

CONSOLIDATED TEST RESULTS SUMMARY

Please see the following pages for full test results.

BULK SKU CHOC.FS.DS4

BATCH # DK05

PRODUCT NAME Full Spectrum Dark Chocolate Sea Salt **SERVING SIZE** 2 squares(~10 g)

LABORATORY: Columbia Laboratories **OREGON ACCREDITATION:** OR100028

LOQ: Limit Of Quantitation
LOD: Limit Of Detection

1 g = 10⁻³ kg = 10³ mg = 10⁶ µg
1 mg/kg = 1 ppm = 1000 ppb

POTENCY	PER SERVING	PER GRAM	Percent
Cannabidiol (CBD)	45.7 mg/serving	4.57 mg/g	0.457 %
Total THC (d9-THC, THCA)	<LOQ mg/serving	<LOQ mg/g	<LOQ %
Cannabigerol (CBG)	<LOQ mg/serving	<LOQ mg/g	<LOQ %
Cannabinol (CBN)	<LOQ mg/serving	<LOQ mg/g	<LOQ %
Cannabichromene (CBC)	<LOQ mg/serving	<LOQ mg/g	<LOQ %
Tetrahydrocannabinolic Acid (THCA)	<LOQ mg/serving	<LOQ mg/g	<LOQ %
Delta-9-THC (d9-THC)	<LOQ mg/serving	<LOQ mg/g	<LOQ %
Delta-8-THC (d8-THC)	<LOQ mg/serving	<LOQ mg/g	<LOQ %

HEAVY METALS	PER SERVING	PER GRAM	REGULATORY ACTION LEVEL
Arsenic	<LOQ µg/serving	<LOQ µg/g	10 µg/day ^[1]
Cadmium	2.20 µg/serving	0.220 µg/g	4.1 µg/day ^[1]
Lead	<LOQ µg/serving	<LOQ µg/g	3.5 µg/day ^[2]
Mercury	<LOQ µg/serving	<LOQ µg/g	2 µg/day ^[1]

PESTICIDES	REGULATORY ACTION LEVEL
None of the other 59 pesticides tested found above limit of detection in the sample.	10 ppb ^[1]

RESIDUAL SOLVENTS	Results	REGULATORY ACTION LEVEL
Ethanol	<LOQ	
Heptane	<LOQ	
None of the 34 residual solvents tested found above limit of quantitation in the sample.		

MICROBIAL	PASS/FAIL
Yeast & Mold	Pass
Coliform	Pass



1. American Herbal Pharmacopoeia. (2014). Cannabis Inflorescence: Standards of Identity, Analysis, and Quality Control. Washington DC: AHP.

2. US Food and Drug Administration. (2019). Lead in Food, Foodwares, and Dietary Supplements. Washington DC: FDA. US Food and Drug Administration. (2019). Lead in Food, Foodwares, and Dietary Supplements. Washington DC: FDA.



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



Report Number: 21-013871/D002.R000
Report Date: 12/03/2021
ORELAP#: OR100028
Purchase Order:
Received: 11/24/21 14:41

Customer: Etz Hayim Holdings
Product identity: FORM-DK05-CHOC.FS.DS4
Client/Metric ID: .
Laboratory ID: 21-013871-0001

Summary

Potency:

Analyte per 1g	Result	Limits	Units	Status	
CBD per 1g	4.57		mg/1g		CBD Total per 1g 4.57 mg/1g
					THC-Total per 1g <LOQ
(Reported in milligrams per serving)					

Residual Solvents:

All analytes passing and less than LOQ.

Pesticides:

All analytes passing and less than LOQ.

Metals:

Analyte	Result	Units	Limit	Status
Cadmium	0.220	mg/kg		

Microbiology:

Less than LOQ for all analytes.



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Customer: Etz Hayim Holdings
 16427 NE Airport Way
 PORTLAND 97230
 United States of America (USA)

Product identity: FORM-DK05-CHOC.FS.DS4

Client/Metric ID: .

