Columbia LABORATORIES
A Tentamus Company

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

Report Number: 26-000838/D006.R001
Report Date: 02/20/2026
ORELAP#: OR100028
Received: 01/21/26 09:55

This is an amended version of report# 26-000838/D006.R000.
Reason: Updated client information.

Customer: Secret Nature

Huntington Beach, CA 92649
United States of America (USA)

Product identity: Live D9 Gummy, Knockout Oregon Huckleberry 20mg
Material: Cannabinoid Edible
Laboratory ID: 26-000838-0004
Evidence of Cooling: No
Temp: 12.1 °C
Lot #: 3626.1NC_011926
Serving Size #1: 8 g

Sample Results

Potency **Method:** J AOAC 2015 V98-6 (mod)[þ] **Batch:** 2600626 **Analyzed:** 01/26/26

Analyte	Result	Units	LOQ	Notes	Serving Size #1 Result	Units	LOQ
CBC	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
CBC-A	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
CBC-Total	< LOQ	%	0.0144		< LOQ	mg/8g	1.15
CBD[⊥]	0.00870	%	0.0077		0.696	mg/8g	0.613
CBD-A[⊥]	0.0131	%	0.0077		1.05	mg/8g	0.613
CBD-Total[⊥]	0.0202	%	0.0144		1.62	mg/8g	1.15
CBDV	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
CBDV-A	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
CBDV-Total	< LOQ	%	0.0143		< LOQ	mg/8g	1.14
CBE	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
CBG	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
CBG-A	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
CBG-Total	< LOQ	%	0.0143		< LOQ	mg/8g	1.14
CBL	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
CBL-A	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
CBL-Total	< LOQ	%	0.0144		< LOQ	mg/8g	1.15
CBN	0.128	%	0.0077		10.2	mg/8g	0.613
CBT	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
Δ10-THC-9R	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
Δ10-THC-9S	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
Δ10-THC-Total	< LOQ	%	0.0153		< LOQ	mg/8g	1.23
Δ8-THC[⊥]	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
Δ8-THCV	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
Δ9-THC[⊥]	0.260	%	0.0077		20.8	mg/8g	0.613
Δ9-THC-A[⊥]	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
Δ9-THC-Total[⊥]	0.260	%	0.0144		20.8	mg/8g	1.15
Δ9-THCP	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
Δ9-THCV	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
Δ9-THCV-A	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
Δ9-THCV-Total	< LOQ	%	0.0143		< LOQ	mg/8g	1.14

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Potency

Method: J AOAC 2015 V98-6 (mod)[þ] **Batch:** 2600626 **Analyzed:** 01/26/26

Analyte	Result	Units	LOQ	Notes	Serving Size #1 Result	Units	LOQ
exo-THC	< LOQ	%	0.0077		< LOQ	mg/8g	0.613
Total Cannabinoids	0.410	%			32.8	mg/8g	

Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed	Method	Status	Notes
Salmonella spp. ⊥	Negative		/25g		2600524	01/24/26	AOAC 2020.02[þ]		
EHEC including STEC ⊥	Negative		/25g		2600525	01/24/26	AOAC 2020.06[þ]		

Solvents

Method: Residual Solvents by HS-GC-MS[þ] **Units** µg/g **Batch** 2600584 **Analyze:** 01/24/26

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 ⊥	< 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	

Pesticides

Method: AOAC 2007.01 & EN 15662 (mod) **Units** mg/kg **Batch** 2600666 **Analyze:** 01/28/26

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-ethyl ⊥	< LOQ	0.20	0.100	pass	
Clofentezine ⊥	< LOQ	0.20	0.100	pass		Cyfluthrin (sum) ⊥	< LOQ	1.0	0.500	pass	
Cypermethrin (sum) ⊥	< 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	

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Pesticides

Method: AOAC 2007.01 & EN 15662 (mod) **Units** mg/kg **Batch** 2600666 **Analyze:** 01/28/26

Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
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	0.0156	2600556	01/23/26	AOAC 2013.06 (mod.)þ	pass	
Cadmium ⊥	< LOQ	0.200	mg/kg	0.0156	2600556	01/23/26	AOAC 2013.06 (mod.)þ	pass	
Lead ⊥	< LOQ	0.500	mg/kg	0.0156	2600556	01/23/26	AOAC 2013.06 (mod.)þ	pass	
Mercury ⊥	< LOQ	0.100	mg/kg	0.00782	2600556	01/23/26	AOAC 2013.06 (mod.)þ	pass	

