Columbia LABORATORIES
A Tentamus Company

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

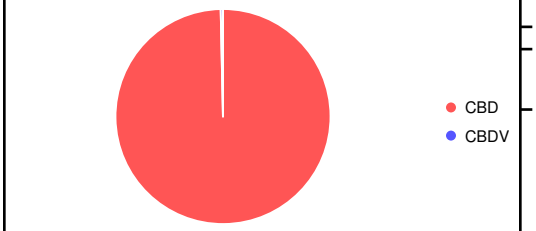
Report Number: 23-001623/D002.R000
Report Date: 02/10/2023
ORELAP#: OR100028
Purchase Order:
Received: 02/07/23 15:18

Customer: IHC LLC
Product identity: 01ISO204
Client/Metrc ID: .
Laboratory ID: 23-001623-0001

Summary

Potency:

Analyte	Result (%)
CBD	91.2
CBDV	0.309



CBD-Total 91.2%

THC-Total <LOQ

(Reported in percent of total sample)

Residual Solvents:

Analyte	Result (µg/g)	Limits (µg/g)	Status
n-Heptane	247	5000	pass

Pesticides:

All analytes passing and less than LOQ.

Metals:

Less than LOQ for all analytes.

Testing in accordance with: OAR 333-007-0430

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Customer: IHC LLC
825 NW 16th Ave
Portland Oregon 97209
United States of America (USA)

Product identity: 01ISO204
Client/Metrc ID: .
Sample Date:
Laboratory ID: 23-001623-0001
Evidence of Cooling: No
Temp: 17.0
Relinquished by: ramos

THE HEMP COLLECT

Sample Results

Potency **Method:** J AOAC 2015 V98-6 (mod)[b] **Units** % **Batch:** 2301306 **Analyze:** 2/9/23 10:44:00 PM

Analyte	As Received	Dry weight	LOQ	Notes
CBC	< LOQ		0.0676	
CBC-A	< LOQ		0.0676	
CBC-Total	< LOQ		0.127	
CBD	91.2		0.676	
CBD-A	< LOQ		0.0676	
CBD-Total	91.2		0.735	
CBDV	0.309		0.0676	
CBDV-A	< LOQ		0.0676	
CBDV-Total	0.309		0.126	
CBE	< LOQ		0.0676	
CBG	< LOQ		0.0676	
CBG-A	< LOQ		0.0676	
CBG-Total	< LOQ		0.126	
CBL	< LOQ		0.0676	
CBL-A	< LOQ		0.0676	
CBL-Total	< LOQ		0.127	
CBN	< LOQ		0.0676	
CBT	< LOQ		0.0676	
Δ10-THC-9R	< LOQ		0.0676	
Δ10-THC-9S	< LOQ		0.0676	
Δ10-THC-Total	< LOQ		0.135	
Δ8-THC	< LOQ		0.0676	
Δ8-THCV	< LOQ		0.0676	
Δ9-THC	< LOQ		0.0676	
exo-THC	< LOQ		0.0676	
THC-A	< LOQ		0.0676	
THC-Total	< LOQ		0.127	
THCV	< LOQ		0.0676	
THCV-A	< LOQ		0.0676	
THCV-Total	< LOQ		0.126	
Total Cannabinoids	91.5			

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Solvents **Method:** Residual Solvents by GC/MS^{b} **Units** µg/g **Batch** 2301280 **Analyze** 02/09/23 03:09 PM

Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
1,4-Dioxane	< LOQ	380	100	pass		2-Butanol	< LOQ	5000	200	pass	
2-Ethoxyethanol	< LOQ	160	30.0	pass		2-Methylbutane (Isopentane)	< LOQ		200		
2-Methylpentane	< LOQ		30.0			2-Propanol (IPA)	< LOQ	5000	200	pass	
2,2-Dimethylbutane	< LOQ		30.0			2,2-Dimethylpropane (neo-pentane)	< LOQ		200		
2,3-Dimethylbutane	< LOQ		30.0			3-Methylpentane	< LOQ		30.0		
Acetone	< LOQ	5000	200	pass		Acetonitrile	< LOQ	410	100	pass	
Benzene	< LOQ	2.00	1.00	pass		Butanes (sum)	< LOQ	5000	400	pass	
Cyclohexane	< LOQ	3880	200	pass		Ethyl acetate	< LOQ	5000	200	pass	
Ethyl benzene	< LOQ		200			Ethyl ether	< LOQ	5000	200	pass	
Ethylene glycol	< LOQ	620	200	pass		Ethylene oxide	< LOQ	50.0	20.0	pass	
Hexanes (sum)	< LOQ	290	150	pass		Isopropyl acetate	< LOQ	5000	200	pass	
Isopropylbenzene (Cumene)	< LOQ	70.0	30.0	pass		m,p-Xylene	< LOQ		200		
Methanol	< LOQ	3000	200	pass		Methylene chloride	< LOQ	600	60.0	pass	
Methylpropane (Isobutane)	< LOQ		200			n-Butane	< LOQ		200		
n-Heptane	247	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)þ **Units** mg/kg **Batch** 2301287 **Analyze** 02/10/23 08:01 AM

Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
Abamectin¥	< LOQ	0.50	0.250	pass		Acephate¥	< LOQ	0.40	0.200	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		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	0.100	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	Analyzed	Method	Status	Notes
Arsenic	< LOQ	0.200	mg/kg	0.0974	2301312	02/10/23	AOAC 2013.06 (mod.)þ	pass	
Cadmium	< LOQ	0.200	mg/kg	0.0974	2301312	02/10/23	AOAC 2013.06 (mod.)þ	pass	
Lead	< LOQ	0.500	mg/kg	0.0974	2301312	02/10/23	AOAC 2013.06 (mod.)þ	pass	
Mercury	< LOQ	0.100	mg/kg	0.0487	2301312	02/10/23	AOAC 2013.06 (mod.)þ	pass	

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Abbreviations

Limits: Action Levels per OAR-333-007-0400, OAR-333-007-0210, OAR-333-007-0220, CCR title 16-division 42. BCC-section 5723

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

þ = ISO/IEC 17025:2017 accredited method.
¥ = TNI accredited analyte.

Units of Measure

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

Approved Signatory

Derrick Tanner
General Manager

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Columbia LABORATORIES
A Tentamus Company

Hemp / Cannabis Usable / Extract / Finished Pro

Chain of Custody Record

Revision: 4.00 Control#: CF023 Rev 02/24/2021 Eff: 03/04/20

ORELAP ID: OR100028

#CLCC: 23-001623

#CLCC:

Company: The Hemp Collect
Contact: kyle@thehempcollect.com
Street: 431 NW Flanders st.
City: Portland State: OR Zip: 97209
☑ Email Results: dropbox (IHC)
Ph: (61) 6U8164 ☐ Fx Results: (___)
Billing (if different): joel@thehempcollect.com

Project Number:
Project Name:
Custom Reporting:
Report to State – ☐ METRC or ☐ Other:
Turnaround time: ☐ 5 Business Day Standard Turnaround
☑ 3 Business Day Rush Turnaround*
☐ 2 Business Day Rush Turnaround*
*Check for availability
Sampled by:

Analysis Requested

Lab ID	Client Sample Identification	Date	Time	Pesticides – OR 59 compounds	Pesticide Multi-Residue – 378 compounds	Potency	Residual Solvents	Moisture & Water Activity	Terpenes	Micro: Yeast and Mold	Micro: E.Coli and Total Coliform	Heavy Metals	Mycotoxins	Other:	Sample Type †	Weight (Units)	Comments/Metrc ID
1	01ISO204			X		X	X					X			S		Sample #1: Alternate client name: Top Notch Medicinals
2																	
3																	
4																	
5																	
6																	
7																	
8																	
9																	
10																	

Relinquished By:	Date	Time	Received By:	Date	Time
Kyle Farook	2/7	11:30 A		2.7.23	1350
	2.7.23	345		2/8/23	518

Lab Use Only:
☐ Shipped Via: _____ or ☐ Client drop
Evidence of cooling: ☐ Yes | ☐ No - Temp (°C): 17.0
Sample in good condition: ☐ Yes | ☐ No
☐ Cash | ☐ Check | ☐ CC | ☐ Net: _____
Prelog storage: _____

† - Sample Type Codes: Vegetation (V) ; Isolates (S) ; Extract/Concentrate (C) ; Tincture/Topical (T) ; Edible (E) ; Beverage (B)

Samples submitted to Columbia Laboratories will assume acceptance constitutes an agreement for services in accordance with the current terms of service associated with the COC. By signing "Relinquished by" you are agreeing to these terms.

