



Customer: NW Natural Goods
Product identity: BEV - BB -B-58-1
Metrc ID: .
Metrc Source ID:
Laboratory ID: 25-002244-0001

Summary

Potency:

Analyte per 355ml	Result	Limits	Units	Status	
CBC per 355ml	0.482		mg/355ml		CBD-Total per Serving Size 51.4 mg/355ml
CBD per 355ml	51.4		mg/355ml		
CBG per 355ml	1.63		mg/355ml		Delta-9-THC-Total per <LOQ
					(Reported in milligrams per serving)

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.



Customer: NW Natural Goods
 United States of America (USA)
Product identity: BEV - BB -B-58-1
Metrc ID: .
Metrc Source ID:
Material: Cannabinoid Beverage
Sample Date:
Laboratory ID: 25-002244-0001
Evidence of Cooling: No
Temp: 17.4 °C
Serving Size #1: 362.1 g
Density: 1.020 g/ml

Sample Results

Potency per 355ml	Method: J AOAC 2015 V98-6 (mod) ^b		Units mg/se	Batch: 2501615	Analyze: 3/6/25	2:32:00 PM
Analyte	Result	Limits	Units	LOQ	Notes	
CBC per 355ml	0.482		mg/355ml	0.352		
CBC-A per 355ml	< LOQ		mg/355ml	0.352		
CBC-Total per 355ml	< LOQ		mg/355ml	0.660		
CBD per 355ml	51.4		mg/355ml	0.352		
CBD-A per 355ml ¹	< LOQ		mg/355ml	0.352		
CBD-Total per 355ml ¹	51.4		mg/355ml	0.660		
CBDV per 355ml	< LOQ		mg/355ml	0.352		
CBDV-A per 355ml	< LOQ		mg/355ml	0.352		
CBDV-Total per 355ml	< LOQ		mg/355ml	0.657		
CBE per 355ml	< LOQ		mg/355ml	0.352		
CBG per 355ml	1.63		mg/355ml	0.352		
CBG-A per 355ml	< LOQ		mg/355ml	0.352		
CBG-Total per 355ml	1.63		mg/355ml	0.657		
CBL per 355ml	< LOQ		mg/355ml	0.352		
CBL-A per 355ml	< LOQ		mg/355ml	0.352		
CBL-Total per 355ml	< LOQ		mg/355ml	0.660		
CBN per 355ml	< LOQ		mg/355ml	0.352		
CBT per 355ml	< LOQ		mg/355ml	0.352		
Δ10-THC-9R per 355ml	< LOQ		mg/355ml	0.352		
Δ10-THC-9S per 355ml	< LOQ		mg/355ml	0.352		
Δ10-THC-Total per 355ml	< LOQ		mg/355ml	0.704		
Δ8-THC per 355ml ¹	< LOQ		mg/355ml	0.352		
Δ8-THCV per 355ml	< LOQ		mg/355ml	0.352		
Δ9-THC per 355ml ¹	< LOQ		mg/355ml	0.352		
Δ9-THC-Total per 355ml	< LOQ		mg/355ml	0.660		
Δ9-THCP per 355ml	< LOQ		mg/355ml	0.352		
Δ9-THCV per 355ml	< LOQ		mg/355ml	0.352		
Δ9-THCV-A per 355ml	< LOQ		mg/355ml	0.352		
Δ9-THCV-Total per 355ml	< LOQ		mg/355ml	0.661		



Potency per 355ml Method: J AOAC 2015 V98-6 (mod)^p Units mg/se Batch: 2501615 Analyze: 3/6/25 2:32:00 PM

Analyte	Result	Limits	Units	LOQ	Notes
exo-THC per 355ml	< LOQ		mg/355ml	0.352	
THC-A per 355ml ¹	< LOQ		mg/355ml	0.352	
Total Cannabinoids per 355ml	53.5		mg/355ml		

Microbiology

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
Aerobic Plate Count	< LOQ		cfu/g	10	2501549	03/07/25 AOAC 990.12 (Petrifilm)		
E.coli	< LOQ		cfu/g	10	2501547	03/07/25 AOAC 991.14 (Petrifilm)		
Total Coliforms	< LOQ		cfu/g	10	2501547	03/07/25 AOAC 991.14 (Petrifilm)		
Mold (RAPID Petrifilm)	< LOQ		cfu/g	10	2501548	03/08/25 AOAC 2014.05 (RAPID)		
Yeast (RAPID Petrifilm)	< LOQ		cfu/g	10	2501548	03/08/25 AOAC 2014.05 (RAPID)		

Solvents Method: Residual Solvents by HS-GC-MS^p Units µg/g Batch 2501655 Analyze 03/07/25 02:52 PM

