



CONSOLIDATED TEST RESULTS SUMMARY

Please see the following pages for full test results

PRODUCT	Live Resin CBD/Delta-8 Tincture	BATCH#	21-012822-0001	LOQ: Limit of Quantification	
SKU	30ML-LR-CBD-D8-TINCTURE	SERVING SIZE	1.0ml	LOD: Limit of Detection	
LABORATORY	Columbia Laboratory	ACCREDITATION	OR1000028	1g = 1000 mg = 100,000 ug 1 mg/kg = 1ppm = 1000 ppb	
CANNABINOID	PERCENTAGE	MG/SERVING	HEAVY METALS	PER/SERVING	ACTION LEVEL
CBD (Cannabidiol)	2.02	20.20	Arsenic	LOQ	10 ug/day *
Delta-8 THC	1.48	14.80	Cadmium	LOQ	4.1 ug/day *
CBT (Cannabicitran)	0.0848	0.85	Lead	LOQ	3.5 ug/day **
CBC (Cannabichromene)	0.0397	0.40	Mercury	LOQ	2 ug/day *
CBG (Cannabigerol)	0.0326	0.33	RESIDUAL SOLVENTS		
CBE (Cannabielsoin)	0.0113	0.11	All analytes less than LOQ	PASS	---
CBDA	0.0098	0.10			
Delta-8 THCV	0.0096	0.10			
CBDV	0.00751	0.08			
TOTAL Cannabinoids	3.61051	36.11			
PESTICIDES	RESULT	LIMIT	MICROBIOLOGY	RESULTS	ACTION LEVEL
All analytes less than LOQ	PASS	---	All analytes less than LOQ	PASS	---
PRODUCT PHOTO(S)					



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

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



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



Report Number: 21-012822/D006.R000
Report Date: 11/08/2021
ORELAP#: OR100028
Purchase Order:
Received: 10/29/21 15:41

Customer: Woodstock Hemp Co.
Product identity: WHC Live Resin 2:1 CBD + D8 Tincture
Client/Metric ID: .
Laboratory ID: 21-012822-0001

Summary

Potency:

Analyte	Result (%)			
CBD	2.02		CBD-Total	2.03%
Δ8-THC†	1.48		THC-Total	<LOQ
CBT†	0.0848		(Reported in percent of total sample)	
CBC	0.0397			
CBG†	0.0326			
CBE†	0.0113			
CBD-A	0.00980			
Δ8-THCV	0.00960			
CBDV†	0.00751			

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: Woodstock Hemp Co.
Product identity: WHC Live Resin 2:1 CBD + D8 Tincture
Client/Metric ID: .
Sample Date:
Laboratory ID: 21-012822-0001
Evidence of Cooling: No
Temp: 19.2 °C
Relinquished by: Client

Sample Results

Potency	Method J AOAC 2015 V98-6 (mod)			Units %	Batch: 2109934	Analyze: 11/3/21 6:33:00 AM
Analyte	As Received	Dry weight	LOQ	Notes		
CBC	0.0397		0.0032			
CBC-A†	< LOQ		0.0032			
CBC-Total†	0.0397		0.0061			
CBD	2.02		0.0324			
CBD-A	0.00980		0.0032			
CBD-Total	2.03		0.0352			
CBDV†	0.00751		0.0032			
CBDV-A†	< LOQ		0.0032			
CBDV-Total†	0.00751		0.0060			
CBE†	0.0113		0.0032			
CBG†	0.0326		0.0032			
CBG-A†	< LOQ		0.0032			
CBG-Total	0.0326		0.0060			
CBL†	< LOQ		0.0032			
CBL-A†	< LOQ		0.0032			
CBL-Total†	< LOQ		0.0061			
CBN	< LOQ		0.0032			
CBT†	0.0848		0.0032			
Δ8-THC†	1.48		0.0324			
Δ8-THCV	0.00960		0.0032			
Δ9-THC	< LOQ		0.0032			
THC-A	< LOQ		0.0032			
THC-Total	< LOQ		0.0061			
THCV†	< LOQ		0.0032			
THCV-A†	< LOQ		0.0032			
THCV-Total†	< LOQ		0.0060			
Total Cannabinoids†	3.70					