12423 NE Whitaker Way
Portland, OR 97230

P: (503) 254-1794 | Fax: (503) 254-1452
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Page ____ of ____
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Purchase Order:
Received: 02/07/23 15:18

Revision: 2 Document ID: 7087
Legacy ID: CFL-E33Effective:

Laboratory Quality Control Results

Residual Solvents **Batch ID:** 2301280

Method Blank					Laboratory Control Sample							
Analyte	**Result**		**LOQ**	**Notes**	**Result**	**Spike**	**Units**	**% Rec**	**Limits**			**Notes**
Propane	ND	<	200		425	572	µg/g	74.3	60	-	120	
Isobutane	ND	<	200		561	731	µg/g	76.7	60	-	120	
Butane	ND	<	200		531	731	µg/g	72.6	60	-	120	
2,2-Dimethylpropane	ND	<	200		649	936	µg/g	69.3	60	-	120	
Methanol	ND	<	200		1260	1620	µg/g	77.8	60	-	120	
Ethylene Oxide	ND	<	30		40.8	56.2	µg/g	72.6	60	-	120	
2-Methylbutane	ND	<	200		1210	1610	µg/g	75.2	60	-	120	
Pentane	ND	<	200		1220	1600	µg/g	76.3	60	-	120	
Ethanol	ND	<	200		1300	1610	µg/g	80.7	70	-	130	
Ethyl Ether	ND	<	200		1300	1630	µg/g	79.8	60	-	120	
2,2-Dimethylbutane	ND	<	30		134	171	µg/g	78.4	60	-	120	
Acetone	ND	<	200		1300	1630	µg/g	79.8	60	-	120	
2-Propanol	ND	<	200		1330	1620	µg/g	82.1	60	-	120	
Ethyl Formate	ND	<	500		1300	1670	µg/g	77.8	70	-	130	
Acetonitrile	ND	<	100		378	498	µg/g	75.9	60	-	120	
Methyl Acetate	ND	<	500		1260	1730	µg/g	72.8	70	-	130	
2,3-Dimethylbutane	ND	<	30		136	171	µg/g	79.5	60	-	120	
Dichloromethane	ND	<	60		403	483	µg/g	83.4	60	-	120	
2-Methylpentane	ND	<	30		128	168	µg/g	76.2	60	-	120	
MTBE	ND	<	500		1290	1650	µg/g	78.2	70	-	130	
3-Methylpentane	ND	<	30		121	167	µg/g	72.5	60	-	120	
Hexane	ND	<	30		174	182	µg/g	95.6	60	-	120	
1-Propanol	ND	<	500		1120	1620	µg/g	69.1	70	-	130	Q6
Methylethylketone	ND	<	500		1180	1620	µg/g	72.8	70	-	130	
Ethyl acetate	ND	<	200		1280	1610	µg/g	79.5	60	-	120	
2-Butanol	ND	<	200		1290	1600	µg/g	80.6	60	-	120	
Tetrahydrofuran	ND	<	100		392	483	µg/g	81.2	60	-	120	
Cyclohexane	ND	<	200		1330	1610	µg/g	82.6	60	-	120	
2-methyl-1-propanol	ND	<	500		1170	1620	µg/g	72.2	70	-	130	
Benzene	ND	<	1		4.48	5.02	µg/g	89.2	60	-	120	
Isopropyl Acetate	ND	<	200		1300	1620	µg/g	80.2	60	-	120	
Heptane	ND	<	200		1240	1610	µg/g	77.0	60	-	120	
1-Butanol	ND	<	500		1160	1630	µg/g	71.2	70	-	130	
Propyl Acetate	ND	<	500		1140	1610	µg/g	70.8	70	-	130	
1,4-Dioxane	ND	<	100		409	491	µg/g	83.3	60	-	120	
2-Ethoxyethanol	ND	<	30		138	181	µg/g	76.2	60	-	120	
Methylisobutylketone	ND	<	500		1190	1620	µg/g	73.5	70	-	130	
3-Methyl-1-butanol	ND	<	500		1150	1630	µg/g	70.6	70	-	130	
Ethylene Glycol	ND	<	200		357	484	µg/g	73.8	60	-	120	
Toluene	ND	<	100		399	485	µg/g	82.3	60	-	120	
Isobutyl Acetate	ND	<	500		1130	1630	µg/g	69.3	70	-	130	Q6
1-Pentanol	ND	<	500		1080	1620	µg/g	66.7	70	-	130	Q6
Butyl Acetate	ND	<	500		1110	1620	µg/g	68.5	70	-	130	Q6
Ethylbenzene	ND	<	200		790	969	µg/g	81.5	60	-	120	
m,p-Xylene	ND	<	200		804	994	µg/g	80.9	60	-	120	
o-Xylene	ND	<	200		792	967	µg/g	81.9	60	-	120	
Cumene	ND	<	30		138	171	µg/g	80.7	60	-	120	
Anisole	ND	<	500		1180	1630	µg/g	72.4	70	-	130	
DMSO	ND	<	500		1130	1680	µg/g	67.3	70	-	130	Q6
1,2-dimethoxyethane	ND	<	50		122	169	µg/g	72.2	70	-	130	
Triethylamine	ND	<	500		1240	1630	µg/g	76.1	70	-	130	
N,N-dimethylformamide	ND	<	150		347	482	µg/g	72.0	70	-	130	
N,N-dimethylacetamide	ND	<	150		328	510	µg/g	64.3	70	-	130	Q6
Pyridine	ND	<	50		144	203	µg/g	70.9	70	-	130	
Sulfolane	ND	<	50		114	172	µg/g	66.3	70	-	130	Q6
1,2-Dichloroethane	ND	<	1		0.701	1	µg/g	70.1	70	-	130	
Chloroform	ND	<	1		0.708	1	µg/g	70.8	70	-	130	
Trichloroethylene	ND	<	1		0.695	1	µg/g	69.5	70	-	130	Q6
1,1-Dichloroethane	ND	<	1		0.672	1	µg/g	67.2	70	-	130	Q6