Analyte	Result	Limits	LOQ	Status	Notes	Analyte	Result	Limits	LOQ	Status	Notes
1,4-Dioxane ¹	< LOQ	380	100	pass		2-Butanol ¹	< LOQ	5000	200	pass	
2-Ethoxyethanol ¹	< LOQ	160	30.0	pass		2-Methylbutane (Isopentane) ¹	< LOQ		200		
2-Methylpentane ¹	< LOQ		30.0			2-Propanol (IPA) ¹	< LOQ	5000	200	pass	
2,2-Dimethylbutane ¹	< LOQ		30.0			2,2-Dimethylpropane (neo-pentane) ¹	< LOQ		200		
2,3-Dimethylbutane ¹	< LOQ		30.0			3-Methylpentane ¹	< LOQ		30.0		
Acetone ¹	< LOQ	5000	200	pass		Acetonitrile ¹	< LOQ	410	100	pass	
Benzene ¹	< LOQ	2.00	1.00	pass		Butanes (sum) ¹	< LOQ	5000	400	pass	
Cyclohexane ¹	< LOQ	3880	200	pass		Ethyl acetate ¹	< LOQ	5000	200	pass	
Ethyl benzene	< LOQ		200			Ethyl ether ¹	< LOQ	5000	200	pass	
Ethylene glycol ¹	< LOQ	620	200	pass		Ethylene oxide ¹	< LOQ	50.0	20.0	pass	
Hexanes (sum) ¹	< LOQ	290	150	pass		Isopropyl acetate ¹	< LOQ	5000	200	pass	
Isopropylbenzene (Cumene) ¹	< LOQ	70.0	30.0	pass		m,p-Xylene ¹	< LOQ		200		
Methanol ¹	< LOQ	3000	200	pass		Methylene chloride ¹	< LOQ	600	60.0	pass	
Methylpropane (Isobutane) ¹	< LOQ		200			n-Butane ¹	< LOQ		200		
n-Heptane ¹	< LOQ	5000	200	pass		n-Hexane ¹	< LOQ		30.0		
n-Pentane ¹	< LOQ		200			o-Xylene ¹	< LOQ		200		
Pentanes (sum)	< LOQ	5000	600	pass		Propane	< LOQ	5000	200	pass	
Tetrahydrofuran ¹	< LOQ	720	100	pass		Toluene ¹	< LOQ	890	100	pass	
Total Xylenes ¹	< LOQ		400			Total Xylenes and Ethyl benzene	< LOQ	2170	600	pass	

Pesticides Method: AOAC 2007.01 Units mg/kg Batch 2501693 Analyze 03/10/25 12:01 PM

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



Metals

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
Arsenic [±]	< LOQ	0.200	mg/kg	0.00390	2501604	03/06/25 AOAC 2013.06 (mod.) ^Þ	pass	
Cadmium [±]	< LOQ	0.200	mg/kg	0.00390	2501604	03/06/25 AOAC 2013.06 (mod.) ^Þ	pass	
Lead [±]	< LOQ	0.500	mg/kg	0.00390	2501604	03/06/25 AOAC 2013.06 (mod.) ^Þ	pass	
Mercury [±]	< LOQ	0.100	mg/kg	0.00195	2501604	03/06/25 AOAC 2013.06 (mod.) ^Þ	pass	

Mycotoxins

Analyte	Result	Limits	Units	LOQ	Batch	Analyzed Method	Status	Notes
Aflatoxin B1 [±]	< LOQ		µg/kg	5.00	2501729	03/11/25 Mycotoxins by AOAC 2007.01		
Aflatoxin B2 [±]	< LOQ		µg/kg	5.00	2501729	03/11/25 Mycotoxins by AOAC 2007.01		
Aflatoxin G1 [±]	< LOQ		µg/kg	5.00	2501729	03/11/25 Mycotoxins by AOAC 2007.01		
Aflatoxin G2 [±]	< LOQ		µg/kg	5.00	2501729	03/11/25 Mycotoxins by AOAC 2007.01		
Ochratoxin A	< LOQ	20.0	µg/kg	5.00	2501729	03/11/25 Mycotoxins by AOAC 2007.01 ^Þ	pass	
Total Aflatoxins	< LOQ	20.0	µg/kg	20.0		03/11/25 Mycotoxins by AOAC 2007.01 ^Þ	pass	

**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.

▷ = ISO/IEC 17025:2017 accredited method.

⊥ = TNI accredited analyte.