Sample Date:

Laboratory ID: 21-013871-0001

Evidence of Cooling: No

Temp: 20.6 °C

Relinquished by: Client

Serving Size #1: 1 g

Density: 0.8000 g/ml

Sample Results

Potency per 1g					
Method J AOAC 2015 V98-6 (mod)Units mg/se Batch: 2110747 Analyze: 11/30/21 8:08:00 PM					
Analyte	Result	Limits	Units	LOQ	Notes
CBC per 1g [†]	< LOQ		mg/1g	0.0320	
CBC-A per 1g [†]	< LOQ		mg/1g	0.0320	
CBC-Total per 1g [†]	< LOQ		mg/1g	0.0600	
CBD per 1g	4.57		mg/1g	0.0320	
CBD-A per 1g	< LOQ		mg/1g	0.0320	
CBD-Total per 1g	4.57		mg/1g	0.0600	
CBDV per 1g [†]	< LOQ		mg/1g	0.0320	
CBDV-A per 1g [†]	< LOQ		mg/1g	0.0320	
CBDV-Total per 1g [†]	< LOQ		mg/1g	0.0597	
CBE per 1g [†]	< LOQ		mg/1g	0.0320	
CBG per 1g [†]	< LOQ		mg/1g	0.0320	
CBG-A per 1g [†]	< LOQ		mg/1g	0.0320	
CBG-Total per 1g [†]	< LOQ		mg/1g	0.0597	
CBL per 1g [†]	< LOQ		mg/1g	0.0320	
CBL-A per 1g [†]	< LOQ		mg/1g	0.0320	
CBL-Total per 1g [†]	< LOQ		mg/1g	0.0600	
CBN per 1g	< LOQ		mg/1g	0.0320	
CBT per 1g [†]	< LOQ		mg/1g	0.0320	
Δ8-THCV per 1g [†]	< LOQ		mg/1g	0.0320	
Δ8-THC per 1g [†]	< LOQ		mg/1g	0.0320	
Δ9-THC per 1g	< LOQ		mg/1g	0.0320	
exo-THC per 1g [†]	< LOQ		mg/1g	0.0320	
THC-A per 1g	< LOQ		mg/1g	0.0320	
THC-Total per 1g	< LOQ		mg/1g	0.0600	
THCV per 1g [†]	< LOQ		mg/1g	0.0320	
THCV-A per 1g [†]	< LOQ		mg/1g	0.0320	
THCV-Total per 1g [†]	< LOQ		mg/1g	0.0600	
Total Cannabinoids per 1g	4.57		mg/1g		



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Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyze	Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2110622	11/29/21	AOAC 991.14 (Petrifilm)		X
Total Coliforms	< LOQ		cfu/g	10	2110622	11/29/21	AOAC 991.14 (Petrifilm)		X
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2110623	11/30/21	AOAC 2014.05 (RAPID)		X
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2110623	11/30/21	AOAC 2014.05 (RAPID)		X

Solvents Method Residual Solvents by GC/MS Units µg/g Batch 2110727 Analyze 12/01/21 08:48 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 (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		Ethanol ^l	< LOQ		200		
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							



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Pesticides											
Method AOAC 2007.01 & EN 15662 (mod) Units mg/kg Batch 2110676 Analyze 11/29/21 05:28 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	0.200	pass		Carbaryl	< LOQ	0.20	0.100	pass	
Carbofuran	< LOQ	0.20	0.100	pass		Chlorantraniliprole	< LOQ	0.20	0.100	pass	
Chlorfenapyr	< LOQ	1.0	0.500	pass		Chlorpyrifos	< LOQ	0.20	0.100	pass	
Clofentezine	< LOQ	0.20	0.100	pass		Cyfluthrin	< LOQ	1.0	0.500	pass	
Cypermethrin	< LOQ	1.0	0.500	pass		Daminozide	< LOQ	1.0	0.500	pass	
Diazinon	< LOQ	0.20	0.100	pass		Dichlorvos	< LOQ	1.0	0.500	pass	
Dimethoate	< LOQ	0.20	0.100	pass		Ethoprophos	< LOQ	0.20	0.100	pass	
Etofenprox	< LOQ	0.40	0.200	pass		Etoxazole	< LOQ	0.20	0.100	pass	
Fenoxycarb	< LOQ	0.20	0.100	pass		Fenpyroximate	< LOQ	0.40	0.200	pass	
Fipronil	< LOQ	0.40	0.200	pass		Fonicamid	< 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	0.200	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	Status	Notes	
Arsenic	< LOQ		mg/kg	0.00949	2110789	12/02/21	AOAC 2013.06 (mod.)		X	
Cadmium	0.220		mg/kg	0.00949	2110789	12/02/21	AOAC 2013.06 (mod.)		X	
Lead	< LOQ		mg/kg	0.00949	2110789	12/02/21	AOAC 2013.06 (mod.)		X	
Mercury	< LOQ		mg/kg	0.00475	2110789	12/02/21	AOAC 2013.06 (mod.)		X	



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These test results are representative of the individual sample selected and submitted by the client.