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Revision: 2 Document ID: 7087
Legacy ID: CFL-E33Effective:

QC - Sample Duplicate **Sample ID: 23-001411-0001**

Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Accept/Fail	Notes
Propane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Isobutane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Butane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2,2-Dimethylpropane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Methanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethylene Oxide	ND	ND	30	µg/g	0.0	< 20	Acceptable	
2-Methylbutane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Pentane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethyl Ether	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2,2-Dimethylbutane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Acetone	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2-Propanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethyl Formate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Acetonitrile	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Methyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
2,3-Dimethylbutane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Dichloromethane	ND	ND	60	µg/g	0.0	< 20	Acceptable	
2-Methylpentane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
MTBE	ND	ND	500	µg/g	0.0	< 20	Acceptable	
3-Methylpentane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Hexane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
1-Propanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
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	100	µg/g	0.0	< 20	Acceptable	
Isobutyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
1-Pentanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Butyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Ethylbenzene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
m,p-Xylene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
o-Xylene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Cumene	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Anisole	ND	ND	500	µg/g	0.0	< 20	Acceptable	
DMSO	ND	ND	500	µg/g	0.0	< 20	Acceptable	
1,2-dimethoxyethane	ND	ND	50	µg/g	0.0	< 20	Acceptable	
Triethylamine	ND	ND	500	µg/g	0.0	< 20	Acceptable	
N,N-dimethylformamide	ND	ND	150	µg/g	0.0	< 20	Acceptable	
N,N-dimethylacetamide	ND	ND	150	µg/g	0.0	< 20	Acceptable	
Pyridine	ND	ND	50	µg/g	0.0	< 20	Acceptable	
Sulfolane	ND	ND	50	µg/g	0.0	< 20	Acceptable	
1,2-Dichloroethane	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Chloroform	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Trichloroethylene	ND	ND	1	µg/g	0.0	< 20	Acceptable	
1,1-Dichloroethane	ND	ND	1	µg/g	0.0	< 20	Acceptable	

Abbreviations

ND - None Detected at or above MRL
RPD - Relative Percent Difference
LOQ - Limit of Quantitation
Q6 - Quality control outside QC limits. Data acceptable based on remaining QC.

Units of Measure:

µg/g- Microgram per gram or ppm

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12423 NE Whitaker Way
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503-254-1794

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Laboratory Pesticide Quality Control Results