Units of Measure

cfu/g = Colony forming units per gram

g = Gram

g/ml = Gram per milliliter

µg/g = Microgram per gram

µg/kg = Micrograms per kilogram = parts per billion (ppb)

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

mg/355ml = Milligram per 355ml

% = Percentage of sample

% wt = µg/g divided by 10,000



Residue List

Method AOAC 2007.01

Units mg/kg

Analyzed 3/10/25

Parameter	LOQ	Parameter	LOQ	Parameter	LOQ	Parameter	LOQ
2,4-D	0.10	2,4-DB	0.10	2,4-DP	0.10	2,4,5-T	0.10
2,4,5-TP	0.10	Abamectin (Avermectin)	0.10	Acephate	0.20	Acequinocyl	0.10
Acetamiprid	0.10	Acetochlor	0.20	Acibenzolar-s-methyl	0.10	Acifluorfen	0.10
Acrinathrin	0.10	Afidopyropen	0.10	Alachlor	0.20	Aldicarb	0.10
Aldicarb-sulfone	0.10	Aldicarb-sulfoxide	0.10	Aldrin	0.10	Ametoctradin	0.10
Ametryn	0.10	Aminocyclopyrachlor	0.10	Anilazine	0.30	Aspon	0.10
Asulam	0.10	Atrazine	0.10	Atrazine-desethyl	0.10	Azadirachtin	0.10
Azinphos-ethyl	0.10	Azinphos-methyl	0.10	Azoxystrobin	0.10	Benalaxyl	0.10
Bendiocarb	0.10	Benfluralin	0.10	Benoxacor	0.10	Bensulide	0.10
Bentazone	0.10	Benzovindiflupyr	0.10	BHC (α , β , γ , δ isomers)	0.10	Bifenazate	0.10
Bifenox	0.10	Bifenthrin	0.10	Binapacryl	0.40	Bioresmethrin	0.10
Bitertanol	0.20	Boscalid	0.10	Broflanilide	0.10	Bromacil	0.20
Bromophos-ethyl	0.20	Bromophos-methyl	0.10	Bromopropylate	0.10	Bromoxynil	0.10
Bromuconazole	0.10	Bupirimate	0.10	Buprofezin	0.10	Butachlor	0.10
Butoxycarboxim	0.10	Butralin	0.20	Butylate	0.10	Cadusafos	0.10
Captafol	1.00	Captan	0.20	Carbaryl	0.10	Carbendazim	0.10
Carbofuran	0.10	Carbofuran-3-hydroxy	0.10	Carbophenothion	0.10	Carbophenothion-methyl	0.10
Carboxin	0.10	Carfentrazone-ethyl	0.10	Chlorantraniliprole	0.10	Chlordane	0.10
Chlordimeform	0.10	Chlorfenapyr	0.20	Chlorfenson	0.10	Chlorfenvinphos	0.10
Chlorimuron-ethyl	0.10	Chlormitrofen	0.20	Chlorobenzilate	0.10	Chloroneb	0.10
Chlorothalonil	0.40	Chlorpropham (CIPC)	0.10	Chlorpyrifos-ethyl	0.10	Chlorpyrifos-methyl	0.10
Chlorsulfuron	0.10	Chlorthal-dimethyl (Dacthal, D)	0.10	Chlorthion	0.20	Chlorthiophos	0.10
Cinerin I	0.10	Clethodim	0.10	Clethodim-sulfone	0.10	Clethodim-sulfoxide	0.10
Clofentezine	0.10	Clomazone	0.10	Clopyralid	0.10	Clothianidin	0.10
Coumaphos	0.10	Crotoxyphos	0.10	Cyanazine	0.10	Cyanofenphos	0.10
Cyanophos	0.40	Cyantraniliprole	0.10	Cyazofamid	0.10	Cycloate	0.10
Cycloxydim	0.10	Cyflufenamid	0.10	Cyflumetofen	0.10	Cyfluthrin (incl. Beta-Cyfluthrin	0.20
Cyhalothrin, lambda	0.10	Cymoxanil	0.10	Cypermethrin and isomers (su	0.10	Cyprodinil	0.10
Cyromazine	0.10	DDD-o,p'	0.10	DDD-p,p'	0.10	DDE-o,p'	0.10
DDE-p,p'	0.10	DDT-o,p'	0.10	DDT-p,p'	0.10	DEF (Tribufos)	0.10
Deltamethrin	0.10	Demeton	0.20	Demeton-s-methyl	0.20	Demeton-s-methyl sulfone	0.20
Desmedipham	0.10	Diallate	0.10	Diazinon	0.10	Diazoxon (Diazinon OA)	0.10
Dicamba	0.10	Dichlobenil	0.10	Dichlofenthion	0.10	Dichlofluanid	0.10
Dichlorbenzamide	0.10	Dichlorvos	0.10	Diclobutrazol	0.10	Diclofop	0.10



