



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



Report Number: 23-002684/D046.R000
Report Date: 03/13/2023
ORELAP#: OR100028
Purchase Order:
Received: 03/06/23 16:17

Customer: Seventh Hill Releaf LLC
Product identity: Hemp Collect LRD9
Client/Metric ID: HCLR40CBD20231
Laboratory ID: 23-002684-0015

Summary

Potency:

| Analyte | Result | Limits | Units | Status | |
|-----------------|---------|--------|--------|--------|--|
| CBD | 0.00370 | | % | | THC-Total per Serving Size 42.1 mg/16g |
| CBD-A | 0.00980 | | % | | |
| Δ8-THC | 0.00407 | | % | | CBD-Total per Serving Size 1.97 mg/16g |
| Δ9-THC | 0.263 | | % | | (Reported in milligrams per serving) |
| Analyte per 16g | Result | Limits | Units | Status | |
| CBD per 16g | 0.592 | | mg/16g | | |
| CBD-A per 16g | 1.57 | | mg/16g | | |
| Δ8-THC per 16g | 0.651 | | mg/16g | | |
| Δ9-THC per 16g | 42.1 | | mg/16g | | |

Residual Solvents:

All analytes passing and less than LOQ.

Metals:

| Analyte | Result | Units | Limit | Status |
|----------|--------|-------|-------|--------|
| Arsenic* | 0.0206 | mg/kg | 0.200 | pass |



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Received: 03/06/23 16:17

Customer: Seventh Hill Releaf LLC
215 S 19th St
Springfield Oregon 97477
United States of America (USA)

Product identity: Hemp Collect LRD9
Client/Metric ID: HCLR40CBD20231

Sample Date:
Laboratory ID: 23-002684-0015

Evidence of Cooling: No
Temp: 19.8

Relinquished by: client
Serving Size #1: 16 g

Sample Results

| Potency | Method: J AOAC 2015 V98-6 (mod) ^b | Units % | Batch: 2302154 | Analyze: 3/9/23 8:23:00 PM | |
|---------------------------|--|---------|----------------|----------------------------|-------|
| Analyte | Result | Limits | Units | LOQ | Notes |
| CBD | 0.00370 | | % | 0.00317 | |
| CBD-A | 0.00980 | | % | 0.00317 | |
| CBD-Total | 0.0123 | | % | 0.00595 | |
| CBG | < LOQ | | % | 0.00317 | |
| CBG-A | < LOQ | | % | 0.00317 | |
| CBG-Total | < LOQ | | % | 0.00592 | |
| CBN | < LOQ | | % | 0.00317 | |
| Δ10-THC-9R | < LOQ | | % | 0.00317 | |
| Δ10-THC-9S | < LOQ | | % | 0.00317 | |
| Δ10-THC-Total | < LOQ | | % | 0.00634 | |
| Δ8-THC | 0.00407 | | % | 0.00317 | |
| Δ9-THC | 0.263 | | % | 0.00317 | |
| THC-A | < LOQ | | % | 0.00317 | |
| THC-Total | 0.263 | | % | 0.00595 | |
| Total Cannabinoids | 0.281 | | % | | |

| Potency per 16g | Method: J AOAC 2015 V98-6 (mod) ^b | Units mg/se | Batch: 2302154 | Analyze: 3/9/23 8:23:00 PM | |
|-----------------------|--|-------------|----------------|----------------------------|-------|
| Analyte | Result | Limits | Units | LOQ | Notes |
| CBD per 16g | 0.592 | | mg/16g | 0.507 | |
| CBD-A per 16g | 1.57 | | mg/16g | 0.507 | |
| CBD-Total per 16g | 1.97 | | mg/16g | 0.952 | |
| CBG per 16g | < LOQ | | mg/16g | 0.571 | |
| CBG-A per 16g | < LOQ | | mg/16g | 0.571 | |
| CBG-Total per 16g | < LOQ | | mg/16g | 1.88 | |
| CBN per 16g | < LOQ | | mg/16g | 0.571 | |
| Δ10-THC-9R per 16g | < LOQ | | mg/16g | 0.571 | |
| Δ10-THC-9S per 16g | < LOQ | | mg/16g | 0.571 | |
| Δ10-THC-Total per 16g | < LOQ | | mg/16g | 2.00 | |
| Δ8-THC per 16g | 0.651 | | mg/16g | 0.507 | |
| Δ9-THC per 16g | 42.1 | | mg/16g | 0.507 | |