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



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Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyze	Method	Status	Notes
E.coli	< LOQ		cfu/g	10	2109823	11/03/21	AOAC 991.14 (Petrifilm)		X
Total Coliforms	< LOQ		cfu/g	10	2109823	11/03/21	AOAC 991.14 (Petrifilm)		X
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2109811	11/03/21	AOAC 2014.05 (RAPID)		X
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2109811	11/03/21	AOAC 2014.05 (RAPID)		X

Solvents Method Residual Solvents by GC/MS Units µg/g Batch 2110022 Analyze 11/05/21 03:42 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	< 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 2109929 Analyze 11/03/21 04:08 PM

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

Metals

Analyte	Result	Limits	Units	LOQ	Batch	Analyze	Method	Status	Notes
Arsenic	< LOQ	0.200	mg/kg	0.0478	2109948	11/03/21	AOAC 2013.06 (mod.)	pass	X
Cadmium	< LOQ	0.200	mg/kg	0.0478	2109948	11/03/21	AOAC 2013.06 (mod.)	pass	X
Lead	< LOQ	0.500	mg/kg	0.0478	2109948	11/03/21	AOAC 2013.06 (mod.)	pass	X
Mercury	< LOQ	0.100	mg/kg	0.0239	2109948	11/03/21	AOAC 2013.06 (mod.)	pass	X



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Mycotoxins

Analyte	Result	Limits	Units	LOQ	Batch	Analyze	Method	Status	Notes
Aflatoxin B2 [†]	< LOQ		µg/kg	5.00	2109863	11/04/21	AOAC 2007.01 & EN 15		
Aflatoxin B1 [†]	< LOQ		µg/kg	5.00	2109863	11/04/21	AOAC 2007.01 & EN 15		
Aflatoxin G1 [†]	< LOQ		µg/kg	5.00	2109863	11/04/21	AOAC 2007.01 & EN 15		
Aflatoxin G2 [†]	< LOQ		µg/kg	5.00	2109863	11/04/21	AOAC 2007.01 & EN 15		
Deoxynivalenol [†]	< LOQ		µg/kg	200	2109863	11/04/21	AOAC 2007.01 & EN 15		
Fumonisin B1 [†]	< LOQ		µg/kg	200	2109863	11/04/21	AOAC 2007.01 & EN 15		
Fumonisin B2 [†]	< LOQ		µg/kg	200	2109863	11/04/21	AOAC 2007.01 & EN 15		
HT2-Toxin [†]	< LOQ		µg/kg	40.0	2109863	11/04/21	AOAC 2007.01 & EN 15		
Nivalenol [†]	< LOQ		µg/kg	400	2109863	11/04/21	AOAC 2007.01 & EN 15		
Ochratoxin A [†]	< LOQ		µg/kg	5.00	2109863	11/04/21	AOAC 2007.01 & EN 15		
Ochratoxin B [†]	< LOQ		µg/kg	2.00	2109863	11/04/21	AOAC 2007.01 & EN 15		
T2-Toxin [†]	< LOQ		µg/kg	20.0	2109863	11/04/21	AOAC 2007.01 & EN 15		
Zearalenone [†]	< LOQ		µg/kg	200	2109863	11/04/21	AOAC 2007.01 & EN 15		



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

Abbreviations

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

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

† = Analyte not NELAP accredited.

Units of Measure

cfu/g = Colony forming units per gram

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

X: Not ORELAP accredited.

