



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

Report Number: 23-003182/D009.R000
Report Date: 04/24/2023
ORELAP#: OR100028
Purchase Order:
Received: 03/31/23 13:40

Customer:
Product identity: Full Spectrum Compliant CBD Distillate V3 GVL-TST582
Client/Metric ID: .
Laboratory ID: 23-003182-0005

Summary

Potency:

Analyte	Result (%)		
CBD	89.1	<ul style="list-style-type: none"> ● CBD ● CBC ● CBT ● CBG ● CBDV ● CBE ● CBN ● 9-THC 	CBD-Total 89.1%
CBC	0.828		THC-Total 0.215%
CBT	0.409		(Reported in percent of total sample)
CBG	0.398		
CBDV	0.346		
CBE	0.312		
CBN	0.258		
Δ9-THC	0.215		

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:

Product identity: Full Spectrum Compliant CBD Distillate V3 GVL-TST582 .
Client/Metric ID:
Sample Date:
Laboratory ID: 23-003182-0005
Evidence of Cooling: No
Temp: 18.9
Relinquished by: courier

Sample Results

Potency	Method: J AOAC 2015 V98-6 (mod) ^p			Units %	Batch: 2305629	Analyze: 4/4/23 7:01:00 PM
Analyte	As Received	Dry weight	LOQ	Notes		
CBC	0.828		0.0691			
CBC-A	< LOQ		0.0691			
CBC-Total	0.828		0.130			
CBD	89.1		0.691			
CBD-A	< LOQ		0.0691			
CBD-Total	89.1		0.752			
CBDV	0.346		0.0691			
CBDV-A	< LOQ		0.0691			
CBDV-Total	0.346		0.129			
CBE	0.312		0.0691			
CBG	0.398		0.0691			
CBG-A	< LOQ		0.0691			
CBG-Total	0.398		0.129			
CBL	< LOQ		0.0691			
CBL-A	< LOQ		0.0691			
CBL-Total	< LOQ		0.130			
CBN	0.258		0.0691			
CBT	0.409		0.0691			
Δ10-THC-9R	< LOQ		0.0691			
Δ10-THC-9S	< LOQ		0.0691			
Δ10-THC-Total	< LOQ		0.138			
Δ8-THC	< LOQ		0.0691			
Δ8-THCV	< LOQ		0.0691			
Δ9-THC	0.215		0.0691			
exo-THC	< LOQ		0.0691			
THC-A	< LOQ		0.0691			
THC-Total	0.215		0.130			
THCV	< LOQ		0.0691			
THCV-A	< LOQ		0.0691			
THCV-Total	< LOQ		0.129			
Total Cannabinoids	91.9					

*
 COA

- CBD
- CBC
- CBT
- CBG
- CBDV
- CBE
- CBN
- 9-THC



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Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2306447	04/20/23 AOAC 991.14 (Petrifilm) ^P		
Total Coliforms	< LOQ		cfu/g	10	2306447	04/20/23 AOAC 991.14 (Petrifilm) ^P		
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2306448	04/21/23 AOAC 2014.05 (RAPID) ^P		
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2306448	04/21/23 AOAC 2014.05 (RAPID) ^P		

Solvents Method: Residual Solvents by GC/MS^P Units µg/g Batch 2306600 Analyze 04/21/23 01:40 PM

Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
1,4-Dioxane	< LOQ	380	100	pass		2-Butanol	< LOQ	5000	200	pass	
2-Ethoxyethanol	< LOQ	160	30.0	pass		2-Methylbutane (Isopentane)	< LOQ		200		
2-Methylpentane	< LOQ		30.0			2-Propanol (IPA)	< LOQ	5000	200	pass	
2,2-Dimethylbutane	< LOQ		30.0			2,2-Dimethylpropane (neo-pentane)	< LOQ		200		
2,3-Dimethylbutane	< LOQ		30.0			3-Methylpentane	< LOQ		30.0		
Acetone	< LOQ	5000	200	pass		Acetonitrile	< LOQ	410	100	pass	
Benzene	< LOQ	2.00	1.00	pass		Butanes (sum)	< LOQ	5000	400	pass	
Cyclohexane	< LOQ	3880	200	pass		Ethyl acetate	< LOQ	5000	200	pass	
Ethyl benzene	< LOQ		200			Ethyl ether	< LOQ	5000	200	pass	
Ethylene glycol	< LOQ	620	200	pass		Ethylene oxide	< LOQ	50.0	20.0	pass	
Hexanes (sum)	< LOQ	290	150	pass		Isopropyl acetate	< LOQ	5000	200	pass	
Isopropylbenzene (Cumene)	< LOQ	70.0	30.0	pass		m,p-Xylene	< LOQ		200		
Methanol	< LOQ	3000	200	pass		Methylene chloride	< LOQ	600	60.0	pass	
Methylpropane (Isobutane)	< LOQ		200			n-Butane	< LOQ		200		
n-Heptane	< LOQ	5000	200	pass		n-Hexane	< LOQ		30.0		
n-Pentane	< LOQ		200			o-Xylene	< LOQ		200		
Pentanes (sum)	< LOQ	5000	600	pass		Propane	< LOQ	5000	200	pass	
Tetrahydrofuran	< LOQ	720	100	pass		Toluene	< LOQ	890	100	pass	
Total Xylenes	< LOQ		400			Total Xylenes and Ethyl benzene	< LOQ	2170	600	pass	

*SAMPLE



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Pesticides											
Method: AOAC 2007.01 & EN 15662 (mod) ^b											
Units mg/kg Batch 2306597 Analyze 04/21/23 12:33 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		Acetamidiprid [‡]	< LOQ	0.20	0.100	pass	
Aldicarb [‡]	< LOQ	0.40	0.200	pass		Azoxystrobin [‡]	< LOQ	0.20	0.100	pass	
Bifentazate [‡]	< LOQ	0.20	0.100	pass		Bifenthrin [‡]	< LOQ	0.20	0.100	pass	
Boscalid [‡]	< LOQ	0.40	0.200	pass		Carbaryl [‡]	< LOQ	0.20	0.100	pass	
Carbofuran [‡]	< LOQ	0.20	0.100	pass		Chlorantraniliprole [‡]	< LOQ	0.20	0.100	pass	
Chlorfenapyr [‡]	< LOQ	1.0	0.500	pass		Chlorpyrifos [‡]	< LOQ	0.20	0.100	pass	
Clofentezine [‡]	< LOQ	0.20	0.100	pass		Cyfluthrin [‡]	< LOQ	1.0	0.500	pass	
Cypermethrin [‡]	< LOQ	1.0	0.500	pass		Daminozide [‡]	< LOQ	1.0	0.500	pass	
Diazinon [‡]	< LOQ	0.20	0.100	pass		Dichlorvos [‡]	< LOQ	1.0	0.500	pass	
Dimethoate [‡]	< LOQ	0.20	0.100	pass		Ethoprophos [‡]	< LOQ	0.20	0.100	pass	
Etofenprox [‡]	< LOQ	0.40	0.200	pass		Etoxazole [‡]	< LOQ	0.20	0.100	pass	
Fenoxycarb [‡]	< LOQ	0.20	0.100	pass		Fenpyroximate [‡]	< LOQ	0.40	0.200	pass	
Fipronil [‡]	< LOQ	0.40	0.200	pass		Flonicamid [‡]	< LOQ	1.0	0.400	pass	
Fludioxonil [‡]	< LOQ	0.40	0.200	pass		Hexythiazox [‡]	< LOQ	1.0	0.400	pass	
Imazalil [‡]	< LOQ	0.20	0.100	pass		Imidacloprid [‡]	< LOQ	0.40	0.200	pass	
Kresoxim-methyl [‡]	< LOQ	0.40	0.200	pass		Malathion [‡]	< LOQ	0.20	0.100	pass	
Metalaxyl [‡]	< LOQ	0.20	0.100	pass		Methiocarb [‡]	< LOQ	0.20	0.100	pass	
Methomyl [‡]	< LOQ	0.40	0.200	pass		MGK-264 [‡]	< LOQ	0.20	0.100	pass	
Myclobutanil [‡]	< LOQ	0.20	0.100	pass		Naled [‡]	< LOQ	0.50	0.250	pass	
Oxamyl [‡]	< LOQ	1.0	0.500	pass		Pacllobutrazole [‡]	< LOQ	0.40	0.200	pass	
Parathion-Methyl [‡]	< LOQ	0.20	0.100	pass		Permethrin [‡]	< LOQ	0.20	0.100	pass	
Phosmet [‡]	< LOQ	0.20	0.100	pass		Piperonyl butoxide [‡]	< LOQ	2.0	1.00	pass	
Prallethrin [‡]	< LOQ	0.20	0.100	pass		Propiconazole [‡]	< LOQ	0.40	0.200	pass	
Propoxur [‡]	< LOQ	0.20	0.100	pass		Pyrethrin I (total) [‡]	< LOQ	1.0	0.500	pass	
Pyridaben [‡]	< LOQ	0.20	0.100	pass		Spinosad [‡]	< LOQ	0.20	0.100	pass	
Spiromesifen [‡]	< LOQ	0.20	0.100	pass		Spirotetramat [‡]	< LOQ	0.20	0.100	pass	
Spiroxamine [‡]	< LOQ	0.40	0.200	pass		Tebuconazole [‡]	< LOQ	0.40	0.200	pass	
Thiacloprid [‡]	< LOQ	0.20	0.100	pass		Thiamethoxam [‡]	< LOQ	0.20	0.100	pass	
Trifloxystrobin [‡]	< LOQ	0.20	0.100	pass							

