

## CONSOLIDATED TEST RESULTS SUMMARY

Please see the following pages for full test results.

|  |                                       |  |
|--|---------------------------------------|--|
| <b>BULK SKU</b> SG50                     | <b>BATCH #</b> EC04                   | <b>LOQ:</b> Limit Of Quantitation<br><b>LOD:</b> Limit Of Detection<br><br>1 g = 10 <sup>-3</sup> kg = 10 <sup>3</sup> mg = 10 <sup>6</sup> µg<br>1 mg/kg = 1 ppm = 1000 ppb |
| <b>PRODUCT NAME</b> CBD Softgels - 50 mg | <b>SERVING SIZE</b> 1 softgel (~0.4g) |  |
| <b>LABORATORY:</b> Columbia Laboratories | <b>OREGON ACCREDITATION:</b> OR100028 |  |

| POTENCY                            | PER SERVING      | PER GRAM   | Percent |
|------------------------------------|------------------|------------|---------|
| Cannabidiol (CBD)                  | 51.60 mg/serving | 129.0 mg/g | 12.90 % |
| Total THC (d9-THC, THCA)           | 0.40 mg/serving  | 1.01 mg/g  | 0.101 % |
| Cannabigerol (CBG)                 | <LOQ mg/serving  | <LOQ mg/g  | <LOQ %  |
| Cannabinol (CBN)                   | <LOQ mg/serving  | <LOQ mg/g  | <LOQ %  |
| Cannabichromene (CBC)              | 0.35 mg/serving  | 0.89 mg/g  | 0.089 % |
| Tetrahydrocannabinolic Acid (THCA) | <LOQ mg/serving  | <LOQ mg/g  | <LOQ %  |
| Delta-9-THC (d9-THC)               | 0.40 mg/serving  | 1.01 mg/g  | 0.101 % |
| Delta-8-THC (d8-THC)               | <LOQ mg/serving  | <LOQ mg/g  | <LOQ %  |

| HEAVY METALS | PER SERVING     | PER GRAM  | REGULATORY ACTION LEVEL   |
|--------------|-----------------|-----------|---------------------------|
| Arsenic      | <LOQ µg/serving | <LOQ µg/g | 10 µg/day <sup>[1]</sup>  |
| Cadmium      | <LOQ µg/serving | <LOQ µg/g | 4.1 µg/day <sup>[1]</sup> |
| Lead         | <LOQ µg/serving | <LOQ µg/g | 3.5 µg/day <sup>[2]</sup> |
| Mercury      | <LOQ µg/serving | <LOQ µg/g | 2 µg/day <sup>[1]</sup>   |

| PESTICIDES   | REGULATORY ACTION LEVEL |
|--|-------------------------|
| None of the other 59 pesticides tested found above limit of detection in the sample. | 10 ppb <sup>[1]</sup>   |

| RESIDUAL SOLVENTS | Results | REGULATORY ACTION LEVEL |
|-------------------|---------|-------------------------|
| Ethanol           | <LOQ    | 50,000 mg/day           |
| Heptane           | <LOQ    | 50,000 mg/day           |

None of the 34 residual solvents tested found above limit of quantitation in the sample.

| MICROBIAL    | PASS/FAIL |
|--------------|-----------|
| Yeast & Mold | Pass      |
| Coliform     | Pass      |

| TERPENES            | % OF SAMPLE |
|---------------------|-------------|
| Farnesene           | <LOQ %      |
| β-Caryophyllene     | 1.98 %      |
| α-Bisabolol         | 0.0248 %    |
| Guaiol              | <LOQ %      |
| Humulene            | 0.164 %     |
| Caryophyllene Oxide | 0.020 %     |



1. American Herbal Pharmacopoeia. (2014). Cannabis Inflorescence: Standards of Identity, Analysis, and Quality Control. Washington DC: AHP.

2. US Food and Drug Administration. (2019). Lead in Food, Foodwares, and Dietary Supplements. Washington DC: FDA. US Food and Drug Administration. (2019). Lead in Food, Foodwares, and Dietary Supplements. Washington DC: FDA.



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



**Report Number:** 22-003508/D002.R000  
**Report Date:** 04/01/2022  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 03/28/22 14:43

**Customer:** Etz Hayim Holdings  
**Product identity:** FORM-SG50-EC04  
**Client/Metric ID:** .  
**Laboratory ID:** 22-003508-0001

### Summary

**Potency:**

| Analyte per 1g           | Result | Limits | Units | Status |                                      |
|--------------------------|--------|--------|-------|--------|--------------------------------------|
| CBC per 1g <sup>†</sup>  | 0.885  |        | mg/1g |        | CBD Total per 1g 129 mg/1g           |
| CBD per 1g               | 129    |        | mg/1g |        |                                      |
| CBDV per 1g <sup>†</sup> | 0.398  |        | mg/1g |        | THC-Total per 1g 1.01 mg/1g          |
| CBT per 1g <sup>†</sup>  | 1.14   |        | mg/1g |        |                                      |
| Δ9 THC per 1g            | 1.01   |        | mg/1g |        | (Reported in milligrams per serving) |



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**Purchase Order:**  
**Received:** 03/28/22 14:43



**Customer:** Etz Hayim Holdings  
 16427 NE Airport Way  
 PORTLAND 97230  
 United States of America (USA)

**Product identity:** FORM-SG50-EC04

**Client/Metric ID:** .

**Sample Date:**

**Laboratory ID:** 22-003508-0001

**Evidence of Cooling:** No

**Temp:** 21.7 °C

**Relinquished by:** Client

**Serving Size #1:** 1 g

### Sample Results

| Potency per 1g  |        |        |       |        |       |
|---|--------|--------|-------|--------|-------|
| Method J AOAC 2015 V98-6 (mod)Units mg/se Batch: 2202732 Analyze: 3/30/22 10:22:00 AM |        |        |       |        |       |
| Analyte   | Result | Limits | Units | LOQ    | Notes |
| CBC per 1g <sup>†</sup>   | 0.885  |        | mg/1g | 0.0329 |       |
| CBC-A per 1g <sup>†</sup>   | < LOQ  |        | mg/1g | 0.0329 |       |
| CBC-Total per 1g <sup>†</sup>   | 0.885  |        | mg/1g | 0.0618 |       |
| CBD per 1g  | 129    |        | mg/1g | 3.29   |       |
| CBD-A per 1g  | < LOQ  |        | mg/1g | 0.0329 |       |
| CBD-Total per 1g  | 129    |        | mg/1g | 3.32   |       |
| CBDV per 1g <sup>†</sup>  | 0.398  |        | mg/1g | 0.0329 |       |
| CBDV-A per 1g <sup>†</sup>  | < LOQ  |        | mg/1g | 0.0329 |       |
| CBDV-Total per 1g <sup>†</sup>  | 0.398  |        | mg/1g | 0.0615 |       |
| CBE per 1g <sup>†</sup>   | < LOQ  |        | mg/1g | 0.0329 |       |
| CBG per 1g <sup>†</sup>   | < LOQ  |        | mg/1g | 0.0329 |       |
| CBG-A per 1g <sup>†</sup>   | < LOQ  |        | mg/1g | 0.0329 |       |
| CBG-Total per 1g <sup>†</sup>   | < LOQ  |        | mg/1g | 0.0615 |       |
| CBL per 1g <sup>†</sup>   | < LOQ  |        | mg/1g | 0.0329 |       |
| CBL-A per 1g <sup>†</sup>   | < LOQ  |        | mg/1g | 0.0329 |       |
| CBL-Total per 1g <sup>†</sup>   | < LOQ  |        | mg/1g | 0.0618 |       |
| CBN per 1g  | < LOQ  |        | mg/1g | 0.0329 |       |
| CBT per 1g <sup>†</sup>   | 1.14   |        | mg/1g | 0.0329 |       |
| Δ8-THCV per 1g <sup>†</sup>   | < LOQ  |        | mg/1g | 0.0329 |       |
| Δ8-THC per 1g <sup>†</sup>  | < LOQ  |        | mg/1g | 0.0329 |       |
| Δ9-THC per 1g   | 1.01   |        | mg/1g | 0.0329 |       |
| exo-THC per 1g <sup>†</sup>   | < LOQ  |        | mg/1g | 0.0329 |       |
| THC-A per 1g  | < LOQ  |        | mg/1g | 0.0329 |       |
| THC-Total per 1g  | 1.01   |        | mg/1g | 0.0618 |       |
| THCV per 1g <sup>†</sup>  | < LOQ  |        | mg/1g | 0.0329 |       |
| THCV-A per 1g <sup>†</sup>  | < LOQ  |        | mg/1g | 0.0329 |       |
| THCV-Total per 1g <sup>†</sup>  | < LOQ  |        | mg/1g | 0.0618 |       |
| Total Cannabinoids per 1g   | 132    |        | mg/1g |        |       |



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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.

† = Analyte not NELAP accredited.

