



 Job Number:
 19-002293

 Report Number:
 19-002293-00

 Report Date:
 03/07/2019

Purchase Order:

ORELAP#:

Received: 03/05/19 15:31

OR100028

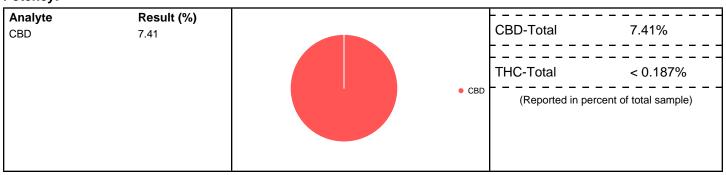
This report cannot be used for ODA, OHA or OLCC compliance requirements.

Product identity: S Cap -Bal Client/Metrc ID:

Laboratory ID: 19-002293-0001 **Sample Date:** 03/05/19 11:00

Summary

Potency:



Residual Solvents:

All analytes passing and less than LOQ.

Pesticides:

All analytes passing and less than LOQ.





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Customer: Cura Wellness

3931 NE Columbia Blvd Portland Oregon 97211

United States

Product identity: S Cap -Bal

Client/Metrc ID:

 Sample Date:
 03/05/19 11:00

 Laboratory ID:
 19-002293-0001

 Relinquished by:
 Erin Harbacek

Temp: 20.7 °C

Sample Results

Potency	Method J AOAC	2015 V98-6		Units %	Batch 1901935	Analyze 03/06/19	10:07 PM
Analyte		Dry LOQ weight	Notes				
CBC [†]	< LOQ	0.0914					
CBC-A [†]	< LOQ	0.0914					
CBC-Total [†]	< LOQ	0.188					
CBD	7.41	0.0914					CBD
CBD-A	< LOQ	0.0914					
CBD-Total	7.41	0.188					
CBDV [†]	< LOQ	0.0914					
CBDV-A [†]	< LOQ	0.0914					
CBDV-Total [†]	< LOQ	0.187					
CBG [†]	< LOQ	0.0914					
CBG-A [†]	< LOQ	0.0914					
CBG-Total [†]	< LOQ	0.188					
CBL [†]	< LOQ	0.0914					
CBN	< LOQ	0.0914					
$\Delta 8\text{-THC}^{\dagger}$	< LOQ	0.0914					
Δ9-THC	< LOQ	0.0914					
THC-A	< LOQ	0.0914					
THC-Total	< LOQ	0.187					
THCV [†]	< LOQ	0.0914					
THCV-A [†]	< LOQ	0.0914					
THCV-Total [†]	< LOQ	0.187					





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Solvents	Method	EPA502	21A			Units µg/g Batch 19	901872	Analyz	ze 03/0	06/19 09:24 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	1 15662 (mod)	Units mg/kg Batc	h 1901924	Analy	ze 03/07/19(9:49 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		Etoxazol	< LOQ	0.20	0.100 pass	
Fenoxycarb	< LOQ	0.20	0.100 pass		Fenpyroximat	< 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	
Spiroxamin	< LOQ	0.40	0.200 pass		Tebuconazol	< 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.

Units of Measure

μg/g = Microgram per gram
 mg/kg = Milligram per kilogram
 % = Percentage of sample
 % wt = μg/g divided by 10,000

Approved Signatory

Derrick Tanner General Manager





Job Number: 19

19-002293

Report Number: Report Date:

