



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



Report Number: 22-011232/D002.R000
Report Date: 09/27/2022
ORELAP#: OR100028
Purchase Order:
Received: 09/19/22 16:00

Customer: Brothers Apothecary
Product identity: Hemp Distillate
Client/Metric ID: .
Laboratory ID: 22-011232-0003

Summary

Potency:

Analyte	Result (%)		
CBD	69.9	<ul style="list-style-type: none"> ● CBD ● CBN ● CBG ● Δ9-THC 	CBD-Total 69.9%
CBN	2.66		THC-Total 0.283%
CBG	1.01		(Reported in percent of total sample)
Δ9-THC	0.283		

Residual Solvents:

All analytes passing and less than LOQ.

Pesticides:

All analytes passing and less than LOQ.

Metals:

Less than LOQ for all analytes.

Microbiology:

Less than LOQ for all analytes.



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Customer: Brothers Apothecary
3423 SE18th Ave
Portland OR, 97202
United States of America (USA)

Product identity: Hemp Distillate

Client/Metric ID: .

Sample Date:

Laboratory ID: 22-011232-0003

Evidence of Cooling: No

Temp: 29 °C

Relinquished by: client

Sample Results

Potency	Method: J AOAC 2015 V98-6 (mod) ^p			Units %	Batch: 2207941	Analyze: 9/20/22 9:56:00 PM
Analyte	As Received	Dry weight	LOQ	Notes		
CBD	69.9		0.734			
CBD-A	< LOQ		0.0734			
CBD-Total	69.9		0.798			
CBG	1.01		0.0734			
CBG-A	< LOQ		0.0734			
CBG-Total	1.01		0.137			
CBN	2.66		0.0734			
Δ10-THC	< LOQ		0.0734			
Δ8-THC	< LOQ		0.0734			
Δ9-THC	0.283		0.0734			
THC-A	< LOQ		0.0734			
THC-Total	0.283		0.138			
Total Cannabinoids	73.9					

Microbiology								
Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ	100.00	cfu/g	10	2207922	09/23/22 AOAC 991.14 (Petrifilm) ^p	pass	
Total Coliforms	< LOQ	100.00	cfu/g	10	2207922	09/23/22 AOAC 991.14 (Petrifilm) ^p	pass	
Mold (RAPID Petrifilm)	< LOQ	1,000.	cfu/g	10	2207923	09/24/22 AOAC 2014.05 (RAPID) ^p	pass	
Yeast (RAPID Petrifilm)	< LOQ	1,000.	cfu/g	10	2207923	09/24/22 AOAC 2014.05 (RAPID) ^p	pass	
Salmonella spp. by PCR	Negative		/15g		2207930	09/22/22 AOAC 2020.02 ^b		
EHEC including STEC	Negative		/15g		2207932	09/22/22 AOAC RI 121806 ^b		



