Acceptable	
CBC	<LOQ	0.00328	%	< 0.00328	Acceptable	
THCA	<LOQ	0.00328	%	< 0.00328	Acceptable	
CBCA	<LOQ	0.00328	%	< 0.00328	Acceptable	
CBLA	<LOQ	0.00328	%	< 0.00328	Acceptable	
d9THCP	<LOQ	0.00328	%	< 0.00328	Acceptable	
CBT	<LOQ	0.00328	%	< 0.00328	Acceptable	

Abbreviations

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

Units of Measure:

% - Percent



Laboratory Quality Control Results

AOAC 2015 V98-6		Batch ID: 2403711						
Sample Duplicate		Sample ID: 24-005147-0002						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
CBDV	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
CBE	0.0243	0.0246	0.00317	%	1.46	< 20	Acceptable	
CBDA	0.0666	0.0677	0.00317	%	1.64	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
CBG	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
CBD	0.162	0.165	0.00317	%	1.63	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
CBN	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
d9THC	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
d8THC	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
9S-d10THC	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
9R-d10THC	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
CBC	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
CBCA	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
d9THCP	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	
CBT	<LOQ	<LOQ	0.00317	%	NA	< 20	Acceptable	

Abbreviations

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

Units of Measure:

% - Percent


Laboratory Quality Control Results

Residual Solvents				Batch ID: 2403762					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		460	584	µg/g	78.8	60 - 120	
Isobutane	ND	< 200		618	767	µg/g	80.6	60 - 120	
Butane	ND	< 200		646	782	µg/g	82.6	60 - 120	
2,2-Dimethylpropane	ND	< 200		794	939	µg/g	84.6	60 - 120	
Methanol	ND	< 200		1140	1600	µg/g	71.3	60 - 120	
Ethylene Oxide	ND	< 30		47.8	57.1	µg/g	83.7	60 - 120	
2-Methylbutane	ND	< 200		1210	1600	µg/g	75.6	60 - 120	
Pentane	ND	< 200		1230	1610	µg/g	76.4	60 - 120	
Ethanol	ND	< 200		1230	1600	µg/g	76.9	70 - 130	
Ethyl Ether	ND	< 200		1250	1600	µg/g	78.1	60 - 120	
2,2-Dimethylbutane	ND	< 30		129	162	µg/g	79.6	60 - 120	
Acetone	ND	< 200		1200	1600	µg/g	75.0	60 - 120	
2-Propanol	ND	< 200		1250	1600	µg/g	78.1	60 - 120	
Ethyl Formate	ND	< 500		1350	1630	µg/g	82.8	70 - 130	
Acetonitrile	ND	< 100		379	487	µg/g	77.8	60 - 120	
Methyl Acetate	ND	< 500		1420	1610	µg/g	88.2	70 - 130	
2,3-Dimethylbutane	ND	< 30		123	161	µg/g	76.4	60 - 120	
Dichloromethane	ND	< 60		373	483	µg/g	77.2	60 - 120	
2-Methylpentane	ND	< 30		125	164	µg/g	76.2	60 - 120	
MTBE	ND	< 500		1450	1610	µg/g	90.1	70 - 130	
3-Methylpentane	ND	< 30		124	160	µg/g	77.5	60 - 120	
Hexane	ND	< 30		134	171	µg/g	78.4	60 - 120	
1-Propanol	ND	< 500		1470	1610	µg/g	91.3	70 - 130	
Methylethylketone	ND	< 500		1450	1610	µg/g	90.1	70 - 130	
Ethyl acetate	ND	< 200		1300	1620	µg/g	80.2	60 - 120	
2-Butanol	ND	< 200		1310	1600	µg/g	81.9	60 - 120	
Tetrahydrofuran	ND	< 100		387	481	µg/g	80.5	60 - 120	
Cyclohexane	ND	< 200		1370	1610	µg/g	85.1	60 - 120	
2-methyl-1-propanol	ND	< 500		1350	1610	µg/g	83.9	70 - 130	
Benzene	ND	< 1		4.32	5.17	µg/g	83.6	60 - 120	
Isopropyl Acetate	ND	< 200		1380	1600	µg/g	86.3	60 - 120	
Heptane	ND	< 200		1370	1620	µg/g	84.6	60 - 120	
1-Butanol	ND	< 500		1320	1610	µg/g	82.0	70 - 130	
Propyl Acetate	ND	< 500		1370	1610	µg/g	85.1	70 - 130	
1,4-Dioxane	ND	< 100		410	497	µg/g	82.5	60 - 120	
2-Ethoxyethanol	ND	< 30		130	160	µg/g	81.3	60 - 120	
Methylisobutylketone	ND	< 500		1320	1620	µg/g	81.5	70 - 130	
3-Methyl-1-butanol	ND	< 500		1250	1610	µg/g	77.6	70 - 130	
Ethylene Glycol	ND	< 200		355	483	µg/g	73.5	60 - 120	
Toluene	ND	< 100		411	482	µg/g	85.3	60 - 120	
Isobutyl Acetate	ND	< 500		1320	1620	µg/g	81.5	70 - 130	
1-Pentanol	ND	< 500		1240	1610	µg/g	77.0	70 - 130	
Butyl Acetate	ND	< 500		1360	1650	µg/g	82.4	70 - 130	
Ethylbenzene	ND	< 200		808	970	µg/g	83.3	60 - 120	
m,p-Xylene	ND	< 200		799	963	µg/g	83.0	60 - 120	
o-Xylene	ND	< 200		819	961	µg/g	85.2	60 - 120	
Cumene	ND	< 30		140	164	µg/g	85.4	60 - 120	
Anisole	ND	< 500		1120	1610	µg/g	69.6	70 - 130	Q6
DMSO	ND	< 500		1280	1610	µg/g	79.5	70 - 130	
1,2-dimethoxyethane	ND	< 50		151	170	µg/g	88.8	70 - 130	
Triethylamine	ND	< 500		1120	1620	µg/g	69.1	70 - 130	Q6
N,N-dimethylformamide	ND	< 150		424	499	µg/g	85.0	70 - 130	
N,N-dimethylacetamide	ND	< 150		333	489	µg/g	68.1	70 - 130	Q6
Pyridine	ND	< 50		120	167	µg/g	71.9	70 - 130	
Sulfolane	ND	< 50		97.9	169	µg/g	57.9	70 - 130	Q6
1,2-Dichloroethane	ND	< 1		0.822	1	µg/g	82.2	70 - 130	
Chloroform	ND	< 1		0.865	1	µg/g	86.5	70 - 130	
Trichloroethylene	ND	< 1		0.846	1	µg/g	84.6	70 - 130	
1,1-Dichloroethane	ND	< 1		0.851	1	µg/g	85.1	70 - 130	



