

SUMMARY OF ANALYSIS (SAMPLE ID: SA36516)

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: OR10727 Lot Number: M00065C13225221421 Production Run: 20231120straw26	Sample Type: Primary Matrix: Edible Mass: 1bag Date Collected: 11/27/2023 Date Received: 11/27/2023
Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries			Date Completed: 11/29/2023

*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 (%)

0.000

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
CBD	0.000	0.000
CBDA	0.000	0.000
CBDV	0.000	0.000
TOTAL CBD	0.000	0.000
TOTAL THC	0.000	0.000
TOTAL CANNABINOIDS	0.000	0.000

Terpenes (Top 5)	(%)	µg/g
β-Caryophyllene	0.00811	81.1
Guaiol	0.00695	69.5
α-Bisabolol	0.00502	50.2
α-Humulene	0.00347	34.7
Camphene		
TOTAL TERPENES	0.0235	235

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



Scan the QR code to verify results.

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REPORT OF LABORATORY ANALYSIS

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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:** 11/29/2023

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

Analysis Date/Time: 11/28/2023 1253 **Method:** HPLC/DAD **Moisture Content (%):** 0.000
Analyst: PW **Instrument:** Agilent 1100 **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	0.000	0.000	0.00289	0.00674	-	-	-
CBCA	0.000	0.000	0.00897	0.0209	-	-	-
CBD	0.000	0.000	0.0204	0.0475	-	-	-
CBDA	0.000	0.000	0.00750	0.0175	-	-	-
CBDV	0.000	0.000	0.00327	0.00763	-	-	-
CBDVA	0.000	0.000	0.00871	0.0203	-	-	-
CBG	0.000	0.000	0.0132	0.0309	-	0.0887	2.66
CBGA	0.000	0.000	0.0187	0.0221	-	-	-
CBL	0.000	0.000	0.0153	0.0357	-	-	-
CBN	0.000	0.000	0.00702	0.0164	-	0.0903	2.71
CBNA	0.000	0.000	0.00757	0.0177	-	-	-
Δ9-THC	0.000	0.000	0.00841	0.0196	-	9.74	292
Δ8-THC	0.000	0.000	0.0131	0.0306	-	0.484	14.5
THCA	0.000	0.000	0.00456	0.0107	-	-	-
THCV	0.000	0.000	0.0109	0.0255	-	0.0613	1.84
THCVA	0.000	0.000	0.00350	0.00813	-	-	-
TOTAL	0.000	0.000			-	10.5	314
TOTAL CBC	0.000	0.000			-	-	-
TOTAL CBD	0.000	0.000			-	-	-
TOTAL CBDV	0.000	0.000			-	-	-
TOTAL CBG	0.000	0.000			-	0.0887	2.66
TOTAL CBN	0.000	0.000			-	0.0903	2.71
TOTAL THC	0.000	0.000			-	9.74	292
TOTAL THC V	0.000	0.000			-	0.0613	1.84



SERVING MASS (g): 0.760
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 THC V = (THCVA x 0.867) + THC V

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:** 11/29/2023

TERPENOID PROFILE

Analysis Date/Time: 11/28/2023 2048
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	50.2	0.00502	
Camphene	ND	-	
δ-3-Carene	ND	-	
β-Caryophyllene	81.1	0.00811	
Caryophyllene oxide	ND	-	
p-Cymene	ND	-	
Eucalyptol	ND	-	
Geraniol	ND	-	
Guaiol	69.5	0.00695	
α-Humulene	34.7	0.00347	
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	235	0.0235	



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): 9.65

"-" Not detected above LOD.