AOAC 2007.1 &EN 15662			Units: mg/Kg					Batch ID: 2301287	
Method Blank				Laboratory Control Sample					
Analyte	Blank Result	Blank Limits	Notes	LCS Result	LCS Spike	LCS % Rec	Limits		Notes
Abamectin	0.000	< 0.250		0.929	1.000	92.9	50.0	150	
Acephate	0.000	< 0.200		0.682	0.800	85.2	60.0	120	
Acequinocyl	0.000	< 1.000		3.758	4.000	93.9	40.0	160	
Acetamiprid	0.000	< 0.100		0.390	0.400	97.5	60.0	120	
Aldicarb	0.000	< 0.200		0.765	0.800	95.7	60.0	120	
Azoxystrobin	0.000	< 0.100		0.370	0.400	92.6	60.0	120	
Bifenazate	0.000	< 0.100		0.355	0.400	88.9	60.0	120	
Bifenthrin	0.000	< 0.100		0.379	0.400	94.9	50.0	150	
Boscalid	0.000	< 0.200		0.765	0.800	95.6	60.0	120	
Carbaryl	0.000	< 0.100		0.389	0.400	97.2	60.0	120	
Carbofuran	0.000	< 0.100		0.385	0.400	96.2	60.0	120	
Chlorantraniliprole	0.000	< 0.100		0.400	0.400	99.9	60.0	120	
Chlorfenapyr	0.000	< 0.500		2.044	2.000	102.2	60.0	120	
Chlorpyrifos	0.000	< 0.100		0.380	0.400	95.0	60.0	120	
Clofentezine	0.000	< 0.100		0.371	0.400	92.7	60.0	120	
Cyfluthrin	0.000	< 0.500		1.965	2.000	98.3	50.0	150	
Cypermethrin	0.000	< 0.500		1.951	2.000	97.6	50.0	150	
Daminozide	0.000	< 0.500		1.808	2.000	90.4	60.0	120	
Diazinon	0.000	< 0.100		0.409	0.400	102.1	60.0	120	
Dichlorvos	0.000	< 0.500		1.948	2.000	97.4	60.0	120	
Dimethoate	0.000	< 0.100		0.392	0.400	98.1	60.0	120	
Ethoprophos	0.000	< 0.100		0.383	0.400	95.7	60.0	120	
Etofenprox	0.000	< 0.200		0.775	0.800	96.9	50.0	150	
Etoxazole	0.000	< 0.100		0.369	0.400	92.3	60.0	120	
Fenoxycarb	0.000	< 0.100		0.398	0.400	99.4	60.0	120	
Fenpyroximate	0.000	< 0.200		0.928	0.800	115.9	60.0	120	
Fipronil	0.000	< 0.200		0.810	0.800	101.3	60.0	120	
Flonicamid	0.000	< 0.250		1.050	1.000	105.0	60.0	120	
Fludioxonil	0.000	< 0.200		0.771	0.800	96.3	50.0	150	
Hexythiazox	0.000	< 0.250		0.952	1.000	95.2	60.0	120	
Imazalil	0.000	< 0.100		0.368	0.400	92.1	60.0	120	
Imidacloprid	0.000	< 0.200		0.795	0.800	99.4	60.0	120	
Kresoxim-methyl	0.000	< 0.200		0.757	0.800	94.6	60.0	120	
Malathion	0.000	< 0.100		0.389	0.400	97.3	60.0	120	
Metalaxyl	0.000	< 0.100		0.388	0.400	96.9	60.0	120	
Methiocarb	0.000	< 0.100		0.373	0.400	93.4	60.0	120	
Methomyl	0.000	< 0.200		0.858	0.800	107.2	60.0	120	
MGK-264	0.000	< 0.100		0.386	0.400	96.6	50.0	150	
Myclobutanil	0.000	< 0.100		0.381	0.400	95.2	60.0	120	
Naled	0.000	< 0.250		0.973	1.000	97.3	50.0	150	
Oxamyl	0.000	< 0.500		1.971	2.000	98.6	60.0	120	
Paclobutrazole	0.000	< 0.200		0.804	0.800	100.5	60.0	120	
Parathion-Methyl	0.000	< 0.100		0.399	0.400	99.6	50.0	150	
Permethrin	0.003	< 0.100		0.389	0.400	97.3	50.0	150	
Phosmet	0.000	< 0.100		0.399	0.400	99.7	50.0	150	
Piperonyl butoxide	0.000	< 0.500		1.888	2.000	94.4	60.0	120	
Prallethrin	0.000	< 0.100		0.385	0.400	96.4	60.0	120	
Propiconazole	0.000	< 0.200		0.786	0.800	98.2	60.0	120	
Propoxur	0.000	< 0.100		0.382	0.400	95.4	60.0	120	
Pyrethrin (Summe)	0.000	< 0.100		0.486	0.488	99.5	60.0	120	
Pyridaben	0.000	< 0.100		0.383	0.400	95.7	50.0	150	
Spinosad	0.000	< 0.100		0.377	0.388	97.2	50.0	150	
Spiromesifen	0.000	< 0.100		0.376	0.400	94.0	60.0	120	
Spirotetramat	0.000	< 0.100		0.402	0.400	100.4	60.0	120	
Spiroxamine	0.000	< 0.200		0.738	0.800	92.2	60.0	120	
Tebuconazole	0.000	< 0.200		0.786	0.800	98.3	60.0	120	
Thiacloprid	0.000	< 0.100		0.392	0.400	98.1	60.0	120	
Thiamethoxam	0.000	< 0.100		0.421	0.400	105.3	60.0	120	
Trifloxystrobin	0.000	< 0.100		0.380	0.400	95.0	60.0	120	

