

SUMMARY OF ANALYSIS (SAMPLE ID: SA35855)

Testing Location: Arkansas 232 S. Broadview St. Greenbrier, AR 72058 License: ADH 113	Customer ID: 2168 River Valley Relief MIPS 5601 Old Greenwood Rd Suite C Fort Smith, AR 72903 License: 00065C	Order ID: OR10579 Lot Number: M00065C13226286865 Production Run: 20230817STRAW13	Sample Type: Primary Matrix: Edible Mass: 1bag Date Collected: 08/21/2023 Date Received: 08/21/2023
Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries			Date Completed: 08/25/2023

*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 (%)

Not Tested

Water Activity (aw)

Not Tested

PASS/FAIL

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.

Cannabinoids (Top 3)	(%)	mg/g
Δ9-THC	1.84	18.4
CBG	0.0768	0.768
CBN	0.0285	0.285
TOTAL CBD	0.00880	0.0880
TOTAL THC	1.84	18.4
TOTAL CANNABINOIDS	1.97	19.7

Terpenes (Top 5)	(%)	µg/g
α-Bisabolol		
Camphene		
δ-3-Carene		
β-Caryophyllene		
Caryophyllene oxide		
TOTAL TERPENES	-	-

Contaminants	PASS/FAIL
Heavy Metals:	PASS
Microbiology:	PASS
Pesticides:	PASS
Residual Solvents:	PASS



Scan the QR code to verify results.

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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35855)

Testing Location: Arkansas 232 S. Broadview St. Greenbrier, AR 72058 License: ADH 113	Customer ID: 2168 River Valley Relief MIPS 5601 Old Greenwood Rd Suite C Fort Smith, AR 72903 License: 00065C	Order ID: OR10579 Lot Number: M00065C13226286865 Production Run: 20230817STRAW13	Sample Type: Primary Matrix: Edible Mass: 1bag Date Collected: 08/21/2023 Date Received: 08/21/2023
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Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries **Date Completed:** 08/25/2023

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

Analysis Date/Time: 08/22/2023 1655
Analyst: PW

Method: HPLC/DAD
Instrument: Agilent 1100

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

<u>Cannabinoid</u>	<u>Result (%)</u>	<u>Result (mg/g)</u>	<u>LOD (mg/g)</u>	<u>LOQ (mg/g)</u>	<u>Result (mg/mL)</u>	<u>Per Serving (mg)</u>	<u>Per Unit (mg)</u>
CBC	ND	ND	0.00430	0.0100	-	-	-
CBCA	ND	ND	0.0134	0.0312	-	-	-
CBD	0.00880	0.0880	0.0303	0.0708	-	0.0466	1.40
CBDa	ND	ND	0.0112	0.0261	-	-	-
CBDV	ND	ND	0.00487	0.0114	-	-	-
CBDVA	ND	ND	0.0130	0.0303	-	-	-
CBG	0.0768	0.768	0.0197	0.0460	-	0.406	12.2
CBGA	ND	ND	0.0279	0.0330	-	-	-
CBL	ND	ND	0.0228	0.0531	-	-	-
CBN	0.0285	0.285	0.0105	0.0244	-	0.151	4.52
CBNA	ND	ND	0.0113	0.0263	-	-	-
Δ9-THC	1.84	18.4	0.0125	0.0292	-	9.76	293
Δ8-THC	ND	ND	0.0196	0.0456	-	-	-
THCA	ND	ND	0.00679	0.0159	-	-	-
THCV	0.0129	0.129	0.0163	0.0380	-	0.0680	2.04
THCVA	ND	ND	0.00521	0.0121	-	-	-
TOTAL	1.97	19.7				10.4	313
TOTAL CBC	-	-				-	-
TOTAL CBD	0.00880	0.0880				0.0466	1.40
TOTAL CBDV	-	-				-	-
TOTAL CBG	0.0768	0.768				0.406	12.2
TOTAL CBN	0.0285	0.285				0.151	4.52
TOTAL THC	1.84	18.4				9.76	293
TOTAL THCv	0.0129	0.129				0.0680	2.04



SERVING MASS (g): 0.529
SERVINGS/UNIT: 30

"-" 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) + Δ9-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.