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Solvents											Method: Residual Solvents by GC/MS ^b					Units µg/g		Batch 2208097		Analyze 09/26/22 12:54 PM			
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes												
1-Butanol	< LOQ		500			1-Pentanol	< LOQ		500														
1,1-Dichloroethane	< LOQ		1.00			1,2-Dichloroethane	< LOQ		1.00														
1,2-Dimethoxyethane	< LOQ		50.0			1,4-Dioxane	< LOQ		100														
2-Butanol	< LOQ		200			2-Ethoxyethanol	< LOQ		30.0														
2-methyl-1-propanol	< LOQ		500			2-Methylbutane (Isopentane)	< LOQ	1000	200	pass													
2-Methylpentane	< LOQ	60.0	30.0	pass		2-Propanol (IPA)	< LOQ	1000	200	pass													
2,2-Dimethylbutane	< LOQ	60.0	30.0	pass		2,2-Dimethylpropane (neo-pentane)	< LOQ	1000	200	pass													
2,3-Dimethylbutane	< LOQ	60.0	30.0	pass		3-Methyl-(1)-Butanol	< LOQ		500														
3-Methylpentane	< LOQ	60.0	30.0	pass		Acetic Acid	< LOQ		250														
Acetone	< LOQ	1000	200	pass		Acetonitrile	< LOQ		100														
Anisole	< LOQ		500			Benzene	< LOQ	2.00	1.00	pass													
Butanes (sum)	< LOQ	1000	400	pass		Butyl acetate	< LOQ		500														
Chloroform	< LOQ		1.00			Cyclohexane	< LOQ		200														
DMSO	< LOQ		500			Ethanol	< LOQ	1000	200	pass													
Ethyl acetate	< LOQ	1000	200	pass		Ethyl benzene	< LOQ		200														
Ethyl ether	< LOQ		200			Ethyl Formate	< LOQ		500														
Ethylene glycol	< LOQ		200			Ethylene oxide	< LOQ		1.00														
Formic Acid	< LOQ		250			Hexanes (sum)	< LOQ	60.0	150	pass													
Isobutyl acetate	< LOQ		500			Isopropyl acetate	< LOQ		200														
Isopropylbenzene (Cumene)	< LOQ		30.0			m,p-Xylene	< LOQ	430	200	pass													
Methanol	< LOQ	600	200	pass		Methyl-t-butyl ether	< LOQ		500														
Methylacetat	< LOQ		500			Methylene chloride	< LOQ		1.00														
Methylethylketone	< LOQ		500			Methylisobutylketone	< LOQ		500														
Methylpropane (Isobutane)	< LOQ	1000	200	pass		n-Butane	< LOQ	1000	200	pass													
n-Heptane	< LOQ	1000	200	pass		n-Hexane	< LOQ	60.0	30.0	pass													
n-Pentane	< LOQ	1000	200	pass		n-Propanol	< LOQ		500														
N,N-dimethylacetamide	< LOQ		200			N,N-dimethylformamide	< LOQ		200														
o-Xylene	< LOQ	430	200	pass		Pentanes (sum)	< LOQ	1000	600	pass													
Propane	< LOQ	1000	200	pass		Propyl Acetate	< LOQ		500														
Pyridine	< LOQ		50.0			Sulfolane	pending		50.0														
Tetrahydrofuran	< LOQ		100			Toluene	< LOQ	180	100	pass													
Total Residual Solvents	< LOQ		5,000			Total Xylenes	< LOQ	430	400	pass													
Total Xylenes and Ethyl benzene	< LOQ		600			Trichloroethylene	< LOQ		1.00														
Triethylamine	< LOQ		500																				



