



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



Report Number: 23-005965/D002.R000
Report Date: 05/25/2023
ORELAP#: OR100028
Purchase Order:
Received: 05/18/23 11:01

Customer: NW Natural Goods
Product identity: HEMP - PCH 0007
Client/Metric ID: .
Laboratory ID: 23-005965-0001

Summary

Potency:

Analyte per 4g	Result	Limits	Units	Status	
CBC per 4g	10.0		mg/4g		CBD-Total per Serving Size 21.4 mg/4g
CBD per 4g	21.4		mg/4g		
CBDV per 4g	0.124		mg/4g		THC-Total per Serving Size <LOQ
CBG per 4g	0.616		mg/4g		(Reported in milligrams per serving)
CBT per 4g	0.159		mg/4g		

Residual Solvents:

All analytes passing and less than LOQ.

Pesticides:

Analyte	Result (mg/kg)	Limits (mg/kg)	Status
Multi-Residue Pesticide Profile	< LOQ for all analytes		

Metals:

Less than LOQ for all analytes.

Microbiology:

Less than LOQ for all analytes.



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Customer: NW Natural Goods

Product identity: HEMP - PCH 0007

Client/Metric ID: .

Sample Date:

Laboratory ID: 23-005965-0001

Evidence of Cooling: No

Temp: 18.6 °C

Relinquished by: Ramos

Serving Size #1: 4 g

Sample Results

Potency per 4g		Method: J AOAC 2015 V98-6 (mod) ^b		Units mg/se Batch: 2307525		Analyze: 5/20/23 4:31:00 AM
Analyte	Result	Limits	Units	LOQ	Notes	
CBC per 4g	10.0		mg/4g	0.124		
CBC-A per 4g	< LOQ		mg/4g	0.124		
CBC-Total per 4g	10.0		mg/4g	0.232		
CBD per 4g	21.4		mg/4g	0.124		
CBD-A per 4g	< LOQ		mg/4g	0.124		
CBD-Total per 4g	21.4		mg/4g	0.232		
CBDV per 4g	0.124		mg/4g	0.124		
CBDV-A per 4g	< LOQ		mg/4g	0.124		
CBDV-Total per 4g	< LOQ		mg/4g	0.231		
CBE per 4g	< LOQ		mg/4g	0.124		
CBG per 4g	0.616		mg/4g	0.124		
CBG-A per 4g	< LOQ		mg/4g	0.124		
CBG-Total per 4g	0.616		mg/4g	0.231		
CBL per 4g	< LOQ		mg/4g	0.124		
CBL-A per 4g	< LOQ		mg/4g	0.124		
CBL-Total per 4g	< LOQ		mg/4g	0.232		
CBN per 4g	< LOQ		mg/4g	0.124		
CBT per 4g	0.159		mg/4g	0.124		
Δ8-THCV per 4g	< LOQ		mg/4g	0.124		
Δ10-THC-9R per 4g	< LOQ		mg/4g	0.124		
Δ10-THC-9S per 4g	< LOQ		mg/4g	0.124		
Δ10-THC-Total per 4g	< LOQ		mg/4g	0.248		
Δ8-THC per 4g	< LOQ		mg/4g	0.124		
Δ9-THC per 4g	< LOQ		mg/4g	0.124		
delta-9-THCP per 4g	< LOQ		mg/4g	0.124		
exo-THC per 4g	< LOQ		mg/4g	0.124		
THC-A per 4g	< LOQ		mg/4g	0.124		
THC-Total per 4g	< LOQ		mg/4g	0.232		
THCV per 4g	< LOQ		mg/4g	0.124		
THCV-A per 4g	< LOQ		mg/4g	0.124		
THCV-Total per 4g	< LOQ		mg/4g	0.232		
Total Cannabinoids per 4g	32.3		mg/4g			



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Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2307452	05/21/23 AOAC 991.14 (Petrifilm) ^P		
Total Coliforms	< LOQ		cfu/g	10	2307452	05/21/23 AOAC 991.14 (Petrifilm) ^P		
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2307453	05/21/23 AOAC 2014.05 (RAPID) ^P		
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2307453	05/21/23 AOAC 2014.05 (RAPID) ^P		