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.

† = Analyte not NELAP accredited.

Units of Measure

cfu/g = Colony forming units per gram

g = Gram

g/ml = Gram per milliliter

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



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12423 NE Whitaker Way Portland OR 97230 p.503-254-1794

Cannabis Chain of Custody Record

ORELAP ID: OR100028

Field ID		Date/Time Collected	Pesticides - OR 59 compounds	Pesticide Multi-Residue - 379 compounds	Potency	Residual Solvents	Water Activity	Moisture	Terpenes	Micro: Yeast and Mold	Micro: E.Coli and Total Coliform	Heavy Metals	Mycotoxins	Other	Matrix	Weight	Serving size for edibles	Comments/Metric ID
FORM-DK05-CHOC.FS.DS4		11/24 1350			X													Laz Nat Discount
FORM-DK05-CHOC.FS.DS4		11/24 1350							X	X								
FORM-DK05-CHOC.FS.DS4		11/24 1350	X			X						X						<i>run in parallel</i>

Collected By:	Relinquished By:	Date	Time	Received by:	Date	Time	Lab Use Only:
<input checked="" type="checkbox"/> Standard (5 day)							Client Alias:
<input type="checkbox"/> Rush (3-4 day) (1.5x Standard)							Order Number:
<input type="checkbox"/> Priority Rush (2 day) (2x Standard)							Proper Container
							Sample Condition
							Temperature: <i>20.6°C</i>
							Shipped Via: <i>Client</i>
							Evidence of cooling: <input type="checkbox"/> Yes <input checked="" type="checkbox"/> 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

Test results relate only to the parameters tested and to the samples as received by the laboratory. Test results meet all requirements of NELAP and the Columbia Laboratories quality assurance plan unless otherwise noted. This report shall not be reproduced, except in full, without the written consent of this laboratory. Samples will be retained for a maximum of 30 days from the receipt date unless prior arrangements have been made.
 Testing in accordance with: OAR 333-007-0390 OAR 333-007-0400 OAR 333-007-0410 OAR 333-007-0430



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Revision: 2 Document ID: 3120
Legacy ID: CFL-C21Effective:

Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662		Units: mg/Kg				Batch ID: 2110676				
Matrix Spike/Matrix Spike Duplicate Recoveries		Sample ID: 21-013572-0001								
Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes
Accephate	0.000	0.645	0.668	1.000	3.5%	< 30	64.5%	66.8%	50 - 150	
Acetaminiprid	0.000	0.496	0.498	0.400	0.4%	< 30	124.0%	124.5%	50 - 150	
Aldicarb	0.000	3.136	1.545	0.800	68.0%	< 30	392.0%	193.1%	50 - 150	R, Q1
Abamectin	0.000	0.867	0.913	1.000	5.2%	< 30	86.7%	91.3%	50 - 150	
Azoxystrobin	0.000	0.515	0.480	0.400	7.0%	< 30	128.8%	120.0%	50 - 150	
Bifenazate	0.000	0.419	0.457	0.400	8.5%	< 30	104.9%	114.1%	50 - 150	
Bifenthrin	0.000	0.536	0.574	0.400	6.8%	< 30	134.0%	143.5%	50 - 150	
Boscalid	0.000	0.870	0.974	0.800	11.3%	< 30	108.7%	121.8%	50 - 150	
Carbaryl	0.000	0.482	0.517	0.400	7.1%	< 30	120.4%	129.3%	50 - 150	
Carbofuran	0.000	0.529	0.604	0.400	13.3%	< 30	132.3%	151.1%	50 - 150	Q1
Chlorantraniliprol	0.000	0.383	0.369	0.400	3.9%	< 30	95.8%	92.1%	50 - 150	
Chlorfenapyr	0.000	2.479	2.573	2.000	3.7%	< 30	124.0%	128.7%	50 - 150	
Chlorpyrifos	0.000	0.853	0.904	0.400	5.9%	< 30	213.2%	226.1%	50 - 150	Q1
Clofentezine	0.000	0.301	0.300	0.400	0.3%	< 30	75.1%	74.9%	50 - 150	
Cyfluthrin	0.000	2.073	2.046	2.000	1.3%	< 30	103.7%	102.3%	30 - 150	
Cypermethrin	0.000	2.056	2.081	2.000	1.2%	< 30	102.8%	104.1%	50 - 150	
Daminozide	0.000	0.537	0.536	2.000	0.2%	< 30	26.8%	26.8%	30 - 150	Q
Diazinon	0.000	0.414	0.429	0.400	3.4%	< 30	103.6%	107.2%	50 - 150	
Dichlorvos	0.000	2.579	2.420	2.000	6.4%	< 30	129.0%	121.0%	50 - 150	
Dimethoat	0.000	0.492	0.478	0.400	2.9%	< 30	122.9%	119.5%	50 - 150	
Ethoprophos	0.000	0.420	0.465	0.400	10.3%	< 30	104.8%	116.1%	50 - 150	
Etofenprox	0.000	1.087	1.269	0.800	15.4%	< 30	135.9%	158.6%	50 - 150	Q1
Etoxazol	0.000	0.363	0.437	0.400	18.5%	< 30	90.7%	109.2%	50 - 150	
Fenoxycarb	0.000	0.412	0.409	0.400	0.8%	< 30	103.1%	102.3%	50 - 150	
Fenpyroximat	0.000	0.886	0.897	0.800	1.2%	< 30	110.8%	112.1%	50 - 150	
Fipronil	0.000	1.207	1.067	0.800	12.4%	< 30	150.9%	133.3%	50 - 150	Q1
Fonicamid	0.000	0.891	0.885	1.000	0.7%	< 30	89.1%	88.5%	50 - 150	