Residue List

Method AOAC 2007.01

Units mg/kg

Analyzed 3/10/25

Parameter	LOQ	Parameter	LOQ	Parameter	LOQ	Parameter	LOQ
Diclofop-methyl	0.10	Dicloran	0.40	Dicofol o,p	0.20	Dicofol-p,p	0.20
Dicrotophos	0.10	Dieldrin	0.10	Diethofencarb	0.10	Diethyltoluamide (DEET)	0.10
Difenoconazole	0.10	Diffubenzuron	0.10	Diffufenzopyr	0.10	Dimethenamid	0.10
Dimethoate	0.10	Dimethomorph	0.10	Diniconazole	0.10	Dinocap	0.10
Dinoseb	0.10	Dinotefuran	0.10	Dioxathion	0.10	Diphenamid	0.10
Diphenylamine	0.10	Disulfoton	0.20	Disulfoton-sulfone	0.10	Disulfoton-sulfoxide	0.10
Dithianon	0.10	Dithiopyr	0.10	Diuron	0.10	Diuron metabolite (DCPMU)	0.10
DNOC (Dinitrocresol)	0.10	Edifenphos	0.10	Endosulfan I (alpha)	0.20	Endosulfan II (beta)	0.20
Endosulfan sulfate	0.10	Endrin	0.20	Endrin Aldehyde	0.20	EPN	0.10
EPTC	0.10	Esfenvalerate	0.20	Etaconazole	0.10	Ethaboxam	0.10
Ethalfuralin	0.10	Ethiofencarb	0.10	Ethion	0.10	Ethirimol	0.10
Ethofumesate	0.10	Ethoprophos	0.10	Ethoxyquin	0.20	Etofenprox	0.10
Etoxazole	0.10	Etridiazole	0.10	Etrimfos	0.10	Famoxadone	0.20
Famphur	0.10	Fenamidone	0.10	Fenamiphos	0.10	Fenamiphos-sulfone	0.10
Fenamiphos-sulfoxide	0.10	Fenarimol	0.10	Fenazaquin	0.10	Fenbuconazole	0.10
Fenbutatin oxide	0.10	Fenchlorphos	0.10	Fenchlorphos-oxon	0.10	Fenhexamid	0.10
Fenitrothion	0.10	Fenobucarb	0.10	Fenoxaprop-p-ethyl	0.10	Fenoxycarb	0.10
Fenpropathrin	0.10	Fenpyroximate	0.10	Fenson	0.20	Fensulfthion	0.10
Fenthion	0.10	Fenuron	0.10	Fipronil	0.10	Fonicamid	0.10
Fluazifop	0.10	Fluazinam	0.10	Fluchloralin	0.10	Flucythrinate	0.30
Fludioxonil	0.10	Flufenacet	0.10	Flumetsulam	0.10	Flumioxazin	0.10
Fluometuron	0.10	Fluopicolide	0.10	Fluopyram	0.10	Fluoxastrobin	0.10
Fluprimidol	0.10	Flupyradifurone	0.10	Fluridone	0.10	Fluroxypyr	0.10
Flusilazole	0.10	Fluthiacet-methyl	0.10	Flutianil	0.10	Flutolanil	0.10
Flutriafol	0.10	Fluxapyroxad	0.10	Folpet	0.10	Fomesafen	0.10
Fonofos	0.10	Foramsulfuron	0.10	Forchlorfenuron	0.10	Formetanate	0.10
Furathiocarb	0.10	Halosulfuron-methyl	0.10	Haloxypop	0.10	Heptachlor	0.10
Heptachlor epoxide	0.10	Hexachlorobenzene	0.10	Hexaconazole	0.10	Hexazinone	0.10
Hexythiazox	0.10	Hydroprene	0.10	Imazalil	0.10	Imazamox	0.10
Imazapic	0.10	Imazapyr	0.10	Imazaquin	0.10	Imazethapyr	0.10
Imidacloprid	0.10	Indaziflam	0.10	Indoxacarb	0.10	Iprobenfos	0.10
Iprodione	0.10	Isazophos	0.10	Isobenzan	0.10	Isocarbophos	0.10
Isodrin	0.10	Isofenphos	0.10	Isofenphos-methyl	0.10	Isofenphos-oxon	0.10
Isofetamid	0.10	Isoprocab	0.10	Isopropalin	0.10	Isoprothiolane	0.10



