







# **SUMMARY OF ANALYSIS (SAMPLE ID: SA35466)**

Order ID: OR10498 **Testing Location:** Customer ID: 2161 Sample Type: Primary Arkansas River Valley Relief Cultivation Lot Number: Matrix: Flower 232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13133898357 **Mass:** 23g

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 06/19/2023 6034H20230504O43-3 License: ADH 113 License: 00065C **Date Received:** 06/20/2023

Cultivar (Strain) or Sample Description: Orange 43 **Date Completed:** 06/21/2023

\*This page is simply a summary of the analysis performed. For analytical details, please consult the individual Certificate(s) of Analysis for each of the specific test(s) performed. All contaminant action levels are referenced from the State of Arkansas MMJ testing guidelines

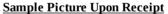
> PASS/FAIL **Moisture Content (%)** Water Activity (aw) 0.474 @ 24.0°C **PASS** 13.0

Moisture content/water activity action levels are referenced from the State of Arkansas MMJ testing guidelines.

Moisture content levels less than 15% are recommended but the sample does not fail. Water activity levels must be less than 0.65aw.

| <u>Cannabino</u>        | ids (Top 3)            | <u>(%)</u> | mg/g           |
|-------------------------|------------------------|------------|----------------|
| TH                      | CA                     | 27.6       | 276            |
| Δ9-٦                    | THC                    | 1.26       | 12.6           |
| CBG                     | GA                     | 1.19       | 11.9           |
| TOTAL                   | CBD                    | -          | -              |
| TOTAL                   | THC                    | 25.5       | 255            |
| TOTAL CAN               | NABINOIDS              | 30.4       | 304            |
| <u>Terpenes (Top 5)</u> |                        | <u>(%)</u> | <u>µg/g</u>    |
| Terpin                  | olene                  | 0.368      | 3680           |
| β-Муг                   | β-Myrcene              |            | 2440           |
| d-Lim                   | onene                  | 0.114      | 1140           |
| β-Ocimene               |                        | 0.0659     | 659            |
| β-Pir                   | nene                   | 0.0555     | 555            |
| TOTAL TI                | ERPENES                | 1.00       | 10000          |
| Contaminants            | Contaminants PASS/FAIL |            | e Upon Receipt |

| <b>Contaminants</b> | PASS/FAIL |
|---------------------|-----------|
| Heavy Metals:       | PASS      |
| Microbiology:       | PASS      |
| Pesticides:         | PASS      |







Scan the QR code to verify results.

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Page 1 of 6

06/21/2023









# **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35466)**

Testing Location:Customer ID: 2161Order ID: OR10498Sample Type: PrimaryArkansasRiver Valley Relief CultivationLot Number:Matrix: Flower232 S. Broadviow St.5601 Old Croopwood Pd Suito C.M00065C13133898357Mass: 23g

 232 S. Broadview St.
 5601 Old Greenwood Rd Suite C
 M00065C13133898357
 Mass: 23g

 Greenbrier, AR 72058
 Fort Smith, AR 72903
 Production Run:
 Date Collected: 06/19/2023