**Units of Measure**

- g = g
- mg/1g = Milligram per 1g
- % = Percentage of sample
- % wt =  $\mu\text{g/g}$  divided by 10,000

Approved Signatory

Derrick Tanner  
General Manager



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503-254-1794



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Received: 03/28/22 14:43



12423 NE Whitaker Way Portland OR 97230 p.503-254-1794

**Cannabis Chain of Custody Record**

ORELAP ID: OR100028

| Field ID       |  | Date/Time Collected |         | Analysis Requested           |   |         |                   |                |          |          |                       |                                  |              | Matrix     |       |        | Serving size for edibles | Comments/Metric ID |                              |  |
|----------------|--|---------------------|---------|------------------------------|---|---------|-------------------|----------------|----------|----------|-----------------------|----------------------------------|--------------|------------|-------|--------|--------------------------|--------------------|------------------------------|--|
|                |  |                     |         | Pesticides - OR 59 compounds | Pesticide Multi-Residue - 379 compounds | Potency | Residual Solvents | Water Activity | Moisture | Terpenes | Micro: Yeast and Mold | Micro: E.Coli and Total Coliform | Heavy Metals | Mycotoxins | Other |        |                          |                    |                              |  |
| Form-SGSD-EC04 |  | 3/28                | 11:40 A |                              |   | X       |                   |                |          |          |                       |                                  |              |            |       | Liquid | TN-                      | mg/g               | LaZ-Nat Discount potency 1st |  |
| Form-SGSD-EC04 |  | 3/28                | 11:40 A | X                            |   |         | X                 |                |          | X        | X                     | X                                | X            |            |       |        |                          |                    |                              |  |
|                |  |                     |         |                              |   |         |                   |                |          |          |                       |                                  |              |            |       |        |                          |                    |                              |  |
|                |  |                     |         |                              |   |         |                   |                |          |          |                       |                                  |              |            |       |        |                          |                    |                              |  |
|                |  |                     |         |                              |   |         |                   |                |          |          |                       |                                  |              |            |       |        |                          |                    |                              |  |

Purchase Order Number:  
Project Number:  
Project Name:  
 Report Instructions:  
 Send to State - METRC  
 Email Final Results:  
 Fax Final Results  
 Cash/Check/CC/Net 30  
Other:

| Collected By:  | Relinquished By: | Date | Time | Received by: | Date | Time | Lab Use Only:  |
|--|------------------|------|------|--------------|------|------|--|
| <input checked="" type="checkbox"/> Standard (5 day)         |                  |      |      |              |      |      | Client Alias:  |
| <input type="checkbox"/> Rush (3-4 day) (1.5x Standard)      |                  |      |      |              |      |      | Order Number:  |
| <input type="checkbox"/> Priority Rush (2 day) (2x Standard) |                  |      |      |              |      |      | Proper Container <input checked="" type="checkbox"/>                                     |
|  |                  |      |      |              |      |      | Sample Condition <input checked="" type="checkbox"/>                                     |
|  |                  |      |      |              |      |      | Temperature: 21.7 °C   |
|  |                  |      |      |              |      |      | Shipped Via: Courier   |
|  |                  |      |      |              |      |      | Evidence of cooling: <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No |

SUBMISSION OF SAMPLES WITH TESTING REQUIREMENTS TO PIXIS WILL BE UNDERSTOOD TO BE AN AGREEMENT FOR SERVICES IN ACCORDANCE WITH THE CONDITIONS LISTED ON THE BACK OF THIS FORM

Revision: 1.02 Control#: CF023  
Effective 01/31/2019 Revised 01/31/2019

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

**Report Number:** 22-003508/D002.R000  
**Report Date:** 04/01/2022  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 03/28/22 14:43



Revision 1 Document D 7148  
Legacy D Workshee Validated 04/20/2021

**Laboratory Quality Control Results**

| J AOAC 2015 V98-6         |        |       |       |       |            |            |       |  |
|---------------------------|--------|-------|-------|-------|------------|------------|-------|--|
| Batch ID: 2202732         |        |       |       |       |            |            |       |  |
| Laboratory Control Sample |        |       |       |       |            |            |       |  |
| Analyte                   | Result | Spike | Units | % Rec | Limits     | Evaluation | Notes |  |
| CBDVA                     | 0.0318 | 0.033 | %     | 95.5  | 80.0 - 120 | Acceptable |       |  |
| CBDV                      | 0.0439 | 0.033 | %     | 132   | 80.0 - 120 | Acceptable | Q6    |  |
| CBE                       | 0.0314 | 0.033 | %     | 94.1  | 80.0 - 120 | Acceptable |       |  |
| CBDA                      | 0.0345 | 0.033 | %     | 104   | 80.0 - 120 | Acceptable |       |  |
| CBGA                      | 0.0322 | 0.033 | %     | 96.5  | 80.0 - 120 | Acceptable |       |  |
| CBG                       | 0.0336 | 0.033 | %     | 101   | 80.0 - 120 | Acceptable |       |  |
| CBD                       | 0.0328 | 0.033 | %     | 98.3  | 80.0 - 120 | Acceptable |       |  |
| THCV                      | 0.0327 | 0.033 | %     | 98.0  | 80.0 - 120 | Acceptable |       |  |
| d8THCV                    | 0.0330 | 0.033 | %     | 98.9  | 80.0 - 120 | Acceptable |       |  |
| THCVA                     | 0.0310 | 0.033 | %     | 92.9  | 80.0 - 120 | Acceptable |       |  |
| CBN                       | 0.0346 | 0.033 | %     | 104   | 80.0 - 120 | Acceptable |       |  |
| exo-THC                   | 0.0308 | 0.033 | %     | 92.5  | 80.0 - 120 | Acceptable |       |  |
| d9THC                     | 0.0339 | 0.033 | %     | 102   | 80.0 - 120 | Acceptable |       |  |
| d8THC                     | 0.0339 | 0.033 | %     | 102   | 80.0 - 120 | Acceptable |       |  |
| CBL                       | 0.0319 | 0.033 | %     | 95.7  | 80.0 - 120 | Acceptable |       |  |
| CBC                       | 0.0344 | 0.033 | %     | 103   | 80.0 - 120 | Acceptable |       |  |
| THCA                      | 0.0325 | 0.033 | %     | 97.6  | 80.0 - 120 | Acceptable |       |  |
| CBCA                      | 0.0328 | 0.033 | %     | 98.5  | 80.0 - 120 | Acceptable |       |  |
| CBLA                      | 0.0308 | 0.033 | %     | 92.4  | 80.0 - 120 | Acceptable |       |  |
| CBT                       | 0.0333 | 0.033 | %     | 99.9  | 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 |       |
| CBDA    | < LOQ  | 0.003 | %     | < 0.003 | Acceptable |       |
| CBGA    | < LOQ  | 0.003 | %     | < 0.003 | Acceptable |       |
| CBG     | < LOQ  | 0.003 | %     | < 0.003 | Acceptable |       |
| CBD     | < 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 |       |
| CBL     | < LOQ  | 0.003 | %     | < 0.003 | Acceptable |       |
| CBC     | < LOQ  | 0.003 | %     | < 0.003 | Acceptable |       |
| THCA    | < LOQ  | 0.003 | %     | < 0.003 | Acceptable |       |
| CBCA    | < 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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 Portland, OR 97230  
 503-254-1794



**Report Number:** 22-003508/D002.R000  
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**Purchase Order:**  
**Received:** 03/28/22 14:43

Revision 1 Document D 7148  
 Legacy D Workshee Validated 04/20/2021

**Laboratory Quality Control Results**

| J AOAC 2015 V98-6       |         |             |       |       |      |        |            |       |
|-------------------------|---------|-------------|-------|-------|------|--------|------------|-------|
| Batch ID: 2202732       |         |             |       |       |      |        |            |       |
| Sample Duplicate        |         |             |       |       |      |        |            |       |
| Sample D 22-003445-0002 |         |             |       |       |      |        |            |       |
| 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                     | 0.00420 | 0.00437     | 0.003 | %     | 3.89 | < 20   | Acceptable |       |
| THCV                    | < LOQ   | < LOQ       | 0.003 | %     | NA   | < 20   | Acceptable |       |
| d8THCV                  | 0.108   | 0.107       | 0.003 | %     | 1.00 | < 20   | Acceptable |       |
| THCVA                   | < LOQ   | < LOQ       | 0.003 | %     | NA   | < 20   | Acceptable |       |
| CBN                     | < LOQ   | < LOQ       | 0.003 | %     | NA   | < 20   | Acceptable |       |
| exo-THC                 | < LOQ   | < LOQ       | 0.003 | %     | NA   | < 20   | Acceptable |       |
| d9THC                   | < LOQ   | < LOQ       | 0.003 | %     | NA   | < 20   | Acceptable |       |
| d8THC                   | < LOQ   | < LOQ       | 0.003 | %     | NA   | < 20   | Acceptable |       |
| CBL                     | < 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                     | 0.0512  | 0.0528      | 0.003 | %     | 3.03 | < 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.                      |





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**Report Number:** 22-003508/D004.R000  
**Report Date:** 04/07/2022  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 03/28/22 14:43

**Customer:** Etz Hayim Holdings  
**Product identity:** FORM-SG50-EC04  
**Client/Metric ID:** .  
**Laboratory ID:** 22-003508-0002

