



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



**Report Number:** 23-000929/D008.R000  
**Report Date:** 01/30/2023  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 01/23/23 11:13

**Customer:** NW Natural Goods  
**Product identity:** Bev - LM 022335 11-1  
**Client/Metric ID:** .  
**Laboratory ID:** 23-000929-0002

### Summary

**Potency:**

| Analyte per 355ml | Result | Limits | Units    | Status |  |
|-------------------|--------|--------|----------|--------|--|
| CBD per 355ml     | 27.1   |        | mg/355ml |        | CBD-Total per Serving Size 27.1 mg/355ml |
| CBG per 355ml     | 0.753  |        | mg/355ml |        |  |
|                   |        |        |          |        | THC-Total per Serving Size <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.*



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**Product identity:** Bev - LM 022335 11-1  
**Client/Metric ID:** .  
**Sample Date:**  
**Laboratory ID:** 23-000929-0002  
**Evidence of Cooling:** No  
**Temp:** 18 °C  
**Relinquished by:** Hinton  
**Serving Size #1:** 362.1 g  
**Density:** 1.020 g/ml

### Sample Results

| Potency per 355ml            | Method: J AOAC 2015 V98-6 (mod) <sup>b</sup> | Units mg/se | Batch: 2300772 | Analyze: 1/24/23 11:36:00 PM |       |
|------------------------------|--|-------------|----------------|------------------------------|-------|
| Analyte                      | Result                                       | Limits      | Units          | LOQ                          | Notes |
| CBC per 355ml                | < LOQ  |             | mg/355ml       | 0.362                        |       |
| CBC-A per 355ml              | < LOQ  |             | mg/355ml       | 0.362                        |       |
| CBC-Total per 355ml          | < LOQ  |             | mg/355ml       | 0.679                        |       |
| CBD per 355ml                | 27.1   |             | mg/355ml       | 0.362                        |       |
| CBD-A per 355ml              | < LOQ  |             | mg/355ml       | 0.362                        |       |
| CBD-Total per 355ml          | 27.1   |             | mg/355ml       | 0.679                        |       |
| CBDV per 355ml               | < LOQ  |             | mg/355ml       | 0.362                        |       |
| CBDV-A per 355ml             | < LOQ  |             | mg/355ml       | 0.362                        |       |
| CBDV-Total per 355ml         | < LOQ  |             | mg/355ml       | 0.675                        |       |
| CBE per 355ml                | < LOQ  |             | mg/355ml       | 0.362                        |       |
| CBG per 355ml                | 0.753  |             | mg/355ml       | 0.362                        |       |
| CBG-A per 355ml              | < LOQ  |             | mg/355ml       | 0.362                        |       |
| CBG-Total per 355ml          | 0.753  |             | mg/355ml       | 0.675                        |       |
| CBL per 355ml                | < LOQ  |             | mg/355ml       | 0.362                        |       |
| CBL-A per 355ml              | < LOQ  |             | mg/355ml       | 0.362                        |       |
| CBL-Total per 355ml          | < LOQ  |             | mg/355ml       | 0.679                        |       |
| CBN per 355ml                | < LOQ  |             | mg/355ml       | 0.362                        |       |
| CBT per 355ml                | < LOQ  |             | mg/355ml       | 0.362                        |       |
| Δ8-THCV per 355ml            | < LOQ  |             | mg/355ml       | 0.362                        |       |
| Δ10-THC-9R per 355ml         | < LOQ  |             | mg/355ml       | 0.362                        |       |
| Δ10-THC-9S per 355ml         | < LOQ  |             | mg/355ml       | 0.362                        |       |
| Δ8-THC per 355ml             | < LOQ  |             | mg/355ml       | 0.362                        |       |
| Δ9-THC per 355ml             | < LOQ  |             | mg/355ml       | 0.362                        |       |
| exo-THC per 355ml            | < LOQ  |             | mg/355ml       | 0.362                        |       |
| THC-A per 355ml              | < LOQ  |             | mg/355ml       | 0.362                        |       |
| THC-Total per 355ml          | < LOQ  |             | mg/355ml       | 0.679                        |       |
| THCV per 355ml               | < LOQ  |             | mg/355ml       | 0.362                        |       |
| THCV-A per 355ml             | < LOQ  |             | mg/355ml       | 0.362                        |       |
| THCV-Total per 355ml         | < LOQ  |             | mg/355ml       | 0.679                        |       |
| Total Cannabinoids per 355ml | 28.0   |             | mg/355ml       |                              |       |



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**Microbiology**

| Analyte                 | Result | Limits | Units | LOQ | Batch   | Analyzed Method                               | Status | Notes |
|-------------------------|--------|--------|-------|-----|---------|---|--------|-------|
| E.coli                  | < LOQ  |        | cfu/g | 10  | 2300700 | 01/26/23 AOAC 991.14 (Petrifilm) <sup>®</sup> |        |       |
| Total Coliforms         | < LOQ  |        | cfu/g | 10  | 2300700 | 01/26/23 AOAC 991.14 (Petrifilm) <sup>®</sup> |        |       |
| Mold (RAPID Petrifilm)  | < LOQ  |        | cfu/g | 10  | 2300701 | 01/27/23 AOAC 2014.05 (RAPID) <sup>®</sup>    |        |       |
| Yeast (RAPID Petrifilm) | < LOQ  |        | cfu/g | 10  | 2300701 | 01/27/23 AOAC 2014.05 (RAPID) <sup>®</sup>    |        |       |