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| Potency per 16g | | | | | |
|--|--------|-------------|--------|----------------|----------------------------|
| Method: J AOAC 2015 V98-6 (mod) ^b | | Units mg/se | | Batch: 2302154 | Analyze: 3/9/23 8:23:00 PM |
| Analyte | Result | Limits | Units | LOQ | Notes |
| THC-A per 16g | < LOQ | | mg/16g | 0.571 | |
| THC-Total per 16g | 42.1 | | mg/16g | 0.952 | |

| Solvents | | | | | | | | | | | |
|---|--------|------------|-------|---------------|-------|--------------------------------------|--------|--------|------|--------|-------|
| Method: Residual Solvents by GC/MS ^b | | Units µg/g | | Batch 2302150 | | Analyze 03/10/23 10:42 AM | | | | | |
| Analyte | Result | Limits | LOQ | Status | Notes | Analyte | Result | Limits | LOQ | Status | Notes |
| 1-Butanol | < LOQ | 5000 | 500 | pass | | 1-Pentanol | < LOQ | 5000 | 500 | pass | |
| 1,1-Dichloroethane | < LOQ | | 1.00 | | | 1,2-Dichloroethane | < LOQ | | 1.00 | | |
| 1,2-Dimethoxyethane | < LOQ | | 50.0 | | | 1,4-Dioxane | < LOQ | 380 | 100 | pass | |
| 2-Butanol | < LOQ | 5000 | 200 | pass | | 2-Ethoxyethanol | < LOQ | 160 | 30.0 | pass | |
| 2-methyl-1-propanol | < LOQ | | 500 | | | 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-Methyl-(1)-Butanol | < LOQ | | 500 | | |
| 3-Methylpentane | < LOQ | | 30.0 | | | Acetic Acid | < LOQ | | 250 | | |
| Acetone | < LOQ | 5000 | 200 | pass | | Acetonitrile | < LOQ | 410 | 100 | pass | |
| Anisole | < LOQ | | 500 | | | Benzene | < LOQ | 2.00 | 1.00 | pass | |
| Butanes (sum) | < LOQ | 5000 | 400 | pass | | Butyl acetate | < LOQ | | 500 | | |
| Chloroform | < LOQ | | 1.00 | | | Cyclohexane | < LOQ | 3880 | 200 | pass | |
| DMSO | < LOQ | 5000 | 500 | pass | | Ethanol | < LOQ | | 200 | | |
| Ethyl acetate | < LOQ | 5000 | 200 | pass | | Ethyl benzene | < LOQ | | 200 | | |
| Ethyl ether | < LOQ | 5000 | 200 | pass | | Ethyl Formate | < LOQ | | 500 | | |
| Ethylene glycol | < LOQ | 620 | 200 | pass | | Ethylene oxide | < LOQ | 50.0 | 1.00 | pass | |
| Formic Acid | < LOQ | | 250 | | | Hexanes (sum) | < LOQ | 290 | 150 | pass | |
| Isobutyl acetate | < LOQ | 5000 | 500 | 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 | | Methyl-t-butyl ether | < LOQ | | 500 | | |
| Methylacetat | < LOQ | | 500 | | | Methylene chloride | < LOQ | 600 | 1.00 | pass | |
| Methylethylketone | < LOQ | | 500 | | | Methylisobutylketone | < LOQ | | 500 | | |
| Methylpropane (Isobutane) | < LOQ | | 200 | | | n-Butane | < LOQ | | 200 | | |
| n-Heptane | < LOQ | 5000 | 200 | pass | | n-Hexane | < LOQ | | 30.0 | | |
| n-Pentane | < LOQ | | 200 | | | n-Propanol | < LOQ | | 500 | | |
| N,N-dimethylacetamide | < LOQ | 1090 | 200 | pass | | N,N-dimethylformamide | < LOQ | | 200 | | |
| o-Xylene | < LOQ | | 200 | | | Pentanes (sum) | < LOQ | 5000 | 600 | pass | |
| Propane | < LOQ | 5000 | 200 | pass | | Propyl Acetate | < LOQ | | 500 | | |
| Pyridine | < LOQ | 200 | 50.0 | pass | | Sulfolane | < LOQ | 160 | 50.0 | pass | |
| Tetrahydrofuran | < LOQ | 720 | 100 | pass | | Toluene | < LOQ | 890 | 100 | pass | |
| Total Residual Solvents | < LOQ | | 5,000 | | | Total Xylenes | < LOQ | | 400 | | |
| Total Xylenes and Ethyl benzene | < LOQ | 2170 | 600 | pass | | Trichloroethylene | < LOQ | | 1.00 | | |
| Triethylamine | < LOQ | | 500 | | | | | | | | |