Mycotoxins

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed	Method	Status	Notes
Aflatoxin B1 ⊥	< LOQ		µg/kg	5.00	2600663	01/28/26	Mycotoxins by AOAC 2007.01		
Aflatoxin B2 ⊥	< LOQ		µg/kg	5.00	2600663	01/28/26	Mycotoxins by AOAC 2007.01		
Aflatoxin G1 ⊥	< LOQ		µg/kg	5.00	2600663	01/28/26	Mycotoxins by AOAC 2007.01		
Aflatoxin G2 ⊥	< LOQ		µg/kg	5.00	2600663	01/28/26	Mycotoxins by AOAC 2007.01		
Ochratoxin A ⊥	< LOQ	20.0	µg/kg	5.00	2600663	01/28/26	Mycotoxins by AOAC 2007.01	pass	
Total Aflatoxins	< LOQ	20.0	µg/kg	20.0		01/28/26	Mycotoxins by AOAC 2007.01þ	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.

Threshold Note: OAR 333-007-0400

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

Units of Measure

% wt = µg/g divided by 10,000
/25g = Per 25g
µg/g = Microgram per gram
µg/kg = Micrograms per kilogram = parts per billion (ppb)
mg/kg = Milligram per kilogram = parts per million (ppm)
% = Percentage of sample
mg/8g = Milligram per 8g

Approved Signatory

Derrick Tanner
General Manager

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

Laboratory Quality Control Results

Residual Solvents **Batch ID:** 2600584

Method Blank					Laboratory Control Sample					
Analyte	**Result**		**LOQ**	**Notes**	**Result**	**Spike**	**Units**	**% Rec**	**Limits**	**Notes**
1,1,1-Trichloroethane	ND	<	5		3.69	5	µg/g	73.8	50-150	
1,2-Dichloroethene, trans-	ND	<	1		0.668	1	µg/g	66.8	50-150	
1,2-Dichloroethene,cis-	ND	<	1		0.704	1	µg/g	70.4	50-150	
1,4-Dioxane	ND	<	100		418	509	µg/g	82.1	60-120	
1-Butanol	ND	<	500		1460	1610	µg/g	90.7	50-150	
2,2-Dimethylbutane	ND	<	30		147	188	µg/g	78.2	60-120	
2,2-Dimethylpropane	ND	<	200		852	956	µg/g	89.1	60-120	
2,3-Dimethylbutane	ND	<	30		147	188	µg/g	78.2	60-120	
2-Butanol	ND	<	200		1310	1640	µg/g	79.9	60-120	
2-Ethoxyethanol	ND	<	30		163	188	µg/g	86.7	60-120	
2-Methylbutane	ND	<	200		1330	1660	µg/g	80.1	60-120	
2-Methylpentane	ND	<	30		153	189	µg/g	81.0	60-120	
2-Propanol	ND	<	200		1400	1680	µg/g	83.3	60-120	
3-Methylpentane	ND	<	30		155	188	µg/g	82.4	60-120	
Acetone	ND	<	200		1400	1670	µg/g	83.8	60-120	
Acetonitrile	ND	<	100		422	511	µg/g	82.6	60-120	
Benzene	ND	<	1		0.696	1	µg/g	69.6	50-150	
Butane	ND	<	200		695	769	µg/g	90.4	60-120	
Chloroform	ND	<	1		0.752	1	µg/g	75.2	50-150	
Cumene	ND	<	30		164	192	µg/g	85.4	60-120	
Cyclohexane	ND	<	200		1390	1650	µg/g	84.2	60-120	
Dichloromethane	ND	<	1		0.671	1	µg/g	67.1	50-150	
DMSO	ND	<	500		1320	1620	µg/g	81.5	50-150	
Ethanol	ND	<	200		1370	1650	µg/g	83.0	60-120	
Ethyl acetate	ND	<	200		1350	1630	µg/g	82.8	60-120	
Ethyl Ether	ND	<	200		1330	1630	µg/g	81.6	60-120	
Ethylbenzene	ND	<	200		845	996	µg/g	84.8	60-120	
Ethylene Glycol	ND	<	200		426	520	µg/g	81.9	60-120	
Ethylene Oxide	ND	<	1		0.688	1	µg/g	68.8	50-150	
Heptane	ND	<	200		1360	1630	µg/g	83.4	60-120	
Hexane	ND	<	30		150	191	µg/g	78.5	60-120	
Isobutane	ND	<	200		682	770	µg/g	88.6	60-120	
Isobutyl Acetate	ND	<	500		1530	1640	µg/g	93.3	50-150	
Isopropyl Acetate	ND	<	200		1430	1660	µg/g	86.1	60-120	
m,p-Xylene	ND	<	200		892	1030	µg/g	86.6	60-120	
Methanol	ND	<	200		1360	1660	µg/g	81.9	60-120	
Methylethylketone	ND	<	500		1510	1620	µg/g	93.2	50-150	
Methylisobutylketone	ND	<	500		1540	1620	µg/g	95.1	50-150	
MTBE	ND	<	500		1430	1610	µg/g	88.8	50-150	
N,N-dimethylformamide	ND	<	150		484	492	µg/g	98.4	50-150	
o-Xylene	ND	<	200		867	996	µg/g	87.0	60-120	
Pentane	ND	<	200		1300	1630	µg/g	79.8	60-120	
Propane	ND	<	200		538	585	µg/g	92.0	60-120	
Tetrahydrofuran	ND	<	100		448	519	µg/g	86.3	60-120	
Toluene	ND	<	100		452	518	µg/g	87.3	60-120	
Triethylamine	ND	<	500		1380	1610	µg/g	85.7	50-150	

