



Job Number: 19-008626

Report Number: 19-008626-00

Report Date: 07/29/2019

ORELAP#: OR100028

Purchase Order:

Received: 07/22/19 14:57

This report cannot be used for ODA, OHA or OLCC compliance requirements.

Product identity: HDTO 1159 Sentia Lem Gin 1000mg Client/Metrc ID: Laboratory ID: 19-008626-0003 Sample Date:

Summary

Potency:

Analyte CBD	Result 3.42	Limits	Units %	LOQ 0.09	CBD-Total per 1g	34.2 mg/1g
Analyte per 1g	Result	Limits	Units	LOQ	THC-Total (%)	< 0.172 %
CBD per 1g	34.2		mg/1g	1.00		

Residual Solvents:

All analytes passing and less than LOQ.

Pesticides:

All analytes passing and less than LOQ.





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Customer: Sentia Wellness

3931 NE Columbia Blvd Portland Oregon 97211

United States

Product identity: HDTD 1159 Sentia Lem Gin 1000mg

Client/Metrc ID: .

Sample Date:

Laboratory ID: 19-008626-0003
Relinquished by: Received By Mail

Temp: 19.5 °C **Serving Size #1:** 1 g

Sample Results

Potency			Batch: 190	6616			
Analyte	Result	Limits	Units	LOQ	Analyze	Method	Notes
CBC [†]	< LOQ		%	0.0917	07/24/19	J AOAC 2015 V98-6	
CBC-A [†]	< LOQ		%	0.0917	07/24/19	J AOAC 2015 V98-6	
CBC-Total [†]	< LOQ		%	0.172	07/26/19	J AOAC 2015 V98-6	
CBD	3.42		%	0.0917	07/24/19	J AOAC 2015 V98-6	
CBD-A	< LOQ		%	0.0917	07/24/19	J AOAC 2015 V98-6	
CBD-Total	3.42		%	0.172	07/26/19	J AOAC 2015 V98-6	
CBDV†	< LOQ		%	0.0917	07/24/19	J AOAC 2015 V98-6	
CBDV-A [†]	< LOQ		%	0.0917	07/24/19	J AOAC 2015 V98-6	
CBDV-Total [†]	< LOQ		%	0.171	07/26/19	J AOAC 2015 V98-6	
CBG [†]	< LOQ		%	0.0917	07/24/19	J AOAC 2015 V98-6	
CBG-A [†]	< LOQ		%	0.0917	07/24/19	J AOAC 2015 V98-6	
CBG-Total [†]	< LOQ		%	0.171	07/26/19	J AOAC 2015 V98-6	
CBL†	< LOQ		%	0.0917	07/24/19	J AOAC 2015 V98-6	
CBN	< LOQ		%	0.0917	07/24/19	J AOAC 2015 V98-6	
∆8-THC [†]	< LOQ		%	0.0917	07/24/19	J AOAC 2015 V98-6	
19-THC	< LOQ		%	0.0917	07/24/19	J AOAC 2015 V98-6	
ГНС-А	< LOQ		%	0.0917	07/24/19	J AOAC 2015 V98-6	
ΓHC-Total	< LOQ		%	0.172	07/26/19	J AOAC 2015 V98-6	
ΓHCV [†]	< LOQ		%	0.0917	07/24/19	J AOAC 2015 V98-6	
ΓHCV-A [†]	< LOQ		%	0.0917	07/24/19	J AOAC 2015 V98-6	
ΓHCV-Total [†]	< LOQ		%	0.171	07/26/19	J AOAC 2015 V98-6	
Potency per 1g			Batch: 190	6616			
Analyte	Result	Limits	Units	LOQ	Analyze	Method	Notes
CBC per 1g [†]	< LOQ		mg/1g	1.00	07/26/19	J AOAC 2015 V98-6	
CBC-A per 1g [†]	< LOQ		mg/1g	1.00	07/26/19	J AOAC 2015 V98-6	
CBC-Total per 1g [†]	< LOQ		mg/1g	1.88	07/26/19	J AOAC 2015 V98-6	

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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 Pixis 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 kept a maximum of 15 days from the report date unless prior arrangements have been made.