Approved Signatory

Derrick Tanner
General Manager



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Cannabis Multi-Residue Profile, Limits of Quantitation

Compound	LOQ (mg/kg)	Compound	LOQ (mg/kg)	Compound	LOQ (mg/kg)
Abamectin	0.100	Cbthodim	0.050	Endrin	0.100
Acephate	0.100	Cbthodim Sulfone	0.050	EPN	0.050
Acequinocyl	0.100	Cbthodim Sulfoxide	0.050	EPTC	0.100
Acetamiprid	0.020	Cbfentazine	0.020	Esfenvalerate/ Fenvalerate	0.200
Acetochlor	0.100	Cbmazone	0.020	Etaconazole	0.100
Acrinathrin	0.100	Cbthianidin	0.200	Ethalfuralin	0.100
Alachlor	0.100	Cumaphos	0.050	Ethiofencarb	0.050
Aldicarb	0.100	Crdoxyphos	0.020	Ethion	0.200
Aldicarb sulfoxide	0.100	Cyarazine	0.020	Ethirimol	0.100
Aldoxycarb (Aldicarb-sulfone)	0.100	Cyarofenphos	0.020	Ethofumesate	0.050
Aldrin	0.100	Cyatr anilprole	0.050	Ethoprophos	0.020
Ametoctradin	0.020	Cyazfamid	0.020	Etofenprox	0.020
Ametryn	0.500	Cydoate	0.100	Etoxazole	0.020
Aspon	0.100	Cyfluthrin	0.200	Etridiazole	0.100
Asulam	0.100	Cyhalothrin, lambda	0.200	Etrinfos	0.020
Atrazine	0.100	Cymoxanil	0.050	Famoxadone	0.200
Atrazine-desethyl	0.100	Cypermethrin	0.200	Famphur	0.100
Azinphos-ethyl	0.020	Cyprodinil	0.100	Fenamidone	0.020
Azinphos-methyl	0.020	Dadhal	0.100	Fenamiphos	0.020
Azoxystrobin	0.020	Damhozide	0.100	Fenamiphos sulfone	0.020
Beralaxyl	0.020	DCPMU	0.050	Fenamiphos sulfoxide	0.020
Bendiocarb	0.020	DDD, op'-	0.100	Fenazaquin	0.100
Berfluralin	0.100	DDD, p,p'-	0.100	Fenbuconazole	0.100
Berxacor	0.050	DCE, o,p'-	0.100	Fenchlorphos	0.100
Bersulide	0.050	DCE, p,p'-	0.100	Fenchlorphos-oxon	0.100
BHC alpha isomer	0.100	DDT, o,p'-	0.100	Fenhexamid	0.100
BHC beta isomer	0.100	DDT, p,p'-	0.100	Fenitrothion	0.100
BHC delta isomer	0.500	DEF (Tribufos)	0.100	Fenobucarb	0.050
Bifenazate	0.020	Detamethrin	0.100	Fenoxycarb	0.020
Bifenthrin	0.020	Desmedipham	0.100	Fenpropathrin	0.050
Boscalid	0.020	Diallate	0.100	Fenpyroximate	0.020
Bromophos-ethyl	0.100	Diazinon	0.020	Fenson	0.100
Bromophos-methyl	0.200	Diazoxon	0.100	Fensulfiothion	0.020
Bromopropylate	0.100	Dichlobenil	0.100	Fensulfiothion oxon	0.020
Bromuconazole	0.100	Dichlofluanid	0.100	Fensulfiothion sulfone	0.100
Bupirimate	0.020	Dichlorvos	0.100	Fensulfiothion-oxon-sulfone	0.020
Buprofezin	0.050	Diclobutrazol	0.050	Fenthion	0.050
Butadlor	0.500	Dicofol	0.100	Fenthion oxon	0.020
Butralin	0.200	Dicrotophos	0.050	Fenthion oxon sulfone	0.100
Butylate	0.100	Dieldrin	0.100	Fenthion sulfone	0.050
Cadusafos	0.020	Diethofencarb	0.020	Fenuron	0.020
Captan	1.000	Diethyltoluamide (DEET)	0.050	Fipronil	0.100
Carbaryl	0.050	Difenoconazole	0.100	Fonicamid	0.100
Carbendazim	0.100	Dimethenamid	0.050	Fluchloralin	0.100
Carbendiazim	0.100	Dimethoate	0.050	Flucythrinate	0.100
Carbendiazim	0.020	Dimethomorph	0.050	Fludioxonil	0.200
Carbophenothion	0.200	Diniconazole	0.200	Flufenacet	0.020
Carboxin	0.020	Dinotefuran	0.200	Fumioxazin	0.100
Carfentrazone-ethyl	0.100	Dioxathion	0.100	Fuometuron	0.020
Chlorantrilprole	0.020	Diphenamid	0.020	Fluopicolide	0.050
Chordane, cis-	0.200	Diphenylamine	0.100	Fluopyram	0.020
Chordane, trans-	0.200	Disulfoton	0.100	Fuoxastrobin	0.050
Chlorfenapyr	0.500	Disulfoton sulfone	0.100	Flupyradifurone	0.020
Chlorfenson	0.200	Disulfoton sulfoxide	0.100	Furidone	0.100
Chlorfenvinphos	0.050	Diuron	0.050	Fusilazole	0.020
Chlorobenzilate	0.100	Edifenphos	0.050	Flutolanil	0.020
Chloroneb	0.200	Endosulfan alpha	0.200	Flutriafol	0.020
Chlorpyrifos	0.050	Endosulfan beta	0.200	Fluvalinate, tau-	0.100
Chlorpyrifos-methyl	0.200	Endosulfan sulfate	0.100	Fluxapyroxad	0.020
CIPC	1.000				