**Solvents Method: Residual Solvents by GC/MS<sup>®</sup> Units µg/g Batch 2300804 Analyze 01/26/23 11:39 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-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)<sup>®</sup> Units mg/kg Batch 2300812 Analyze 01/26/23 02:23 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 | 2300864 | 01/27/23 AOAC 2013.06 (mod.) <sup>®</sup> | pass   |       |
| Cadmium | < LOQ  | 0.200  | mg/kg | 0.00390 | 2300864 | 01/27/23 AOAC 2013.06 (mod.) <sup>®</sup> | pass   |       |
| Lead    | < LOQ  | 0.500  | mg/kg | 0.00390 | 2300864 | 01/27/23 AOAC 2013.06 (mod.) <sup>®</sup> | pass   |       |
| Mercury | < LOQ  | 0.100  | mg/kg | 0.00195 | 2300864 | 01/27/23 AOAC 2013.06 (mod.) <sup>®</sup> | pass   |       |



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**Mycotoxins**

| Analyte                       | Result | Limits | Units | LOQ  | Batch   | Analyzed Method                                     | Status | Notes |
|-------------------------------|--------|--------|-------|------|---------|---|--------|-------|
| Aflatoxin B2 <sup>‡</sup>     | < LOQ  |        | µg/kg | 5.00 | 2300825 | 01/27/23 AOAC 2007.01 & EN 15662 (mod) <sup>‡</sup> |        |       |
| Aflatoxin B1 <sup>‡</sup>     | < LOQ  |        | µg/kg | 5.00 | 2300825 | 01/27/23 AOAC 2007.01 & EN 15662 (mod) <sup>‡</sup> |        |       |
| Aflatoxin G1 <sup>‡</sup>     | < LOQ  |        | µg/kg | 5.00 | 2300825 | 01/27/23 AOAC 2007.01 & EN 15662 (mod) <sup>‡</sup> |        |       |
| Aflatoxin G2 <sup>‡</sup>     | < LOQ  |        | µg/kg | 5.00 | 2300825 | 01/27/23 AOAC 2007.01 & EN 15662 (mod) <sup>‡</sup> |        |       |
| Ochratoxin A <sup>‡</sup>     | < LOQ  | 20.0   | µg/kg | 5.00 | 2300825 | 01/27/23 AOAC 2007.01 & EN 15662 (mod) <sup>‡</sup> | pass   |       |
| Total Aflatoxins <sup>‡</sup> | 0.000  | 20.0   | µg/kg | 20.0 |         | 01/30/23 AOAC 2007.01 & EN 15662 (mod) <sup>‡</sup> | pass   |       |

**Nutrition**

| Analyte                   | Result | Limits | Units  | LOQ   | Batch   | Analyzed Method                          | Status | Notes |
|---------------------------|--------|--------|--------|-------|---------|--|--------|-------|
| Moisture (Loss on Drying) | 99.8   |        | g/100g | 0.10  | 2300809 | 01/25/23 AOAC 925.10 (mod.) <sup>‡</sup> |        |       |
| Water Activity            | 0.990  |        | Aw     | 0.030 | 2300763 | 01/25/23 AOAC 978.18 <sup>‡</sup>        |        |       |



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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.

Ⓟ = ISO/IEC 17025:2017 accredited method.

\* = TNI accredited analyte.

