



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

This is an amended version of the report# 077724-00.  
Reason: Updated Batch ID.

**Product identity:** Sentia Pet Dist Drops 7.9.19 750mg      **Client/Metric ID:** HDTO-1146  
**Laboratory ID:** 19-008079-0001      **Sample Date:** 07/09/19 14:00

**Summary**

**Potency:**

Analyte	Result	Limits	Units	LOQ	
CBC <sup>1</sup>	0.0891		%	0.09	CBD-Total per 1g      32.0 mg/1g
CBD	3.19		%	0.09	THC-Total per 1g      < 1.886 mg/1g
					(Reported in milligrams per serving)
Analyte per 1g	Result	Limits	Units	LOQ	
CBD per 1g	32.0		mg/1g	1.00	

**Residual Solvents:**

All analytes passing and less than LOQ.

**Pesticides:**

All analytes passing and less than LOQ.

**Metals:**

Less than LOQ for all analytes.



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**Customer:** Sentia Wellness  
3931 NE Columbia Blvd  
Portland Oregon 97211  
United States

**Product identity:** Sentia Pet Dist Drops 7.9.19 750mg

**Client/Metric ID:** HDTO-1146

**Sample Date:** 07/09/19 14:00

**Laboratory ID:** 19-008079-0001

**Relinquished by:** Sentia Wellness

**Temp:** 25.4 °C

**Weight Received:** 16 g

**Serving Size #1:** 1.003 g

### Sample Results

Potency		Batch: 1906200					
Analyte	Result	Limits	Units	LOQ	Analyze	Method	Notes
CBC <sup>†</sup>	0.0891		%	0.0871	07/10/19	J AOAC 2015 V98-6	
CBC-A <sup>†</sup>	< LOQ		%	0.0871	07/10/19	J AOAC 2015 V98-6	
CBC-Total <sup>†</sup>	0.0891		%	0.164	07/16/19	J AOAC 2015 V98-6	
CBD	3.19		%	0.0871	07/10/19	J AOAC 2015 V98-6	
CBD-A	< LOQ		%	0.0871	07/10/19	J AOAC 2015 V98-6	
CBD-Total	3.19		%	0.164	07/16/19	J AOAC 2015 V98-6	
CBDV <sup>†</sup>	< LOQ		%	0.0871	07/10/19	J AOAC 2015 V98-6	
CBDV-A <sup>†</sup>	< LOQ		%	0.0871	07/10/19	J AOAC 2015 V98-6	
CBDV-Total <sup>†</sup>	< LOQ		%	0.163	07/16/19	J AOAC 2015 V98-6	
CBG <sup>†</sup>	< LOQ		%	0.0871	07/10/19	J AOAC 2015 V98-6	
CBG-A <sup>†</sup>	< LOQ		%	0.0871	07/10/19	J AOAC 2015 V98-6	
CBG-Total <sup>†</sup>	< LOQ		%	0.163	07/16/19	J AOAC 2015 V98-6	
CBL <sup>†</sup>	< LOQ		%	0.0871	07/10/19	J AOAC 2015 V98-6	
CBN	< LOQ		%	0.0871	07/10/19	J AOAC 2015 V98-6	
Δ8-THC <sup>†</sup>	< LOQ		%	0.0871	07/10/19	J AOAC 2015 V98-6	
Δ9-THC	< LOQ		%	0.0871	07/10/19	J AOAC 2015 V98-6	
THC-A	< LOQ		%	0.0871	07/10/19	J AOAC 2015 V98-6	
THC-Total	< LOQ		%	0.164	07/16/19	J AOAC 2015 V98-6	
THCV <sup>†</sup>	< LOQ		%	0.0871	07/10/19	J AOAC 2015 V98-6	
THCV-A <sup>†</sup>	< LOQ		%	0.0871	07/10/19	J AOAC 2015 V98-6	
THCV-Total <sup>†</sup>	< LOQ		%	0.163	07/16/19	J AOAC 2015 V98-6	

Potency per 1g		Batch: 1906200					
Analyte	Result	Limits	Units	LOQ	Analyze	Method	Notes
CBC per 1g <sup>†</sup>	< LOQ		mg/1g	1.00	07/16/19	J AOAC 2015 V98-6	
CBC-A per 1g <sup>†</sup>	< LOQ		mg/1g	1.00	07/16/19	J AOAC 2015 V98-6	



