



SUMMARY OF ANALYSIS (SAMPLE ID: SA36325)

| Testing Location: | Customer ID: 2168 | Order ID: OR10680 | Sample Type: Primary |
|-------------------------------|-----------------------------------|--------------------|----------------------------|
| Arkansas | River Valley Relief MIPS | Lot Number: | Matrix: Edible |
| 232 S. Broadview St. | 5601 Old Greenwood Rd Suite C | M00065C13222435363 | Mass: 10pcs |
| Greenbrier, AR 72058 | Fort Smith, AR 72903 | Production Run: | Date Collected: 10/30/2023 |
| License: ADH 113 | License: 00065C | P20231023WAT14 | Date Received: 10/31/2023 |
| Cultivar (Strain) or Sample D | Date Completed: 11/01/2023 | | |

*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.

| Moisture Content (%) | Water Activity (aw) | PASS/FAIL | | |
|----------------------|---------------------|-----------|--|--|
| Not Tested | Not Tested | PASS | | |

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>Cannabinoi</u> | <u>ds (Top 3)</u> | <u>(%)</u> | <u>mg</u> /g | | |
|---------------------|-------------------|----------------|---------------------|--|--|
| Δ9-Τ | HC | 0.197 | 1.97 | | |
| Δ8-Τ | HC | 0.0113 | 0.113 | | |
| CB | D | | - | | |
| TOTAL | CBD | - | - | | |
| TOTAL | THC | 0.197 | 1.97 | | |
| TOTAL CANN | NABINOIDS | 0.208 | 2.08 | | |
| <u>Terpenes</u> | <u>(Top 5)</u> | <u>(%)</u> | µg∕g | | |
| β-Caryop | hyllene | 0.000941 | 9.41 | | |
| α-Bisa | bolol | | | | |
| Campl | nene | | | | |
| δ-3-Ca | rene | | | | |
| Caryophyllo | ene oxide | | | | |
| TOTAL TE | RPENES | 0.000941 9.41 | | | |
| Contaminants | PASS/FAIL | Sample Picture | <u>Upon Receipt</u> | | |
| Heavy Metals: | PASS | | | | |
| Microbiology: | PASS | | | | |
| Pesticides: | PASS | | | | |
| Residual Solvents: | PASS | Sec. | | | |
| | | | | | |
| | | I III | | | |



Scan the QR code to verify results.

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Kyle W. Felling, Ph.D. Laboratory Director

REPORT OF LABORATORY ANALYSIS

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| Cultivar (Strain) or Sample | Date Completed: 11/01/2023 | | |

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

Analysis Date/Time: 10/31/2023 1531 Analyst: PW Method: HPLC/DAD Instrument: Agilent 1100

| <u>Cannabinoid</u> | <u>Result</u> (%) | <u>Result</u> (mg/g) | <u>LOD</u> (mg/g) | LOQ (mg/g) | <u>Result</u> (<u>mg/</u> <u>mL</u>) | <u>Per</u> Serving (mg) | <u>Per</u> <u>Unit</u> (mg) |
|--------------------|----------------------|-------------------------|----------------------|---------------|----------------------------------------------|-------------------------------|-----------------------------------|
| CBC | ND | ND | 0.00215 | 0.00501 | - | - | - |
| CBCA | ND | ND | 0.00667 | 0.0156 | - | - | - |
| CBD | ND | ND | 0.0151 | 0.0353 | - | - | - |
| CBDA | ND | ND | 0.00557 | 0.0130 | - | - | - |
| CBDV | ND | ND | 0.00243 | 0.00567 | - | - | - |
| CBDVA | ND | ND | 0.00648 | 0.0151 | - | - | - |
| CBG | ND | ND | 0.00983 | 0.0230 | - | - | - |
| CBGA | ND | ND | 0.0139 | 0.0164 | - | - | - |
| CBL | ND | ND | 0.0114 | 0.0265 | - | - | - |
| CBN | ND | ND | 0.00522 | 0.0122 | - | - | - |
| CBNA | ND | ND | 0.00563 | 0.0131 | - | - | - |
| Δ9-ΤΗC | 0.197 | 1.97 | 0.00625 | 0.0146 | - | 9.62 | 96.2 |
| $\Delta 8$ -THC | 0.0113 | 0.113 | 0.00975 | 0.0228 | - | 0.549 | 5.49 |
| THCA | ND | ND | 0.00339 | 0.00793 | - | - | - |
| THCV | ND | ND | 0.00813 | 0.0190 | - | - | - |
| THCVA | ND | ND | 0.00260 | 0.00604 | - | - | - |
| TOTAL | 0.208 | 2.08 | | , | | 10.2 | 102 |
| TOTAL CBC | - | - | | | | - | - |
| TOTAL CBD | - | - | | | - | - | - |
| TOTAL CBDV | - | - | | | - | - | - |
| TOTAL CBG | - | - | | | - | - | - |
| TOTAL CBN | - | - | | | - | - | - |
| TOTAL THC | 0.197 | 1.97 | | | - | 9.62 | 96.2 |
| TOTAL THCV | - | - | | | - | - | - |

