



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



Report Number: 23-010635/D002.R000
Report Date: 09/14/2023
ORELAP#: OR100028
Purchase Order:
Received: 09/07/23 11:13

Customer: NW Natural Goods
Product identity: GF 023249-1
Client/Metric ID: .
Laboratory ID: 23-010635-0001

Summary

Potency:

| Analyte per 355ml | Result | Limits | Units | Status | |
|-------------------|--------|--------|----------|--------|--|
| CBD per 355ml | 21.5 | | mg/355ml | | CBD-Total per Serving Size 21.5 mg/355ml |
| CBG per 355ml | 11.5 | | 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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Customer: NW Natural Goods

Product identity: GF 023249-1

Client/Metric ID: .

Sample Date:

Laboratory ID: 23-010635-0001

Evidence of Cooling: No

Temp: 19.5 °C

Relinquished by: Ramos

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: 2310817 | Analyze: 9/11/23 5:04:00 PM |
|-------------------------|--------|--|----------|-------------|----------------|-----------------------------|
| Analyte | Result | Limits | Units | LOQ | Notes | |
| CBC per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| CBC-A per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| CBC-Total per 355ml | < LOQ | | mg/355ml | 0.667 | | |
| CBD per 355ml | 21.5 | | mg/355ml | 0.355 | | |
| CBD-A per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| CBD-Total per 355ml | 21.5 | | mg/355ml | 0.667 | | |
| CBDV per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| CBDV-A per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| CBDV-Total per 355ml | < LOQ | | mg/355ml | 0.663 | | |
| CBE per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| CBG per 355ml | 11.5 | | mg/355ml | 0.355 | | |
| CBG-A per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| CBG-Total per 355ml | 11.5 | | mg/355ml | 0.663 | | |
| CBL per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| CBL-A per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| CBL-Total per 355ml | < LOQ | | mg/355ml | 0.667 | | |
| CBN per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| CBT per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| Δ8-THCV per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| Δ10-THC-9R per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| Δ10-THC-9S per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| Δ10-THC-Total per 355ml | < LOQ | | mg/355ml | 0.711 | | |
| Δ8-THC per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| Δ9-THC per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| delta-9-THCP per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| exo-THC per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| THC-A per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| THC-Total per 355ml | < LOQ | | mg/355ml | 0.667 | | |
| THCV per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| THCV-A per 355ml | < LOQ | | mg/355ml | 0.355 | | |
| THCV-Total per 355ml | < LOQ | | mg/355ml | 0.667 | | |



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Potency per 355ml Method: J AOAC 2015 V98-6 (mod)^P Units mg/se Batch: 2310817 Analyze: 9/11/23 5:04:00 PM

| Analyte | Result | Limits | Units | LOQ | Notes |
|------------------------------|--------|--------|----------|-----|-------|
| Total Cannabinoids per 355ml | 33.0 | | mg/355ml | | |

Microbiology

| Analyte | Result | Limits | Units | LOQ | Batch | Analyzed Method | Status | Notes |
|-------------------------|--------|--------|-------|-----|---------|---|--------|-------|
| Aerobic Plate Count | < LOQ | | cfu/g | 10 | 2310707 | 09/10/23 AOAC 990.12 (Petrifilm) ^P | | |
| E.coli | < LOQ | | cfu/g | 10 | 2310703 | 09/10/23 AOAC 991.14 (Petrifilm) ^P | | |
| Total Coliforms | < LOQ | | cfu/g | 10 | 2310703 | 09/10/23 AOAC 991.14 (Petrifilm) ^P | | |
| Mold (RAPID Petrifilm) | < LOQ | | cfu/g | 10 | 2310706 | 09/10/23 AOAC 2014.05 (RAPID) ^P | | |
| Yeast (RAPID Petrifilm) | < LOQ | | cfu/g | 10 | 2310706 | 09/10/23 AOAC 2014.05 (RAPID) ^P | | |