Revision: 2 Document ID: 7087

Legacy ID: CFL-E33Effective:

QC - Sample Duplicate
Sample ID: 24-004985-0001-01

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



Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662		Units: mg/Kg			Batch ID: 2403785			
Method Blank		Laboratory Control Sample						
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes
Abamectin	0.000	< 0.250		0.802	1.000	80.2	50.0	150
Acephate	0.072	< 0.200		0.646	0.800	80.7	60.0	120
Acequinocyl	0.000	< 1.000		3.844	4.000	96.1	40.0	160
Acetamiprid	0.000	< 0.100		0.355	0.400	88.8	60.0	120
Aldicarb	0.000	< 0.200		0.762	0.800	95.3	60.0	120
Azoxystrobin	0.010	< 0.100		0.360	0.400	90.0	60.0	120
Bifenazate	0.000	< 0.100		0.361	0.400	90.2	60.0	120
Bifenthrin	0.000	< 0.100		0.379	0.400	94.8	50.0	150
Boscalid	0.000	< 0.200		0.742	0.800	92.8	60.0	120
Carbaryl	0.000	< 0.100		0.354	0.400	88.4	60.0	120
Carbofuran	0.000	< 0.100		0.335	0.400	83.8	60.0	120
Chlorantraniliprole	0.000	< 0.100		0.381	0.400	95.2	60.0	120
Chlorfenapyr	0.000	< 0.500		1.531	2.000	76.6	60.0	120
Chlorpyrifos	0.000	< 0.100		0.356	0.400	89.0	60.0	120
Clofentezine	0.000	< 0.100		0.328	0.400	82.1	60.0	120
Cyfluthrin	0.000	< 0.500		1.963	2.000	98.2	50.0	150
Cypermethrin	0.000	< 0.500		1.939	2.000	97.0	50.0	150
Daminozide	0.000	< 0.500		0.790	2.000	39.5	60.0	120
Diazinon	0.000	< 0.100		0.366	0.400	91.6	60.0	120
Dichlorvos	0.000	< 0.500		1.940	2.000	97.0	60.0	120
Dimethoate	0.000	< 0.100		0.367	0.400	91.8	60.0	120
Ethoprophos	0.000	< 0.100		0.369	0.400	92.4	60.0	120
Etofenprox	0.000	< 0.200		0.764	0.800	95.5	50.0	150
Etoazole	0.000	< 0.100		0.392	0.400	97.9	60.0	120
Fenoxycarb	0.000	< 0.100		0.354	0.400	88.6	60.0	120
Fenpyroximate	0.000	< 0.200		0.714	0.800	89.3	60.0	120
Fipronil	0.000	< 0.200		0.843	0.800	105.3	60.0	120
Fonicamid	0.000	< 0.250		0.920	1.000	92.0	60.0	120
Fludioxonil	0.000	< 0.200		0.746	0.800	93.3	50.0	150
Hexythiazox	0.000	< 0.250		0.894	1.000	89.4	60.0	120
Imazalil	0.000	< 0.100		0.368	0.400	92.1	60.0	120
Imidacloprid	0.000	< 0.200		0.705	0.800	88.1	60.0	120
Kresoxim-methyl	0.000	< 0.200		0.720	0.800	89.9	60.0	120
Malathion	0.000	< 0.100		0.360	0.400	90.0	60.0	120
Metalaxyl	0.000	< 0.100		0.366	0.400	91.6	60.0	120
Methiocarb	0.000	< 0.100		0.365	0.400	91.4	60.0	120
Methomyl	0.000	< 0.200		0.722	0.800	90.3	60.0	120
MGK-264	0.000	< 0.100		0.338	0.400	84.5	50.0	150
Myclobutanil	0.000	< 0.100		0.386	0.400	96.6	60.0	120
Naled	0.000	< 0.250		0.921	1.000	92.1	50.0	150
Oxamyl	0.000	< 0.500		2.071	2.000	103.6	60.0	120
Paclobutrazole	0.000	< 0.200		0.766	0.800	95.8	60.0	120
Parathion-Methyl	0.000	< 0.100		0.426	0.400	106.6	50.0	150
Permethrin	0.000	< 0.100		0.421	0.400	105.3	50.0	150
Phosmet	0.000	< 0.100		0.346	0.400	86.6	50.0	150
Piperonyl butoxide	0.000	< 0.500		1.973	2.000	98.7	60.0	120
Prallethrin	0.000	< 0.100		0.356	0.400	89.1	60.0	120
Propiconazole	0.000	< 0.200		0.738	0.800	92.3	60.0	120
Propoxur	0.000	< 0.100		0.366	0.400	91.6	60.0	120
Pyrethrin (Summe)	0.000	< 0.100		0.445	0.488	91.1	60.0	120
Pyridaben	0.000	< 0.100		0.387	0.400	96.8	50.0	150
Spinosad	0.000	< 0.100		0.367	0.388	94.5	50.0	150
Spiromesifen	0.000	< 0.100		0.337	0.400	84.2	60.0	120
Spirotetramat	0.000	< 0.100		0.356	0.400	89.1	60.0	120
Spiroxamine	0.000	< 0.200		0.723	0.800	90.4	60.0	120