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Potency per 1g			Batch: 19	06616				
Analyte	Result	Limits	Units	L	oq	Analyze	Method	Notes
CBD per 1g	34.2		mg/1g	1.	00	07/26/19	J AOAC 2015 V98-6	
CBD-A per 1g	< LOQ		mg/1g	1.	00	07/26/19	J AOAC 2015 V98-6	
CBD-Total per 1g	34.2		mg/1g	1.	88	07/26/19	J AOAC 2015 V98-6	
CBDV per 1g [†]	< LOQ		mg/1g	1.	00	07/26/19	J AOAC 2015 V98-6	
CBDV-A per 1g [†]	< LOQ		mg/1g	1.	00	07/26/19	J AOAC 2015 V98-6	
CBDV-Total per 1g [†]	< LOQ		mg/1g	1.	87	07/26/19	J AOAC 2015 V98-6	
CBG per 1g [†]	< LOQ		mg/1g	1.	00	07/26/19	J AOAC 2015 V98-6	
CBG-A per 1g [†]	< LOQ		mg/1g	1.	00	07/26/19	J AOAC 2015 V98-6	
CBG-Total per 1g [†]	< LOQ		mg/1g	1.	88	07/26/19	J AOAC 2015 V98-6	
CBL per 1g [†]	< LOQ		mg/1g	1.	00	07/26/19	J AOAC 2015 V98-6	
CBN per 1g	< LOQ		mg/1g	1.	00	07/26/19	J AOAC 2015 V98-6	
$\Delta 8$ -THC per 1g †	< LOQ		mg/1g	1.	00	07/26/19	J AOAC 2015 V98-6	
Δ9-THC per 1g	< LOQ		mg/1g	1.	00	07/26/19	J AOAC 2015 V98-6	
THC-A per 1g	< LOQ		mg/1g	1.	00	07/26/19	J AOAC 2015 V98-6	
THC-Total per 1g	< LOQ		mg/1g	1.	88	07/26/19	J AOAC 2015 V98-6	
THCV per 1g [†]	< LOQ		mg/1g	1.	00	07/26/19	J AOAC 2015 V98-6	
THCV-A per 1g [†]	< LOQ		mg/1g	1.	00	07/26/19	J AOAC 2015 V98-6	
THCV-Total per 1g [†]	< LOQ		mg/1g	1.	87	07/26/19	J AOAC 2015 V98-6	

Solvents	Method	EPA502	21A			Units µg/g Batch 19	906654	Analyz	ze 07/2	25/19 08:51 A	M
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	< LOQ		200		
2-Methylpentane	< LOQ		30.0			2-Propanol (IPA)	< LOQ	5000	200	pass	
2,2-Dimethylbutane	< LOQ		30.0			2,2-Dimethylpropane	< 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	30.0	pass	
Hexanes (sum)	< LOQ	290	150	pass		Isopropyl acetate	< LOQ	5000	200	pass	
Isopropylbenzene	< LOQ	70.0	30.0	pass		m,p-Xylene	< LOQ		200		
Methanol	< LOQ	3000	200	pass		Methylene chloride	< LOQ	600	200	pass	
Methylpropane	< 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	< LOQ	2170	600	pass	