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Cannabis Multi-Residue Profile, Limits of Quantitation

Compound	LOQ(mg/kg)	Compound	LOQ(mg/kg)	Compound	LOQ(mg/kg)
Fomesafen	0.100	Mexacarbate	0.020	Prpamacarb	0.050
Fonofos	0.100	MGK 264	0.020	Prparil	0.050
Forchlorfenuron	0.050	Mirex	0.100	Prpargite	0.050
Formetanate	0.050	Molinate	0.050	Prpazine	0.020
Furathiocarb	0.020	Monocrotophos	0.100	Prpzetamphos	0.050
Heptachlor	0.100	Monolinuron	0.020	Prppham	0.050
Heptachlor epoxide	0.100	Myclobutanil	0.050	Prppiconazole	0.050
Heptenophos	0.100	Naled	0.100	Prppoxur	0.050
Hexachlorobenzene	0.100	Napropamide	0.050	Prppoxycarbazone Na	0.050
Hexaconazole	0.100	Neburon	0.020	Prppyzamide	0.050
Hexazinone	0.100	Nitrapyrin	0.100	Prrothiofos	0.100
Hexythiazox	0.020	Norflurazon	0.050	Prpyraclostrobin	0.020
Imazalil	0.100	Omethoate	0.100	Prpyazophos	0.050
Imidacoprid	0.100	O-Phenylphenol	0.100	Prpyrethrins	0.050
Indaziflam	0.020	Oxadixyl	0.100	Prpyridaben	0.020
Indoxacarb	0.020	Oxamyl	0.100	Prpyridafol	0.100
Iprobenfos	0.100	Oxamyl-oxime	0.100	Prpyridate	0.020
Iprodione	0.100	Oxychloridane	0.100	Prpyrimetharil	0.050
Isobenzan	0.100	Oxydemeton-Methyl	0.100	Prpyriproxifen	0.020
Isocarbophos	0.500	Oxythioquinox	0.200	Prpyroxasulfone	0.020
Isodrin	0.100	Padlobutrazol	0.050	Prpyroxulam	0.020
Isfenphos	0.050	Paraaxon-ethyl	0.020	Quinalphos	0.050
Isfenphos-methyl	0.020	Paraaxon methyl	0.100	Quinoxifen	0.050
Isfenphos oxon	0.050	Parathion ethyl	0.100	Quintozene (PQNB)	0.200
Isoprocarb	0.020	Parathion methyl	0.200	Resmethrin	0.050
Isopropalin	0.200	Perconazole	0.050	Rotenone	0.050
Isoprothiolane	0.050	Perdimethalin	0.050	S421	0.100
Isoproturon	0.050	Perflufen	0.020	Smaazine	0.100
Isoxaben	0.050	Pertachloroaniline	0.100	Smetryn	0.200
Isoxaflutole	0.050	Pertachloroanisole	0.100	Spinetoram	0.020
Kresoxim-methyl	0.050	Pentachlorobenzene (PCB)	0.100	Spinosad	0.050
Ladofen	0.500	Pentachlorothioanisole (PCTA)	0.100	Spirodiclofen	0.100
Lenadi	0.100	Perthiopyrad	0.020	Spiromesifen	0.050
Lindane (gammaBHC)	0.100	Permethrin	0.050	Spirotetramat	0.050
Linuron	0.020	Perthane	0.100	Spiroxamine	0.020
Malaoxon	0.050	Phenmedipham	0.050	Sulfotep	0.050
Malathion	0.050	Phenthoate	0.050	Sulfoxaflof	0.050
Mandipropamid	0.020	Phorate	0.050	Sulprofos	0.020
Mecarbam	0.020	Phorate Sulfone	0.050	Tebuconazole	0.100
Mepanipyrim	0.050	Phorate Sulfoxide	0.050	Tebufenozide	0.020
Merphos	0.500	Phosalone	0.050	Tebuthiuron	0.020
Metalaxyl	0.050	Phosmet	0.100	Tecnazene	0.100
Metaldehyde	0.050	Phosphamidon	0.050	Tefluthrin	0.100
Metconazole	0.100	Phoxim	0.050	Terbufos	0.020
Methacifos	0.100	Pinoxaden	0.020	Terbufos sulfone	0.050
Methamidophos	0.050	Piperonyl butoxide	0.050	Terbufos sulfoxide	0.050
Methidathion	0.050	Pirimicarb	0.020	Terbuthylazine	0.020
Methiocarb	0.050	Pirimiphos-methyl	0.050	Terbutryn	0.020
Methiocarb sulfone	0.100	Pirimiphos-ethyl	0.020	Tetrachlorvinphos	0.050
Methiocarb sulfoxide	0.100	Prallethrin	0.100	Tetraconazole	0.050
Methomyl	0.100	Prochloraz	0.020	Tetradfon	0.200
Methoxychlor	0.100	Proxymidone	0.100	Tetramethrin	0.050
Methoxyfenozide	0.020	Prfenofos	0.100	Tetrasul	0.100
Metobromuron	0.050	Prfluralin	0.100	Thiabendazole	0.100
Metolachlor	0.100	Promecarb	0.050	Thiabendazole, 5-hydroxy	0.100
Metolcarb	0.050	Prometon	0.100	Thiadoprid	0.050
Metraferone	0.050	Prometryn	0.020	Thiamethoxam	0.100
Metribuzin	0.100	Prpadhlor	0.020	Thiobencarb	0.050
Mevinphos	0.100			Thiodicarb	0.050
				Thiophanate-methyl	0.050