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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Moisture Content (%): -Water Activity (aw): -



SERVING MASS (g): 4.88 SERVINGS/UNIT: 10

"-" 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) + Δ 9-THC Total THCV = (THCVA x 0.867) + THCV

Percentage results are reported by mass.

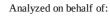
mg/g results are reported as mass component per mass material.

Kyle W. Felling, Ph.D. atory Directo











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|-----------------------------|-----------------------------------|------------------------|----------------------------|
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| License: ADH 113 | License: 00065C | P20231023WAT14 | Date Received: 10/31/2023 |
| Cultivar (Strain) or Sample | Date Completed: 11/01/2023 | | |

| Analysis Date/Time:10/31/2023 2250 | | Met | hod: GC/MS | Deviations from SOP: | | | |
|------------------------------------|-------------------------------------------------------------------------------------------------------------------|-----------------------------|-----------------------------------|-------------------------------------------------------------------------------------|--|--|--|
| Analyst: KF | | Inst | r ument: Agilent 7890/5975 | None | | | |
| <u>Terpene</u> | <u>Result</u> (µg/g) | <u>Result</u> <u>(%)</u> | | | | | |
| α-Bisabolol | ND | - | | | | | |
| Camphene | ND | - | | | | | |
| δ-3-Carene | ND | - | | | | | |
| β-Caryophyllene | 9.41 | 0.000941 | | MA I I | | | |
| Caryophyllene oxide | ND | - | | SOUR GUMMIES WITERALINA- | | | |
| p-Cymene | ND | - | | So will be a | | | |
| Eucalyptol | ND | - | | · AN HETMELER LEVE | | | |
| Geraniol | ND | - | | | | | |
| Guaiol | ND | - | | Abhumistismu CC Car | | | |
| α-Humulene | ND | - | | Abbreviations: GC - Gas Chromatography, MS - Mass | | | |
| Isopulegol | ND | - | | Spectrometry, RL - Reporting Limit | | | |
| d-Limonene | ND | - | | <i>Abbreviations:</i> ND - Not Detected, , LOD - Limit of Detection, LOQ - Limit | | | |
| Linalool | ND | - | | of Quantitation | | | |
| β-Myrcene | ND | - | | This information is provided as a service and makes no claims of efficacy and/or | | | |
| cis-Nerolidol | ND | - | | safety of this product. | | | |
| trans-Nerolidol | ND | - | | Results are applicable only for the sample(s) analyzed and for the specific | | | |
| α-Ocimene | ND | - | | analysis conducted. | | | |
| β-Ocimene | ND | - | | This report is for informational purposes | | | |
| α-Pinene | <loq< td=""><td>-</td><td></td><td>only and should not be used to diagnose, treat, or prevent any</td></loq<> | - | | only and should not be used to diagnose, treat, or prevent any | | | |
| β-Pinene | ND | - | | medical-related symptoms. | | | |
| α-Terpinene | ND | - | | The statements and results herein have not been approved and/or endorsed by | | | |
| γ-Terpinene | ND | - | | the FDA. | | | |
| Terpinolene | ND | - | | | | | |
| TOTAL | 9.41 | 0.000941 | | Reporting Limit (µg/g): 7 | | | |

Reporting Limit (µg/g): 7.85

"-" Not detected above LOD.













