



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



Report Number: 22-012186/D015.R000
Report Date: 10/17/2022
ORELAP#: OR100028
Purchase Order:
Received: 10/10/22 10:10

Customer: Hawaiian Heart/31 Labs
Product identity: CBD REM
Project Number: CBD100322
Client/Metric ID: .
Laboratory ID: 22-012186-0007

Summary

Potency:

| Analyte | Result (%) | |
|---------|------------|--|
| CBD | 49.7 | |
| CBG | 11.0 | |
| CBE | 8.51 | |
| CBDV | 2.41 | |
| Δ9-THC | 0.264 | |
| CBN | 0.158 | |
| | | |

Residual Solvents:

All analytes passing and less than LOQ.

Pesticides:

All analytes passing and less than LOQ.

Metals:

Less than LOQ for all analytes.



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Purchase Order:
Received: 10/10/22 10:10

Customer: Hawaiian Heart/31 Labs
3505 Cadillac Ave BLDG E
Costa Mesa 92626
United States of America (USA)

Product identity: CBD REM
Project Number: CBD100322
Client/Metric ID: .
Sample Date:
Laboratory ID: 22-012186-0007
Evidence of Cooling: No
Temp: 21 °C
Relinquished by: UPS

Sample Results

| Potency | Method: J AOAC 2015 V98-6 (mod) ^b | | | Units % | Batch: 2208681 | Analyze: 10/12/22 9:18:00 AM |
|---------|--|------------|--------|---------|----------------|------------------------------|
| Analyte | As Received | Dry weight | LOQ | Notes | | |
| CBC | < LOQ | | 0.0720 | | | |
| CBC-A | < LOQ | | 0.0720 | | | |
| CBD | 49.7 | | 0.720 | | | |
| CBD-A | < LOQ | | 0.0720 | | | |
| CBDV | 2.41 | | 0.0720 | | | |
| CBDV-A | < LOQ | | 0.0720 | | | |
| CBE | 8.51 | | 0.0720 | | | |
| CBG | 11.0 | | 0.0720 | | | |
| CBG-A | < LOQ | | 0.0720 | | | |
| CBL | < LOQ | | 0.0720 | | | |
| CBL-A | < LOQ | | 0.0720 | | | |
| CBN | 0.158 | | 0.0720 | | | |
| CBT | < LOQ | | 0.0720 | | | |
| Δ10-THC | < LOQ | | 0.0720 | | | |
| Δ8-THC | < LOQ | | 0.0720 | | | |
| Δ8-THCV | < LOQ | | 0.0720 | | | |
| Δ9-THC | 0.264 | | 0.0720 | | | |
| exo-THC | < LOQ | | 0.0720 | | | |
| THC-A | < LOQ | | 0.0720 | | | |
| THCV | < LOQ | | 0.0720 | | | |
| THCV-A | < LOQ | | 0.0720 | | | |

| Potency | Method: J AOAC 2015 V98-6 (mod) ^b | | | Units | Batch: 2208681 | Analyze: 10/12/22 9:18:00 AM |
|------------|--|------------|------|-------|----------------|------------------------------|
| Analyte | As Received | Dry weight | LOQ | Notes | | |
| CBC-Total | pending | | 1.88 | | | |
| CBD-Total | pending | | 1.88 | | | |
| CBDV-Total | pending | | 1.87 | | | |
| CBG-Total | pending | | 1.88 | | | |
| CBL-Total | pending | | 1.88 | | | |
| THC-Total | pending | | 1.88 | | | |



| Potency | Method: J AOAC 2015 V98-6 (mod) ^p | Units | Batch: 2208681 | Analyze: 10/12/22 9:18:00 AM |
|---------------------------|--|------------|----------------|------------------------------|
| Analyte | As Received | Dry weight | LOQ | Notes |
| THCV-Total | pending | | 1.87 | |
| Total Cannabinoids | pending | | | |

| Solvents | Method: Residual Solvents by GC/MS ^b | Units µg/g | Batch 2208661 | Analyze 10/12/22 11:45 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) | pending | 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) | pending | 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) | pending | 5000 | 600 | pass | | Propane | < LOQ | 5000 | 200 | pass | |
| Tetrahydrofuran | < LOQ | 720 | 100 | pass | | Toluene | < LOQ | 890 | 100 | pass | |
| Total Xylenes | pending | | 400 | | | Total Xylenes and Ethyl benzene | pending | 2170 | 600 | pass | |