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Cannabis Multi-Residue Profile, Limits of Quantitation

Compound	LOQ(mg/kg)	Compound	LOQ(mg/kg)	Compound	LOQ(mg/kg)
Tolclofos-methyl	0.100	Triazophos	0.020	Trifloxystrobin	0.020
Triforin	0.100	Tolyfluarid	0.050	Triticonazole	0.050
Tralkoxydim	0.100	Tridiphane	0.500	Vindozolin	0.100
Triadimefon	0.050	Triflumizde	0.020	Zoxamide	0.020
Triallate	0.100	Trifluralin	0.100		

LOQ=Limit of Quantitation, mg/kg

Factors affecting the LOQ include instrumentation sensitivity for 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 Products

Chain of Custody Record

Document Control ID: 2732 Revision: 1
 Effective: 05/04/2021
 ORELAP ID: OR100028

Company: <u>Woodstock Hemp Company</u> Contact: <u>Jason Peterson / Jerrod Peterson</u> Street: <u>23355 S Kamrath rd</u> City: <u>Oregon City</u> State: <u>OR</u> Zip: <u>97045</u> <input checked="" type="checkbox"/> Email Results: <u>jason@woodstockhempcompany.com</u> Ph: <u>(503) 4074500</u> <input type="checkbox"/> Fx Results: () Billing (if different): <u>14757 Glen Oaks rd</u> <u>Oregon City, OR 97045</u>				Analysis Requested Pesticides - OR 59 compounds Pesticide Multi-Residue - 379 compounds Potency Residual Solvents Moisture & Water Activity Terpenes Micro: Yeast and Mold Micro: E.Coli and Total Coliform Heavy Metals Mycotoxins Other:										PO Number: _____ Project Number: _____ Project Name: _____ Custom Reporting: _____ Report to State - <input type="checkbox"/> METRC or <input type="checkbox"/> Other: _____ Turnaround time: <input checked="" type="checkbox"/> 5 Business Day Standard Turnaround <input type="checkbox"/> 3 Business Day Rush Turnaround* <input type="checkbox"/> 2 Business Day Rush Turnaround* <i>*Check for availability</i>			
Lab ID	Client Sample Identification	Date	Time	Pesticides - OR 59 compounds	Pesticide Multi-Residue - 379 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
	{ Pura Vida Honeycomb Ginger CBD Tincture Batch # 10272101	10/29/21			X	X	X			X	X	X	X	*	T	30 ml	* Mycotoxins analysis canceled via email. -PG 10/29
	{ THC Free Broad Spectrum Distillate	10/29/21				X									C	20 ml	
	{ Perform Oil # PO150921	10/29/21				X	X			X	X	X	X		T	30 ml	
	{ WHC Live Resin CBD Tincture 1,500mg				X	X	X			X	X	X	X		T	30 ml	
Relinquished By:		Date	Time	Received By:		Date	Time	Lab Use Only:									
		10/29/21	15:41	BD		10/29/21	15:41	<input type="checkbox"/> Shipped Via: _____ or <input checked="" type="checkbox"/> Client drop Evidence of cooling: <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No - Temp (°C): <u>19.2</u> Sample in good condition: <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> Cash <input type="checkbox"/> Check <input type="checkbox"/> CC <input type="checkbox"/> 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 with testing requirements constitute an agreement for services in accordance with the [current terms of service](#) 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 _____