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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Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries **Date Completed:** 08/25/2023

TERPENOID PROFILE

Analysis Date/Time: 08/22/2023 1834
Analyst: KF

Method: GC/MS
Instrument: Agilent 7890/5975

Deviations from SOP:
None

<u>Terpene</u>	<u>Result (µg/g)</u>	<u>Result (%)</u>
α-Bisabolol	ND	-
Camphene	ND	-
δ-3-Carene	ND	-
β-Caryophyllene	ND	-
Caryophyllene oxide	ND	-
p-Cymene	ND	-
Eucalyptol	ND	-
Geraniol	ND	-
Guaiol	ND	-
α-Humulene	ND	-
Isopulegol	ND	-
d-Limonene	ND	-
Linalool	ND	-
β-Myrcene	ND	-
cis-Nerolidol	ND	-
trans-Nerolidol	ND	-
α-Ocimene	ND	-
β-Ocimene	ND	-
α-Pinene	ND	-
β-Pinene	ND	-
α-Terpinene	ND	-
γ-Terpinene	ND	-
Terpinolene	ND	-
TOTAL	0.000	0.000



Abbreviations: GC - Gas Chromatography, MS - Mass Spectrometry, RL - Reporting Limit
Abbreviations: ND - Not Detected, LOD - Limit of Detection, LOQ - Limit of Quantitation

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Reporting Limit (µg/g): 9.44

"-" Not detected above LOD.

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Kyle W. Felling
Kyle W. Felling, Ph.D.
Laboratory Director



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Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries			Date Completed: 08/25/2023

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

Analysis Date/Time: 08/23/2023 0250	Method: HS/GC/MS	Deviations from SOP:
Analyst: KF	Instrument: Agilent 7890/5975	None

Solvent	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)	Solvent	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
Acetone (67-64-1)	-	131	262	5000	n-Heptane (142-82-5)	-	131	262	5000
Acetonitrile (75-5-8)	-	131	262	410	n-Hexane (110-54-3)	-	45.8	91.7	290
Benzene (71-43-2)	-	1.31	2.62	2	Isobutane (75-28-5)	-	131	262	5000
n-Butane (106-97-2)	-	131	262	5000	Isopropanol (67-63-0)	-	131	262	5000
1-Butanol (71-36-3)	-	131	262	5000	Isopropyl acetate (108-21-4)	-	131	262	5000
2-Butanol (78-92-2)	-	131	262	5000	Isopropyl benzene (98-82-8)	-	13.1	26.2	70
2-Butanone (78-93-3)	-	131	262	5000	Methanol (67-56-1)	-	131	262	3000
Cyclohexane (110-82-7)	-	131	262	3880	2-Methylbutane (78-78-4)	-	131	262	5000
1,2-Dimethoxyethane (110-71-4)	-	13.1	26.2	100	Methylene chloride (75-9-2)	-	131	262	600
N,N-Dimethylacetamide (127-19-5)	-	131	262	1090	2-Methylpentane (107-83-5)	-	45.8	91.7	290
2,2-Dimethylbutane (75-83-2)	-	45.8	91.7	290	3-Methylpentane (96-10-0)	-	45.8	91.7	290
2,3-Dimethylbutane (79-29-8)	-	45.8	91.7	290	n-Pentane (109-66-0)	-	131	262	5000
N,N-Dimethylformamide (68-12-2)	-	131	262	880	1-Pentanol (71-41-0)	-	131	262	5000
Dimethylsulfoxide (67-68-5)	-	131	262	5000	n-Propane (74-98-6)	-	131	262	5000
1,4-Dioxane (123-91-1)	-	131	262	380	1-Propanol (71-23-8)	-	131	262	5000
Ethanol (64-17-5)	-	131	262	5000	Pyridine (110-86-1)	-	45.8	91.7	200
2-Ethoxyethanol (110-80-5)	-	45.8	91.7	160	Tetrahydrofuran (109-99-9)	-	131	262	720
Ethyl ether (60-29-7)	-	131	262	5000	Tetramethylene sulfone (126-33-0)	-	45.8	91.7	160
Ethyl acetate (141-78-6)	-	131	262	5000	Toluene (108-88-3)	-	131	262	890
Ethyl benzene (100-41-4)	-	131	262	2170	o-Xylene (95-47-6)	-	131	262	2170
Ethylene glycol (107-21-1)	-	131	262	620	m,p-Xylene (108-38-3 or 106-42-3)	-	131	262	2170
Ethylene oxide (75-21-8)	-	13.1	26.2	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170



