





## SUMMARY OF ANALYSIS (SAMPLE ID: SA41318)

Testing Location:	Customer ID: 2168	<b>Order ID:</b> OR11630	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13229894153	Mass: 1ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 06/05/2025
License: ADH 113	License: 00065C	P20250603CAR05	Date Received: 06/05/2025
Cultivar (Strain) or Sample Des	Date Completed:06/09/2025		

\*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.

\*Where provided, statements of conformity (e.g. Pass/Fail) are made in accordance with ILAC G8, Binary Statement for Simple Acceptance Rule (w=0, AL=TL). PASS: when the result is within the acceptance interval. FAIL: when the result is outside the acceptance interval

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.

<u>Cannabinoi</u>	<u>ds (Top 3)</u>	<u>(%)</u>	<u>mg/g</u>		
Δ9-Τ		0.246	2.46		
CB	G	0.00363	0.0363		
CB	N	0.00283	0.0283		
TOTAL	CBD	0.000	0.000		
TOTAL	THC	0.246	2.46		
TOTAL CANN	NABINOIDS	0.252	2.52		
<u>Terpenes</u>	<u>(Top 5)</u>	<u>(%)</u>	<u>ppm</u>		
α-Hum	ulene	0.0675	675		
trans-Ne	rolidol	0.00289	28.9		
α-Bisa	bolol				
Campl	nene				
δ-3-Ca	rene				
TOTAL TE	RPENES	0.0704	704		
<b>Contaminants</b>	PASS/FAIL	Sample Picture	e Upon Receipt		
Heavy Metals:	PASS				
Microbiology:	PASS				
Pesticides:	PASS				
Residual Solvents:	PASS				



Scan the QR code to verify results.

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The statements and results herein have not been approved and/or endorsed by the FDA.

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

REPORT OF LABORATORY ANALYSIS

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

Analysis Date/Time: 06/06/2025 1505 Analyst: PW Method: HPLC/DAD Instrument: Agilent 1100

<u>Cannabinoid</u>	<u>Result</u> (%)	<u>Result</u> (mg/g)	<u>LOD</u> (mg/g)	LOQ (mg/g)	<u>Result</u> (mg/ mL)	<u>Per</u> Serving (mg)	<u>Per</u> <u>Unit</u> (mg)
CBC	ND	ND	0.00225	0.00524	-	0.000	0.000
CBCA	ND	ND	0.00698	0.0163	-	0.000	0.000
CBD	ND	ND	0.0159	0.0370	-	0.000	0.000
CBDA	ND	ND	0.00584	0.0136	-	0.000	0.000
CBDV	ND	ND	0.00254	0.00593	-	0.000	0.000
CBDVA	ND	ND	0.00678	0.0158	-	0.000	0.000
CBG	0.00363	0.0363	0.0103	0.0240	-	0.148	1.48
CBGA	ND	ND	0.0146	0.0172	-	0.000	0.000
CBL	ND	ND	0.0119	0.0278	-	0.000	0.000
CBN	0.00283	0.0283	0.00546	0.0128	-	0.115	1.15
CBNA	ND	ND	0.00590	0.0137	-	0.000	0.000
Δ9-ΤΗC	0.246	2.46	0.00655	0.0153	-	10.0	100
$\Delta 8$ -THC	ND	ND	0.0102	0.0238	-	0.000	0.000
THCA	ND	ND	0.00355	0.00830	-	0.000	0.000
THCV	ND	ND	0.00852	0.0199	-	0.000	0.000
THCVA	ND	ND	0.00272	0.00633	-	0.000	0.000
TOTAL	0.252	2.52			-	10.3	103
TOTAL CBC	0.000	0.000				0.000	0.000
TOTAL CBD	0.000	0.000			-	0.000	0.000
TOTAL CBDV	0.000	0.000			-	0.000	0.000
TOTAL CBG	0.00363	0.0363			-	0.148	1.48
TOTAL CBN	0.00283	0.0283			-	0.115	1.15
TOTAL THC	0.246	2.46			-	10.0	100
TOTAL THCV	0.000	0.000			-	0.000	0.000

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, ND - Not Detected (less than LOD)

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



SERVING MASS (g): 4.07 SERVINGS/UNIT: 10

"-" Not reported for this sample.