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| Pesticides | | | | | | | | | | | |
|---|--------|--------|-------|--------|-------|----------------------------------|--------|--------|-------|--------|-------|
| Method: AOAC 2007.01 & EN 15662 (mod) ^b | | | | | | | | | | | |
| Units mg/kg Batch 2208761 Analyze 10/14/22 10:49 AM | | | | | | | | | | | |
| Analyte | Result | Limits | LOQ | Status | Notes | Analyte | Result | Limits | LOQ | Status | Notes |
| Abamectin [‡] | < LOQ | 0.50 | 0.250 | pass | | Acephate [‡] | < LOQ | 0.40 | 0.250 | pass | |
| Acequinocyl [‡] | < LOQ | 2.0 | 1.00 | pass | | Acetamiprid [‡] | < LOQ | 0.20 | 0.100 | pass | |
| Aldicarb [‡] | < LOQ | 0.40 | 0.200 | pass | | Azoxystrobin [‡] | < LOQ | 0.20 | 0.100 | pass | |
| Bifenazate [‡] | < LOQ | 0.20 | 0.100 | pass | | Bifenthrin [‡] | < LOQ | 0.20 | 0.100 | pass | |
| Boscalid [‡] | < LOQ | 0.40 | 0.200 | pass | | Carbaryl [‡] | < LOQ | 0.20 | 0.100 | pass | |
| Carbofuran [‡] | < LOQ | 0.20 | 0.100 | pass | | Chlorantraniliprole [‡] | < LOQ | 0.20 | 0.100 | pass | |
| Chlorfenapyr [‡] | < LOQ | 1.0 | 0.500 | pass | | Chlorpyrifos [‡] | < LOQ | 0.20 | 0.100 | pass | |
| Clofentezine [‡] | < LOQ | 0.20 | 0.100 | pass | | Cyfluthrin [‡] | < LOQ | 1.0 | 0.500 | pass | |
| Cypermethrin [‡] | < LOQ | 1.0 | 0.500 | pass | | Daminozide [‡] | < LOQ | 1.0 | 0.500 | pass | |
| Diazinon [‡] | < LOQ | 0.20 | 0.100 | pass | | Dichlorvos [‡] | < LOQ | 1.0 | 0.500 | pass | |
| Dimethoate [‡] | < LOQ | 0.20 | 0.100 | pass | | Ethoprophos [‡] | < LOQ | 0.20 | 0.100 | pass | |
| Etofenprox [‡] | < LOQ | 0.40 | 0.200 | pass | | Etoxazole [‡] | < LOQ | 0.20 | 0.100 | pass | |
| Fenoxycarb [‡] | < LOQ | 0.20 | 0.100 | pass | | Fenpyroximate [‡] | < LOQ | 0.40 | 0.200 | pass | |
| Fipronil [‡] | < LOQ | 0.40 | 0.200 | pass | | Flonicamid [‡] | < LOQ | 1.0 | 0.400 | pass | |
| Fludioxonil [‡] | < LOQ | 0.40 | 0.200 | pass | | Hexythiazox [‡] | < LOQ | 1.0 | 0.400 | pass | |
| Imazalil [‡] | < LOQ | 0.20 | 0.100 | pass | | Imidacloprid [‡] | < LOQ | 0.40 | 0.200 | pass | |
| Kresoxim-methyl [‡] | < LOQ | 0.40 | 0.200 | pass | | Malathion [‡] | < LOQ | 0.20 | 0.100 | pass | |
| Metalaxyl [‡] | < LOQ | 0.20 | 0.100 | pass | | Methiocarb [‡] | < LOQ | 0.20 | 0.100 | pass | |
| Methomyl [‡] | < LOQ | 0.40 | 0.200 | pass | | MGK-264 [‡] | < LOQ | 0.20 | 0.100 | pass | |
| Myclobutanil [‡] | < LOQ | 0.20 | 0.100 | pass | | Naled [‡] | < LOQ | 0.50 | 0.250 | pass | |
| Oxamyl [‡] | < LOQ | 1.0 | 0.500 | pass | | Pacllobutrazole [‡] | < LOQ | 0.40 | 0.200 | pass | |
| Parathion-Methyl [‡] | < LOQ | 0.20 | 0.200 | pass | | Permethrin [‡] | < LOQ | 0.20 | 0.100 | pass | |
| Phosmet [‡] | < LOQ | 0.20 | 0.100 | pass | | Piperonyl butoxide [‡] | < LOQ | 2.0 | 1.00 | pass | |
| Prallethrin [‡] | < LOQ | 0.20 | 0.200 | pass | | Propiconazole [‡] | < LOQ | 0.40 | 0.200 | pass | |
| Propoxur [‡] | < LOQ | 0.20 | 0.100 | pass | | Pyrethrin I (total) [‡] | < LOQ | 1.0 | 0.500 | pass | |
| Pyridaben [‡] | < LOQ | 0.20 | 0.100 | pass | | Spinosad [‡] | < LOQ | 0.20 | 0.100 | pass | |
| Spiromesifen [‡] | < LOQ | 0.20 | 0.100 | pass | | Spirotetramat [‡] | < LOQ | 0.20 | 0.100 | pass | |
| Spiroxamine [‡] | < LOQ | 0.40 | 0.200 | pass | | Tebuconazole [‡] | < LOQ | 0.40 | 0.200 | pass | |
| Thiacloprid [‡] | < LOQ | 0.20 | 0.100 | pass | | Thiamethoxam [‡] | < LOQ | 0.20 | 0.100 | pass | |
| Trifloxystrobin [‡] | < LOQ | 0.20 | 0.100 | pass | | | | | | | |

| Metals | | | | | | | | | | |
|---------|--------|--------|-------|--------|---------|---|--------|-------|--|--|
| Analyte | Result | Limits | Units | LOQ | Batch | Analyzed Method | Status | Notes | | |
| Arsenic | < LOQ | 0.200 | mg/kg | 0.0838 | 2208721 | 10/12/22 AOAC 2013.06 (mod.) ^b | pass | | | |
| Cadmium | < LOQ | 0.200 | mg/kg | 0.0838 | 2208721 | 10/12/22 AOAC 2013.06 (mod.) ^b | pass | | | |
| Lead | < LOQ | 0.500 | mg/kg | 0.0838 | 2208721 | 10/12/22 AOAC 2013.06 (mod.) ^b | pass | | | |
| Mercury | < LOQ | 0.100 | mg/kg | 0.0419 | 2208721 | 10/12/22 AOAC 2013.06 (mod.) ^b | pass | | | |



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Mycotoxins

| Analyte | Result | Limits | Units | LOQ | Batch | Analyzed Method | Status | Notes |
|-------------------------------|----------------|--------|-------|------|---------|---|--------|-------|
| Aflatoxin B2 [‡] | < LOQ | | µg/kg | 5.00 | 2208731 | 10/13/22 AOAC 2007.01 & EN 15662 (mod) [‡] | | |
| Aflatoxin B1 [‡] | < LOQ | | µg/kg | 5.00 | 2208731 | 10/13/22 AOAC 2007.01 & EN 15662 (mod) [‡] | | |
| Aflatoxin G1 [‡] | < LOQ | | µg/kg | 5.00 | 2208731 | 10/13/22 AOAC 2007.01 & EN 15662 (mod) [‡] | | |
| Aflatoxin G2 [‡] | < LOQ | | µg/kg | 5.00 | 2208731 | 10/13/22 AOAC 2007.01 & EN 15662 (mod) [‡] | | |
| Ochratoxin A [‡] | < LOQ | 20.0 | µg/kg | 5.00 | 2208731 | 10/13/22 AOAC 2007.01 & EN 15662 (mod) [‡] | pass | |
| Total Aflatoxins [‡] | <i>pending</i> | 20.0 | | | | AOAC 2007.01 & EN 15662 (mod) [‡] | pass | |



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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

µg/g = Microgram per gram

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

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

% = Percentage of sample

% wt = µg/g divided by 10,000

Approved Signatory

Derrick Tanner
General Manager



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**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