19-002293-00 03/07/2019

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12423 NE Whitaker Way Portland OR 972	30 p.503	J-254-17	94	Ca	nnal	ois C					ecor	d			ORELAP	PIXIS Labs (D: OR100028	
Company: Wa Wellness							А	naiys	is ke	quest	.ea					Purchase Order Number:	
Contact: Colin Gallison Address: 3931 Necolumbia	Blia	spunod														Project Number: Project Name:	
Email: CBP COMPLIANCE @ CURSCI Phone: CF addams 7 Fax: Processor's License:	OR 59 compounds	ulti-Residue – 379 compounds		vents	ity			t and Mold	Micro: E.Coli and Total Coliform	s						Report Instructions: Send to State - METRC Email Final Results: Gash/Check/CC/Net 30 Other:	
Field ID Date/Tin		/	Potency	Residual Solvents	Water Activity	Moisture	Terpenes	Micro: Yeast and Mold	Micro: E.Co	Heavy Metals	Mycotoxins	Other	Matrix	Weight	Serving size for edibles		
Scap-BOD 3/5 Scap-Rec 3/5	a V		V										Oil/nia			inside the capsule capsules do not no	
SCap-BOO 3/5 SCap-Rec 3/5	a V		/	V	/								oilfris			inside the capsule.	5
S Cap - Rec / 3/5 1	a 1/		1	1									oil/mix			capsules do not ne	ed testing
S Cap - Res / 3/5 1	a /		_	/									DiVmix				
Collected By: Relinquish	ed By:			Date		Time		Recei	ved by	·:			Date	Tim	e La	b Use Only:	
□Standard (5 day) EY ₩	Har	pac	CRK	3.	5.19	3:	310	4	Mi	10		~	3-	5-19 13	5)3/ or	ent Alias: der Number: oper Container	
(1.5x Standard)									U							mple Condition	
Priority Rush (2 day) (2x Standard)															Sh	mperature: 20.) ipped Via: Clent idence of cooling: □Yes □No	

SUBMISSION OF SAMPLES WITH TESTING REQUIREMENTS TO PIXIS WILL BE UNDERSTOOD TO BE AN AGREEMENT FOR SERVICES IN ACCORDANCE WITH THE CONDITIONS LISTED ON THE BACK OF THIS FORM

Revision: 1.02 Control#: CF023 Effective 01/31/2019 Revised 01/31/2019 www.pixislabs.com

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	Labo	oratory	Quali	ty Contro	Results							
EPA 5021						Ba	tch ID:	190187	2			
Method Blank					Laborator	y Cont	rol San	nple				
Analyte	Result			Notes	Result	Spike	Units	%Rec	L	imi	ts	Notes
Propane	ND	<	200		1290	1170	µg/g	110.3	70	-	130	
Isobutane	ND	<	200		1610	1560	μg/g	103.2	70	-	130	
n-Butane	ND	<	200		1610	1560	μg/g	103.2	70	-	130	
neo-pentane	ND	<	200		2090	1930	μg/g	108.3	70	-	130	
Methanol	ND	<	200		2340	2390	μg/g	97.9	70	-	130	
Ethylene Oxide	ND	<	30		40.2	38.7	μg/g	103.9	70	-	130	
Isopentane	ND	<	200		2390	2430	μg/g	98.4	70	-	130	
Pentane	ND	<	200		2410	2380	μg/g	101.3	70	-	130	
Bhanol	ND	<	200		2270	2400	цд/д	94.6	70	-	130	
Bhyl Bher	ND	<	200		2410	2430	цд/д	99.2	70	-	130	
2,2-Dimethylbutane	ND	<	30		605	620	цд/д	97.6	70	-	130	
Acetone	ND	<	200		2310	2380	цд/д	97.1	70	-	130	
2-Propanol	ND	<	200		2210	2380	цд/д	92.9	70	-	130	
Acetonitrile	ND	<	100		901	919	μg/g	98.0	70	-	130	
2,3-Dimethylbutane	ND	<	30		286	303	μg/g	94.4	70	-	130	
Dichloromethane	ND	<	200		895	948	μg/g	94.4	70	-	130	
2-methylpentane	ND	<	30		290	293	μg/g	99.0	70	-	130	
3-methylpentane	ND	<	30		291	314	μg/g	92.7	70	-	130	
Hexane	ND	<	30		282	297	μg/g	94.9	70	-	130	
Ethyl Acetate	ND	<	200		2240	2370	μg/g	94.5	70	-	130	
2-Butanol	ND	<	200		2280	2410	μg/g	94.6	70	-	130	
Tetrahydrofuran	ND	<	100		907	943	μg/g	96.2	70	-	130	
Cyclohexane	ND	<	200		2230	2370	μg/g	94.1	70	-	130	
Benzene	ND	<	1.0		35.8	38.4	μg/g	93.2	70	-	130	
Isopropyl acetate	ND	٧	200		2270	2420	μg/g	93.8	70		130	
Heptane	ND	٧	200		2210	2380	μg/g	92.9	70		130	
1,4-Dioxane	ND	٧	100		896	933	μg/g	96.0	70		130	
2-Ethoxyethanol	ND	<	30		2210	2370	μg/g	93.2	70	-	130	
Toluene	ND	<	100		874	937	μg/g	93.3	70	-	130	
Ethylene glycol	ND	<	200		689	934	μg/g	73.8	70	-	130	
Ethylbenzene	ND	<	200		1770	1920	μg/g	92.2	70	-	130	
m,p-Xylene	ND	<	200		1770	1880	μg/g	94.1	70		130	
o-Xylene	ND	<	200		1810	1910	μg/g	94.8	70		130	
Qumene Qumene	ND	<	30		350	368	μg/g	95.1	70	-	130	