**Units of Measure**

cfu/g = Colony forming units per gram

g = g

g/ml = Gram per milliliter

g/100g = Grams per 100 Grams

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

Aw = Water Activity

% wt = µg/g divided by 10,000

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       | Clethodim               | 0.050       | Endrin                     | 0.100       |
| Acephate                      | 0.100       | Clethodim Sulfone       | 0.050       | EPN                        | 0.050       |
| Acequinocyl                   | 0.100       | Clethodim Sulfoxide     | 0.050       | EPTC                       | 0.100       |
| Acetamiprid                   | 0.020       | Clofentezine            | 0.020       | Esfenvalerate/Fenvalerate  | 0.200       |
| Acetochlor                    | 0.100       | Clomazone               | 0.020       | Etaconazole                | 0.100       |
| Acrinathrin                   | 0.100       | Clothianidin            | 0.200       | Ethalfuralin               | 0.100       |
| Alachlor                      | 0.100       | Coumaphos               | 0.050       | Ethiofencarb               | 0.050       |
| Aldicarb                      | 0.100       | Crotoxypfos             | 0.020       | Ethion                     | 0.200       |
| Aldicarb sulfoxide            | 0.100       | Cyanazine               | 0.020       | Ethirimol                  | 0.100       |
| Aldoxycarb (Aldicarb-sulfone) | 0.100       | Cyanofenphos            | 0.020       | Ethofumesate               | 0.050       |
| Aldrin                        | 0.100       | Cyantranilprole         | 0.050       | Ethoprophos                | 0.020       |
| Ametocrtadin                  | 0.020       | Cyazofamid              | 0.020       | Etofenprox                 | 0.020       |
| Ametryn                       | 0.500       | Cycloate                | 0.100       | Etozazole                  | 0.020       |
| Aspon                         | 0.100       | Cyfluthrin              | 0.200       | Etridiazole                | 0.100       |
| Asulam                        | 0.100       | Cyhalothrin, lambda     | 0.200       | Etrimfos                   | 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       | Dacthal                 | 0.100       | Fenamiphos                 | 0.020       |
| Azoxystrobin                  | 0.020       | Daminozide              | 0.100       | Fenamiphos sulfone         | 0.020       |
| Benalaxyl                     | 0.020       | DCPMU                   | 0.050       | Fenamiphos sulfoxide       | 0.020       |
| Bendiocarb                    | 0.020       | DDD, o,p'-              | 0.100       | Fenazaquin                 | 0.100       |
| Benfluralin                   | 0.100       | DDD, p,p'-              | 0.100       | Fenbuconazole              | 0.100       |
| Benoxacor                     | 0.050       | DDE, o,p'-              | 0.100       | Fenchlorphos               | 0.100       |
| Bensulide                     | 0.050       | DDE, 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       | Deltamethrin            | 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       | Fensulfothion              | 0.020       |
| Bromopropylate                | 0.100       | Dichlobenil             | 0.100       | Fensulfothion oxon         | 0.020       |
| Bromuconazole                 | 0.100       | Dichlofluanid           | 0.100       | Fensulfothion sulfone      | 0.100       |
| Bupirimate                    | 0.020       | Dichlorvos              | 0.100       | Fensulfothion-oxon-sulfone | 0.020       |
| Buprofezin                    | 0.050       | Diclobutrazol           | 0.050       | Fenthion                   | 0.050       |
| Butachlor                     | 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       | Flonicamid                 | 0.100       |
| Carbendazim                   | 0.100       | Dimethenamid            | 0.050       | Fluchloralin               | 0.100       |
| Carbofuran                    | 0.020       | Dimethoate              | 0.050       | Flucythrinate              | 0.100       |
| Carbophenothion               | 0.200       | Dimethomorph            | 0.050       | Fludioxonil                | 0.200       |
| Carboxin                      | 0.020       | Diniconazole            | 0.200       | Flufenacet                 | 0.020       |
| Carfentrazone-ethyl           | 0.100       | Dinotefuran             | 0.200       | Flumioxazin                | 0.100       |
| Chlorantranilprole            | 0.020       | Dioxathion              | 0.100       | Fluometuron                | 0.020       |
| Chlordane, cis-               | 0.200       | Diphenamid              | 0.020       | Fluopicolide               | 0.050       |
| Chlordane, trans-             | 0.200       | Diphenylamine           | 0.100       | Fluopyram                  | 0.020       |
| Chlorfenapyr                  | 0.500       | Disulfoton              | 0.100       | Fluoxastrobin              | 0.050       |
| Chlorfenson                   | 0.200       | Disulfoton sulfone      | 0.100       | Flupyradifurone            | 0.020       |
| Chlorfenvinphos               | 0.050       | Disulfoton sulfoxide    | 0.100       | Fluridone                  | 0.100       |
| Chlorobenzilate               | 0.100       | Diuron                  | 0.050       | Flusilazole                | 0.020       |
| Chloroneb                     | 0.200       | Edifenphos              | 0.050       | Flutolanil                 | 0.020       |
| Chlorpyrifos                  | 0.050       | Endosulfan alpha        | 0.200       | Flutriafol                 | 0.020       |
| Chlorpyrifos-methyl           | 0.200       | Endosulfan beta         | 0.200       | Fluxalinate, tau-          | 0.100       |
| CIPC                          | 1.000       | Endosulfan sulfate      | 0.100       | Fluxapyroxad               | 0.020       |



