







SUMMARY OF ANALYSIS (SAMPLE ID: SA36123)

Testing Location:Customer ID: 2168Order ID: OR10628Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible

 232 S. Broadview St.
 5601 Old Greenwood Rd Suite C
 M00065C13229275793
 Mass: 1ea

 Greenbrier, AR 72058
 Fort Smith, AR 72903
 Production Run:
 Date Collected: 10/02/2023

 License: ADH 113
 License: 00065C
 20230925STRAW22
 Date Received: 10/03/2023

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

Date Completed: 10/04/2023

*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/FAILNot TestedNot TestedPASS

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>	ds (Top 3)	<u>(%)</u>	mg/g	
Δ9-Τ	HC	0.551	5.51	
CB	D	0.0430	0.430	
CBo	\hat{G}	0.0194	0.194	
TOTAL	CBD	0.0430	0.430	
TOTAL	THC	0.551	5.51	
TOTAL CANN	IABINOIDS	0.623	6.23	
<u>Terpenes</u>	(Top 5)	(%)	<u>µg</u> /g	
α-Bisa	polol			
Campl	nene			
δ-3-Ca	rene			
β-Caryop	nyllene			
Caryophyllo	ene oxide			
TOTAL TE	RPENES	-	-	
Contaminants	PASS/FAIL	Sample Picture Upon Receipt		
Hoarn Motale	DACC			

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





Scan the QR code to verify results.

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License: ADH 113







CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36123)

Order ID: OR10628 **Testing Location:** Customer ID: 2168 Arkansas River Valley Relief MIPS Lot Number: 232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13229275793 Greenbrier, AR 72058

Fort Smith, AR 72903 **Production Run:** 20230925STRAW22 License: 00065C

Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries **Date Completed:** 10/04/2023

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

Analysis Date/Time: 10/03/2023 1454 Method: HPLC/DAD **Analyst: PW Instrument:** Agilent 1100

<u>Cannabinoid</u>	Result (%)	Result (mg/g)	LOD (mg/g)	LOQ (mg/g)	Result (mg/ mL)	<u>Per</u> Serving (mg)	Per Unit (mg)
CBC	ND	ND	0.00232	0.00542	-	-	-
CBCA	ND	ND	0.00721	0.0168	-	-	-
CBD	0.0430	0.430	0.0164	0.0383	-	0.718	21.5
CBDA	ND	ND	0.00603	0.0141	-	-	-
CBDV	ND	ND	0.00263	0.00613	-	-	-
CBDVA	ND	ND	0.00701	0.0164	-	-	-
CBG	0.0194	0.194	0.0106	0.0248	-	0.324	9.72
CBGA	ND	ND	0.0151	0.0178	-	-	-
CBL	ND	ND	0.0123	0.0287	-	-	-
CBN	0.00571	0.0571	0.00565	0.0132	-	0.0953	2.86
CBNA	ND	ND	0.00609	0.0142	-	-	-
Δ9-ΤΗС	0.551	5.51	0.00677	0.0158	-	9.20	276
Δ8-ΤΗС	ND	ND	0.0106	0.0246	-	-	-
THCA	ND	ND	0.00367	0.00858	-	-	-
THCV	0.00432	0.0432	0.00880	0.0205	-	0.0722	2.17
THCVA	ND	ND	0.00281	0.00654	-	-	-
TOTAL	0.623	6.23			-	10.4	312
TOTAL CBC	-	-			-	_	-
TOTAL CBD	0.0430	0.430			-	0.718	21.5
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	0.0194	0.194			-	0.324	9.72
TOTAL CBN	0.00571	0.0571			-	0.0953	2.86
TOTAL THC	0.551	5.51			-	9.20	276
TOTAL THCV	0.00432	0.0432			-	0.0722	2.17

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): -



Sample Type: Primary

Date Collected: 10/02/2023

Date Received: 10/03/2023

Matrix: Edible

Mass: 1ea

SERVING MASS (g): 1.67 **SERVINGS/UNIT: 30**

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 = $(CBCA \times 0.877) + CBC$ Total CBD = $(CBDA \times 0.877) + CBD$ Total CBDV = $(CBDVA \times 0.867) + CBDV$ Total CBG = $(CBGA \times 0.878) + CBG$ Total CBN = $(CBNA \times 0.876) + CBN$ Total THC = (THCA x 0.877) + $\Delta 9$ -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.





Page 2 of 7

10/04/2023

[&]quot;-" Not detected above LOD.