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Potency per 1g		Batch: 1906200					
Analyte	Result	Limits	Units	LOQ	Analyze	Method	Notes
CBC-Total per 1g <sup>†</sup>	< LOQ		mg/1g	1.89	07/16/19	J AOAC 2015 V98-6	
CBD per 1g	32.0		mg/1g	1.00	07/16/19	J AOAC 2015 V98-6	
CBD-A per 1g	< LOQ		mg/1g	1.00	07/16/19	J AOAC 2015 V98-6	
CBD-Total per 1g	32.0		mg/1g	1.89	07/16/19	J AOAC 2015 V98-6	
CBDV per 1g <sup>†</sup>	< LOQ		mg/1g	1.00	07/16/19	J AOAC 2015 V98-6	
CBDV-A per 1g <sup>†</sup>	< LOQ		mg/1g	1.00	07/16/19	J AOAC 2015 V98-6	
CBDV-Total per 1g <sup>†</sup>	< LOQ		mg/1g	1.88	07/16/19	J AOAC 2015 V98-6	
CBG per 1g <sup>†</sup>	< LOQ		mg/1g	1.00	07/16/19	J AOAC 2015 V98-6	
CBG-A per 1g <sup>†</sup>	< LOQ		mg/1g	1.00	07/16/19	J AOAC 2015 V98-6	
CBG-Total per 1g <sup>†</sup>	< LOQ		mg/1g	1.89	07/16/19	J AOAC 2015 V98-6	
CBL per 1g <sup>†</sup>	< LOQ		mg/1g	1.00	07/16/19	J AOAC 2015 V98-6	
CBN per 1g	< LOQ		mg/1g	1.00	07/16/19	J AOAC 2015 V98-6	
Δ8-THC per 1g <sup>†</sup>	< LOQ		mg/1g	1.00	07/16/19	J AOAC 2015 V98-6	
Δ9-THC per 1g	< LOQ		mg/1g	1.00	07/16/19	J AOAC 2015 V98-6	
THC-A per 1g	< LOQ		mg/1g	1.00	07/16/19	J AOAC 2015 V98-6	
THC-Total per 1g	< LOQ		mg/1g	1.89	07/16/19	J AOAC 2015 V98-6	
THCV per 1g <sup>†</sup>	< LOQ		mg/1g	1.00	07/16/19	J AOAC 2015 V98-6	
THCV-A per 1g <sup>†</sup>	< LOQ		mg/1g	1.00	07/16/19	J AOAC 2015 V98-6	
THCV-Total per 1g <sup>†</sup>	< LOQ		mg/1g	1.88	07/16/19	J AOAC 2015 V98-6	

Solvents		Method EPA5021A				Units μg/g	Batch 1906109	Analyze 07/10/19 11:20 AM			
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
1,4-Dioxane	< LOQ	380	100	pass		2-Butanol	< LOQ	5000	200	pass	
2-Ethoxyethanol	< LOQ	160	30.0	pass		2-Methylbutane	< 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 Batch 1906201 Analyze 07/12/19 09:36 AM											
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	< 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	< 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	Analyze	Method	Notes	
Arsenic	< LOQ		mg/kg	0.0493	1906294	07/15/19	AOAC 2013.06		X
Cadmium	< LOQ		mg/kg	0.0493	1906294	07/15/19	AOAC 2013.06		X
Lead	< LOQ		mg/kg	0.0493	1906294	07/15/19	AOAC 2013.06		X
Mercury	< LOQ		mg/kg	0.0247	1906294	07/15/19	AOAC 2013.06		X

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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**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 = µg/g divided by 10,000

**Glossary of Qualifiers**

X: Not ORELAP accredited.