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RESIDUAL SOLVENT PROFILE (SOP: SOP-RS-001)

| Analysis Date/Time: Analyst: KF | 10/31/20 |)23 1933 | 1933 Method: HS/GC/MS Instrument: Agilent 7890/5975 | | | Deviations from SOP: 75 None | | | | |
|-------------------------------------|-------------------------|----------------------|--------------------------------------------------------|----------------------------------|--------------------------------------|----------------------------------------|----------------------|----------------------|----------------------------------|---------------------------------------------------|
| Solvent | <u>Result</u> (µg/g) | <u>LOD</u> (µg/g) | <u>LOQ</u> (µg/g) | <u>Action</u> Level (µg/g) | Solvent | <u>Result</u> (µg/g) | <u>LOD</u> (µg/g) | <u>LOQ</u> (µg/g) | <u>Action</u> Level (µg/g) | |
| Acetone (67-64-1) | - | 126 | 251 | 5000 | n-Heptane (142-82-5) | _ | 126 | 251 | 5000 | |
| Acetonitrile (75-5-8) | - | 126 | 251 | 410 | n-Hexane (110-54-3) | - | 44.0 | 87.9 | 290 | |
| Benzene (71-43-2) | - | 1.26 | 2.51 | 2 | Isobutane (75-28-5) | - | 126 | 251 | 5000 | MA IN |
| n-Butane (106-97-2) | - | 126 | 251 | 5000 | Isopropanol (67-63-0) | _ | 126 | 251 | 5000 | |
| 1-Butanol (71-36-3) | - | 126 | 251 | 5000 | Isopropyl acetate | | 100 | 251 | 5000 | SOURIE |
| 2-Butanol (78-92-2) | - | 126 | 251 | 5000 | (108-21-4) | - | 126 | 251 | 5000 | No antimical |
| 2-Butanone (78-93-3) | - | 126 | 251 | 5000 | Isopropyl benzene | _ | 12.6 | 25.1 | 70 | Pun - |
| Cyclohexane (110-82-7) | - | 126 | 251 | 3880 | (98-82-8) | | | | | Color Key |
| 1,2-Dimethoxyethane | | 10.0 | 05.4 | 100 | Methanol (67-56-1) | 288 | 126 | 251 | 3000 | <u>coor Rey</u> |
| (110-71-4) | - | 12.6 | 25.1 | 100 | 2-Methylbutane (78-78-4) | - | 126 | 251 | 5000 | RESULT < AL |
| N,N-Dimethylacetamide (127-19-5) | - | 126 | 251 | 1090 | Methylene chloride (75-9-2) | - | 126 | 251 | 600 | RESULT > AL |
| 2,2-Dimethylbutane (75-83-2) | - | 44.0 | 87.9 | 290 | 2-Methylpentane (107-83-5) | - | 44.0 | 87.9 | 290 | "DET" detected less than LOQ |
| 2,3-Dimethylbutane | | 44.0 | 87.9 | 290 | 3-Methylpentane (96-10-0) | - | 44.0 | 87.9 | 290 | "-" not detected above |
| (79-29-8) | _ | 44.0 | 07.9 | 290 | n-Pentane (109-66-0) | - | 126 | 251 | 5000 | LOD |
| N,N-Dimethylformamide | _ | 126 | 251 | 880 | 1-Pentanol (71-41-0) | - | 126 | 251 | 5000 | "*" - o,m,p-Xylene and |
| (68-12-2) | | | | | n-Propane (74-98-6) | - | 126 | 251 | 5000 | Ethylbenzene |
| Dimethylsulfoxide (67-68-5) | - | 126 | 251 | 5000 | 1-Propanol (71-23-8) | - | 126 | 251 | 5000 | |
| 1,4-Dioxane (123-91-1) | _ | 126 | 251 | 380 | Pyridine (110-86-1) | - | 44.0 | 87.9 | 200 | Action levels are referenced from the State of |
| Ethanol (64-17-5) | _ | 126 | 251 | 5000 | Tetrahydrofuran (109-99-9) | - | 126 | 251 | 720 | Arkansas |
| 2-Ethoxyethanol (110-80-5) | _ | 44.0 | 87.9 | 160 | Tetramethylene sulfone | _ | 44.0 | 87.9 | 160 | MMJ testing |
| Ethyl ether (60-29-7) | _ | 126 | 251 | 5000 | (126-33-0) | | | | | guidelines. |
| Ethyl acetate (141-78-6) | _ | 126 | 251 | 5000 | Toluene (108-88-3) | - | 126 | 251 | 890 | A value of "-" |
| Ethyl benzene (100-41-4) | _ | 126 | 251 | 2170 | o-Xylene (95-47-6) | - | 126 | 251 | 2170 | for the action level |
| Ethylene glycol (107-21-1) | _ | 126 | 251 | 620 | m,p-Xylene (108-38-3 or 106-42-3) | - | 126 | 251 | 2170 | means that analyte |
| Ethylene oxide (75-21-8) | | 12.6 | 25.1 | 50 | Xylenes* (1330-20-7) | | 43.3 | 86.7 | 2170 | is not currently |
| , | | | | 50 | | | - | 00.7 | 2170 | regulated by the regulations referenced above. |
| <u>Solvent</u> | | Synonym(s | | | <u>Solvent</u> | | ynonym(s) | | | |
| Acetonitrile 1-Butanol | | Methyl Cya | | L - 1 | Ethylene glycol | | ,2-Ethanediol | | | |
| | | n-Butanol, | 5 | 001 | Isobutane | | -Methylpropa | | | |
| 2-Butanol | | sec-Butyl a | | (E) I | Isopropanol | | -Propanol, IP/ | | | |
| 2-Butanone | | Methyl eth | · · | 1EK | Isopropyl Acetate | | cetic acid iso | | r | |
| 1,2-Dimethoxyethane | | Monoglym | | | Methanol | | fethyl alcoho | l | | |
| 2,3-Dimethylbutane | | Neohexane | | | 2-Methylbutane | | sopentane | | | |
| 2,3-Dimethylbutane | | Diisopropy | 71 | | Methylene chloride | | Dichlorometha | ine | | |
| N,N-Dimethylformamide | | DMF | | | 2-Methylpentane | | sohexane | | | |
| Dimethysufoxide | | DMSO | | | 1-Pentanol | | -Amyl alcoho | | | |
| 2-Ethoxyethanol | | Cellosolve | | 101 | 1-Propanol | | ropyl alcohol | | | |
| Ethyl ether | | Diethyl eth | er, Ether | | Tetrahydrofuran | | ΉF | | | |
| Ethyl acetate | | EtOAc | | | Tetramethylene sulfone | | ulfolane | | | |
| Ethyl benzene | | Phenyletha | ne | | Xylene | D | Dimethylbenze | ene | | |