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

Report Number: 21-012822/D006.R000
 Report Date: 11/08/2021
 ORELAP#: OR100028
 Purchase Order:
 Received: 10/29/21 15:41



Hemp / Cannabis Usable / Extract / Finished Products

Chain of Custody Record

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Company: <u>Woodstock Hemp Company</u> Contact: <u>Jason Peterson / Jerrod Peterson</u> Street: <u>23355 S Kamrath rd</u> City: <u>Oregon City</u> State: <u>OR</u> Zip: <u>97045</u> <input checked="" type="checkbox"/> Email Results: <u>jason@woodstockhempcompany.com</u> Ph: <u>(503) 4074500</u> <input type="checkbox"/> Fx Results: () Billing (if different): <u>14757 Glen Oaks rd</u> <u>Oregon City, OR 97045</u>				Analysis Requested <input type="checkbox"/> Pesticides - OR 59 compounds <input type="checkbox"/> Pesticide Multi-Residue - 379 compounds <input type="checkbox"/> Potency <input type="checkbox"/> Residual Solvents <input type="checkbox"/> Moisture & Water Activity <input type="checkbox"/> Terpenes <input type="checkbox"/> Micro: Yeast and Mold <input type="checkbox"/> Micro: E.Coli and Total Coliform <input type="checkbox"/> Heavy Metals <input type="checkbox"/> Mycotoxins <input type="checkbox"/> Other:										PO Number: _____ Project Number: _____ Project Name: _____ Custom Reporting: _____ Report to State - <input type="checkbox"/> METRC or <input type="checkbox"/> Other: _____ Turnaround time: <input checked="" type="checkbox"/> 5 Business Day Standard Turnaround <input type="checkbox"/> 3 Business Day Rush Turnaround* <input type="checkbox"/> 2 Business Day Rush Turnaround* *Check for availability Sampled by: _____			
Lab ID	Client Sample Identification	Date	Time	Pesticides - OR 59 compounds	Pesticide Multi-Residue - 379 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
	{ WHC Live Resin CBD Vapor Cartridge	10/29/21		X	X	X				X	X	X	X		C	1.0g	> Not enough sample submit
	{ WHC Full Spectrum DB Vapor Cartridge	10/29/21		X	X	X				X	X	X	X		C	1.0g	> Not enough sample submit
	{ WHC DB Vapor Cartridge	10/29/21		X	X	X				X	X	X	X		C	1.0g	> Not enough sample submit
	{ WHC Live Resin 2:1 CBD + DB Tincture			X	X	X				X	X	X	X		T	30 ml	
Relinquished By:		Date	Time	Received By:		Date	Time	Lab Use Only:									
<i>JWP</i>		10/29/21	15:41	<i>BD</i>		10/29/21	15:41	<input type="checkbox"/> Shipped Via: _____ or <input checked="" type="checkbox"/> Client drop Evidence of cooling: <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No - Temp (°C): <u>19.2</u> Sample in good condition: <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No <input type="checkbox"/> Cash <input type="checkbox"/> Check <input type="checkbox"/> CC <input type="checkbox"/> Net: _____ Prelog storage: _____									