Solvents Method: Residual Solvents by GC/MS^P Units µg/g Batch 2307514 Analyze 05/22/23 11:29 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-Dimethyl butane	< LOQ		30.0			2,2-Dimethylpropane (neo-pentane)	< LOQ		200		
2,3-Dimethyl butane	< 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	
Isopropyl benzene (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)^P Units mg/kg Batch 2307527 Analyze 05/23/23 12:00 PM

Analyte	Result	Limits	Status	Notes
Multi-Residue Pesticide Profile	< LOQ for all analytes			



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Metals

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
Arsenic*	< LOQ	0.200	mg/kg	0.0163	2307620	05/24/23 AOAC 2013.06 (mod.) ^p	pass	
Cadmium*	< LOQ	0.200	mg/kg	0.0163	2307620	05/24/23 AOAC 2013.06 (mod.) ^p	pass	
Lead*	< LOQ	0.500	mg/kg	0.0163	2307620	05/24/23 AOAC 2013.06 (mod.) ^p	pass	
Mercury*	< LOQ	0.100	mg/kg	0.00815	2307620	05/24/23 AOAC 2013.06 (mod.) ^p	pass	

Nutrition

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
Moisture (Loss on Drying)	18.4		g/100g	0.10	2307539	05/23/23 AOAC 925.10 (mod.) ^p		
Water Activity	0.675		Aw	0.030	2307473	05/19/23 AOAC 978.18 ^p		



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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.

^p = ISO/IEC 17025:2017 accredited method.

[¥] = TNI accredited analyte.

Units of Measure

cfu/g = Colony forming units per gram

g = g

g/100g = Grams per 100 Grams

µg/g = Microgram per gram

mg/kg = Milligram per kilogram = parts per million (ppm)