Q7



Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662		Units: mg/Kg					Batch ID: 2403785			
Matrix Spike/Matrix Spike Duplicate Recoveries							Sample ID: 24-005230-0001			
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes
Abamectin	0.000	0.886	0.949	1.000	6.9%	< 30	88.6%	94.9%	50 - 150	
Acephate	0.000	0.524	0.478	0.800	9.2%	< 30	65.5%	59.8%	50 - 150	
Acequinocyl	0.000	4.326	4.086	4.000	5.7%	< 30	108.1%	102.2%	50 - 150	
Acetamiprid	0.000	0.386	0.372	0.400	3.7%	< 30	96.6%	93.1%	50 - 150	
Aldicarb	0.000	0.721	0.683	0.800	5.5%	< 30	90.2%	85.4%	50 - 150	
Azoxystrobin	0.028	0.405	0.381	0.400	6.7%	< 30	94.2%	88.1%	50 - 150	
Bifenazate	0.000	0.297	0.294	0.400	1.0%	< 30	74.3%	73.6%	50 - 150	
Bifenthrin	0.000	0.443	0.387	0.400	13.5%	< 30	110.8%	96.8%	50 - 150	
Boscalid	0.000	0.573	0.612	0.800	6.6%	< 30	71.6%	76.5%	50 - 150	
Carbaryl	0.000	0.359	0.363	0.400	0.9%	< 30	89.8%	90.7%	50 - 150	
Carbofuran	0.000	0.334	0.333	0.400	0.4%	< 30	83.6%	83.3%	50 - 150	
Chlorantraniliprole	0.000	0.338	0.342	0.400	1.0%	< 30	84.6%	85.4%	50 - 150	
Chlorfenapyr	0.000	0.750	2.161	2.000	96.9%	< 30	37.5%	108.0%	50 - 150	R, Q
Chlorpyrifos	0.000	0.408	0.381	0.400	6.6%	< 30	101.9%	95.4%	50 - 150	
Clofentezine	0.000	0.351	0.333	0.400	5.0%	< 30	87.6%	83.4%	50 - 150	
Cyfluthrin	0.000	1.882	1.909	2.000	1.4%	< 30	94.1%	95.5%	30 - 150	
Cypermethrin	0.018	2.168	2.201	2.000	1.5%	< 30	107.5%	109.1%	50 - 150	
Daminozide	0.000	0.693	0.633	2.000	9.0%	< 30	34.6%	31.7%	30 - 150	
Diazinon	0.000	0.354	0.347	0.400	2.0%	< 30	88.5%	86.7%	50 - 150	
Dichlorvos	0.000	2.000	2.001	2.000	0.0%	< 30	100.0%	100.0%	50 - 150	
Dimethoate	0.000	0.333	0.358	0.400	7.0%	< 30	83.4%	89.4%	50 - 150	
Ethoprophos	0.000	0.332	0.328	0.400	1.4%	< 30	83.1%	81.9%	50 - 150	
Etofenprox	0.000	0.728	0.732	0.800	0.7%	< 30	90.9%	91.5%	50 - 150	
Etoxazole	0.000	0.394	0.379	0.400	3.9%	< 30	98.5%	94.7%	50 - 150	
Fenoxycarb	0.000	0.365	0.352	0.400	3.6%	< 30	91.3%	88.1%	50 - 150	
Fenpyroximate	0.000	0.700	0.742	0.800	5.9%	< 30	87.5%	92.8%	50 - 150	