Abbreviations: HS-Headspace, GC-Gas Chromatography, MS-Mass Spectrometry, RL-Reporting Limit, AL-Action Level, CAS-Chemical Abstract Services, LOD - Limit of Detection, LOQ - Limit of Quantification

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Deviations from SOP:

None



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36325)

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Analysis Date/Time: 10/31/2023 1609 Analyst: KF

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Method: LC/MS/MS Instrument: Shimadzu LC-8050

| | <u>Result</u> (µg/g) | <u>LOD</u> (µg/g) | <u>LOQ</u> (µg/g) | <u>Action</u> Level (µg/g) | <u>Pesticide</u> | <u>Result</u> (µg/g) | <u>LOD</u> (µg/g) | <u>LOQ</u> (µg/g) | <u>Action</u> Level (µg/g) | |
|--------------------------------------|-------------------------|----------------------|----------------------|----------------------------------|--------------------------------------------------|-------------------------|----------------------|----------------------|----------------------------------|----------------------------------------------------------------------------------------|
| Abamectin (71751-41-2) | - | 0.00946 | 0.0757 | 0.5 | Kresoxim-methyl | _ | 0.00946 | 0.0757 | 0.4 | |
| Acephate (30560-19-1) | - | 0.00946 | 0.0757 | 0.4 | (143390-89-0) | | | | | |
| Acequinocyl (57960-19-7) | - | 0.00946 | 0.0757 | 2 | Malathion (121-75-5) | - | 0.00946 | 0.0757 | 0.2 | |
| Acetamiprid (135410-20-7) | - | 0.00946 | 0.0757 | 0.2 | Metalaxyl (57837-19-1) Methiocarb (2032-65-7) | - | 0.00946 0.00946 | 0.0757 0.0757 | 0.2 0.2 | SOUR CUMMES SOUR CUMMES SOUR CUMMES SOUR CUMMES |
| Aldicarb (116-06-3) | - | 0.00946 | 0.0757 | 0.4 | Methomyl (16752-77-5) | - | 0.00946 | 0.0757 | 0.4 | |
| Azoxystrobin (131860-33-8) | - | 0.00946 | 0.0757 | 0.2 | Methyl parathion (298-0-0) | - | 0.00946 | 0.0757 | 0.2 | Air anni |
| Bifenazate (149877-41-8) | - | 0.00946 | 0.0757 | 0.2 | MGK 264 (113-48-4) | - | 0.00946 | 0.0757 | 0.2 | <u>Color Key</u> |
| Bifenthrin (82657-04-3) | - | 0.00946 | 0.0757 | 0.2 | Myclobutanil | _ | 0.00946 | 0.0757 | 0.2 | RESULT < AL |
| Boscalid (188425-85-6) | - | 0.00946 | 0.0757 | 0.4 | (88671-89-0) | | | | | RESULT > AL |
| Carbaryl (63-25-2) | - | 0.00946 | 0.0757 | 0.2 | Naled (300-76-5) | - | 0.00946 | 0.0757 | 0.5 | |
| Carbofuran (1563-66-2) | - | 0.00946 | 0.0757 | 0.2 | Oxamyl (23135-22-0) | - | 0.00946 | 0.0757 | 1 | "DET" detected less than LOQ |
| Chlorantraniliprole (800008-45-7) | - | 0.00946 | 0.0757 | 0.2 | Paclobutrazol (76738-62-0) | - | 0.00946 | 0.0757 | 0.4 | "-" not detected above LOD |
| Chlorfenapyr (122453-73-0) | - | 0.00946 | 0.0757 | 1 | Permethrins (52645-53-1) | - | 0.00946 | 0.0757 | 0.2 | Permethrins measured as the |