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Pesticides											
Method: AOAC 2007.01 & EN 15662 (mod)						Units mg/kg		Batch 2207934		Analyze 09/21/22 09:07 AM	
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
Abamectin	< LOQ	0.25	0.070	pass		Acephate	< LOQ	0.050	0.020	pass	
Acequinocyl	< LOQ	0.030	0.025	pass		Acetamidiprid	< LOQ	0.050	0.050	pass	
Aldicarb	< LOQ	0.50	0.100	pass		Allethrin	< LOQ	0.10	0.100	pass	
Atrazine	< LOQ	0.0250	0.025	pass		Azadirachtin	< LOQ	1.0	0.500	pass	
Azoxystrobin	< LOQ	0.010	0.010	pass		Benzovindiflupyr	< LOQ	0.010	0.010	pass	
Bifenazate	< LOQ	0.010	0.010	pass		Bifenthrin	< LOQ	1.0	0.100	pass	
Boscalid	< LOQ	0.010	0.010	pass		Buprofezin	< LOQ	0.020	0.010	pass	
Captan	< LOQ		0.700			Carbaryl	< LOQ	0.025	0.025	pass	
Carbofuran	< LOQ	0.010	0.010	pass		Chlorantraniliprole	< LOQ	0.020	0.010	pass	
Chlordane (cis+trans)	< LOQ		0.100			Chlorfenapyr	< LOQ	1.5	0.100	pass	
Chlorpyrifos	< LOQ	0.50	0.010	pass		Clofentezine	< LOQ	0.010	0.010	pass	
Clothianidin	< LOQ	0.025	0.025	pass		Coumaphos	< LOQ	0.010	0.010	pass	
Cyantraniliprole	< LOQ	0.010	0.010	pass		Cyfluthrin	< LOQ	0.20	0.400	pass	
Cyhalothrin,lambda	< LOQ	0.0200	0.250	pass		Cypermethrin	< LOQ	0.30	0.300	pass	
Cyprodinil	< LOQ	0.010	0.010	pass		Daminozide	< LOQ	0.10	0.050	pass	
Deltamethrin	< LOQ	0.50	0.500	pass		Diazinon	< LOQ	0.020	0.010	pass	
Dichlorvos	< LOQ	0.050	0.050	pass		Dimethoate	< LOQ	0.010	0.010	pass	
Dimethomorph	< LOQ	0.050	0.050	pass		Dinotefuran	< LOQ	0.050	0.050	pass	
Diuron	< LOQ	0.125	0.125	pass		Dodemorph	< LOQ	0.050	0.050	pass	
Endosulfan I (alpha)	< LOQ	2.5	0.050	pass		Endosulfan II (beta)	< LOQ	2.5	0.050	pass	
Endosulfan sulfate	< LOQ	2.5	0.050	pass		Ethoprophos	< LOQ	0.010	0.010	pass	
Etofenprox	< LOQ	0.050	0.010	pass		Etoxazole	< LOQ	0.020	0.010	pass	
Etridiazole	< LOQ	0.15	0.050	pass		Fenhexamid	< LOQ	0.13	0.100	pass	
Fenoxycarb	< LOQ	0.010	0.010	pass		Fenpyroximate	< LOQ	0.020	0.020	pass	
Fensulfothion	< LOQ	0.010	0.010	pass		Fenthion	< LOQ	0.010	0.010	pass	
Fenvalerate	< LOQ		0.200			Fipronil	< LOQ	0.010	0.010	pass	
Flonicamid	< LOQ	0.025	0.025	pass		Fludioxonil	< LOQ	0.010	0.010	pass	
Fluopyram	< LOQ	0.010	0.010	pass		Hexythiazox	< LOQ	0.010	0.010	pass	
Imazalil	< LOQ	0.010	0.010	pass		Imidacloprid	< LOQ	0.010	0.010	pass	
Iprodione	< LOQ	0.50	0.500	pass		Kinoprene	< LOQ	1.3	0.050	pass	
Kresoxim-methyl	< LOQ	0.15	0.010	pass		Malathion	< LOQ	0.010	0.010	pass	
Metalaxyl	< LOQ	0.010	0.010	pass		Methiocarb	< LOQ	0.010	0.010	pass	
Methomyl	< LOQ	0.025	0.025	pass		Methoprene	< LOQ	2.0	1.00	pass	
Mevinphos	< LOQ	0.025	0.025	pass		MGK-264	< LOQ	0.050	0.050	pass	
Myclobutanil	< LOQ	0.010	0.010	pass		Naled	< LOQ	0.10	0.100	pass	
Novaluron	< LOQ	0.025	0.025	pass		Oxamyl	< LOQ	1.5	0.500	pass	
Paclbutrazole	< LOQ	0.010	0.010	pass		Parathion-Methyl	< LOQ	0.050	0.030	pass	
Permethrin	< LOQ	0.50	0.040	pass		Phenothrin	< LOQ	0.050	0.025	pass	
Phosmet	< LOQ	0.020	0.010	pass		Piperonyl butoxide	< LOQ	1.3	0.200	pass	
Pirimicarb	< LOQ	0.010	0.010	pass		Prallethrin	< LOQ	0.050	0.050	pass	
Propiconazole	< LOQ	0.10	0.010	pass		Propoxur	< LOQ	0.010	0.010	pass	
Pyraclostrobin	< LOQ	0.010	0.010	pass		Pyrethrins (total)	< LOQ	0.050	0.025	pass	
Pyridaben	< LOQ	0.020	0.020	pass		Pyriproxyfen	< LOQ	0.0100	0.010	pass	
Quintozene	< LOQ	0.020	0.020	pass		Resmethrin	< LOQ	0.050	0.020	pass	
Spinetoram	< LOQ	0.010	0.010	pass		Spinosad	< LOQ	0.010	0.010	pass	
Spirodiclofen	< LOQ	0.25	0.250	pass		Spiromesifen	< LOQ	3.0	0.030	pass	



