F.A.S.T. LABORATORIES/RESEARCH

ACCREDITED CERTIFICATE #4473.01

ASA PFC Laboratory

Analyzed on behalf of:

RVR RIVER VALLEY RELIEF

SUMMARY OF ANALYSIS (SAMPLE ID: SA35822)

Testing Location:	**Customer ID:** 2168	**Order ID:** OR10570	**Sample Type:** Primary
Arkansas	River Valley Relief MIPS	**Lot Number:**	**Matrix:** Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13226761233	**Mass:** 1bag
Greenbrier, AR 72058	Fort Smith, AR 72903	**Production Run:**	**Date Collected:** 08/14/2023
License: ADH 113	License: 00065C	20230811STRAW12	**Date Received:** 08/14/2023

Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries — **Date Completed:** 08/17/2023

*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

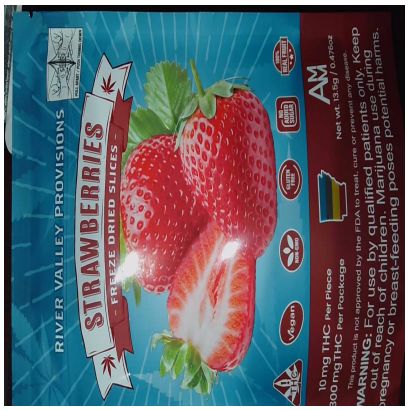
Moisture content/water activity action levels are referenced from the State of Arkansas MMJ testing guidelines.
Moisture content levels less than 15% are recommended but the sample does not fail. Water activity levels must be less than 0.65aw.

Cannabinoids (Top 3)	(%)	mg/g
Δ9-THC	0.987	9.87
CBG	0.0380	0.380
CBN	0.0153	0.153
TOTAL CBD	0.00482	0.0482
TOTAL THC	0.987	9.87
TOTAL CANNABINOIDS	1.05	10.5

Terpenes (Top 5)	(%)	µg/g
β-Caryophyllene	0.0102	102
α-Bisabolol		
Camphene		
δ-3-Carene		
Caryophyllene oxide		
TOTAL TERPENES	0.0102	102

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

Sample Picture Upon Receipt



Scan the QR code to verify results.

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Kyle W. Felling, Ph.D.
Laboratory Director

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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35822)

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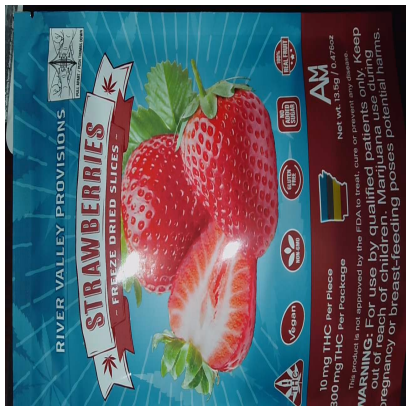
License: ADH 113	License: 00065C	20230811STRAW12	**Date Received:** 08/14/2023

Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries — **Date Completed:** 08/17/2023

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

Analysis Date/Time: 08/15/2023 1437
Analyst: PW
Method: HPLC/DAD
Instrument: Agilent 1100
Moisture Content (%): -
Water Activity (aw): -

Cannabinoid	Result (%)	Result (mg/g)	LOD (mg/g)	LOQ (mg/g)	Result (mg/mL)	Per Serving (mg)	Per Unit (mg)
CBC	ND	ND	0.00237	0.00552	-	-	-
CBCA	ND	ND	0.00735	0.0171	-	-	-
CBD	0.00482	0.0482	0.0167	0.0389	-	0.0466	1.40
CBDA	ND	ND	0.00614	0.0143	-	-	-
CBDV	ND	ND	0.00268	0.00625	-	-	-
CBDVA	ND	ND	0.00714	0.0167	-	-	-
CBG	0.0380	0.380	0.0108	0.0253	-	0.367	11.0
CBGA	ND	ND	0.0154	0.0181	-	-	-
CBL	ND	ND	0.0125	0.0292	-	-	-
CBN	0.0153	0.153	0.00575	0.0134	-	0.148	4.45
CBNA	ND	ND	0.00620	0.0145	-	-	-
Δ9-THC	0.987	9.87	0.00689	0.0161	-	9.53	286
Δ8-THC	ND	ND	0.0107	0.0251	-	-	-
THCA	ND	ND	0.00373	0.00874	-	-	-
THCV	0.00627	0.0627	0.00896	0.0209	-	0.0606	1.82
THCVA	ND	ND	0.00286	0.00666	-	-	-
TOTAL	1.05	10.5			-	10.2	305
TOTAL CBC	-	-			-	-	-
TOTAL CBD	0.00482	0.0482			-	0.0466	1.40
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	0.0380	0.380			-	0.367	11.0
TOTAL CBN	0.0153	0.153			-	0.148	4.45
TOTAL THC	0.987	9.87			-	9.53	286
TOTAL THCV	0.00627	0.0627			-	0.0606	1.82