Metals										
Analyte	Result	Limits	Units	LOQ	Batch	Analyzed	Method	Status	Notes	
Arsenic [‡]	< LOQ	0.200	mg/kg	0.0874	2306543	04/19/23	AOAC 2013.06 (mod.) ^b	pass		
Cadmium [‡]	< LOQ	0.200	mg/kg	0.0874	2306543	04/19/23	AOAC 2013.06 (mod.) ^b	pass		
Lead [‡]	< LOQ	0.500	mg/kg	0.0874	2306543	04/19/23	AOAC 2013.06 (mod.) ^b	pass		
Mercury [‡]	< LOQ	0.100	mg/kg	0.0437	2306543	04/19/23	AOAC 2013.06 (mod.) ^b	pass		



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Abbreviations

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

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

Ⓐ = ISO/IEC 17025:2017 accredited method.

Ⓜ = TNI accredited analyte.

Units of Measure

cfu/g = Colony forming units per gram

µg/g = Microgram per gram

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

% = Percentage of sample

% wt = µg/g divided by 10,000

Approved Signatory

Derrick Tanner
General Manager



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

Laboratory Quality Control Results
 Batch ID: 2305629

Laboratory Control Sample									
Analyte	LCS	Result	Spike	Units	% Rec	Limits		Evaluation	Notes
CBDVA	2	0.102	0.100	%	102	80.0	- 120	Acceptable	
CBDV	2	0.0998	0.100	%	99.8	80.0	- 120	Acceptable	
CBE	2	0.101	0.100	%	101	80.0	- 120	Acceptable	
CBDA	1	0.0931	0.093	%	100	90.0	- 110	Acceptable	
CBGA	1	0.0791	0.077	%	102	80.0	- 120	Acceptable	
CBG	1	0.0950	0.093	%	102	80.0	- 120	Acceptable	
CBD	1	0.0831	0.082	%	101	90.0	- 110	Acceptable	
THCV	2	0.0989	0.100	%	98.9	80.0	- 120	Acceptable	
d8THCV	2	0.0996	0.100	%	99.6	80.0	- 120	Acceptable	
THCVA	2	0.102	0.100	%	102	80.0	- 120	Acceptable	
CBN	1	0.0814	0.081	%	101	80.0	- 120	Acceptable	
exo-THC	2	0.0935	0.100	%	93.5	80.0	- 120	Acceptable	
d9THC	1	0.0902	0.093	%	96.5	90.0	- 110	Acceptable	
d8THC	1	0.0935	0.094	%	99.9	90.0	- 110	Acceptable	
9S-d10THC	1	0.0914	0.094	%	96.9	80.0	- 120	Acceptable	
CBL	2	0.0952	0.100	%	95.2	80.0	- 120	Acceptable	
9R-d10THC	1	0.0892	0.096	%	93.0	80.0	- 120	Acceptable	
CBC	2	0.0991	0.100	%	99.1	80.0	- 120	Acceptable	
THCA	1	0.103	0.108	%	95.2	90.0	- 110	Acceptable	
CBCA	2	0.105	0.100	%	105	80.0	- 120	Acceptable	
CBLA	2	0.105	0.100	%	105	80.0	- 120	Acceptable	
CBT	2	0.0863	0.100	%	86.3	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	
9S-d10THC	<LOQ	0.077	%	< 0.077	Acceptable	
CBL	<LOQ	0.077	%	< 0.077	Acceptable	
9R-d10THC	<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