| Company: Hawaiian Heart/31 Labs Contact: Paisley Byrnes Address: 3505 Cadillac Ave BLDG E City: Costa Mesa State: CA Zip Code: 92626 <input checked="" type="checkbox"/> Email Results: team@purehawaiianheart.com <input type="checkbox"/> Ph: (949) - 4637643 <i>Billing Contact (if different)</i> Name: DNG LLC Email: team@purehawaiianheart.com Address: 24040 Camino Del Avion City: Dana Point State: CA Zip: 92629 Ph: (949) - 4637643 | | | Analysis Requested <table border="1"> <tr> <td>Potency</td> <td>Pesticides</td> <td>Residual Solvents</td> <td>Heavy Metals</td> <td>Mycotoxins</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> </tr> <tr> <td>✓</td> <td>✓</td> <td>✓</td> <td>✓</td> <td>✓</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> </tr> </table> | | | | | | Potency | Pesticides | Residual Solvents | Heavy Metals | Mycotoxins | | | | | | | | | | | | | | | | | ✓ | ✓ | ✓ | ✓ | ✓ | | | | | | | | | | | | | | | | | | PO Number: Project ID: CBD/D8 Testing Batch ID: CBD100322 Sampled by: Paisley Byrnes Custom Reporting: Source Material: <input checked="" type="checkbox"/> - Ind. Hemp product <input checked="" type="checkbox"/> - Rec. Cannabis Reporting Type: <input type="checkbox"/> - Compliance <input checked="" type="checkbox"/> - R&D Report to: <input type="checkbox"/> - METRC <input type="checkbox"/> - ODA <input type="checkbox"/> - USDA <input type="checkbox"/> - Other: Turnaround time (TAT - Business Days): <input checked="" type="checkbox"/> - 5BD <input type="checkbox"/> - 3BD* <input type="checkbox"/> - 2BD* <i>*Check for availability</i> | | |
|--|------------------------------|-------------------|---|--------------------------|-------------------|--------------|------------|--|---------|------------|-------------------|--------------|------------|--|--|--|--|--|--|--|--|-----------------|----------------|--------------------|--|--|--|--|--|---|---|---|---|---|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|
| Potency | Pesticides | Residual Solvents | Heavy Metals | Mycotoxins | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| ✓ | ✓ | ✓ | ✓ | ✓ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Lab ID | Client Sample Identification | Sample date | Potency | Pesticides | Residual Solvents | Heavy Metals | Mycotoxins | | | | | | | | | | | | | | | Material Type † | Weight (Units) | Comments/Metric ID | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | CBD BDAY | 09/29/22 | ✓ | ✓ | ✓ | ✓ | ✓ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | CBD GELATO | 09/29/22 | ✓ | ✓ | ✓ | ✓ | ✓ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | CBD GRNC | 09/29/22 | ✓ | ✓ | ✓ | ✓ | ✓ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | CBD MNGHZ | 09/29/22 | ✓ | ✓ | ✓ | ✓ | ✓ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | CBD NLTS | 09/29/22 | ✓ | ✓ | ✓ | ✓ | ✓ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | CBD NYC | 09/29/22 | ✓ | ✓ | ✓ | ✓ | ✓ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | CBD REM | 09/29/22 | ✓ | ✓ | ✓ | ✓ | ✓ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | D8 BDAY | 09/29/22 | ✓ | ✓ | ✓ | ✓ | ✓ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | D8 BLUD | 09/29/20 | ✓ | ✓ | ✓ | ✓ | ✓ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | D8 GEL | 09/29/22 | ✓ | ✓ | ✓ | ✓ | ✓ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Signature - Relinquished By: | | Date | Time | Signature - Received By: | | Date | Time | Lab Use Only: | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | | 10/03/20 | 9:30am | RBS | | 10/10/22 | 10:10 | <input checked="" type="checkbox"/> Shipped Via: UPS or <input type="checkbox"/> Client drop off Evidence of cooling: <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No - Temp (°C): 21.0 Sample in good condition: <input type="checkbox"/> Yes <input type="checkbox"/> No Payment: <input type="checkbox"/> Cash <input type="checkbox"/> Check <input type="checkbox"/> CC <input type="checkbox"/> Net: 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

12423 NE Whitaker Way
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info@columbialaboratories.com

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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

| | | | | | | | | | | |
|---|-------------------------------------|--------------------|---------------------------|---------------------------------|--------------------------|---------------------|-------------------|--|-----------------------|---------------------------|
| Company: <u>Hawaiian Heart/31 Labs</u> Contact: <u>Paisley Byrnes</u> Address: <u>3505 Cadillac Ave BLDG E</u> City: <u>Costa Mesa</u> State: <u>CA</u> Zip Code: <u>92626</u> <input checked="" type="checkbox"/> Email Results: <u>team@purehawaiianheart.com</u> <input type="checkbox"/> Ph: <u>(949) - 4637643</u> <i>Billing Contact (if different)</i> Name: <u>DNG LLC</u> Email: <u>team@purehawaiianheart.com</u> Address: <u>24040 Camino Del Avion</u> City: <u>Dana Point</u> State: <u>CA</u> Zip: <u>92629</u> Ph: <u>(949) - 4637643</u> | | | Analysis Requested | | | | | PO Number: _____ Project ID: <u>CBD/D8 Testing</u> Batch ID: <u>CBD100322</u> Sampled by: <u>Paisley Byrnes</u> Custom Reporting: _____ Source Material: <input type="checkbox"/> - Ind. Hemp product <input checked="" type="checkbox"/> - Rec. Cannabis Reporting Type: <input type="checkbox"/> - Compliance <input checked="" type="checkbox"/> - R&D Report to: <input type="checkbox"/> - METRC <input type="checkbox"/> - ODA <input type="checkbox"/> - USDA <input type="checkbox"/> - Other: _____ Turnaround time (TAT - Business Days): <input checked="" type="checkbox"/> - 5BD <input type="checkbox"/> - 3BD* <input type="checkbox"/> - 2BD* <i>*Check for availability</i> | | |
| Lab ID | Client Sample Identification | Sample date | Potency | Pesticides | Residual Solvents | Heavy Metals | Mycotoxins | Material Type † | Weight (Units) | Comments/Metric ID |
| | D8 MNGHZ | 09/29/22 | ✓ | ✓ | ✓ | ✓ | ✓ | | | |
| | D8 MWOW | 09/29/22 | ✓ | ✓ | ✓ | ✓ | ✓ | | | |
| | D8 NLTS | 09/29/22 | ✓ | ✓ | ✓ | ✓ | ✓ | | | |
| | D8 NYCD | 09/29/22 | ✓ | ✓ | ✓ | ✓ | ✓ | | | |
| | D8 REM | 09/29/22 | ✓ | ✓ | ✓ | ✓ | ✓ | | | |
| | D8 ZTZ | 09/29/22 | ✓ | ✓ | ✓ | ✓ | ✓ | | | |
| Signature - Relinquished By: | | Date | Time | Signature - Received By: | | Date | Time | Lab Use Only: | | |
| | | 10/03/20 | 9:30am | | | 10/10/22 | 10:11 | Shipped Via: <u>UPS</u> or <input type="checkbox"/> Client drop off Evidence of cooling: <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No - Temp (°C): <u>21.0</u> Sample in good condition: <input type="checkbox"/> Yes <input type="checkbox"/> No Payment: <input type="checkbox"/> Cash <input type="checkbox"/> Check <input type="checkbox"/> CC <input type="checkbox"/> Net: _____ Prelog storage: _____ | | |