| Chlorpyrifos (2921-88-2) | - | 0.00946 | 0.0757 | 0.2 | Phosmet (732-11-6) | - | 0.00946 | 0.0757 | 0.2 | cumulative residue of the <i>cis</i> - and <i>trans</i> - permethrin isomers. |
| Clofentezine (74115-24-5) | - | 0.00946 | 0.0757 | 0.2 | Piperonyl butoxide (51-03-6) | - | 0.00946 | 0.0757 | 2 | L |
| Cyfluthrin (68359-37-5) | - | 0.00946 | 0.0757 | 1 | (31-03-0) Prallethrins (2331-36-9) | _ | 0.00946 | 0.0757 | 0.2 | Pyrethrins measured as the cumulative residue of the |
| Cypermethrin (52315-07-8) | - | 0.00946 | 0.0757 | 1 | Propiconazole (60207-90-1)) | _ | 0.00946 | 0.0757 | 0.4 | pyrethrin I, cinerin I, and jasmolin I isomers. |
| Daminozide (1596-84-5) | - | 0.00946 | 0.0757 | 1 | Propoxur (114-26-1) | _ | 0.00946 | 0.0757 | 0.2 | Action levels are referenced from |
| DDVP (62-73-7) | - | 0.00946 | 0.0757 | 0.1 | Pyrethrins (8003-34-7) | _ | 0.00946 | 0.0757 | 1 | the |
| Diazinon (333-41-5) | - | 0.00946 | 0.0757 | 0.2 | Pyridaben (96489-71-3) | _ | 0.00946 | 0.0757 | 0.2 | State of Arkansas MMJ testing guidelines. |
| Dimethoate (60-51-5) | - | 0.00946 | 0.0757 | 0.2 | Spinosad (168316-95-8) | _ | 0.00946 | 0.0757 | 0.2 | guidennes. |
| Ethoprophos (13194-48-4) | - | 0.00946 | 0.0757 | 0.2 | Spiromesifen (283594-90-1) | - | 0.00946 | 0.0757 | 0.2 | A value of "-" for the action level means that analyte is not |
| Etofenprox (80844-07-1) | - | 0.00946 | 0.0757 | 0.4 | (283394-90-1) Spirotetramat | | | | | currently regulated by the regulations referenced above. |
| Etoxazole (153233-91-1) | - | 0.00946 | 0.0757 | 0.2 | (203313-25-1) | - | 0.00946 | 0.0757 | 0.2 | legulations referenced above. |
| Fenoxycarb (72490-01-8) | - | 0.00946 | 0.0757 | 0.2 | Spiroxamine | | 0.00946 | 0.0757 | 0.4 | <i>Disclaimer:</i> This information is provided as a service and makes |
| (E)-Fenpyroximate (134098-61-6) | - | 0.00946 | 0.0757 | 0.4 | (118134-30-8) Tebuconazole | - | | | | no claims of efficacy and/or safety of this product. Results are |
| Fipronil (120068-37-3) | - | 0.00946 | 0.0757 | 0.4 | (80443-41-0) | - | 0.00946 | 0.0757 | 0.4 | applicable only for the sample(s) |
| Flonicamid (158062-67-0) | - | 0.00946 | 0.0757 | 1 | Thiacloprid | _ | 0.00946 | 0.0757 | 0.2 | analyzed and for the specific analysis conducted. This report is |
| Fludioxinil (131341-86-1) | - | 0.00946 | 0.0757 | 0.4 | (111988-49-9) | | | | | for informational purposes only |
| Hexythiazox (78587-05-0) | - | 0.00946 | 0.0757 | 1 | Thiamethoxam (153719-23-4) | - | 0.00946 | 0.0757 | 0.2 | and should not be used to |
| Imazalil (35554-44-0) | - | 0.00946 | 0.0757 | 0.2 | Trifloxystrobin | | 0.000.45 | 0.0775 | 0.0 | diagnose, treat, or prevent any medical-related symptoms. The |
| Imidacloprid (138261-41-3) | - | 0.00946 | 0.0757 | 0.4 | (141517-21-7) | | 0.00946 | 0.0757 | 0.2 | statements and results herein have not been approved and/or endorsed by the FDA. |
| | | | | | | | | | | |