Residue List

Method AOAC 2007.01

Units mg/kg

Analyzed 3/10/25

Parameter	LOQ	Parameter	LOQ	Parameter	LOQ	Parameter	LOQ
Isoproturon	0.10	Isoxaben	0.10	Isoxaflutole	0.10	Jasmodin I	0.10
Kresoxim-methyl	0.10	Lactofen	0.20	Lenacil	0.10	Linuron	0.10
Malaaxon	0.10	Malathion	0.10	Mandestrobin	0.10	Mandipropamid	0.10
MCPA	0.10	MCPB	0.10	MCPP (Mecoprop)	0.10	MCPP-P	0.10
Mecarbam	0.10	Mefentrifluconazole	0.10	Mepanipyrim	0.10	Mesosulfuron-methyl	0.10
Mesotrione	0.10	Metaxyl	0.10	Metalddehyde	0.10	Metconazole	0.10
Methacrifos	0.10	Methamidophos	0.10	Methidathion	0.10	Methiocarb	0.10
Methiocarb-sulfone	0.10	Methiocarb-sulfoxide	0.10	Methiozolin	0.10	Methomyl	0.10
Methoxychlor	0.10	Methoxyfenozide	0.10	Metobromuron	0.10	Metolacarb	0.10
Metolachlor	0.10	Metrafenone	0.10	Metribuzin	0.10	Metsulfuron-methyl	0.10
Mevinphos	0.10	Mexacarbate	0.10	MGK-264	0.10	Mirex	0.10
Molinate	0.10	Monocrotophos	0.10	Monolinuron	0.10	Myclobutanil	0.10
Naled	0.10	Napropamide	0.10	Neburon	0.10	Nicosulfuron	0.10
Nitrapyrin	0.10	Nitrofen	0.20	Norflurazon	0.10	Novaluron	0.10
Nuarimol	0.20	O-Phenylphenol	0.50	Omethoate	0.10	Oryzalin	0.10
Oxadiazon	0.10	Oxadixyl	0.10	Oxamyl	0.10	Oxamyl-oxime	0.10
Oxathiapiprolin	0.10	Oxychlorane	0.10	Oxydemeton-methyl	0.10	Oxyfluorfen	0.10
Oxythioquinox	0.20	Paclobutrazole	0.10	Paraaxon-ethyl	0.10	Paraaxon-methyl	0.10
Parathion-ethyl	0.10	Parathion-methyl	0.30	Penconazole	0.10	Pendimethalin	0.10
Penflufen	0.10	Pentachloroaniline	0.10	Pentachloroanisole	0.10	Pentachlorobenzene	0.10
Pentachlorophenol	0.10	Pentachloroethoxyanisole	0.30	Penthiopyrad	0.10	Permethrin	0.10
Perthane	0.10	Phenmedipham	0.10	Phenothrin	0.10	Phenthoate	0.10
Phorate	0.10	Phorate OA	0.10	Phorate-sulfone	0.10	Phorate-sulfoxide	0.10
Phosalone	0.10	Phosmet	0.10	Phosmet oxon	0.10	Phosphamidon	0.10
Phoxim	0.10	Picloram	0.10	Pinoxaden	0.10	Piperonyl butoxide	0.10
Pirimicarb	0.10	Pirimiphos-ethyl	0.10	Pirimiphos-methyl	0.10	Prallethrin	0.10
Primisulfuron-methyl	0.10	Prochloraz	0.10	Procymidone	0.10	Prodiamine	0.10
Profenofos	0.10	Profluralin	0.10	Promecarb	0.10	Prometon	0.10
Prometryn	0.10	Pronamide (Propyzamid)	0.10	Propachlor	0.10	Propamocarb	0.10
Propanil	0.10	Propargite	0.10	Propazine	0.10	Propetamphos	0.10
Propham	0.10	Propiconazole	0.10	Propoxur	0.10	Propoxycarbazon sodium	0.10
Prosulfuron	0.10	Prothioconazole	0.10	Prothiofos	0.10	Pydiflumetofen	0.10
Pymetrozine	0.10	Pyraclostrobin	0.10	Pyraflufen-ethyl	0.10	Pyrazophos	0.10
Pyrethrins (total)	0.10	Pyridaben	0.10	Pyridate	0.10	Pyrifluquinazon	0.10