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

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Portland, OR 97230

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

Report Number: 22-012186/D015.R000
Report Date: 10/17/2022
ORELAP#: OR100028
Purchase Order:
Received: 10/10/22 10:10



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

Laboratory Quality Control Results

| Residual Solvents | | | | Batch ID: 2208661 | | | | | |
|-----------------------|--------|-------|-------|---------------------------|-------|-------|-------|----------|-------|
| Method Blank | | | | Laboratory Control Sample | | | | | |
| Analyte | Result | LOQ | Notes | Result | Spike | Units | % Rec | Limits | Notes |
| Propane | ND | < 200 | | 493 | 572 | µg/g | 86.2 | 60 - 120 | |
| Isobutane | ND | < 200 | | 582 | 731 | µg/g | 79.6 | 60 - 120 | |
| Butane | ND | < 200 | | 575 | 731 | µg/g | 78.7 | 60 - 120 | |
| 2,2-Dimethylpropane | ND | < 200 | | 877 | 936 | µg/g | 93.7 | 60 - 120 | |
| Methanol | ND | < 200 | | 1710 | 1650 | µg/g | 103.6 | 60 - 120 | |
| Ethylene Oxide | ND | < 30 | | 54.3 | 56.2 | µg/g | 96.6 | 60 - 120 | |
| 2-Methylbutane | ND | < 200 | | 1580 | 1650 | µg/g | 95.8 | 60 - 120 | |
| Pentane | ND | < 200 | | 1600 | 1650 | µg/g | 97.0 | 60 - 120 | |
| Ethanol | ND | < 200 | | 1610 | 1660 | µg/g | 97.0 | 70 - 130 | |
| Ethyl Ether | ND | < 200 | | 1530 | 1630 | µg/g | 93.9 | 60 - 120 | |
| 2,2-Dimethylbutane | ND | < 30 | | 173 | 189 | µg/g | 91.5 | 60 - 120 | |
| Acetone | ND | < 200 | | 1640 | 1650 | µg/g | 99.4 | 60 - 120 | |
| 2-Propanol | ND | < 200 | | 1610 | 1650 | µg/g | 97.6 | 60 - 120 | |
| Ethyl Formate | ND | < 500 | | 1420 | 1610 | µg/g | 88.2 | 70 - 130 | |
| Acetonitrile | ND | < 100 | | 490 | 504 | µg/g | 97.2 | 60 - 120 | |
| Methyl Acetate | ND | < 500 | | 1620 | 1630 | µg/g | 99.4 | 70 - 130 | |
| 2,3-Dimethylbutane | ND | < 30 | | 168 | 174 | µg/g | 96.6 | 60 - 120 | |
| Dichloromethane | ND | < 60 | | 467 | 521 | µg/g | 89.6 | 60 - 120 | |
| 2-Methylpentane | ND | < 30 | | 165 | 187 | µg/g | 88.2 | 60 - 120 | |
| MTBE | ND | < 500 | | 1560 | 1600 | µg/g | 97.5 | 70 - 130 | |
| 3-Methylpentane | ND | < 30 | | 176 | 188 | µg/g | 93.6 | 60 - 120 | |
| Hexane | ND | < 30 | | 176 | 182 | µg/g | 96.7 | 60 - 120 | |
| 1-Propanol | ND | < 500 | | 1560 | 1610 | µg/g | 96.9 | 70 - 130 | |
| Methylethylketone | ND | < 500 | | 1520 | 1600 | µg/g | 95.0 | 70 - 130 | |
| Ethyl acetate | ND | < 200 | | 1580 | 1630 | µg/g | 96.9 | 60 - 120 | |
| 2-Butanol | ND | < 200 | | 1530 | 1630 | µg/g | 93.9 | 60 - 120 | |
| Tetrahydrofuran | ND | < 100 | | 453 | 506 | µg/g | 89.9 | 60 - 120 | |
| Cyclohexane | ND | < 200 | | 1480 | 1640 | µg/g | 90.2 | 60 - 120 | |
| 2-methyl-1-propanol | ND | < 500 | | 1530 | 1620 | µg/g | 94.4 | 70 - 130 | |
| Benzene | ND | < 1 | | 4.3 | 4.93 | µg/g | 87.2 | 60 - 120 | |
| Isopropyl Acetate | ND | < 200 | | 1590 | 1640 | µg/g | 97.0 | 60 - 120 | |
| Heptane | ND | < 200 | | 1580 | 1630 | µg/g | 96.9 | 60 - 120 | |
| 1-Butanol | ND | < 500 | | 1600 | 1600 | µg/g | 100.0 | 70 - 130 | |
| Propyl Acetate | ND | < 500 | | 1610 | 1620 | µg/g | 99.4 | 70 - 130 | |
| 1,4-Dioxane | ND | < 100 | | 426 | 493 | µg/g | 86.4 | 60 - 120 | |
| 2-Ethoxyethanol | ND | < 30 | | 173 | 171 | µg/g | 101.2 | 60 - 120 | |
| Methylisobutylketone | ND | < 500 | | 1600 | 1620 | µg/g | 98.8 | 70 - 130 | |
| 3-Methyl-1-butanol | ND | < 500 | | 1610 | 1610 | µg/g | 100.0 | 70 - 130 | |
| Ethylene Glycol | ND | < 200 | | 443 | 494 | µg/g | 89.7 | 60 - 120 | |
| Toluene | ND | < 100 | | 435 | 506 | µg/g | 86.0 | 60 - 120 | |
| Isobutyl Acetate | ND | < 500 | | 1640 | 1620 | µg/g | 101.2 | 70 - 130 | |
| 1-Pentanol | ND | < 500 | | 1660 | 1610 | µg/g | 103.1 | 70 - 130 | |
| Butyl Acetate | ND | < 500 | | 1670 | 1610 | µg/g | 103.7 | 70 - 130 | |
| Ethylbenzene | ND | < 200 | | 867 | 998 | µg/g | 87.0 | 60 - 120 | |
| m,p-Xylene | ND | < 200 | | 885 | 1010 | µg/g | 87.6 | 60 - 120 | |
| o-Xylene | ND | < 200 | | 854 | 979 | µg/g | 87.2 | 60 - 120 | |
| Cumene | ND | < 30 | | 160 | 188 | µg/g | 85.1 | 60 - 120 | |
| Anisole | ND | < 500 | | 1530 | 1610 | µg/g | 95.0 | 70 - 130 | |
| DMSO | ND | < 500 | | 1520 | 1600 | µg/g | 95.0 | 70 - 130 | |
| 1,2-dimethoxyethane | ND | < 50 | | 189 | 190 | µg/g | 99.5 | 70 - 130 | |
| Triethylamine | ND | < 500 | | 1550 | 1610 | µg/g | 96.3 | 70 - 130 | |
| N,N-dimethylformamide | ND | < 150 | | 518 | 496 | µg/g | 104.4 | 70 - 130 | |
| N,N-dimethylacetamide | ND | < 150 | | 482 | 483 | µg/g | 99.8 | 70 - 130 | |
| Pyridine | ND | < 50 | | 155 | 167 | µg/g | 92.8 | 70 - 130 | |
| 1,2-Dichloroethane | ND | < 1 | | 1.04 | 1 | µg/g | 104.0 | 70 - 130 | |
| Chloroform | ND | < 1 | | 1.01 | 1 | µg/g | 101.0 | 70 - 130 | |
| Trichloroethylene | ND | < 1 | | 1.01 | 1 | µg/g | 101.0 | 70 - 130 | |
| 1,1-Dichloroethane | ND | < 1 | | 0.96 | 1 | µg/g | 96.0 | 70 - 130 | |