### Summary

**Residual Solvents:**

*All analytes passing and less than LOQ.*

**Pesticides:**

*All analytes passing and less than LOQ.*

**Terpenes:**

| Analyte                           | Percent by weight | Percent of Total | Analyte                              | Percent by weight | Percent of Total |
|-----------------------------------|-------------------|------------------|--------------------------------------|-------------------|------------------|
| (R)-(+)-Limonene <sup>†</sup>     | 10.9              | 60.89%           | β-Myrcene <sup>†</sup>               | 4.73              | 26.42%           |
| β-Caryophyllene <sup>†</sup>      | 1.98              | 11.06%           | Humulene <sup>†</sup>                | 0.164             | 0.92%            |
| d-3-Carene <sup>†</sup>           | 0.0345            | 0.19%            | α-Bisabolol <sup>†</sup>             | 0.0248            | 0.14%            |
| Geranyl acetate <sup>†</sup>      | 0.0225            | 0.13%            | (-)-caryophyllene oxide <sup>†</sup> | 0.0200            | 0.11%            |
| <b>Total Terpenes<sup>†</sup></b> | <b>17.9</b>       | <b>100.00%</b>   |                                      |                   |                  |

**Metals:**

*Less than LOQ for all analytes.*

**Microbiology:**

*Less than LOQ for all analytes.*



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**Purchase Order:**  
**Received:** 03/28/22 14:43

**Customer:** Etz Hayim Holdings  
16427 NE Airport Way  
PORTLAND 97230  
United States of America (USA)

**Product identity:** FORM-SG50-EC04

**Client/Metric ID:** .

**Sample Date:**

**Laboratory ID:** 22-003508-0002

**Evidence of Cooling:** No

**Temp:** 21.7 °C

**Relinquished by:** Client

### Sample Results

#### Microbiology

| Analyte                 | Result | Limits | Units | LOQ | Batch   | Analyze  | Method                  | Status | Notes |
|-------------------------|--------|--------|-------|-----|---------|----------|-------------------------|--------|-------|
| E.coli                  | < LOQ  |        | cfu/g | 10  | 2202828 | 04/04/22 | AOAC 991.14 (Petrifilm) | X, I   |       |
| Total Coliforms         | < LOQ  |        | cfu/g | 10  | 2202828 | 04/04/22 | AOAC 991.14 (Petrifilm) | X, I   |       |
| Mold (RAPID Petrifilm)  | < LOQ  |        | cfu/g | 10  | 2202829 | 04/05/22 | AOAC 2014.05 (RAPID)    | X, I   |       |
| Yeast (RAPID Petrifilm) | < LOQ  |        | cfu/g | 10  | 2202829 | 04/05/22 | AOAC 2014.05 (RAPID)    | X, I   |       |

#### Solvents

| Residual Solvents by GC/MS      |        |        |      |        |       | Units µg/g                           | Batch 2202942 | Analyze 04/06/22 08:55 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<br>(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<br>(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   |       | Ethanol <sup>l</sup>                 | < LOQ         |                           | 200  |        |       |
| Ethyl acetate                   | < LOQ  | 5000   | 200  | pass   |       | Ethyl benzene                        | < LOQ         |                           | 200  |        |       |
| Ethyl ether                     | < LOQ  | 5000   | 200  | pass   |       | Ethylene glycol                      | < LOQ         | 620                       | 200  | pass   |       |
| Ethylene oxide                  | < LOQ  | 50.0   | 20.0 | pass   |       | Hexanes (sum)                        | < LOQ         | 290                       | 150  | pass   |       |
| Isopropyl acetate               | < LOQ  | 5000   | 200  | pass   |       | Isopropylbenzene<br>(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<br>(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   |       |                                      |               |                           |      |        |       |



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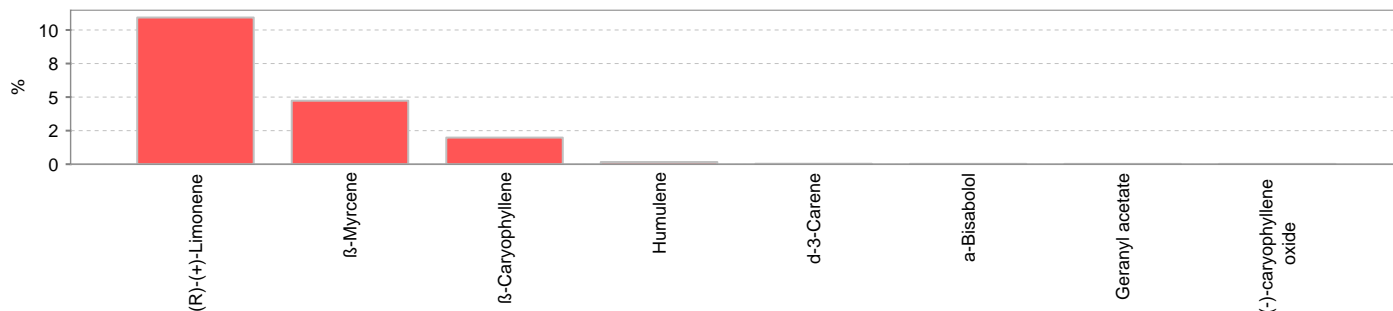


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**Report Date:** 04/07/2022  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 03/28/22 14:43

| Pesticides   |        |        |       |        |       |                     |        |        |       |        |       |
|--|--------|--------|-------|--------|-------|---------------------|--------|--------|-------|--------|-------|
| Method AOAC 2007.01 & EN 15662 (mod) Units mg/kg Batch 2202924 Analyze 04/05/22 02:43 PM |        |        |       |        |       |                     |        |        |       |        |       |
| 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   |       | Fonicamid           | < 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   |       | Paclotrazole        | < 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   |       |                     |        |        |       |        |       |



| Terpenes                         |             |       |            | Method J AOAC 2015 V98-6 | Units %                              | Batch 2202887 | Analyze 04/05/22 | 05:40 AM   |       |
|----------------------------------|-------------|-------|------------|--------------------------|--------------------------------------|---------------|------------------|------------|-------|
| Analyte                          | Result      | LOQ   | % of Total | Notes                    | Analyte                              | Result        | LOQ              | % of Total | Notes |
| (R)-(+)-Limonene <sup>†</sup>    | 10.9        | 0.194 | 60.89%     |                          | β-Myrcene <sup>†</sup>               | 4.73          | 0.019            | 26.42%     |       |
| β-Caryophyllene <sup>†</sup>     | 1.98        | 0.019 | 11.06%     |                          | Humulene <sup>†</sup>                | 0.164         | 0.019            | 0.916%     |       |
| d-3-Carene <sup>†</sup>          | 0.0345      | 0.019 | 0.1927%    |                          | a-Bisabolol <sup>†</sup>             | 0.0248        | 0.019            | 0.1385%    |       |
| Geranyl acetate <sup>†</sup>     | 0.0225      | 0.019 | 0.1257%    |                          | (-)-caryophyllene oxide <sup>†</sup> | 0.0200        | 0.019            | 0.1117%    |       |
| α-Terpinene <sup>†</sup>         | < LOQ       | 0.019 | 0.00%      |                          | p-Cymene <sup>†</sup>                | < LOQ         | 0.019            | 0.00%      |       |
| Geraniol <sup>†</sup>            | < LOQ       | 0.019 | 0.00%      |                          | gamma-Terpinene <sup>†</sup>         | < LOQ         | 0.019            | 0.00%      |       |
| nerol <sup>†</sup>               | < LOQ       | 0.019 | 0.00%      |                          | Sabinene <sup>†</sup>                | < LOQ         | 0.019            | 0.00%      |       |
| farnesene <sup>†</sup>           | < LOQ       | 0.019 | 0.00%      |                          | (+)-Pulegone <sup>†</sup>            | < LOQ         | 0.019            | 0.00%      |       |
| (-)-β-Pinene <sup>†</sup>        | < LOQ       | 0.019 | 0.00%      |                          | Menthol <sup>†</sup>                 | < LOQ         | 0.019            | 0.00%      |       |
| (±)-Camphor <sup>†</sup>         | < LOQ       | 0.019 | 0.00%      |                          | (-)-a-Terpineol <sup>†</sup>         | < LOQ         | 0.019            | 0.00%      |       |
| Sabinene hydrate <sup>†</sup>    | < LOQ       | 0.019 | 0.00%      |                          | (+)-Cedrol <sup>†</sup>              | < LOQ         | 0.019            | 0.00%      |       |
| (-)-Guaiol <sup>†</sup>          | < LOQ       | 0.019 | 0.00%      |                          | (-)-Isopulegol <sup>†</sup>          | < LOQ         | 0.019            | 0.00%      |       |
| (+)-Borneol <sup>†</sup>         | < LOQ       | 0.019 | 0.00%      |                          | (+)-fenchol <sup>†</sup>             | < LOQ         | 0.019            | 0.00%      |       |
| (±)-cis-Nerolidol <sup>†</sup>   | < LOQ       | 0.019 | 0.00%      |                          | (±)-fenchone <sup>†</sup>            | < LOQ         | 0.019            | 0.00%      |       |
| (±)-trans-Nerolidol <sup>†</sup> | < LOQ       | 0.019 | 0.00%      |                          | a-cedrene <sup>†</sup>               | < LOQ         | 0.019            | 0.00%      |       |
| a-phellandrene <sup>†</sup>      | < LOQ       | 0.019 | 0.00%      |                          | a-pinene <sup>†</sup>                | < LOQ         | 0.019            | 0.00%      |       |
| Camphene <sup>†</sup>            | < LOQ       | 0.019 | 0.00%      |                          | cis-β-Ocimene <sup>†</sup>           | < LOQ         | 0.006            | 0.00%      |       |
| Eucalyptol <sup>†</sup>          | < LOQ       | 0.019 | 0.00%      |                          | Isoborneol <sup>†</sup>              | < LOQ         | 0.019            | 0.00%      |       |
| Linalool <sup>†</sup>            | < LOQ       | 0.019 | 0.00%      |                          | Terpinolene <sup>†</sup>             | < LOQ         | 0.019            | 0.00%      |       |
| trans-β-Ocimene <sup>†</sup>     | < LOQ       | 0.012 | 0.00%      |                          | valencene <sup>†</sup>               | < LOQ         | 0.019            | 0.00%      |       |
| <b>Total Terpenes</b>            | <b>17.9</b> |       |            |                          |                                      |               |                  |            |       |