Approved Signatory

Derrick Tanner  
General Manager



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

Laboratory Quality Control Results									
EPA 5021				Batch ID: 1906109					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		1070	1200	µg/g	89.2	70 - 130	
Isobutane	ND	< 200		1450	1570	µg/g	92.4	70 - 130	
Butane	ND	< 200		1440	1570	µg/g	91.7	70 - 130	
2,2-dimethylpropane	ND	< 200		1740	1980	µg/g	87.9	70 - 130	
Methanol	ND	< 200		2150	2390	µg/g	90.0	70 - 130	
Ethylene Oxide	ND	< 30		104	119	µg/g	87.4	70 - 130	
2-Methylbutane	ND	< 200		2340	2430	µg/g	96.3	70 - 130	
n-Pentane	ND	< 200		2240	2380	µg/g	94.1	70 - 130	
Ethanol	ND	< 200		2180	2400	µg/g	90.8	70 - 130	
Ethyl Ether	ND	< 200		2260	2430	µg/g	93.0	70 - 130	
2,2-Dimethylbutane	ND	< 30		626	620	µg/g	101.0	70 - 130	
Acetone	ND	< 200		2250	2380	µg/g	94.5	70 - 130	
Isopropyl alcohol	ND	< 200		2150	2380	µg/g	90.3	70 - 130	
Ethyl Formate	ND	< 500		2360	2440	µg/g	96.7	70 - 130	
Acetonitrile	ND	< 100		820	919	µg/g	89.2	70 - 130	
Methyl Acetate	ND	< 500		2370	2450	µg/g	96.7	70 - 130	
2,3-Dimethylbutane	ND	< 30		253	303	µg/g	83.5	70 - 130	
Dichloromethane	ND	< 200		905	948	µg/g	95.5	70 - 130	
2-Methylpentane	ND	< 30		278	293	µg/g	94.9	70 - 130	
MTBE	ND	< 500		2320	2440	µg/g	95.1	70 - 130	
3-Methylpentane	ND	< 30		280	314	µg/g	89.2	70 - 130	
Hexane	ND	< 30		278	297	µg/g	93.6	70 - 130	
1-Propanol	ND	< 500		1990	2350	µg/g	84.7	70 - 130	
Methylethylketone	ND	< 500		2360	2400	µg/g	98.3	70 - 130	
Ethyl acetate	ND	< 200		2230	2370	µg/g	94.1	70 - 130	
2-Butanol	ND	< 200		2270	2410	µg/g	94.2	70 - 130	
Tetrahydrofuran	ND	< 100		680	943	µg/g	72.1	70 - 130	
Cyclohexane	ND	< 200		1840	2370	µg/g	77.6	70 - 130	
2-methyl-1-propanol	ND	< 500		1950	2400	µg/g	81.3	70 - 130	
Benzene	ND	< 1		27.7	38.4	µg/g	72.1	70 - 130	
Isopropyl Acetate	ND	< 200		1750	2420	µg/g	72.3	70 - 130	
Heptane	ND	< 200		1720	2380	µg/g	72.3	70 - 130	
1-Butanol	ND	< 500		1900	2370	µg/g	80.2	70 - 130	
Propyl Acetate	ND	< 500		2190	2470	µg/g	88.7	70 - 130	
1,4-Dioxane	ND	< 100		918	933	µg/g	98.4	70 - 130	
2-Ethoxyethanol	ND	< 30		2090	2370	µg/g	88.2	70 - 130	
Methylisobutylketone	ND	< 500		2110	2460	µg/g	85.8	70 - 130	
3-Methyl-1-butanol	ND	< 500		1950	2400	µg/g	81.3	70 - 130	
Ethylene Glycol	ND	< 200		1040	934	µg/g	111.3	70 - 130	
Toluene	ND	< 200		912	937	µg/g	97.3	70 - 130	
Isobutyl Acetate	ND	< 500		2120	2450	µg/g	86.5	70 - 130	
1-Pentanol	ND	< 500		1860	2440	µg/g	76.2	70 - 130	
Butyl Acetate	ND	< 500		2080	2750	µg/g	75.6	70 - 130	
Ethylbenzene	ND	< 200		1580	1920	µg/g	82.3	70 - 130	
m,p-Xylene	ND	< 200		1790	1880	µg/g	95.2	70 - 130	
o-Xylene	ND	< 200		1780	1910	µg/g	93.2	70 - 130	
Cumene	ND	< 30		338	368	µg/g	91.8	70 - 130	
Anisole	ND	< 500		2020	2450	µg/g	82.4	70 - 130	



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QC - Sample Duplicate Sample ID: 19-007807-0004

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	
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	
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	200 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	200 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	

**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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Laboratory Quality Control Results