Fludioxonil	0.000	0.823	0.917	0.800	10.9%	< 30	102.8%	114.7%	50 - 150	
Hexythiazox	0.000	1.074	1.113	1.000	3.6%	< 30	107.4%	111.3%	50 - 150	
Imazalil	0.000	0.440	0.443	0.400	0.6%	< 30	110.0%	110.8%	50 - 150	
Imidacloprid	0.000	0.769	0.747	0.800	2.9%	< 30	96.1%	93.3%	50 - 150	
Kresoxim-Methyl	0.000	0.850	0.884	0.800	4.0%	< 30	106.2%	110.5%	50 - 150	
Malathion	0.000	0.464	0.447	0.400	3.7%	< 30	115.9%	111.8%	50 - 150	
Metaxyl	0.000	0.483	0.454	0.400	6.2%	< 30	120.8%	113.6%	50 - 150	
Methiocarb	0.000	0.495	0.409	0.400	19.2%	< 30	123.8%	102.1%	50 - 150	
Methomyl	0.000	0.859	0.793	0.800	8.0%	< 30	107.4%	99.1%	50 - 150	
MGK 264	0.000	0.377	0.406	0.400	7.5%	< 30	94.2%	101.6%	50 - 150	
Myclobutanil	0.000	0.481	0.440	0.400	8.9%	< 30	120.2%	109.9%	50 - 150	
Naled	0.000	1.227	1.115	1.000	9.6%	< 30	122.7%	111.5%	50 - 150	
Oxamyl	0.000	2.087	2.056	2.000	1.5%	< 30	104.4%	102.8%	50 - 150	
Paclobutrazol	0.000	0.857	0.872	0.800	1.8%	< 30	107.1%	109.0%	50 - 150	
Parathion Methyl	0.000	1.065	1.042	0.800	2.2%	< 30	133.2%	130.3%	30 - 150	
Permethrin	0.000	0.406	0.396	0.400	2.5%	< 30	101.4%	99.0%	50 - 150	
Phosmet	0.000	0.478	0.478	0.400	0.0%	< 30	119.5%	119.5%	50 - 150	
Piperonyl butoxide	0.000	2.672	2.743	2.000	2.6%	< 30	133.6%	137.2%	50 - 150	
Prallethrin	0.000	0.306	0.327	0.400	6.4%	< 30	76.6%	81.6%	50 - 150	
Propiconazole	0.000	0.818	0.901	0.800	9.6%	< 30	102.2%	112.6%	50 - 150	
Propoxur	0.000	0.491	0.541	0.400	9.6%	< 30	122.8%	135.3%	50 - 150	
Pyrethrins	0.000	0.028	0.031	0.413	8.8%	< 30	6.8%	7.4%	50 - 150	Q
Pyridaben	0.000	0.379	0.405	0.400	6.6%	< 30	94.8%	101.3%	50 - 150	
Spinosad	0.000	0.514	0.479	0.388	6.9%	< 30	132.4%	123.5%	50 - 150	
Spiromesifen	0.000	0.374	0.420	0.400	11.6%	< 30	93.6%	105.1%	50 - 150	
Spirotetramat	0.000	0.366	0.360	0.400	1.5%	< 30	91.5%	90.1%	50 - 150	
Spiroxamine	0.000	0.819	0.877	0.800	6.8%	< 30	102.4%	109.6%	50 - 150	
Tebuconazol	0.000	0.902	0.963	0.800	6.5%	< 30	112.8%	120.3%	50 - 150	
Thiadoprid	0.000	0.482	0.494	0.400	2.5%	< 30	120.5%	123.5%	50 - 150	
Thiamethoxam	0.000	0.328	0.311	0.400	5.2%	< 30	81.9%	77.7%	50 - 150	
Trifloxystrobin	0.000	0.458	0.494	0.400	7.5%	< 30	114.6%	123.5%	50 - 150	