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Received: 10/10/22 10:10



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

QC- Sample Duplicate Sample ID: 22-012140-0001

| 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 | |
| 1,2-Dichloroethane | ND | ND | 1 µg/g | 0.0 | < 20 | Acceptable | |
| Chloroform | ND | ND | 1 µg/g | 0.0 | < 20 | Acceptable | |
| Trichloroethylene | ND | ND | 1 µg/g | 0.0 | < 20 | Acceptable | |
| 1,1-Dichloroethane | ND | ND | 1 µg/g | 0.0 | < 20 | Acceptable | |

Abbreviations

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

Units of Measure:

µg/g- Microgram per gram or ppm



12423 NE Whitaker Way
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Report Number: 22-012186/D015.R000
Report Date: 10/17/2022
ORELAP#: OR100028
Purchase Order:
Received: 10/10/22 10:10

Revision: 1 Document ID: 7148
Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

JAOAC2015 V986 Batch ID: 2208681

| Analyte | LCS | Result | Spike | Units | % Rec | Limits | Evaluation | Notes |
|---------|-----|--------|-------|-------|-------|------------|------------|-------|
| CBDA | 2 | 0.114 | 0.103 | % | 110 | 80.0 - 120 | Acceptable | |
| CBDV | 2 | 0.120 | 0.110 | % | 110 | 80.0 - 120 | Acceptable | |
| CBE | 2 | 0.115 | 0.105 | % | 110 | 80.0 - 120 | Acceptable | |
| CBDA | 1 | 0.0972 | 0.100 | % | 96.8 | 90.0 - 110 | Acceptable | |
| CBSA | 1 | 0.0983 | 0.101 | % | 97.7 | 80.0 - 120 | Acceptable | |
| CBS | 1 | 0.102 | 0.103 | % | 99.0 | 80.0 - 120 | Acceptable | |
| CB | 1 | 0.102 | 0.103 | % | 98.9 | 90.0 - 110 | Acceptable | |
| THCV | 2 | 0.123 | 0.113 | % | 109 | 80.0 - 120 | Acceptable | |
| δ8THCV | 2 | 0.121 | 0.110 | % | 109 | 80.0 - 120 | Acceptable | |
| THCVA | 2 | 0.111 | 0.101 | % | 110 | 80.0 - 120 | Acceptable | |
| CBN | 1 | 0.101 | 0.101 | % | 99.6 | 90.0 - 110 | Acceptable | |
| exo-THC | 2 | 0.111 | 0.103 | % | 109 | 80.0 - 120 | Acceptable | |
| δ9THC | 1 | 0.104 | 0.104 | % | 100 | 90.0 - 110 | Acceptable | |
| δ8THC | 1 | 0.106 | 0.100 | % | 106 | 90.0 - 110 | Acceptable | |
| CB | 2 | 0.108 | 0.099 | % | 110 | 80.0 - 120 | Acceptable | |
| δ10THC | 1 | 0.0939 | 0.096 | % | 98.1 | 80.0 - 120 | Acceptable | |
| CB | 2 | 0.120 | 0.108 | % | 111 | 80.0 - 120 | Acceptable | |
| THCA | 1 | 0.0965 | 0.099 | % | 97.0 | 90.0 - 110 | Acceptable | |
| CBA | 2 | 0.116 | 0.105 | % | 110 | 80.0 - 120 | Acceptable | |
| CBA | 2 | 0.0613 | 0.056 | % | 109 | 80.0 - 120 | Acceptable | |
| CB | 2 | 0.122 | 0.112 | % | 109 | 80.0 - 120 | Acceptable | |

| Analyte | Result | LOQ | Units | Limits | Evaluation | Notes |
|---------|--------|--------|-------|----------|------------|-------|
| CBDA | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| CBDV | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| CBE | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| CBDA | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| CBSA | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| CBS | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| CB | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| THCV | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| δ8THCV | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| THCVA | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| CBN | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| exo-THC | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| δ9THC | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| δ8THC | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| CB | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| δ10THC | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| CB | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| THCA | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| CBA | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| CBA | <LOQ | 0.0077 | % | < 0.0077 | Acceptable | |
| CB | <LOQ | 0.0077 | % | < 0.0077 | 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
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Report Number: 22-012186/D015.R000
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ORELAP#: OR100028
Purchase Order:
Received: 10/10/22 10:10



Revision: 1 Document ID: 7148
 Legacy ID: Worksheet Validated 04/20/2021

Laboratory Quality Control Results

| JAOAC2015 V986 | | Batch ID: 2208681 | | | | | | |
|------------------|--------|--------------------------|--------|-------|-------|--------|------------|-------|
| Sample Duplicate | | Sample ID: 22-0121770001 | | | | | | |
| Analyte | Result | Org. Result | LOQ | Units | RPD | Limits | Evaluation | Notes |
| CBDA | <LOQ | <LOQ | 0.0077 | % | NA | < 20 | Acceptable | |
| CBDV | 0.283 | 0.289 | 0.0077 | % | 1.98 | < 20 | Acceptable | |
| CBF | <LOQ | <LOQ | 0.0077 | % | NA | < 20 | Acceptable | |
| CBDA | <LOQ | <LOQ | 0.0077 | % | NA | < 20 | Acceptable | |
| CBGA | <LOQ | <LOQ | 0.0077 | % | NA | < 20 | Acceptable | |
| CBG | 1.23 | 1.23 | 0.0077 | % | 0.136 | < 20 | Acceptable | |
| CB | 58.2 | 57.6 | 0.0077 | % | 1.15 | < 20 | Acceptable | |
| THCV | <LOQ | <LOQ | 0.0077 | % | NA | < 20 | Acceptable | |
| δ8THCV | <LOQ | <LOQ | 0.0077 | % | NA | < 20 | Acceptable | |
| THCVA | <LOQ | <LOQ | 0.0077 | % | NA | < 20 | Acceptable | |
| CBN | 0.356 | 0.355 | 0.0077 | % | 0.517 | < 20 | Acceptable | |
| exo-THC | <LOQ | <LOQ | 0.0077 | % | NA | < 20 | Acceptable | |
| δ9THC | 1.15 | 1.15 | 0.0077 | % | 0.135 | < 20 | Acceptable | |
| δ8THC | <LOQ | <LOQ | 0.0077 | % | NA | < 20 | Acceptable | |
| CB | <LOQ | <LOQ | 0.0077 | % | NA | < 20 | Acceptable | |
| δ10THC | <LOQ | <LOQ | 0.0077 | % | NA | < 20 | Acceptable | |
| CB | 2.32 | 2.19 | 0.0077 | % | 5.62 | < 20 | Acceptable | |
| THCA | <LOQ | <LOQ | 0.0077 | % | NA | < 20 | Acceptable | |
| CBGA | <LOQ | <LOQ | 0.0077 | % | NA | < 20 | Acceptable | |
| CBGA | <LOQ | <LOQ | 0.0077 | % | NA | < 20 | Acceptable | |
| CB | 0.631 | 0.626 | 0.0077 | % | 0.829 | < 20 | Acceptable | |