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Laboratory Pesticide Quality Control Results

AOAC2007.1 &EN 15662 | Units: mg/Kg | Batch ID: 2301287

Matrix Spike/Matrix Spike Duplicate Recoveries | Sample ID: 23-0016510001

Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Re	MSD % Re	Limits	Notes
Abamectin	0.000	0.653	0.764	1.000	15.6%	< 30	65.3%	76.4%	50 - 150	
Acephate	0.000	0.568	0.483	0.800	16.1%	< 30	71.0%	60.4%	50 - 150	
Acequinocyl	0.000	2.348	2.370	4.000	0.9%	< 30	58.7%	59.2%	50 - 150	
Acetamiprid	0.000	0.293	0.307	0.400	4.6%	< 30	73.2%	76.6%	50 - 150	
Aldicarb	0.000	0.660	0.694	0.800	5.0%	< 30	82.5%	86.7%	50 - 150	
Azoxystrobin	0.000	0.239	0.247	0.400	3.1%	< 30	59.8%	61.7%	50 - 150	
Bifenazate	0.000	0.284	0.312	0.400	9.3%	< 30	71.0%	78.0%	50 - 150	
Bifenthrin	0.000	0.221	0.223	0.400	0.8%	< 30	55.2%	55.6%	50 - 150	
Boscalid	0.044	0.467	0.487	0.800	4.5%	< 30	52.9%	55.3%	50 - 150	
Carbaryl	0.000	0.212	0.222	0.400	4.6%	< 30	53.1%	55.6%	50 - 150	
Carbofuran	0.000	0.240	0.245	0.400	2.3%	< 30	59.9%	61.3%	50 - 150	
Chlorantraniliprole	0.000	0.265	0.290	0.400	9.1%	< 30	66.2%	72.5%	50 - 150	
Chlorfenapyr	0.000	1.294	1.108	2.000	15.5%	< 30	64.7%	55.4%	50 - 150	
Chlorpyrifos	0.017	0.365	0.341	0.400	7.1%	< 30	87.1%	81.1%	50 - 150	
Clofentezine	0.010	0.191	0.180	0.400	6.2%	< 30	45.2%	42.4%	50 - 150	Q
Cyfluthrin	0.000	0.647	0.717	2.000	10.2%	< 30	32.4%	35.9%	30 - 150	
Cypermethrin	0.000	0.635	0.636	2.000	0.2%	< 30	31.7%	31.8%	50 - 150	Q
Daminozide	0.201	1.317	1.410	2.000	8.0%	< 30	55.8%	60.4%	30 - 150	
Diazinon	0.000	0.127	0.133	0.400	4.0%	< 30	31.9%	33.2%	50 - 150	Q
Dichlorvos	0.206	1.454	1.598	2.000	10.9%	< 30	62.4%	69.6%	50 - 150	
Dimethoate	0.000	0.340	0.358	0.400	5.4%	< 30	84.9%	89.6%	50 - 150	
Ethoprophos	0.000	0.210	0.217	0.400	2.9%	< 30	52.6%	54.2%	50 - 150	
Etofenprox	0.000	0.497	0.502	0.800	1.0%	< 30	62.1%	62.7%	50 - 150	
Etoxazole	0.000	0.323	0.319	0.400	1.2%	< 30	80.7%	79.7%	50 - 150	
Fenoxycarb	0.000	0.250	0.284	0.400	12.6%	< 30	62.5%	70.9%	50 - 150	
Fenpyroximate	0.000	0.219	0.242	0.800	10.2%	< 30	27.3%	30.3%	50 - 150	Q
Fipronil	0.000	0.357	0.371	0.800	3.8%	< 30	44.6%	46.4%	50 - 150	Q