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Pesticides	Method	AOAC	2007.01 & EN	15662 (mod)	Units mg/kg Ba	atch 1906643	Analy	ze 07/24/19 04:14 PM
Analyte	Result	Limits	LOQ Status	Notes	Analyte	Result	Limits	LOQ Status Notes
Abamectin	< LOQ	0.50	0.250 pass		Acephate	< LOQ	0.40	0.250 pass
Acequinocyl	< LOQ	2.0	1.00 pass		Acetamiprid	< LOQ	0.20	0.100 pass
Aldicarb	< LOQ	0.40	0.200 pass		Azoxystrobin	< LOQ	0.20	0.100 pass
Bifenazate	< LOQ	0.20	0.100 pass		Bifenthrin	< LOQ	0.20	0.100 pass
Boscalid	< LOQ	0.40	0.100 pass		Carbaryl	< LOQ	0.20	0.100 pass
Carbofuran	< LOQ	0.20	0.100 pass		Chlorantraniliprole	e < 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 (incl.	< LOQ	1.0	0.500 pass
Cypermethrin	< LOQ	1.0	0.500 pass		Daminozide	< LOQ	1.0	0.500 pass
Diazinon	< LOQ	0.20	0.100 pass		Dichlorvos	< LOQ	1.0	0.500 pass
Dimethoate	< LOQ	0.20	0.100 pass		Ethoprophos	< LOQ	0.20	0.100 pass
Etofenprox	< LOQ	0.40	0.200 pass		Etoxazole	< LOQ	0.20	0.100 pass
Fenoxycarb	< LOQ	0.20	0.100 pass		Fenpyroximate	< LOQ	0.40	0.200 pass
Fipronil	< LOQ	0.40	0.200 pass		Flonicamid	< LOQ	1.0	0.400 pass
Fludioxonil	< LOQ	0.40	0.200 pass		Hexythiazox	< LOQ	1.0	0.400 pass
Imazalil	< LOQ	0.20	0.100 pass		Imidacloprid	< LOQ	0.40	0.200 pass
Kresoxim-methyl	< LOQ	0.40	0.200 pass		Malathion	< LOQ	0.20	0.100 pass
Metalaxyl	< LOQ	0.20	0.100 pass		Methiocarb	< LOQ	0.20	0.100 pass
Methomyl	< LOQ	0.40	0.200 pass		MGK-264	< LOQ	0.20	0.100 pass
Myclobutanil	< LOQ	0.20	0.100 pass		Naled	< LOQ	0.50	0.250 pass
Oxamyl	< LOQ	1.0	0.500 pass		Paclobutrazole	< LOQ	0.40	0.200 pass
Parathion-Methyl	< LOQ	0.20	0.200 pass		Permethrin	< LOQ	0.20	0.100 pass
Phosmet	< LOQ	0.20	0.100 pass		Piperonyl butoxide	e < 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					





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Abbreviations

Limits: Action Levels per OAR-333-007-0400, OAR-333-007-0210, OAR-333-007-0220

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

† = Analyte not NELAP accredited.

Units of Measure

g = Gram

μg/g = Microgram per gram

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

mg/1g = Milligram per 1g

% = Percentage of sample

% wt = μ g/g divided by 10,000

Approved Signatory

Derrick Tanner General Manager





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Company: Scotia V	Vellness												19-0			PIXIS Labs
Contact: 611) Hurt	cet				T	1	_	Anar	ysis i	Reque	ested				OI.	RELAP ID: OR100028
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Laboratory Quality Control Results

J AOAC 2015	5 V98-6			Bat	ch ID: 1906616		
Laboratory C	Control Sample						
Analyte	Result	Spike	Units	% Rec	Limits	Evaluation	Notes
CBDV-A	0.194	0.2	%	97.0	85 - 115	Acceptable	
CBDV	0.202	0.2	%	101	85 - 115	Acceptable	
CBD-A	0.200	0.2	%	100	85 - 115	Acceptable	
CBG-A	0.193	0.2	%	96.5	85 - 115	Acceptable	
CBG	0.192	0.2	%	96.0	85 - 115	Acceptable	
CBD	0.200	0.2	%	100	85 - 115	Acceptable	
THCV	0.194	0.2	%	97.0	85 - 115	Acceptable	
THCVA	0.192	0.2	%	96.0	85 - 115	Acceptable	
CBN	0.201	0.2	%	101	85 - 115	Acceptable	
THC	0.192	0.2	%	96.0	85 - 115	Acceptable	
D8THC	0.181	0.2	%	90.5	85 - 115	Acceptable	
CBL	0.187	0.2	%	93.5	85 - 115	Acceptable	
CBC	0.191	0.2	%	95.5	85 - 115	Acceptable	
THCA	0.207	0.2	%	104	85 - 115	Acceptable	
CBCA	0.182	0.2	%	91.0	85 - 115	Acceptable	