Color Key

RESULT < AL
RESULT > AL

"DET" detected less than LOQ

"-" not detected above LOD

"*" - o,m,p-Xylene and Ethylbenzene

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.

Solvent	Synonym(s)	Solvent	Synonym(s)
Acetonitrile	Methyl Cyanide, ACN	Ethylene glycol	1,2-Ethanediol
1-Butanol	n-Butanol, Butyl Alcohol	Isobutane	2-Methylpropane
2-Butanol	sec-Butyl alcohol	Isopropanol	2-Propanol, IPA
2-Butanone	Methyl ethyl ketone, MEK	Isopropyl Acetate	Acetic acid isopropyl ester
1,2-Dimethoxyethane	Monoglyme	Methanol	Methyl alcohol
2,3-Dimethylbutane	Neohexane	2-Methylbutane	Isopentane
2,3-Dimethylbutane	Diisopropyl	Methylene chloride	Dichloromethane
N,N-Dimethylformamide	DMF	2-Methylpentane	Isohexane
Dimethylsulfoxide	DMSO	1-Pentanol	n-Amyl alcohol
2-Ethoxyethanol	Cellosolve, Ethyl glycol	1-Propanol	Propyl alcohol
Ethyl ether	Diethyl ether, Ether	Tetrahydrofuran	THF
Ethyl acetate	EtOAc	Tetramethylene sulfone	Sulfolane
Ethyl benzene	Phenylethane	Xylene	Dimethylbenzene

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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Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries **Date Completed:** 08/25/2023

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 01/01/1900 1200
Analyst: KF