mg/4g = Milligram per 4g

% = Percentage of sample

Aw = Water Activity

% wt = µg/g divided by 10,000

Approved Signatory

Derrick Tanner
General Manager



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Compound	LOQ (mg/kg)	Compound	LOQ (mg/kg)	Compound	LOQ (mg/kg)
Abamec in	0.100	Cle hodim	0.050	ndrin	0.100
Acepha e	0.100	Cle hodim Sul one	0.050	PN	0.050
Acequinocyl	0.100	Cle hodim Sul oxide	0.050	PTC	0.100
Ace amiprid	0.020	Clo en ezine	0.020	s envalera e/ envalera e	0.200
Ace ochlor	0.100	Clomazone	0.020	aconazole	0.100
Acrina hrin	0.100	Clo hianidin	0.200	hal luralin	0.100
Alachlor	0.100	Coumaphos	0.050	hio encarb	0.050
Aldicarb	0.100	Cro oxyphos	0.020	hion	0.200
Aldicarb sul oxide	0.100	Cyanazine	0.020	hirimol	0.100
Aldoxycarb (Aldicarb-sul one)	0.100	Cyano enphos	0.020	ho umesa e	0.050
Aldrin	0.100	Cyan raniliprole	0.050	hoprophos	0.020
Ame ocr radin	0.020	Cyazo amid	0.020	o enprox	0.020
Ame ryn	0.500	Cycloa e	0.100	oxazole	0.020
Aspon	0.100	Cy lu hrin	0.200	ridiazole	0.100
Asulam	0.100	Cyhalo hrin, lambda	0.200	rim os	0.020
A razine	0.100	Cymoxanil	0.050	amoxadone	0.200
A razine-dese hyl	0.100	Cyperme hrin	0.200	amphur	0.100
Azinphos-e hyl	0.020	Cyprodinil	0.100	enamidone	0.020
Azinphos-me hyl	0.020	Dac hal	0.100	enamiphos	0.020
Azoxys robin	0.020	Daminozide	0.100	enamiphos sul one	0.020
Benalaxyl	0.020	DCPMU	0.050	enamiphos sul oxide	0.020
Bendiocarb	0.020	DDD, o,p'-	0.100	enazaquin	0.100
Ben luralin	0.100	DDD, p,p'-	0.100	enbuconazole	0.100
Benoxacor	0.050	DD , o,p'-	0.100	enchlorphos	0.100
Bensulide	0.050	DD , p,p'-	0.100	enchlorphos-oxon	0.100
B C alpha isomer	0.100	DDT, o,p'-	0.100	enhexamid	0.100
B C be a isomer	0.100	DDT, p,p'-	0.100	eni ro hion	0.100
B C del a isomer	0.500	D (Tribu os)	0.100	enobucarb	0.050
Bi enaza e	0.020	Del ame hrin	0.100	enoxycarb	0.020
Bi en hrin	0.020	Desmedipham	0.100	enpropa hrin	0.050
Boscalid	0.020	Dialla e	0.100	enpyroxima e	0.020
Bromophos-e hyl	0.100	Diazinon	0.020	enson	0.100
Bromophos-me hyl	0.200	Diazoxon	0.100	ensul o hion	0.020
Bromopropyla e	0.100	Dichlobenil	0.100	ensul o hion oxon	0.020
Bromuconazole	0.100	Dichlo luanid	0.100	ensul o hion sul one	0.100
Bupirima e	0.020	Dichlorvos	0.100	Fensulfothion-oxon-sulfone	0.020
Bupro ezin	0.050	Diclobu razol	0.050	en hion	0.050
Bu achlor	0.500	Dico ol	0.100	en hion oxon	0.020
Bu ralin	0.200	Dicro ophos	0.050	en hion oxon sul one	0.100
Bu yla e	0.100	Dieldrin	0.100	en hion sul one	0.050
Cadusa os	0.020	Die ho encarb	0.020	enuron	0.020
Cap an	1.000	Die hyl oluamide (D T)	0.050	ipronil	0.100
Carbaryl	0.050	Di enoconazole	0.100	lonicamid	0.100
Carbendazim	0.100	Dime henamid	0.050	luchloralin	0.100
Carbo uran	0.020	Dime hoa e	0.050	lucy hrina e	0.100
Carbopheno hion	0.200	Dime homorph	0.050	ludioxonil	0.200
Carboxin	0.020	Diniconazole	0.200	lu enace	0.020
Car en razone-e hyl	0.100	Dino e uran	0.200	lumioxazin	0.100
Chloran raniliprole	0.020	Dioxa hion	0.100	luome uron	0.020
Chlordane, cis-	0.200	Diphenamid	0.020	luopicolide	0.050
Chlordane, rans-	0.200	Diphenylamine	0.100	luopyram	0.020
Chlor enapyr	0.500	Disul o on	0.100	luoxas robin	0.050
Chlor enson	0.200	Disul o on sul one	0.100	lupyradi urone	0.020
Chlor envinphos	0.050	Disul o on sul oxide	0.100	luridone	0.100
Chlorobenzila e	0.100	Diuron	0.050	lusilazole	0.020
Chloroneb	0.200	di enphos	0.050	lu olanil	0.020
Chlorpyri os	0.050	ndosul an alpha	0.200	lu ria ol	0.020
Chlorpyri os-me hyl	0.200	ndosul an be a	0.200	luvalina e, au-	0.100
C PC	1.000	ndosul an sul a e	0.100	luxapyroxad	0.020