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Pesticides											
Method: AOAC 2007.01 & EN 15662 (mod)						Units mg/kg	Batch 2207934	Analyze 09/21/22 09:07 AM			
Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
Spirotetramat	< LOQ	0.010	0.010	pass		Spiroxamine	< LOQ	0.10	0.010	pass	
Tebuconazole	< LOQ	0.010	0.010	pass		Tebufenozide	< LOQ	0.010	0.010	pass	
Teflubenzuron	< LOQ	0.025	0.025	pass		Tetrachlorvinphos	< LOQ	0.010	0.010	pass	
Tetramethrin	< LOQ	0.10	0.050	pass		Thiabendazole	< LOQ	0.0200	0.020	pass	
Thiacloprid	< LOQ	0.010	0.010	pass		Thiamethoxam	< LOQ	0.010	0.010	pass	
Thiophanate-Methyl	< LOQ	0.050	0.030	pass		Trifloxystrobin	< LOQ	0.010	0.010	pass	

Metals										
Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method		Status	Notes	
Arsenic	< LOQ	1.50	mg/kg	0.0979	2208028	09/22/22	AOAC 2013.06 (mod.) ^P	pass		
Cadmium	< LOQ	0.50	mg/kg	0.0979	2208028	09/22/22	AOAC 2013.06 (mod.) ^P	pass		
Lead	< LOQ	0.50	mg/kg	0.0979	2208028	09/22/22	AOAC 2013.06 (mod.) ^P	pass		
Mercury	< LOQ	1.50	mg/kg	0.0489	2208028	09/22/22	AOAC 2013.06 (mod.) ^P	pass		

Mycotoxins										
Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method		Status	Notes	
Aflatoxin B2 ^Y	< LOQ	5.00	µg/kg	5.00	2207994	09/22/22	AOAC 2007.01 & EN 15662 (mod) ^P	pass		
Aflatoxin B1 ^Y	< LOQ	5.00	µg/kg	5.00	2207994	09/22/22	AOAC 2007.01 & EN 15662 (mod) ^P	pass		
Aflatoxin G1 ^Y	< LOQ	5.00	µg/kg	5.00	2207994	09/22/22	AOAC 2007.01 & EN 15662 (mod) ^P	pass		
Aflatoxin G2 ^Y	< LOQ	5.00	µg/kg	5.00	2207994	09/22/22	AOAC 2007.01 & EN 15662 (mod) ^P	pass		
Ochratoxin A ^Y	< LOQ	5.00	µg/kg	5.00	2207994	09/22/22	AOAC 2007.01 & EN 15662 (mod) ^P	pass		
Total Aflatoxins ^Y	0.000		µg/kg	20.0		09/27/22	AOAC 2007.01 & EN 15662 (mod) ^P			



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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.

Threshold Note: Action levels per 6 CCR 1010-21 CDPHE requirements

Ⓟ = ISO/IEC 17025:2017 accredited method.

Ⓢ = TNI accredited analyte.

Units of Measure

/15g = Per 15g

cfu/g = Colony forming units per gram

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

% wt = µg/g divided by 10,000

Glossary of Qualifiers

I: Insufficient sample received to meet method requirements.

Approved Signatory

Derrick Tanner
General Manager



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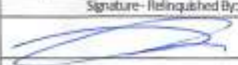
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**Hemp & Cannabis: Usable / Extract / Finished Product
Chain of Custody Record**