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QC- Sample Duplicate Sample ID: 19-002165-0001

QC- cample Duplicate	,					cample ib.	13-002 100-0001	
Analyte	Result	Org. Result	LOQ L	Jnits	RPD	Limits	Accept/Fail	Notes
Propane	ND	ND	200 L	ıg/g	0.0	< 20	Acceptable	
Isobutane	ND	ND		ıg/g	0.0	< 20	Acceptable	İ
n-Butane	ND	ND	200 j	ıg/g	0.0	< 20	Acceptable	
neo-pentane	ND	ND	200 j	ıg/g	0.0	< 20	Acceptable	
Methanol	ND	ND	200 j	ıg/g	0.0	< 20	Acceptable	
Ethylene Oxide	ND	ND	30 µ	ıg/g	0.0	< 20	Acceptable	
Isopentane	ND	ND	200 µ	ıg/g	0.0	< 20	Acceptable	
Pentane	ND	ND	200 µ	ıg/g	0.0	< 20	Acceptable	
Ethanol	ND	ND	200 µ	ıg/g	0.0	< 20	Acceptable	
Ethyl Ether	ND	ND	200 µ	ıg/g	0.0	< 20	Acceptable	
2,2-Dimethylbutane	ND	ND	30 j	ıg/g	0.0	< 20	Acceptable	
Acetone	ND	ND	200 J	ıg/g	0.0	< 20	Acceptable	
2-Propanol	ND	ND	200 J	ıg/g	0.0	< 20	Acceptable	
Acetonitrile	ND	ND	100 յ	ıg/g	0.0	< 20	Acceptable	
2,3-Dimethylbutane	ND	ND	30 j	ıg/g	0.0	< 20	Acceptable	
Dichloromethane	ND	ND	200 µ	ıg/g	0.0	< 20	Acceptable	
2-methylpentane	ND	ND	30 j	ıg/g	0.0	< 20	Acceptable	
3-methylpentane	ND	ND	30 j	ıg/g	0.0	< 20	Acceptable	
Hexane	ND	ND	30 j	ıg/g	0.0	< 20	Acceptable	
Ethyl Acetate	ND	ND	200 µ	ıg/g	0.0	< 20	Acceptable	
2-Butanol	ND	ND	200 µ	ıg/g	0.0	< 20	Acceptable	
Tetrahydrofuran	ND	ND		ıg/g	0.0	< 20	Acceptable	
Cyclohexane	ND	ND	200 լ	ıg/g	0.0	< 20	Acceptable	
Benzene	ND	ND	1.0 į	ıg/g	0.0	< 20	Acceptable	
Isopropyl acetate	ND	ND	200 լ	ıg/g	0.0	< 20	Acceptable	
Heptane	ND	ND	200 լ	ıg/g	0.0	< 20	Acceptable	
1,4-Dioxane	ND	ND	100 լ	ıg/g	0.0	< 20	Acceptable	
2-Ethoxyethanol	ND	ND		ıg/g	0.0	< 20	Acceptable	
Toluene	ND	ND	100 լ	ıg/g	0.0	< 20	Acceptable	
Ethylene glycol	ND	ND	200 µ	ıg/g	0.0	< 20	Acceptable	
Ethylbenzene	ND	ND	200 µ	ıg/g	0.0	< 20	Acceptable	
m,p-Xylene	ND	ND	200 µ	ıg/g	0.0	< 20	Acceptable	
o-Xylene	ND	ND	200 µ	ıg/g	0.0	< 20	Acceptable	
Cumene	ND	ND	30 į	лд/д	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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Revision: 0.01 Control: CFL-C22 Revised: 12/4/2018 Effective: 12/4/2018

Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN	15662	62 Units: mg/Kg Batch ID: 1								
Method Blank				Laboratory Cor	ntrol Sample					
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes		
Acephate	ND	< 0.200		1.020	1.000	102.0	70 - 130			
Acequinocyl	ND	< 1.000		3.480	4.000	87.0	70 - 130			
Acetamiprid	ND	< 0.100		0.404	0.400	101.0	70 - 130			
Aldicarb	ND	< 0.200		0.749	0.800	93.6	70 - 130			
Abamectin	ND	< 0.288		1.020	1.000	102.0	70 - 130			
Azoxystrobin	ND	< 0.100		0.423	0.400	105.8	70 - 130			
Bifenazate	ND	< 0.100		0.382	0.400	95.5	70 - 130			
Bifenthrin	ND	< 0.100		0.409	0.400	102.3	70 - 130			
Boscalid	ND	< 0.100		0.692	0.800	86.5	70 - 130	l		
Carbaryl	ND	< 0.100		0.409	0.400	102.3	70 - 130			
Carbofuran	ND	< 0.100		0.432	0.400	108.0	70 - 130			
Chlorantraniliprol	ND	< 0.100		0.400	0.400	100.0	70 - 130	l .		
Chlorfenapyr	ND	< 1.000	1	1.840	2.000	92.0	70 - 130			
Chlorpyrifos	ND	< 0.100		0.355	0.400	88.8	70 - 130			
Clofentezine	ND	< 0.100		0.281	0.400	70.3	70 - 130			
Cyfluthrin	ND	< 1.000		2.070	2.000	103.5	30 - 150			
Cypermethrin	ND	< 1.000		2.130	2.000	106.5	70 - 130			
Daminozide	ND	< 1.000		2.040	2.000	102.0	30 - 150			
Diazinon	ND	< 0.100		0.393	0.400	98.3	70 - 130			
Dichlorvos	ND	< 0.500		1.950	2.000	97.5	70 - 130			
Dimethoat	ND	< 0.100	1	0.397	0.400	99.3	70 - 130			
Ethoprophos	ND	< 0.100	1	0.395	0.400	98.8	70 - 130			
Etofenprox	ND	< 0.100	1	0.924	0.800	115.5	70 - 130			
Etoxazol	ND	< 0.100	1	0.384	0.400	96.0	70 - 130			
Fenoxycarb	ND	< 0.100		0.396	0.400	99.0	70 - 130			
Fenpyroximat	ND	< 0.100	1	0.800	0.800	100.0	70 - 130			
Fipronil	ND	< 0.100		0.786	0.800	98.3	70 - 130			
Flonicamid	ND	< 0.400	1	1.030	1.000	103.0	70 - 130			
Fludioxonil	ND	< 0.100	1	0.838	0.800	104.8	70 - 130			
Hexythiazox	ND	< 0.400	_	0.953	1.000	95.3	70 - 130			
mazalil	ND	< 0.100	1	0.403	0.400	100.8	70 - 130	l –		
Imidacloprid	ND	< 0.200	1	0.774	0.800	96.8	70 - 130			
Kresoxim-Methyl	ND	< 0.100	1	0.786	0.800	98.3	70 - 130	l		
Malathion	ND	< 0.100	1	0.374	0.400	93.5	70 - 130			
Metalaxyl	ND	< 0.100	1	0.383	0.400	95.8	70 - 130			
Methiocarb	ND	< 0.100	1	0.421	0.400	105.3	70 - 130	 		
Methomyl	ND	< 0.200	+	0.856	0.800	107.0	70 - 130	-		
MGK 264	ND	< 0.100	1	0.359	0.400	89.8	70 - 130			
Myclobutanil	ND I	< 0.100	+	0.392	0.400	98.0	70 - 130	 		
Naled	ND	< 0.200	1	1.010	1.000	101.0	70 - 130			
Oxamyl	ND ND	< 0.400	1	2.070	2.000	103.5	70 - 130			
Paclobutrazol	ND ND	< 0.200	1	0.784	0.800	98.0	70 - 130			
Parathion Methyl	ND ND	< 0.200	+	0.805	0.800	100.6	30 - 150	<u> </u>		
Permethrin	ND ND	< 0.100	1	0.422	0.400	105.5	70 - 130			
Phosmet	ND ND	< 0.100	1	0.387	0.400	96.8	70 - 130			
Piperonyl butoxide	ND ND	< 1.000	+	1.480	2.000	74.0	70 - 130			
Prallethrin	ND I	< 0.200	1	0.358	0.400	89.5	70 - 130	1		
Propiconazole	ND	< 0.200	1	0.787	0.800	98.4	70 - 130			
Propoxur	ND	< 0.100	1	0.385	0.400	96.3	70 - 130			
Pyrethrins	ND	< 0.500	+	0.292	0.284	102.8	70 - 130			
Pyridaben	ND ND	< 0.100	1	0.631	0.400	157.8	70 - 130	Q1		
Spinosad	ND ND	< 0.100	1	0.454	0.388	117.0	70 - 130	"		
Spiromesifen	ND ND	< 0.100	+	0.375	0.400	93.8	70 - 130			
Spirotetramat	ND ND	< 0.100	+	0.399	0.400	99.8	70 - 130	<u> </u>		
Spiroxamine	ND ND	< 0.100	+	0.830	0.800	103.8	70 - 130	 		
rebuconazol	ND ND	< 0.200	-	0.803	0.800	100.4	70 - 130	-		
Thiacloprid	ND ND	< 0.100	-	0.402	0.400	100.4	70 - 130	-		
Thiaciopriu Thiamethoxam	ND ND	< 0.100	-	0.402	0.400	100.5	70 - 130	-		
Frifloxystrobin	ND I	< 0.100	-	0.412	0.400	94.0	70 - 130			
THIOXYSTIODIII	טאו	< 0.100	1	0.370	0.400	94.0	70 - 130			