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Compound	LOQ (mg/kg)	Compound	LOQ (mg/kg)	Compound	LOQ (mg/kg)
omosa en	0.100	Mexacarba e	0.020	Propamocarb	0.050
ono os	0.100	MGK 264	0.020	Propanil	0.050
orchlor enuron	0.050	Mirex	0.100	Propargi e	0.050
orme ana e	0.050	Molina e	0.050	Propazine	0.020
ura hiocarb	0.020	Monocro ophos	0.100	Prope amphos	0.050
ep achlor	0.100	Monolinuron	0.020	Propham	0.050
ep achlor epoxide	0.100	Myclobu anil	0.050	Propiconazole	0.050
ep enophos	0.100	Naled	0.100	Propoxur	0.050
exachlorobenzene	0.100	Napropamide	0.050	Propoxycarbazona Na	0.050
exaconazole	0.100	Neburon	0.020	Propyzamide	0.050
exazinone	0.100	Ni rapyrin	0.100	Pro hio os	0.100
exy hiazox	0.020	Nor lurazon	0.050	Pyraclos robin	0.020
mazalil	0.100	Ome hoa e	0.100	Pyrazophos	0.050
midacloprid	0.100	O-Phenylphenol	0.100	Pyre hrins	0.050
ndazi lam	0.020	Oxadixyl	0.100	Pyridaben	0.020
ndoxacarb	0.020	Oxamyl	0.100	Pyrida ol	0.100
proben os	0.100	Oxamyl-oxime	0.100	Pyrida e	0.020
prodione	0.100	Oxychlorthane	0.100	Pyrima hanil	0.050
sobenzan	0.100	Oxydeme on-Me hyl	0.100	Pyriproxi en	0.020
socarbophos	0.500	Oxy hioquinox	0.200	Pyroxasul one	0.020
sodrin	0.100	Paclobu razol	0.050	Pyroxulam	0.020
so enphos	0.050	Paraaxon-e hyl	0.020	Quinalphos	0.050
so enphos-me hyl	0.020	Paraaxon me hyl	0.100	Quinoxy en	0.050
so enphos oxon	0.050	Para hion e hyl	0.100	Quin ozene (PCNB)	0.200
soproc carb	0.020	Para hion me hyl	0.200	Resme hrin	0.050
sopropalin	0.200	Penconazole	0.050	Ro enone	0.050
sopro hiolane	0.050	Pendime halin	0.050	S421	0.100
sopro uron	0.050	Pen lu en	0.020	Simazine	0.100
soxaben	0.050	Pen achloroaniline	0.100	Sime ryn	0.200
soxa lu ole	0.050	Pen achloroanisole	0.100	Spine oram	0.020
Kresoxim-me hyl	0.050	Pen achlorobenzene (PCB)	0.100	Spinosad	0.050
ac o en	0.500	Pentachlorothioanisole (PCTA)	0.100	Spirodiclo en	0.100
enacil	0.100	Pen hiopyrad	0.020	Spiromesi en	0.050
indane (gamma B C)	0.100	Perme hrin	0.050	Spiro e rama	0.050
inuron	0.020	Per hane	0.100	Spiroxamine	0.020
Malaaxon	0.050	Phenmedipham	0.050	Sul o ep	0.050
Mala hion	0.050	Phen hoa e	0.050	Sul oxa lor	0.050
Mandipropamid	0.020	Phora e	0.050	Sulpro os	0.020
Mecarbam	0.020	Phora e Sul one	0.050	Tebuconazole	0.100
Mepanipyrim	0.050	Phora e Sul oxide	0.050	Tebu enozide	0.020
Merphos	0.500	Phosalone	0.050	Tebu hiuron	0.020
Me alaxyl	0.050	Phosme	0.100	Tecnazene	0.100
Me aldehyde	0.050	Phosphamidon	0.050	Te lu hrin	0.100
Me conazole	0.100	Phoxim	0.050	Terbu os	0.020
Me hacri os	0.100	Pinoxaden	0.020	Terbu os sul one	0.050
Me hamidophos	0.050	Piperonyl bu oxide	0.050	Terbu os sul oxide	0.050
Me hida hion	0.050	Pirimicarb	0.020	Terbu hylazine	0.020
Me hiocarb	0.050	Pirimiphos-me hyl	0.050	Terbu ryn	0.020
Me hiocarb sul one	0.100	Pirimiphos-e hyl	0.020	Te rachlorvinphos	0.050
Me hiocarb sul oxide	0.100	Pralle hrin	0.100	Te raconazole	0.050
Me homyl	0.100	Prochloraz	0.020	Te radi on	0.200
Me hoxychlor	0.100	Procymidone	0.100	Te rame hrin	0.050
Me hoxy enozide	0.020	Pro eno os	0.100	Te rasul	0.100
Me obromuron	0.050	Pro luralin	0.100	Thiabendazole	0.100
Me olachlor	0.100	Promecarb	0.050	Thiabendazole, 5-hydroxy	0.100
Me olcarb	0.050	Prome on	0.100	Thiacloprid	0.050
Me ra enone	0.050	Prome ryn	0.020	Thiame hoxam	0.100
Me ribuzin	0.100	Propachlor	0.020	Thiobencarb	0.050
Mevinphos	0.100			Thiodicarb	0.050
				Thiophana e-me hyl	0.050



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Compound	LOQ (mg/kg)	Compound	LOQ (mg/kg)	Compound	LOQ (mg/kg)
Tolclo os-me hyl	0.100	Triazophos	0.020	Tri loxys robin	0.020
Tri orin	0.100	Tolyl luanid	0.050	Tri iconazole	0.050
Tralkoxydim	0.100	Tridiphane	0.500	Vinclozolin	0.100
Triadime on	0.050	Tri lumizole	0.020	Zoxamide	0.020
Trialla e	0.100	Tri luralin	0.100		

LOQ= Limit of Quantitation, mg/kg

Factors affecting the LOQ include instrument sensitivity or a particular analyte, sample size, moisture content (percent solids) of the sample, effectiveness of the cleanup on the sample extract, and especially the type of sample matrix.