JAOAC2015 V98-6		Batch ID: 1906200								
Laboratory Control Sample										
Analyte	Result	Spike	Units	% Rec	Limits		Evaluation	Notes		
CBDV-A	0.202	0.2	%	101	85	- 115	Acceptable			
CBDV	0.201	0.2	%	101	85	- 115	Acceptable			
CBD-A	0.195	0.2	%	98	85	- 115	Acceptable			
CBG-A	0.206	0.2	%	103	85	- 115	Acceptable			
CBG	0.209	0.2	%	105	85	- 115	Acceptable			
CBD	0.211	0.2	%	106	85	- 115	Acceptable			
THCV	0.198	0.2	%	99	85	- 115	Acceptable			
THCV-A	0.194	0.2	%	97	85	- 115	Acceptable			
CBN	0.209	0.2	%	105	85	- 115	Acceptable			
THC	0.201	0.2	%	101	85	- 115	Acceptable			
D8THC	0.191	0.2	%	96	85	- 115	Acceptable			
CBL	0.195	0.2	%	98	85	- 115	Acceptable			
CBC	0.201	0.2	%	101	85	- 115	Acceptable			
THCA	0.207	0.2	%	104	85	- 115	Acceptable			
CBCA	0.186	0.2	%	93	85	- 115	Acceptable			

Method Blank

Analyte	Result	LOC	Units	Limits		Evaluation	Notes			
CBDV-A	ND	0.0878	%	< 0.0878		Acceptable				
CBDV	ND	0.0878	%	< 0.0878		Acceptable				
CBD-A	ND	0.0878	%	< 0.0878		Acceptable				
CBG-A	ND	0.0878	%	< 0.0878		Acceptable				
CBG	ND	0.0878	%	< 0.0878		Acceptable				
CBD	ND	0.0878	%	< 0.0878		Acceptable				
THCV	ND	0.0878	%	< 0.0878		Acceptable				
THCV-A	ND	0.0878	%	< 0.0878		Acceptable				
CBN	ND	0.0878	%	< 0.0878		Acceptable				
THC	ND	0.0878	%	< 0.0878		Acceptable				
D8THC	ND	0.0878	%	< 0.0878		Acceptable				
CBL	ND	0.0878	%	< 0.0878		Acceptable				
CBC	ND	0.0878	%	< 0.0878		Acceptable				
THCA	ND	0.0878	%	< 0.0878		Acceptable				
CBCA	ND	0.0878	%	< 0.0878		Acceptable				

**Abbreviations**

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

**Units of Measure:**

%- Percent





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

JAOAC2015 V98-6		Batch ID: 1906200						
Sample Duplicate		Sample ID: 19-0080790001						
Analyte	Result	Org. Result	LOC	Units	RPD	Limits	Evaluation	Notes
CBDV-A	ND	ND	0.0871	%	0	< 20	Acceptable	
CBDV	ND	ND	0.0871	%	0	< 20	Acceptable	
CBD-A	ND	ND	0.0871	%	0	< 20	Acceptable	
CBG-A	ND	ND	0.0871	%	0	< 20	Acceptable	
CBG	ND	ND	0.0871	%	0	< 20	Acceptable	
CBD	3.18	3.19	0.0871	%	0.31	< 20	Acceptable	
THCV	ND	ND	0.0871	%	0	< 20	Acceptable	
THCV-A	ND	ND	0.0871	%	0	< 20	Acceptable	
CBN	ND	ND	0.0871	%	0	< 20	Acceptable	
THC	ND	ND	0.0871	%	0	< 20	Acceptable	
Δ8THC	ND	ND	0.0871	%	0	< 20	Acceptable	
CBL	ND	ND	0.0871	%	0	< 20	Acceptable	
CBC	0.0891	0.0891	0.0871	%	0	< 20	Acceptable	
THCA	ND	ND	0.0871	%	0	< 20	Acceptable	
CBCA	ND	ND	0.0871	%	0	< 20	Acceptable	

