

SUMMARY OF ANALYSIS (SAMPLE ID: SA35964)

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: OR10599 Lot Number: M00065C13226721777 Production Run: 20230830straw17	Sample Type: Primary Matrix: Edible Mass: 1bag Date Collected: 09/05/2023 Date Received: 09/05/2023
Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries			Date Completed: 09/08/2023

*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	1.60	16.0
CBG	0.0611	0.611
CBN	0.0257	0.257
TOTAL CBD	0.0165	0.165
TOTAL THC	1.60	16.0
TOTAL CANNABINOIDS	1.71	17.1

Terpenes (Top 5)	(%)	µg/g
γ-Terpinene	0.0358	358
β-Caryophyllene	0.0328	328
β-Pinene	0.0159	159
α-Humulene	0.0108	108
α-Bisabolol		
TOTAL TERPENES	0.0952	952

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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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35964)

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: OR10599 Lot Number: M00065C13226721777 Production Run: 20230830straw17	Sample Type: Primary Matrix: Edible Mass: 1bag Date Collected: 09/05/2023 Date Received: 09/05/2023
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Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries **Date Completed:** 09/08/2023

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

Analysis Date/Time: 09/06/2023 1523
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	0.00379	0.00884	-	-	-
CBCA	ND	ND	0.0118	0.0275	-	-	-
CBD	0.0165	0.165	0.0267	0.0624	-	0.0991	2.97
CBDa	ND	ND	0.00984	0.0230	-	-	-
CBDV	ND	ND	0.00429	0.0100	-	-	-
CBDVA	ND	ND	0.0114	0.0267	-	-	-
CBG	0.0611	0.611	0.0174	0.0405	-	0.368	11.0
CBGA	ND	ND	0.0246	0.0290	-	-	-
CBL	ND	ND	0.0201	0.0468	-	-	-
CBN	0.0257	0.257	0.00921	0.0215	-	0.155	4.65
CBNA	ND	ND	0.00994	0.0232	-	-	-
Δ9-THC	1.60	16.0	0.0110	0.0257	-	9.63	289
Δ8-THC	ND	ND	0.0172	0.0402	-	-	-
THCA	ND	ND	0.00599	0.0140	-	-	-
THCV	0.00938	0.0938	0.0144	0.0335	-	0.0565	1.69
THCVA	ND	ND	0.00459	0.0107	-	-	-
TOTAL	1.71	17.1				10.3	309
TOTAL CBC	-	-				-	-
TOTAL CBD	0.0165	0.165				0.0991	2.97
TOTAL CBDV	-	-				-	-
TOTAL CBG	0.0611	0.611				0.368	11.0
TOTAL CBN	0.0257	0.257				0.155	4.65
TOTAL THC	1.60	16.0				9.63	289
TOTAL THCv	0.00938	0.0938				0.0565	1.69



SERVING MASS (g): 0.602
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:** 09/08/2023

TERPENOID PROFILE

Analysis Date/Time: 09/06/2023 1818
Analyst: KF

Method: GC/MS
Instrument: Agilent 7890/5975

Deviations from SOP:
None

Terpene	Result (µg/g)	Result (%)	
α-Bisabolol	ND	-	
Camphene	ND	-	
δ-3-Carene	ND	-	
β-Caryophyllene	328	0.0328	Orange bar
Caryophyllene oxide	ND	-	
p-Cymene	ND	-	
Eucalyptol	ND	-	
Geraniol	ND	-	
Guaiol	ND	-	
α-Humulene	108	0.0108	Green bar
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	159	0.0159	Green bar
α-Terpinene	ND	-	
γ-Terpinene	358	0.0358	Yellow bar
Terpinolene	ND	-	
TOTAL	952	0.0952	



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): 12.8

"-" Not detected above LOD.

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

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

Analysis Date/Time: 09/07/2023 0204	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)	-	163	326	5000	n-Heptane (142-82-5)	-	163	326	5000
Acetonitrile (75-5-8)	-	163	326	410	n-Hexane (110-54-3)	-	57.0	114	290
Benzene (71-43-2)	-	1.63	3.26	2	Isobutane (75-28-5)	-	163	326	5000
n-Butane (106-97-2)	-	163	326	5000	Isopropanol (67-63-0)	-	163	326	5000
1-Butanol (71-36-3)	-	163	326	5000	Isopropyl acetate (108-21-4)	-	163	326	5000
2-Butanol (78-92-2)	-	163	326	5000	Isopropyl benzene (98-82-8)	-	16.3	32.6	70
2-Butanone (78-93-3)	-	163	326	5000	Methanol (67-56-1)	-	163	326	3000
Cyclohexane (110-82-7)	-	163	326	3880	2-Methylbutane (78-78-4)	-	163	326	5000
1,2-Dimethoxyethane (110-71-4)	-	16.3	32.6	100	Methylene chloride (75-9-2)	-	163	326	600
N,N-Dimethylacetamide (127-19-5)	-	163	326	1090	2-Methylpentane (107-83-5)	-	57.0	114	290
2,2-Dimethylbutane (75-83-2)	-	57.0	114	290	3-Methylpentane (96-10-0)	-	57.0	114	290
2,3-Dimethylbutane (79-29-8)	-	57.0	114	290	n-Pentane (109-66-0)	-	163	326	5000
N,N-Dimethylformamide (68-12-2)	-	163	326	880	1-Pentanol (71-41-0)	-	163	326	5000
Dimethylsulfoxide (67-68-5)	-	163	326	5000	n-Propane (74-98-6)	-	163	326	5000
1,4-Dioxane (123-91-1)	-	163	326	380	1-Propanol (71-23-8)	-	163	326	5000
Ethanol (64-17-5)	-	163	326	5000	Pyridine (110-86-1)	-	57.0	114	200
2-Ethoxyethanol (110-80-5)	-	57.0	114	160	Tetrahydrofuran (109-99-9)	-	163	326	720
Ethyl ether (60-29-7)	-	163	326	5000	Tetramethylene sulfone (126-33-0)	-	57.0	114	160
Ethyl acetate (141-78-6)	-	163	326	5000	Toluene (108-88-3)	-	163	326	890
Ethyl benzene (100-41-4)	-	163	326	2170	o-Xylene (95-47-6)	-	163	326	2170
Ethylene glycol (107-21-1)	-	163	326	620	m,p-Xylene (108-38-3 or 106-42-3)	-	163	326	2170
Ethylene oxide (75-21-8)	-	16.3	32.6	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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Kyle W. Felling, Ph.D.
Laboratory Director