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

Northeast-Natural-Goods-1684355179

OREIAP ID OR1000028 ANAB ISO 17025 IDAT1508

Project Information Project Name <u>HEMP - PCH 0007</u> PO Number <u>N/A</u> Turnaround Time <u>5 Business Days (standard) (required for microbial testing)</u> Samples Delivered to Laboratory <u>Schedule Pick-Up</u> Cannabis Type <u>Industrial</u> Pick-Up Location Street Address <u>11781 SE HWY 212 #404</u> City, State, Zip <u>Clackamas, Oregon 97015</u> Pick-Up Location Phone <u>3866699210</u>			Testing						
			Heavy Metals Profile OR (As, Cd, Pb & Hg)	Moisture as Loss on Drying	Pesticide - Multi-Residue Profile	Potency Cannabinoid Basic + Extended Profile	Residual Solvents - OR	Total Coliforms + E. Coli	Water Activity
#	Sample Name	Sample Material	Amount Provided						
1	HEMP - PCH 0007	Edible	20 units for sale	✓	✓	✓	✓	✓	✓

Relinquished By	Date	Time	Temp., °C	Received By	Date	Time	Received Temp., °C	Evidence of Cooling?
<i>Kristen Johnson</i>	<i>5/17/2023</i>	<i>13:26</i>		<i>BR</i>	<i>5/19/2023</i>	<i>10:27</i>		<i>No</i>
<i>BR</i>	<i>5/18/2023</i>	<i>10:51</i>	<i>18.6</i>	<i>MRH</i>	<i>5/18/2023</i>	<i>11:01</i>		<i>No</i>

Samples submitted to Columbia Laboratories with testing requirements constitute an agreement for services in accordance with the [current terms of services](#) associated with this COC. By signing "Relinquished by" you are agreeing to these terms.

Columbia Laboratories
 12423 NE Whitaker Way
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P (503) 254-1794 | Fx (503) 254-1452
info@columbiaboratories.com