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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 15	662		Units:	mg/Kg				Batch	ID: 19019	124
Matrix Spike/Matrix	Spike Duplic	ate Reco	veries			S	ample ID:	19-00219	1-0002	
Analyte	Result	MS Res	MSD Res	Spike	RP	PD%		MSD % Rec		Notes
Acephate	0.000	0.967	0.983	1.000	1.6	< 30	96.7	98.3	50 - 150	
Acequinocyl	0.000	4.730	4.580	4.000	3.2	< 30	118.3	114.5	50 - 150	
Acetamiprid	0.000	0.408	0.407	0.400	0.2	< 30	102.0	101.8	50 - 150	
Aldicarb	0.000	0.766	0.786	0.800	2.6	< 30	95.8	98.3	50 - 150	
Abamectin	0.000	1.090	1.110	1.000	1.8	< 30	109.0	111.0	50 - 150	
Azoxystrobin	0.000	0.435	0.456	0.400	4.7	< 30	108.8	114.0	50 - 150	
Bifenazate	0.000	0.369	0.380	0.400	2.9	< 30	92.3	95.0	50 - 150	
Bifenthrin	0.245	1.040	1.120	0.400	7.4	< 30	198.8	218.8	50 - 150	Q1
Boscalid	0.000	0.729	0.922	0.800	23.4	< 30	91.1	115.3	50 - 150	
Carbaryl	0.000	0.396	0.390	0.400	1.5	< 30	99.0	97.5	50 - 150	
Carbofuran	0.000	0.430	0.423	0.400	1.6	< 30	107.5	105.8	50 - 150	
Chlorantraniliprol	0.000	0.382	0.399	0.400	4.4	< 30	95.5	99.8	50 - 150	
Chlorfenapyr	1.110	2.100	1.860	2.000	12.1	< 30	49.5	37.5	50 - 150	Q
Chlorpyrifos	0.000	0.115	0.137	0.400	17.5	< 30	28.8	34.3	50 - 150	Q
Clofentezine	0.000	0.352	0.420	0.400	17.6	< 30	88.0	105.0	50 - 150	
Cyfluthrin	0.000	2.040	1.940	2.000	5.0	< 30	102.0	97.0	30 - 150	
Cypermethrin	0.126	2.040	2.120	2.000	3.8	< 30	95.7	99.7	50 - 150	
Daminozide	0.000	1.440	1.530	2.000	6.1	< 30	72.0	76.5	30 - 150	
Diazinon	0.000	0.356	0.408	0.400	13.6	< 30	89.0	102.0	50 - 150	
Dichlorvos	0.000	2.020	2.160	2.000	6.7	< 30	101.0	108.0	50 - 150	
Dimethoat	0.000	0.415	0.412	0.400	0.7	< 30	103.8	103.0	50 - 150	
Ethoprophos	0.000	0.383	0.405	0.400	5.6	< 30	95.8	101.3	50 - 150	
Etofenprox	0.000	0.706	0.813	0.800	14.1	< 30	88.3	101.6	50 - 150	
Etoxazol	0.000	0.367	0.413	0.400	11.8	< 30	91.8	103.3	50 - 150	
Fenoxycarb	0.000	0.362	0.391	0.400	7.7	< 30	90.5	97.8	50 - 150	
Fenpyroximat	0.000	0.786	0.849	0.800	7.7	< 30	98.3	106.1	50 - 150	
Fipronil	0.000	0.624	0.747	0.800	17.9	< 30	78.0	93.4	50 - 150	