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

Deviations from SOP:
None

Pesticide	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)	Pesticide	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
Abamectin (71751-41-2)	-	0.0189	0.151	0.5	Kresoxim-methyl (143390-89-0)	-	0.0189	0.151	0.4
Acephate (30560-19-1)	-	0.0189	0.151	0.4	Malathion (121-75-5)	-	0.0189	0.151	0.2
Acetamiprid (135410-20-7)	-	0.0189	0.151	0.2	Metalaxyl (57837-19-1)	-	0.0189	0.151	0.2
Aldicarb (116-06-3)	-	0.0189	0.151	0.4	Methiocarb (2032-65-7)	-	0.0189	0.151	0.2
Azoxystrobin (131860-33-8)	-	0.0189	0.151	0.2	Methomyl (16752-77-5)	-	0.0189	0.151	0.4
Bifenazate (149877-41-8)	-	0.0189	0.151	0.2	Methyl parathion (298-0-0)	-	0.0189	0.151	0.2
Bifenthrin (82657-04-3)	-	0.0189	0.151	0.2	MGK 264 (113-48-4)	-	0.0189	0.151	0.2
Boscalid (188425-85-6)	-	0.0189	0.151	0.4	Myclobutanil (88671-89-0)	-	0.0189	0.151	0.2
Carbaryl (63-25-2)	-	0.0189	0.151	0.2	Naled (300-76-5)	-	0.0189	0.151	0.5
Carbofuran (1563-66-2)	-	0.0189	0.151	0.2	Oxamyl (23135-22-0)	-	0.0189	0.151	1
Chlorantraniliprole (800008-45-7)	-	0.0189	0.151	0.2	Pacllobutrazol (76738-62-0)	-	0.0189	0.151	0.4
Chlorfenapyr (122453-73-0)	-	0.0189	0.151	1	Permethrins (52645-53-1)	-	0.0189	0.151	0.2
Chlorpyrifos (2921-88-2)	-	0.0189	0.151	0.2	Phosmet (732-11-6)	-	0.0189	0.151	0.2
Clofentezine (74115-24-5)	-	0.0189	0.151	0.2	Piperonyl butoxide (51-03-6)	-	0.0189	0.151	2
Cyfluthrin (68359-37-5)	-	0.0189	0.151	1	Prallethrins (2331-36-9)	-	0.0189	0.151	0.2
Cypermethrin (52315-07-8)	-	0.0189	0.151	1	Propiconazole (60207-90-1)	-	0.0189	0.151	0.4
Daminozide (1596-84-5)	-	0.0189	0.151	1	Propoxur (114-26-1)	-	0.0189	0.151	0.2
DDVP (62-73-7)	-	0.0189	0.151	0.1	Pyrethrins (8003-34-7)	-	0.0189	0.151	1
Diazinon (333-41-5)	-	0.0189	0.151	0.2	Pyridaben (96489-71-3)	-	0.0189	0.151	0.2
Dimethoate (60-51-5)	-	0.0189	0.151	0.2	Spinosad (168316-95-8)	-	0.0189	0.151	0.2
Ethoprophos (13194-48-4)	-	0.0189	0.151	0.2	Spiromesifen (283594-90-1)	-	0.0189	0.151	0.2
Etofenprox (80844-07-1)	-	0.0189	0.151	0.4	Spirotetramat (203313-25-1)	-	0.0189	0.151	0.2
Etozazole (153233-91-1)	-	0.0189	0.151	0.2	Spiroxamine (118134-30-8)	-	0.0189	0.151	0.4
Fenoxycarb (72490-01-8)	-	0.0189	0.151	0.2	Tebuconazole (80443-41-0)	-	0.0189	0.151	0.4
(E)-Fenpyroximate (134098-61-6)	-	0.0189	0.151	0.4	Thiacloprid (111988-49-9)	-	0.0189	0.151	0.2
Fipronil (120068-37-3)	-	0.0189	0.151	0.4	Thiamethoxam (153719-23-4)	-	0.0189	0.151	0.2
Fonicamid (158062-67-0)	-	0.0189	0.151	1	Trifloxystrobin (141517-21-7)	-	0.0189	0.151	0.2
Fludioxinil (131341-86-1)	-	0.0189	0.151	0.4					
Hexythiazox (78587-05-0)	-	0.0189	0.151	1					
Imazalil (35554-44-0)	-	0.0189	0.151	0.2					
Imidacloprid (138261-41-3)	-	0.0189	0.151	0.4					



Color Key

RESULT < AL
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.
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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

Pesticide	Synonym(s)	Pesticide	Synonym(s)	Pesticide	Synonym(s)
Cyfluthrin	Baythroid	Myclobutanil	Sythane	Propiconazole	Tilt
DDVP	Dichlorvos	Naled	Dibrom	Propoxur	Baygon
Ethoprophos	Prophos	Phosmet	Imidan		

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Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries			Date Completed: 08/25/2023

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

Analysis Date/Time: 08/24/2023 2102 (ICP/OES) **Method:** ICP/MS **Deviations from SOP:**
Analysis Date/Time: - (DMA) **Instrument:** Agilent 7500ce None
Analyst: KF

<u>Heavy Metal</u>	<u>Result (µg/kg)</u>	<u>LOD (µg/kg)</u>	<u>LOQ (µg/kg)</u>	<u>Action Level (µg/kg)</u>
Arsenic (As)	-	58.8	93.1	200
Cadmium (Cd)	-	58.8	93.1	200
Lead (Pb)	-	58.8	93.1	500
Mercury (Hg)	-	58.8	93.1	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



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 "-" not detected above LOD

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

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Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries			

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 08/24/2023 1021 **Method:** Hardy Diagnostics CompactDry **Deviations from SOP:** None
Analyst: PW **Instrument:** Thermo Incubator

Bacteria/Microbe	Result (CFU/g)	Action Level (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	-



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)

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.

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.

REPORT OF LABORATORY ANALYSIS

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