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.

Dry percent = Wet percent / (1-(Moisture Content/ 100))

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) +  $\Delta$ 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.

W. Felling, Ph.D. rv Dir

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F.A.S.T. Personnel perform sampling following the Sampling SOP (SOP-02).







<b>Testing Location:</b>	Customer ID: 2168	<b>Order ID:</b> OR11630	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13229894153	Mass: 1ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 06/05/2025
License: ADH 113	License: 00065C	P20250603CAR05	Date Received: 06/05/2025
Cultivar (Strain) or Sample	Date Completed:06/09/2025		

## A A

<b>TERPENOID PROFILE</b>
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Analysis Date/Time:06/06/2025 1625 Analyst: KF		Method: GC	/MS Agilent 7890/5975	<b>Deviations from SOP:</b> None			
-			Agricia 7050/5575	110116			
<u>Terpene</u>	<u>Result</u> (ppm)	<u>Result</u> (%)					
α-Bisabolol	<loq< td=""><td>-</td><td></td><td>and the second sec</td></loq<>	-		and the second sec			
Camphene	ND	-					
δ-3-Carene	ND	-					
β-Caryophyllene	ND	-		The second			
Caryophyllene oxide	ND	-					
p-Cymene	ND	-					
Eucalyptol	<loq< td=""><td>-</td><td></td><td></td></loq<>	-					
Geraniol	ND	-					
Guaiol	ND	-		Abbreviations: GC - Gas			
α-Humulene	675	0.0675		Chromatography, MS - Mass			
Isopulegol	ND	-		Spectrometry, RL - Reporting Limit Abbreviations: ND - Not Detected, ,			
d-Limonene	ND	-		LOD - Limit of Detection, LOQ - Limit			
Linalool	ND	-		of Quantitation			
β-Myrcene	ND	-		This information is provided as a service and makes no claims of efficacy and/or			
cis-Nerolidol	ND	-		safety of this product.			
trans-Nerolidol	28.9	0.00289		Results are applicable only for the sample(s) analyzed and for the specific			
α-Ocimene	ND	-		analysis conducted.			
β-Ocimene	ND	-		This report is for informational purposes only and should not be used to diagnose,			
α-Pinene	ND	-		treat, or prevent any			
β-Pinene	ND	-		medical-related symptoms.			
α-Terpinene	ND	-		The statements and results herein have not been approved and/or endorsed by			
γ-Terpinene	ND	-		the FDA.			
Terpinolene	ND	-					
TOTAL	704	0.0704		Reporting Limit (ppm): 4.93			

"-" Not detected above LOD.











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License: ADH 113	License: 00065C	P20250603CAR05	Date Received: 06/05/2025
Cultivar (Strain) or Sample	Date Completed:06/09/2025		