**Abbreviations**

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

**Units of Measure:**

%- Percent



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

Revision: 0.01 Control: CFL-C2  
Revised: 12/4/2018 Effective: 12/4/2018

Laboratory Pesticide Quality Control Results

AOAC2007.1 & EN 15662		Units: mg/Kg		Batch ID: 1906201				
Method Blank			Laboratory Control Sample					
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spke	LCS% Re	Limits	Notes
Acephate	ND	< 0.200		0.856	1.000	85.6	70 - 130	
Acequinocyl	ND	< 1.000		4.280	4.000	106.5	70 - 130	
Acetamiprid	ND	< 0.100		0.422	0.400	105.5	70 - 130	
Aldicarb	ND	< 0.200		0.844	0.800	105.5	70 - 130	
Abamectin	ND	< 0.288		1.000	1.000	100.0	70 - 130	
Azoxystrobin	ND	< 0.100		0.474	0.400	118.5	70 - 130	
Bifenazate	ND	< 0.100		0.431	0.400	107.8	70 - 130	
Bifenthrin	ND	< 0.100		0.483	0.400	120.8	70 - 130	
Boscalid	ND	< 0.100		0.800	0.800	100.0	70 - 130	
Carbaryl	ND	< 0.100		0.413	0.400	103.3	70 - 130	
Carbofuran	ND	< 0.100		0.446	0.400	111.5	70 - 130	
Chlorantraniliprol	ND	< 0.100		0.454	0.400	113.5	70 - 130	
Chlorfenapyr	ND	< 1.000		1.930	2.000	96.5	70 - 130	
Chlorpyrifos	ND	< 0.100		0.500	0.400	125.0	70 - 130	
Cofentazine	ND	< 0.100		0.416	0.400	104.0	70 - 130	
Cyfluthrin	ND	< 1.000		1.980	2.000	99.0	30 - 150	
Cypermethrin	ND	< 1.000		2.300	2.000	115.0	70 - 130	
Daminozide	ND	< 1.000		2.120	2.000	106.0	30 - 150	
Diazinon	ND	< 0.100		0.433	0.400	108.3	70 - 130	
Dichlorvos	ND	< 0.500		2.180	2.000	109.0	70 - 130	
Dimethoat	ND	< 0.100		0.411	0.400	102.8	70 - 130	
Ethiofoprofos	ND	< 0.100		0.427	0.400	106.8	70 - 130	
Etofenprox	ND	< 0.100		0.933	0.800	116.6	70 - 130	
Etoxaol	ND	< 0.100		0.432	0.400	108.0	70 - 130	
Fenoxycarb	ND	< 0.100		0.434	0.400	108.5	70 - 130	
Fenpyroximat	ND	< 0.100		0.923	0.800	115.4	70 - 130	
Fipronil	ND	< 0.100		0.886	0.800	110.8	70 - 130	
Fonicamid	ND	< 0.400		1.010	1.000	101.0	70 - 130	
Fludioxonil	ND	< 0.100		0.831	0.800	103.9	70 - 130	
Hexythiazox	ND	< 0.400		1.140	1.000	114.0	70 - 130	
Imazali	ND	< 0.100		0.428	0.400	107.0	70 - 130	
Imidacloprid	ND	< 0.200		0.875	0.800	109.4	70 - 130	
Kiesoxim-Methyl	ND	< 0.100		0.757	0.800	94.6	70 - 130	
Malathion	ND	< 0.100		0.419	0.400	104.8	70 - 130	
Metaxalyl	ND	< 0.100		0.424	0.400	106.0	70 - 130	
Methiocarb	ND	< 0.100		0.427	0.400	106.8	70 - 130	
Methomyl	ND	< 0.200		0.831	0.800	103.9	70 - 130	
MGK 264	ND	< 0.100		0.419	0.400	104.8	70 - 130	
Myclobutanil	ND	< 0.100		0.412	0.400	103.0	70 - 130	
Naled	ND	< 0.200		1.050	1.000	105.0	70 - 130	
Oxamyl	ND	< 0.400		2.030	2.000	101.5	70 - 130	
Padobutrazol	ND	< 0.200		0.900	0.800	112.5	70 - 130	
Parathion Methyl	ND	< 0.200		0.805	0.800	100.6	30 - 150	
Permethrin	ND	< 0.100		0.480	0.400	120.0	70 - 130	
Phosmet	ND	< 0.100		0.455	0.400	113.8	70 - 130	
Piperonyl butoxide	ND	< 1.000		2.400	2.000	120.0	70 - 130	
Prallethrin	ND	< 0.200		0.418	0.400	104.5	70 - 130	
Propiconazole	ND	< 0.200		0.864	0.800	108.0	70 - 130	
Propoxur	ND	< 0.100		0.423	0.400	105.8	70 - 130	
Pyrethrins	ND	< 0.500		0.323	0.284	113.7	70 - 130	
Pyridaben	ND	< 0.100		0.446	0.400	111.5	70 - 130	
Spinosad	ND	< 0.100		0.452	0.388	116.5	70 - 130	
Spiromesifen	ND	< 0.100		0.450	0.400	112.5	70 - 130	
Spirotetramat	ND	< 0.100		0.413	0.400	103.3	70 - 130	
Spiroxamine	ND	< 0.100		0.939	0.800	117.4	70 - 130	
Tebuconazol	ND	< 0.200		0.855	0.800	107.0	70 - 130	
Thiadoprid	ND	< 0.100		0.416	0.400	104.0	70 - 130	
Thiamethoxam	ND	< 0.100		0.421	0.400	105.3	70 - 130	
Trifloxystrobin	ND	< 0.100		0.415	0.400	103.8	70 - 130	