| Metals  |        |        |       |        |         |          |                     |        |       |
|---------|--------|--------|-------|--------|---------|----------|---------------------|--------|-------|
| Analyte | Result | Limits | Units | LOQ    | Batch   | Analyze  | Method              | Status | Notes |
| Arsenic | < LOQ  | 0.200  | mg/kg | 0.0840 | 2202970 | 04/06/22 | AOAC 2013.06 (mod.) | pass   | X     |
| Cadmium | < LOQ  | 0.200  | mg/kg | 0.0840 | 2202970 | 04/06/22 | AOAC 2013.06 (mod.) | pass   | X     |
| Lead    | < LOQ  | 0.500  | mg/kg | 0.0840 | 2202970 | 04/06/22 | AOAC 2013.06 (mod.) | pass   | X     |
| Mercury | < LOQ  | 0.100  | mg/kg | 0.0420 | 2202970 | 04/06/22 | AOAC 2013.06 (mod.) | pass   | X     |



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**Purchase Order:**  
**Received:** 03/28/22 14:43

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.

† = Analyte not NELAP accredited.

**Units of Measure**

cfu/g = Colony forming units per gram

µg/g = Microgram per gram

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

% = Percentage of sample

% wt = µg/g divided by 10,000

**Glossary of Qualifiers**

I: Insufficient sample received to meet method requirements.

X: Not ORELAP accredited.

Approved Signatory

Derrick Tanner  
General Manager



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**Cannabis Chain of Custody Record**

ORELAP ID: OR100028

| Field ID       |  | Date/Time Collected |  | Pesticides - OR 59 compounds | Pesticide Multi-Residue - 379 compounds | Potency | Residual Solvents | Water Activity | Moisture | Terpenes | Micro: Yeast and Mold | Micro: E.Coli and Total Coliform | Heavy Metals | Mycotoxins | Other | Matrix | Weight | Serving size for edibles | Comments/Metric ID            |
|----------------|--|---------------------|--|------------------------------|---|---------|-------------------|----------------|----------|----------|-----------------------|----------------------------------|--------------|------------|-------|--------|--------|--------------------------|-------------------------------|
| Form-SG6D-EC04 |  | 3/28 11:40 A        |  |                              |   | X       |                   |                |          |          |                       |                                  |              |            |       | Liquid | TN-    | mg/g                     | La-Z-Nat Discount potency 1st |
| Form-SG6D-EC04 |  | 3/28 11:40 A        |  | X                            |   |         | X                 |                |          | X        | X                     | X                                | X            |            |       |        |        |                          |                               |

Purchase Order Number:  
Project Number:  
Project Name:  
 Report Instructions:  
 Send to State - METRC  
 Email Final Results:  
 Fax Final Results  
 Cash/Check/CC/Net 30  
Other:

|   |                  |      |      |              |      |      |   |
|---|------------------|------|------|--------------|------|------|---|
| Collected By:   | Relinquished By: | Date | Time | Received by: | Date | Time | Lab Use Only:   |
| <input checked="" type="checkbox"/> Standard (5 day)<br><input type="checkbox"/> Rush (3-4 day) (1.5x Standard)<br><input type="checkbox"/> Priority Rush (2 day) (2x Standard) | [Redacted]       |      |      | [Redacted]   |      |      | Client Alias:<br>Order Number:<br>Proper Container <input checked="" type="checkbox"/><br>Sample Condition <input checked="" type="checkbox"/><br>Temperature: 21.7 °C<br>Shipped Via: <u>Courier</u><br>Evidence of cooling: <input type="checkbox"/> Yes <input checked="" type="checkbox"/> No |

SUBMISSION OF SAMPLES WITH TESTING REQUIREMENTS TO PIXIS WILL BE UNDERSTOOD TO BE AN AGREEMENT FOR SERVICES IN ACCORDANCE WITH THE CONDITIONS LISTED ON THE BACK OF THIS FORM  
 Revision: 1.02 Control#: CF023 [www.pixislabs.com](http://www.pixislabs.com)  
 Effective 01/31/2019 Revised 01/31/2019 Page 1 of 2



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Revision 1 Document D 7086  
 Legacy D CFL-E57Worksheet Validated 11/04/2020

**Terpenes Quality Control Results**

| Method Reference EPA 5035 |        |       |       | Batch ID 2202887          |     |       |           |         |       |
|---------------------------|--------|-------|-------|---------------------------|-----|-------|-----------|---------|-------|
| Method Blank              |        |       |       | Laboratory Control Sample |     |       |           |         |       |
| Analyte                   | Result | LOQ   | Notes | Result                    | LCS | Units | LCS % Rec | Limits  | Notes |
| a-pinene                  | < OQ   | < 200 |       | 497                       | 500 | µg/g  | 99%       | 70 - 30 |       |
| Camphene                  | < OQ   | < 200 |       | 49                        | 500 | µg/g  | 98%       | 70 - 30 |       |
| Sabinene                  | < OQ   | < 200 |       | 496                       | 500 | µg/g  | 99%       | 70 - 30 |       |
| b-Pinene                  | < OQ   | < 200 |       | 482                       | 500 | µg/g  | 96%       | 70 - 30 |       |
| b-Myrcene                 | < OQ   | < 200 |       | 492                       | 500 | µg/g  | 98%       | 70 - 30 |       |
| a-phellandrene            | < OQ   | < 200 |       | 40                        | 500 | µg/g  | 80%       | 70 - 30 |       |
| d-3-Carene                | < OQ   | < 200 |       | 547                       | 500 | µg/g  | 109%      | 70 - 30 |       |
| a-Terpinene               | < OQ   | < 200 |       | 449                       | 500 | µg/g  | 90%       | 70 - 30 |       |
| p-Cymene                  | < OQ   | < 200 |       | 487                       | 500 | µg/g  | 97%       | 70 - 30 |       |
| D-imonene                 | < OQ   | < 200 |       | 45                        | 500 | µg/g  | 90%       | 70 - 30 |       |
| α-caryophyllol            | < OQ   | < 200 |       | 466                       | 500 | µg/g  | 93%       | 70 - 30 |       |
| b-cis-Ocimene             | < OQ   | < 67  |       | 49                        | 67  | µg/g  | 90%       | 70 - 30 |       |
| b-trans-Ocimene           | < OQ   | < 33  |       | 279                       | 333 | µg/g  | 84%       | 70 - 30 |       |
| g-Terpinene               | < OQ   | < 200 |       | 4                         | 500 | µg/g  | 82%       | 70 - 30 |       |
| Sabinene hydrate          | < OQ   | < 200 |       | 454                       | 500 | µg/g  | 91%       | 70 - 30 |       |
| Terpinolene               | < OQ   | < 200 |       | 443                       | 500 | µg/g  | 89%       | 70 - 30 |       |
| D-enchone                 | < OQ   | < 200 |       | 433                       | 500 | µg/g  | 87%       | 70 - 30 |       |
| linalool                  | < OQ   | < 200 |       | 446                       | 500 | µg/g  | 89%       | 70 - 30 |       |
| α-enchol                  | < OQ   | < 200 |       | 467                       | 500 | µg/g  | 93%       | 70 - 30 |       |
| Camphor                   | < OQ   | < 200 |       | 482                       | 500 | µg/g  | 96%       | 70 - 30 |       |
| sopulego                  | < OQ   | < 200 |       | 485                       | 500 | µg/g  | 97%       | 70 - 30 |       |
| soborneol                 | < OQ   | < 200 |       | 48                        | 500 | µg/g  | 96%       | 70 - 30 |       |
| Borneol                   | < OQ   | < 200 |       | 474                       | 500 | µg/g  | 95%       | 70 - 30 |       |
| D-Menthol                 | < OQ   | < 200 |       | 489                       | 500 | µg/g  | 98%       | 70 - 30 |       |
| Terpineol                 | < OQ   | < 200 |       | 404                       | 500 | µg/g  | 81%       | 70 - 30 |       |
| Nerol                     | < OQ   | < 200 |       | 429                       | 500 | µg/g  | 86%       | 70 - 30 |       |
| Pulegone                  | < OQ   | < 200 |       | 476                       | 500 | µg/g  | 95%       | 70 - 30 |       |
| Geraniol                  | < OQ   | < 200 |       | 377                       | 500 | µg/g  | 75%       | 70 - 30 |       |
| Geranyl Acetate           | < OQ   | < 200 |       | 496                       | 500 | µg/g  | 99%       | 70 - 30 |       |
| α-Cedrene                 | < OQ   | < 200 |       | 49                        | 500 | µg/g  | 98%       | 70 - 30 |       |
| b-Caryophyllene           | < OQ   | < 200 |       | 500                       | 500 | µg/g  | 100%      | 70 - 30 |       |
| α-Humulene                | < OQ   | < 200 |       | 483                       | 500 | µg/g  | 97%       | 70 - 30 |       |
| Valenene                  | < OQ   | < 200 |       | 495                       | 500 | µg/g  | 99%       | 70 - 30 |       |
| cis-Nerolidol             | < OQ   | < 200 |       | 503                       | 500 | µg/g  | 101%      | 70 - 30 |       |
| α-arnesene                | < OQ   | < 200 |       | 564                       | 500 | µg/g  | 113%      | 70 - 30 |       |
| trans-Nerolidol           | < OQ   | < 200 |       | 55                        | 500 | µg/g  | 103%      | 70 - 30 |       |
| Caryophyllene Oxide       | < OQ   | < 200 |       | 498                       | 500 | µg/g  | 100%      | 70 - 30 |       |
| Guaiol                    | < OQ   | < 200 |       | 529                       | 500 | µg/g  | 106%      | 70 - 30 |       |
| Cedrol                    | < OQ   | < 200 |       | 53                        | 500 | µg/g  | 103%      | 70 - 30 |       |
| α-Bisabolol               | < OQ   | < 200 |       | 55                        | 500 | µg/g  | 103%      | 70 - 30 |       |