Solvents Method: Residual Solvents by GC/MS^P Units µg/g Batch 2310820 Analyze 09/12/23 01:46 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-Dimethyl butane | < LOQ | | 30.0 | | | 2,2-Dimethylpropane (neo-pentane) | < LOQ | | 200 | | |
| 2,3-Dimethyl butane | < LOQ | | 30.0 | | | 3-Methylpentane | < LOQ | | 30.0 | | |
| Acetone | < LOQ | 5000 | 200 | pass | | Acetonitrile | < LOQ | 410 | 100 | pass | |
| Benzene | < LOQ | 2.00 | 1.00 | pass | | Butanes (sum) | < LOQ | 5000 | 400 | pass | |
| Cyclohexane | < LOQ | 3880 | 200 | pass | | Ethyl acetate | < LOQ | 5000 | 200 | pass | |
| Ethyl benzene | < LOQ | | 200 | | | Ethyl ether | < LOQ | 5000 | 200 | pass | |
| Ethylene glycol | < LOQ | 620 | 200 | pass | | Ethylene oxide | < LOQ | 50.0 | 20.0 | pass | |
| Hexanes (sum) | < LOQ | 290 | 150 | pass | | Isopropyl acetate | < LOQ | 5000 | 200 | pass | |
| Isopropyl benzene (Cumene) | < LOQ | 70.0 | 30.0 | pass | | m,p-Xylene | < LOQ | | 200 | | |
| Methanol | < LOQ | 3000 | 200 | pass | | Methylene chloride | < LOQ | 600 | 60.0 | pass | |
| Methylpropane (Isobutane) | < LOQ | | 200 | | | n-Butane | < LOQ | | 200 | | |
| n-Heptane | < LOQ | 5000 | 200 | pass | | n-Hexane | < LOQ | | 30.0 | | |
| n-Pentane | < LOQ | | 200 | | | o-Xylene | < LOQ | | 200 | | |
| Pentanes (sum) | < LOQ | 5000 | 600 | pass | | Propane | < LOQ | 5000 | 200 | pass | |
| Tetrahydrofuran | < LOQ | 720 | 100 | pass | | Toluene | < LOQ | 890 | 100 | pass | |
| Total Xylenes | < LOQ | | 400 | | | Total Xylenes and Ethyl benzene | < LOQ | 2170 | 600 | pass | |

Pesticides Method: AOAC 2007.01 & EN 15662 (mod)^P Units mg/kg Batch 2310775 Analyze 09/12/23 11:47 AM

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



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Metals

| Analyte | Result | Limits | Units | LOQ | Batch | Analyzed Method | Status | Notes |
|----------------------|--------|--------|-------|---------|---------|---|--------|-------|
| Arsenic [‡] | < LOQ | 0.200 | mg/kg | 0.00371 | 2310725 | 09/08/23 AOAC 2013.06 (mod.) [‡] | pass | |
| Cadmium [‡] | < LOQ | 0.200 | mg/kg | 0.00371 | 2310725 | 09/08/23 AOAC 2013.06 (mod.) [‡] | pass | |
| Lead [‡] | < LOQ | 0.500 | mg/kg | 0.00371 | 2310725 | 09/08/23 AOAC 2013.06 (mod.) [‡] | pass | |
| Mercury [‡] | < LOQ | 0.100 | mg/kg | 0.00185 | 2310725 | 09/08/23 AOAC 2013.06 (mod.) [‡] | pass | |

Mycotoxins

| Analyte | Result | Limits | Units | LOQ | Batch | Analyzed Method | Status | Notes |
|-------------------------------|--------|--------|-------|------|---------|---|--------|-------|
| Aflatoxin B2 [‡] | < LOQ | | µg/kg | 5.00 | 2310759 | 09/11/23 AOAC 2007.01 & EN 15662 (mod) [‡] | | |
| Aflatoxin B1 [‡] | < LOQ | | µg/kg | 5.00 | 2310759 | 09/11/23 AOAC 2007.01 & EN 15662 (mod) [‡] | | |
| Aflatoxin G1 [‡] | < LOQ | | µg/kg | 5.00 | 2310759 | 09/11/23 AOAC 2007.01 & EN 15662 (mod) [‡] | | |
| Aflatoxin G2 [‡] | < LOQ | | µg/kg | 5.00 | 2310759 | 09/11/23 AOAC 2007.01 & EN 15662 (mod) [‡] | | |
| Ochratoxin A [‡] | < LOQ | 20.0 | µg/kg | 5.00 | 2310759 | 09/11/23 AOAC 2007.01 & EN 15662 (mod) [‡] | pass | |
| Total Aflatoxins [‡] | 0.000 | 20.0 | µg/kg | 20.0 | | 09/14/23 AOAC 2007.01 & EN 15662 (mod) [‡] | pass | |



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Abbreviations

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

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

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

[¥] = TNI accredited analyte.