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

Report Number: 21-013871/D002.R000
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 Received: 11/24/21 14:41



Revision: 2 Document ID: 3120
 Legacy ID: CFL-C21Effective:

Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662		Units: mg/Kg		Batch ID: 2110676				
Method Blank		Laboratory Control Sample						
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits	Notes
Accephate	0.010	< 0.250		0.913	1.000	91.3	71.9 - 134	
Acetaminocyl	0.000	< 1.000		3.308	4.000	82.7	73.8 - 137	
Acetamidiprid	0.000	< 0.100		0.406	0.400	101.5	75.0 - 139	
Aldicarb	0.000	< 0.200		1.579	0.800	197.4	87.3 - 162	Q1
Abamectin	0.000	< 0.250		0.968	1.000	96.8	77.1 - 143	
Azoxystrobin	0.000	< 0.100		0.380	0.400	94.9	72.6 - 135	
Bifenazate	0.000	< 0.100		0.372	0.400	93.0	75.1 - 140	
Bifenthrin	0.000	< 0.100		0.365	0.400	91.4	72.0 - 134	
Boscalid	0.000	< 0.200		0.739	0.800	92.4	75.3 - 140	
Carbaryl	0.000	< 0.100		0.421	0.400	105.4	74.7 - 139	
Carbofuran	0.000	< 0.100		0.408	0.400	102.0	74.2 - 138	
Chlorantraniliprol	0.000	< 0.100		0.379	0.400	94.8	70.3 - 130	
Chlorfenapyr	0.000	< 0.500		1.894	2.000	94.7	74.0 - 137	
Chlorpyrifos	0.000	< 0.100		0.457	0.400	114.2	71.5 - 133	
Clofentezine	0.000	< 0.100		0.167	0.400	41.8	42.9 - 79.6	Q6
Cyfluthrin	0.000	< 0.500		1.959	2.000	98.0	79.2 - 147	
Cypermethrin	0.000	< 0.500		1.917	2.000	95.8	77.4 - 144	
Daminozide	0.016	< 0.500		0.716	2.000	35.8	27.3 - 50.7	
Diazinon	0.000	< 0.100		0.388	0.400	96.9	76.0 - 141	
Dichlorvos	0.000	< 0.500		2.354	2.000	117.7	81.5 - 151	
Dimethoat	0.000	< 0.100		0.418	0.400	104.4	73.8 - 137	
Ethoprophos	0.000	< 0.100		0.380	0.400	94.9	74.7 - 139	
Etofenprox	0.000	< 0.200		0.745	0.800	93.1	73.3 - 136	
Etoxazol	0.000	< 0.100		0.318	0.400	79.5	88.0 - 163	Q6
Fenoxycarb	0.000	< 0.100		0.371	0.400	92.8	73.3 - 136	
Fenpyroximat	0.000	< 0.200		0.831	0.800	103.8	72.8 - 135	
Fipronil	0.000	< 0.200		0.813	0.800	101.6	73.9 - 137	
Fonicamid	0.000	< 0.250		1.116	1.000	111.6	75.9 - 141	
Fludioxonil	0.000	< 0.200		0.841	0.800	105.2	76.4 - 142	
Hexythiazox	0.000	< 0.250		0.989	1.000	98.9	70.8 - 132	
Imazali	0.009	< 0.100		0.395	0.400	98.7	72.2 - 134	
Imidacloprid	0.000	< 0.200		0.812	0.800	101.5	74.6 - 139	
Kresoxim-Methyl	0.000	< 0.200		0.764	0.800	95.5	74.3 - 138	
Malathion	0.000	< 0.100		0.367	0.400	91.9	74.1 - 138	
Metaxyl	0.009	< 0.100		0.411	0.400	102.8	74.0 - 137	
Methiocarb	0.000	< 0.100		0.361	0.400	90.3	74.9 - 139	
Methomyl	0.000	< 0.200		0.859	0.800	107.3	72.3 - 134	
MGK 264	0.000	< 0.100		0.412	0.400	102.9	76.0 - 141	
Myclobutanil	0.000	< 0.100		0.383	0.400	95.7	73.7 - 137	
Naled	0.000	< 0.250		0.494	1.000	49.4	38.6 - 71.7	
Oxamyl	0.000	< 0.500		2.151	2.000	107.5	73.3 - 136	
Paclobutrazol	0.000	< 0.200		0.756	0.800	94.6	74.4 - 138	
Parathion Methyl	0.000	< 0.200		0.828	0.800	103.6	72.5 - 135	
Permethrin	0.000	< 0.100		0.383	0.400	95.7	74.9 - 139	
Phosmet	0.000	< 0.100		0.361	0.400	90.2	74.0 - 137	
Piperonyl butoxide	0.000	< 0.500		2.143	2.000	107.2	75.1 - 140	
Prallethrin	0.001	< 0.100		0.388	0.400	97.1	72.5 - 135	
Propiconazole	0.085	< 0.200		0.712	0.800	89.0	75.5 - 140	
Propoxur	0.012	< 0.100		0.395	0.400	98.7	73.7 - 137	
Pyrethrins	0.009	< 0.100		0.032	0.413	7.8	72.2 - 134	Q6
Pyridaben	0.000	< 0.100		0.322	0.400	80.4	76.4 - 142	
Spinosad	0.000	< 0.100		0.422	0.388	108.7	75.8 - 141	
Spiromesifen	0.000	< 0.100		0.358	0.400	89.4	75.4 - 140	
Spirotetramat	0.000	< 0.100		0.375	0.400	93.8	72.8 - 135	
Spiroxamine	0.000	< 0.200		0.765	0.800	95.6	72.7 - 135	
Tebuconazol	0.000	< 0.200		0.742	0.800	92.8	74.1 - 138	
Thiadoprid	0.000	< 0.100		0.381	0.400	95.3	74.4 - 138	
Thiamethoxam	0.000	< 0.100		0.414	0.400	103.5	74.0 - 137	
Trifloxystrobin	0.000	< 0.100		0.434	0.400	108.4	73.9 - 137	