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QC - Sample Duplicate **Sample ID: 26-000872-0001**

Analyte	SR Result	SD Result	LOQ	Units	RPD	Limits	Accept/Fail	Notes
1,1,1-Trichloroethane	ND	ND	5	µg/g	0.0	< 20	Acceptable	
1,2-Dichloroethene, trans-	ND	ND	1	µg/g	0.0	< 20	Acceptable	
1,2-Dichloroethene,cis-	ND	ND	1	µg/g	0.0	< 20	Acceptable	
1,4-Dioxane	ND	ND	100	µg/g	0.0	< 20	Acceptable	
1-Butanol	ND	ND	500	µg/g	0.0	< 20	Acceptable	
2,2-Dimethylbutane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
2,2-Dimethylpropane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2,3-Dimethylbutane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
2-Butanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2-Ethoxyethanol	ND	ND	30	µg/g	0.0	< 20	Acceptable	
2-Methylbutane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
2-Methylpentane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
2-Propanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
3-Methylpentane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Acetone	846	1120	200	µg/g	27.9	< 20	FAIL	R, E2
Acetonitrile	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Benzene	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Butane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Chloroform	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Cumene	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Cyclohexane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Dichloromethane	ND	ND	1	µg/g	0.0	< 20	Acceptable	
DMSO	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Ethanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethyl acetate	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethyl Ether	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethylbenzene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethylene Glycol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Ethylene Oxide	ND	ND	1	µg/g	0.0	< 20	Acceptable	
Heptane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Hexane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Isobutane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Isobutyl Acetate	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Isopropyl Acetate	ND	ND	200	µg/g	0.0	< 20	Acceptable	
m,p-Xylene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Methanol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Methylethylketone	ND	ND	500	µg/g	0.0	< 20	Acceptable	
Methylisobutylketone	ND	ND	500	µg/g	0.0	< 20	Acceptable	
MTBE	ND	ND	500	µg/g	0.0	< 20	Acceptable	
N,N-dimethylformamide	ND	ND	150	µg/g	0.0	< 20	Acceptable	
o-Xylene	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Pentane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Propane	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Tetrahydrofuran	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Toluene	ND	ND	100	µg/g	0.0	< 20	Acceptable	
Triethylamine	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

Units of Measure:

µg/g- Microgram per gram or ppm

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Portland, OR 97230
503-254-1794

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Laboratory Quality Control Results