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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       | Propamocarb              | 0.050       |
| Fonofos              | 0.100       | MGK 264                       | 0.020       | Propanil                 | 0.050       |
| Forchlorfenuron      | 0.050       | Mirex                         | 0.100       | Propargite               | 0.050       |
| Formetanate          | 0.050       | Molinate                      | 0.050       | Propazine                | 0.020       |
| Furathiocarb         | 0.020       | Monocrotophos                 | 0.100       | Propetamphos             | 0.050       |
| Heptachlor           | 0.100       | Monolinuron                   | 0.020       | Propham                  | 0.050       |
| Heptachlor epoxide   | 0.100       | Myclobutanil                  | 0.050       | Propiconazole            | 0.050       |
| Heptenophos          | 0.100       | Naled                         | 0.100       | Propoxur                 | 0.050       |
| Hexachlorobenzene    | 0.100       | Napropamide                   | 0.050       | Propoxycarbazone Na      | 0.050       |
| Hexaconazole         | 0.100       | Neburon                       | 0.020       | Propyzamide              | 0.050       |
| Hexazinone           | 0.100       | Nitrapyrin                    | 0.100       | Prothiofos               | 0.100       |
| Hexythiazox          | 0.020       | Norflurazon                   | 0.050       | Pyraclostrobin           | 0.020       |
| Imazalil             | 0.100       | Omethoate                     | 0.100       | Pyrazophos               | 0.050       |
| Imidacloprid         | 0.100       | O-Phenylphenol                | 0.100       | Pyrethrins               | 0.050       |
| Indaziflam           | 0.020       | Oxadixyl                      | 0.100       | Pyridaben                | 0.020       |
| Indoxacarb           | 0.020       | Oxamyl                        | 0.100       | Pyridafol                | 0.100       |
| Iprobenfos           | 0.100       | Oxamyl-oxime                  | 0.100       | Pyridate                 | 0.020       |
| Iprodione            | 0.100       | Oxychlorane                   | 0.100       | Pyrimethanil             | 0.050       |
| Isobenzan            | 0.100       | Oxydemeton-Methyl             | 0.100       | Pyriproxifen             | 0.020       |
| Isocarbophos         | 0.500       | Oxythioquinox                 | 0.200       | Pyroxasulfone            | 0.020       |
| Isodrin              | 0.100       | Paclobutrazol                 | 0.050       | Pyroxulam                | 0.020       |
| Isofenphos           | 0.050       | Paraoxon-ethyl                | 0.020       | Quinalphos               | 0.050       |
| Isofenphos-methyl    | 0.020       | Paraoxon methyl               | 0.100       | Quinoxyfen               | 0.050       |
| Isofenphos oxon      | 0.050       | Parathion ethyl               | 0.100       | Quintozene (PCNB)        | 0.200       |
| Isoprocarb           | 0.020       | Parathion methyl              | 0.200       | Resmethrin               | 0.050       |
| Isopropalin          | 0.200       | Penconazole                   | 0.050       | Rotenone                 | 0.050       |
| Isoprothiolane       | 0.050       | Pendimethalin                 | 0.050       | S421                     | 0.100       |
| Isoproturon          | 0.050       | Penflufen                     | 0.020       | Simazine                 | 0.100       |
| Isoxaben             | 0.050       | Pentachloroaniline            | 0.100       | Simetryn                 | 0.200       |
| Isoxaflutole         | 0.050       | Pentachloroanisole            | 0.100       | Spinetoram               | 0.020       |
| Kresoxim-methyl      | 0.050       | Pentachlorobenzene (PCB)      | 0.100       | Spinosad                 | 0.050       |
| Lactofen             | 0.500       | Pentachlorothioanisole (PCTA) | 0.100       | Spirodiclofen            | 0.100       |
| Lenacil              | 0.100       | Penthiopyrad                  | 0.020       | Spiromesifen             | 0.050       |
| Lindane (gamma BHC)  | 0.100       | Permethrin                    | 0.050       | Spirotetramat            | 0.050       |
| Linuron              | 0.020       | Perthane                      | 0.100       | Spiroxamine              | 0.020       |
| Malaaxon             | 0.050       | Phenmedipham                  | 0.050       | Sulfotep                 | 0.050       |
| Malathion            | 0.050       | Phenthoate                    | 0.050       | Sulfoxaflor              | 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       |
| Methacrifos          | 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       | Terbutylazine            | 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       | Tetradifon               | 0.200       |
| Methoxychlor         | 0.100       | Procymidone                   | 0.100       | Tetramethrin             | 0.050       |
| Methoxyfenozide      | 0.020       | Profenofos                    | 0.100       | Tetrasul                 | 0.100       |
| Metobromuron         | 0.050       | Profluralin                   | 0.100       | Thiabendazole            | 0.100       |
| Metolachlor          | 0.100       | Promecarb                     | 0.050       | Thiabendazole, 5-hydroxy | 0.100       |
| Metolcarb            | 0.050       | Prometon                      | 0.100       | Thiacloprid              | 0.050       |
| Metrafenone          | 0.050       | Prometryn                     | 0.020       | Thiamethoxam             | 0.100       |
| Metribuzin           | 0.100       | Propachlor                    | 0.020       | Thiobencarb              | 0.050       |
| Mevinphos            | 0.100       |                               |             | Thiodicarb               | 0.050       |
|                      |             |                               |             | Thiophanate-methyl       | 0.050       |



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**Report Number:** 23-000929/D008.R000  
**Report Date:** 01/30/2023  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 01/23/23 11:13



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       | Tolyfluanid  | 0.050       | Triticonazole   | 0.050       |
| Tralkoxydim      | 0.100       | Tridiphane   | 0.500       | Vinclozolin     | 0.100       |
| Triadimefon      | 0.050       | Triflumizole | 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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**Received:** 01/23/23 11:13