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Laboratory Quality Control Results

Batch ID: 2 0727

Residual Solvents		Method Blank		Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	200		0	9.8	µg/g	0	0	0
Isobutane	ND	200		350	260	µg/g	0	0	0
Butane	ND	200		350	260	µg/g	0	0	0
2,2-Dimethylpropane	ND	200		9.0	600	µg/g	23	0	0
Methanol	ND	200		660	6.0	µg/g	03	0	0
Ethylene Oxide	ND	30		3	95	µg/g	8	0	0
2-Methylbutane	ND	200		650	6.0	µg/g	02.5	0	0
Pentane	ND	200		600	6.0	µg/g	99	0	0
Ethanol	ND	200		560	6.0	µg/g	96.9	0	0
Ethyl Ether	ND	200		0	6.0	µg/g	9.3	0	0
2,2-Dimethylbutane	ND	30		53	6	µg/g	93.2	0	0
Acetone	ND	200		530	6.0	µg/g	95.0	0	0
2-Propanol	ND	200		580	6.0	µg/g	98	0	0
Ethyl Formate	ND	500		5.0	6.0	µg/g	93.8	0	0
Acetonitrile	ND	00		66	8	µg/g	96.3	0	0
Methyl Acetate	ND	500		5.0	6.0	µg/g	9.5	0	0
2,3-Dimethylbutane	ND	30		59	6	µg/g	95.2	0	0
Dichloromethane	ND	60		39	9	µg/g	89	0	0
2-Methylpentane	ND	30		53	65	µg/g	92	0	0
M. BE	ND	500		5.0	600	µg/g	98	0	0
3-Methylpentane	ND	30		5	2	µg/g	9.3	0	0
Hexane	ND	30		5	6	µg/g	90	0	0
Propanol	ND	500		5.0	6.0	µg/g	95	0	0
Methylethylketone	ND	500		530	620	µg/g	9	0	0
Ethyl acetate	ND	200		50	6.0	µg/g	90	0	0
2-Butanol	ND	200		30	6.0	µg/g	88.8	0	0
tetrahydrofuran	ND	00		25	83	µg/g	88.2	0	0
Cyclohexane	ND	200		00	6.0	µg/g	8.0	0	0
2-methyl propanol	ND	500		5.0	620	µg/g	95	0	0
Benzene	ND	200		29	5.36	µg/g	80.0	0	0
Isopropyl Acetate	ND	200		0	620	µg/g	90	0	0
Heptane	ND	200		30	6.0	µg/g	88.8	0	0
Butanol	ND	500		5.0	6.0	µg/g	93.8	0	0
Propyl Acetate	ND	500		3.0	620	µg/g	8.6	0	0
Dioxane	ND	00		00	89	µg/g	8.8	0	0
2-Ethoxyethanol	ND	30		8	6	µg/g	88.6	0	0
Methyl isobutylketone	ND	500		60	6.0	µg/g	90	0	0
3-Methyl butanol	ND	500		0	6.0	µg/g	89	0	0
Ethylene Glycol	ND	200		0	50	µg/g	82	0	0
o-xylene	ND	200		38	8	µg/g	93	0	0
Isobutyl Acetate	ND	500		500	6.0	µg/g	93.2	0	0
Pentanol	ND	500		350	6.0	µg/g	83.9	0	0
Butyl Acetate	ND	500		500	620	µg/g	92.6	0	0
Ethyl benzene	ND	200		33	68	µg/g	5	0	0
m,p Xylene	ND	200		5	9	µg/g	2	0	0
o-Xylene	ND	200		0	982	µg/g	5.8	0	0
Cumene	ND	30		2	69	µg/g	3	0	0
Anisole	ND	500		350	630	µg/g	82.8	0	0
DMSO	ND	500		320	630	µg/g	8.0	0	0
2,2-dimethoxyethane	ND	50		66	62	µg/g	03	0	0
diethylamine	ND	500		590	6.0	µg/g	95.2	0	0
N,N-dimethylformamide	ND	50		53	502	µg/g	90.2	0	0
N,N-dimethylacetamide	ND	50		05	85	µg/g	83.5	0	0
Pyridine	ND	50		50	66	µg/g	90	0	0
2-Dichloroethane	ND	0.926		0.926		µg/g	92.6	0	0
Chloroform	ND	0.88		0.88		µg/g	88	0	0
richloroethylene	ND	0.9		0.9		µg/g	9	0	0
Ethylene Oxide	ND	0.93		0.93		µg/g	93	0	0
Dichloromethane	ND	0.863		0.863		µg/g	86.3	0	0
Benzene	ND	0.83		0.83		µg/g	83.0	0	0



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QC Sample Duplicate Sample ID: 21 013830 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	
Ethyl 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	00	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	
M.E.	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	
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	00	µg/g	0.0	20	Acceptable	
Cyclohexane	ND	ND	200	µg/g	0.0	20	Acceptable	
2 methyl propanol	ND	ND	500	µg/g	0.0	20	Acceptable	
Benzene	ND	ND	100	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	
Butanol	ND	ND	500	µg/g	0.0	20	Acceptable	
Propyl Acetate	ND	ND	500	µg/g	0.0	20	Acceptable	
Dioxane	ND	ND	00	µg/g	0.0	20	Acceptable	
2 Ethoxyethanol	ND	ND	30	µg/g	0.0	20	Acceptable	
Methyl isobutyl ketone	ND	ND	500	µg/g	0.0	20	Acceptable	
3 Methyl butanol	ND	ND	500	µg/g	0.0	20	Acceptable	
Ethyl Glycol	ND	ND	200	µg/g	0.0	20	Acceptable	
o-xylene	ND	ND	200	µg/g	0.0	20	Acceptable	
Isobutyl Acetate	ND	ND	500	µg/g	0.0	20	Acceptable	
Pentanol	ND	ND	500	µg/g	0.0	20	Acceptable	
Butyl Acetate	ND	ND	500	µg/g	0.0	20	Acceptable	
Ethyl benzene	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	
2,4-dimethoxyethane	ND	ND	50	µg/g	0.0	20	Acceptable	
riethylamine	ND	ND	500	µg/g	0.0	20	Acceptable	
N,N-dimethylformamide	ND	ND	50	µg/g	0.0	20	Acceptable	
N,N-dimethylacetamide	ND	ND	50	µg/g	0.0	20	Acceptable	
Pyridine	ND	ND	50	µg/g	0.0	20	Acceptable	
2,2-Dichloroethane	ND	ND	100	µg/g	0.0	20	Acceptable	
Chloroform	ND	ND	100	µg/g	0.0	20	Acceptable	
trichloroethylene	ND	ND	100	µg/g	0.0	20	Acceptable	
Ethyl Oxide	ND	ND	30	µg/g	0.0	20	Acceptable	
Dichloromethane	ND	ND	100	µg/g	0.0	20	Acceptable	
Benzene	ND	ND	100	µg/g	0.0	20	Acceptable	