Definitions

|       |                           |
|-------|---------------------------|
| LOQ   | Limit of Quantitation     |
| LCS   | Laboratory Control Sample |
| % REC | Percent Recovery          |





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**Report Number:** 22-003508/D004.R000  
**Report Date:** 04/07/2022  
**ORELAP#:** OR100028  
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**Received:** 03/28/22 14:43

Revision 1 Document D 7086  
 Legacy D CFL-E57Worksheet Validated 11/04/2020

**Terpenes Quality Control Results**

| Method Reference EPA 5035 |        | Batch ID 2202887         |      |       |       |       |       |
|---------------------------|--------|--------------------------|------|-------|-------|-------|-------|
| Sample/Sample Duplicate   |        | Sample ID 22-003397-0002 |      |       |       |       |       |
| Analyte                   | Result | Org. Result              | LOQ  | Units | % RPD | LIMIT | Notes |
| a-pinene                  | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| Camphene                  | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| Sabinene                  | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| b-Pinene                  | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| b-Myrcene                 | 46600  | 44 00                    | 83   | µg/g  | 6%    | < 20  |       |
| a-phellandrene            | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| d-3-Carene                | 523    | 524                      | 83   | µg/g  | 0%    | < 20  |       |
| a-Terpinene               | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| p-Cymene                  | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| D-imonene                 | 0000   | 08000                    | 83   | µg/g  | 2%    | < 20  |       |
| α-calyptol                | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| b-cis-Ocimene             | < OQ   | < OQ                     | 60.9 | µg/g  | 0%    | < 20  |       |
| b-trans-Ocimene           | < OQ   | < OQ                     | 22   | µg/g  | 0%    | < 20  |       |
| g-Terpinene               | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| Sabinene hydrate          | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| Terpinolene               | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| D-enchone                 | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| α-inalool                 | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| α-enchol                  | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| Camphor                   | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| α-sopulego                | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| α-soborneol               | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| Borneol                   | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| D-Menthol                 | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| Terpineol                 | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| Nerol                     | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| Pulegone                  | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| Geraniol                  | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| Geranyl Acetate           | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| α-Cedrene                 | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| b-Caryophyllene           | 9900   | 9900                     | 83   | µg/g  | 0%    | < 20  |       |
| α-Humulene                | 420    | 4 0                      | 83   | µg/g  | 1%    | < 20  |       |
| Valenene                  | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| cis-Nerolidol             | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| α-arnesene                | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| trans-Nerolidol           | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| Caryophyllene Oxide       | 228    | 235                      | 83   | µg/g  | 3%    | < 20  |       |
| Guaiol                    | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| Cedrol                    | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |
| α-Bisabolol               | < OQ   | < OQ                     | 83   | µg/g  | 0%    | < 20  |       |

Definitions

RPD Relative Percent Difference





12423 NE Whitaker Way  
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 503-254-1794



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 Legacy D C L C2 Wo ks ee Va da ed 0/30/2020