Flonicamid	0.000	0.969	0.981	1.000	1.2	< 30	96.9	98.1	50 - 150	
Fludioxonil	0.000	0.759	0.729	0.800	4.0	< 30	94.9	91.1	50 - 150	
Hexythiazox	0.000	0.392	0.438	1.000	11.1	< 30	39.2	43.8	50 - 150	Q
Imazalil	0.000	0.442	0.445	0.400	0.7	< 30	110.5	111.3	50 - 150	
Imidacloprid	0.000	0.868	0.896	0.800	3.2	< 30	108.5	112.0	50 - 150	
Kresoxim-Methyl	0.000	0.754	0.815	0.800	7.8	< 30	94.3	101.9	50 - 150	
Malathion	0.000	0.371	0.381	0.400	2.7	< 30	92.8	95.3	50 - 150	
Metalaxyl	0.000	0.396	0.404	0.400	2.0	< 30	99.0	101.0	50 - 150	
Methiocarb	0.000	0.381	0.417	0.400	9.0	< 30	95.3	104.3	50 - 150	
Methomyl	0.000	0.788	0.846	0.800	7.1	< 30	98.5	105.8	50 - 150	
MGK 264	0.000	0.350	0.354	0.400	1.1	< 30	87.5	88.5	50 - 150	
Myclobutanil	0.000	0.368	0.406	0.400	9.8	< 30	92.0	101.5	50 - 150	
Naled	0.000	1.020	1.020	1.000	0.0	< 30	102.0	102.0	50 - 150	
Oxamyl	0.000	2.010	2.170	2.000	7.7	< 30	100.5	108.5	50 - 150	
Paclobutrazol	0.000	0.746	0.818	0.800	9.2	< 30	93.3	102.3	50 - 150	
Parathion Methyl	0.000	0.632	0.685	0.800	8.0	< 30	79.0	85.6	30 - 150	
Permethrin	0.015	0.491	0.513	0.400	4.4	< 30	119.0	124.5	50 - 150	
Phosmet	0.000	0.415	0.419	0.400	1.0	< 30	103.8	104.8	50 - 150	
Piperonyl butoxide	0.000	1.760	2.090	2.000	17.1	< 30	88.0	104.5	50 - 150	
Prallethrin	0.000	0.407	0.479	0.400	16.3	< 30	101.8	119.8	50 - 150	
Propiconazole	0.000	0.749	0.805	0.800	7.2	< 30	93.6	100.6	50 - 150	
Propoxur	0.000	0.398	0.391	0.400	1.8	< 30	99.5	97.8	50 - 150	
Pyrethrins	0.000	0.303	0.342	0.284	12.1	< 30	106.7	120.4	50 - 150	
Pyridaben	0.000	0.291	0.319	0.400	9.2	< 30	72.8	79.8	50 - 150	
Spinosad	0.000	0.425	0.444	0.388	4.4	< 30	109.5	114.4	50 - 150	
Spiromesifen	0.000	0.304	0.336	0.400	10.0	< 30	76.0	84.0	50 - 150	
Spirotetramat	0.000	0.348	0.376	0.400	7.7	< 30	87.0	94.0	50 - 150	
Spiroxamine	0.000	0.896	0.893	0.800	0.3	< 30	112.0	111.6	50 - 150	
Tebuconazol	0.000	0.686	0.792	0.800	14.3	< 30	85.8	99.0	50 - 150	
Thiacloprid	0.000	0.405	0.412	0.400	1.7	< 30	101.3	103.0	50 - 150	
Thiamethoxam	0.000	0.373	0.418	0.400	11.4	< 30	93.3	104.5	50 - 150	
Trifloxystrobin	0.000	0.300	0.346	0.400	9.7	< 30	75.0	86.5	50 - 150	