Document Control ID: 2832 Revision: 5
Effective: 01/04/2022

ORELAP ID: OR100028 ANAB ISO 17025 ID: AT-1506

Company: Brothers Apothecary Contact: Jesse Richardson Address: 3423 SE 18th Ave City: Portland State: OR Zip Code: 97202 <input type="checkbox"/> Email Results: jesse@brotherstea.com <input type="checkbox"/> Ph: (503) 653.4634 <i>(Billing Contact (if different))</i> Name: _____ Email: _____ Address: _____ City: _____ State: _____ Zip: _____ Ph: (____) _____			Analysis Requested <table border="1"> <tr> <td>H0014 - basic potency</td> <td>H0024 cannabis solvents</td> <td>P2140 cannabis pesticides</td> <td>M264 aspergillus</td> <td>M7000 Colorado Profile</td> <td>H0013 Metals</td> <td>Mycotoxins (H0042)</td> </tr> <tr> <td><input checked="" type="checkbox"/></td> <td><input checked="" type="checkbox"/></td> <td><input checked="" type="checkbox"/></td> <td><input checked="" type="checkbox"/></td> <td><input checked="" type="checkbox"/></td> <td><input checked="" type="checkbox"/></td> <td><input checked="" type="checkbox"/></td> </tr> </table>						H0014 - basic potency	H0024 cannabis solvents	P2140 cannabis pesticides	M264 aspergillus	M7000 Colorado Profile	H0013 Metals	Mycotoxins (H0042)	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	PO Number: _____ Project ID: _____ Batch ID: COL-BR-01 Sampled by: Richardson-Random Custom Reporting: Source Material: <input type="checkbox"/> - Ind. Hemp product <input type="checkbox"/> - Rec. Cannabis Reporting Type: <input type="checkbox"/> - Compliance <input type="checkbox"/> - R&D Report to: <input type="checkbox"/> - METRC <input type="checkbox"/> - CGA <input type="checkbox"/> - USDA <input type="checkbox"/> - Other: Turnaround time (TAT - Business Days): <input checked="" type="checkbox"/> - 5BD <input type="checkbox"/> - 3BD* <input type="checkbox"/> - 2BD* <i>*Check for availability</i>		
H0014 - basic potency	H0024 cannabis solvents	P2140 cannabis pesticides	M264 aspergillus	M7000 Colorado Profile	H0013 Metals	Mycotoxins (H0042)																			
<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>																			
Lab ID	Client Sample Identification	Sample date					Material Type	Weight (Units)	Comments/Metric ID																
	Brothers Hemp Base (Tea, Capsules, Protein)		<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	P																		
	Hemp Isolate		<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	I																		
	Hemp Distillate		<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	<input checked="" type="checkbox"/>	C																		
Signature - Relinquished By: 			Date: 9/19/22	Time: 3:08 pm	Signature - Received By: JF	Date: 9/19	Time: 15:07	Lab Use Only: <input type="checkbox"/> Shipped Via: _____ or <input checked="" type="checkbox"/> Client drop off Evidence of cooling: <input type="checkbox"/> Yes <input type="checkbox"/> No - Temp (°C): 29.0 Sample in good condition: <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No Payment: <input type="checkbox"/> Cash <input type="checkbox"/> Check <input type="checkbox"/> CC <input type="checkbox"/> Net: Prelog storage: _____																	

- Material Type Codes: Plant Material [] ; Isolate [] ; Concentrate/Extract [] ; Tincture/Topical [] ; Edible [] ; Beverage [] ; Vapor Product []
 Samples submitted to Columbia Laboratories with testing requirements constitute an agreement for services in accordance with the _____ associated with this COC. By signing "Relinquished By" you are agreeing to these terms.
 12423 NE Whitaker Way P: (503) 254-1794 | Fax: (503) 254-1452 Page _____ of _____
 Portland, OR 97230



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Revision: 1 Document ID: 7148
Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