License: ADH 113







CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36123)

Customer ID: 2168 **Testing Location:** Arkansas River Valley Relief MIPS 232 S. Broadview St. 5601 Old Greenwood Rd Suite C Fort Smith, AR 72903 Greenbrier, AR 72058

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

Order ID: OR10628 Lot Number: M00065C13229275793 **Production Run:** 20230925STRAW22

Matrix: Edible Mass: 1ea **Date Collected:** 10/02/2023 **Date Received:** 10/03/2023

Sample Type: Primary

Date Completed: 10/04/2023

TERPENOID PROFILE

Analysis Date/Time:10/04/2023 0056 Method: GC/MS **Deviations from SOP:**

Analyst: KF Instrument: Agilent 7890/5975 None

License: 00065C

<u>Terpene</u>	<u>Result</u> (μg/g)	Result (%)	RIVER VALLEY PROVISION
α-Bisabolol	ND	-	KIVER VALLET PROVISION
Camphene	<loq< td=""><td>-</td><td>TO AMPEDOLE</td></loq<>	-	TO AMPEDOLE
δ-3-Carene	ND	-	CLKAMRFKKIEG J
β-Caryophyllene	<loq< td=""><td>-</td><td>of fittee a sectiff.</td></loq<>	-	of fittee a sectiff.
Caryophyllene oxide	ND	-	FREEZE DRIED SLICES.
p-Cymene	ND	-	
Eucalyptol	ND	-	0.0000
Geraniol	ND	-	
Guaiol	ND	-	
α-Humulene	<loq< td=""><td>-</td><td>Abbreviations: GC - Gas Chromatography, MS - Mass</td></loq<>	-	Abbreviations: GC - Gas Chromatography, MS - Mass
Isopulegol	ND	-	Spectrometry, RL - Reporting Limit
d-Limonene	ND	-	Abbreviations: ND - Not Detected, , LOD - Limit of Detection, LOQ - Lin
Linalool	ND	-	of Quantitation
β-Myrcene	ND	-	This information is provided as a serv and makes no claims of efficacy and/o
cis-Nerolidol	ND	-	safety of this product.
trans-Nerolidol	ND	-	Results are applicable only for the
α-Ocimene	ND	-	sample(s) analyzed and for the specifi analysis conducted.
β-Ocimene	ND	-	This report is for informational purpo
α-Pinene	<loq< td=""><td>-</td><td>only and should not be used to diagnot treat, or prevent any</td></loq<>	-	only and should not be used to diagnot treat, or prevent any
β-Pinene	ND	-	medical-related symptoms.
α-Terpinene	ND	-	The statements and results herein have
γ-Terpinene	ND	-	not been approved and/or endorsed by the FDA.
Terpinolene	ND	-	
TOTAL	0.000	0.000	Reporting Limit (µ

Reporting Limit (µg/g): 8.49

"-" Not detected above LOD.













CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36123)

Testing Location:Customer ID: 2168Order ID: OR10628Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13229275793 Mass: 1ea

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run:** Date Collected: 10/02/2023 License: ADH 113 License: 00065C 20230925STRAW22 Date Received: 10/03/2023

Cultivar (Strain) or Sample Description: ARV-Provisions Strawberries **Date Completed:** 10/04/2023

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

Analysis Date/Time: 10/03/2023 2013 Method: HS/GC/MS Deviations from SOP:

Analyst: KF Instrument: Agilent 7890/5975 None

Solvent	Result (μg/g)	<u>LOD</u> (μg/g)	LOQ (µg/g)	<u>Action</u> <u>Level</u> (μg/g)	Solvent	Result (μg/g)	<u>LOD</u> (µg/g)	LOQ (µg/g)	<u>Action</u> <u>Level</u> (μg/g)
Acetone (67-64-1)	-	150	300	5000	n-Heptane (142-82-5)	-	150	300	5000
Acetonitrile (75-5-8)	-	150	300	410	n-Hexane (110-54-3)	-	52.5	105	290
Benzene (71-43-2)	-	1.50	3.00	2	Isobutane (75-28-5)	-	150	300	5000
n-Butane (106-97-2)	-	150	300	5000	Isopropanol (67-63-0)	-	150	300	5000
1-Butanol (71-36-3)	-	150	300	5000	Isopropyl acetate	_	150	300	5000
2-Butanol (78-92-2)	-	150	300	5000	(108-21-4)		150	500	5000
2-Butanone (78-93-3)	-	150	300	5000	Isopropyl benzene (98-82-8)	-	15.0	30.0	70
Cyclohexane (110-82-7)	-	150	300	3880	Methanol (67-56-1)	_	150	300	3000
1,2-Dimethoxyethane (110-71-4)	-	15.0	30.0	100	2-Methylbutane (78-78-4)	-	150	300	5000
N,N-Dimethylacetamide (127-19-5)	-	150	300	1090	Methylene chloride (75-9-2)	-	150	300	600
2,2-Dimethylbutane (75-83-2)	-	52.5	105	290	2-Methylpentane (107-83-5)	-	52.5	105	290
2,3-Dimethylbutane (79-29-8)	-	52.5	105	290	3-Methylpentane (96-10-0) n-Pentane (109-66-0)	-	52.5 150	105 300	290 5000
N,N-Dimethylformamide (68-12-2)	-	150	300	880	1-Pentanol (71-41-0) n-Propane (74-98-6)	-	150 150	300 300	5000 5000
Dimethylsulfoxide (67-68-5)	-	150	300	5000	1-Propanol (71-23-8)		150	300	5000
1,4-Dioxane (123-91-1)	_	150	300	380	Pyridine (110-86-1)	-	52.5	105	200
Ethanol (64-17-5)	4070	150	300	5000	Tetrahydrofuran (109-99-9)	-	150	300	720
2-Ethoxyethanol (110-80-5)	-	52.5	105	160	Tetramethylene sulfone (126-33-0)	-	52.5	105	160
Ethyl ether (60-29-7)	-	150	300	5000	Toluene (108-88-3)	_	150	300	890
Ethyl acetate (141-78-6)	-	150	300	5000	o-Xylene (95-47-6)		150	300	2170
Ethyl benzene (100-41-4)	-	150	300	2170	m,p-Xylene (108-38-3 or				
Ethylene glycol (107-21-1)		150	300	620	106-42-3)		150	300	2170
Ethylene oxide (75-21-8)	-	15.0	30.0	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170

Solvent	Synonym(s)	Solvent	Synonym(s)
Solvent	<u>Synonym(s)</u>		<u>Synonym(s)</u>
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



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.

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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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36123)

Testing Location:Customer ID: 2168Order ID: OR10628Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13229275793 Mass: 1ea
Greenbrier, AR 72058 Fort Smith, AR 72903 Production Run: Date Collected: 10/02/2023

License: ADH 113 License: 00065C 20230925STRAW22 **Date Received:** 10/03/2023 **Cultivar (Strain) or Sample Description:** ARV-Provisions Strawberries **Date Completed:** 10/04/2023

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 10/03/2023 1525 Method: LC/MS/MS Deviations from SOP:

Analyst: KF Instrument: Shimadzu LC-8050 None

<u>Pesticide</u>	<u>Result</u> (μg/g)	LOD (µg/g)	<u>LOQ</u> (µg/g)	Action Level (µg/g)	<u>Pesticide</u>	<u>Result</u> (μg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> <u>Level</u> (μg/g)
Abamectin (71751-41-2)	-	0.0101	0.0810	0.5	Kresoxim-methyl	_	0.0101	0.0810	0.4
Acephate (30560-19-1)	-	0.0101	0.0810	0.4	(143390-89-0)				
Acequinocyl (57960-19-7)	-	0.0101	0.0810	2	Malathion (121-75-5)	-	0.0101	0.0810	0.2
Acetamiprid (135410-20-7)	-	0.0101	0.0810	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.0101 0.0101	0.0810 0.0810	0.2 0.2
Aldicarb (116-06-3)	-	0.0101	0.0810	0.4	Methomyl (16752-77-5)	-	0.0101	0.0810	0.4
Azoxystrobin (131860-33-8)	-	0.0101	0.0810	0.2	Methyl parathion (298-0-0)	-	0.0101	0.0810	0.2
Bifenazate (149877-41-8)	DET	0.0101	0.0810	0.2	MGK 264 (113-48-4)	-	0.0101	0.0810	0.2
Bifenthrin (82657-04-3)	-	0.0101	0.0810	0.2	Myclobutanil		0.0101	0.0810	0.2
Boscalid (188425-85-6)	DET	0.0101	0.0810	0.4	(88671-89-0)		0.0101	0.0010	0.2
Carbaryl (63-25-2)	-	0.0101	0.0810	0.2	Naled (300-76-5)	-	0.0101	0.0810	0.5
Carbofuran (1563-66-2)	-	0.0101	0.0810	0.2	Oxamyl (23135-22-0)	-	0.0101	0.0810	1
Chlorantraniliprole (800008-45-7)	-	0.0101	0.0810	0.2	Paclobutrazol (76738-62-0)	-	0.0101	0.0810	0.4
Chlorfenapyr		0.0101	0.0810	1	Permethrins (52645-53-1)	-	0.0101	0.0810	0.2
(122453-73-0)	-	0.0101	0.0810	1	Phosmet (732-11-6)	-	0.0101	0.0810	0.2
Chlorpyrifos (2921-88-2)	-	0.0101	0.0810	0.2	Piperonyl butoxide	_	0.0101	0.0810	2
Clofentezine (74115-24-5)	-	0.0101	0.0810	0.2	(51-03-6)		0.0404	0.0010	0.0
Cyfluthrin (68359-37-5)	-	0.0101	0.0810	1	Prallethrins (2331-36-9)	-	0.0101	0.0810	0.2
Cypermethrin (52315-07-8)	-	0.0101	0.0810	1	Propiconazole (60207-90-1))	-	0.0101	0.0810	0.4
Daminozide (1596-84-5)	-	0.0101	0.0810	1	Propoxur (114-26-1)	-	0.0101	0.0810	0.2
DDVP (62-73-7)	-	0.0101	0.0810	0.1	Pyrethrins (8003-34-7)	-	0.0101	0.0810	1
Diazinon (333-41-5)	-	0.0101	0.0810	0.2	Pyridaben (96489-71-3)	-	0.0101	0.0810	0.2
Dimethoate (60-51-5)	-	0.0101	0.0810	0.2	Spinosad (168316-95-8)	-	0.0101	0.0810	0.2
Ethoprophos (13194-48-4)	DET	0.0101	0.0810	0.2	Spiromesifen (283594-90-1)	-	0.0101	0.0810	0.2
Etofenprox (80844-07-1)	-	0.0101	0.0810	0.4	Spirotetramat				
Etoxazole (153233-91-1)	-	0.0101	0.0810	0.2	(203313-25-1)	-	0.0101	0.0810	0.2
Fenoxycarb (72490-01-8)	-	0.0101	0.0810	0.2	Spiroxamine		0.0101	0.0810	0.4
(E)-Fenpyroximate (134098-61-6)	-	0.0101	0.0810	0.4	(118134-30-8) Tebuconazole		0.0101	0.0810	0.4
Fipronil (120068-37-3)	-	0.0101	0.0810	0.4	(80443-41-0)		0.0101	0.0610	0.4
Flonicamid (158062-67-0)	-	0.0101	0.0810	1	Thiacloprid		0.0101	0.0810	0.2
Fludioxinil (131341-86-1)	-	0.0101	0.0810	0.4	(111988-49-9)				
Hexythiazox (78587-05-0)	-	0.0101	0.0810	1	Thiamethoxam (153719-23-4)	-	0.0101	0.0810	0.2
Imazalil (35554-44-0)	-	0.0101	0.0810	0.2	Trifloxystrobin				
Imidacloprid (138261-41-3)	-	0.0101	0.0810	0.4	(141517-21-7)		0.0101	0.0810	0.2



Color Key

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

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

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













Date Received: 10/03/2023

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Testing Location:Customer ID: 2168Order ID: OR10628Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13229275793 Mass: 1ea

Greenbrier, AR 72058 Fort Smith, AR 72903 Production Run: Date Collected: 10/02/2023

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

Date Completed: 10/04/2023

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

20230925STRAW22

Analysis Date/Time: 10/03/2023 1826 (ICP/OES) Method: ICP/MS Deviations from SOP:

Analysis Date/Time: - (DMA) Instrument: Agilent 7500ce None

Analyst: KF

License: ADH 113

Heavy Metal	<u>Result</u> (μg/kg)	<u>LOD</u> (µg/kg)	<u>LOQ</u> (μg/kg)	<u>Action Level</u> (μg/kg)
Arsenic (As)	DET	59.7	94.5	200
Cadmium (Cd)	-	59.7	94.5	200
Lead (Pb)	-	59.7	94.5	500
Mercury (Hg)	_	59.7	94.5	100

License: 00065C



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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Date Completed: 10/04/2023

CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36123)

Testing Location:Customer ID: 2168Sample ID: SA36123Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible232 S. Broadview St.5601 Old Greenwood Rd Suite CM00065C13229275793Mass: 1ea

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 10/02/2023 License: ADA 05_H273 License: 00065C 20230925STRAW22 **Date Received:** 10/03/2023

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 10/04/2023 0838 **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	-

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



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)

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.