 License: ADH 113
 License: 00065C
 6034H20230504O43-3
 Date Received: 06/20/2023

Cultivar (Strain) or Sample Description: Orange 43

Date Completed: 06/21/2023

### CANNABINOID (POTENCY) PROFILE (SOP: SOP-CANN-001)

Analysis Date/Time: 06/20/2023 1709 Method: HPLC/DAD Moisture Content (%): 13.0

Analyst: PW Instrument: Agilent 1100 Water Activity (aw): 0.474 @ 24.0°C

| Cannabinoid | Result<br>(%)   | Result<br>(mg/g)  | LOD<br>(mg/g) | LOQ<br>(mg/g) | Result<br>(mg/ S<br>mL) | Per<br>Serving<br>(mg) | Per<br>Unit<br>(mg) |
|-------------|---|---|---------------|---------------|-------------------------|------------------------|---------------------|
| CBC         | ND  | ND  | 0.0438        | 0.102         | -                       | -                      | -                   |
| CBCA        | ND  | ND  | 0.136         | 0.318         | -                       | -                      | -                   |
| CBD         | ND  | ND  | 0.309         | 0.722         | -                       | -                      | -                   |
| CBDA        | ND  | ND  | 0.114         | 0.266         | -                       | -                      | -                   |
| CBDV        | ND  | ND  | 0.0496        | 0.116         | -                       | -                      | -                   |
| CBDVA       | ND  | ND  | 0.132         | 0.309         | -                       | -                      | -                   |
| CBG         | <loq< td=""><td><loq< td=""><td>0.201</td><td>0.469</td><td>-</td><td>-</td><td>-</td></loq<></td></loq<> | <loq< td=""><td>0.201</td><td>0.469</td><td>-</td><td>-</td><td>-</td></loq<> | 0.201         | 0.469         | -                       | -                      | -                   |
| CBGA        | 1.19  | 11.9  | 0.284         | 0.336         | -                       | 11.9                   | 11.9                |
| CBL         | ND  | ND  | 0.232         | 0.541         | -                       | -                      | -                   |
| CBN         | ND  | ND  | 0.106         | 0.249         | -                       | -                      | -                   |
| CBNA        | ND  | ND  | 0.115         | 0.268         | -                       | -                      | -                   |
| Δ9-ΤΗС      | 1.26  | 12.6  | 0.128         | 0.298         | -                       | 12.6                   | 12.6                |
| Δ8-ΤΗС      | ND  | ND  | 0.199         | 0.465         | -                       | -                      | -                   |
| THCA        | 27.6  | 276   | 0.0692        | 0.162         | -                       | 276                    | 276                 |
| THCV        | ND  | ND  | 0.166         | 0.387         | -                       | -                      | -                   |
| THCVA       | 0.356   | 3.56  | 0.0530        | 0.123         | -                       | 3.56                   | 3.56                |
| TOTAL       | 30.4  | 304   |               |               | -                       | 304                    | 304                 |
| TOTAL CBC   | -   | -   |               |               | -                       | -                      | -                   |
| TOTAL CBD   | -   | -   |               |               | -                       | -                      | -                   |
| TOTAL CBDV  | -   | -   |               |               | -                       | -                      | -                   |
| TOTAL CBG   | 1.04  | 10.4  |               |               | -                       | 10.4                   | 10.4                |
| TOTAL CBN   | -   | -   |               |               | -                       | -                      | -                   |
| TOTAL THC   | 25.5  | 255   |               |               | -                       | 255                    | 255                 |
| TOTAL THCV  | 0.308   | 3.08  |               |               | -                       | 3.08                   | 3.08                |

Abbreviations: DAD - Diode Array Detector, HPLC - High Pressure Liquid Chromatography, RL - Reporting Limit, RPD - Relative Percent Difference, RSD - Relative Standard Deviation, DET - Detected (less than LOQ), LOD - Limit of Detection, LOQ - Limit of Quantitation, UM - Measurement Uncertainty

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SERVING MASS (g): 1.00 SERVINGS/UNIT: 1

"-" Not detected above LOD.

Deviations from standard operating procedure: None

Recoveries for all analyte standards: 90-110% Replicate Uncertainties: <5% RSD, <20% RPD Sample/Reagent Blanks: <RL for all analytes

Values for plant matter are adjusted for moisture content.

Total CBC = (CBCA x 0.877) + CBC Total CBD = (CBDA x 0.877) + CBD Total CBDV = (CBDVA x 0.867) + CBDV Total CBG = (CBGA x 0.878) + CBG Total CBN = (CBNA x 0.876) + CBN Total THC = (THCA x 0.877) +  $\Delta$ 9-THC Total THCV = (THCVA x 0.867) + THCV

Percentage results are reported by mass.

mg/g results are reported as mass component per mass material.







Cultivar (Strain) or Sample Description: Orange 43

License: ADH 113







### **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35466)**

**Testing Location:** Customer ID: 2161 Arkansas River Valley Relief Cultivation 232 S. Broadview St. 5601 Old Greenwood Rd Suite C Greenbrier, AR 72058

Fort Smith, AR 72903

License: 00065C

Order ID: OR10498

Lot Number: M00065C13133898357

**Production Run:** 6034H20230504O43-3 Sample Type: Primary

Matrix: Flower **Mass:** 23g

**Date Collected:** 06/19/2023 **Date Received:** 06/20/2023

**Date Completed:** 06/21/2023

### TERPENOID PROFILE

Analysis Date/Time:06/21/2023 0738 Method: GC/MS **Deviations from SOP: Analyst: KF** Instrument: Agilent 7890/5975 None