Flonicamid	0.000	0.836	0.917	1.000	9.2%	< 30	83.6%	91.7%	50 - 150	
Fludioxonil	0.000	1.076	1.152	0.800	6.8%	< 30	134.5%	144.0%	50 - 150	
Hexythiazox	0.000	0.611	0.598	1.000	2.2%	< 30	61.1%	59.8%	50 - 150	
Imazalil	0.020	0.291	0.318	0.400	9.5%	< 30	67.8%	74.6%	50 - 150	
Imidacloprid	0.027	0.775	0.803	0.800	3.6%	< 30	93.5%	97.0%	50 - 150	
Kresoxim-methyl	0.000	0.495	0.540	0.800	8.7%	< 30	61.9%	67.5%	50 - 150	
Malathion	0.014	0.232	0.257	0.400	10.7%	< 30	54.7%	60.8%	50 - 150	
Metalaxyl	0.000	0.268	0.287	0.400	6.8%	< 30	66.9%	71.6%	50 - 150	
Methiocarb	0.000	0.228	0.240	0.400	5.2%	< 30	56.9%	59.9%	50 - 150	
Methomyl	0.000	0.765	0.801	0.800	4.6%	< 30	95.6%	100.1%	50 - 150	
MGK-264	0.000	0.114	0.125	0.400	8.6%	< 30	28.6%	31.1%	50 - 150	Q
Myclobutanil	0.000	0.120	0.119	0.400	0.7%	< 30	30.0%	29.8%	50 - 150	Q
Naled	0.000	0.528	0.542	1.000	2.6%	< 30	52.8%	54.2%	50 - 150	
Oxamyl	0.000	1.890	1.926	2.000	1.9%	< 30	94.5%	96.3%	50 - 150	
Paclobutrazole	0.000	0.406	0.469	0.800	14.3%	< 30	50.7%	58.6%	50 - 150	
Parathion-Methyl	0.000	0.089	0.091	0.400	2.0%	< 30	22.2%	22.6%	30 - 150	Q
Permethrin	0.005	0.236	0.229	0.400	2.9%	< 30	57.8%	56.1%	50 - 150	
Phosmet	0.000	0.212	0.244	0.400	13.8%	< 30	53.1%	60.9%	50 - 150	
Piperonyl butoxide	0.147	1.583	1.664	2.000	5.5%	< 30	71.8%	75.8%	50 - 150	
Prallethrin	0.054	0.158	0.167	0.400	8.3%	< 30	26.0%	28.2%	50 - 150	Q
Propiconazole	0.000	0.524	0.552	0.800	5.1%	< 30	65.5%	69.0%	50 - 150	
Propoxur	0.017	0.264	0.281	0.400	6.8%	< 30	61.6%	66.0%	50 - 150	
Pyrethrin (Summe)	0.178	0.485	0.490	0.488	1.6%	< 30	63.0%	64.0%	50 - 150	
Pyridaben	0.001	0.280	0.275	0.400	1.7%	< 30	69.8%	68.6%	50 - 150	
Spinosad	0.000	0.312	0.318	0.388	1.7%	< 30	80.5%	81.9%	50 - 150	
Spiromesifen	0.000	0.266	0.288	0.400	8.0%	< 30	66.5%	72.0%	50 - 150	
Spirotetramat	0.000	0.401	0.444	0.400	10.2%	< 30	100.3%	111.1%	50 - 150	
Spiroxamine	0.000	0.606	0.658	0.800	8.3%	< 30	75.8%	82.3%	50 - 150	
Tebuconazole	0.000	0.488	0.528	0.800	7.8%	< 30	61.0%	66.0%	50 - 150	
Thiacloprid	0.000	0.291	0.303	0.400	4.2%	< 30	72.6%	75.8%	50 - 150	
Thiamethoxam	0.000	0.364	0.397	0.400	8.7%	< 30	91.0%	99.3%	50 - 150	
Trifloxystrobin	0.000	0.275	0.287	0.400	4.2%	< 30	68.7%	71.6%	50 - 150	