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

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



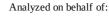
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| Testing Location: | Customer ID: 2168 | Order ID: OR10680 | Sample Type: Primary |
|-------------------------------------------|-----------------------------------|--------------------|----------------------------|
| Arkansas | River Valley Relief MIPS | Lot Number: | Matrix: Edible |
| 232 S. Broadview St. | 5601 Old Greenwood Rd Suite C | M00065C13222435363 | Mass: 10pcs |
| Greenbrier, AR 72058 Fort Smith, AR 72903 | | Production Run: | Date Collected: 10/30/2023 |
| License: ADH 113 | Date Received: 10/31/2023 | | |
| Cultivar (Strain) or Sample | Date Completed: 11/01/2023 | | |

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

| Analysis Date/T Analysis Date/T Analyst: KF | ime: 10/31/2023 1 ime: - (DMA) | 709 (ICP/OES) | Method: ICP/MS Instrument: Agilent 7500ce | | Deviations from SOP: None |
|---------------------------------------------------|-------------------------------------------------|-----------------------|------------------------------------------------------------|--------------------------------|-------------------------------------------------------------------------------------------------------------|
| <u>Heavy Metal</u> | <u>Result</u> (µg/kg) | <u>LOD</u> (µg/kg) | <u>LOQ</u> (µg/kg) | <u>Action Level</u> (µg/kg) | |
| Arsenic (As) | - | 57.5 | 91.1 | 200 | |
| Cadmium (Cd) | - | 57.5 | 91.1 | 200 | |
| Lead (Pb) | - | 57.5 | 91.1 | 500 | |
| Mercury (Hg) | - | 57.5 | 91.1 | 100 | CURCHANES BURGUMANES BURGUMANES BURGUMANES BURGUMANES BURGUMANES BURGUMANES BURGUMANES |

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

<u>Color Key</u>



"DET" detected less than LOQ

"-" not detected above LOD

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

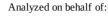
Disclaimer: This information is provided as a service and makes no claims of efficacy and/or safety of this product. Results are applicable only for the sample(s) analyzed and for the specific analysis conducted. This report is for informational purposes only and should not be used to diagnose, treat, or prevent any medical-related symptoms. The statements and results herein have not been approved and/or endorsed by the FDA.













| Testing Location: | Customer ID: 2168 | Sample ID: SA36325 | Sample Type: Primary |
|-----------------------------------------------------|-----------------------------------|--------------------|----------------------------|
| Arkansas | River Valley Relief MIPS | Lot Number: | Matrix: Edible |
| 232 S. Broadview St. | 5601 Old Greenwood Rd Suite C | M00065C13222435363 | Mass: 10pcs |
| Greenbrier, AR 72058 | Fort Smith, AR 72903 | Production Run: | Date Collected: 10/30/2023 |
| License: ADA 05_H273 License: 00065C P20231023WAT14 | | | Date Received: 10/31/2023 |
| Cultivar (Strain) or Sample | Date Completed: 11/01/2023 | | |

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

| Analysis Date/Time: 11/01/202 Analyst: PW | | d: Hardy Diagnostics CompactDry ment: Thermo Incubator | Deviations from SOP: None |
|----------------------------------------------|--------------------------|-----------------------------------------------------------|----------------------------------------|
| Bacteria/Microbe | <u>Result</u> (CFU/g) | <u>Action Level</u> (CFU/g) | |
| Aerobic Plate Count | NT | - | |
| Coliforms, Total | Absent | 1 | |
| Escherichia Coli (E. Coli) | Absent | 100 | |
| Mold/Yeast | NT | - | MA I E |
| Pseudomonas aeruginosa | NT | - | |
| Salmonella spp. | NT | - | Ual miles Sourcements Sourcement |
| Staphylococcus aureus | NT | - | Avi arminated |
| | | | |

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





1

Action levels for microbiology 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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