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

Laboratory Quality Control Results

Residual Solvents				Batch D: 2307514					
Method Blank				Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec	Limits	Notes
Propane	ND	< 200		532	584	µg/g	91.1	60 - 120	
Isobutane	ND	< 200		671	767	µg/g	87.5	60 - 120	
Butane	ND	< 200		676	782	µg/g	86.3	60 - 120	
2,2-Dimethylpropane	ND	< 200		941	939	µg/g	100.2	60 - 120	
Methanol	ND	< 200		1340	1610	µg/g	83.2	60 - 120	
Ethylene Oxide	ND	< 30		52	57.1	µg/g	91.1	60 - 120	
2-Methylbutane	ND	< 200		1230	1600	µg/g	76.9	60 - 120	
Pentane	ND	< 200		1240	1610	µg/g	77.0	60 - 120	
Ethanol	ND	< 200		1260	1600	µg/g	78.8	70 - 130	
Ethyl Ether	ND	< 200		1260	1610	µg/g	78.3	60 - 120	
2,2-Dimethylbutane	ND	< 30		134	173	µg/g	77.5	60 - 120	
Acetone	ND	< 200		1280	1620	µg/g	79.0	60 - 120	
2-Propanol	ND	< 200		1250	1600	µg/g	78.1	60 - 120	
Ethyl Formate	ND	< 500		1510	1610	µg/g	93.8	70 - 130	
Acetonitrile	ND	< 100		380	488	µg/g	77.9	60 - 120	
Methyl Acetate	ND	< 500		1370	1610	µg/g	85.1	70 - 130	
2,3-Dimethylbutane	ND	< 30		127	165	µg/g	77.0	60 - 120	
Dichloromethane	ND	< 60		382	487	µg/g	78.4	60 - 120	
2-Methylpentane	ND	< 30		121	160	µg/g	75.6	60 - 120	
MTBE	ND	< 500		1420	1600	µg/g	88.8	70 - 130	
3-Methylpentane	ND	< 30		126	161	µg/g	78.3	60 - 120	
Hexane	ND	< 30		126	162	µg/g	77.8	60 - 120	
1-Propanol	ND	< 500		1470	1620	µg/g	90.7	70 - 130	
Methyl ethyl ketone	ND	< 500		1440	1610	µg/g	89.4	70 - 130	
Ethyl acetate	ND	< 200		1240	1600	µg/g	77.5	60 - 120	
2-Butanol	ND	< 200		1250	1610	µg/g	77.6	60 - 120	
Tetrahydrofuran	ND	< 100		377	483	µg/g	78.1	60 - 120	
Cyclohexane	ND	< 200		1270	1610	µg/g	78.9	60 - 120	
2-methyl-1-propanol	ND	< 500		1480	1630	µg/g	90.8	70 - 130	
Benzene	ND	< 1		3.85	4.98	µg/g	77.3	60 - 120	
Isopropyl Acetate	ND	< 200		1240	1610	µg/g	77.0	60 - 120	
Heptane	ND	< 200		1280	1620	µg/g	77.8	60 - 120	
1-Butanol	ND	< 500		1460	1600	µg/g	91.3	70 - 130	
Propyl Acetate	ND	< 500		1460	1620	µg/g	90.1	70 - 130	
1,4-Dioxane	ND	< 100		383	494	µg/g	77.5	60 - 120	
2-Ethoxyethanol	ND	< 30		135	165	µg/g	81.8	60 - 120	
Methylisobutylketone	ND	< 500		1450	1610	µg/g	90.1	70 - 130	
3-Methyl-1-butanol	ND	< 500		1380	1610	µg/g	85.3	70 - 130	
Ethylene Glycol	ND	< 200		310	488	µg/g	63.8	60 - 120	
Toluene	ND	< 100		374	513	µg/g	72.9	60 - 120	
Isobutyl Acetate	ND	< 500		1430	1600	µg/g	89.4	70 - 130	
1-Pentanol	ND	< 500		1400	1610	µg/g	87.0	70 - 130	
Butyl Acetate	ND	< 500		1450	1610	µg/g	90.1	70 - 130	
Ethylbenzene	ND	< 200		742	967	µg/g	76.7	60 - 120	
m,p-Xylene	ND	< 200		798	994	µg/g	80.3	60 - 120	
o-Xylene	ND	< 200		751	992	µg/g	75.7	60 - 120	
Cumene	ND	< 30		132	171	µg/g	77.2	60 - 120	
Anisole	ND	< 500		1420	1610	µg/g	88.2	70 - 130	
DMSO	ND	< 500		1180	1610	µg/g	73.3	70 - 130	
1,2-dimethoxyethane	ND	< 50		157	172	µg/g	91.3	70 - 130	
Triethylamine	ND	< 500		1510	1620	µg/g	93.2	70 - 130	
N,N-dimethylformamide	ND	< 150		440	499	µg/g	88.2	70 - 130	
N,N-dimethylacetamide	ND	< 150		367	491	µg/g	74.7	70 - 130	
Pyridine	ND	< 50		153	171	µg/g	89.5	70 - 130	
Silolane	ND	< 50		145	160	µg/g	90.6	70 - 130	
1,2-Dichloroethane	ND	< 1		0.957	1	µg/g	95.7	70 - 130	
Chloroform	ND	< 1		0.978	1	µg/g	97.8	70 - 130	
Trichloroethylene	ND	< 1		0.903	1	µg/g	90.3	70 - 130	
1,1-Dichloroethane	ND	< 1		0.963	1	µg/g	96.3	70 - 130	



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Report Number: 23-005965/D002.R000
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Received: 05/18/23 11:01

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

QC- Sample Duplicate		Sample ID: 23-005910-0001						
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

Units of Measure:

µg/g- Microgram per gram or ppm



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 Portland, OR 97230
 503-254-1794



Report Number: 23-005965/D002.R000
Report Date: 05/25/2023
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Purchase Order:
Received: 05/18/23 11:01