Units of Measure

cfu/g = Colony forming units per gram

g = g

g/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

Approved Signatory

Derrick Tanner
General Manager



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Cannab s Mu t -Res due Prof e, L m ts of Quant tat on

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



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Cannab s Mu t -Res due Prof e, L m ts of Quant tat on

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



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Cannab s Mu t-Res due Prof e, L m ts of Quant tat on

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

LOQ= Limit of Quantitation, mg/kg

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



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



Report Number: 23-010635/D002.R000
Report Date: 09/14/2023
ORELAP#: OR100028
Purchase Order:
Received: 09/07/23 11:13



Hemp & Cannabis
 Chain of Custody

NW-Natural-Goods-
 1694025625

ORELAP ID: CR1000028 ANAB ISO 17025 ID: AT1608

| Project Information | | | | | Testing | | | | | | |
|--|------------------|----------|------------------|--------------------|---|-------------------------------|--|--------------------------------|------------------------|---|---|
| Project Name <i>N/A</i> PO Number <i>N/A</i> Turnaround Time <i>5 Business Days (Req. For Micro Testing) Standard</i> Samples Delivered to Laboratory <i>Schedule Pick-Up</i> | | | | | H0010 - Potency Cannabis/oid Basic + Extended Profile | H0042 - Mycotoxins Compliance | F20200 - Pesticide - Multi-Residue Profile | H0008 - Residual Solvents - CR | M010 - Micro Profile D | H0013 - Heavy Metals Profile CR (As, Cd, Pb & Hg) | |
| # | Sample Name Test | Material | Amount Provided | Reporting Unit | Serving Size | | | | | | |
| 1 | GF 023249-1 | Beverage | 4 units for sale | mg/g & mg/serveing | 365 ml | ✓ | ✓ | ✓ | ✓ | ✓ | ✓ |

| Relinquished By | Date | Time | Temp., °C | Received By | Date | Time | Received Temp., °C | Evidence of Cooling? |
|------------------|-----------------|--------------|-------------|-------------|-----------------|--------------|--------------------|----------------------|
| <i>Jack Hood</i> | <i>9/6/2023</i> | <i>11:40</i> | | <i>BR</i> | <i>9/7/2023</i> | <i>10:23</i> | | <i>No</i> |
| <i>BR</i> | <i>9/7/2023</i> | <i>11:07</i> | <i>19.5</i> | <i>MRH</i> | <i>9/7/2023</i> | <i>11:13</i> | | <i>No</i> |

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

Columbia Laboratories
 12423 NE Whitaker Way
 Portland, OR 97230

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info@columbialaboratories.com

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



Report Number: 23-010635/D002.R000
 Report Date: 09/14/2023
 ORELAP#: OR100028
 Purchase Order:
 Received: 09/07/23 11:13