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

Laboratory Quality Control Results

J AOAC 2015 V98-6		Batch ID: 2305629						
Sample Duplicate		Sample ID: 23-003141-0001						
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	
CBDA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBG	2.75	2.79	0.077	%	1.23	< 20	Acceptable	
CBD	0.117	0.132	0.077	%	12.0	< 20	Acceptable	
THCV	0.354	0.362	0.077	%	2.08	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBN	0.205	0.207	0.077	%	1.15	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
d9THC	70.1	70.8	0.077	%	1.05	< 20	Acceptable	
d8THC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
9S-d10THC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
9R-d10THC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBC	1.55	1.57	0.077	%	1.28	< 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	<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:

% - Percent



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

Laboratory Pesticide Quality Control Results

AOAC2007.1 & EN 15662		Units: mg/Kg			Batch ID 2306597			
Method Blank	Blank Result	Blank Limits	Notes	LCS Result	LCS Spk	LCS % Re	Limits	Notes
Abamectin	0.00	< 0.250		1.094	1.00	109.4	50.0	150
Acephate	0.00	< 0.200		0.800	0.80	100.0	60.0	120
Acetamiprid	0.00	< 1.000		3.545	4.00	88.6	40.0	160
Acetamiprid	0.00	< 0.100		0.433	0.40	108.3	60.0	120
Aldicarb	0.00	< 0.200		0.844	0.80	105.5	60.0	120
Azoxystrobin	0.00	< 0.100		0.429	0.40	107.3	60.0	120
Bifenazate	0.00	< 0.100		0.490	0.40	122.5	60.0	120
Bifenthrin	0.00	< 0.100		0.396	0.40	99.1	50.0	150
Boscalid	0.00	< 0.200		0.920	0.80	115.0	60.0	120
Carbaryl	0.00	< 0.100		0.420	0.40	104.9	60.0	120
Carbendazim	0.00	< 0.100		0.431	0.40	107.9	60.0	120
Chlorantraniliprole	0.00	< 0.100		0.419	0.40	104.7	60.0	120
Chlorfenapyr	0.00	< 0.500		2.346	2.00	117.3	60.0	120
Chlorpyrifos	0.00	< 0.100		0.408	0.40	100.7	60.0	120
Clofentezine	0.00	< 0.100		0.422	0.40	105.4	60.0	120
Cyfluthrin	0.00	< 0.500		2.172	2.00	108.6	50.0	150
Cypermethrin	0.00	< 0.500		2.093	2.00	104.7	50.0	150
Daminozide	0.215	< 0.500		2.409	2.00	120.5	60.0	120
Diazinon	0.00	< 0.100		0.467	0.40	116.8	60.0	120
Dichlorvos	0.00	< 0.500		2.060	2.00	103.0	60.0	120
Dimethoate	0.00	< 0.100		0.414	0.40	103.6	60.0	120
Ethiofencarb	0.00	< 0.100		0.427	0.40	106.7	60.0	120
Ethofenprox	0.00	< 0.200		0.822	0.80	102.8	50.0	150
Etoxazole	0.00	< 0.100		0.426	0.40	106.4	60.0	120
Fenoxycarb	0.00	< 0.100		0.421	0.40	105.3	60.0	120
Fenpyroximate	0.00	< 0.200		0.904	0.80	113.1	60.0	120
Fipronil	0.00	< 0.200		0.822	0.80	102.7	60.0	120
Fonicamid	0.00	< 0.250		1.019	1.00	101.9	60.0	120
Fludioxonil	0.00	< 0.200		0.906	0.80	113.2	50.0	150
Hexythiazox	0.00	< 0.250		1.037	1.00	103.7	60.0	120
Imazalil	0.00	< 0.100		0.426	0.40	106.5	60.0	120
Imidacloprid	0.00	< 0.200		0.839	0.80	104.9	60.0	120
Kiesoxim-methyl	0.00	< 0.200		0.857	0.80	107.1	60.0	120
Malathion	0.00	< 0.100		0.430	0.40	107.4	60.0	120
Metaxyl	0.00	< 0.100		0.442	0.40	110.5	60.0	120
Methiocarb	0.00	< 0.100		0.421	0.40	105.3	60.0	120
Methomyl	0.00	< 0.200		0.849	0.80	106.1	60.0	120
MCK-264	0.00	< 0.100		0.433	0.40	108.3	50.0	150
Myclobutanil	0.00	< 0.100		0.443	0.40	110.8	60.0	120
Naled	0.00	< 0.250		1.021	1.00	102.1	50.0	150
Oxaryl	0.00	< 0.500		2.166	2.00	108.3	60.0	120
Padobutrazole	0.00	< 0.200		0.898	0.80	112.2	60.0	120
Parathion-Methyl	0.00	< 0.100		0.508	0.40	126.9	50.0	150
Permethrin	0.002	< 0.100		0.424	0.40	106.0	50.0	150
Phosmet	0.00	< 0.100		0.416	0.40	104.1	50.0	150
Piperonyl butoxide	0.00	< 0.500		2.250	2.00	112.5	60.0	120
Prallethrin	0.00	< 0.100		0.418	0.40	104.4	60.0	120
Propiconazole	0.00	< 0.200		0.863	0.80	107.9	60.0	120
Propoxur	0.00	< 0.100		0.426	0.40	106.9	60.0	120
Pyrethrin (Summe)	0.00	< 0.100		0.522	0.48	108.9	60.0	120
Pyridaben	0.00	< 0.100		0.419	0.40	104.8	50.0	150
Spinosad	0.00	< 0.100		0.413	0.38	106.6	50.0	150
Spiromesfen	0.00	< 0.100		0.446	0.40	111.1	60.0	120
Spirotetramat	0.00	< 0.100		0.435	0.40	108.7	60.0	120
Spiroxamine	0.00	< 0.200		0.864	0.80	107.9	60.0	120
Tebuconazole	0.00	< 0.200		0.900	0.80	112.5	60.0	120
Thiadoprid	0.00	< 0.100		0.425	0.40	106.2	60.0	120
Thiamethoxam	0.00	< 0.100		0.437	0.40	109.2	60.0	120
Trifloxystrobin	0.00	< 0.100		0.427	0.40	106.9	60.0	120