Residue List

Method AOAC 2007.01

Units mg/kg

Analyzed 3/10/25

Parameter	LOQ	Parameter	LOQ	Parameter	LOQ	Parameter	LOQ
Pyrimethanil	0.10	Pyriproxyfen	0.10	Pyroxasulfone	0.10	Pyroxulam	0.10
Quinalphos	0.10	Quinclorac	0.10	Quinoxifen	0.10	Quintozene (PCNB)	0.10
Quizalofop	0.10	Resmethrin	0.10	Rimsulfuron	0.10	Rotenone	0.10
S-421	0.10	Saflufenacil	0.10	Sebuthylazine	0.10	Sedaxane	0.10
Sethoxydim	0.10	Siduron	0.10	Simazine	0.10	Simetryn	0.10
Spinetoram	0.10	Spinosad	0.10	Spirodiclofen	0.10	Spiromesifen	0.10
Spirotetramat	0.10	Spirotetramat enol	0.10	Spiroxamine	0.10	Sulfallate	0.10
Sulfentrazone	0.30	Sulfometuron-methyl	0.10	Sulfosulfuron	0.10	Sulfotep	0.10
Sulfoxaflor	0.10	Sulprofos	0.10	Tau-fluvalinate	0.10	Tebuconazole	0.10
Tebufenozide	0.10	Tebuthiuron	0.10	Tecnazene	0.10	Tefluthrin	0.10
Tembotrione	0.10	Terbacil	0.40	Terbufos	0.10	Terbufos-sulfone	0.10
Terbufos-sulfoxide	0.10	Terbuthylazine	0.10	Terbutryn	0.10	Tetrachlorvinphos	0.10
Tetraconazole	0.10	Tetradifon	0.10	Tetramethrin	0.10	Tetrasul	0.10
Thiabendazol 5 hydroxy	0.10	Thiabendazole	0.10	Thiacloprid	0.10	Thiamethoxam	0.10
Thifensulfuron-methyl	0.10	Thiobencarb	0.10	Thiodicarb	0.10	Thiometon	0.20
Thionazin	0.10	Thiophanate-methyl	0.10	Tolclofos-methyl	0.10	Tolfenpyrad	0.10
Tolyfluanid	0.10	Topramezone	0.10	Tralkoxydim	0.10	Triadimefon	0.10
Triadimenol	0.10	Triallate	0.10	Triasulfuron	0.10	Triazophos	0.10
Tribenuron-methyl	0.10	Trichlorfon (Metrifonate)	0.10	Triclopyr	0.20	Trifloxystrobin	0.10
Trifloxysulfuron	0.10	Triflumizole	0.10	Trifluralin	0.10	Triflurosulfuron-methyl	0.10
Triforine	0.10	Trinexapac-ethyl	0.10	Triticonazole	0.10	Vinclozolin	0.10
Zoxamide	0.10						



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



Report Number: 25-002244/D001.R000
Report Date: 03/11/2025
ORELAP#: OR100028
Purchase Order:
Received: 03/04/25 13:55



Hemp & Cannabis
Chain of Custody

Northwest-Natural-
Goods-1741030220

Company Details Company: <u>Northwest Natural Goods</u> [Redacted] [Redacted] [Redacted] [Redacted] [Redacted] [Redacted] Billing Information Billing Email: [Redacted]			Project Details Turnaround Time: <u>5 Business Days</u> Req. For Micro Testing Standard Relinquishment Sampling, Courier & Shipping Options: <u>Pick-Up Courier Service</u> Project Name / ID: <u>BEV - BB - B-58-1</u> Pick-Up Details Pick-Up Location Name: <u>Northwest Natural Goods</u> [Redacted] [Redacted] [Redacted] Receipt Information Evidence of Cooling?: No Sample Condition: Satisfactory Prelog Storage: Canna Shelves			Testing					
						H0013 - Cannabis Heavy Metals Pro le OR	H0010 - Potency Cannabis (Basic+Expanded)	H0008 - Residual Solvents (Cannabis - Oregon)	H0042S - A atoxine+Ochratoxin CLCC	M010 - Micro Pro le D	P2320 - Multi-Residue Pesticide Pro le (Cannabis)
#	Sample Name	Material	Amount Provided	Reporting Unit	Serving Size						
1	BEV - BB - B-58-1	Cannabinoid Beverage	4 each	mg/g & mg/serving	362.1g	✓	✓	✓	✓	✓	✓

Relinquished By	Date	Time	Temp., °C	Received By	Date	Time	Received Temp., °C	IRTherm. CL#
Kristen Johnson	03/03/2025	11:30		BCR	03/04/2025	11:41	NA	NA
BCR	03/04/2025	11:21	16.9	new	03/04/2025	13:55	17.4	CL-1196

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
Portland, OR 97230

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Page 1 of 1
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Laboratory Quality Control Results