Abbreviations

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

Units of Measure:



12423 NE Whitaker Way
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Report Number: 22-012186/D015.R000
Report Date: 10/17/2022
ORELAP#: OR100028
Purchase Order:
Received: 10/10/22 10:10

Revision: 3 Document ID: 3120
Legacy ID: CFL-C21 Worksheet Validated 10/30/2020

Laboratory Pesticide Quality Control Results

| AOAC2007.1 &EN 15662 | | Units: mg/Kg | | | Batch ID 2208761 | | | |
|----------------------|---------------------------|--------------|-------|------------|------------------|----------|--------|-------|
| Method Blank | Laboratory Control Sample | | | | | | | |
| Analyte | Blank Result | Blank Limits | Notes | LCS Result | LCS Spk | LCS % Re | Limits | Notes |
| Abamectin | 0.000 | < 0.250 | | 0.880 | 1.000 | 88.0 | 50.0 | 150 |
| Acephate | 0.000 | < 0.250 | | 0.824 | 1.000 | 82.4 | 60.0 | 120 |
| Acetamiprid | 0.000 | < 1.000 | | 3.353 | 4.000 | 83.8 | 40.0 | 160 |
| Acetamiprid | 0.000 | < 0.100 | | 0.368 | 0.400 | 91.9 | 60.0 | 120 |
| Aldicarb | 0.000 | < 0.200 | | 0.717 | 0.800 | 89.7 | 60.0 | 120 |
| Azoxystrobin | 0.000 | < 0.100 | | 0.363 | 0.400 | 90.8 | 60.0 | 120 |
| Bifenazate | 0.000 | < 0.100 | | 0.362 | 0.400 | 90.6 | 60.0 | 120 |
| Bifenthrin | 0.000 | < 0.100 | | 0.350 | 0.400 | 87.5 | 50.0 | 150 |
| Boscalid | 0.000 | < 0.200 | | 0.719 | 0.800 | 89.8 | 60.0 | 120 |
| Carbaryl | 0.000 | < 0.100 | | 0.367 | 0.400 | 91.8 | 60.0 | 120 |
| Carbofuran | 0.000 | < 0.100 | | 0.371 | 0.400 | 92.7 | 60.0 | 120 |
| Chlorantraniliprole | 0.000 | < 0.100 | | 0.350 | 0.400 | 87.4 | 60.0 | 120 |
| Chlorfenapyr | 0.000 | < 0.500 | | 1.607 | 2.000 | 80.3 | 60.0 | 120 |
| Chlorpyrifos | 0.000 | < 0.100 | | 0.368 | 0.400 | 92.0 | 60.0 | 120 |
| Clofentazine | 0.000 | < 0.100 | | 0.357 | 0.400 | 89.2 | 60.0 | 120 |
| Cyfluthrin | 0.000 | < 0.500 | | 1.919 | 2.000 | 95.9 | 50.0 | 150 |
| Cypermethrin | 0.000 | < 0.500 | | 1.806 | 2.000 | 90.3 | 50.0 | 150 |
| Daminozide | 0.000 | < 0.500 | | 1.886 | 2.000 | 94.3 | 60.0 | 120 |
| Diazinon | 0.000 | < 0.100 | | 0.372 | 0.400 | 93.1 | 60.0 | 120 |
| Dichlorvos | 0.000 | < 0.500 | | 1.842 | 2.000 | 92.1 | 60.0 | 120 |
| Dimethoate | 0.000 | < 0.100 | | 0.369 | 0.400 | 92.4 | 60.0 | 120 |
| Ethoprophos | 0.000 | < 0.100 | | 0.347 | 0.400 | 86.6 | 60.0 | 120 |
| Etofenprox | 0.000 | < 0.200 | | 0.707 | 0.800 | 88.4 | 50.0 | 150 |
| Etoxazole | 0.000 | < 0.100 | | 0.361 | 0.400 | 90.2 | 60.0 | 120 |
| Fenoxycarb | 0.000 | < 0.100 | | 0.355 | 0.400 | 88.8 | 60.0 | 120 |
| Fenpyroximate | 0.000 | < 0.200 | | 0.708 | 0.800 | 88.5 | 60.0 | 120 |
| Fipronil | 0.000 | < 0.200 | | 0.719 | 0.800 | 89.9 | 60.0 | 120 |
| Fonicamid | 0.000 | < 0.250 | | 0.944 | 1.000 | 94.4 | 60.0 | 120 |
| Fludioxonil | 0.000 | < 0.200 | | 0.805 | 0.800 | 100.6 | 50.0 | 150 |
| Hexythiazox | 0.000 | < 0.250 | | 0.893 | 1.000 | 89.3 | 60.0 | 120 |
| Imazalil | 0.000 | < 0.100 | | 0.371 | 0.400 | 92.6 | 60.0 | 120 |
| Imidacloprid | 0.000 | < 0.200 | | 0.727 | 0.800 | 90.8 | 60.0 | 120 |
| Kresoxim-methyl | 0.000 | < 0.200 | | 0.749 | 0.800 | 93.7 | 60.0 | 120 |
| Malathion | 0.000 | < 0.100 | | 0.364 | 0.400 | 91.0 | 60.0 | 120 |
| Metaxalyl | 0.000 | < 0.100 | | 0.369 | 0.400 | 92.1 | 60.0 | 120 |
| Methiocarb | 0.000 | < 0.100 | | 0.367 | 0.400 | 91.8 | 60.0 | 120 |
| Methomyl | 0.000 | < 0.200 | | 0.666 | 0.800 | 83.2 | 60.0 | 120 |
| MGK-264 | 0.000 | < 0.100 | | 0.364 | 0.400 | 91.1 | 50.0 | 150 |
| Myclobutanil | 0.000 | < 0.100 | | 0.360 | 0.400 | 90.1 | 60.0 | 120 |
| Naled | 0.000 | < 0.250 | | 0.877 | 1.000 | 87.7 | 50.0 | 150 |
| Oxamyl | 0.000 | < 0.500 | | 1.911 | 2.000 | 95.5 | 60.0 | 120 |
| Paclotrazole | 0.000 | < 0.200 | | 0.715 | 0.800 | 89.4 | 60.0 | 120 |
| Parathion-Methyl | 0.000 | < 0.200 | | 0.728 | 0.800 | 91.0 | 50.0 | 150 |
| Permethrin | 0.000 | < 0.100 | | 0.355 | 0.400 | 88.8 | 50.0 | 150 |
| Phosmet | 0.000 | < 0.100 | | 0.351 | 0.400 | 87.9 | 50.0 | 150 |
| Piperonyl butoxide | 0.000 | < 0.500 | | 1.739 | 2.000 | 87.0 | 60.0 | 120 |
| Prallethrin | 0.000 | < 0.100 | | 0.368 | 0.400 | 91.9 | 60.0 | 120 |
| Propiconazole | 0.000 | < 0.200 | | 0.734 | 0.800 | 91.7 | 60.0 | 120 |
| Propoxur | 0.000 | < 0.100 | | 0.374 | 0.400 | 93.5 | 60.0 | 120 |
| Pyrethrin (Summe) | 0.000 | < 0.100 | | 0.377 | 0.413 | 91.2 | 60.0 | 120 |
| Pyridaben | 0.000 | < 0.100 | | 0.349 | 0.400 | 87.3 | 50.0 | 150 |
| Spirosad | 0.000 | < 0.100 | | 0.319 | 0.388 | 82.2 | 50.0 | 150 |
| Spiromesifen | 0.000 | < 0.100 | | 0.363 | 0.400 | 90.7 | 60.0 | 120 |
| Spirotetramat | 0.000 | < 0.100 | | 0.369 | 0.400 | 92.2 | 60.0 | 120 |
| Spiroxamine | 0.000 | < 0.200 | | 0.732 | 0.800 | 91.4 | 60.0 | 120 |
| Tebuconazole | 0.000 | < 0.200 | | 0.723 | 0.800 | 90.4 | 60.0 | 120 |
| Thiacloprid | 0.000 | < 0.100 | | 0.374 | 0.400 | 93.5 | 60.0 | 120 |
| Thiamethoxam | 0.000 | < 0.100 | | 0.359 | 0.400 | 89.8 | 60.0 | 120 |
| Trifloxystrobin | 0.000 | < 0.100 | | 0.360 | 0.400 | 89.9 | 60.0 | 120 |