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Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries **Date Completed:** 09/08/2023

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 09/06/2023 1511
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)	-	0.0166	0.133	0.5	Kresoxim-methyl (143390-89-0)	-	0.0166	0.133	0.4
Acephate (30560-19-1)	-	0.0166	0.133	0.4	Malathion (121-75-5)	-	0.0166	0.133	0.2
Acequinocyl (57960-19-7)	-	0.0166	0.133	2	Metalaxyl (57837-19-1)	-	0.0166	0.133	0.2
Acetamiprid (135410-20-7)	-	0.0166	0.133	0.2	Methiocarb (2032-65-7)	-	0.0166	0.133	0.2
Aldicarb (116-06-3)	-	0.0166	0.133	0.4	Methomyl (16752-77-5)	-	0.0166	0.133	0.4
Azoxystrobin (131860-33-8)	-	0.0166	0.133	0.2	Methyl parathion (298-0-0)	-	0.0166	0.133	0.2
Bifenazate (149877-41-8)	-	0.0166	0.133	0.2	MGK 264 (113-48-4)	-	0.0166	0.133	0.2
Bifenthrin (82657-04-3)	-	0.0166	0.133	0.2	Myclobutanil (88671-89-0)	-	0.0166	0.133	0.2
Boscalid (188425-85-6)	-	0.0166	0.133	0.4	Naled (300-76-5)	-	0.0166	0.133	0.5
Carbaryl (63-25-2)	-	0.0166	0.133	0.2	Oxamyl (23135-22-0)	-	0.0166	0.133	1
Carbofuran (1563-66-2)	-	0.0166	0.133	0.2	Pacllobutrazol (76738-62-0)	-	0.0166	0.133	0.4
Chlorantraniliprole (800008-45-7)	-	0.0166	0.133	0.2	Permethrins (52645-53-1)	-	0.0166	0.133	0.2
Chlorfenapyr (122453-73-0)	-	0.0166	0.133	1	Phosmet (732-11-6)	-	0.0166	0.133	0.2
Chlorpyrifos (2921-88-2)	-	0.0166	0.133	0.2	Piperonyl butoxide (51-03-6)	-	0.0166	0.133	2
Clofentezine (74115-24-5)	-	0.0166	0.133	0.2	Prallethrins (2331-36-9)	-	0.0166	0.133	0.2
Cyfluthrin (68359-37-5)	-	0.0166	0.133	1	Propiconazole (60207-90-1)	-	0.0166	0.133	0.4
Cypermethrin (52315-07-8)	-	0.0166	0.133	1	Propoxur (114-26-1)	-	0.0166	0.133	0.2
Daminozide (1596-84-5)	-	0.0166	0.133	1	Pyrethrins (8003-34-7)	-	0.0166	0.133	1
DDVP (62-73-7)	-	0.0166	0.133	0.1	Pyridaben (96489-71-3)	-	0.0166	0.133	0.2
Diazinon (333-41-5)	-	0.0166	0.133	0.2	Spinosad (168316-95-8)	-	0.0166	0.133	0.2
Dimethoate (60-51-5)	-	0.0166	0.133	0.2	Spiromesifen (283594-90-1)	-	0.0166	0.133	0.2
Ethoprophos (13194-48-4)	-	0.0166	0.133	0.2	Spirotetramat (203313-25-1)	-	0.0166	0.133	0.2
Etofenprox (80844-07-1)	-	0.0166	0.133	0.4	Spiroxamine (118134-30-8)	-	0.0166	0.133	0.4
Etoxazole (153233-91-1)	-	0.0166	0.133	0.2	Tebuconazole (80443-41-0)	-	0.0166	0.133	0.4
Fenoxycarb (72490-01-8)	-	0.0166	0.133	0.2	Thiacloprid (111988-49-9)	-	0.0166	0.133	0.2
(E)-Fenpyroximate (134098-61-6)	-	0.0166	0.133	0.4	Thiamethoxam (153719-23-4)	-	0.0166	0.133	0.2
Fipronil (120068-37-3)	-	0.0166	0.133	0.4	Trifloxystrobin (141517-21-7)	-	0.0166	0.133	0.2
Fonicamid (158062-67-0)	-	0.0166	0.133	1					
Fludioxinil (131341-86-1)	-	0.0166	0.133	0.4					
Hexythiazox (78587-05-0)	-	0.0166	0.133	1					
Imazalil (35554-44-0)	-	0.0166	0.133	0.2					
Imidacloprid (138261-41-3)	-	0.0166	0.133	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.
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

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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Laboratory Director



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Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries			Date Completed: 09/08/2023

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

Analysis Date/Time: 09/06/2023 1826 (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)	-	59.5	94.2	200
Cadmium (Cd)	-	59.5	94.2	200
Lead (Pb)	-	59.5	94.2	500
Mercury (Hg)	-	59.5	94.2	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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Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries			

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 09/07/2023 0935 **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



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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