Revision 4 Documen D 7148
 Legacy D Workshee Valida ed 04/20/2021

Laboratory Quality Control Results

JACAC2015 V98-6 Batch ID: 2310817

| Laboratory Control Sample | | | | | | | | | |
|---------------------------|-----|---------|---------|-------|-------|--------|-------|------------|-------|
| Analyte | LCS | Result | Spike | Units | % Rec | Limits | | | Notes |
| CBDVA | 2 | 0.00101 | 0.0009 | % | 109 | 80.0 | - 120 | Acceptable | |
| CBDV | 2 | 0.00106 | 0.0009 | % | 112 | 80.0 | - 120 | Acceptable | |
| CBE | 2 | 0.00111 | 0.0010 | % | 112 | 80.0 | - 120 | Acceptable | |
| CBD | 1 | 0.00118 | 0.00109 | % | 109 | 90.0 | - 110 | Acceptable | |
| CBGA | 1 | 0.00122 | 0.00110 | % | 110 | 80.0 | - 120 | Acceptable | |
| CBG | 1 | 0.00125 | 0.00118 | % | 107 | 80.0 | - 120 | Acceptable | |
| THCV | 1 | 0.00123 | 0.00114 | % | 108 | 90.0 | - 110 | Acceptable | |
| THCVA | 2 | 0.0007 | 0.0006 | % | 109 | 80.0 | - 120 | Acceptable | |
| THCV | 2 | 0.0007 | 0.0006 | % | 109 | 80.0 | - 120 | Acceptable | |
| THCVA | 2 | 0.0009 | 0.0008 | % | 114 | 80.0 | - 120 | Acceptable | |
| THCVA | 2 | 0.00100 | 0.0009 | % | 109 | 80.0 | - 120 | Acceptable | |
| CBN | 1 | 0.00123 | 0.00111 | % | 111 | 80.0 | - 120 | Acceptable | |
| exo-THC | 2 | 0.00101 | 0.0009 | % | 111 | 80.0 | - 120 | Acceptable | |
| d9THC | 1 | 0.00123 | 0.00115 | % | 106 | 90.0 | - 110 | Acceptable | |
| d8THC | 1 | 0.00130 | 0.00117 | % | 110 | 90.0 | - 110 | Acceptable | Q1 |
| 9S-d10THC | 1 | 0.00122 | 0.00115 | % | 106 | 80.0 | - 120 | Acceptable | |
| CBL | 2 | 0.00105 | 0.0009 | % | 112 | 80.0 | - 120 | Acceptable | |
| 9R-d10THC | 1 | 0.0004 | 0.0004 | % | 112 | 80.0 | - 120 | Acceptable | |
| CBG | 2 | 0.00110 | 0.00100 | % | 110 | 80.0 | - 120 | Acceptable | |
| THCA | 1 | 0.00122 | 0.00111 | % | 111 | 90.0 | - 110 | Acceptable | Q1 |
| CBGA | 2 | 0.00103 | 0.0010 | % | 108 | 80.0 | - 120 | Acceptable | |
| CBLA | 2 | 0.00216 | 0.00194 | % | 111 | 80.0 | - 120 | Acceptable | |
| d9THCP | 2 | 0.00106 | 0.0009 | % | 112 | 80.0 | - 120 | Acceptable | |
| CBT | 2 | 0.00104 | 0.0009 | % | 113 | 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 | |
| CBGA | <LOQ | 0.0001 | % | < 0.0001 | Acceptable | |
| CBG | <LOQ | 0.0001 | % | < 0.0001 | Acceptable | |
| THCV | <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 | |
| CBG | <LOQ | 0.0001 | % | < 0.0001 | Acceptable | |
| THCA | <LOQ | 0.0001 | % | < 0.0001 | Acceptable | |
| CBGA | <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



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



Report Number: 23-010635/D002.R000
Report Date: 09/14/2023
ORELAP#: OR100028
Purchase Order:
Received: 09/07/23 11:13

Revision 4 Documen D 7148
Legacy D Workshee Valida ed 04/20/2021

Laboratory Quality Control Results

| JAOAC2015 V98-6 | | Batch ID: 2310817 | | | | | | |
|------------------|---------|---------------------------|--------|-------|-------|--------|------------|-------|
| Sample Duplicate | | Sample ID: 23-010635-0001 | | | | | | |
| 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 | |
| CBD | <LOQ | <LOQ | 0.0001 | % | NA | < 20 | Acceptable | |
| CBDA | <LOQ | <LOQ | 0.0001 | % | NA | < 20 | Acceptable | |
| CBGA | <LOQ | <LOQ | 0.0001 | % | NA | < 20 | Acceptable | |
| CBG | 0.00318 | 0.00317 | 0.0001 | % | 0.368 | < 20 | Acceptable | |
| CBD | 0.00597 | 0.00594 | 0.0001 | % | 0.462 | < 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