Test results relate only to the parameters tested and to the samples as received by the laboratory. Test results meet all requirements of NELAP and the Columbia Laboratories quality assurance plan unless otherwise noted. This report shall not be reproduced, except in full, without the written consent of this laboratory. Samples will be retained for a maximum of 30 days from the receipt date unless prior arrangements have been made.

Testing in accordance with: OAR 333-007-0390 OAR 333-007-0400 OAR 333-007-0410 OAR 333-007-0430



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

Report Number: 23-003182/D009.R000
 Report Date: 04/24/2023
 ORELAP#: OR100028
 Purchase Order:
 Received: 03/31/23 13:40

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

Laboratory Pesticide Quality Control Results

AOAC2007.1 & EN 15662		Units: mg/Kg				Batch ID 2306597				
Matrix Spke/Matrix Spke Duplicate Recoveries	Result	MS Res	MSD Res	Spike	RFD%	Limit	MS% Re	MSD % Re	Limits	Notes
Abamectin	0.00	0.968	1.049	1.00	8.0%	< 30	96.8%	104.9%	50 - 150	
Acephate	0.00	0.730	0.740	0.800	1.3%	< 30	91.3%	92.5%	50 - 150	
Acetamiprid	0.00	0.379	0.398	0.400	4.8%	< 30	94.7%	99.4%	50 - 150	
Aldicarb	0.00	0.759	0.864	0.800	12.9%	< 30	94.9%	108.0%	50 - 150	
Azoxystrobin	0.00	0.338	0.367	0.400	8.4%	< 30	84.0%	91.4%	50 - 150	
Bifenazate	0.106	0.491	0.560	0.400	16.9%	< 30	96.3%	113.6%	50 - 150	
Bifenthrin	0.00	0.214	0.251	0.400	16.0%	< 30	53.8%	62.8%	50 - 150	
Boscalid	0.745	1.234	1.400	0.800	29.0%	< 30	61.1%	81.9%	50 - 150	
Carbaryl	0.00	0.313	0.331	0.400	5.7%	< 30	78.2%	82.8%	50 - 150	
Carbofuran	0.00	0.350	0.380	0.400	8.0%	< 30	87.8%	94.9%	50 - 150	
Chlorantraniliprole	0.00	0.374	0.401	0.400	7.0%	< 30	93.8%	100.4%	50 - 150	
Chlorfenapyr	0.00	1.788	1.671	2.000	6.8%	< 30	89.4%	83.5%	50 - 150	
Chlorpyrifos	0.062	0.453	0.409	0.400	12.1%	< 30	97.9%	86.7%	50 - 150	
Clofentezine	0.011	0.225	0.265	0.400	17.0%	< 30	53.8%	63.8%	50 - 150	
Cyfluthrin	0.00	1.153	1.215	2.000	5.2%	< 30	57.8%	60.7%	30 - 150	
Cypermethrin	0.00	0.992	1.010	2.000	1.8%	< 30	49.6%	50.5%	50 - 150	Q
Daminozide	0.316	2.415	2.869	2.000	19.6%	< 30	104.9%	127.7%	30 - 150	
Diazinon	0.00	0.162	0.193	0.400	17.0%	< 30	40.4%	48.3%	50 - 150	Q
Dichlorvos	0.00	1.691	1.700	2.000	0.5%	< 30	84.8%	85.0%	50 - 150	
Dimethoate	0.00	0.355	0.366	0.400	3.0%	< 30	88.8%	91.9%	50 - 150	
Ethionphos	0.003	0.343	0.364	0.400	6.0%	< 30	85.1%	90.3%	50 - 150	
Etofenprox	0.036	0.500	0.560	0.800	12.2%	< 30	58.0%	65.5%	50 - 150	