J AOAC 2015 V98-6 **Batch ID: 2600626**

Laboratory Control Sample

Analyte	LCS	Result	Spike	Units	% Rec	Limits	Evaluation	Notes
CBDVA	2	0.0303	0.0295	%	103	80.0 - 120	Acceptable	
CBDV	2	0.0318	0.0311	%	102	80.0 - 120	Acceptable	
CBE	2	0.0344	0.0341	%	101	80.0 - 120	Acceptable	
CBDA	1	0.0293	0.0288	%	102	90.0 - 110	Acceptable	
CBGA	1	0.0304	0.0300	%	101	80.0 - 120	Acceptable	
CBG	1	0.0283	0.0283	%	99.8	80.0 - 120	Acceptable	
CBD	1	0.0273	0.0264	%	103	90.0 - 110	Acceptable	
THCV	2	0.0315	0.0313	%	101	80.0 - 120	Acceptable	
d8THCV	2	0.0322	0.0328	%	98.1	80.0 - 120	Acceptable	
THCVA	2	0.0576	0.0558	%	103	80.0 - 120	Acceptable	
CBN	1	0.0282	0.0277	%	102	80.0 - 120	Acceptable	
exo-THC	2	0.0304	0.0298	%	102	80.0 - 120	Acceptable	
d9THC	1	0.0277	0.0274	%	101	90.0 - 110	Acceptable	
d8THC	1	0.0293	0.0293	%	100	90.0 - 110	Acceptable	
9S-d10THC	1	0.0310	0.0308	%	101	80.0 - 120	Acceptable	
CBL	2	0.0298	0.0292	%	102	80.0 - 120	Acceptable	
9R-d10THC	1	0.0332	0.0330	%	101	80.0 - 120	Acceptable	
CBC	2	0.0319	0.0314	%	102	80.0 - 120	Acceptable	
THCA	1	0.0287	0.0289	%	99.5	90.0 - 110	Acceptable	
CBCA	2	0.0308	0.0306	%	101	80.0 - 120	Acceptable	
CBLA	2	0.0309	0.0306	%	101	80.0 - 120	Acceptable	
d9THCP	2	0.0298	0.0293	%	101	80.0 - 120	Acceptable	
CBT	2	0.0313	0.0316	%	99.0	80.0 - 120	Acceptable	

Method Blank

Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.00776	%	< 0.00776	Acceptable	
CBDV	<LOQ	0.00776	%	< 0.00776	Acceptable	
CBE	<LOQ	0.00776	%	< 0.00776	Acceptable	
CBDA	<LOQ	0.00776	%	< 0.00776	Acceptable	
CBGA	<LOQ	0.00776	%	< 0.00776	Acceptable	
CBG	<LOQ	0.00776	%	< 0.00776	Acceptable	
CBD	<LOQ	0.00776	%	< 0.00776	Acceptable	
THCV	<LOQ	0.00776	%	< 0.00776	Acceptable	
d8THCV	<LOQ	0.00776	%	< 0.00776	Acceptable	
THCVA	<LOQ	0.00776	%	< 0.00776	Acceptable	
CBN	<LOQ	0.00776	%	< 0.00776	Acceptable	
exo-THC	<LOQ	0.00776	%	< 0.00776	Acceptable	
d9THC	<LOQ	0.00776	%	< 0.00776	Acceptable	
d8THC	<LOQ	0.00776	%	< 0.00776	Acceptable	
9S-d10THC	<LOQ	0.00776	%	< 0.00776	Acceptable	
CBL	<LOQ	0.00776	%	< 0.00776	Acceptable	
9R-d10THC	<LOQ	0.00776	%	< 0.00776	Acceptable	
CBC	<LOQ	0.00776	%	< 0.00776	Acceptable	
THCA	<LOQ	0.00776	%	< 0.00776	Acceptable	
CBCA	<LOQ	0.00776	%	< 0.00776	Acceptable	
CBLA	<LOQ	0.00776	%	< 0.00776	Acceptable	
d9THCP	<LOQ	0.00776	%	< 0.00776	Acceptable	
CBT	<LOQ	0.00776	%	< 0.00776	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: 2600626**

Sample Duplicate Sample ID: **26-000835-0001**

Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
CBDV	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
CBE	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
CBDA	0.0101	0.0102	0.00770	%	0.416	< 10	Acceptable	
CBGA	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
CBG	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
CBD	<LOQ	<LOQ	0.00770	%	NA	< 10	Acceptable	
THCV	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
CBN	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
d9THC	0.220	0.217	0.00770	%	1.15	< 10	Acceptable	
d8THC	<LOQ	<LOQ	0.00770	%	NA	< 10	Acceptable	
9S-d10THC	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
9R-d10THC	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
CBC	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.00770	%	NA	< 10	Acceptable	
CBCA	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
d9THCP	<LOQ	<LOQ	0.00770	%	NA	< 20	Acceptable	
CBT	<LOQ	<LOQ	0.00770	%	NA	< 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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Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662 **Units:** mg/Kg **Batch ID:** 2600666