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Revision: 0.01 Control: CFL-C2  
Revised: 12/4/2018 Effective: 12/4/2018

Laboratory Pesticide Quality Control Results

AOAC2007.1 & EN 15662		Units: mg/Kg				Batch ID: 1906201				
Matrix Spike/ Matrix Spike Duplicate Recoveries					Sample ID: 19-008079-0001					
Analyte	Result	MS Res	MSD Res	Spike	RPD%	MS % Re	MSD % Re	Limits	Notes	
Acephate	0.00	0.269	0.257	1.000	4.6	< 30	26.9	25.7	50 - 150	Q
Acequinocyl	0.00	4.930	4.860	4.000	1.4	< 30	123.3	121.5	50 - 150	
Acetamiprid	0.00	0.411	0.434	0.400	5.4	< 30	102.8	108.5	50 - 150	
Aldicarb	0.00	0.841	0.863	0.800	2.6	< 30	105.1	107.9	50 - 150	
Abamectin	0.00	1.260	1.240	1.000	1.6	< 30	126.0	124.0	50 - 150	
Azoxystrobin	0.00	0.478	0.519	0.400	8.2	< 30	119.5	129.8	50 - 150	
Bifenazate	0.00	0.430	0.449	0.400	4.3	< 30	107.5	112.3	50 - 150	
Bifenthrin	0.00	1.290	1.290	0.400	0.0	< 30	322.5	322.5	50 - 150	Q1
Boscalid	0.00	0.857	0.886	0.800	3.3	< 30	107.1	110.8	50 - 150	
Carbaryl	0.00	0.438	0.437	0.400	0.2	< 30	109.5	109.3	50 - 150	
Carbofuran	0.00	0.455	0.464	0.400	2.0	< 30	113.8	116.0	50 - 150	
Chlorantraniliprol	0.00	0.422	0.413	0.400	2.2	< 30	105.5	103.3	50 - 150	
Chlorfenapyr	0.00	2.450	2.230	2.000	9.4	< 30	122.5	111.5	50 - 150	
Chlorpyrifos	0.00	0.735	0.730	0.400	0.6	< 30	198.8	197.5	50 - 150	Q1
Cofentazine	0.00	0.492	0.507	0.400	3.0	< 30	123.0	126.8	50 - 150	
Cyfluthrin	0.00	3.590	3.570	2.000	0.6	< 30	179.5	178.5	30 - 150	Q1
Cypermethrin	0.00	2.420	2.290	2.000	5.5	< 30	121.0	114.5	50 - 150	
Daminozide	0.00	2.100	2.150	2.000	2.4	< 30	105.0	107.5	30 - 150	
Diazinon	0.00	0.434	0.448	0.400	3.2	< 30	108.5	112.0	50 - 150	
Dichlorvos	0.00	2.250	2.210	2.000	1.8	< 30	112.5	110.5	50 - 150	
Dimethoat	0.00	0.409	0.437	0.400	6.6	< 30	102.3	109.3	50 - 150	
Ethoprophos	0.00	0.432	0.440	0.400	1.8	< 30	108.0	110.0	50 - 150	
Etofenprox	0.00	0.986	0.960	0.800	2.7	< 30	123.3	120.0	50 - 150	
Etoxazol	0.00	0.451	0.473	0.400	4.8	< 30	112.8	118.3	50 - 150	
Fenoxycarb	0.00	0.422	0.432	0.400	2.3	< 30	105.5	108.0	50 - 150	
Fenpyroximat	0.00	0.838	0.868	0.800	3.5	< 30	104.8	108.5	50 - 150	
Fipronil	0.00	1.060	1.100	0.800	3.7	< 30	132.5	137.5	50 - 150	
Fonicamid	0.00	0.963	1.120	1.000	15.1	< 30	96.3	112.0	50 - 150	