Method Blank

Wicking Blank						
Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDV-A	ND	0.1	%	< 0.1	Acceptable	
CBDV	ND	0.1	%	< 0.1	Acceptable	
CBD-A	ND	0.1	%	< 0.1	Acceptable	
CBG-A	ND	0.1	%	< 0.1	Acceptable	
CBG	ND	0.1	%	< 0.1	Acceptable	
CBD	ND	0.1	%	< 0.1	Acceptable	
THCV	ND	0.1	%	< 0.1	Acceptable	
THCVA	ND	0.1	%	< 0.1	Acceptable	
CBN	ND	0.1	%	< 0.1	Acceptable	
THC	ND	0.1	%	< 0.1	Acceptable	
D8THC	ND	0.1	%	< 0.1	Acceptable	
CBL	ND	0.1	%	< 0.1	Acceptable	
CBC	ND	0.1	%	< 0.1	Acceptable	
THCA	ND	0.1	%	< 0.1	Acceptable	
CBCA	ND	0.1	%	< 0.1	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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J AOAC 2015	V98-6				Bat	ch ID: 190661	6	
Sample Dupli	icate				Sam	ple ID: 19-008 4	79-0001	
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDV-A	ND	ND	0.1	%	0	< 20	Acceptable	
CBDV	ND	ND	0.1	%	0	< 20	Acceptable	
CBD-A	0.194	0.195	0.1	%	0.514	< 20	Acceptable	
CBG-A	0.921	0.921	0.1	%	0	< 20	Acceptable	
CBG	0.324	0.325	0.1	%	0.308	< 20	Acceptable	
CBD	ND	ND	0.1	%	0	< 20	Acceptable	
THCV	ND	ND	0.1	%	0	< 20	Acceptable	
THCVA	0.409	0.404	0.1	%	1.23	< 20	Acceptable	
CBN	0.203	0.195	0.1	%	4.02	< 20	Acceptable	
THC	11.7	11.8	0.1	%	0.851	< 20	Acceptable	
D8THC	ND	ND	0.1	%	0	< 20	Acceptable	
CBL	ND	ND	0.1	%	0	< 20	Acceptable	
CBC	0.205	0.209	0.1	%	1.93	< 20	Acceptable	
THCA	67.9	69.7	0.1	%	2.62	< 20	Acceptable	
CBCA	0.951	0.963	0.1	%	1.25	< 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: 0.01 Control: CFL-C22 Revised: 12/4/2018 Effective: 12/4/2018

Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN	15662	Unit	s: mg/Kg			Batch ID: 1906643					
Method Blank				Laboratory Cor	ntrol Samp	ole					
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes			
Acephate	ND	< 0.200		0.892	1.000	89.2	70 - 130	П			
Acequinocyl	ND	< 1.000		3.960	4.000	99.0	70 - 130				
Acetamiprid	ND	< 0.100	1	0.428	0.400	107.0	70 - 130				
Aldicarb	ND	< 0.200	1	0.847	0.800	105.9	70 - 130	1			
Abamectin	ND	< 0.288	1	1.020	1.000	102.0	70 - 130				
Azoxystrobin	ND	< 0.100	1	0.401	0.400	100.3	70 - 130	1			
Bifenazate	ND	< 0.100	+	0.424	0.400	106.0	70 - 130				
Bifenthrin	ND	< 0.100	1	0.396	0.400	99.0	70 - 130				
Boscalid	ND	< 0.100	1	0.768	0.800	96.0	70 - 130	1			
Carbaryl	ND ND	< 0.100	-	0.423	0.400	105.8	70 - 130	1			
Carbofuran	ND ND	< 0.100	-	0.445	0.400	111.3	70 - 130	1			
Chlorantraniliprol	ND ND	< 0.100	+	0.414	0.400	103.5	70 - 130	 			
Chlorfenapyr	ND ND	< 1.000		1.660	2.000	83.0	70 - 130	-			
Chlorpyrifos	ND ND	< 0.100	-	0.403	0.400	100.8	70 - 130				
Clofentezine	ND ND	< 0.100		0.405	0.400	100.8	70 - 130				
Liorentezine Cvfluthrin	0.070					87.0					
A Committee of the Comm	ND ND	< 1.000	-	1.740	2.000	100000000000000000000000000000000000000		-			
Cypermethrin	ND	< 1.000		2.110	2.000	105.5	70 - 130				
Daminozide	ND	< 1.000		2.050	2.000	102.5	30 - 150				
Diazinon	ND	< 0.100		0.399	0.400	99.8	70 - 130				
Dichlorvos	ND	< 0.500		1.850	2.000	92.5	70 - 130				
Dimethoat	ND	< 0.100		0.410	0.400	102.5	70 - 130				
Ethoprophos	ND	< 0.100	ı	0.425	0.400	106.3	70 - 130				
Etofenprox	ND	< 0.100		0.890	0.800	111.3	70 - 130				
Etoxazol	ND	< 0.100	1	0.421	0.400	105.3	70 - 130				
enoxycarb	ND	< 0.100	1	0.432	0.400	108.0	70 - 130	1			
enpyroximat	ND	< 0.100		0.867	0.800	108.4	70 - 130				
ipronil	ND	< 0.100		0.803	0.800	100.4	70 - 130				
lonicamid	ND	< 0.400		1.040	1.000	104.0	70 - 130	1			
Fludioxonil	ND	< 0.100	1	0.885	0.800	110.6	70 - 130				
Hexythiazox	ND	< 0.400		1.060	1.000	106.0	70 - 130				
mazalil	ND I	< 0.100	1	0.399	0.400	99.8	70 - 130	1			
midacloprid	ND	< 0.200	1	0.845	0.800	105.6	70 - 130	1			
Kresoxim-Methyl	ND	< 0.100	1	0.865	0.800	108.1	70 - 130	1			
Malathion	ND	< 0.100		0.408	0.400	102.0	70 - 130				
Metalaxyl	ND	< 0.100	-	0.426	0.400	106.5	70 - 130	1			
Methiocarb	ND ND	< 0.100	+	0.418	0.400	104.5	70 - 130	 			
Methomyl	ND ND	< 0.200	+	0.838	0.800	104.8	70 - 130	1			
MGK 264	ND ND	< 0.100	-	0.414	0.400	103.5	70 - 130	-			
Myclobutanil	ND ND	< 0.100	-	0.421	0.400	105.3	70 - 130	-			
Naled	ND ND	< 0.200	-	1.150	1.000	115.0	70 - 130	l -			
Oxamyl	ND ND	< 0.400	-	1.900	2.000	95.0	70 - 130	1			
Paclobutrazol	ND ND	< 0.400		0.865	0.800	108.1	70 - 130	-			
	ND ND	< 0.200	-	1.230	0.800	108.1 153.8	30 - 150	Q1			
Parathion Methyl Permethrin	ND ND	< 0.200		0.448	0.800	112.0	70 - 130	Į QI			
Permethrin	ND ND	< 0.100	-	0.448	0.400	105.8					
								-			
Piperonyl butoxide	ND ND	< 1.000		2.330	2.000	116.5	70 - 130				
Prallethrin	ND ND	< 0.200		0.857	0.800	107.1	70 - 130				
Propiconazole	ND	< 0.200		0.810	0.800	101.3	70 - 130				
ropoxur	ND	< 0.100		0.406	0.400	101.5	70 - 130				
Pyrethrins	ND	< 0.500		0.326	0.284	114.8	70 - 130				
Pyridaben	ND	< 0.100		0.498	0.400	124.5	70 - 130				
pinosad	ND	< 0.100		0.424	0.388	109.3	70 - 130				
piromesifen	ND	< 0.100		0.413	0.400	103.3	70 - 130				
Spirotetramat	ND	< 0.100		0.410	0.400	102.5	70 - 130				
piroxamine	ND	< 0.100		0.897	0.800	112.1	70 - 130				
l'ebuconazol	ND	< 0.200		0.844	0.800	105.5	70 - 130	1			
hiacloprid	ND	< 0.100		0.411	0.400	102.8	70 - 130	1			
Thiamethoxam	ND	< 0.100		0.428	0.400	107.0	70 - 130				
Frifloxystrobin	ND ND	< 0.100	1	0.431	0.400	107.8	70 - 130	1			