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

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

Report Number: 21-012822/D006.R000
Report Date: 11/08/2021
ORELAP#: OR100028
Purchase Order:
Received: 10/29/21 15:41



Revision #: 0.00 Control : CFL-D06
Revision Date: 05/31/2019 Effective Date: 05/31/2019

Laboratory Quality Control Results

JAOAC2015 V986 Batch ID: 2109934

Laboratory Control Sample							
Analyte	Result	Spike	Units	% Rec	Limits	Evaluation	Notes
CBDVA	0.00959	0.01	%	95.9	85.0 - 115	Acceptable	
CBDV	0.0103	0.01	%	103	85.0 - 115	Acceptable	
CBE	0.00998	0.01	%	99.8	85.0 - 115	Acceptable	
CBDA	0.00990	0.01	%	99.0	85.0 - 115	Acceptable	
CBGA	0.00961	0.01	%	96.1	85.0 - 115	Acceptable	
CBG	0.0100	0.01	%	100	85.0 - 115	Acceptable	
CBD	0.0101	0.01	%	101	85.0 - 115	Acceptable	
THCV	0.00962	0.01	%	96.2	85.0 - 115	Acceptable	
d8THCV	0.00967	0.01	%	96.7	85.0 - 115	Acceptable	
THCVA	0.00941	0.01	%	94.1	85.0 - 115	Acceptable	
CBN	0.0103	0.01	%	103	85.0 - 115	Acceptable	
exo-THC	0.0102	0.01	%	102	85.0 - 115	Acceptable	
d9THC	0.00991	0.01	%	99.1	85.0 - 115	Acceptable	
d8THC	0.00951	0.01	%	95.1	85.0 - 115	Acceptable	
CBL	0.00935	0.01	%	93.5	85.0 - 115	Acceptable	
CBC	0.00999	0.01	%	99.9	85.0 - 115	Acceptable	
THCA	0.00995	0.01	%	99.5	85.0 - 115	Acceptable	
CBCA	0.00981	0.01	%	98.1	85.0 - 115	Acceptable	
CBLA	0.00995	0.01	%	99.5	85.0 - 115	Acceptable	
CBT	0.0112	0.01	%	112	85.0 - 115	Acceptable	

Method Blank

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

Abbreviations

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

Units of Measure:

% - Percent



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Report Number: 21-012822/D006.R000
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Received: 10/29/21 15:41

Revision #: 0.00 Control : CFL-D06
 Revision Date: 05/31/2019 Effective Date: 05/31/2019

Laboratory Quality Control Results

JAOAC2015 V986		Batch ID: 2109934						
Sample Duplicate		Sample ID: 21-012818-0001						
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBDV	0.00455	0.00454	0.003	%	0.334	< 20	Acceptable	
CBE	0.00805	0.00828	0.003	%	2.82	< 20	Acceptable	
CBDA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBG	0.0127	0.0126	0.003	%	0.601	< 20	Acceptable	
CBD	0.912	0.911	0.003	%	0.0379	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBN	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
d9THC	0.0248	0.0246	0.003	%	0.659	< 20	Acceptable	
d8THC	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBC	0.0474	0.0476	0.003	%	0.321	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBCA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.003	%	NA	< 20	Acceptable	
CBT	0.00545	0.00534	0.003	%	2.09	< 20	Acceptable	