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Revision: 3 Document ID: 3120
Legacy ID: CFL-C21 Worksheet Validated 10/30/2020

Laboratory Pesticide Quality Control Results

| AOAC2007.1 & EN 15662 | | | | | | | | | | |
|--|--------|--------|---------|-------|-------------------------|-------|--------|---------|----------|------------------|
| Units: mg/Kg | | | | | | | | | | Batch ID 2208761 |
| Matrix Spke/Matrix Spke Duplicate Recoveries | | | | | Sample ID: 220121770001 | | | | | |
| Analyte | Result | MS Res | MSD Res | Spike | RFD% | Limit | MS% Re | MSD% Re | Limits | Notes |
| Abamectin | 0.000 | 0.638 | 0.648 | 1.000 | 1.6% | < 30 | 63.8% | 64.8% | 50 - 150 | |
| Acephate | 0.000 | 0.605 | 0.654 | 1.000 | 7.8% | < 30 | 60.5% | 65.4% | 50 - 150 | |
| Acequinocyl | 0.000 | 1.501 | 1.570 | 4.000 | 4.5% | < 30 | 37.5% | 39.2% | 50 - 150 | Q |
| Acetamiprid | 0.000 | 0.347 | 0.354 | 0.400 | 2.0% | < 30 | 86.8% | 88.5% | 50 - 150 | |
| Aldicarb | 0.000 | 0.669 | 0.718 | 0.800 | 7.1% | < 30 | 83.6% | 89.7% | 50 - 150 | |
| Azoxystrobin | 0.000 | 0.269 | 0.268 | 0.400 | 0.7% | < 30 | 67.3% | 66.9% | 50 - 150 | |
| Bifenazate | 0.000 | 0.300 | 0.319 | 0.400 | 6.2% | < 30 | 74.9% | 79.7% | 50 - 150 | |
| Bifenthrin | 0.000 | 0.200 | 0.217 | 0.400 | 8.3% | < 30 | 49.9% | 54.2% | 50 - 150 | Q |
| Boscalid | 0.069 | 0.683 | 0.645 | 0.800 | 6.4% | < 30 | 76.8% | 72.0% | 50 - 150 | |
| Carbaryl | 0.000 | 0.300 | 0.297 | 0.400 | 1.1% | < 30 | 74.9% | 74.1% | 50 - 150 | |
| Carbofuran | 0.000 | 0.278 | 0.280 | 0.400 | 0.9% | < 30 | 69.4% | 70.0% | 50 - 150 | |
| Chlorantraniliprole | 0.000 | 0.303 | 0.312 | 0.400 | 3.0% | < 30 | 75.7% | 78.0% | 50 - 150 | |
| Chlorfenapyr | 0.000 | 1.231 | 1.067 | 2.000 | 14.2% | < 30 | 61.5% | 53.4% | 50 - 150 | |
| Chlorpyrifos | 0.000 | 2.487 | 2.665 | 0.400 | 6.9% | < 30 | 621.8% | 666.2% | 50 - 150 | Q |
| Clofentezine | 0.000 | 0.229 | 0.235 | 0.400 | 2.7% | < 30 | 57.1% | 58.7% | 50 - 150 | |
| Cyfluthrin | 0.000 | 0.871 | 0.940 | 2.000 | 7.6% | < 30 | 43.5% | 47.0% | 30 - 150 | |
| Cypermethrin | 0.000 | 0.759 | 0.817 | 2.000 | 7.5% | < 30 | 37.9% | 40.9% | 50 - 150 | Q |
| Daminozide | 0.000 | 0.290 | 0.318 | 2.000 | 9.3% | < 30 | 14.5% | 15.9% | 30 - 150 | Q |
| Diazinon | 0.000 | 0.205 | 0.205 | 0.400 | 0.4% | < 30 | 51.3% | 51.1% | 50 - 150 | |
| Dichlorvos | 0.000 | 1.564 | 1.589 | 2.000 | 1.6% | < 30 | 78.2% | 79.5% | 50 - 150 | |
| Dimethoate | 0.000 | 0.352 | 0.358 | 0.400 | 1.7% | < 30 | 88.1% | 89.6% | 50 - 150 | |
| Ethoprophos | 0.000 | 0.250 | 0.250 | 0.400 | 0.2% | < 30 | 62.4% | 62.5% | 50 - 150 | |
| Etofenprox | 0.000 | 0.347 | 0.351 | 0.800 | 1.3% | < 30 | 43.3% | 43.9% | 50 - 150 | Q |
| Etoxazole | 0.000 | 0.275 | 0.271 | 0.400 | 1.6% | < 30 | 68.8% | 67.7% | 50 - 150 | |
| Fenoxycarb | 0.000 | 0.259 | 0.269 | 0.400 | 3.7% | < 30 | 64.7% | 67.2% | 50 - 150 | |
| Fenpyroximate | 0.000 | 0.308 | 0.343 | 0.800 | 10.6% | < 30 | 38.5% | 42.8% | 50 - 150 | Q |
| Fipronil | 0.000 | 0.519 | 0.550 | 0.800 | 5.7% | < 30 | 64.9% | 68.7% | 50 - 150 | |