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



Report Number: 23-010635/D002.R000
 Report Date: 09/14/2023
 ORELAP#: OR100028
 Purchase Order:
 Received: 09/07/23 11:13

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

Laboratory Quality Control Results

| Residual Solvents | | | | Batch D: 2310820 | | | | | |
|-----------------------|--------|-------|-------|---------------------------|-------|-------|-------|----------|-------|
| Method Blank | | | | Laboratory Control Sample | | | | | |
| Analyte | Result | LOQ | Notes | Result | Spike | Units | % Rec | Limits | Notes |
| Propane | ND | < 200 | | 560 | 584 | µg/g | 95.9 | 60 - 120 | |
| Isobutane | ND | < 200 | | 711 | 767 | µg/g | 92.7 | 60 - 120 | |
| Butane | ND | < 200 | | 732 | 782 | µg/g | 93.6 | 60 - 120 | |
| 2,2-Dimethylpropane | ND | < 200 | | 892 | 939 | µg/g | 95.0 | 60 - 120 | |
| Methanol | ND | < 200 | | 1690 | 1670 | µg/g | 101.2 | 60 - 120 | |
| Ethylene Oxide | ND | < 30 | | 57.9 | 57.1 | µg/g | 101.4 | 60 - 120 | |
| 2-Methylbutane | ND | < 200 | | 1550 | 1680 | µg/g | 92.3 | 60 - 120 | |
| Pentane | ND | < 200 | | 1540 | 1670 | µg/g | 92.2 | 60 - 120 | |
| Ethanol | ND | < 200 | | 1670 | 1660 | µg/g | 100.6 | 70 - 130 | |
| Ethyl Ether | ND | < 200 | | 1600 | 1670 | µg/g | 95.8 | 60 - 120 | |
| 2,2-Dimethylbutane | ND | < 30 | | 176 | 189 | µg/g | 93.1 | 60 - 120 | |
| Acetone | ND | < 200 | | 1630 | 1670 | µg/g | 97.6 | 60 - 120 | |
| 2-Propanol | ND | < 200 | | 1620 | 1630 | µg/g | 99.4 | 60 - 120 | |
| Ethyl Formate | ND | < 500 | | 1380 | 1600 | µg/g | 86.3 | 70 - 130 | |
| Acetonitrile | ND | < 100 | | 471 | 492 | µg/g | 95.7 | 60 - 120 | |
| Methyl Acetate | ND | < 500 | | 1570 | 1600 | µg/g | 98.1 | 70 - 130 | |
| 2,3-Dimethylbutane | ND | < 30 | | 186 | 180 | µg/g | 103.3 | 60 - 120 | |
| Dichloromethane | ND | < 60 | | 475 | 488 | µg/g | 97.3 | 60 - 120 | |
| 2-Methylpentane | ND | < 30 | | 173 | 182 | µg/g | 95.1 | 60 - 120 | |
| MTBE | ND | < 500 | | 1560 | 1610 | µg/g | 96.9 | 70 - 130 | |
| 3-Methylpentane | ND | < 30 | | 160 | 177 | µg/g | 90.4 | 60 - 120 | |
| Hexane | ND | < 30 | | 169 | 177 | µg/g | 95.5 | 60 - 120 | |
| 1-Propanol | ND | < 500 | | 1510 | 1600 | µg/g | 94.4 | 70 - 130 | |
| Methyl ethyl ketone | ND | < 500 | | 1550 | 1610 | µg/g | 96.3 | 70 - 130 | |
| Ethyl acetate | ND | < 200 | | 1570 | 1630 | µg/g | 96.3 | 60 - 120 | |
| 2-Butanol | ND | < 200 | | 1540 | 1630 | µg/g | 94.5 | 60 - 120 | |
| Tetrahydrofuran | ND | < 100 | | 454 | 488 | µg/g | 93.0 | 60 - 120 | |
| Cyclohexane | ND | < 200 | | 1510 | 1610 | µg/g | 93.8 | 60 - 120 | |
| 2-methyl-1-propanol | ND | < 500 | | 1480 | 1610 | µg/g | 91.9 | 70 - 130 | |
| Benzene | ND | < 1 | | 4.69 | 4.79 | µg/g | 97.9 | 60 - 120 | |
