



SUMMARY OF ANALYSIS (SAMPLE ID: SA38965)

Testing Location:	Customer ID: 2168	Order ID: OR11215	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13220003538	Mass: 1ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 09/26/2024
License: ADH 113	License: 00065C	P20240925GHEE10	Date Received: 09/27/2024
Cultivar (Strain) or Sample Do	Date Completed: 09/29/2024		

*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.

Cannabino	ids (Top 3)	<u>(%)</u>	<u>mg/g</u>		
Δ9-1	ΉС	0.661	6.61		
Δ8-1	ΉС	0.0291	0.291		
CE	G	0.0268	0.268		
TOTAI	CBD	0.0257	0.257		
TOTAI	L THC	0.661	6.61		
TOTAL CAN	NABINOIDS	0.760	7.60		
Terpenes	<u>s (Top 5)</u>	<u>(%)</u>	µg∕g		
d-Lim	onene	0.000889	8.89		
β-Caryop	hyllene	0.000719	7.19		
β-Муг	cene	0.000595	5.95		
α-Bisa	bolol	0.000573	5.73		
Camp	hene				
TOTAL TH	ERPENES	0.00278	27.8		
Contaminants	PASS/FAIL	Sample Picture Upon Receipt			
Heavy Metals:	PASS				

Heavy Metals:	PASS
Microbiology:	PASS
Pesticides:	PASS
Residual Solvents:	PASS





Scan the QR code to verify results.

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Kyle W. Felling, Ph.D. Laboratory Director

REPORT OF LABORATORY ANALYSIS

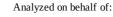
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Cultivar (Strain) or Sample	Date Completed: 09/29/2024		

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

Analysis Date/Time: 09/27/2024 1458 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> Unit (mg)	
CBC	ND	ND	0.00215	0.00501	-	-	-	
CBCA	ND	ND	0.00666	0.0155	-	-	-	
CBD	0.0257	0.257	0.0151	0.0353	-	0.386	23.1	
CBDA	ND	ND	0.00557	0.0130	-	-	-	
CBDV	ND	ND	0.00243	0.00566	-	-	-	
CBDVA	ND	ND	0.00647	0.0151	-	-	-	
CBG	0.0268	0.268	0.00982	0.0229	-	0.402	24.1	
CBGA	ND	ND	0.0139	0.0164	-	-	-	
CBL	ND	ND	0.0113	0.0265	-	-	-	
CBN	0.0176	0.176	0.00521	0.0122	-	0.264	15.8	
CBNA	ND	ND	0.00563	0.0131	-	-	-	
Δ9-ΤΗC	0.661	6.61	0.00625	0.0146	-	9.92	595	
Δ8-THC	0.0291	0.291	0.00975	0.0228	-	0.437	26.2	
THCA	ND	ND	0.00339	0.00792	-	-	-	
THCV	ND	ND	0.00813	0.0190	-	-	-	
THCVA	ND	ND	0.00260	0.00604	-	-	-	
TOTAL	0.760	7.60	·	,		11.4	684	
TOTAL CBC	-	-		,			_	
TOTAL CBD	0.0257	0.257			-	0.386	23.1	
TOTAL CBDV	-	-			-	-	-	
TOTAL CBG	0.0268	0.268			-	0.402	24.1	
TOTAL CBN	0.0176	0.176			-	0.264	15.8	
TOTAL THC	0.661	6.61			-	9.92	595	
TOTAL THCV	-	-			-	-	-	

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): -



SERVING MASS (g): 1.50 SERVINGS/UNIT: 60

"-" 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) + Δ 9-THC Total THCV = (THCVA x 0.867) + THCV

Percentage results are reported by mass.

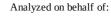
mg/g results are reported as mass component per mass material.