SERVING MASS (g): 0.966
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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Kyle W. Felling, Ph.D.
Laboratory Director

www.FASTLaboratories.com

Analyzed on behalf of:

CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35822)

Testing Location:	**Customer ID:** 2168	**Order ID:** OR10570	**Sample Type:** Primary
Arkansas	River Valley Relief MIPS	**Lot Number:**	**Matrix:** Edible
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Greenbrier, AR 72058	Fort Smith, AR 72903	**Production Run:**	**Date Collected:** 08/14/2023
License: ADH 113	License: 00065C	20230811STRAW12	**Date Received:** 08/14/2023

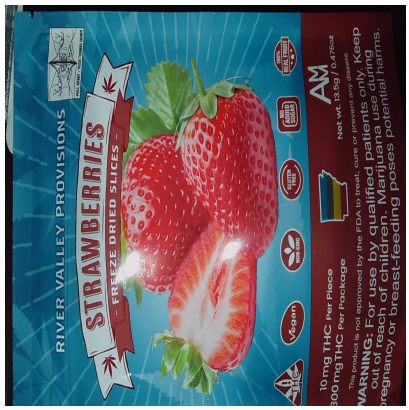
Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries — **Date Completed:** 08/17/2023

TERPENOID PROFILE

Analysis Date/Time:08/17/2023 0541 — **Method:** GC/MS — **Deviations from SOP:**

Analyst: KF — **Instrument:** Agilent 7890/5975 — None

Terpene	Result (µg/g)	Result (%)
α-Bisabolol	ND	-
Camphene	ND	-
δ-3-Carene	ND	-
β-Caryophyllene	102	0.0102
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	102	0.0102



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): 5.19

"-" Not detected above LOD.

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Laboratory Director

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ACCREDITED CERTIFICATE #4473.01

ASA PFC Laboratory

Analyzed on behalf of:

RVR RIVER VALLEY RELIEF

CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35822)

Testing Location:	**Customer ID:** 2168	**Order ID:** OR10570	**Sample Type:** Primary
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Greenbrier, AR 72058	Fort Smith, AR 72903	**Production Run:**	**Date Collected:** 08/14/2023
License: ADH 113	License: 00065C	20230811STRAW12	**Date Received:** 08/14/2023

Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries — **Date Completed:** 08/17/2023

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

Analysis Date/Time: 08/16/2023 1303 — **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)
Acetone (67-64-1)	-	159	317	5000
Acetonitrile (75-5-8)	-	159	317	410
Benzene (71-43-2)	-	1.59	3.17	2
n-Butane (106-97-2)	-	159	317	5000
1-Butanol (71-36-3)	-	159	317	5000
2-Butanol (78-92-2)	-	159	317	5000
2-Butanone (78-93-3)	-	159	317	5000
Cyclohexane (110-82-7)	-	159	317	3880
1,2-Dimethoxyethane (110-71-4)	-	15.9	31.7	100
N,N-Dimethylacetamide (127-19-5)	-	159	317	1090
2,2-Dimethylbutane (75-83-2)	-	55.6	111	290
2,3-Dimethylbutane (79-29-8)	-	55.6	111	290
N,N-Dimethylformamide (68-12-2)	-	159	317	880
Dimethylsulfoxide (67-68-5)	-	159	317	5000
1,4-Dioxane (123-91-1)	-	159	317	380
Ethanol (64-17-5)	-	159	317	5000
2-Ethoxyethanol (110-80-5)	-	55.6	111	160
Ethyl ether (60-29-7)	-	159	317	5000
Ethyl acetate (141-78-6)	-	159	317	5000
Ethyl benzene (100-41-4)	-	159	317	2170
Ethylene glycol (107-21-1)	-	159	317	620
Ethylene oxide (75-21-8)	-	15.9	31.7	50