Fludioxonil	0.00	0.788	0.764	0.800	3.1	< 30	98.5	95.5	50 - 150	
Hexythiazox	0.00	2.670	2.690	1.000	0.7	< 30	267.0	269.0	50 - 150	Q1
Imazali	0.00	0.395	0.387	0.400	2.3	< 30	99.0	96.8	50 - 150	
Imidacloprid	0.00	0.855	0.873	0.800	2.1	< 30	106.9	109.1	50 - 150	
Kiesoxim-Methyl	0.043	0.910	0.922	0.800	1.3	< 30	108.4	109.9	50 - 150	
Malathion	0.00	0.450	0.457	0.400	1.5	< 30	112.5	114.3	50 - 150	
Metaxalyl	0.00	0.418	0.437	0.400	4.4	< 30	104.5	109.3	50 - 150	
Methiocarb	0.008	0.439	0.451	0.400	2.7	< 30	107.6	110.6	50 - 150	
Methomyl	0.00	0.774	0.855	0.800	9.9	< 30	96.8	106.9	50 - 150	
MKG 264	0.00	0.429	0.463	0.400	7.6	< 30	107.3	115.8	50 - 150	
Mydobutanil	0.00	0.446	0.461	0.400	3.5	< 30	111.3	115.3	50 - 150	
Naled	0.00	1.080	1.130	1.000	4.5	< 30	108.0	113.0	50 - 150	
Oxaryl	0.00	2.000	2.120	2.000	5.8	< 30	100.0	106.0	50 - 150	
Padobutrazol	0.00	1.020	1.010	0.800	1.0	< 30	127.5	126.3	50 - 150	
Parathion Methyl	0.00	0.612	0.753	0.800	20.7	< 30	76.5	94.1	30 - 150	
Permethrin	0.008	0.556	0.546	0.400	2.0	< 30	136.9	134.2	50 - 150	
Phosmet	0.00	0.440	0.462	0.400	4.9	< 30	110.0	115.5	50 - 150	
Piperonyl butoxide	0.00	2.430	2.550	2.000	4.8	< 30	121.5	127.5	50 - 150	
Prallethrin	0.005	0.586	0.560	0.400	4.5	< 30	145.2	138.7	50 - 150	
Propiconazole	0.00	0.951	0.935	0.800	1.7	< 30	118.9	116.9	50 - 150	
Propoxur	0.00	0.428	0.421	0.400	1.6	< 30	107.0	105.3	50 - 150	
Pyrethrins	0.011	0.354	0.399	0.284	1.4	< 30	121.0	122.7	50 - 150	
Pyridaben	0.00	0.355	0.357	0.400	0.6	< 30	88.8	89.3	50 - 150	
Spinosad	0.00	0.391	0.399	0.388	2.0	< 30	100.8	102.8	50 - 150	
Spiromesifen	0.00	0.563	0.578	0.400	2.6	< 30	140.8	144.5	50 - 150	
Spirotetramet	0.00	0.380	0.372	0.400	2.1	< 30	95.0	93.0	50 - 150	
Spiroxamine	0.00	0.914	0.909	0.800	0.5	< 30	114.3	113.6	50 - 150	
Tebuconazol	0.00	0.923	0.954	0.800	3.3	< 30	115.4	119.3	50 - 150	
Thiadoprid	0.00	0.417	0.443	0.400	6.0	< 30	104.3	110.8	50 - 150	
Thiamethoxam	0.00	0.399	0.403	0.400	1.0	< 30	99.8	100.8	50 - 150	
Trifloxystrobin	0.00	0.440	0.440	0.400	0.0	< 30	110.0	110.0	50 - 150	

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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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.
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.