Abbreviations

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

Units of Measure:

%- Percent



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Report Number: 21-012822/D006.R000
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Received: 10/29/21 15:41

Revision: Document ID:
Legacy ID: Effective:

Laboratory Quality Control Results

Batch ID: 2110022

Residual Solvents		Method Blank		Laboratory Control Sample					
Analyte	Result	LOQ	Notes	Result	Spike	Units	%Rec	Limits	Notes
Propane	ND	< 200		378	407	µg/g	92.8	70	130
Isobutane	ND	< 200		467	491	µg/g	95.1	70	130
Butane	ND	< 200		464	491	µg/g	94.5	70	130
2,2-Dimethylpropane	ND	< 200		609	609	µg/g	100.0	70	130
Methanol	ND	< 200		1870	1810	µg/g	116.1	70	130
Ethylene Oxide	ND	< 30		35.9	38.9	µg/g	92.3	70	130
2-Methylbutane	ND	< 200		1650	1610	µg/g	102.5	70	130
Pentane	ND	< 200		1640	1610	µg/g	101.9	70	130
Ethanol	ND	< 200		1720	1610	µg/g	106.8	70	130
Ethyl Ether	ND	< 200		1480	1610	µg/g	91.9	70	130
2,2-Dimethylbutane	ND	< 30		143	164	µg/g	87.2	70	130
Acetone	ND	< 200		1630	1610	µg/g	101.2	70	130
2-Propanol	ND	< 200		1740	1610	µg/g	108.1	70	130
Acetonitrile	ND	< 100		507	484	µg/g	104.8	70	130
2,3-Dimethylbutane	ND	< 30		173	167	µg/g	103.6	70	130
Dichloromethane	ND	< 60		429	491	µg/g	87.4	70	130
2-Methylpentane	ND	< 30		158	169	µg/g	95.8	70	130
3-Methylpentane	ND	< 30		166	172	µg/g	96.5	70	130
Hexane	ND	< 30		160	167	µg/g	95.8	70	130
Ethyl acetate	ND	< 200		1650	1610	µg/g	102.5	70	130
2-Butanol	ND	< 200		1660	1610	µg/g	103.1	70	130
Tetrahydrofuran	ND	< 100		440	483	µg/g	91.1	70	130
Cyclohexane	ND	< 200		1340	1610	µg/g	83.2	70	130
Benzene	ND	< 1		4.5	5.36	µg/g	84.0	70	130
Isopropyl Acetate	ND	< 200		1620	1620	µg/g	112.3	70	130
Heptane	ND	< 200		1770	1610	µg/g	109.9	70	130
1,4-Dioxane	ND	< 100		405	488	µg/g	82.8	70	130
2-Ethoxyethanol	ND	< 30		183	167	µg/g	109.6	70	130
Ethylene Glycol	ND	< 200		527	504	µg/g	104.6	70	130
Toluene	ND	< 200		394	484	µg/g	81.4	70	130
Ethylbenzene	ND	< 200		771	968	µg/g	79.6	70	130
m,p-Xylene	ND	< 200		797	977	µg/g	81.6	70	130
o-Xylene	ND	< 200		787	982	µg/g	80.1	70	130
Cumene	ND	< 30		124	169	µg/g	73.4	70	130



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Received: 10/29/21 15:41

Revision: Document ID:
 Legacy ID: Effective:

CC - Simple Duplicate Sample ID: 21-012816-0001

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	
Acetonitrile	ND	ND	100	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	
3-Methylpentane	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Hexane	ND	ND	30	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	
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	
1,4-Dioxane	ND	ND	100	µg/g	0.0	< 20	Acceptable	
2-Ethoxyethanol	ND	ND	30	µg/g	0.0	< 20	Acceptable	
Ethylene Glycol	ND	ND	200	µg/g	0.0	< 20	Acceptable	
Toluene	ND	ND	200	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	

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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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.