**Laboratory Pesticide Quality Control Results**

| AOAC 2007.1 & EN 15662 |              | Units: mg/Kg              |       |            | Batch ID: 2202924 |              |        |       |
|------------------------|--------------|---------------------------|-------|------------|-------------------|--------------|--------|-------|
| Method Blank           |              | Laboratory Control Sample |       |            |                   |              |        |       |
| Analyte                | Blank Result | Blank Limits              | Notes | LCS Result | LCS Spike         | LCS % Rec    | Limits | Notes |
| Abamectin              | 0.000        | < 0.250                   |       | 1.889      | 1.000             | <b>188.9</b> | 50.0   | 150   |
| Acephate               | 0.000        | < 0.250                   |       | 1.017      | 1.000             | 101.7        | 60.0   | 120   |
| Acetaminocyl           | 0.000        | < 1.000                   |       | 4.263      | 4.000             | 106.6        | 40.0   | 160   |
| Acetamiprid            | 0.000        | < 0.100                   |       | 0.395      | 0.400             | 98.8         | 60.0   | 120   |
| Aldicarb               | 0.000        | < 0.200                   |       | 0.779      | 0.800             | 97.4         | 60.0   | 120   |
| Azoxystrobin           | 0.000        | < 0.100                   |       | 0.385      | 0.400             | 96.3         | 60.0   | 120   |
| Bifenazate             | 0.000        | < 0.100                   |       | 0.387      | 0.400             | 96.7         | 60.0   | 120   |
| Bifenthrin             | 0.000        | < 0.100                   |       | 0.414      | 0.400             | 103.5        | 50.0   | 150   |
| Boscalid               | 0.000        | < 0.200                   |       | 0.841      | 0.800             | 105.2        | 60.0   | 120   |
| Carbaryl               | 0.000        | < 0.100                   |       | 0.397      | 0.400             | 99.4         | 60.0   | 120   |
| Carbofuran             | 0.000        | < 0.100                   |       | 0.391      | 0.400             | 97.6         | 60.0   | 120   |
| Chlorantraniliprole    | 0.000        | < 0.100                   |       | 0.384      | 0.400             | 95.9         | 60.0   | 120   |
| Chlorfenapyr           | 0.000        | < 0.500                   |       | 2.217      | 2.000             | 110.8        | 60.0   | 120   |
| Chlorpyrifos           | 0.000        | < 0.100                   |       | 0.404      | 0.400             | 101.1        | 60.0   | 120   |
| Clofentazine           | 0.000        | < 0.100                   |       | 0.400      | 0.400             | 100.0        | 60.0   | 120   |
| Cyfluthrin             | 0.000        | < 0.500                   |       | 1.642      | 2.000             | 82.1         | 50.0   | 150   |
| Cypermethrin           | 0.000        | < 0.500                   |       | 1.997      | 2.000             | 99.8         | 50.0   | 150   |
| Daminozide             | 0.000        | < 0.500                   |       | 2.435      | 2.000             | <b>121.8</b> | 60.0   | 120   |
| Diazinon               | 0.000        | < 0.100                   |       | 0.404      | 0.400             | 100.9        | 60.0   | 120   |
| Dichlorvos             | 0.000        | < 0.500                   |       | 2.057      | 2.000             | 102.8        | 60.0   | 120   |
| Dimethoate             | 0.000        | < 0.100                   |       | 0.404      | 0.400             | 101.0        | 60.0   | 120   |
| Ethoprophos            | 0.000        | < 0.100                   |       | 0.403      | 0.400             | 100.8        | 60.0   | 120   |
| Etofenprox             | 0.000        | < 0.200                   |       | 0.819      | 0.800             | 102.4        | 50.0   | 150   |
| Etoxazole              | 0.000        | < 0.100                   |       | 0.412      | 0.400             | 102.9        | 60.0   | 120   |
| Fenoxycarb             | 0.000        | < 0.100                   |       | 0.393      | 0.400             | 98.2         | 60.0   | 120   |
| Fenpyroximate          | 0.000        | < 0.200                   |       | 0.796      | 0.800             | 99.6         | 60.0   | 120   |
| Fipronil               | 0.000        | < 0.200                   |       | 0.787      | 0.800             | 98.4         | 60.0   | 120   |
| Fonicamid              | 0.000        | < 0.250                   |       | 1.065      | 1.000             | 106.5        | 60.0   | 120   |
| Fludioxonil            | 0.000        | < 0.200                   |       | 0.794      | 0.800             | 99.3         | 50.0   | 150   |
| Hexythiazox            | 0.000        | < 0.250                   |       | 1.027      | 1.000             | 102.7        | 60.0   | 120   |
| Imazalil               | 0.000        | < 0.100                   |       | 0.395      | 0.400             | 98.8         | 60.0   | 120   |
| Imidacloprid           | 0.000        | < 0.200                   |       | 0.784      | 0.800             | 98.0         | 60.0   | 120   |
| Kresoxim methyl        | 0.000        | < 0.200                   |       | 0.840      | 0.800             | 105.1        | 60.0   | 120   |
| Malathion              | 0.000        | < 0.100                   |       | 0.403      | 0.400             | 100.6        | 60.0   | 120   |
| Metaxalyl              | 0.000        | < 0.100                   |       | 0.395      | 0.400             | 98.8         | 60.0   | 120   |
| Methiocarb             | 0.000        | < 0.100                   |       | 0.385      | 0.400             | 96.3         | 60.0   | 120   |
| Methomyl               | 0.000        | < 0.200                   |       | 0.728      | 0.800             | 90.9         | 60.0   | 120   |
| MGK 264                | 0.000        | < 0.100                   |       | 0.401      | 0.400             | 100.2        | 50.0   | 150   |
| Myclobutanil           | 0.000        | < 0.100                   |       | 0.394      | 0.400             | 98.4         | 60.0   | 120   |
| Naled                  | 0.000        | < 0.250                   |       | 0.976      | 1.000             | 97.6         | 50.0   | 150   |
| Oxamyl                 | 0.000        | < 0.500                   |       | 1.994      | 2.000             | 99.7         | 60.0   | 120   |
| Pacllobutrazole        | 0.000        | < 0.200                   |       | 0.782      | 0.800             | 97.8         | 60.0   | 120   |
| Parathion Methyl       | 0.000        | < 0.200                   |       | 0.893      | 0.800             | 111.6        | 50.0   | 150   |
| Permethrin             | 0.000        | < 0.100                   |       | 0.401      | 0.400             | 100.2        | 50.0   | 150   |
| Phosmet                | 0.000        | < 0.100                   |       | 0.396      | 0.400             | 99.0         | 50.0   | 150   |
| Piperonyl butoxide     | 0.000        | < 0.500                   |       | 2.376      | 2.000             | 118.8        | 60.0   | 120   |
| Prallethrin            | 0.000        | < 0.100                   |       | 0.411      | 0.400             | 102.8        | 60.0   | 120   |
| Propiconazole          | 0.000        | < 0.200                   |       | 0.819      | 0.800             | 102.3        | 60.0   | 120   |
| Propoxur               | 0.000        | < 0.100                   |       | 0.397      | 0.400             | 99.3         | 60.0   | 120   |
| Pyrethrin (Summe)      | 0.000        | < 0.100                   |       | 0.418      | 0.413             | 101.1        | 60.0   | 120   |
| Pyridaben              | 0.000        | < 0.100                   |       | 0.409      | 0.400             | 102.3        | 50.0   | 150   |
| Spirosad               | 0.000        | < 0.100                   |       | 0.409      | 0.388             | 105.4        | 50.0   | 150   |
| Spiromesifen           | 0.000        | < 0.100                   |       | 0.422      | 0.400             | 105.4        | 60.0   | 120   |
| Spirotetramat          | 0.000        | < 0.100                   |       | 0.404      | 0.400             | 101.1        | 60.0   | 120   |
| Spiroxamine            | 0.000        | < 0.200                   |       | 0.806      | 0.800             | 100.8        | 60.0   | 120   |
| ebuconazole            | 0.000        | < 0.200                   |       | 0.801      | 0.800             | 100.1        | 60.0   | 120   |
| hiacloprid             | 0.000        | < 0.100                   |       | 0.395      | 0.400             | 98.8         | 60.0   | 120   |
| hiamethoxam            | 0.000        | < 0.100                   |       | 0.424      | 0.400             | 105.9        | 60.0   | 120   |
| rifloxystrobin         | 0.000        | < 0.100                   |       | 0.402      | 0.400             | 100.4        | 60.0   | 120   |



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Laboratory Pesticide Quality Control Results