929 10f2



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

ORELAP ID: OR100028 ANAB ISO 17025 ID: AT-1508

NWNATURALGOODS 23-000929



|  |                    |          |  |   |   |  |   |   |   |   |   |   |   |  |        |  |
|--|--------------------|----------|--|---|---|--|---|---|---|---|---|---|---|--|--------|--|
| <b>Company:</b> Northwest Natural Goods<br><b>Contact:</b> Annie Nair<br><b>Address:</b> 11791 SE HWY 212<br><b>City:</b> Clackamas <b>State:</b> OR <b>Zip Code:</b> 97015<br><input checked="" type="checkbox"/> <b>Email Results:</b> annienair@nwnaturalgoods.com<br><input type="checkbox"/> <b>Ph:</b> ( ) -<br><i>Billing Contact (if different)</i><br><b>Name:</b> <b>Email:</b><br><b>Address:</b><br><b>City:</b> <b>State:</b> <b>Zip:</b><br><b>Ph:</b> ( ) - |                    |          | <b>Analysis Requested</b><br>Pesticides - OR 59 Compounds<br>Pesticide Multi-Residue - 379 compounds<br>Potency<br>Residual Solvents<br>Water Activity<br>Moisture<br>Micro: Yeast and Mold<br>Micro: E. Coli and Total Coliform<br>Heavy Metals<br>Mycotoxins |   |   |  |   |   |   |   |   |   | <b>PO Number:</b><br><b>Project ID:</b><br><b>Batch ID:</b><br><b>Sampled by:</b><br><b>Custom Reporting:</b> |  |        |  |
| <b>Lab ID</b> <b>Client Sample Identification</b> <b>Sample date</b>   |                    |          | <b>Material Type †</b> <b>Weight (Units)</b> <b>Comments/Metric ID</b>   |   |   |  |   |   |   |   |   |   |   |  |        |  |
|  | Bev - 88 023020 17 | 01/20/23 | ✓  | ✓ | ✓ | ✓  | ✓ | ✓ | ✓ | ✓ | ✓ | ✓ |   |  | 362.1g |  |
| <b>Signature - Relinquished By:</b> <b>Date</b> <b>Time</b>  |                    |          | <b>Signature - Received By:</b> <b>Date</b> <b>Time</b>  |   |   | <b>Lab Use Only:</b>   |   |   |   |   |   |   |   |  |        |  |
| Annie Nair 01/23/23  |                    |          | MRA 1/23 10:31   |   |   | <input type="checkbox"/> Shipped Via: _____ or <input type="checkbox"/> Client drop off<br>Evidence of cooling: <input type="checkbox"/> Yes   <input type="checkbox"/> No - Temp (°C): 18<br>Sample in good condition: <input type="checkbox"/> Yes   <input type="checkbox"/> No<br>Payment: <input type="checkbox"/> Cash   <input type="checkbox"/> Check   <input type="checkbox"/> CC   <input type="checkbox"/> Net:<br>Prelog storage: |   |   |   |   |   |   |   |  |        |  |
| MRA 1/23 10:55   |                    |          | RBS 01/23/23 11:3  |   |   |  |   |   |   |   |   |   |   |  |        |  |

† - Material Type Codes: Plant Material (P) ; Isolate (I) ; Concentrate/Extract (C) ; Tincture/Topical (T) ; Edible (E) ; Beverage (B) ; Vapor Product (V)

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  
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**Report Number:** 23-000929/D008.R000  
**Report Date:** 01/30/2023  
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929 20f2



**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-1508

| <b>Company:</b> Northwest Natural Goods<br><b>Contact:</b> Annie Nair<br><b>Address:</b> 11791 SE HWY 212<br><b>City:</b> Clackamas <b>State:</b> OR <b>Zip Code:</b> 97015<br><input checked="" type="checkbox"/> <b>Email Results:</b> annienair@nwnaturalgoods.com<br><input type="checkbox"/> <b>Ph:</b> ( ) -<br><i>Billing Contact (if different)</i><br><b>Name:</b> <b>Email:</b><br><b>Address:</b><br><b>City:</b> <b>State:</b> <b>Zip:</b><br><b>Ph:</b> ( ) - |                              |             | <b>Analysis Requested</b>  |                           |   |   |   |                                  |                               |  |   |   | <b>PO Number:</b><br><b>Project ID:</b><br><b>Batch ID:</b><br><b>Sampled by:</b><br><b>Custom Reporting:</b>  |                |                    |
|--|------------------------------|-------------|--|---------------------------|---|---|---|----------------------------------|-------------------------------|--|---|---|--|----------------|--------------------|
|  |                              |             | Pesticides - OR 59 Compounds<br>Pesticide Multi-Residue - 379 compounds<br>Potency<br>Residual Solvents<br>Water Activity<br>Moisture<br>Micro: Yeast and Mold<br>Micro: E.Coli and Total Coliform<br>Heavy Metals<br>Mycotoxins |                           |   |   |   |                                  |                               |  |   |   | <b>Source Material:</b> <input type="checkbox"/> - Ind. Hemp product   <input type="checkbox"/> - Rec. Cannabis<br><b>Reporting Type:</b> <input type="checkbox"/> - Compliance   <input type="checkbox"/> - R&D<br><b>Report to:</b> <input type="checkbox"/> - METRC   <input type="checkbox"/> - ODA   <input type="checkbox"/> - USDA  <br><input type="checkbox"/> - Other: |                |                    |
|  |                              |             | Turnaround time (TAT - Business Days):<br><input type="checkbox"/> - 5BD   <input checked="" type="checkbox"/> - 3BD*   <input type="checkbox"/> - 2BD*<br>*Check for availability   |                           |   |   |   |                                  |                               |  |   |   |  |                |                    |
| Lab ID   | Client Sample Identification | Sample date |  |                           |   |   |   |                                  |                               |  |   |   | Material Type †  | Weight (Units) | Comments/Metric ID |
|  | Bev - LM 022335 11-1         | 01/20/23    | ✓  | ✓                         | ✓   | ✓ | ✓ | ✓                                | ✓                             | ✓  | ✓ | ✓ |  | 362.1g         |                    |
| <b>Signature - Relinquished By:</b><br>Annie Nair<br>met   |                              |             | <b>Date:</b><br>01/23/23<br>1/23   | <b>Time:</b><br><br>10:55 | <b>Signature - Received By:</b><br>MRA<br>RBS |   |   | <b>Date:</b><br>1/23<br>01/23/23 | <b>Time:</b><br>10:31<br>1113 | <b>Lab Use Only:</b><br><input type="checkbox"/> Shipped Via: _____ or <input type="checkbox"/> Client drop off<br>Evidence of cooling: <input type="checkbox"/> Yes   <input type="checkbox"/> No - Temp (°C): 19<br>Sample in good condition: <input type="checkbox"/> Yes   <input type="checkbox"/> No<br>Payment: <input type="checkbox"/> Cash   <input type="checkbox"/> Check   <input type="checkbox"/> CC   <input type="checkbox"/> Net:<br>Prelog storage: |   |   |  |                |                    |

