



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:** GVB Oregon  
**Product identity:** Comp BSD V3 GVL-TST582  
**Client/Metric ID:** .  
**Laboratory ID:** 23-003182-0005

### Summary

**Potency:**

Analyte	Result (%)		
CBD	89.1	<ul style="list-style-type: none"> <li><span style="color: red;">●</span> CBD</li> <li><span style="color: blue;">●</span> CBC</li> <li><span style="color: green;">●</span> CBT</li> <li><span style="color: yellow;">●</span> CBG</li> <li><span style="color: magenta;">●</span> CBDV</li> <li><span style="color: cyan;">●</span> CBE</li> <li><span style="color: pink;">●</span> CBN</li> <li><span style="color: red;">●</span> 9-THC</li> </ul>	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:** GVB Oregon  
United States of America (USA)  
**Product identity:** Comp BSD 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) <sup>P</sup>			Units %	Batch: 2305629	Analyze: 4/4/23 7:01:00 PM
Analyte	As Received	Dry weight	LOQ	Notes		
CBC	0.828		0.0691		<ul style="list-style-type: none"> <li><span style="color: red;">●</span> CBD</li> <li><span style="color: blue;">●</span> CBC</li> <li><span style="color: green;">●</span> CBT</li> <li><span style="color: yellow;">●</span> CBG</li> <li><span style="color: purple;">●</span> CBDV</li> <li><span style="color: cyan;">●</span> CBE</li> <li><span style="color: pink;">●</span> CBN</li> <li><span style="color: red;">●</span> 9-THC</li> </ul>	
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			
<b>Total Cannabinoids</b>	91.9					



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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) <sup>P</sup>		
Total Coliforms	< LOQ		cfu/g	10	2306447	04/20/23 AOAC 991.14 (Petrifilm) <sup>P</sup>		
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2306448	04/21/23 AOAC 2014.05 (RAPID) <sup>P</sup>		
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2306448	04/21/23 AOAC 2014.05 (RAPID) <sup>P</sup>		