Etoxazole	0.00	0.330	0.363	0.400	9.5%	< 30	82.4%	90.7%	50 - 150	
Fenoxycarb	0.00	0.305	0.377	0.400	21.1%	< 30	76.2%	94.1%	50 - 150	
Fenpyroximate	0.00	0.209	0.259	0.800	21.3%	< 30	26.1%	32.4%	50 - 150	Q
Fipronil	0.00	0.448	0.494	0.800	9.8%	< 30	56.0%	61.7%	50 - 150	
Fonicamid	0.00	0.923	0.955	1.000	3.4%	< 30	92.3%	95.5%	50 - 150	
Fludioxonil	0.00	1.308	1.184	0.800	10.0%	< 30	163.5%	148.0%	50 - 150	Q
Hexythiazox	0.00	0.814	0.904	1.000	10.8%	< 30	81.4%	90.4%	50 - 150	
Imazalil	0.00	0.356	0.399	0.400	11.4%	< 30	89.1%	99.9%	50 - 150	
Imidacloprid	0.00	0.718	0.744	0.800	3.6%	< 30	89.7%	93.0%	50 - 150	
Kiesoxim-methyl	0.00	0.561	0.663	0.800	16.7%	< 30	70.1%	82.9%	50 - 150	
Malathion	0.00	0.309	0.367	0.400	16.9%	< 30	77.4%	91.7%	50 - 150	
Metolaxyl	0.00	0.351	0.436	0.400	21.8%	< 30	87.9%	109.1%	50 - 150	
Methiocarb	0.00	0.322	0.387	0.400	18.1%	< 30	80.8%	96.7%	50 - 150	
Methomyl	0.00	0.771	0.855	0.800	10.3%	< 30	96.3%	106.8%	50 - 150	
MCK-264	0.00	0.196	0.242	0.400	21.2%	< 30	48.9%	60.5%	50 - 150	Q
Mydobutani	0.046	0.315	0.343	0.400	9.7%	< 30	67.4%	74.2%	50 - 150	
Naled	0.00	0.764	0.836	1.000	9.1%	< 30	76.4%	83.6%	50 - 150	
Oxamyl	0.00	1.904	1.986	2.000	4.2%	< 30	95.2%	99.3%	50 - 150	
Padbutrazole	0.003	0.659	0.751	0.800	13.1%	< 30	82.0%	93.8%	50 - 150	
Parathion-Methyl	0.00	0.257	0.353	0.400	32.8%	< 30	64.3%	89.8%	30 - 150	R
Permethrin	0.011	0.281	0.309	0.400	9.8%	< 30	67.7%	74.7%	50 - 150	
Phosmet	0.002	0.327	0.406	0.400	21.8%	< 30	81.4%	101.0%	50 - 150	
Piperonyl butoxide	0.00	1.911	2.179	2.000	13.1%	< 30	95.8%	109.0%	50 - 150	
Prallethrin	0.00	0.222	0.276	0.400	21.5%	< 30	55.8%	69.0%	50 - 150	
Propiconazole	0.059	0.655	0.783	0.800	21.1%	< 30	69.7%	86.1%	50 - 150	
Propoxur	0.007	0.353	0.373	0.400	5.8%	< 30	86.3%	91.5%	50 - 150	
Pyrethrin (Summe)	0.106	0.442	0.501	0.488	16.2%	< 30	68.8%	80.9%	50 - 150	
Pyridaben	0.012	0.266	0.287	0.400	7.6%	< 30	63.7%	68.7%	50 - 150	
Spirosad	0.00	0.303	0.336	0.388	10.3%	< 30	78.1%	86.6%	50 - 150	
Spiromesfen	0.00	0.347	0.371	0.400	6.6%	< 30	86.7%	92.7%	50 - 150	
Spirotetramat	0.00	0.462	0.504	0.400	8.6%	< 30	115.5%	125.9%	50 - 150	
Spiroxamine	0.00	0.731	0.901	0.800	20.8%	< 30	91.4%	112.7%	50 - 150	
Tebuconazole	0.00	0.629	0.760	0.800	18.9%	< 30	78.8%	95.1%	50 - 150	
Thiadoprid	0.00	0.355	0.404	0.400	12.8%	< 30	89.1%	101.1%	50 - 150	
Thiamethoxam	0.00	0.384	0.408	0.400	6.1%	< 30	95.9%	102.0%	50 - 150	
Trifloxystrobin	0.005	0.299	0.336	0.400	12.0%	< 30	73.9%	82.8%	50 - 150	