† - Material Type Codes: Plant Material (P) ; Isolate (I) ; Concentrate/Extract (C) ; Tincture/Topical (T) ; Edible (E) ; Beverage (B) ; Vapor Product (V)  
 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  
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**Report Number:** 23-000929/D008.R000  
**Report Date:** 01/30/2023  
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Revision: 1 Document ID: 7148  
 Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

JAOAC2015 V98-6 Batch ID: 2300772

| Laboratory Control Sample |     |        |       |       |       |        |       |            |       |
|---------------------------|-----|--------|-------|-------|-------|--------|-------|------------|-------|
| Analyte                   | LCS | Result | Spike | Units | % Rec | Limits |       | Evaluation | Notes |
| CBDVA                     | 2   | 0.0009 | 0.001 | %     | 92.7  | 80.0   | - 120 | Acceptable |       |
| CBDV                      | 2   | 0.0011 | 0.001 | %     | 99.2  | 80.0   | - 120 | Acceptable |       |
| CBE                       | 2   | 0.0010 | 0.001 | %     | 98.5  | 80.0   | - 120 | Acceptable |       |
| CBD                       | 1   | 0.0009 | 0.001 | %     | 93.1  | 90.0   | - 110 | Acceptable |       |
| CBG <sup>A</sup>          | 1   | 0.0009 | 0.001 | %     | 95.3  | 80.0   | - 120 | Acceptable |       |
| CBG                       | 1   | 0.0010 | 0.001 | %     | 99.2  | 80.0   | - 120 | Acceptable |       |
| CBD                       | 1   | 0.0009 | 0.001 | %     | 95.7  | 90.0   | - 110 | Acceptable |       |
| THCV                      | 2   | 0.0010 | 0.001 | %     | 97.7  | 80.0   | - 120 | Acceptable |       |
| δ8THCV                    | 2   | 0.0011 | 0.001 | %     | 102   | 80.0   | - 120 | Acceptable |       |
| THCV/A                    | 2   | 0.0009 | 0.001 | %     | 92.0  | 80.0   | - 120 | Acceptable |       |
| CBN                       | 1   | 0.0010 | 0.001 | %     | 99.8  | 80.0   | - 120 | Acceptable |       |
| exo-THC                   | 2   | 0.0010 | 0.001 | %     | 99.7  | 80.0   | - 120 | Acceptable |       |
| δ9THC                     | 1   | 0.0010 | 0.001 | %     | 98.9  | 90.0   | - 110 | Acceptable |       |
| δ8THC                     | 1   | 0.0010 | 0.001 | %     | 95.2  | 90.0   | - 110 | Acceptable |       |
| 9SΔ10THC                  | 1   | 0.0005 | 0.000 | %     | 113   | 80.0   | - 120 | Acceptable |       |
| CBL                       | 2   | 0.0010 | 0.001 | %     | 99.8  | 90.0   | - 110 | Acceptable |       |
| 9RΔ10THC                  | 1   | 0.0005 | 0.000 | %     | 98.9  | 80.0   | - 120 | Acceptable |       |
| CBC                       | 2   | 0.0010 | 0.001 | %     | 97.9  | 80.0   | - 120 | Acceptable |       |
| THCA                      | 1   | 0.0009 | 0.001 | %     | 91.9  | 80.0   | - 120 | Acceptable |       |
| CBCA                      | 2   | 0.0009 | 0.001 | %     | 88.9  | 80.0   | - 120 | Acceptable |       |
| CBLA                      | 2   | 0.0010 | 0.001 | %     | 94.2  | 90.0   | - 110 | Acceptable |       |
| CBT                       | 2   | 0.0010 | 0.001 | %     | 99.4  | 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 |       |
| CBD              | <LOQ   | 0.0001 | %     | < 0.0001 | Acceptable |       |
| CBG <sup>A</sup> | <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 |       |
| δ8THCV           | <LOQ   | 0.0001 | %     | < 0.0001 | Acceptable |       |
| THCV/A           | <LOQ   | 0.0001 | %     | < 0.0001 | Acceptable |       |
| CBN              | <LOQ   | 0.0001 | %     | < 0.0001 | Acceptable |       |
| exo-THC          | <LOQ   | 0.0001 | %     | < 0.0001 | Acceptable |       |
| δ9THC            | <LOQ   | 0.0001 | %     | < 0.0001 | Acceptable |       |
| δ8THC            | <LOQ   | 0.0001 | %     | < 0.0001 | Acceptable |       |
| 9SΔ10THC         | <LOQ   | 0.0001 | %     | < 0.0001 | Acceptable |       |
| CBL              | <LOQ   | 0.0001 | %     | < 0.0001 | Acceptable |       |
| 9RΔ10THC         | <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 |       |
| CBT              | <LOQ   | 0.0001 | %     | < 0.0001 | Acceptable |       |