Method Blank				Laboratory Control Sample					
Analyte	**Blank Result**	**Blank Limits**	**Notes**	**LCS Result**	**LCS Spike**	**LCS % Rec**	**Limits**		**Notes**
Abamectin	0.000	< 0.070		0.290	0.280	103.5	50.0	150	
Acephate	0.000	< 0.020		0.074	0.080	92.7	60.0	120	
Acequinocyl	0.000	< 0.020		0.107	0.080	133.7	40.0	160	
Acetamiprid	0.000	< 0.020		0.068	0.080	84.4	60.0	120	
Aldicarb	0.000	< 0.100		0.379	0.400	94.7	60.0	120	
Azoxystrobin	0.001	< 0.010		0.037	0.040	92.0	60.0	120	
Bifenazate	0.001	< 0.010		0.041	0.040	103.1	60.0	120	
Bifenthrin	0.000	< 0.100		0.345	0.400	86.3	50.0	150	
Boscalid	0.001	< 0.010		0.037	0.040	92.9	60.0	120	
Carbaryl	0.000	< 0.020		0.069	0.080	86.7	60.0	120	
Carbofuran	0.000	< 0.010		0.038	0.040	94.2	60.0	120	
Chlorantraniliprole	0.000	< 0.010		0.036	0.040	90.4	60.0	120	
Chlorfenapyr	0.000	< 0.050		0.230	0.200	115.1	60.0	120	
Chlorpyrifos	0.000	< 0.010		0.033	0.040	82.4	60.0	120	
Clofentezine	0.001	< 0.010		0.036	0.040	89.5	60.0	120	
Cyfluthrin	0.000	< 0.200		0.643	0.800	80.4	50.0	150	
Cypermethrin	0.009	< 0.200		0.701	0.800	87.6	50.0	150	
Daminozide	0.050	< 0.050		0.186	0.200	93.1	60.0	120	
Diazinon	0.002	< 0.010		0.036	0.040	90.9	60.0	120	
Dichlorvos	0.000	< 0.050		0.179	0.200	89.5	60.0	120	
Dimethoate	0.001	< 0.010		0.036	0.040	89.1	60.0	120	
Ethoprophos	0.001	< 0.010		0.036	0.040	90.4	60.0	120	
Etofenprox	0.002	< 0.010		0.036	0.040	89.9	50.0	150	
Etoxazole	0.000	< 0.010		0.035	0.040	88.7	60.0	120	
Fenoxycarb	0.000	< 0.010		0.041	0.040	103.0	60.0	120	
Fenpyroximate	0.003	< 0.020		0.073	0.080	91.1	60.0	120	
Fipronil	0.000	< 0.010		0.040	0.040	101.2	60.0	120	
Flonicamid	0.000	< 0.020		0.076	0.080	95.1	60.0	120	
Fludioxonil	0.000	< 0.010		0.036	0.040	89.3	50.0	150	
Hexythiazox	0.000	< 0.010		0.036	0.040	88.9	60.0	120	
Imazalil	0.001	< 0.010		0.038	0.040	96.2	60.0	120	
Imidacloprid	0.000	< 0.010		0.035	0.040	88.1	60.0	120	
Kresoxim-methyl	0.000	< 0.010		0.036	0.040	91.0	60.0	120	
Malathion	0.000	< 0.010		0.037	0.040	92.7	60.0	120	
Metalaxyl	0.001	< 0.010		0.035	0.040	88.6	60.0	120	
Methiocarb	0.000	< 0.010		0.036	0.040	90.0	60.0	120	
Methomyl	0.000	< 0.020		0.069	0.080	86.9	60.0	120	
MGK-264	0.002	< 0.050		0.179	0.200	89.6	50.0	150	
Myclobutanil	0.000	< 0.010		0.036	0.040	91.0	60.0	120	
Naled	0.012	< 0.100		0.379	0.400	94.6	50.0	150	
Oxamyl	0.000	< 0.200		0.737	0.800	92.2	60.0	120	
Paclobutrazole	0.001	< 0.010		0.037	0.040	92.2	60.0	120	
Parathion-Methyl	0.010	< 0.030		0.113	0.120	93.8	50.0	150	
Permethrin	0.000	< 0.040		0.132	0.160	82.3	50.0	150	
Phosmet	0.000	< 0.010		0.047	0.040	118.1	50.0	150	
Piperonyl butoxide	0.000	< 0.200		0.754	0.800	94.3	60.0	120	
Prallethrin	0.002	< 0.050		0.173	0.200	86.4	60.0	120	
Propiconazole	0.001	< 0.010		0.036	0.040	89.9	60.0	120	
Propoxur	0.000	< 0.010		0.037	0.040	92.3	60.0	120	
Pyrethrin (Summe)	0.005	< 0.100		0.099	0.100	99.0	60.0	120	
Pyridaben	0.002	< 0.020		0.071	0.080	88.5	50.0	150	
Spinosad	0.000	< 0.100		0.035	0.040	87.7	50.0	150	
Spiromesifen	0.002	< 0.030		0.106	0.120	88.1	60.0	120	
Spirotetramat	0.000	< 0.010		0.039	0.040	96.6	60.0	120	
Spiroxamine	0.000	< 0.010		0.036	0.040	90.3	60.0	120	