COA*

*SAMPLE



12423 NE Whitaker Way
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 503-254-1794

Report Number: 23-003182/D009.R000
 Report Date: 04/24/2023
 ORELAP#: OR100028
 Purchase Order:
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Revision: 2 Document ID: 7087
 Legacy ID: CFL-E33Effective:

Laboratory Quality Control Results

Residual Solvents				Batch ID: 2306600						
Method Blank				Laboratory Control Sample						
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes	
Propane	ND	< 200		444	584	µg/g	76.0	60 - 120		
Isobutane	ND	< 200		553	767	µg/g	72.1	60 - 120		
Butane	ND	< 200		586	782	µg/g	74.9	60 - 120		
2,2-Dimethylpropane	ND	< 200		764	939	µg/g	81.4	60 - 120		
Methanol	ND	< 200		1620	1610	µg/g	100.6	60 - 120		
Ethylene Oxide	ND	< 30		46.4	57.1	µg/g	81.3	60 - 120		
2-Methylbutane	ND	< 200		1570	1600	µg/g	98.1	60 - 120		
Pentane	ND	< 200		1620	1610	µg/g	100.6	60 - 120		
Ethanol	ND	< 200		1600	1600	µg/g	100.0	70 - 130		
Ethyl Ether	ND	< 200		1570	1610	µg/g	97.5	60 - 120		
2,2-Dimethylbutane	ND	< 30		173	173	µg/g	100.0	60 - 120		
Acetone	ND	< 200		1600	1620	µg/g	98.8	60 - 120		
2-Propanol	ND	< 200		1590	1600	µg/g	99.4	60 - 120		
Ethyl Formate	ND	< 500		1690	1610	µg/g	105.0	70 - 130		
Acetonitrile	ND	< 100		470	488	µg/g	96.3	60 - 120		
Methyl Acetate	ND	< 500		1450	1610	µg/g	90.1	70 - 130		
2,3-Dimethylbutane	ND	< 30		167	165	µg/g	101.2	60 - 120		
Dichloromethane	ND	< 60		482	487	µg/g	99.0	60 - 120		
2-Methylpentane	ND	< 30		156	160	µg/g	97.5	60 - 120		
MTBE	ND	< 500		1500	1600	µg/g	93.8	70 - 130		
3-Methylpentane	ND	< 30		153	161	µg/g	95.0	60 - 120		
Hexane	ND	< 30		157	162	µg/g	96.9	60 - 120		
1-Propanol	ND	< 500		1500	1620	µg/g	92.6	70 - 130		
Methylethylketone	ND	< 500		1450	1610	µg/g	90.1	70 - 130		
Ethyl acetate	ND	< 200		1550	1600	µg/g	96.9	60 - 120		
2-Butanol	ND	< 200		1580	1610	µg/g	98.1	60 - 120		
Tetrahydrofuran	ND	< 100		464	483	µg/g	96.1	60 - 120		
Cyclohexane	ND	< 200		1570	1610	µg/g	97.5	60 - 120		
2-methyl-1-propanol	ND	< 500		1470	1630	µg/g	90.2	70 - 130		
Benzene	ND	< 1		4.84	4.98	µg/g	97.2	60 - 120		
Isopropyl Acetate	ND	< 200		1580	1610	µg/g	98.1	60 - 120		
Heptane	ND	< 200		1550	1620	µg/g	95.7	60 - 120		