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

Units of Measure:  
 %- Percent



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**Report Number:** 23-000929/D008.R000  
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 Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

| JAOAC2015 V98-6  |        | Batch ID: 2300772         |        |       |      |        |            |       |
|------------------|--------|---------------------------|--------|-------|------|--------|------------|-------|
| Sample Duplicate |        | Sample ID: 23-000912-0001 |        |       |      |        |            |       |
| 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 |       |
| CBDAA            | <LOQ   | <LOQ                      | 0.0001 | %     | NA   | < 20   | Acceptable |       |
| CBD <sup>A</sup> | <LOQ   | <LOQ                      | 0.0001 | %     | NA   | < 20   | Acceptable |       |
| CBC              | 0.0002 | 0.0002                    | 0.0001 | %     | 1.86 | < 20   | Acceptable |       |
| CBD              | 0.0073 | 0.0074                    | 0.0001 | %     | 1.63 | < 20   | Acceptable |       |
| THCV             | <LOQ   | <LOQ                      | 0.0001 | %     | NA   | < 20   | Acceptable |       |
| δ8THCV           | <LOQ   | <LOQ                      | 0.0001 | %     | NA   | < 20   | Acceptable |       |
| THCV/A           | <LOQ   | <LOQ                      | 0.0001 | %     | NA   | < 20   | Acceptable |       |
| CBN              | <LOQ   | <LOQ                      | 0.0001 | %     | NA   | < 20   | Acceptable |       |
| exo-THC          | <LOQ   | <LOQ                      | 0.0001 | %     | NA   | < 20   | Acceptable |       |
| δ9THC            | <LOQ   | <LOQ                      | 0.0001 | %     | NA   | < 20   | Acceptable |       |
| δ8THC            | <LOQ   | <LOQ                      | 0.0001 | %     | NA   | < 20   | Acceptable |       |
| 9Sδ10THC         | <LOQ   | <LOQ                      | 0.0001 | %     | NA   | < 20   | Acceptable |       |
| CBL              | <LOQ   | <LOQ                      | 0.0001 | %     | NA   | < 20   | Acceptable |       |
| 9Rδ10THC         | <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 |       |
| 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



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**Report Number:** 23-000929/D008.R000  
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**Received:** 01/23/23 11:13