**Solvents** Method: Residual Solvents by GC/MS<sup>B</sup> 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	



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

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

J AOAC 2015 V98-6 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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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	
CBDV	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBD	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	Laboratory Control Sample							
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spk	LCS % Re	Limits	Notes
Abamectin	0.00	< 0.250		1.094	1.000	109.4	50.0	150
Acephate	0.00	< 0.200		0.800	0.800	100.0	60.0	120
Acequinocyl	0.00	< 1.000		3.545	4.000	88.6	40.0	160
Acetamiprid	0.00	< 0.100		0.433	0.400	108.3	60.0	120
Aldicarb	0.00	< 0.200		0.844	0.800	105.5	60.0	120
Azoxystrobin	0.00	< 0.100		0.429	0.400	107.3	60.0	120
Bifenazate	0.00	< 0.100		0.490	0.400	122.5	60.0	120
Bifenthrin	0.00	< 0.100		0.396	0.400	99.1	50.0	150
Boscalid	0.00	< 0.200		0.920	0.800	115.0	60.0	120
Carbaryl	0.00	< 0.100		0.420	0.400	104.9	60.0	120
Carbendazim	0.00	< 0.100		0.431	0.400	107.9	60.0	120
Chlorantraniliprole	0.00	< 0.100		0.419	0.400	104.7	60.0	120
Chlorfenapyr	0.00	< 0.500		2.346	2.000	117.3	60.0	120
Chlorpyrifos	0.00	< 0.100		0.403	0.400	100.7	60.0	120
Clofentezine	0.00	< 0.100		0.422	0.400	105.4	60.0	120
Cyfluthrin	0.00	< 0.500		2.172	2.000	108.6	50.0	150
Cypermethrin	0.00	< 0.500		2.093	2.000	104.7	50.0	150
Daminozide	0.215	< 0.500		2.409	2.000	120.5	60.0	120
Diazinon	0.00	< 0.100		0.467	0.400	116.8	60.0	120
Dichlorvos	0.00	< 0.500		2.060	2.000	103.0	60.0	120
Dimethoate	0.00	< 0.100		0.414	0.400	103.6	60.0	120
Ethoprophos	0.00	< 0.100		0.427	0.400	106.7	60.0	120
Etofenprox	0.00	< 0.200		0.822	0.800	102.8	50.0	150
Etoxazole	0.00	< 0.100		0.425	0.400	106.4	60.0	120
Fenoxycarb	0.00	< 0.100		0.421	0.400	105.3	60.0	120
Fenpyroximate	0.00	< 0.200		0.904	0.800	113.1	60.0	120
Fipronil	0.00	< 0.200		0.822	0.800	102.7	60.0	120
Fonicamid	0.00	< 0.250		1.019	1.000	101.9	60.0	120
Fludioxonil	0.00	< 0.200		0.905	0.800	113.2	50.0	150
Hexythiazox	0.00	< 0.250		1.037	1.000	103.7	60.0	120
Imazalil	0.00	< 0.100		0.425	0.400	106.5	60.0	120
Imidacloprid	0.00	< 0.200		0.839	0.800	104.9	60.0	120
Kiesoxim-methyl	0.00	< 0.200		0.857	0.800	107.1	60.0	120
Malathion	0.00	< 0.100		0.430	0.400	107.4	60.0	120
Metaxyl	0.00	< 0.100		0.442	0.400	110.5	60.0	120
Methiocarb	0.00	< 0.100		0.421	0.400	105.3	60.0	120
Methomyl	0.00	< 0.200		0.849	0.800	106.1	60.0	120
MCK-264	0.00	< 0.100		0.433	0.400	108.3	50.0	150
Mydobutani	0.00	< 0.100		0.443	0.400	110.8	60.0	120
Naled	0.00	< 0.250		1.021	1.000	102.1	50.0	150
Oxamyl	0.00	< 0.500		2.166	2.000	108.3	60.0	120
Padobutrazole	0.00	< 0.200		0.898	0.800	112.2	60.0	120
Parathion-Methyl	0.00	< 0.100		0.508	0.400	126.9	50.0	150
Permethrin	0.002	< 0.100		0.424	0.400	106.0	50.0	150
Phosmet	0.00	< 0.100		0.416	0.400	104.1	50.0	150
Piperonyl butoxide	0.00	< 0.500		2.250	2.000	112.5	60.0	120
Prallethrin	0.00	< 0.100		0.418	0.400	104.4	60.0	120
Propiconazole	0.00	< 0.200		0.863	0.800	107.9	60.0	120
Propoxur	0.00	< 0.100		0.425	0.400	106.9	60.0	120
Pyrethrin (Summe)	0.00	< 0.100		0.522	0.488	106.9	60.0	120
Pyridaben	0.00	< 0.100		0.419	0.400	104.8	50.0	150
Spinosad	0.00	< 0.100		0.413	0.388	106.6	50.0	150
Spiromesfen	0.00	< 0.100		0.445	0.400	111.1	60.0	120
Spirotetramat	0.00	< 0.100		0.435	0.400	108.7	60.0	120
Spiroxamine	0.00	< 0.200		0.864	0.800	107.9	60.0	120
Tebuconazole	0.00	< 0.200		0.900	0.800	112.5	60.0	120
Thiadoprid	0.00	< 0.100		0.425	0.400	106.2	60.0	120
Thiamethoxam	0.00	< 0.100		0.437	0.400	109.2	60.0	120
Trifloxystrobin	0.00	< 0.100		0.427	0.400	106.9	60.0	120





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



Report Number: 23-003182/D009.R000  
 Report Date: 04/24/2023  
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 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	
Acequinocyl	0.00	2.682	3.218	4.000	18.1%	< 30	67.1%	80.4%	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	49.8%	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	0.856	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	
Padobutrazole	0.003	0.659	0.751	0.800	13.1%	< 30	82.0%	93.5%	50 - 150	
Parathion-Methyl	0.00	0.257	0.358	0.400	32.8%	< 30	64.3%	89.5%	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	0.099	0.655	0.788	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	
Spinosad	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	



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Report Number: 23-003182/D009.R000  
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 Received: 03/31/23 13:40



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	



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**Received:** 03/31/23 13:40

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



12423 NE Whitaker Way  
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**Report Number:** 23-003182/D009.R000  
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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.