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Laboratory Pesticide Quality Control Results

AOAC 2007.1 & EN 15662 **Units:** mg/Kg **Batch ID:** 2600666

Matrix Spike/Matrix Spike Duplicate Recoveries **Sample ID: 26-000838-0001**

Analyte	Result	MS Res	MSD Res	Spike	RPD%	Limit	MS % Rec	MSD % Rec	Limits	Notes
Abamectin	0.000	0.243	0.225	0.280	7.6%	< 30	86.7%	80.3%	50 - 150	
Acephate	0.000	0.082	0.076	0.080	7.9%	< 30	102.2%	94.4%	50 - 150	
Acequinocyl	0.027	0.103	0.105	0.080	2.4%	< 30	94.8%	97.1%	50 - 150	
Acetamiprid	0.000	0.075	0.072	0.080	3.4%	< 30	93.6%	90.5%	50 - 150	
Aldicarb	0.000	0.355	0.343	0.400	3.3%	< 30	88.7%	85.8%	50 - 150	
Azoxystrobin	0.001	0.033	0.033	0.040	1.6%	< 30	81.5%	80.2%	50 - 150	
Bifenazate	0.001	0.039	0.037	0.040	4.8%	< 30	95.0%	90.5%	50 - 150	
Bifenthrin	0.000	0.345	0.334	0.400	3.4%	< 30	86.3%	83.5%	50 - 150	
Boscalid	0.001	0.034	0.040	0.040	16.9%	< 30	83.2%	98.6%	50 - 150	
Carbaryl	0.000	0.071	0.063	0.080	12.4%	< 30	88.7%	78.3%	50 - 150	
Carbofuran	0.000	0.033	0.034	0.040	2.0%	< 30	82.6%	84.3%	50 - 150	
Chlorantraniliprole	0.000	0.032	0.034	0.040	5.7%	< 30	79.6%	84.2%	50 - 150	
Chlorfenapyr	0.000	0.345	0.428	0.200	21.4%	< 30	172.7%	214.2%	50 - 150	Q
Chlorpyrifos	0.000	0.035	0.032	0.040	10.1%	< 30	88.5%	80.0%	50 - 150	
Clofentezine	0.001	0.035	0.034	0.040	4.4%	< 30	87.0%	83.3%	50 - 150	
Cyfluthrin	0.000	0.666	0.650	0.800	2.4%	< 30	83.3%	81.3%	30 - 150	
Cypermethrin	0.009	0.760	0.761	0.800	0.1%	< 30	93.8%	94.0%	50 - 150	
Daminozide	0.112	0.240	0.240	0.200	0.4%	< 30	63.7%	64.0%	30 - 150	
Diazinon	0.002	0.035	0.036	0.040	1.6%	< 30	83.7%	85.1%	50 - 150	
Dichlorvos	0.000	0.170	0.174	0.200	2.1%	< 30	85.2%	87.0%	50 - 150	
Dimethoate	0.001	0.036	0.035	0.040	1.8%	< 30	86.9%	85.3%	50 - 150	
Ethoprophos	0.001	0.035	0.036	0.040	2.1%	< 30	85.7%	87.5%	50 - 150	
Etofenprox	0.000	0.035	0.037	0.040	5.8%	< 30	86.4%	91.6%	50 - 150	
Etoxazole	0.000	0.035	0.034	0.040	4.7%	< 30	88.2%	84.1%	50 - 150	
Fenoxycarb	0.001	0.036	0.032	0.040	12.2%	< 30	87.2%	77.2%	50 - 150	
Fenpyroximate	0.003	0.071	0.071	0.080	1.0%	< 30	85.3%	86.2%	50 - 150	