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Revision: 0.01 Control: CFL-C22 Revised: 12/4/2018 Effective: 12/4/2018

Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 1	15662		Units:	mg/Kg	Batch ID: 1906643						
Matrix Spike/Matrix	Spike Duplic	ate Reco	veries			Si	ample ID:	19-008698-0004			
Analyte	Result	MS Res	MSD Res	Spike	RP	D%		MSD % Rec	Limits	Notes	
Acephate	0.000	0.962	0.919	1.000	4.6	< 30	96.2	91.9	50 - 150	T	
Acequinocyl	0.000	4.750	4.310	4.000	9.7	< 30	118.8	107.8	50 - 150		
Acetamiprid	0.000	0.426	0.361	0.400	16.5	< 30	106.5	90.3	50 - 150		
Aldicarb	0.000	0.844	0.744	0.800	12.6	< 30	105.5	93.0	50 - 150	i –	
Abamectin	0.000	0.915	0.867	1.000	5.4	< 30	91.5	86.7	50 - 150		
Azoxystrobin	0.000	0.371	0.356	0.400	4.1	< 30	92.8	89.0	50 - 150		
Bifenazate	0.000	0.406	0.375	0.400	7.9	< 30	101.5	93.8	50 - 150		
Bifenthrin	0.034	0.538	0.478	0.400	11.8	< 30	126.1	111.1	50 - 150		
Boscalid	0.000	0.747	0.624	0.800	17.9	< 30	93.4	78.0	50 - 150		
Carbaryl	0.000	0.425	0.354	0.400	18.2	< 30	106.3	88.5	50 - 150	<u> </u>	
Carbofuran	0.000	0.446	0.385	0.400	14.7	< 30	111.5	96.3	50 - 150		
Chlorantraniliprol	0.000	0.421	0.362	0.400	15.1	< 30	105.3	90.5	50 - 150	 	
Chlorfenapyr	0.000	1.780	1.450	2.000	20.4	< 30	89.0	72.5	50 - 150	-	
Chlorpyrifos	0.028	0.369	0.341	0.400	7.9	< 30	85.3	78.3	50 - 150	-	
Clofentezine	0.000	0.435	0.341	0.400	13.2	< 30	108.8	95.3	50 - 150	-	
Cyfluthrin	0.000	2.370	2.010	2.000	16.4	< 30	118.5	100.5	30 - 150	<u> </u>	
Cypermethrin	0.000	2.030	1.770	2.000	13.7	< 30	99.9	86.9	50 - 150	-	
Daminozide	0.032	2.030	1.820	2.000	12.9	< 30	103.5	91.0	30 - 150	-	
Diazinon	0.000	0.418	0.347	0.400	18.6	< 30	103.5	86.8	50 - 150	-	
Diazinon Dichlorvos	0.000	1.940	1.680	2.000	14.4	< 30	97.0	84.0	50 - 150	-	
										_	
Dimethoat	0.000	0.409	0.366	0.400	11.1	< 30	102.3	91.5			
Ethoprophos	0.000	0.436	0.371	0.400	16.1	< 30	109.0	92.8	50 - 150		
tofenprox	0.000	0.928	0.831	0.800	11.0	< 30	116.0	103.9	50 - 150		
toxazol	0.000	0.440	0.394	0.400	11.0	< 30	110.0	98.5	50 - 150		
enoxycarb	0.000	0.427	0.381	0.400	11.4	< 30	106.8	95.3	50 - 150		
enpyroximat	0.000	0.827	0.707	0.800	15.6	< 30	103.4	88.4	50 - 150		
ipronil	0.000	0.954	0.881	0.800	8.0	< 30	119.3	110.1	50 - 150		