Kyle W. Felling, Ph.D. ory Dire











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Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13220003538	Mass: 1ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 09/26/2024
License: ADH 113	License: 00065C	P20240925GHEE10	Date Received: 09/27/2024
Cultivar (Strain) or Sample D	Date Completed: 09/29/2024		

Analysis Date/Time:09/27/2024 2112 Analyst: KF

Analyst: KF		Instrument: Agilent 7890/5975			
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)			
α-Bisabolol	5.73	0.000573			
Camphene	ND	-			
δ-3-Carene	ND	-			
β-Caryophyllene	7.19	0.000719			
Caryophyllene oxide	<loq< td=""><td>-</td><td></td></loq<>	-			
p-Cymene	ND	-			
Eucalyptol	ND	-			
Geraniol	ND	-			
Guaiol	ND	-			
α-Humulene	<loq< td=""><td>-</td><td></td></loq<>	-			
Isopulegol	ND	-			
d-Limonene	8.89	0.000889			
Linalool	<loq< td=""><td>-</td><td></td></loq<>	-			
β-Myrcene	5.95	0.000595			
cis-Nerolidol	ND	-			
trans-Nerolidol	ND	-			
α-Ocimene	ND	-			
β-Ocimene	<loq< td=""><td>-</td><td></td></loq<>	-			
α-Pinene	ND	-			
β-Pinene	ND	-			
α-Terpinene	ND	-			
γ-Terpinene	ND	-			
Terpinolene	ND	-			
TOTAL	27.8	0.00278			

TERPENOID PROFILE

Method: GC/MS



Deviations from SOP:

None

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): 4.70

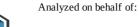
"-" Not detected above LOD.













Testing Location:	Customer ID: 2168	Order ID: OR11215	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13220003538	Mass: 1ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 09/26/2024
License: ADH 113	License: 00065C	P20240925GHEE10	Date Received: 09/27/2024
Cultivar (Strain) or Sample	Date Completed: 09/29/2024		

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

Analysis Date/Time:	09/27/20	24 2211		Method: HS/GC/MS				D	eviations	s from SOP:
Analyst: KF					Instrument: Agilent 7890/5975			Ν	one	
<u>Solvent</u>	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	<u>Solvent</u>	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	
Acetone (67-64-1)	-	39.6	79.2	5000	n-Heptane (142-82-5)	-	39.6	79.2	5000	
Acetonitrile (75-5-8)	-	39.6	79.2	410	n-Hexane (110-54-3)	-	13.9	27.7	290	
Benzene (71-43-2)	-	0.396	0.792	2	Isobutane (75-28-5)	-	39.6	79.2	5000	
n-Butane (106-97-2)	-	39.6	79.2	5000	Isopropanol (67-63-0)	-	39.6	79.2	5000	
1-Butanol (71-36-3)	-	39.6	79.2	5000	Isopropyl acetate	_	39.6	79.2	5000	
2-Butanol (78-92-2)	-	39.6	79.2	5000	(108-21-4)		55.0	/ 5.2	5000	
2-Butanone (78-93-3)	-	39.6	79.2	5000	Isopropyl benzene (98-82-8)	-	3.96	7.92	70	
Cyclohexane (110-82-7)	-	39.6	79.2	3880	(98-82-8) Methanol (67-56-1)		39.6	79.2	3000	Color Key
1,2-Dimethoxyethane (110-71-4)	-	3.96	7.92	100	2-Methylbutane (78-78-4)	-	39.6	79.2	5000	RESULT < AL
N,N-Dimethylacetamide (127-19-5)	-	39.6	79.2	1090	Methylene chloride (75-9-2)	-	39.6	79.2	600	RESULT > AL
2,2-Dimethylbutane (75-83-2)	-	13.9	27.7	290	2-Methylpentane (107-83-5)	-	13.9	27.7	290	"-" not detected above 1/2 Action Level
2,3-Dimethylbutane					3-Methylpentane (96-10-0)	-	13.9	27.7	290	
(79-29-8)	-	13.9	27.7	290	n-Pentane (109-66-0)	-	39.6	79.2	5000	"*" - o,m,p-Xylene and
N,N-Dimethylformamide (68-12-2)	-	39.6	79.2	880	1-Pentanol (71-41-0)	-	39.6	79.2	5000	Ethylbenzene
Dimethylsulfoxide					n-Propane (74-98-6)	-	39.6	79.2	5000	Action levels are referenced from the State of
(67-68-5)	-	39.6	79.2	5000	1-Propanol (71-23-8)	-	39.6	79.2 27.7	5000	Arkansas
1,4-Dioxane (123-91-1)	-	39.6	79.2	380	Pyridine (110-86-1)	-	13.9		200	MMJ testing
Ethanol (64-17-5)	-	39.6	79.2	5000	Tetrahydrofuran (109-99-9)	-	39.6	79.2	720	guidelines.
2-Ethoxyethanol (110-80-5)	-	13.9	27.7	160	Tetramethylene sulfone (126-33-0)	-	13.9	27.7	160	A value of "-"
Ethyl ether (60-29-7)	-	39.6	79.2	5000	Toluene (108-88-3)	-	39.6	79.2	890	for the action level
Ethyl acetate (141-78-6)	-	39.6	79.2	5000	o-Xylene (95-47-6)	-	39.6	79.2	2170	means that analyte
Ethyl benzene (100-41-4)	-	39.6	79.2	2170	m,p-Xylene (108-38-3 or					is not currently
Ethylene glycol (107-21-1)	-	39.6	79.2	620	106-42-3)	-	39.6	79.2	2170	regulated by the
Ethylene oxide (75-21-8)	-	3.96	7.92	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170	regulations referenced above
Solvent		Synonym(s	<u>5)</u>		Solvent	S	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/	ł		
2-Butanone		Methyl eth	yl ketone, N	⁄IEK	Isopropyl Acetate	А	cetic acid iso	propyl ester		
1,2-Dimethoxyethane		Monoglym	e		Methanol	Μ	lethyl alcoho	l		
2,3-Dimethylbutane		Neohexane			2-Methylbutane	Is	opentane			
2,3-Dimethylbutane		Diisopropy	7l		Methylene chloride	D	ichlorometha	ne		
N,N-Dimethylformamide		DMF			2-Methylpentane	Is	ohexane			
Dimethysufoxide		DMSO			1-Pentanol	n·	-Amyl alcoho	1		
2-Ethoxyethanol		Cellosolve	, Ethyl glyc	ol	1-Propanol	Pi	ropyl alcohol			
Ethyl ether		Diethyl eth	ier, Ether		Tetrahydrofuran	Т	HF			