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

Units of Measure:
 µg/g Microgram per gram or ppm



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Laboratory Quality Control Results

J AOAC 2015 V98-6							
Batch ID: 2110747							
Laboratory Control Sample							
Analyte	Result	Spike	Units	% Rec	Limits	Evaluation	Notes
CBDVA	0.00978	0.01	%	97.8	85.0 - 115	Acceptable	
CBDV	0.0104	0.01	%	104	85.0 - 115	Acceptable	
CBE	0.0106	0.01	%	106	85.0 - 115	Acceptable	
CBDA	0.0107	0.01	%	107	85.0 - 115	Acceptable	
CBGA	0.00989	0.01	%	98.9	85.0 - 115	Acceptable	
CBG	0.0103	0.01	%	103	85.0 - 115	Acceptable	
CBD	0.0110	0.01	%	110	85.0 - 115	Acceptable	
THCV	0.0103	0.01	%	103	85.0 - 115	Acceptable	
d8THCV	0.00992	0.01	%	99.2	85.0 - 115	Acceptable	
THCVA	0.00935	0.01	%	93.5	85.0 - 115	Acceptable	
CBN	0.0109	0.01	%	109	85.0 - 115	Acceptable	
exo-THC	0.00964	0.01	%	96.4	85.0 - 115	Acceptable	
d9THC	0.0106	0.01	%	106	85.0 - 115	Acceptable	
d8THC	0.00988	0.01	%	98.8	85.0 - 115	Acceptable	
CBL	0.00928	0.01	%	92.8	85.0 - 115	Acceptable	
CBC	0.0103	0.01	%	103	85.0 - 115	Acceptable	
THCA	0.0106	0.01	%	106	85.0 - 115	Acceptable	
CBCA	0.00937	0.01	%	93.7	85.0 - 115	Acceptable	
CBLA	0.00973	0.01	%	97.3	85.0 - 115	Acceptable	
CBT	0.00921	0.01	%	92.1	85.0 - 115	Acceptable	

Method Blank

Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.003	%	< 0.003	Acceptable	
CBDV	<LOQ	0.003	%	< 0.003	Acceptable	
CBE	<LOQ	0.003	%	< 0.003	Acceptable	
CBDA	<LOQ	0.003	%	< 0.003	Acceptable	
CBGA	<LOQ	0.003	%	< 0.003	Acceptable	
CBG	<LOQ	0.003	%	< 0.003	Acceptable	
CBD	<LOQ	0.003	%	< 0.003	Acceptable	
THCV	<LOQ	0.003	%	< 0.003	Acceptable	
d8THCV	<LOQ	0.003	%	< 0.003	Acceptable	
THCVA	<LOQ	0.003	%	< 0.003	Acceptable	
CBN	<LOQ	0.003	%	< 0.003	Acceptable	
exo-THC	<LOQ	0.003	%	< 0.003	Acceptable	
d9THC	<LOQ	0.003	%	< 0.003	Acceptable	
d8THC	<LOQ	0.003	%	< 0.003	Acceptable	
CBL	<LOQ	0.003	%	< 0.003	Acceptable	
CBC	<LOQ	0.003	%	< 0.003	Acceptable	
THCA	<LOQ	0.003	%	< 0.003	Acceptable	
CBCA	<LOQ	0.003	%	< 0.003	Acceptable	
CBLA	<LOQ	0.003	%	< 0.003	Acceptable	
CBT	<LOQ	0.003	%	< 0.003	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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Laboratory Quality Control Results

J AOAC 2015 V98-6								
Batch ID: 2110747								
Sample Duplicate								
Sample ID: 21-013826-0001								
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBDV	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBE	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBDA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBG	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBD	0.595	0.592	0.003	%	0.580	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBN	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
d9THC	0.135	0.135	0.003	%	0.566	< 20	Acceptable	
d8THC	0.0300	0.0307	0.003	%	2.00	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBC	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBCA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBT	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	

Abbreviations

- ND - None Detected at or above MRL
- RPD - Relative Percent Difference
- LOQ - Limit of Quantitation
- NA - Calculation Not Applicable given non-numerical results

Units of Measure:

% - Percent



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Report Number: 21-013871/D002.R000
Report Date: 12/03/2021
ORELAP#: OR100028
Purchase Order:
Received: 11/24/21 14:41

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.