Flonicamid	0.000	1.090	0.912	1.000	17.8	< 30	109.0	91.2	50 - 150	l	
ludioxonil	0.000	0.820	0.769	0.800	6.4	< 30	102.5	96.1	50 - 150		
Hexythiazox	0.000	1.090	0.920	1.000	16.9	< 30	109.0	92.0	50 - 150		
mazalil	0.000	0.356	0.305	0.400	15.4	< 30	89.0	76.3	50 - 150		
midacloprid	0.000	0.828	0.733	0.800	12.2	< 30	103.5	91.6	50 - 150		
Kresoxim-Methyl	0.000	0.861	0.763	0.800	12.1	< 30	107.6	95.4	50 - 150		
Malathion	0.000	0.439	0.393	0.400	11.1	< 30	109.8	98.3	50 - 150		
Metalaxyl	0.000	0.418	0.361	0.400	14.6	< 30	104.5	90.3	50 - 150		
Methiocarb	0.000	0.426	0.371	0.400	13.8	< 30	106.5	92.8	50 - 150	l	
Methomyl	0.000	0.827	0.657	0.800	22.9	< 30	103.4	82.1	50 - 150	1	
MGK 264	0.000	0.486	0.424	0.400	13.6	< 30	121.5	106.0	50 - 150		
Myclobutanil	0.000	0.417	0.358	0.400	15.2	< 30	104.3	89.5	50 - 150		
Naled	0.000	1.160	0.954	1.000	19.5	< 30	116.0	95.4	50 - 150	l	
Oxamyl	0.000	1.910	1.670	2.000	13.4	< 30	95.5	83.5	50 - 150		
Paclobutrazol	0.000	0.852	0.740	0.800	14.1	< 30	106.5	92.5	50 - 150		
Parathion Methyl	0.002	0.880	0.859	0.800	2.4	< 30	109.8	107.2	30 - 150	1	
Permethrin	0.000	0.435	0.348	0.400	22.2	< 30	108.8	87.0	50 - 150		
Phosmet	0.000	0.421	0.374	0.400	11.8	< 30	105.3	93.5	50 - 150		
Piperonyl butoxide	0.000	2.330	2.020	2.000	14.3	< 30	116.5	101.0	50 - 150		
Prallethrin	0.000	1.170	1.110	0.800	5.3	< 30	146.3	138.8	50 - 150	 	
Propiconazole	0.000	0.849	0.759	0.800	11.2	< 30	106.1	94.9	50 - 150	l	
Propoxur	0.006	0.419	0.369	0.400	12.7	< 30	103.2	90.7	50 - 150	-	
Pyrethrins	0.014	0.413	0.303	0.284	13.8	< 30	90.5	78.2	50 - 150	\vdash	
Pyridaben	0.000	0.453	0.230	0.400	17.8	< 30	113.3	94.8	50 - 150	<u> </u>	
Spinosad	0.000	0.455	0.379	0.400	12.6	< 30	102.6	90.5	50 - 150	<u> </u>	
	0.000	0.523	0.351	0.400	13.9	< 30	130.8	113.8	50 - 150	-	
piromesifen					8.4					-	
pirotetramat	0.000	0.485	0.446	0.400		< 30	121.3	111.5			
piroxamine	0.000	0.887	0.797	0.800	10.7	< 30	110.9	99.6	50 - 150		
ebuconazol	0.011	0.875	0.743	0.800	16.3	< 30	108.1	91.6	50 - 150		
hiacloprid	0.000	0.402	0.368	0.400	8.8	< 30	100.5	92.0	50 - 150		
hiamethoxam	0.000	0.432	0.380	0.400	12.8	< 30	108.0	95.0	50 - 150		
Trifloxystrobin	0.000	0.429	0.369	0.400	9.8	< 30	107.3	92.3	50 - 150	1	