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Laboratory Quality Control Results

J AOAC 2015	V98-6			Bat			
Laboratory C	ontrol Sample						
Analyte	Result	Spike	Units	% Rec	Limits	Evaluation	Notes
CBDV-A	0.192	0.2	%	96.0	85 - 115	Acceptable	
CBDV	0.207	0.2	%	104	85 - 115	Acceptable	
CBD-A	0.190	0.2	%	95.0	85 - 115	Acceptable	
CBG-A	0.187	0.2	%	93.5	85 - 115	Acceptable	
CBG	0.204	0.2	%	102	85 - 115	Acceptable	
CBD	0.192	0.2	%	96.0	85 - 115	Acceptable	
THCV	0.192	0.2	%	96.0	85 - 115	Acceptable	
THCVA	0.181	0.2	%	90.5	85 - 115	Acceptable	
CBN	0.207	0.2	%	104	85 - 115	Acceptable	
THC	0.194	0.2	%	97.0	85 - 115	Acceptable	
D8THC	0.190	0.2	%	95.0	85 - 115	Acceptable	
CBL	0.178	0.2	%	89.0	85 - 115	Acceptable	
CBC	0.208	0.2	%	104	85 - 115	Acceptable	
THCA	0.188	0.2	%	94.0	85 - 115	Acceptable	
CBCA	0.179	0.2	%	89.5	85 - 115	Acceptable	

Method Blank

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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 V	98-6							
Sample Duplica	ite				Sam	ple ID: 19-0021	91-0001	
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	0.227	0.230	0.1	%	1.31	< 20	Acceptable	
CBG-A	1.86	1.82	0.1	%	2.17	< 20	Acceptable	
CBG	0.516	0.523	0.1	%	1.35	< 20	Acceptable	
CBD	ND	ND	0.1	%	0	< 20	Acceptable	
THCV	0.174	0.181	0.1	%	3.94	< 20	Acceptable	
THCVA	0.305	0.300	0.1	%	1.65	< 20	Acceptable	
CBN	ND	ND	0.1	%	0	< 20	Acceptable	
THC	27.7	28.0	0.1	%	1.08	< 20	Acceptable	
D8THC	ND	ND	0.1	%	0	< 20	Acceptable	
CBL	ND	ND	0.1	%	0	< 20	Acceptable	
CBC	0.289	0.291	0.1	%	0.690	< 20	Acceptable	
THCA	42.1	41.3	0.1	%	1.92	< 20	Acceptable	
CBCA	0.758	0.724	0.1	%	4.59	< 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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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.