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

Analysis Date/Time: 06/06/2025 0430		Method: HS/GC/MS				<b>Deviations from SOP:</b>				
Analyst: KF				I	nstrument: Agilent 78	90/5975	5 None			
<u>Solvent</u>	<u>Result</u> (ppm)	<u>LOD</u> (ppm)	<u>LOQ</u> (ppm)	<u>Action</u> Level (ppm)	<u>Solvent</u>	<u>Result</u> (ppm)	<u>LOD</u> (ppm)	<u>LOQ</u> (ppm)	<u>Action</u> <u>Level</u> (ppm)	
Acetone (67-64-1)	-	38.1	76.2	5000	n-Heptane (142-82-5)	-	38.1	76.2	5000	S YEAR /-
Acetonitrile (75-5-8)	-	38.1	76.2	410	n-Hexane (110-54-3)	-	13.3	26.7	290	
Benzene (71-43-2)	-	0.381	0.762	2	Isobutane (75-28-5)	-	38.1	76.2	5000	
n-Butane (106-97-2)	-	38.1	76.2	5000	Isopropanol (67-63-0)	-	38.1	76.2	5000	
1-Butanol (71-36-3)	-	38.1	76.2	5000	Isopropyl acetate		38.1	76.2	5000	
2-Butanol (78-92-2)	-	38.1	76.2	5000	(108-21-4)		50.1	70.2	5000	
2-Butanone (78-93-3)	-	38.1	76.2	5000	Isopropyl benzene	-	3.81	7.62	70	
Cyclohexane (110-82-7)	-	38.1	76.2	3880	(98-82-8) Methanol (67-56-1)		38.1	76.2	3000	Color Key
1,2-Dimethoxyethane (110-71-4)	-	3.81	7.62	100	2-Methylbutane (78-78-4)	-	38.1	76.2	5000	RESULT < AL
N,N-Dimethylacetamide (127-19-5)	-	38.1	76.2	1090	Methylene chloride (75-9-2)	-	38.1	76.2	600	<b>RESULT</b> > AL
2,2-Dimethylbutane (75-83-2)	-	13.3	26.7	290	2-Methylpentane (107-83-5)	-	13.3	26.7	290	"-" not detected above 1/2 Action Level
2,3-Dimethylbutane					3-Methylpentane (96-10-0)	-	13.3	26.7	290	
(79-29-8)	-	13.3	26.7	290	n-Pentane (109-66-0)	-	38.1	76.2	5000	"*" - o,m,p-Xylene and
N,N-Dimethylformamide	_	38.1	76.2	880	1-Pentanol (71-41-0)	-	38.1	76.2	5000	Ethylbenzene
(68-12-2)		50.1	70.2	000	n-Propane (74-98-6)	-	38.1	76.2	5000	Action levels are
Dimethylsulfoxide (67-68-5)	-	38.1	76.2	5000	1-Propanol (71-23-8)	-	38.1	76.2	5000	referenced from the State of
(07-00-5) 1,4-Dioxane (123-91-1)		38.1	76.2	380	Pyridine (110-86-1)	-	13.3	26.7	200	Arkansas
Ethanol (64-17-5)	_	38.1	76.2	5000	Tetrahydrofuran (109-99-9)	-	38.1	76.2	720	MMJ testing guidelines.
2-Ethoxyethanol (110-80-5)	_	13.3	26.7	160	Tetramethylene sulfone	_	13.3	26.7	160	0
Ethyl ether (60-29-7)	_	38.1	76.2	5000	(126-33-0)		20.1	76.2	000	A value of "-"
Ethyl acetate (141-78-6)	_	38.1	76.2	5000	Toluene (108-88-3)	-	38.1	76.2	890	for the action level means that analyte
Ethyl benzene (100-41-4)	_	38.1	76.2	2170	o-Xylene (95-47-6)	-	38.1	76.2	2170	is not currently
Ethylene glycol (107-21-1)	_	38.1	76.2	620	m,p-Xylene (108-38-3 or 106-42-3)	-	38.1	76.2	2170	regulated by the
Ethylene oxide (75-21-8)	-	3.81	7.62	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170	regulations referenced above.
<u>Solvent</u>		Synonym(s	<u>s)</u>		<u>Solvent</u>	<u>S</u>	ynonym(s)			
Acetonitrile		Methyl Cya	anide, ACN		Ethylene glycol	1,	,2-Ethanediol			
1-Butanol		n-Butanol,	Butyl Alco	hol	Isobutane	2.	-Methylpropa	ne		
2-Butanol		sec-Butyl a	lcohol		Isopropanol	2.	-Propanol, IP/	4		
2-Butanone		Methyl eth	yl ketone, N	1EK	Isopropyl Acetate	А	cetic acid iso	propyl este	r	
1,2-Dimethoxyethane		Monoglym	e		Methanol	Ν	fethyl alcoho	l		
2,3-Dimethylbutane		Neohexane			2-Methylbutane	Is	opentane			
2,3-Dimethylbutane		Diisopropy	71		Methylene chloride	D	ichlorometha	ne		
N,N-Dimethylformamide		DMF			2-Methylpentane	Is	ohexane			
Dimethysufoxide		DMSO			1-Pentanol	n·	-Amyl alcoho	1		

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

1-Propanol

Xylene

Tetrahydrofuran

Tetramethylene sulfone

Cellosolve, Ethyl glycol

Diethyl ether, Ether

EtOAc

Phenylethane

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2-Ethoxyethanol

Ethyl ether

Ethyl acetate

Ethyl benzene

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Propyl alcohol

Dimethylbenzene

THF

Sulfolane

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**Deviations from SOP:** 

None



# **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA41318)**

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Cultivar (Strain) or Sample	Date Completed:06/09/2025		