| <u>Terpene</u>      | <u>Result</u><br>(µg/g) | Result<br>(%) |   |
|---------------------|-------------------------|---------------|---|
| α-Bisabolol         | 8.26                    | 0.000826      |   |
| Camphene            | 54.2                    | 0.00542       |   |
| δ-3-Carene          | 137                     | 0.0137        |   |
| β-Caryophyllene     | 267                     | 0.0267        |   |
| Caryophyllene oxide | -                       | -             |   |
| p-Cymene            | -                       | -             |   |
| Eucalyptol          | 22.4                    | 0.00224       |   |
| Geraniol            | 9.85                    | 0.000985      |   |
| Guaiol              | -                       | -             |   |
| α-Humulene          | 87.1                    | 0.00871       |   |
| Isopulegol          | -                       | -             |   |
| d-Limonene          | 1140                    | 0.114         |   |
| Linalool            | 177                     | 0.0177        |   |
| β-Myrcene           | 2440                    | 0.244         |   |
| cis-Nerolidol       | -                       | -             |   |
| trans-Nerolidol     | 35.7                    | 0.00357       |   |
| α-Ocimene           | -                       | -             |   |
| β-Ocimene           | 659                     | 0.0659        |   |
| α-Pinene            | 421                     | 0.0421        |   |
| β-Pinene            | 555                     | 0.0555        |   |
| α-Terpinene         | 188                     | 0.0188        | 1 |
| γ-Terpinene         | 125                     | 0.0125        |   |
| Terpinolene         | 3680                    | 0.368         |   |
| TOTAL               | 10000                   | 1.00          |   |



Abbreviations: GC - Gas Chromatography, MS - Mass Spectrometry, RL - Reporting Limit

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"-" Not detected above RL

Reporting Limit (µg/g): 133 10000 TOTAL 1.00







License: ADH 113





6034H20230504O43-3



# **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35466)**

Testing Location:Customer ID: 2161Order ID: OR10498Sample Type: PrimaryArkansasRiver Valley Relief CultivationLot Number:Matrix: Flower

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13133898357 **Mass:** 23g
Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 06/19/2023