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Laboratory Quality Control Results

J AOAC 2015 V98-6 **Batch ID: 2301306**

Laboratory Control Sample

Analyte	LCS	Result	Spike	Units	% Rec	Limits	Evaluation	Notes
CBDVA	2	0.117	0.101	%	115	80.0 - 120	Acceptable	
CBDV	2	0.128	0.108	%	118	80.0 - 120	Acceptable	
CBE	2	0.121	0.105	%	115	80.0 - 120	Acceptable	
CBDA	1	0.101	0.099	%	103	90.0 - 110	Acceptable	
CBGA	1	0.0987	0.096	%	102	80.0 - 120	Acceptable	
CBG	1	0.0977	0.094	%	104	80.0 - 120	Acceptable	
CBD	1	0.0979	0.096	%	102	90.0 - 110	Acceptable	
THCV	2	0.116	0.103	%	112	80.0 - 120	Acceptable	
d8THCV	2	0.119	0.107	%	112	80.0 - 120	Acceptable	
THCVA	2	0.112	0.097	%	116	80.0 - 120	Acceptable	
CBN	1	0.103	0.101	%	102	80.0 - 120	Acceptable	
exo-THC	2	0.112	0.101	%	111	80.0 - 120	Acceptable	
d9THC	1	0.101	0.100	%	100	90.0 - 110	Acceptable	
d8THC	1	0.104	0.100	%	103	90.0 - 110	Acceptable	
9S-d10THC	1	0.103	0.102	%	101	80.0 - 120	Acceptable	
CBL	2	0.112	0.102	%	109	80.0 - 120	Acceptable	
9S-HHC	3	0.0918	0.100	%	91.8	80.0 - 120	Acceptable	
9R-d10THC	1	0.0985	0.097	%	101	80.0 - 120	Acceptable	
CBC	2	0.119	0.106	%	112	80.0 - 120	Acceptable	
9R-HHC	3	0.0881	0.100	%	88.1	80.0 - 120	Acceptable	
THCA	1	0.0960	0.096	%	99.5	90.0 - 110	Acceptable	
CBCA	2	0.118	0.103	%	115	80.0 - 120	Acceptable	
CBLA	2	0.119	0.104	%	115	80.0 - 120	Acceptable	
d8THCO	3	0.0957	0.100	%	95.7	80.0 - 120	Acceptable	
CBT	2	0.117	0.106	%	110	80.0 - 120	Acceptable	
d9THCO	3	0.0970	0.100	%	97.0	80.0 - 120	Acceptable	

Method Blank

Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.077	%	< 0.077	Acceptable	
CBDV	<LOQ	0.077	%	< 0.077	Acceptable	
CBE	<LOQ	0.077	%	< 0.077	Acceptable	
CBDA	<LOQ	0.077	%	< 0.077	Acceptable	
CBGA	<LOQ	0.077	%	< 0.077	Acceptable	
CBG	<LOQ	0.077	%	< 0.077	Acceptable	
CBD	<LOQ	0.077	%	< 0.077	Acceptable	
THCV	<LOQ	0.077	%	< 0.077	Acceptable	
d8THCV	<LOQ	0.077	%	< 0.077	Acceptable	
THCVA	<LOQ	0.077	%	< 0.077	Acceptable	
CBN	<LOQ	0.077	%	< 0.077	Acceptable	
exo-THC	<LOQ	0.077	%	< 0.077	Acceptable	
d9THC	<LOQ	0.077	%	< 0.077	Acceptable	
d8THC	<LOQ	0.077	%	< 0.077	Acceptable	
9S-d10THC	<LOQ	0.077	%	< 0.077	Acceptable	
CBL	<LOQ	0.077	%	< 0.077	Acceptable	
9S-HHC	<LOQ	0.077	%	< 0.077	Acceptable	
9R-d10THC	<LOQ	0.077	%	< 0.077	Acceptable	
CBC	<LOQ	0.077	%	< 0.077	Acceptable	
9R-HHC	<LOQ	0.077	%	< 0.077	Acceptable	
THCA	<LOQ	0.077	%	< 0.077	Acceptable	
CBCA	<LOQ	0.077	%	< 0.077	Acceptable	
CBLA	<LOQ	0.077	%	< 0.077	Acceptable	
d8THCO	<LOQ	0.077	%	< 0.077	Acceptable	
CBT	<LOQ	0.077	%	< 0.077	Acceptable	
d9THCO	<LOQ	0.077	%	< 0.077	Acceptable	

Abbreviations

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

Units of Measure:

% - Percent

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Laboratory Quality Control Results

J AOAC 2015 V98-6 **Batch ID: 2301306**

Sample Duplicate Sample ID: **23-001332-0002**

Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBDV	0.413	0.413	0.077	%	0.0332	< 20	Acceptable	
CBE	4.02	4.04	0.077	%	0.435	< 20	Acceptable	
CBDA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBG	1.11	1.12	0.077	%	0.363	< 20	Acceptable	
CBD	67.8	67.8	0.077	%	0.0737	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBN	1.12	1.13	0.077	%	0.436	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
d9THC	0.0785	0.0798	0.077	%	1.69	< 20	Acceptable	
d8THC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
9S-d10THC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBL	0.359	0.354	0.077	%	1.38	< 20	Acceptable	
9S-HHC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
9R-d10THC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBC	1.42	1.43	0.077	%	1.10	< 20	Acceptable	
9R-HHC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBCA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
d8THCO	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBT	3.82	3.88	0.077	%	1.50	< 20	Acceptable	
d9THCO	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	

Abbreviations

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

Units of Measure:

% - Percent

Columbia LABORATORIES
A Tentamus Company

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

Report Number:	23-001623/D002.R000
Report Date:	02/10/2023
ORELAP#:	OR100028
Purchase Order:	
Received:	02/07/23 15:18

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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.