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

Laboratory Quality Control Results

| Residual Solvents     |        |       |       | Batch ID: 2300804         |       |       |       |          |       |
|-----------------------|--------|-------|-------|---------------------------|-------|-------|-------|----------|-------|
| Method Blank          |        |       |       | Laboratory Control Sample |       |       |       |          |       |
| Analyte               | Result | LOQ   | Notes | Result                    | Spike | Units | % Rec | Limits   | Notes |
| Propane               | ND     | < 200 |       | 535                       | 572   | µg/g  | 93.5  | 60 - 120 |       |
| Isobutane             | ND     | < 200 |       | 629                       | 731   | µg/g  | 86.0  | 60 - 120 |       |
| Butane                | ND     | < 200 |       | 614                       | 731   | µg/g  | 84.0  | 60 - 120 |       |
| 2,2-Dimethylpropane   | ND     | < 200 |       | 784                       | 936   | µg/g  | 83.8  | 60 - 120 |       |
| Methanol              | ND     | < 200 |       | 1600                      | 1620  | µg/g  | 98.8  | 60 - 120 |       |
| Ethylene Oxide        | ND     | < 30  |       | 48.5                      | 56.2  | µg/g  | 86.3  | 60 - 120 |       |
| 2-Methylbutane        | ND     | < 200 |       | 1410                      | 1610  | µg/g  | 87.6  | 60 - 120 |       |
| Pentane               | ND     | < 200 |       | 1410                      | 1600  | µg/g  | 88.1  | 60 - 120 |       |
| Ethanol               | ND     | < 200 |       | 1510                      | 1610  | µg/g  | 93.8  | 70 - 130 |       |
| Ethyl Ether           | ND     | < 200 |       | 1400                      | 1630  | µg/g  | 85.9  | 60 - 120 |       |
| 2,2-Dimethylbutane    | ND     | < 30  |       | 164                       | 171   | µg/g  | 95.9  | 60 - 120 |       |
| Acetone               | ND     | < 200 |       | 1500                      | 1630  | µg/g  | 92.0  | 60 - 120 |       |
| 2-Propanol            | ND     | < 200 |       | 1510                      | 1620  | µg/g  | 93.2  | 60 - 120 |       |
| Ethyl Formate         | ND     | < 500 |       | 1640                      | 1670  | µg/g  | 98.2  | 70 - 130 |       |
| Acetonitrile          | ND     | < 100 |       | 435                       | 498   | µg/g  | 87.3  | 60 - 120 |       |
| Methyl Acetate        | ND     | < 500 |       | 1680                      | 1730  | µg/g  | 97.1  | 70 - 130 |       |
| 2,3-Dimethylbutane    | ND     | < 30  |       | 140                       | 171   | µg/g  | 81.9  | 60 - 120 |       |
| Dichloromethane       | ND     | < 60  |       | 433                       | 483   | µg/g  | 89.6  | 60 - 120 |       |
| 2-Methylpentane       | ND     | < 30  |       | 146                       | 168   | µg/g  | 86.9  | 60 - 120 |       |
| MTBE                  | ND     | < 500 |       | 1560                      | 1650  | µg/g  | 94.5  | 70 - 130 |       |
| 3-Methylpentane       | ND     | < 30  |       | 136                       | 167   | µg/g  | 81.4  | 60 - 120 |       |
| Hexane                | ND     | < 30  |       | 204                       | 182   | µg/g  | 112.1 | 60 - 120 |       |
| 1-Propanol            | ND     | < 500 |       | 1480                      | 1620  | µg/g  | 91.4  | 70 - 130 |       |
| Methylethylketone     | ND     | < 500 |       | 1650                      | 1620  | µg/g  | 101.9 | 70 - 130 |       |
| Ethyl acetate         | ND     | < 200 |       | 1540                      | 1610  | µg/g  | 95.7  | 60 - 120 |       |
| 2-Butanol             | ND     | < 200 |       | 1510                      | 1600  | µg/g  | 94.4  | 60 - 120 |       |
| Tetrahydrofuran       | ND     | < 100 |       | 400                       | 483   | µg/g  | 82.8  | 60 - 120 |       |
| Cyclohexane           | ND     | < 200 |       | 1600                      | 1610  | µg/g  | 99.4  | 60 - 120 |       |
| 2-methyl-1-propanol   | ND     | < 500 |       | 1420                      | 1620  | µg/g  | 87.7  | 70 - 130 |       |
| Benzene               | ND     | < 1   |       | 4.48                      | 5.02  | µg/g  | 89.2  | 60 - 120 |       |
| Isopropyl Acetate     | ND     | < 200 |       | 1560                      | 1620  | µg/g  | 96.3  | 60 - 120 |       |
| Heptane               | ND     | < 200 |       | 1590                      | 1610  | µg/g  | 98.8  | 60 - 120 |       |
| 1-Butanol             | ND     | < 500 |       | 1110                      | 1630  | µg/g  | 68.1  | 70 - 130 | Q6    |
| Propyl Acetate        | ND     | < 500 |       | 1420                      | 1610  | µg/g  | 88.2  | 70 - 130 |       |
| 1,4-Dioxane           | ND     | < 100 |       | 474                       | 491   | µg/g  | 96.5  | 60 - 120 |       |
| 2-Ethoxyethanol       | ND     | < 30  |       | 103                       | 181   | µg/g  | 56.9  | 60 - 120 | Q6    |
| Methylisobutylketone  | ND     | < 500 |       | 1580                      | 1620  | µg/g  | 97.5  | 70 - 130 |       |
| 3-Methyl-1-butanol    | ND     | < 500 |       | 1210                      | 1630  | µg/g  | 74.2  | 70 - 130 |       |
| Ethylene Glycol       | ND     | < 200 |       | 452                       | 484   | µg/g  | 93.4  | 60 - 120 |       |
| Toluene               | ND     | < 100 |       | 417                       | 485   | µg/g  | 86.0  | 60 - 120 |       |
| Isobutyl Acetate      | ND     | < 500 |       | 1670                      | 1630  | µg/g  | 102.5 | 70 - 130 |       |
| 1-Pentanol            | ND     | < 500 |       | 1240                      | 1620  | µg/g  | 76.5  | 70 - 130 |       |
| Butyl Acetate         | ND     | < 500 |       | 1390                      | 1620  | µg/g  | 85.8  | 70 - 130 |       |
| Ethylbenzene          | ND     | < 200 |       | 936                       | 969   | µg/g  | 96.6  | 60 - 120 |       |
| m,p-Xylene            | ND     | < 200 |       | 964                       | 994   | µg/g  | 97.0  | 60 - 120 |       |
| o-Xylene              | ND     | < 200 |       | 917                       | 967   | µg/g  | 94.8  | 60 - 120 |       |
| Cumene                | ND     | < 30  |       | 147                       | 171   | µg/g  | 86.0  | 60 - 120 |       |
| Anisole               | ND     | < 500 |       | 1170                      | 1630  | µg/g  | 71.8  | 70 - 130 |       |
| DMISO                 | ND     | < 500 |       | 1210                      | 1680  | µg/g  | 72.0  | 70 - 130 |       |
| 1,2-dimethoxyethane   | ND     | < 50  |       | 168                       | 169   | µg/g  | 99.4  | 70 - 130 |       |
| Triethylamine         | ND     | < 500 |       | 1650                      | 1630  | µg/g  | 101.2 | 70 - 130 |       |
| N,N-dimethylformamide | ND     | < 150 |       | 420                       | 482   | µg/g  | 87.1  | 70 - 130 |       |
| N,N-dimethylacetamide | ND     | < 150 |       | 375                       | 510   | µg/g  | 73.5  | 70 - 130 |       |
| Pyridine              | ND     | < 50  |       | 193                       | 203   | µg/g  | 95.1  | 70 - 130 |       |
| Silfolane             | ND     | < 50  |       | 130                       | 172   | µg/g  | 75.6  | 70 - 130 |       |
| 1,2-Dichloroethane    | ND     | < 1   |       | 1.05                      | 1     | µg/g  | 105.0 | 70 - 130 |       |
| Chloroform            | ND     | < 1   |       | 1.08                      | 1     | µg/g  | 108.0 | 70 - 130 |       |
| Trichloroethylene     | ND     | < 1   |       | 1.11                      | 1     | µg/g  | 111.0 | 70 - 130 |       |
| 1,1,1-Trichloroethane | ND     | < 1   |       | 1.01                      | 1     | µg/g  | 101.0 | 70 - 130 |       |



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**Report Number:** 23-000929/D008.R000  
**Report Date:** 01/30/2023  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 01/23/23 11:13

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

QC- Sample Duplicate Sample ID: 23-000757-0001

| 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  
 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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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.                      |