1-Butanol	ND	< 500		1410	1600	µg/g	88.1	70 - 130		
Propyl Acetate	ND	< 500		1350	1620	µg/g	83.3	70 - 130		
1,4-Dioxane	ND	< 100		439	494	µg/g	88.9	60 - 120		
2-Ethoxyethanol	ND	< 30		166	165	µg/g	100.6	60 - 120		
Methylisobutylketone	ND	< 500		1430	1610	µg/g	88.8	70 - 130		
3-Methyl-1-butanol	ND	< 500		1350	1610	µg/g	83.9	70 - 130		
Ethylene Glycol	ND	< 200		411	486	µg/g	84.6	60 - 120		
Toluene	ND	< 100		447	513	µg/g	87.1	60 - 120		
Isobutyl Acetate	ND	< 500		1390	1600	µg/g	86.9	70 - 130		
1-Pentanol	ND	< 500		1340	1610	µg/g	83.2	70 - 130		
Butyl Acetate	ND	< 500		1370	1610	µg/g	85.1	70 - 130		
Ethylbenzene	ND	< 200		874	967	µg/g	90.4	60 - 120		
m,p-Xylene	ND	< 200		886	994	µg/g	89.1	60 - 120		
o-Xylene	ND	< 200		877	992	µg/g	88.4	60 - 120		
Cumene	ND	< 30		144	171	µg/g	84.2	60 - 120		
Anisole	ND	< 500		1200	1610	µg/g	74.5	70 - 130		
DMSO	ND	< 500		1230	1610	µg/g	76.4	70 - 130		
1,2-dimethoxyethane	ND	< 50		157	172	µg/g	91.3	70 - 130		
Triethylamine	ND	< 500		1380	1620	µg/g	85.2	70 - 130		
N,N-dimethylformamide	ND	< 150		416	499	µg/g	83.4	70 - 130		
N,N-dimethylacetamide	ND	< 150		401	491	µg/g	81.7	70 - 130		
Pyridine	ND	< 50		139	171	µg/g	81.3	70 - 130		
Sulfone	ND	< 50		121	160	µg/g	75.6	70 - 130		
1,2-Dichloroethane	ND	< 1		0.844	1	µg/g	84.4	70 - 130		
Chloroform	ND	< 1		0.904	1	µg/g	90.4	70 - 130		
Trichloroethylene	ND	< 1		0.927	1	µg/g	92.7	70 - 130		
1,1-Dichloroethane	ND	< 1		0.87	1	µg/g	87.0	70 - 130		

SAMPLE ONLY



12423 NE Whitaker Way
 Portland, OR 97230
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Revision: 2 Document ID: 7087
 Legacy ID: CFL-E33Effective:

QC - Sample Duplicate Sample ID: 23-004569-0001

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	
1,2-Dichloroethane	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Chloroform	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Trichloroethylene	ND	ND	1	µg/g	0.0	< 20	Acceptable	
1,1-Dichloroethane	ND	ND	1	µg/g	0.0	< 20	Acceptable	

Abbreviations

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

Units of Measure:

µg/g - Microgram per gram or ppm



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SAMPLE COA

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.