Cultivar (Strain) or Sample Description: Orange 43

Date Completed: 06/21/2023

### PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 06/20/2023 1822 Method: LC/MS/MS Deviations from SOP:

Analyst: KF Instrument: Shimadzu LC-8050 None

License: 00065C

| <u>Pesticide</u>                   | Result<br>(μg/g) | <u>LOD</u><br>(µg/g) | <u>LOQ</u><br>(µg/g) | <u>Action</u><br><u>Level</u><br>(μg/g) | <u>Pesticide</u>                                 | Result<br>(μg/g) | <u>LOD</u><br>(µg/g) | <u>LOQ</u><br>(µg/g) | <u>Action</u><br><u>Level</u><br>(μg/g) |
|------------------------------------|------------------|----------------------|----------------------|---|--|------------------|----------------------|----------------------|---|
| Abamectin (71751-41-2)             | -                | 0.0479               | 0.383                | 0.5                                     | Kresoxim-methyl                                  | _                | 0.0479               | 0.383                | 0.4                                     |
| Acephate (30560-19-1)              | -                | 0.0479               | 0.383                | 0.4                                     | (143390-89-0)                                    |                  |                      |                      |   |
| Acequinocyl (57960-19-7)           | -                | 0.0479               | 0.383                | 2                                       | Malathion (121-75-5)                             | -                | 0.0479               | 0.383                | 0.2                                     |
| Acetamiprid<br>(135410-20-7)       | -                | 0.0479               | 0.383                | 0.2                                     | Metalaxyl (57837-19-1)<br>Methiocarb (2032-65-7) | -                | 0.0479<br>0.0479     | 0.383<br>0.383       | 0.2<br>0.2                              |
| Aldicarb (116-06-3)                | _                | 0.0479               | 0.383                | 0.4                                     | Methomyl (16752-77-5)                            | _                | 0.0479               | 0.383                | 0.4                                     |
| Azoxystrobin<br>(131860-33-8)      | -                | 0.0479               | 0.383                | 0.2                                     | Methyl parathion (298-0-0)                       | -                | 0.0479               | 0.383                | 0.2                                     |
| Bifenazate (149877-41-8)           | _                | 0.0479               | 0.383                | 0.2                                     | MGK 264 (113-48-4)                               | -                | 0.0479               | 0.383                | 0.2                                     |
| Bifenthrin (82657-04-3)            | _                | 0.0479               | 0.383                | 0.2                                     | Myclobutanil                                     |                  | 0.0479               | 0.383                | 0.2                                     |
| Boscalid (188425-85-6)             | -                | 0.0479               | 0.383                | 0.4                                     | (88671-89-0)                                     | _                | 0.0479               | 0.363                | 0.2                                     |
| Carbaryl (63-25-2)                 | _                | 0.0479               | 0.383                | 0.2                                     | Naled (300-76-5)                                 | -                | 0.0479               | 0.383                | 0.5                                     |
| Carbofuran (1563-66-2)             | _                | 0.0479               | 0.383                | 0.2                                     | Oxamyl (23135-22-0)                              | -                | 0.0479               | 0.383                | 1                                       |
| Chlorantraniliprole (800008-45-7)  | -                | 0.0479               | 0.383                | 0.2                                     | Paclobutrazol<br>(76738-62-0)                    | -                | 0.0479               | 0.383                | 0.4                                     |
| Chlorfenapyr                       |                  | 0.0450               | 0.000                |   | Permethrins (52645-53-1)                         | -                | 0.0479               | 0.383                | 0.2                                     |
| (122453-73-0)                      | -                | 0.0479               | 0.383                | 1                                       | Phosmet (732-11-6)                               | -                | 0.0479               | 0.383                | 0.2                                     |
| Chlorpyrifos (2921-88-2)           | -                | 0.0479               | 0.383                | 0.2                                     | Piperonyl butoxide                               | _                | 0.0479               | 0.383                | 2                                       |
| Clofentezine (74115-24-5)          | -                | 0.0479               | 0.383                | 0.2                                     | (51-03-6)  |                  |                      |                      |   |
| Cyfluthrin (68359-37-5)            | -                | 0.0479               | 0.383                | 1                                       | Prallethrins (2331-36-9)                         | -                | 0.0479               | 0.383                | 0.2                                     |
| Cypermethrin (52315-07-8)          | -                | 0.0479               | 0.383                | 1                                       | Propiconazole<br>(60207-90-1))                   | -                | 0.0479               | 0.383                | 0.4                                     |
| Daminozide (1596-84-5)             | -                | 0.0479               | 0.383                | 1                                       | Propoxur (114-26-1)                              | -                | 0.0479               | 0.383                | 0.2                                     |
| DDVP (62-73-7)                     | -                | 0.0479               | 0.383                | 0.1                                     | Pyrethrins (8003-34-7)                           | -                | 0.0479               | 0.383                | 1                                       |
| Diazinon (333-41-5)                | -                | 0.0479               | 0.383                | 0.2                                     | Pyridaben (96489-71-3)                           | -                | 0.0479               | 0.383                | 0.2                                     |
| Dimethoate (60-51-5)               | -                | 0.0479               | 0.383                | 0.2                                     | Spinosad (168316-95-8)                           | -                | 0.0479               | 0.383                | 0.2                                     |
| Ethoprophos (13194-48-4)           | -                | 0.0479               | 0.383                | 0.2                                     | Spiromesifen (283594-90-1)                       | _                | 0.0479               | 0.383                | 0.2                                     |
| Etofenprox (80844-07-1)            | -                | 0.0479               | 0.383                | 0.4                                     | Spirotetramat                                    |                  |                      |                      |   |
| Etoxazole (153233-91-1)            | -                | 0.0479               | 0.383                | 0.2                                     | (203313-25-1)                                    | -                | 0.0479               | 0.383                | 0.2                                     |
| Fenoxycarb (72490-01-8)            | -                | 0.0479               | 0.383                | 0.2                                     | Spiroxamine                                      |                  | 0.0479               | 0.383                | 0.4                                     |
| (E)-Fenpyroximate<br>(134098-61-6) | -                | 0.0479               | 0.383                | 0.4                                     | (118134-30-8)<br>Tebuconazole                    |                  | 0.0479               | 0.383                | 0.4                                     |
| Fipronil (120068-37-3)             | -                | 0.0479               | 0.383                | 0.4                                     | (80443-41-0)                                     |                  | 0.0479               | 0.363                | 0.4                                     |
| Flonicamid (158062-67-0)           | -                | 0.0479               | 0.383                | 1                                       | Thiacloprid                                      | _                | 0.0479               | 0.383                | 0.2                                     |
| Fludioxinil (131341-86-1)          | -                | 0.0479               | 0.383                | 0.4                                     | (111988-49-9)                                    |                  |                      |                      |   |
| Hexythiazox (78587-05-0)           | -                | 0.0479               | 0.383                | 1                                       | Thiamethoxam (153719-23-4)                       | -                | 0.0479               | 0.383                | 0.2                                     |
| Imazalil (35554-44-0)              | -                | 0.0479               | 0.383                | 0.2                                     | Trifloxystrobin                                  |                  | 00:-0                | 0.000                |   |
| Imidacloprid<br>(138261-41-3)      | -                | 0.0479               | 0.383                | 0.4                                     | (141517-21-7)                                    |                  | 0.0479               | 0.383                | 0.2                                     |



Color Key

**Date Received:** 06/20/2023

# RESULT < AL

"DET" detected less than LOQ

"-" not detected above LOD

Permethrins measured as the cumulative residue of the *cis*- and *trans*- permethrin isomers.