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Metals

| Analyte | Result | Limits | Units | LOQ | Batch | Analyzed Method | Status | Notes |
|----------------------|--------|--------|-------|---------|---------|---|--------|-------|
| Arsenic [§] | 0.0206 | 0.200 | mg/kg | 0.0142 | 2302077 | 03/08/23 AOAC 2013.06 (mod.) [§] | pass | |
| Cadmium [§] | < LOQ | 0.200 | mg/kg | 0.0142 | 2302077 | 03/08/23 AOAC 2013.06 (mod.) [§] | pass | |
| Lead [§] | < LOQ | 0.500 | mg/kg | 0.0142 | 2302077 | 03/08/23 AOAC 2013.06 (mod.) [§] | pass | |
| Mercury [§] | < LOQ | 0.100 | mg/kg | 0.00708 | 2302077 | 03/08/23 AOAC 2013.06 (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.

Ⓟ = ISO/IEC 17025:2017 accredited method.

Ⓜ = TNI accredited analyte.

Units of Measure

g = g

µg/g = Microgram per gram

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

mg/16g = Milligram per 16g

% = Percentage of sample

% wt = µg/g divided by 10,000

Approved Signatory

Derrick Tanner
General Manager



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Purchase Order:
Received: 03/06/23 16:17

Revision: 1 Document ID: TBA

Laboratory Quality Control Results

| Organic Acids | | | | Batch ID: 2302114 | | | |
|---------------|--------|-------|-------|---------------------------|-------|-------|-------|
| Method Blank | | | | Laboratory Control Sample | | | |
| Analyte | Result | LOQ | Notes | Result | Spike | Units | %Rec |
| Acetic Acid | <LOQ | < 250 | | 488 | 522 | µg/g | 93.3 |
| Formic Acid | <LOQ | < 250 | | 564 | 557 | µg/g | 101.2 |

| QC- Sample Duplicate | | | | | | Sample ID: LCS | |
|----------------------|--------|-------------|-----|-------|-----|----------------|-------------|
| Analyte | Result | Org. Result | LOQ | Units | RPD | Limits | Accept/Fail |
| Acetic Acid | <LOQ | <LOQ | 250 | µg/g | 0.0 | < 20 | Acceptable |
| Formic Acid | <LOQ | <LOQ | 250 | µg/g | 0.0 | < 20 | Acceptable |

Abbreviations

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

Units of Measure:

µg/g- Microgram per gram or ppm



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

Laboratory Quality Control Results

| Residual Solvents | | | | Batch ID: 2302150 | | | | | |
|-----------------------|--------|-------|-------|---------------------------|-------|-------|-------|----------|-------|
| Method Blank | | | | Laboratory Control Sample | | | | | |
| Analyte | Result | LOQ | Notes | Result | Spike | Units | % Rec | Limits | Notes |
| Propane | ND | < 200 | | 505 | 584 | µg/g | 86.5 | 60 - 120 | |
| Isobutane | ND | < 200 | | 643 | 767 | µg/g | 83.8 | 60 - 120 | |
| Butane | ND | < 200 | | 656 | 782 | µg/g | 84.1 | 60 - 120 | |
| 2,2-Dimethylpropane | ND | < 200 | | 784 | 939 | µg/g | 83.5 | 60 - 120 | |
| Methanol | ND | < 200 | | 1420 | 1610 | µg/g | 88.2 | 60 - 120 | |
| Ethylene Oxide | ND | < 30 | | 48.7 | 57.1 | µg/g | 85.3 | 60 - 120 | |
| 2-Methylbutane | ND | < 200 | | 1510 | 1600 | µg/g | 94.4 | 60 - 120 | |
| Pentane | ND | < 200 | | 1500 | 1610 | µg/g | 93.2 | 60 - 120 | |
| Ethanol | ND | < 200 | | 1370 | 1600 | µg/g | 85.6 | 70 - 130 | |
| Ethyl Ether | ND | < 200 | | 1430 | 1610 | µg/g | 88.8 | 60 - 120 | |
| 2,2-Dimethylbutane | ND | < 30 | | 156 | 173 | µg/g | 90.2 | 60 - 120 | |
| Acetone | ND | < 200 | | 1440 | 1620 | µg/g | 88.9 | 60 - 120 | |
| 2-Propanol | ND | < 200 | | 1380 | 1600 | µg/g | 86.3 | 60 - 120 | |
| Ethyl Formate | ND | < 500 | | 1600 | 1610 | µg/g | 99.4 | 70 - 130 | |
| Acetonitrile | ND | < 100 | | 432 | 488 | µg/g | 88.5 | 60 - 120 | |
| Methyl Acetate | ND | < 500 | | 1370 | 1610 | µg/g | 85.1 | 70 - 130 | |
| 2,3-Dimethylbutane | ND | < 30 | | 148 | 165 | µg/g | 89.7 | 60 - 120 | |
| Dichloromethane | ND | < 60 | | 425 | 487 | µg/g | 87.3 | 60 - 120 | |
| 2-Methylpentane | ND | < 30 | | 143 | 160 | µg/g | 89.4 | 60 - 120 | |
| MTBE | ND | < 500 | | 1350 | 1600 | µg/g | 84.4 | 70 - 130 | |
| 3-Methylpentane | ND | < 30 | | 131 | 161 | µg/g | 81.4 | 60 - 120 | |
| Hexane | ND | < 30 | | 172 | 162 | µg/g | 106.2 | 60 - 120 | |
| 1-Propanol | ND | < 500 | | 1480 | 1620 | µg/g | 91.4 | 70 - 130 | |
| Methyl ethyl ketone | ND | < 500 | | 1440 | 1610 | µg/g | 89.4 | 70 - 130 | |
| Ethyl acetate | ND | < 200 | | 1380 | 1600 | µg/g | 86.3 | 60 - 120 | |
| 2-Butanol | ND | < 200 | | 1380 | 1610 | µg/g | 85.7 | 60 - 120 | |
| Tetrahydrofuran | ND | < 100 | | 408 | 483 | µg/g | 84.1 | 60 - 120 | |
| Cyclohexane | ND | < 200 | | 1350 | 1610 | µg/g | 83.9 | 60 - 120 | |
| 2-methyl-1-propanol | ND | < 500 | | 1430 | 1630 | µg/g | 87.7 | 70 - 130 | |
| Benzene | ND | < 1 | | 4.85 | 4.98 | µg/g | 97.4 | 60 - 120 | |
| Isopropyl Acetate | ND | < 200 | | 1350 | 1610 | µg/g | 83.9 | 60 - 120 | |
| Heptane | ND | < 200 | | 1370 | 1620 | µg/g | 84.6 | 60 - 120 | |
| 1-Butanol | ND | < 500 | | 1470 | 1600 | µg/g | 91.9 | 70 - 130 | |
| Propyl Acetate | ND | < 500 | | 1430 | 1620 | µg/g | 88.3 | 70 - 130 | |
| 1,4-Dioxane | ND | < 100 | | 401 | 494 | µg/g | 81.2 | 60 - 120 | |
| 2-Ethoxyethanol | ND | < 30 | | 138 | 165 | µg/g | 83.6 | 60 - 120 | |
| Methylisobutylketone | ND | < 500 | | 1340 | 1610 | µg/g | 83.2 | 70 - 130 | |
| 3-Methyl-1-butanol | ND | < 500 | | 1440 | 1610 | µg/g | 89.4 | 70 - 130 | |
| Ethylene Glycol | ND | < 200 | | 423 | 488 | µg/g | 87.0 | 60 - 120 | |
| Toluene | ND | < 100 | | 400 | 513 | µg/g | 78.0 | 60 - 120 | |
| Isobutyl Acetate | ND | < 500 | | 1380 | 1600 | µg/g | 86.3 | 70 - 130 | |
| 1-Pentanol | ND | < 500 | | 1440 | 1610 | µg/g | 89.4 | 70 - 130 | |
| Butyl Acetate | ND | < 500 | | 1370 | 1610 | µg/g | 85.1 | 70 - 130 | |
| Ethylbenzene | ND | < 200 | | 788 | 967 | µg/g | 81.5 | 60 - 120 | |
| m,p-Xylene | ND | < 200 | | 806 | 994 | µg/g | 81.1 | 60 - 120 | |
| o-Xylene | ND | < 200 | | 793 | 992 | µg/g | 79.9 | 60 - 120 | |
| Cumene | ND | < 30 | | 133 | 171 | µg/g | 77.8 | 60 - 120 | |
| Anisole | ND | < 500 | | 1250 | 1610 | µg/g | 77.6 | 70 - 130 | |
| DMSO | ND | < 500 | | 1480 | 1610 | µg/g | 91.9 | 70 - 130 | |
| 1,2-dimethoxyethane | ND | < 50 | | 152 | 172 | µg/g | 88.4 | 70 - 130 | |
| Triethylamine | ND | < 500 | | 1400 | 1620 | µg/g | 86.4 | 70 - 130 | |
| N,N-dimethylformamide | ND | < 150 | | 407 | 499 | µg/g | 81.6 | 70 - 130 | |
| N,N-dimethylacetamide | ND | < 150 | | 381 | 491 | µg/g | 77.6 | 70 - 130 | |
| Pyridine | ND | < 50 | | 146 | 171 | µg/g | 85.4 | 70 - 130 | |
| Silfolane | ND | < 50 | | 108 | 160 | µg/g | 67.5 | 70 - 130 | Q6 |
| 1,2-Dichloroethane | ND | < 1 | | 0.813 | 1 | µg/g | 81.3 | 70 - 130 | |
| Chloroform | ND | < 1 | | 0.812 | 1 | µg/g | 81.2 | 70 - 130 | |
| Trichloroethylene | ND | < 1 | | 0.795 | 1 | µg/g | 79.5 | 70 - 130 | |
| 1,1-Dichloroethane | ND | < 1 | | 0.848 | 1 | µg/g | 84.8 | 70 - 130 | |