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

Laboratory Quality Control Results

JAOAC2015 V986 Batch ID: 2307525

Laboratory Control Sample									
Analyte	LCS	Result	Spike	Units	% Rec	Limits		Evaluation	Notes
CBDA	2	0.0293	0.0283	%	104	80.0	- 120	Acceptable	
CBV	2	0.0307	0.0291	%	105	80.0	- 120	Acceptable	
CBE	2	0.0358	0.0344	%	104	80.0	- 120	Acceptable	
CBDA	1	0.0341	0.0311	%	109	90.0	- 110	Acceptable	
CBSA	1	0.0338	0.0311	%	109	80.0	- 120	Acceptable	
CBS	1	0.0351	0.0322	%	109	80.0	- 120	Acceptable	
CB	1	0.0351	0.0323	%	109	90.0	- 110	Acceptable	
THCV	2	0.0212	0.0201	%	105	80.0	- 120	Acceptable	
δ8THCV	2	0.0280	0.0268	%	104	80.0	- 120	Acceptable	
THCVA	2	0.0311	0.0299	%	104	80.0	- 120	Acceptable	
CBN	1	0.0359	0.0329	%	109	80.0	- 120	Acceptable	
exo-THC	2	0.0303	0.0292	%	104	80.0	- 120	Acceptable	
δ9THC	1	0.0360	0.0341	%	105	90.0	- 110	Acceptable	
δ8THC	1	0.0444	0.0420	%	106	90.0	- 110	Acceptable	
9SaTHC	1	0.0257	0.0240	%	107	80.0	- 120	Acceptable	
CB	2	0.0329	0.0315	%	105	80.0	- 120	Acceptable	
9RaTHC	1	0.0337	0.0310	%	109	80.0	- 120	Acceptable	
CB	2	0.0320	0.0309	%	104	80.0	- 120	Acceptable	
THCA	1	0.0344	0.0314	%	110	90.0	- 110	Acceptable	
CBA	2	0.0333	0.0326	%	102	80.0	- 120	Acceptable	
CBA	2	0.0340	0.0331	%	103	80.0	- 120	Acceptable	
δ9THCP	2	0.0329	0.0321	%	102	80.0	- 120	Acceptable	
CB	2	0.0334	0.0327	%	102	80.0	- 120	Acceptable	

Method Blank							
Analyte	Result	LOQ	Units	Limits		Evaluation	Notes
CBDA	<LOQ	0.00319	%	< 0.00319		Acceptable	
CBV	<LOQ	0.00319	%	< 0.00319		Acceptable	
CBE	<LOQ	0.00319	%	< 0.00319		Acceptable	
CBDA	<LOQ	0.00319	%	< 0.00319		Acceptable	
CBSA	<LOQ	0.00319	%	< 0.00319		Acceptable	
CBS	<LOQ	0.00319	%	< 0.00319		Acceptable	
CB	<LOQ	0.00319	%	< 0.00319		Acceptable	
THCV	<LOQ	0.00319	%	< 0.00319		Acceptable	
δ8THCV	<LOQ	0.00319	%	< 0.00319		Acceptable	
THCVA	<LOQ	0.00319	%	< 0.00319		Acceptable	
CBN	<LOQ	0.00319	%	< 0.00319		Acceptable	
exo-THC	<LOQ	0.00319	%	< 0.00319		Acceptable	
δ9THC	<LOQ	0.00319	%	< 0.00319		Acceptable	
δ8THC	<LOQ	0.00319	%	< 0.00319		Acceptable	
9SaTHC	<LOQ	0.00319	%	< 0.00319		Acceptable	
CB	<LOQ	0.00319	%	< 0.00319		Acceptable	
9RaTHC	<LOQ	0.00319	%	< 0.00319		Acceptable	
CB	<LOQ	0.00319	%	< 0.00319		Acceptable	
THCA	<LOQ	0.00319	%	< 0.00319		Acceptable	
CBA	<LOQ	0.00319	%	< 0.00319		Acceptable	
CBA	<LOQ	0.00319	%	< 0.00319		Acceptable	
δ9THCP	<LOQ	0.00319	%	< 0.00319		Acceptable	
CB	<LOQ	0.00319	%	< 0.00319		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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Report Number: 23-005965/D002.R000
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ORELAP#: OR100028
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Received: 05/18/23 11:01

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

Laboratory Quality Control Results

JAOAC2015 V986		Batch ID: 2307525						
Sample Duplicate		Sample ID: 23-0059570001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
CBDV	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
CBE	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
CBD	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
CBSA	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
CBS	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
CBD	0.260	0.259	0.00308	%	0.298	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
δ8THCV	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
CBN	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
δ9THC	0.269	0.268	0.00308	%	0.420	< 20	Acceptable	
δ8THC	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
9Sa10THC	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
CB	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
9Rd10THC	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
CB	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
CBCA	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
δ9THCP	<LOQ	<LOQ	0.00308	%	NA	< 20	Acceptable	
CB	<LOQ	<LOQ	0.00308	%	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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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.