Pyrethrins measured as the cumulative residue of the pyrethrin I, cinerin I, and jasmolin I isomers.

Action levels are referenced from the

State of Arkansas MMJ testing guidelines.

A value of "-" for the action level means that analyte is not currently regulated by the regulations referenced above.

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Abbreviations: LC - Liquid Chromatography, MS - Mass Spectrometry, RL - Reporting Limit, AL - Action Level, CAS - Chemical Abstract Services, LOD - Limit of Detection, LOQ - Limit of Quantification

| <u>Pesticide</u> | Synonym(s) | <u>Pesticide</u> | Synonym(s) | <u>Pesticide</u> | Synonym(s) |
|------------------|------------|------------------|------------|------------------|------------|
| Cyfluthrin       | Baythroid  | Myclobutanil     | Systhane   | Propiconazole    | Tilt       |
| DDVP             | Dichlorvos | Naled            | Dibrom     | Propoxur         | Baygon     |
| Ethoprophos      | Prophos    | Phosmet          | Imidan     |                  |            |





**Mass:** 23g







M00065C13133898357



# **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35466)**

Testing Location:Customer ID: 2161Order ID: OR10498Sample Type: PrimaryArkansasRiver Valley Relief CultivationLot Number:Matrix: Flower

 Greenbrier, AR 72058
 Fort Smith, AR 72903
 Production Run:
 Date Collected: 06/19/2023

 License: ADH 113
 License: 00065C
 6034H20230504O43-3
 Date Received: 06/20/2023

Cultivar (Strain) or Sample Description: Orange 43

Date Completed: 06/21/2023

### **HEAVY METAL PROFILE (SOP: SOP-ICP-200.6)**

Analysis Date/Time: 06/20/2023 1838 (ICP/OES) Method: ICP/MS Deviations from SOP:

Analysis Date/Time: - (DMA) Instrument: Agilent 7500ce None

5601 Old Greenwood Rd Suite C

**Analyst:** KF

232 S. Broadview St.

| <u>Heavy Metal</u> | <u>Result</u><br>(μg/kg) | <u>LOD</u><br>(µg/kg) | LOQ<br>(µg/kg) | <u>Action Level</u><br>(μg/kg) |
|--------------------|--------------------------|-----------------------|----------------|--------------------------------|
| Arsenic (As)       | -                        | 59.1                  | 93.7           | 200                            |
| Cadmium (Cd)       | -                        | 59.1                  | 93.7           | 200                            |
| Lead (Pb)          | -                        | 59.1                  | 93.7           | 500                            |
| Mercury (Hg)       | -                        | 59.1                  | 93.7           | 100                            |



Abbreviations: ICP - Inductively-Coupled Plasma, OES - Optical Emission Spectroscopy, DMA - Direct Mercury Analyzer, RL - Reporting Limit, AL - Action Level, LOD - Limit of Detection, LOQ - Limit of Quantitation

#### Color Key

RESULT < AL
RESULT > AL

"DET" detected less than LOQ

"-" not detected above LOD

Action levels for heavy metals are referenced from the State of Arkansas MMJ testing guidelines.

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License: ADA 05 H273





6034H20230504O43-3



**Date Received:** 06/20/2023

### **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35466)**

Testing Location:Customer ID: 2161Sample ID: SA35466Sample Type: PrimaryArkansasRiver Valley Relief CultivationLot Number:Matrix: Flower

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13133898357 **Mass:** 23g
Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 06/19/2023

Cultivar (Strain) or Sample Description: Orange 43

Date Completed: 06/21/2023

### MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 06/21/2023 0953 Method: Hardy Diagnostics CompactDry Deviations from SOP:

Analyst: PW Instrument: Thermo Incubator None

| Bacteria/Microbe           | Result<br>(CFU/g) | Action Level<br>(CFU/g) |
|----------------------------|-------------------|-------------------------|
| Aerobic Plate Count        | NT                | -                       |
| Coliforms, Total           | Absent            | 1                       |
| Escherichia Coli (E. Coli) | Absent            | 100                     |
| Mold/Yeast                 | NT                | -                       |
| Pseudomonas aeruginosa     | NT                | -                       |
| Salmonella spp.            | NT                | -                       |
| Staphylococcus aureus      | NT                | -                       |

License: 00065C



Abbreviations: EC - Escherichia Coli, CFU - Colony-Forming Unit, RL - Reporting Limit, AL - Action Level, TNTC - Too Numerous To Count, NT - Not Tested Absent - Not Detected Above RL, Present - Detected Above RL

Color Key

RESULT < AL
RESULT > AL

Reporting Limit (CFU/g)

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