J AOAC 2015 V98-6 Batch ID: 2207941

Laboratory Control Sample									
Analyte	LCS	Result	Spike	Units	% Rec	Limits		Evaluation	Notes
CBDVA	2	0.106	0.100	%	106	80.0	- 120	Acceptable	
CBDV	2	0.112	0.100	%	112	80.0	- 120	Acceptable	
CBE	2	0.106	0.100	%	106	80.0	- 120	Acceptable	
CBDA	1	0.105	0.100	%	105	90.0	- 110	Acceptable	
CBGA	1	0.104	0.0990	%	105	80.0	- 120	Acceptable	
CBG	1	0.116	0.108	%	107	80.0	- 120	Acceptable	
CBD	1	0.116	0.108	%	108	90.0	- 110	Acceptable	
THCV	2	0.107	0.100	%	107	80.0	- 120	Acceptable	
d8THCV	2	0.108	0.100	%	108	80.0	- 120	Acceptable	
THCVA	2	0.107	0.100	%	107	80.0	- 120	Acceptable	
CBN	1	0.114	0.108	%	105	90.0	- 110	Acceptable	
exo-THC	2	0.103	0.100	%	103	80.0	- 120	Acceptable	
d9THC	1	0.118	0.116	%	102	90.0	- 110	Acceptable	
d8THC	1	0.107	0.100	%	107	80.0	- 120	Acceptable	
CBL	2	0.101	0.100	%	101	80.0	- 120	Acceptable	
d10THC	1	0.0967	0.100	%	96.7	80.0	- 120	Acceptable	
CBC	2	0.105	0.100	%	105	80.0	- 120	Acceptable	
THCA	1	0.0962	0.0955	%	101	90.0	- 110	Acceptable	
CBCA	2	0.104	0.100	%	104	80.0	- 120	Acceptable	
CBLA	2	0.108	0.100	%	108	80.0	- 120	Acceptable	
CBT	2	0.0922	0.100	%	92.2	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			
CBL	<LOQ	0.077	%	< 0.077		Acceptable			
d10THC	<LOQ	0.077	%	< 0.077		Acceptable			
CBC	<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			
CBT	<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: 2207941						
Sample Duplicate		Sample ID: 22-011064-0001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBDV	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBE	0.143	0.159	0.077	%	11.0	< 20	Acceptable	
CBD	3.31	3.45	0.077	%	4.14	< 20	Acceptable	
CBDV	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBDV	0.138	0.150	0.077	%	8.43	< 20	Acceptable	
CBD	2.38	2.51	0.077	%	5.22	< 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	0.275	0.287	0.077	%	4.54	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
d9THC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
d8THC	0.493	0.505	0.077	%	2.29	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
d10THC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBC	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
THCA	0.283	0.281	0.077	%	0.502	< 20	Acceptable	
CBCA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.077	%	NA	< 20	Acceptable	
CBT	0.485	0.476	0.077	%	1.94	< 20	Acceptable	

Abbreviations

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

Units of Measure:



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Laboratory Quality Control Results

Organic Acids				Batch ID: 2207970						
Method Blank				Laboratory Control Sample						
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes	
Acetic Acid	<LOQ	<	250	479	522	µg/g	91.9	70 - 130		
Formic Acid	<LOQ	<	250	642	617	µg/g	104.1	70 - 130		

QC - Sample Duplicate						Sample ID: LCS		
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Accept/Fail	Notes
Acetic Acid	<LOQ	<LOQ	250	µg/g	0.0	< 20	Acceptable	
Formic Acid	<LOQ	<LOQ	250	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