JAOAC2015 V986 Batch ID: 2501615

Laboratory Control Sample

Analyte	LCS	Result	Spike	Units	% Rec	Limits	Evaluation	Notes
CBDVA	2	0.0009	0.0010	%	89.5	80.0 - 120	Acceptable	
CBDV	2	0.00103	0.00107	%	96.5	80.0 - 120	Acceptable	
CBE	2	0.00104	0.00109	%	94.8	80.0 - 120	Acceptable	
CBDA	1	0.00107	0.00106	%	101	90.0 - 110	Acceptable	
CBGA	1	0.00102	0.00106	%	96.6	80.0 - 120	Acceptable	
CBG	1	0.0010	0.00101	%	98.2	80.0 - 120	Acceptable	
CBD	1	0.0009	0.0009	%	106	90.0 - 110	Acceptable	
THCV	2	0.00104	0.00107	%	97.4	80.0 - 120	Acceptable	
d8THCV	2	0.00105	0.00111	%	94.5	80.0 - 120	Acceptable	
THCVA	2	0.0009	0.0010	%	91.1	80.0 - 120	Acceptable	
CBN	1	0.00106	0.00101	%	104	80.0 - 120	Acceptable	
exo-THC	2	0.0010	0.00101	%	97.2	80.0 - 120	Acceptable	
d9THC	1	0.0010	0.0009	%	108	90.0 - 110	Acceptable	
d8THC	1	0.00107	0.00106	%	101	90.0 - 110	Acceptable	
9S-d10THC	1	0.00109	0.00107	%	102	80.0 - 120	Acceptable	
CBL	2	0.00100	0.00104	%	96.1	80.0 - 120	Acceptable	
9R-d10THC	1	0.0010	0.0009	%	104	80.0 - 120	Acceptable	
CBC	2	0.00102	0.00105	%	97.3	80.0 - 120	Acceptable	
THCA	1	0.00109	0.00113	%	96.8	90.0 - 110	Acceptable	
CBCA	2	0.0009	0.00103	%	86.6	80.0 - 120	Acceptable	
CBLA	2	0.0009	0.0010	%	90.4	80.0 - 120	Acceptable	
d9THCP	2	0.0010	0.00102	%	98.1	80.0 - 120	Acceptable	
CBT	2	0.00100	0.00104	%	96.1	80.0 - 120	Acceptable	

Method Blank

Analyte	Result	LOQ	Units	Limits	Evaluation	Notes
CBDVA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBDV	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBE	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBDA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBGA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBG	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBD	<LOQ	0.0001	%	< 0.0001	Acceptable	
THCV	<LOQ	0.0001	%	< 0.0001	Acceptable	
d8THCV	<LOQ	0.0001	%	< 0.0001	Acceptable	
THCVA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBN	<LOQ	0.0001	%	< 0.0001	Acceptable	
exo-THC	<LOQ	0.0001	%	< 0.0001	Acceptable	
d9THC	<LOQ	0.0001	%	< 0.0001	Acceptable	
d8THC	<LOQ	0.0001	%	< 0.0001	Acceptable	
9S-d10THC	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBL	<LOQ	0.0001	%	< 0.0001	Acceptable	
9R-d10THC	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBC	<LOQ	0.0001	%	< 0.0001	Acceptable	
THCA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBCA	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBLA	<LOQ	0.0001	%	< 0.0001	Acceptable	
d9THCP	<LOQ	0.0001	%	< 0.0001	Acceptable	
CBT	<LOQ	0.0001	%	< 0.0001	Acceptable	

Abbreviations

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

Units of Measure:

% - Percent



Laboratory Quality Control Results

JAOAC2015 V986			Batch ID: 2501615					
Sample Duplicate			Sample ID: 25-0021770001					
Analyte	Result	Org. Result	LOQ	Units	RPD	Limits	Evaluation	Notes
CBDVA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBDV	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBE	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBDA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBGA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBG	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBD	0.00682	0.00683	0.0001	%	0.122	< 20	Acceptable	
THCV	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
d8THCV	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
THCVA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBN	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
exo-THC	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
d9THC	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
d8THC	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
9S-d10THC	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBL	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
9R-d10THC	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBC	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
THCA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBCA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBLA	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
d9THCP	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	
CBT	<LOQ	<LOQ	0.0001	%	NA	< 20	Acceptable	

Abbreviations

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

Units of Measure:

% - Percent


Laboratory Quality Control Results

Residual Solvents				Batch ID: 2501655			
Method Blank				Laboratory Control Sample			
Analyte	Result	LOQ	Notes	Result	Spike	Units	% Rec Limits Notes
Propane	ND	< 200		613	585	µg/g	104.8 60 - 120
Isobutane	ND	< 200		802	770	µg/g	104.2 60 - 120
Butane	ND	< 200		825	769	µg/g	107.3 60 - 120
2,2-Dimethylpropane	ND	< 200		1040	956	µg/g	108.8 60 - 120
Methanol	ND	< 200		2090	1650	µg/g	126.7 60 - 120 Q1
Ethylene Oxide	ND	< 30		61.9	57.7	µg/g	107.3 60 - 120
2-Methylbutane	ND	< 200		2020	1620	µg/g	124.7 60 - 120 Q1
Pertane	ND	< 200		2020	1620	µg/g	124.7 60 - 120 Q1
Ethanol	ND	< 200		2070	1680	µg/g	123.2 70 - 130
Ethyl Ether	ND	< 200		1890	1640	µg/g	115.2 60 - 120
2,2-Dimethylbutane	ND	< 30		222	182	µg/g	122.0 60 - 120 Q1
Acetone	ND	< 200		1810	1640	µg/g	110.4 60 - 120
2-Propanol	ND	< 200		1900	1630	µg/g	116.6 60 - 120
Ethyl Formate	ND	< 500		1240	1610	µg/g	77.0 70 - 130
Acetonitrile	ND	< 100		667	513	µg/g	130.0 60 - 120 Q1
Methyl Acetate	ND	< 500		1710	1610	µg/g	106.2 70 - 130
2,3-Dimethylbutane	ND	< 30		217	184	µg/g	117.9 60 - 120
Dichloromethane	ND	< 60		648	589	µg/g	110.0 60 - 120
2-Methylpentane	ND	< 30		238	185	µg/g	128.6 60 - 120 Q1
MTBE	ND	< 500		1710	1610	µg/g	106.2 70 - 130
3-Methylpentane	ND	< 30		202	177	µg/g	114.1 60 - 120
Hexane	ND	< 30		201	174	µg/g	115.5 60 - 120
1-Propanol	ND	< 500		1730	1600	µg/g	108.1 70 - 130
Methylethylketone	ND	< 500		1750	1610	µg/g	108.7 70 - 130
Ethyl acetate	ND	< 200		1900	1620	µg/g	117.3 60 - 120
2-Butanol	ND	< 200		1830	1640	µg/g	111.6 60 - 120
Tetrahydrofuran	ND	< 100		530	511	µg/g	103.7 60 - 120
Cyclohexane	ND	< 200		1600	1610	µg/g	99.4 60 - 120
2-methyl-1-propanol	ND	< 500		1590	1610	µg/g	98.8 70 - 130
Benzene	ND	< 1		5.14	5.56	µg/g	92.4 60 - 120
Isopropyl Acetate	ND	< 200		1690	1620	µg/g	104.3 60 - 120
Heptane	ND	< 200		1610	1610	µg/g	100.0 60 - 120
1-Butanol	ND	< 500		1560	1610	µg/g	96.9 70 - 130
Propyl Acetate	ND	< 500		1630	1620	µg/g	100.6 70 - 130
1,4-Dioxane	ND	< 100		446	487	µg/g	91.6 60 - 120
2-Ethoxyethanol	ND	< 30		182	180	µg/g	101.1 60 - 120
Methylisobutylketone	ND	< 500		1550	1620	µg/g	95.7 70 - 130
3-Methyl-1-butanol	ND	< 500		1480	1600	µg/g	92.5 70 - 130
Ethylene Glycol	ND	< 200		442	530	µg/g	83.4 60 - 120
Toluene	ND	< 100		452	497	µg/g	90.9 60 - 120
Isobutyl Acetate	ND	< 500		1570	1620	µg/g	96.9 70 - 130
1-Pentanol	ND	< 500		1580	1600	µg/g	98.8 70 - 130
Butyl Acetate	ND	< 500		1610	1600	µg/g	100.6 70 - 130
Ethylbenzene	ND	< 200		975	1060	µg/g	92.0 60 - 120
m,p-Xylene	ND	< 200		972	1040	µg/g	93.5 60 - 120
o-Xylene	ND	< 200		922	1090	µg/g	84.6 60 - 120
Cumene	ND	< 30		180	227	µg/g	79.3 60 - 120
Anisole	ND	< 500		1430	1610	µg/g	88.8 70 - 130
DMSO	ND	< 500		1390	1600	µg/g	86.9 70 - 130
1,2-dimethoxyethane	ND	< 50		164	162	µg/g	101.2 70 - 130
Triethylamine	ND	< 500		1610	1600	µg/g	100.6 70 - 130
N,N-dimethylformamide	ND	< 150		474	487	µg/g	97.3 70 - 130
N,N-dimethylacetamide	ND	< 150		458	498	µg/g	92.0 70 - 130
Pyridine	ND	< 50		156	162	µg/g	96.3 70 - 130
Sulfolane	ND	< 50		124	173	µg/g	71.7 70 - 130
1,2-Dichloroethane	ND	< 1		1.01	1	µg/g	101.0 70 - 130
Chloroform	ND	< 1		1.02	1	µg/g	102.0 70 - 130
Trichloroethylene	ND	< 1		0.891	1	µg/g	89.1 70 - 130
1,1-Dichloroethane	ND	< 1		1.22	1	µg/g	122.0 70 - 130



QC- Sample Duplicate

Sample ID: 25-002173-0001

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	
Pertane	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	
Sulfone	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



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



Report Number: 25-002244/D001.R000
Report Date: 03/11/2025
ORELAP#: OR100028
Purchase Order:
Received: 03/04/25 13:55





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.