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

| QC- Sample Duplicate | | Sample ID: 23-002824-0001 | | | | | | |
|-----------------------|--------|---------------------------|-----------|------|--------|--------------|-------|--|
| 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 | 293 | 265 | 200 µg/g | 10.0 | < 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 QClimits. Data acceptable based on remaining QC.

Units of Measure:

µg/g- Microgram per gram or ppm



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



Report Number: 23-002684/D046.R000
Report Date: 03/13/2023
ORELAP#: OR100028
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Laboratory Quality Control Results

J AOAC 2015 V98-6 Batch ID: 2302154

| Laboratory Control Sample | | | | | | | | | |
|---------------------------|-----|--------|-------|-------|-------|--------|-------|------------|-------|
| Analyte | LCS | Result | Spike | Units | % Rec | Limits | | Evaluation | Notes |
| CBDVA | 2 | 0.0368 | 0.033 | % | 111 | 80.0 | - 120 | Acceptable | |
| CBDV | 2 | 0.0392 | 0.035 | % | 111 | 80.0 | - 120 | Acceptable | |
| CBE | 2 | 0.0379 | 0.034 | % | 111 | 80.0 | - 120 | Acceptable | |
| CBD | 1 | 0.0376 | 0.036 | % | 104 | 90.0 | - 110 | Acceptable | |
| CBGA | 1 | 0.0372 | 0.036 | % | 104 | 80.0 | - 120 | Acceptable | |
| CBG | 1 | 0.0392 | 0.038 | % | 104 | 80.0 | - 120 | Acceptable | |
| THCV | 2 | 0.0382 | 0.037 | % | 103 | 90.0 | - 110 | Acceptable | |
| d8THCV | 2 | 0.0378 | 0.033 | % | 113 | 80.0 | - 120 | Acceptable | |
| THCVA | 2 | 0.0394 | 0.036 | % | 109 | 80.0 | - 120 | Acceptable | |
| THCA | 2 | 0.0365 | 0.033 | % | 112 | 80.0 | - 120 | Acceptable | |
| CBN | 1 | 0.0395 | 0.038 | % | 104 | 80.0 | - 120 | Acceptable | |
| exo-THC | 2 | 0.0366 | 0.034 | % | 109 | 80.0 | - 120 | Acceptable | |
| d9THC | 1 | 0.0391 | 0.036 | % | 108 | 90.0 | - 110 | Acceptable | |
| d8THC | 1 | 0.0396 | 0.037 | % | 106 | 90.0 | - 110 | Acceptable | |
| 9S-d10THC | 1 | 0.0395 | 0.037 | % | 106 | 80.0 | - 120 | Acceptable | |
| CBL | 2 | 0.0380 | 0.033 | % | 114 | 80.0 | - 120 | Acceptable | |
| 9R-d10THC | 1 | 0.0367 | 0.036 | % | 101 | 80.0 | - 120 | Acceptable | |
| CB | 2 | 0.0392 | 0.036 | % | 110 | 80.0 | - 120 | Acceptable | |
| THCA | 1 | 0.0380 | 0.036 | % | 105 | 90.0 | - 110 | Acceptable | |
| CBGA | 2 | 0.0390 | 0.035 | % | 112 | 80.0 | - 120 | Acceptable | |
| CBLA | 2 | 0.0384 | 0.035 | % | 110 | 80.0 | - 120 | Acceptable | |
| CBT | 2 | 0.0386 | 0.036 | % | 108 | 80.0 | - 120 | Acceptable | |