Fipronil	0.000	0.731	0.763	0.800	4.2%	< 30	91.4%	95.3%	50 - 150	
Fonicamid	0.000	0.835	0.818	1.000	2.0%	< 30	83.5%	81.8%	50 - 150	
Fludioxonil	0.028	0.833	0.828	0.800	0.5%	< 30	100.6%	100.0%	50 - 150	
Hexythiazox	0.000	1.278	1.267	1.000	0.8%	< 30	127.8%	126.7%	50 - 150	
Imazalil	0.000	0.382	0.347	0.400	9.8%	< 30	95.6%	86.6%	50 - 150	
Imidacloprid	0.000	0.659	0.624	0.800	5.3%	< 30	82.3%	78.0%	50 - 150	
Kresoxim-methyl	0.000	0.701	0.651	0.800	7.3%	< 30	87.6%	81.4%	50 - 150	
Malathion	0.000	0.314	0.335	0.400	6.4%	< 30	78.5%	83.7%	50 - 150	
Metalaxyl	0.000	0.354	0.362	0.400	2.1%	< 30	88.5%	90.4%	50 - 150	
Methiocarb	0.000	0.354	0.345	0.400	2.5%	< 30	88.4%	86.2%	50 - 150	
Methomyl	0.000	0.660	0.637	0.800	3.5%	< 30	82.5%	79.6%	50 - 150	
MGK-264	0.000	0.335	0.351	0.400	4.7%	< 30	83.8%	87.8%	50 - 150	
Myclobutanil	0.000	0.323	0.301	0.400	6.9%	< 30	80.7%	75.3%	50 - 150	
Naled	0.000	0.841	0.885	1.000	5.0%	< 30	84.1%	88.5%	50 - 150	
Oxamyl	0.000	2.004	1.909	2.000	4.8%	< 30	100.2%	95.5%	50 - 150	
Paclobutrazole	0.000	0.556	0.544	0.800	2.2%	< 30	69.5%	68.0%	50 - 150	
Parathion-Methyl	0.000	0.438	0.331	0.400	28.0%	< 30	109.5%	82.6%	30 - 150	
Permethrin	0.000	0.372	0.434	0.400	15.4%	< 30	93.0%	108.5%	50 - 150	
Phosmet	0.000	0.325	0.336	0.400	3.5%	< 30	81.1%	84.0%	50 - 150	
Piperonyl butoxide	0.000	2.054	1.897	2.000	8.0%	< 30	102.7%	94.8%	50 - 150	
Prallethrin	0.000	0.364	0.366	0.400	0.3%	< 30	91.1%	91.4%	50 - 150	
Propiconazole	0.020	0.714	0.681	0.800	4.9%	< 30	86.8%	82.7%	50 - 150	
Propoxur	0.000	0.366	0.359	0.400	2.0%	< 30	91.4%	89.6%	50 - 150	
Pyrethrin (Summe)	0.003	0.434	0.440	0.488	1.3%	< 30	88.4%	89.5%	50 - 150	
Pyridaben	0.000	0.380	0.406	0.400	6.5%	< 30	95.1%	101.5%	50 - 150	
Spinosad	0.000	0.363	0.348	0.388	4.2%	< 30	93.5%	89.7%	50 - 150	
Spiromesifen	0.000	0.391	0.392	0.400	0.3%	< 30	97.7%	98.0%	50 - 150	
Spirotetramat	0.000	0.442	0.443	0.400	0.3%	< 30	110.4%	110.7%	50 - 150	
Spiroxamine	0.000	0.719	0.719	0.800	0.1%	< 30	89.9%	89.8%	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.