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Laboratory Quality Control Results

EPA 5021		, ,		ty Contro		Ba	tch ID:	190665	54			
Method Blank					Laborator	ry Cont	rol Sa	mple				
Analyte	Result		LOQ	Notes	Result	Spike		% Rec	-	Lim	its	Notes
Propane	ND	<	200		911	1200	μg/g	75.9	70	-	130	
Isobutane	ND	<	200		1210	1570	μg/g	77.1	70	-	130	
Butane	ND	<	200		1210	1570	μg/g	77.1	70	-	130	
2,2-dimethylpropane	ND	<	200		1550	1980	μg/g	78.3	70	-	130	
Methanol	ND	٧	200		2220	2390	μg/g	92.9	70	-	130	
Ethylene Oxide	ND	٧	30		94.7	119	μg/g	79.6	70	-	130	
2-Methylbutane	ND	<	200		2380	2430	μg/g	97.9	70	-	130	
n-Pentane	ND	«	200		2320	2380	μg/g	97.5	70	-	130	
Ethanol	ND	٧	200		2180	2400	μg/g	90.8	70	-	130	
Ethyl Ether	ND	٧	200		2310	2430	μg/g	95.1	70		130	
2,2-Dimethylbutane	ND	٧	30		648	620	μg/g	104.5	70		130	
Acetone	ND	٧	200		2330	2380	μg/g	97.9	70		130	
Isopropyl alcohol	ND	<	200		2260	2380	μg/g	95.0	70	-	130	
Acetonitrile	ND	<	100		886	919	μg/g	96.4	70	-	130	
2,3-Dimethylbutane	ND	<	30		283	303	μg/g	93.4	70	-	130	
Dichloromethane	ND	<	200		938	948	μg/g	98.9	70	-	130	
2-Methylpentane	ND	<	30		254	293	μg/g	86.7	70	-	130	
3-Methylpentane	ND	٧	30		303	314	μg/g	96.5	70	-	130	
Hexane	ND	<	30		282	297	μg/g	94.9	70		130	
Ethyl acetate	ND	<	200		2210	2370	μg/g	93.2	70	-	130	
2-Butanol	ND	<	200		2100	2410	μg/g	87.1	70		130	
Tetrahydrofuran	ND	<	100		874	943	μg/g	92.7	70	-	130	
Cyclohexane	ND	<	200		2350	2370	μg/g	99.2	70	-	130	
Benzene	ND	<	1		38.3	38.4	μg/g	99.7	70	-	130	
Isopropyl Acetate	ND	<	200		2250	2420	μg/g	93.0	70	-	130	
Heptane	ND	<	200		2320	2380	μg/g	97.5	70	-	130	
1,4-Dioxane	ND	<	100		907	933	μg/g	97.2	70	-	130	
2-Ethoxyethanol	ND	<	30		2110	2370	μg/g	89.0	70	-	130	
Ethylene Glycol	ND	<	200		900	934	μg/g	96.4	70	-	130	
Toluene	ND	<	200		885	937	μg/g	94.5	70	-	130	
Ethylbenzene	ND	<	200		1740	1920	μg/g	90.6	70	-	130	
m,p-Xylene	ND	<	200		1760	1880	μg/g	93.6	70	-	130	
o-Xylene	ND	<	200		1720	1910	μg/g	90.1	70	-	130	
Cumene	ND	٧	30		328	368	μg/g	89.1	70	-	130	





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QC - Sample Duplicate Sample ID: 19-008438-0005

Analyte		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	цд/д	0.0	< 20	Acceptable	İ
2,2-dimethylpropane	ND	ND	200	µg/g	0.0	< 20	Acceptable	İ
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	
n-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	
Isopropyl alcohol	ND	ND	200	μg/g	0.0	< 20	Acceptable	
Acetonitrile	ND	ND	100	μg/g	0.0	< 20	Acceptable	
2,3-Dimethylbutane	ND	ND	30	μg/g	0.0	< 20	Acceptable	
Dichloromethane	ND	ND	200	μg/g	0.0	< 20	Acceptable	
2-Methylpentane	ND	ND	30	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	
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	
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	
1,4-Dioxane	ND	ND	100	μg/g	0.0	< 20	Acceptable	
2-Ethoxyethanol	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Ethylene Glycol	ND	ND	200	μg/g	0.0	< 20	Acceptable	
Toluene	ND	ND	200	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	

Abbreviations

ND - None Detected at or above MRL

RPD - Relative Percent Difference

LOQ - Limit of Quantitation
* Screening only

Q1 Quality Control result biased high. Only non detect samples reported.

Units of Measure:

μg/g- Microgram per gram or ppm

mg/Kg - Milligrams per Kilogram Aw- Water Activity unit





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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	Quantitaion level raised due to matrix interference.					
В	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.					