| Method Blank | | | | | | |
|--------------|--------|-------|-------|---------|------------|-------|
| Analyte | Result | LOQ | Units | Limits | Evaluation | Notes |
| CBDVA | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| CBDV | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| CBE | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| CBD | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| CBDA | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| CBGA | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| CBG | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| THCV | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| d8THCV | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| THCVA | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| CBN | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| exo-THC | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| d9THC | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| d8THC | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| 9S-d10THC | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| CBL | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| 9R-d10THC | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| CB | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| THCA | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| CBGA | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| CBLA | <LOQ | 0.003 | % | < 0.003 | Acceptable | |
| CBT | <LOQ | 0.003 | % | < 0.003 | Acceptable | |

Abbreviations

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

Units of Measure:

% - Percent



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Laboratory Quality Control Results

| J AOAC 2015 V98-6 | | Batch ID: 2302154 | | | | | | |
|-------------------|---------|---------------------------|-------|-------|------|--------|------------|-------|
| Sample Duplicate | | Sample ID: 23-002464-0001 | | | | | | |
| Analyte | Result | Org. Result | LOQ | Units | RPD | Limits | Evaluation | Notes |
| CBDVA | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| CBDV | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| CBE | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| CBDA | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| CBGA | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| CBG | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| CBD | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| THCV | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| d8THCV | 0.00460 | 0.00470 | 0.003 | % | 2.06 | < 20 | Acceptable | |
| THCVA | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| CBN | 0.00328 | 0.00336 | 0.003 | % | 2.47 | < 20 | Acceptable | |
| exo-THC | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| d9THC | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| d8THC | 0.663 | 0.672 | 0.003 | % | 1.33 | < 20 | Acceptable | |
| 9S-d10THC | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| CBL | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| 9R-d10THC | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| CBC | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| THCA | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| CBCA | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| CBLA | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |
| CBT | <LOQ | <LOQ | 0.003 | % | NA | < 20 | Acceptable | |

Abbreviations

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

Units of Measure:

% - Percent



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Explanation of QC Flag Comments:

| Code | Explanation |
|------|---|
| Q | Matrix interferences affecting spike or surrogate recoveries. |
| Q1 | Quality control result biased high. Only non-detect samples reported. |
| Q2 | Quality control outside QC limits. Data considered estimate. |
| Q3 | Sample concentration greater than four times the amount spiked. |
| Q4 | Non-homogenous sample matrix, affecting RPD result and/or % recoveries. |
| Q5 | Spike results above calibration curve. |
| Q6 | Quality control outside QC limits. Data acceptable based on remaining QC. |
| R | Relative percent difference (RPD) outside control limit. |
| R1 | RPD non-calculable, as sample or duplicate results are less than five times the LOQ. |
| R2 | Sample replicates RPD non-calculable, as only one replicate is within the analytical range. |
| LOQ1 | Quantitation level raised due to low sample volume and/or dilution. |
| LOQ2 | Quantitation level raised due to matrix interference. |
| B | Analyte detected in method blank, but not in associated samples. |
| B1 | The sample concentration is greater than 5 times the blank concentration. |
| B2 | The sample concentration is less than 5 times the blank concentration. |