Residual Solvents				Batch ID: 2208097					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		442	572	µg/g	77.3	60 - 120	
Isobutane	ND	< 200		555	731	µg/g	75.9	60 - 120	
Butane	ND	< 200		540	731	µg/g	73.9	60 - 120	
2,2-Dimethylpropane	ND	< 200		730	936	µg/g	78.0	60 - 120	
Methanol	ND	< 200		1360	1650	µg/g	82.4	60 - 120	
Ethylene Oxide	ND	< 30		41.7	56.2	µg/g	74.2	60 - 120	
2-Methylbutane	ND	< 200		1400	1650	µg/g	84.8	60 - 120	
Pentane	ND	< 200		1410	1650	µg/g	85.5	60 - 120	
Ethanol	ND	< 200		1340	1660	µg/g	80.7	70 - 130	
Ethyl Ether	ND	< 200		1360	1630	µg/g	83.4	60 - 120	
2,2-Dimethylbutane	ND	< 30		160	189	µg/g	84.7	60 - 120	
Acetone	ND	< 200		1370	1650	µg/g	83.0	60 - 120	
2-Propanol	ND	< 200		1320	1650	µg/g	80.0	60 - 120	
Ethyl Formate	ND	< 500		1540	1610	µg/g	95.7	70 - 130	
Acetonitrile	ND	< 100		390	504	µg/g	77.4	60 - 120	
Methyl Acetate	ND	< 500		1700	1630	µg/g	104.3	70 - 130	
2,3-Dimethylbutane	ND	< 30		127	174	µg/g	73.0	60 - 120	
Dichloromethane	ND	< 60		422	521	µg/g	81.0	60 - 120	
2-Methylpentane	ND	< 30		164	187	µg/g	87.7	60 - 120	
MTBE	ND	< 500		1650	1600	µg/g	103.1	70 - 130	
3-Methylpentane	ND	< 30		161	188	µg/g	85.6	60 - 120	
Hexane	ND	< 30		156	182	µg/g	85.7	60 - 120	
1-Propanol	ND	< 500		1710	1610	µg/g	106.2	70 - 130	
Methylethylketone	ND	< 500		1670	1600	µg/g	104.4	70 - 130	
Ethyl acetate	ND	< 200		1300	1630	µg/g	79.8	60 - 120	
2-Butanol	ND	< 200		1250	1630	µg/g	76.7	60 - 120	
Tetrahydrofuran	ND	< 100		398	506	µg/g	78.7	60 - 120	
Cyclohexane	ND	< 200		1300	1640	µg/g	79.3	60 - 120	
2-methyl-1-propanol	ND	< 500		1640	1620	µg/g	101.2	70 - 130	
Benzene	ND	< 1		3.67	4.93	µg/g	74.4	60 - 120	
Isopropyl Acetate	ND	< 200		1260	1640	µg/g	76.8	60 - 120	
Heptane	ND	< 200		1280	1630	µg/g	78.5	60 - 120	
1-Butanol	ND	< 500		1730	1600	µg/g	108.1	70 - 130	
Propyl Acetate	ND	< 500		1680	1620	µg/g	103.7	70 - 130	
1,4-Dioxane	ND	< 100		351	493	µg/g	71.2	60 - 120	
2-Ethoxyethanol	ND	< 30		127	171	µg/g	74.3	60 - 120	
Methylisobutylketone	ND	< 500		1720	1620	µg/g	106.2	70 - 130	
3-Methyl-1-butanol	ND	< 500		1720	1610	µg/g	106.8	70 - 130	
Ethylene Glycol	ND	< 200		289	494	µg/g	58.5	60 - 120	Q6
Toluene	ND	< 100		349	506	µg/g	69.0	60 - 120	
Isobutyl Acetate	ND	< 500		1840	1620	µg/g	113.6	70 - 130	
1-Pentanol	ND	< 500		2050	1610	µg/g	127.3	70 - 130	
Butyl Acetate	ND	< 500		1720	1610	µg/g	106.8	70 - 130	
Ethylbenzene	ND	< 200		702	996	µg/g	70.5	60 - 120	
m,p-Xylene	ND	< 200		716	1010	µg/g	70.9	60 - 120	
o-Xylene	ND	< 200		702	979	µg/g	71.7	60 - 120	
Cumene	ND	< 30		139	188	µg/g	73.9	60 - 120	
Anisole	ND	< 500		1660	1610	µg/g	103.1	70 - 130	
DMSO	ND	< 500		1670	1600	µg/g	104.4	70 - 130	
1,2-dimethoxyethane	ND	< 50		199	190	µg/g	104.7	70 - 130	
Triethylamine	ND	< 500		1650	1610	µg/g	102.5	70 - 130	
N,N-dimethylformamide	ND	< 150		520	496	µg/g	104.8	70 - 130	
N,N-dimethylacetamide	ND	< 150		506	483	µg/g	104.8	70 - 130	
Pyridine	ND	< 50		175	167	µg/g	104.8	70 - 130	
1,2-Dichloroethane	ND	< 1		1.13	1	µg/g	113.0	70 - 130	
Chloroform	ND	< 1		1.15	1	µg/g	115.0	70 - 130	
Trichloroethylene	ND	< 1		1.19	1	µg/g	119.0	70 - 130	



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QC - Sample Duplicate			Sample ID: 22-011013-0001					
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	
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	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	
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	

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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Report Number: 22-011232/D002.R000
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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	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.