| Isopropyl Acetate | ND | < 200 | | 1580 | 1650 | µg/g | 95.8 | 60 - 120 | |
| Heptane | ND | < 200 | | 1510 | 1630 | µg/g | 92.6 | 60 - 120 | |
| 1-Butanol | ND | < 500 | | 1330 | 1600 | µg/g | 83.1 | 70 - 130 | |
| Propyl Acetate | ND | < 500 | | 1430 | 1600 | µg/g | 89.4 | 70 - 130 | |
| 1,4-Dioxane | ND | < 100 | | 460 | 523 | µg/g | 88.0 | 60 - 120 | |
| 2-Ethoxyethanol | ND | < 30 | | 166 | 179 | µg/g | 92.7 | 60 - 120 | |
| Methylisobutylketone | ND | < 500 | | 1370 | 1600 | µg/g | 85.6 | 70 - 130 | |
| 3-Methyl-1-butanol | ND | < 500 | | 1210 | 1600 | µg/g | 75.6 | 70 - 130 | |
| Ethylene Glycol | ND | < 200 | | 289 | 508 | µg/g | 53.2 | 60 - 120 | Q6 |
| Toluene | ND | < 100 | | 446 | 496 | µg/g | 89.9 | 60 - 120 | |
| Isobutyl Acetate | ND | < 500 | | 1440 | 1610 | µg/g | 89.4 | 70 - 130 | |
| 1-Pentanol | ND | < 500 | | 1500 | 1600 | µg/g | 93.8 | 70 - 130 | |
| Butyl Acetate | ND | < 500 | | 1340 | 1610 | µg/g | 83.2 | 70 - 130 | |
| Ethylbenzene | ND | < 200 | | 790 | 978 | µg/g | 80.8 | 60 - 120 | |
| m,p-Xylene | ND | < 200 | | 808 | 994 | µg/g | 81.3 | 60 - 120 | |
| o-Xylene | ND | < 200 | | 770 | 982 | µg/g | 78.4 | 60 - 120 | |
| Cumene | ND | < 30 | | 133 | 171 | µg/g | 77.8 | 60 - 120 | |
| Anisole | ND | < 500 | | 1210 | 1600 | µg/g | 75.6 | 70 - 130 | |
| DMF | ND | < 500 | | 1170 | 1620 | µg/g | 72.2 | 70 - 130 | |
| 1,2-dimethoxyethane | ND | < 50 | | 170 | 186 | µg/g | 91.4 | 70 - 130 | |
| Triethylamine | ND | < 500 | | 1380 | 1600 | µg/g | 86.3 | 70 - 130 | |
| N,N-dimethylformamide | ND | < 150 | | 410 | 480 | µg/g | 85.4 | 70 - 130 | |
| N,N-dimethylacetamide | ND | < 150 | | 343 | 483 | µg/g | 71.0 | 70 - 130 | |
| Pyridine | ND | < 50 | | 133 | 168 | µg/g | 79.2 | 70 - 130 | |
| Silfolane | ND | < 50 | | 91.9 | 161 | µg/g | 57.1 | 70 - 130 | Q6 |
| 1,2-Dichloroethane | ND | < 1 | | 0.985 | 1 | µg/g | 98.5 | 70 - 130 | |
| Chloroform | ND | < 1 | | 0.86 | 1 | µg/g | 86.0 | 70 - 130 | |
| Trichloroethylene | ND | < 1 | | 0.862 | 1 | µg/g | 86.2 | 70 - 130 | |
| 1,1-Dichloroethane | ND | < 1 | | 0.923 | 1 | µg/g | 92.3 | 70 - 130 | |



12423 NE Whitaker Way
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Report Number: 23-010635/D002.R000
Report Date: 09/14/2023
ORELAP#: OR100028
Purchase Order:
Received: 09/07/23 11:13

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

QC- Sample Duplicate Sample ID: 23-010289-0002

| 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



12423 NE Whitaker Way
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Report Number: 23-010635/D002.R000
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Received: 09/07/23 11:13

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. |