### **PESTICIDES PROFILE (SOP: SOP-PEST-001)**

Analysis Date/Time: 06/06/2025 1521 Analyst: KF

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

Instrument: Shimadzu LC-8050

<u>Pesticide</u>	<u>Result</u> (ppm)	<u>LOD</u> (ppm)	<u>LOQ</u> (ppm)	<u>Action</u> Level (ppm)	<u>Pesticide</u>	<u>Result</u> (ppm)	<u>LOD</u> (ppm)	<u>LOQ</u> (ppm)	<u>Action</u> Level (ppm)	
Abamectin (71751-41-2)	-	0.00989	0.0791	0.5	Kresoxim-methyl	_	0.00989	0.0791	0.4	
Acephate (30560-19-1)	-	0.00989	0.0791	0.4	(143390-89-0)					
Acequinocyl (57960-19-7)		0.00989	0.0791	2	Malathion (121-75-5)	-	0.00989	0.0791	0.2	
Acetamiprid (135410-20-7)	-	0.00989	0.0791	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.00989 0.00989	0.0791 0.0791	0.2 0.2	<u>Color Key</u>
Aldicarb (116-06-3)	-	0.00989	0.0791	0.4	Methomyl (16752-77-5)	-	0.00989	0.0791	0.4	<b>RESULT &lt; AL</b>
Azoxystrobin (131860-33-8)	-	0.00989	0.0791	0.2	Methyl parathion (298-0-0)	-	0.00989	0.0791	0.2	<b>RESULT &gt; AL</b>
Bifenazate (149877-41-8)	-	0.00989	0.0791	0.2	MGK 264 (113-48-4)	-	0.00989	0.0791	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3)	-	0.00989	0.0791	0.2	Myclobutanil	_	0.00989	0.0791	0.2	"-" not detected above
Boscalid (188425-85-6)	-	0.00989	0.0791	0.4	(88671-89-0)					LOD
Carbaryl (63-25-2)	-	0.00989	0.0791	0.2	Naled (300-76-5)	-	0.00989	0.0791	0.5	Permethrins measured as the
Carbofuran (1563-66-2)	-	0.00989	0.0791	0.2	Oxamyl (23135-22-0)	-	0.00989	0.0791	1	cumulative residue of the <i>cis</i> - and
Chlorantraniliprole (800008-45-7)	-	0.00989	0.0791	0.2	Paclobutrazol (76738-62-0)	-	0.00989	0.0791	0.4	<i>trans-</i> permethrin isomers.
Chlorfenapyr (122453-73-0)	-	0.00989	0.0791	1	Permethrins (52645-53-1)	-	0.00989	0.0791	0.2	Pyrethrins measured as the cumulative residue of the pyrethrin I, cinerin I, and jasmolin
Chlorpyrifos (2921-88-2)	-	0.00989	0.0791	0.2	Phosmet (732-11-6)	-	0.00989	0.0791	0.2	I isomers.
Clofentezine (74115-24-5)	-	0.00989	0.0791	0.2	Piperonyl butoxide (51-03-6)	_	0.00989	0.0791	2	Action levels are
Cyfluthrin (68359-37-5)	-	0.00989	0.0791	1	(51-03-6) Prallethrins (2331-36-9)		0.00989	0.0791	0.2	referenced from the State of
Cypermethrin (52315-07-8)	-	0.00989	0.0791	1	Propiconazole (60207-90-1))	-	0.00989	0.0791	0.2	Arkansas MMJ testing
Daminozide (1596-84-5)	-	0.00989	0.0791	1	(00207-50-1)) Propoxur (114-26-1)		0.00989	0.0791	0.2	guidelines.
DDVP (62-73-7)	-	0.00989	0.0791	0.1	Pyrethrins (8003-34-7)		0.00989	0.0791	1	A value of "-" for the action level
Diazinon (333-41-5)	-	0.00989	0.0791	0.2	Pyridaben (96489-71-3)		0.00989	0.0791	0.2	means that analyte is not currently regulated by the
Dimethoate (60-51-5)	-	0.00989	0.0791	0.2	Spinosad (168316-95-8)	_	0.00989	0.0791	0.2	regulations referenced above.
Ethoprophos (13194-48-4)	-	0.00989	0.0791	0.2	Spiromesifen (283594-90-1)	-	0.00989	0.0791	0.2	Disclaimer: This information is
Etofenprox (80844-07-1)	-	0.00989	0.0791	0.4	Spirotetramat					provided as a service and makes no claims of efficacy and/or safety
Etoxazole (153233-91-1)	-	0.00989	0.0791	0.2	(203313-25-1)	-	0.00989	0.0791	0.2	of this product. Results are
Fenoxycarb (72490-01-8)	-	0.00989	0.0791	0.2	Spiroxamine		0.00989	0.0791	0.4	applicable only for the sample(s) analyzed and for the specific
(E)-Fenpyroximate (134098-61-6)	-	0.00989	0.0791	0.4	(118134-30-8) Tebuconazole	-	0.00989	0.0791	0.4	analysis conducted. This report is for informational purposes only
Fipronil (120068-37-3)	-	0.00989	0.0791	0.4	(80443-41-0)	-	0.00909	0.0791	0.4	and should not be used to
Flonicamid (158062-67-0)	-	0.00989	0.0791	1	Thiacloprid	_	0.00989	0.0791	0.2	diagnose, treat, or prevent any medical-related symptoms. The
Fludioxinil (131341-86-1)	-	0.00989	0.0791	0.4	(111988-49-9)					statements and results herein have
Hexythiazox (78587-05-0)	-	0.00989	0.0791	1	Thiamethoxam (153719-23-4)	-	0.00989	0.0791	0.2	not been approved and/or endorsed by the FDA.
Imazalil (35554-44-0)	-	0.00989	0.0791	0.2	Trifloxystrobin		0.00989	0.0701	0.2	endorsed by the FDA.
Imidacloprid (138261-41-3)	-	0.00989	0.0791	0.4	(141517-21-7)	-	0.00989	0.0791	0.2	