Solvent	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
n-Heptane (142-82-5)	-	159	317	5000
n-Hexane (110-54-3)	-	55.6	111	290
Isobutane (75-28-5)	-	159	317	5000
Isopropanol (67-63-0)	-	159	317	5000
Isopropyl acetate (108-21-4)	-	159	317	5000
Isopropyl benzene (98-82-8)	-	15.9	31.7	70
Methanol (67-56-1)	-	159	317	3000
2-Methylbutane (78-78-4)	-	159	317	5000
Methylene chloride (75-9-2)	-	159	317	600
2-Methylpentane (107-83-5)	-	55.6	111	290
3-Methylpentane (96-10-0)	-	55.6	111	290
n-Pentane (109-66-0)	-	159	317	5000
1-Pentanol (71-41-0)	-	159	317	5000
n-Propane (74-98-6)	-	159	317	5000
1-Propanol (71-23-8)	-	159	317	5000
Pyridine (110-86-1)	-	55.6	111	200
Tetrahydrofuran (109-99-9)	-	159	317	720
Tetramethylene sulfone (126-33-0)	-	55.6	111	160
Toluene (108-88-3)	-	159	317	890
o-Xylene (95-47-6)	-	159	317	2170
m,p-Xylene (108-38-3 or 106-42-3)	-	159	317	2170
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
Dimethysufoxide	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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Analyzed on behalf of: RVR River Valley Relief

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Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries — **Date Completed:** 08/17/2023

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 08/15/2023 1458 — **Method:** LC/MS/MS — **Deviations from SOP:** None

Analyst: KF — **Instrument:** Shimadzu LC-8050

Pesticide	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
Abamectin (71751-41-2)	-	0.0100	0.0803	0.5
Acephate (30560-19-1)	-	0.0100	0.0803	0.4
Acequinocyl (57960-19-7)	-	0.0100	0.0803	2
Acetamiprid (135410-20-7)	-	0.0100	0.0803	0.2
Aldicarb (116-06-3)	-	0.0100	0.0803	0.4
Azoxystrobin (131860-33-8)	-	0.0100	0.0803	0.2
Bifenazate (149877-41-8)	-	0.0100	0.0803	0.2
Bifenthrin (82657-04-3)	-	0.0100	0.0803	0.2
Boscalid (188425-85-6)	-	0.0100	0.0803	0.4
Carbaryl (63-25-2)	-	0.0100	0.0803	0.2
Carbofuran (1563-66-2)	-	0.0100	0.0803	0.2
Chlorantraniliprole (800008-45-7)	-	0.0100	0.0803	0.2
Chlorfenapyr (122453-73-0)	-	0.0100	0.0803	1
Chlorpyrifos (2921-88-2)	-	0.0100	0.0803	0.2
Clofentezine (74115-24-5)	-	0.0100	0.0803	0.2
Cyfluthrin (68359-37-5)	-	0.0100	0.0803	1
Cypermethrin (52315-07-8)	-	0.0100	0.0803	1
Daminozide (1596-84-5)	-	0.0100	0.0803	1
DDVP (62-73-7)	-	0.0100	0.0803	0.1
Diazinon (333-41-5)	-	0.0100	0.0803	0.2
Dimethoate (60-51-5)	-	0.0100	0.0803	0.2
Ethoprophos (13194-48-4)	-	0.0100	0.0803	0.2
Etofenprox (80844-07-1)	-	0.0100	0.0803	0.4
Etoxazole (153233-91-1)	-	0.0100	0.0803	0.2
Fenoxycarb (72490-01-8)	-	0.0100	0.0803	0.2
(E)-Fenpyroximate (134098-61-6)	-	0.0100	0.0803	0.4
Fipronil (120068-37-3)	-	0.0100	0.0803	0.4
Flonicamid (158062-67-0)	-	0.0100	0.0803	1
Fludioxinil (131341-86-1)	-	0.0100	0.0803	0.4
Hexythiazox (78587-05-0)	-	0.0100	0.0803	1
Imazalil (35554-44-0)	-	0.0100	0.0803	0.2
Imidacloprid (138261-41-3)	-	0.0100	0.0803	0.4