| Fonicamid | 0.000 | 0.927 | 0.910 | 1.000 | 1.9% | < 30 | 92.7% | 91.0% | 50 - 150 | |
| Fludioxonil | 0.000 | 0.809 | 0.777 | 0.800 | 4.0% | < 30 | 101.1% | 97.1% | 50 - 150 | |
| Hexythiazox | 0.000 | 0.658 | 0.665 | 1.000 | 1.1% | < 30 | 65.8% | 66.5% | 50 - 150 | |
| Imazalil | 0.000 | 0.326 | 0.340 | 0.400 | 4.3% | < 30 | 81.5% | 85.1% | 50 - 150 | |
| Imidacloprid | 0.000 | 0.773 | 0.808 | 0.800 | 4.4% | < 30 | 96.6% | 101.0% | 50 - 150 | |
| Kresoxim-methyl | 0.000 | 0.477 | 0.486 | 0.800 | 1.8% | < 30 | 59.7% | 60.7% | 50 - 150 | |
| Malathion | 0.000 | 0.236 | 0.241 | 0.400 | 2.3% | < 30 | 58.9% | 60.3% | 50 - 150 | |
| Metaxalyl | 0.000 | 0.275 | 0.284 | 0.400 | 3.3% | < 30 | 68.8% | 71.1% | 50 - 150 | |
| Methiocarb | 0.000 | 0.282 | 0.281 | 0.400 | 0.6% | < 30 | 70.5% | 70.1% | 50 - 150 | |
| Methomyl | 0.000 | 0.748 | 0.630 | 0.800 | 17.2% | < 30 | 93.5% | 78.7% | 50 - 150 | |
| MGK-264 | 0.000 | 0.154 | 0.158 | 0.400 | 2.2% | < 30 | 38.5% | 39.4% | 50 - 150 | Q |
| Myclobutanil | 0.000 | 0.236 | 0.279 | 0.400 | 16.5% | < 30 | 59.0% | 69.7% | 50 - 150 | |
| Naled | 0.000 | 0.634 | 0.644 | 1.000 | 1.5% | < 30 | 63.4% | 64.4% | 50 - 150 | |
| Oxamyl | 0.000 | 1.904 | 1.702 | 2.000 | 11.2% | < 30 | 95.2% | 85.1% | 50 - 150 | |
| Pacllobutrazole | 0.000 | 0.575 | 0.602 | 0.800 | 4.6% | < 30 | 71.8% | 75.2% | 50 - 150 | |
| Parathion-Methyl | 0.000 | 0.454 | 0.359 | 0.800 | 23.4% | < 30 | 56.7% | 44.9% | 30 - 150 | |
| Permethrin | 0.000 | 0.212 | 0.213 | 0.400 | 0.5% | < 30 | 53.0% | 53.2% | 50 - 150 | |
| Phosmet | 0.000 | 0.283 | 0.293 | 0.400 | 3.4% | < 30 | 70.7% | 73.2% | 50 - 150 | |
| Piperonyl butoxide | 0.000 | 1.325 | 1.377 | 2.000 | 3.8% | < 30 | 66.2% | 68.8% | 50 - 150 | |
| Prallethrin | 0.000 | 0.197 | 0.193 | 0.400 | 1.6% | < 30 | 49.1% | 48.3% | 50 - 150 | Q |
| Propiconazole | 0.000 | 0.622 | 0.634 | 0.800 | 2.0% | < 30 | 77.8% | 79.3% | 50 - 150 | |
| Propoxur | 0.000 | 0.297 | 0.298 | 0.400 | 0.3% | < 30 | 74.4% | 74.6% | 50 - 150 | |
| Pyrethrin (Summe) | 0.000 | 0.320 | 0.323 | 0.413 | 0.9% | < 30 | 77.5% | 78.3% | 50 - 150 | |
| Pyridaben | 0.000 | 0.226 | 0.231 | 0.400 | 2.4% | < 30 | 56.5% | 57.9% | 50 - 150 | |
| Spinosad | 0.000 | 0.230 | 0.230 | 0.388 | 0.1% | < 30 | 59.3% | 59.4% | 50 - 150 | |
| Spiromesifen | 0.000 | 0.267 | 0.247 | 0.400 | 7.7% | < 30 | 66.7% | 61.8% | 50 - 150 | |
| Spirotetramat | 0.000 | 0.506 | 0.525 | 0.400 | 3.7% | < 30 | 126.5% | 131.3% | 50 - 150 | |
| Spiroxamine | 0.000 | 0.576 | 0.625 | 0.800 | 8.2% | < 30 | 72.0% | 78.2% | 50 - 150 | |
| Tebuconazole | 0.000 | 0.601 | 0.649 | 0.800 | 7.7% | < 30 | 75.1% | 81.1% | 50 - 150 | |
| Thiacloprid | 0.000 | 0.334 | 0.344 | 0.400 | 3.0% | < 30 | 83.4% | 86.0% | 50 - 150 | |
| Thiamethoxam | 0.000 | 0.358 | 0.330 | 0.400 | 8.2% | < 30 | 89.6% | 82.5% | 50 - 150 | |
| Trifloxystrobin | 0.000 | 0.243 | 0.247 | 0.400 | 1.5% | < 30 | 60.9% | 61.8% | 50 - 150 | |



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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. |