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> Cyfluthrin DDVP Ethoprophos

<u>Pesticide</u> Myclobutanil Naled Phosmet

Synonym(s)

Baythroid

Dichlorvos

Prophos

<u>Synonym(s)</u> Systhane Dibrom Imidan

<u>Pesticide</u> Propiconazole Propoxur **Synonym(s)** Tilt Baygon



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F.A.S.T. Personnel perform sampling following the Sampling SOP (SOP-02).







Testing Location:	Customer ID: 2168	Order ID: OR11630	Sample Type: Primary
Arkansas	<b>River Valley Relief MIPS</b>	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13229894153	Mass: 1ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 06/05/2025
License: ADH 113	License: 00065C	P20250603CAR05	Date Received: 06/05/2025
Cultivar (Strain) or Sample Description: ARV-Sea Salt Caramel Soft Chews 200mg 20pk			Date Completed:06/09/2025

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

Analysis Date/Tir	<b>ne:</b> 06/06/2025 1	1558	Method: I		<b>Deviations from SOP:</b>
Analyst: KF			Instrume	<b>ıt:</b> Agilent 7700x	None
<u>Heavy Metal</u>	<u>Result</u> (ppb)	<u>LOD</u> (ppb)	<u>LOQ</u> (ppb)	<u>Action Level</u> (ppb)	
Arsenic (As)	ND	58.3	92.2	200	
Cadmium (Cd)	ND	58.3	92.2	200	
Lead (Pb)	ND	58.3	92.2	500	
Mercury (Hg)	ND	58.3	92.2	100	

Abbreviations: ICP - Inductively-Coupled Plasma, MS - Mass Spectrometry,

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 "ND" not detected above LOD Reporting limit is 1/2 the action level of each analyte.

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

*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.

W. Felling, Ph.D rv Dir

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Testing Location:	Customer ID: 2168	Sample ID: SA41318	Sample Type: Primary
Arkansas	<b>River Valley Relief MIPS</b>	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13229894153	Mass: 1ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 06/05/2025
-	License: 00065C	P20250603CAR05	Date Received: 06/05/2025
Cultivar (Strain) or Sample	Date Completed:06/09/2025		

#### MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 06/06/20 Analyst: PW		Hardy Diagnostics Compace ent: Thermo Incubator	tDry <b>Deviations from SOP:</b> None
Bacteria/Microbe	<u>Result</u> (CFU/g)	<u>Action Level</u> (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





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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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