| AOAC 2007.1 & EN 15662                         |        | Units: mg/Kg              |         |       |       |       | Batch ID: 2202924 |           |        |       |  |
|--|--------|---------------------------|---------|-------|-------|-------|-------------------|-----------|--------|-------|--|
| Matrix Spike/Matrix Spike Duplicate Recoveries |        | Sample ID: 22-003658-0001 |         |       |       |       |                   |           |        |       |  |
| Analyte  | Result | MS Res                    | MSD Res | Spike | RPD%  | Limit | MS % Rec          | MSD % Rec | Limits | Notes |  |
| Abamectin                                      | 0.000  | 2.116                     | 2.090   | 1.000 | 1.2%  | < 30  | 211.6%            | 209.0%    | 50 150 | Q     |  |
| Acephate                                       | 0.000  | 1.001                     | 0.986   | 1.000 | 1.5%  | < 30  | 100.1%            | 98.6%     | 50 150 |       |  |
| Acetaminocyl                                   | 0.000  | 5.131                     | 4.659   | 4.000 | 9.6%  | < 30  | 128.3%            | 116.5%    | 50 150 |       |  |
| Acetamiprid                                    | 0.000  | 0.388                     | 0.391   | 0.400 | 0.7%  | < 30  | 97.0%             | 97.7%     | 50 150 |       |  |
| Aldicarb                                       | 0.000  | 0.764                     | 0.757   | 0.800 | 0.9%  | < 30  | 95.5%             | 94.7%     | 50 150 |       |  |
| Azoxystrobin                                   | 0.000  | 0.388                     | 0.401   | 0.400 | 3.5%  | < 30  | 96.9%             | 100.4%    | 50 150 |       |  |
| Bifenazate                                     | 0.000  | 0.422                     | 0.412   | 0.400 | 2.5%  | < 30  | 105.6%            | 103.0%    | 50 150 |       |  |
| Bifenthrin                                     | 0.000  | 0.455                     | 0.459   | 0.400 | 0.9%  | < 30  | 113.8%            | 114.8%    | 50 150 |       |  |
| Boscalid                                       | 0.000  | 0.781                     | 0.747   | 0.800 | 4.4%  | < 30  | 97.6%             | 93.4%     | 50 150 |       |  |
| Carbaryl                                       | 0.000  | 0.389                     | 0.389   | 0.400 | 0.0%  | < 30  | 97.2%             | 97.2%     | 50 150 |       |  |
| Carbofuran                                     | 0.000  | 0.382                     | 0.390   | 0.400 | 2.3%  | < 30  | 95.4%             | 97.6%     | 50 150 |       |  |
| Chlorantraniliprole                            | 0.000  | 0.369                     | 0.379   | 0.400 | 2.7%  | < 30  | 92.2%             | 94.7%     | 50 150 |       |  |
| Chlorfenapyr                                   | 0.000  | 1.985                     | 1.761   | 2.000 | 12.0% | < 30  | 99.3%             | 88.1%     | 50 150 |       |  |
| Chlorpyrifos                                   | 0.000  | 0.414                     | 0.360   | 0.400 | 13.9% | < 30  | 103.6%            | 90.1%     | 50 150 |       |  |
| Clofentazine                                   | 0.000  | 0.453                     | 0.453   | 0.400 | 0.1%  | < 30  | 113.2%            | 113.4%    | 50 150 |       |  |
| Cyfluthrin                                     | 0.000  | 2.018                     | 1.862   | 2.000 | 8.1%  | < 30  | 100.9%            | 93.1%     | 30 150 |       |  |
| Cypermethrin                                   | 0.000  | 2.057                     | 1.600   | 2.000 | 25.0% | < 30  | 102.9%            | 80.0%     | 50 150 |       |  |
| Daminozide                                     | 0.173  | 2.300                     | 2.274   | 2.000 | 1.2%  | < 30  | 106.3%            | 105.0%    | 30 150 |       |  |
| Diazinon                                       | 0.000  | 0.392                     | 0.382   | 0.400 | 2.7%  | < 30  | 98.0%             | 95.4%     | 50 150 |       |  |
| Dichlorvos                                     | 0.000  | 1.982                     | 1.908   | 2.000 | 3.8%  | < 30  | 99.1%             | 95.4%     | 50 150 |       |  |
| Dimethoate                                     | 0.000  | 0.390                     | 0.391   | 0.400 | 0.4%  | < 30  | 97.4%             | 97.8%     | 50 150 |       |  |
| Ethoprophos                                    | 0.000  | 0.384                     | 0.411   | 0.400 | 6.8%  | < 30  | 95.9%             | 102.6%    | 50 150 |       |  |
| Etofenprox                                     | 0.000  | 0.962                     | 0.921   | 0.800 | 4.3%  | < 30  | 120.2%            | 115.2%    | 50 150 |       |  |
| Etoxazole                                      | 0.000  | 0.519                     | 0.530   | 0.400 | 2.1%  | < 30  | 129.7%            | 132.4%    | 50 150 |       |  |
| Fenoxycarb                                     | 0.000  | 0.388                     | 0.384   | 0.400 | 1.0%  | < 30  | 97.0%             | 96.1%     | 50 150 |       |  |
| Fenpyroximate                                  | 0.000  | 0.861                     | 0.888   | 0.800 | 3.1%  | < 30  | 107.6%            | 111.0%    | 50 150 |       |  |
| Fipronil                                       | 0.000  | 0.872                     | 0.901   | 0.800 | 3.3%  | < 30  | 108.9%            | 112.6%    | 50 150 |       |  |
| Fonicamid                                      | 0.000  | 1.032                     | 1.012   | 1.000 | 2.0%  | < 30  | 103.2%            | 101.2%    | 50 150 |       |  |
| Fludioxonil                                    | 0.000  | 0.739                     | 0.764   | 0.800 | 3.4%  | < 30  | 92.3%             | 95.6%     | 50 150 |       |  |
| Hexythiazox                                    | 0.000  | 1.054                     | 1.097   | 1.000 | 4.0%  | < 30  | 105.4%            | 109.7%    | 50 150 |       |  |
| Imazalil                                       | 0.000  | 0.314                     | 0.310   | 0.400 | 1.3%  | < 30  | 78.4%             | 77.4%     | 50 150 |       |  |
| Imidacloprid                                   | 0.000  | 0.770                     | 0.765   | 0.800 | 0.7%  | < 30  | 96.3%             | 95.6%     | 50 150 |       |  |
| Kresoxim methyl                                | 0.000  | 0.823                     | 0.796   | 0.800 | 3.4%  | < 30  | 102.9%            | 99.5%     | 50 150 |       |  |
| Malathion                                      | 0.000  | 0.396                     | 0.389   | 0.400 | 1.9%  | < 30  | 99.0%             | 97.2%     | 50 150 |       |  |
| Metaxalyl                                      | 0.000  | 0.373                     | 0.372   | 0.400 | 0.1%  | < 30  | 93.2%             | 93.1%     | 50 150 |       |  |
| Methiocarb                                     | 0.000  | 0.381                     | 0.381   | 0.400 | 0.0%  | < 30  | 95.2%             | 95.3%     | 50 150 |       |  |
| Methomyl                                       | 0.000  | 0.650                     | 0.753   | 0.800 | 14.7% | < 30  | 81.2%             | 94.1%     | 50 150 |       |  |
| MGK 264  | 0.000  | 0.418                     | 0.412   | 0.400 | 1.4%  | < 30  | 104.5%            | 103.1%    | 50 150 |       |  |
| Myclobutanil                                   | 0.000  | 0.371                     | 0.378   | 0.400 | 1.7%  | < 30  | 92.8%             | 94.5%     | 50 150 |       |  |
| Naled  | 0.000  | 0.882                     | 0.919   | 1.000 | 4.1%  | < 30  | 88.2%             | 91.9%     | 50 150 |       |  |
| Oxamyl   | 0.000  | 1.914                     | 1.812   | 2.000 | 5.5%  | < 30  | 95.7%             | 90.6%     | 50 150 |       |  |
| Pacllobutrazole                                | 0.000  | 0.746                     | 0.736   | 0.800 | 1.4%  | < 30  | 93.3%             | 92.0%     | 50 150 |       |  |
| Parathion Methyl                               | 0.000  | 0.758                     | 0.770   | 0.800 | 1.6%  | < 30  | 94.7%             | 96.2%     | 30 150 |       |  |
| Permethrin                                     | 0.000  | 0.538                     | 0.566   | 0.400 | 5.1%  | < 30  | 134.6%            | 141.6%    | 50 150 |       |  |
| Phosmet  | 0.000  | 0.373                     | 0.378   | 0.400 | 1.3%  | < 30  | 93.2%             | 94.4%     | 50 150 |       |  |
| Piperonyl butoxide                             | 0.000  | 2.345                     | 2.435   | 2.000 | 3.8%  | < 30  | 117.2%            | 121.7%    | 50 150 |       |  |
| Prallethrin                                    | 0.000  | 0.477                     | 0.473   | 0.400 | 0.8%  | < 30  | 119.2%            | 118.2%    | 50 150 |       |  |
| Propiconazole                                  | 0.000  | 0.797                     | 0.791   | 0.800 | 0.6%  | < 30  | 99.6%             | 98.9%     | 50 150 |       |  |
| Propoxur                                       | 0.000  | 0.388                     | 0.392   | 0.400 | 1.1%  | < 30  | 96.9%             | 98.0%     | 50 150 |       |  |
| Pyrethrin (Summe)                              | 4.058  | 5.235                     | 5.381   | 0.413 | 11.7% | < 30  | 284.9%            | 320.2%    | 50 150 | Q     |  |
| Pyridaben                                      | 0.000  | 0.544                     | 0.509   | 0.400 | 6.7%  | < 30  | 135.9%            | 127.1%    | 50 150 |       |  |
| Spinosad                                       | 0.000  | 0.440                     | 0.400   | 0.388 | 9.4%  | < 30  | 113.3%            | 103.1%    | 50 150 |       |  |
| Spiromesifen                                   | 0.000  | 0.412                     | 0.414   | 0.400 | 0.4%  | < 30  | 103.0%            | 103.5%    | 50 150 |       |  |
| Spirotetramat                                  | 0.000  | 0.398                     | 0.406   | 0.400 | 2.0%  | < 30  | 99.5%             | 101.5%    | 50 150 |       |  |
| Spiroxamine                                    | 0.000  | 0.785                     | 0.761   | 0.800 | 3.1%  | < 30  | 98.1%             | 95.1%     | 50 150 |       |  |
| ebuconazole                                    | 0.000  | 0.786                     | 0.768   | 0.800 | 2.4%  | < 30  | 98.3%             | 96.0%     | 50 150 |       |  |
| hiacloprid                                     | 0.000  | 0.377                     | 0.385   | 0.400 | 2.1%  | < 30  | 94.1%             | 96.1%     | 50 150 |       |  |
| hiamethoxam                                    | 0.000  | 0.397                     | 0.419   | 0.400 | 5.3%  | < 30  | 99.3%             | 104.7%    | 50 150 |       |  |
| rifloxystrobin                                 | 0.000  | 0.419                     | 0.415   | 0.400 | 1.1%  | < 30  | 104.8%            | 103.7%    | 50 150 |       |  |