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Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries **Date Completed:** 11/29/2023

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

Analysis Date/Time: 11/28/2023 1807 **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)	-	111	221	5000	n-Heptane (142-82-5)	-	111	221	5000
Acetonitrile (75-5-8)	-	111	221	410	n-Hexane (110-54-3)	-	38.7	77.5	290
Benzene (71-43-2)	-	1.11	2.21	2	Isobutane (75-28-5)	-	111	221	5000
n-Butane (106-97-2)	454	111	221	5000	Isopropanol (67-63-0)	-	111	221	5000
1-Butanol (71-36-3)	-	111	221	5000	Isopropyl acetate (108-21-4)	-	111	221	5000
2-Butanol (78-92-2)	-	111	221	5000	Isopropyl benzene (98-82-8)	-	11.1	22.1	70
2-Butanone (78-93-3)	-	111	221	5000	Methanol (67-56-1)	-	111	221	3000
Cyclohexane (110-82-7)	-	111	221	3880	2-Methylbutane (78-78-4)	-	111	221	5000
1,2-Dimethoxyethane (110-71-4)	-	11.1	22.1	100	Methylene chloride (75-9-2)	-	111	221	600
N,N-Dimethylacetamide (127-19-5)	-	111	221	1090	2-Methylpentane (107-83-5)	-	38.7	77.5	290
2,2-Dimethylbutane (75-83-2)	-	38.7	77.5	290	3-Methylpentane (96-10-0)	-	38.7	77.5	290
2,3-Dimethylbutane (79-29-8)	-	38.7	77.5	290	n-Pentane (109-66-0)	-	111	221	5000
N,N-Dimethylformamide (68-12-2)	-	111	221	880	1-Pentanol (71-41-0)	-	111	221	5000
Dimethylsulfoxide (67-68-5)	-	111	221	5000	n-Propane (74-98-6)	-	111	221	5000
1,4-Dioxane (123-91-1)	-	111	221	380	1-Propanol (71-23-8)	-	111	221	5000
Ethanol (64-17-5)	1660	111	221	5000	Pyridine (110-86-1)	-	38.7	77.5	200
2-Ethoxyethanol (110-80-5)	-	38.7	77.5	160	Tetrahydrofuran (109-99-9)	-	111	221	720
Ethyl ether (60-29-7)	-	111	221	5000	Tetramethylene sulfone (126-33-0)	-	38.7	77.5	160
Ethyl acetate (141-78-6)	-	111	221	5000	Toluene (108-88-3)	-	111	221	890
Ethyl benzene (100-41-4)	-	111	221	2170	o-Xylene (95-47-6)	-	111	221	2170
Ethylene glycol (107-21-1)	-	111	221	620	m,p-Xylene (108-38-3 or 106-42-3)	-	111	221	2170
Ethylene oxide (75-21-8)	-	11.1	22.1	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170



Color Key

RESULT < AL (Green background)
RESULT > AL (Red background)