Pesticide	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
Kresoxim-methyl (143390-89-0)	-	0.0100	0.0803	0.4
Malathion (121-75-5)	-	0.0100	0.0803	0.2
Metalaxyl (57837-19-1)	-	0.0100	0.0803	0.2
Methiocarb (2032-65-7)	-	0.0100	0.0803	0.2
Methomyl (16752-77-5)	-	0.0100	0.0803	0.4
Methyl parathion (298-0-0)	-	0.0100	0.0803	0.2
MGK 264 (113-48-4)	-	0.0100	0.0803	0.2
Myclobutanil (88671-89-0)	-	0.0100	0.0803	0.2
Naled (300-76-5)	-	0.0100	0.0803	0.5
Oxamyl (23135-22-0)	-	0.0100	0.0803	1
Paclobutrazol (76738-62-0)	-	0.0100	0.0803	0.4
Permethrins (52645-53-1)	-	0.0100	0.0803	0.2
Phosmet (732-11-6)	-	0.0100	0.0803	0.2
Piperonyl butoxide (51-03-6)	-	0.0100	0.0803	2
Prallethrins (2331-36-9)	-	0.0100	0.0803	0.2
Propiconazole (60207-90-1))	-	0.0100	0.0803	0.4
Propoxur (114-26-1)	-	0.0100	0.0803	0.2
Pyrethrins (8003-34-7)	-	0.0100	0.0803	1
Pyridaben (96489-71-3)	-	0.0100	0.0803	0.2
Spinosad (168316-95-8)	-	0.0100	0.0803	0.2
Spiromesifen (283594-90-1)	-	0.0100	0.0803	0.2
Spirotetramat (203313-25-1)	-	0.0100	0.0803	0.2
Spiroxamine (118134-30-8)	-	0.0100	0.0803	0.4
Tebuconazole (80443-41-0)	-	0.0100	0.0803	0.4
Thiacloprid (111988-49-9)	-	0.0100	0.0803	0.2
Thiamethoxam (153719-23-4)	-	0.0100	0.0803	0.2
Trifloxystrobin (141517-21-7)	-	0.0100	0.0803	0.2

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	Systhane	Propiconazole	Tilt
DDVP	Dichlorvos	Naled	Dibrom	Propoxur	Baygon
Ethoprophos	Prophos	Phosmet	Imidan		

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Laboratory Director

F.A.S.T. LABORATORIES/RESEARCH

ACCREDITED CERTIFICATE #4473.01

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RVR RIVER VALLEY RELIEF

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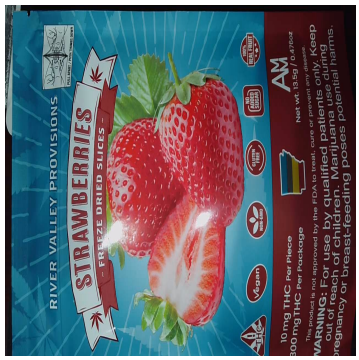
HEAVY METAL PROFILE (SOP: SOP-ICP-200.6)

Analysis Date/Time: 08/15/2023 1907 (ICP/OES)
Analysis Date/Time: - (DMA)
Analyst: KF

Method: ICP/MS
Instrument: Agilent 7500ce

Deviations from SOP:
None

Heavy Metal	Result (µg/kg)	LOD (µg/kg)	LOQ (µg/kg)	Action Level (µg/kg)
Arsenic (As)	-	57.4	90.8	200
Cadmium (Cd)	DET	57.4	90.8	200
Lead (Pb)	-	57.4	90.8	500
Mercury (Hg)	-	57.4	90.8	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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Kyle W. Felling, Ph.D.
Laboratory Director

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F.A.S.T. LABORATORIES/RESEARCH

ACCREDITED CERTIFICATE #4473.01

Analyzed on behalf of:

CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35822)

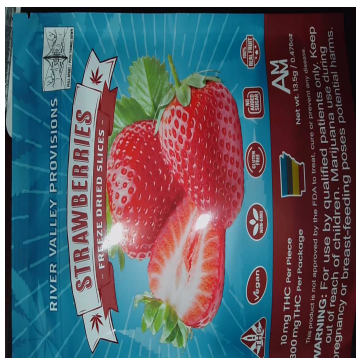
Testing Location:	**Customer ID:** 2168	**Sample ID:** SA35822	**Sample Type:** Primary
Arkansas	River Valley Relief MIPS	**Lot Number:**	**Matrix:** Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13226761233	**Mass:** 1bag
Greenbrier, AR 72058	Fort Smith, AR 72903	**Production Run:**	**Date Collected:** 08/14/2023
License: ADA 05_H273	License: 00065C	20230811STRAW12	**Date Received:** 08/14/2023

Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries — **Date Completed:** 08/17/2023

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 08/16/2023 0933 — **Method:** Hardy Diagnostics CompactDry — **Deviations from SOP:**
Analyst: PW — **Instrument:** Thermo Incubator — None

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.

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