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| Laboratory Quality Control Results |        |     |       |        |                           |       |       |        |       |
|------------------------------------|--------|-----|-------|--------|---------------------------|-------|-------|--------|-------|
| Residual Solvents                  |        |     |       |        |                           |       |       |        |       |
| Batch ID: 2202942                  |        |     |       |        |                           |       |       |        |       |
| Method Blank                       |        |     |       |        | Laboratory Control Sample |       |       |        |       |
| Analyte                            | Result | LOQ | Notes | Result | Spike                     | Units | % Rec | Limits | Notes |
| Propane                            | ND     | 200 |       | 535    | 5 2                       | µg/g  | 93 5  | 60 20  |       |
| Isobutane                          | ND     | 200 |       | 9      | 3                         | µg/g  | 02 5  | 60 20  |       |
| Butane                             | ND     | 200 |       | 66     | 3                         | µg/g  | 0 8   | 60 20  |       |
| 2,2-Dimethylpropane                | ND     | 200 |       | 89     | 596                       | µg/g  | 95 8  | 60 20  |       |
| Methanol                           | ND     | 200 |       | 5 0    | 620                       | µg/g  | 95    | 60 20  |       |
| Ethylene Oxide                     | ND     | 0   |       | 59 5   | 56 2                      | µg/g  | 05 9  | 60 20  |       |
| 2-Methylbutane                     | ND     | 200 |       | 590    | 620                       | µg/g  | 98    | 60 20  |       |
| Pentane                            | ND     | 200 |       | 5 0    | 6 0                       | µg/g  | 95    | 60 20  |       |
| Ethanol                            | ND     | 200 |       | 590    | 630                       | µg/g  | 9 5   | 0 30   |       |
| Ethyl Ether                        | ND     | 200 |       | 80     | 620                       | µg/g  | 9     | 60 20  |       |
| 2,2-Dimethylbutane                 | ND     | 0   |       | 5      |                           | µg/g  | 90 2  | 60 20  |       |
| Acetone                            | ND     | 200 |       | 360    | 650                       | µg/g  | 82    | 60 20  |       |
| 2-Propanol                         | ND     | 200 |       | 0      | 6 0                       | µg/g  | 28 8  | 60 20  |       |
| Ethyl Formate                      | ND     | 500 |       | 0      | 600                       | µg/g  | 28    | 0 30   |       |
| Acetonitrile                       | ND     | 0   |       | 2      | 98                        | µg/g  | 9 8   | 60 20  |       |
| Methyl Acetate                     | ND     | 500 |       | 90     | 6 0                       | µg/g  | 92 5  | 0 30   |       |
| 2,3-Dimethylbutane                 | ND     | 0   |       | 6      | 6                         | µg/g  | 9 5   | 60 20  |       |
| Dichloromethane                    | ND     | 60  |       | 85     | 5 0                       | µg/g  | 95 3  | 60 20  |       |
| 1,2-Methylpentane                  | ND     | 0   |       | 50     | 6                         | µg/g  | 85 2  | 60 20  |       |
| m-NE                               | ND     | 500 |       | 0      | 600                       | µg/g  | 90 0  | 0 30   |       |
| 1,3-Methylpentane                  | ND     | 0   |       | 55     | 5                         | µg/g  | 88 6  | 60 20  |       |
| Hexane                             | ND     | 0   |       | 3      |                           | µg/g  | 9     | 60 20  |       |
| Propanol                           | ND     | 500 |       | 5 0    | 6 0                       | µg/g  | 93 8  | 0 30   |       |
| Methylethylketone                  | ND     | 500 |       | 5 0    | 600                       | µg/g  | 9     | 0 30   |       |
| Ethyl acetate                      | ND     | 200 |       | 500    | 630                       | µg/g  | 92 0  | 60 20  |       |
| 2-Butanol                          | ND     | 200 |       | 0      | 620                       | µg/g  | 88 3  | 60 20  |       |
| Tetrahydrofuran                    | ND     | 0   |       | 3      | 500                       | µg/g  | 8     | 60 20  |       |
| Cyclohexane                        | ND     | 200 |       | 390    | 620                       | µg/g  | 85 8  | 60 20  |       |
| 2-methyl propanol                  | ND     | 500 |       | 3 0    | 620                       | µg/g  | 80 9  | 0 30   |       |
| Benzene                            | ND     |     |       | 5      | 5 32                      | µg/g  | 85 3  | 60 20  |       |
| Isopropyl Acetate                  | ND     | 200 |       | 3 0    | 620                       | µg/g  | 8 6   | 60 20  |       |
| Heptane                            | ND     | 200 |       | 0      |                           | µg/g  | 9     | 60 20  |       |
| Butanol                            | ND     | 500 |       | 250    | 600                       | µg/g  | 8     | 0 30   |       |
| Propyl Acetate                     | ND     | 500 |       | 50     | 600                       | µg/g  | 90 6  | 0 30   |       |
| Dioxane                            | ND     | 0   |       | 23     | 50                        | µg/g  | 83 9  | 60 20  |       |
| 2-Ethoxyethanol                    | ND     | 0   |       | 56     | 8                         | µg/g  | 86 2  | 60 20  |       |
| Methylisobutylketone               | ND     | 500 |       | 360    | 6 0                       | µg/g  | 8 5   | 0 30   |       |
| 1,2-Methyl butanol                 | ND     | 500 |       | 200    | 6 0                       | µg/g  | 5     | 0 30   |       |
| Ethylene Glycol                    | ND     | 200 |       | 36     | 9                         | µg/g  | 3     | 60 20  |       |
| toluene                            | ND     | 200 |       | 0      | 9                         | µg/g  | 8 6   | 60 20  |       |
| Isobutyl Acetate                   | ND     | 500 |       | 500    | 600                       | µg/g  | 93 8  | 0 30   |       |
| Pentanol                           | ND     | 500 |       | 5 0    | 6 0                       | µg/g  | 95 0  | 0 30   |       |
| Butyl Acetate                      | ND     | 500 |       | 0      | 6 0                       | µg/g  | 9 3   | 0 30   |       |
| Ethyl benzene                      | ND     | 200 |       | 86     | 9 3                       | µg/g  | 88 5  | 60 20  |       |
| m-p Xylene                         | ND     | 200 |       | 85     | 596                       | µg/g  | 86 0  | 60 20  |       |
| o-Xylene                           | ND     | 200 |       | 890    | 9 3                       | µg/g  | 9 5   | 60 20  |       |
| Cumene                             | ND     | 0   |       | 0      |                           | µg/g  | 82 9  | 60 20  |       |
| Anisole                            | ND     | 500 |       | 350    | 6 0                       | µg/g  | 83 9  | 0 30   |       |
| DMSO                               | ND     | 500 |       | 560    | 630                       | µg/g  | 95    | 0 30   |       |
| 2-dimethoxyethane                  | ND     | 50  |       | 8      | 6                         | µg/g  | 90 2  | 0 30   |       |
| diethylamine                       | ND     | 500 |       | 360    | 600                       | µg/g  | 85 0  | 0 30   |       |
| N,N-dimethylformamide              | ND     | 50  |       | 0      | 9                         | µg/g  | 86 5  | 0 30   |       |
| N,N-dimethylacetamide              | ND     | 50  |       | 88     | 98                        | µg/g  | 98 0  | 0 30   |       |
| Pyridine                           | ND     | 50  |       | 5      | 80                        | µg/g  | 9 2   | 0 30   |       |
| 1,2-Dichloroethane                 | ND     | 0   |       | 0      |                           | µg/g  | 0     | 0 30   |       |
| Chloroform                         | ND     |     |       | 03     |                           | µg/g  | 03 0  | 0 30   |       |
| trichloroethylene                  | ND     |     |       |        |                           | µg/g  | 0     | 0 30   |       |



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| QC Sample Duplicate   |        | Sample ID: 22 003693 0004 |      |       |     |            |             |       |
|-----------------------|--------|---------------------------|------|-------|-----|------------|-------------|-------|
| 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 Methyl butane       | 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  |       |
| Aceton Nitrile        | ND     | ND                        | 60   | 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 Methyl pentane      | ND     | ND                        | 30   | µg/g  | 0.0 | 20         | Acceptable  |       |
| Mt SE                 | ND     | ND                        | 500  | µg/g  | 0.0 | 20         | Acceptable  |       |
| 3 Methyl pentane      | ND     | ND                        | 30   | µg/g  | 0.0 | 20         | Acceptable  |       |
| Hexane                | ND     | ND                        | 30   | µg/g  | 0.0 | 20         | Acceptable  |       |
| 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                        | 60   | µg/g  | 0.0 | 20         | Acceptable  |       |
| Cyclohexane           | ND     | ND                        | 200  | µg/g  | 0.0 | 20         | Acceptable  |       |
| 2 methyl propanol     | ND     | ND                        | 500  | µg/g  | 0.0 | 20         | Acceptable  |       |
| Benzene               | ND     | ND                        | 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  |       |
| Butanol               | ND     | ND                        | 500  | µg/g  | 0.0 | 20         | Acceptable  |       |
| Propyl Acetate        | ND     | ND                        | 500  | µg/g  | 0.0 | 20         | Acceptable  |       |
| Dioxane               | ND     | ND                        | 60   | 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 butanol      | ND     | ND                        | 500  | µg/g  | 0.0 | 20         | Acceptable  |       |
| Ethylene Glycol       | ND     | ND                        | 200  | µg/g  | 0.0 | 20         | Acceptable  |       |
| okane                 | ND     | ND                        | 200  | µg/g  | 0.0 | 20         | Acceptable  |       |
| Isobutyl Acetate      | ND     | ND                        | 500  | µg/g  | 0.0 | 20         | Acceptable  |       |
| Pentanol              | ND     | ND                        | 500  | µg/g  | 0.0 | 20         | Acceptable  |       |
| Butyl Acetate         | ND     | ND                        | 500  | µg/g  | 0.0 | 20         | Acceptable  |       |
| Ethyl benzene         | 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  |       |
| 2 dimethoxyethane     | ND     | ND                        | 50   | µg/g  | 0.0 | 20         | Acceptable  |       |
| riethylamine          | ND     | ND                        | 500  | µg/g  | 0.0 | 20         | Acceptable  |       |
| N N d methylformam de | ND     | ND                        | 50   | µg/g  | 0.0 | 20         | Acceptable  |       |
| N N d methylacetamide | ND     | ND                        | 50   | µg/g  | 0.0 | 20         | Acceptable  |       |
| Pyrid ne              | ND     | ND                        | 50   | µg/g  | 0.0 | 20         | Acceptable  |       |
| 2 Dichloroethane      | ND     | ND                        | µg/g | 0.0   | 20  | Acceptable |             |       |
| Chloroform            | ND     | ND                        | µg/g | 0.0   | 20  | Acceptable |             |       |
| richloroethylene      | ND     | ND                        | µg/g | 0.0   | 20  | Acceptable |             |       |

**Abbreviations**

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

**Units of Measure:**

µg/g Microgram per gram or ppm



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



**Report Number:** 22-003508/D004.R000  
**Report Date:** 04/07/2022  
**ORELAP#:** OR100028  
**Purchase Order:**  
**Received:** 03/28/22 14:43





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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.                      |