"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:** 11/29/2023

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 11/28/2023 1243
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)	-	0.0128	0.103	0.5	Kresoxim-methyl (143390-89-0)	-	0.0128	0.103	0.4
Acephate (30560-19-1)	-	0.0128	0.103	0.4	Malathion (121-75-5)	-	0.0128	0.103	0.2
Acequinocyl (57960-19-7)	-	0.0128	0.103	2	Metalaxyl (57837-19-1)	-	0.0128	0.103	0.2
Acetamiprid (135410-20-7)	-	0.0128	0.103	0.2	Methiocarb (2032-65-7)	-	0.0128	0.103	0.2
Aldicarb (116-06-3)	-	0.0128	0.103	0.4	Methomyl (16752-77-5)	-	0.0128	0.103	0.4
Azoxystrobin (131860-33-8)	-	0.0128	0.103	0.2	Methyl parathion (298-0-0)	-	0.0128	0.103	0.2
Bifenazate (149877-41-8)	-	0.0128	0.103	0.2	MGK 264 (113-48-4)	-	0.0128	0.103	0.2
Bifenthrin (82657-04-3)	-	0.0128	0.103	0.2	Myclobutanil (88671-89-0)	-	0.0128	0.103	0.2
Boscalid (188425-85-6)	-	0.0128	0.103	0.4	Naled (300-76-5)	-	0.0128	0.103	0.5
Carbaryl (63-25-2)	-	0.0128	0.103	0.2	Oxamyl (23135-22-0)	-	0.0128	0.103	1
Carbofuran (1563-66-2)	-	0.0128	0.103	0.2	Pacllobutrazol (76738-62-0)	-	0.0128	0.103	0.4
Chlorantraniliprole (800008-45-7)	-	0.0128	0.103	0.2	Permethrins (52645-53-1)	-	0.0128	0.103	0.2
Chlorfenapyr (122453-73-0)	-	0.0128	0.103	1	Phosmet (732-11-6)	-	0.0128	0.103	0.2
Chlorpyrifos (2921-88-2)	-	0.0128	0.103	0.2	Piperonyl butoxide (51-03-6)	-	0.0128	0.103	2
Clofentezine (74115-24-5)	-	0.0128	0.103	0.2	Prallethrins (2331-36-9)	-	0.0128	0.103	0.2
Cyfluthrin (68359-37-5)	-	0.0128	0.103	1	Propiconazole (60207-90-1)	-	0.0128	0.103	0.4
Cypermethrin (52315-07-8)	-	0.0128	0.103	1	Propoxur (114-26-1)	-	0.0128	0.103	0.2
Daminozide (1596-84-5)	-	0.0128	0.103	1	Pyrethrins (8003-34-7)	-	0.0128	0.103	1
DDVP (62-73-7)	-	0.0128	0.103	0.1	Pyridaben (96489-71-3)	-	0.0128	0.103	0.2
Diazinon (333-41-5)	-	0.0128	0.103	0.2	Spinosad (168316-95-8)	-	0.0128	0.103	0.2
Dimethoate (60-51-5)	-	0.0128	0.103	0.2	Spiromesifen (283594-90-1)	-	0.0128	0.103	0.2
Ethoprophos (13194-48-4)	-	0.0128	0.103	0.2	Spirotetramat (203313-25-1)	-	0.0128	0.103	0.2
Etofenprox (80844-07-1)	-	0.0128	0.103	0.4	Spiroxamine (118134-30-8)	-	0.0128	0.103	0.4
Etoxazole (153233-91-1)	-	0.0128	0.103	0.2	Tebuconazole (80443-41-0)	-	0.0128	0.103	0.4
Fenoxycarb (72490-01-8)	-	0.0128	0.103	0.2	Thiacloprid (111988-49-9)	-	0.0128	0.103	0.2
(E)-Fenpyroximate (134098-61-6)	-	0.0128	0.103	0.4	Thiamethoxam (153719-23-4)	-	0.0128	0.103	0.2
Fipronil (120068-37-3)	-	0.0128	0.103	0.4	Trifloxystrobin (141517-21-7)	-	0.0128	0.103	0.2
Fonicamid (158062-67-0)	-	0.0128	0.103	1					
Fludioxinil (131341-86-1)	-	0.0128	0.103	0.4					
Hexythiazox (78587-05-0)	-	0.0128	0.103	1					
Imazalil (35554-44-0)	-	0.0128	0.103	0.2					
Imidacloprid (138261-41-3)	-	0.0128	0.103	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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HEAVY METAL PROFILE (SOP: SOP-ICP-200.6)

Analysis Date/Time: 11/28/2023 1830 (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)	-	56.1	88.9	200
Cadmium (Cd)	-	56.1	88.9	200
Lead (Pb)	-	56.1	88.9	500
Mercury (Hg)	-	56.1	88.9	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

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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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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36516)

Testing Location: Arkansas 232 S. Broadview St. Greenbrier, AR 72058 License: ADA 05_H273	Customer ID: 2168 River Valley Relief MIPS 5601 Old Greenwood Rd Suite C Fort Smith, AR 72903 License: 00065C	Sample ID: SA36516 Lot Number: M00065C13225221421 Production Run: 20231120straw26	Sample Type: Primary Matrix: Edible Mass: 1bag Date Collected: 11/27/2023 Date Received: 11/27/2023
Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries			Date Completed: 11/29/2023

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 11/29/2023 1026 **Method:** Hardy Diagnostics CompactDry **Deviations from SOP:**
Analyst: PW **Instrument:** Thermo Incubator None

<u>Bacteria/Microbe</u>	<u>Result (CFU/g)</u>	<u>Action Level (CFU/g)</u>
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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Kyle W. Felling
Kyle W. Felling, Ph.D.
Laboratory Director