Fipronil	0.000	0.033	0.038	0.040	14.1%	< 30	82.4%	94.9%	50 - 150	
Flonicamid	0.000	0.074	0.067	0.080	10.0%	< 30	93.0%	84.2%	50 - 150	
Fludioxonil	0.000	0.029	0.033	0.040	12.3%	< 30	72.3%	81.8%	50 - 150	
Hexythiazox	0.000	0.041	0.039	0.040	4.7%	< 30	102.1%	97.4%	50 - 150	
Imazalil	0.001	0.034	0.038	0.040	12.0%	< 30	80.7%	91.0%	50 - 150	
Imidacloprid	0.000	0.036	0.033	0.040	8.6%	< 30	89.6%	82.2%	50 - 150	
Kresoxim-methyl	0.000	0.032	0.034	0.040	6.0%	< 30	79.0%	83.8%	50 - 150	
Malathion	0.000	0.035	0.034	0.040	2.8%	< 30	87.1%	84.7%	50 - 150	
Metalaxyl	0.000	0.034	0.033	0.040	4.7%	< 30	85.8%	81.9%	50 - 150	
Methiocarb	0.000	0.037	0.034	0.040	6.9%	< 30	90.3%	84.3%	50 - 150	
Methomyl	0.000	0.066	0.072	0.080	9.0%	< 30	82.4%	90.1%	50 - 150	
MGK-264	0.000	0.173	0.172	0.200	0.6%	< 30	86.6%	86.1%	50 - 150	
Myclobutanil	0.000	0.034	0.033	0.040	1.4%	< 30	83.7%	82.5%	50 - 150	
Naled	0.000	0.372	0.388	0.400	4.4%	< 30	92.9%	97.0%	50 - 150	
Oxamyl	0.000	0.715	0.693	0.800	3.2%	< 30	89.4%	86.6%	50 - 150	
Paclobutrazole	0.000	0.035	0.034	0.040	2.8%	< 30	88.6%	86.2%	50 - 150	
Parathion-Methyl	0.015	0.110	0.105	0.120	5.0%	< 30	79.3%	75.4%	30 - 150	
Permethrin	0.000	0.136	0.154	0.160	12.8%	< 30	84.8%	96.4%	50 - 150	
Phosmet	0.000	0.051	0.046	0.040	9.9%	< 30	128.3%	116.2%	50 - 150	
Piperonyl butoxide	0.000	0.725	0.741	0.800	2.2%	< 30	90.7%	92.7%	50 - 150	
Prallethrin	0.000	0.168	0.172	0.200	2.3%	< 30	83.9%	85.8%	50 - 150	
Propiconazole	0.000	0.033	0.037	0.040	12.2%	< 30	82.9%	93.7%	50 - 150	
Propoxur	0.000	0.038	0.036	0.040	4.8%	< 30	95.0%	90.6%	50 - 150	
Pyrethrin (Summe)	0.004	0.111	0.099	0.100	12.1%	< 30	107.8%	95.5%	50 - 150	
Pyridaben	0.002	0.078	0.077	0.080	2.0%	< 30	95.3%	93.4%	50 - 150	
Spinosad	0.000	0.034	0.036	0.040	7.9%	< 30	83.8%	90.6%	50 - 150	
Spiromesifen	0.000	0.099	0.106	0.120	7.8%	< 30	82.1%	88.7%	50 - 150	
Spirotetramat	0.000	0.036	0.036	0.040	0.2%	< 30	90.0%	89.8%	50 - 150	
Spiroxamine	0.000	0.034	0.036	0.040	3.3%	< 30	85.9%	88.8%	50 - 150	

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