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

Sulfolane

Dimethylbenzene

Tetramethylene sulfone

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Xylene



EtOAc

Phenylethane

Ethyl acetate

Ethyl benzene

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Deviations from SOP:

None



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA38965)

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License: ADH 113	License: 00065C	P20240925GHEE10	Date Received: 09/27/2024
Cultivar (Strain) or Sample	Date Completed: 09/29/2024		

Analysis Date/Time: 09/27/2024 1446 Analyst: KF

PESTICIDES PROFILE (SOP: SOP-PEST-001)

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

<u>Pesticide</u>	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> <u>Level</u>	<u>Pesticide</u>	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> <u>Level</u>	
Abamectin (71751-41-2)		0.00934	0.0748	(µg/g) 0.5	Kresoxim-methyl				<u>(µg/g)</u>	
Acephate (30560-19-1)	_	0.00934	0.0748	0.4	(143390-89-0)	-	0.00934	0.0748	0.4	
Acequinocyl (57960-19-7)	_	0.00934	0.0748	2	Malathion (121-75-5)	-	0.00934	0.0748	0.2	
Acetamiprid					Metalaxyl (57837-19-1)	-	0.00934	0.0748	0.2	
(135410-20-7)	-	0.00934	0.0748	0.2	Methiocarb (2032-65-7)	-	0.00934	0.0748	0.2	<u>Color Key</u>
Aldicarb (116-06-3)	-	0.00934	0.0748	0.4	Methomyl (16752-77-5)	-	0.00934	0.0748	0.4	RESULT < AL
Azoxystrobin (131860-33-8)	-	0.00934	0.0748	0.2	Methyl parathion (298-0-0)	-	0.00934	0.0748	0.2	RESULT > AL
Bifenazate (149877-41-8)	-	0.00934	0.0748	0.2	MGK 264 (113-48-4)	-	0.00934	0.0748	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3)	-	0.00934	0.0748	0.2	Myclobutanil	_	0.00934	0.0748	0.2	"-" not detected above
Boscalid (188425-85-6)	-	0.00934	0.0748	0.4	(88671-89-0)					LOD
Carbaryl (63-25-2)	-	0.00934	0.0748	0.2	Naled (300-76-5)	-	0.00934	0.0748	0.5	Permethrins measured as the
Carbofuran (1563-66-2)	-	0.00934	0.0748	0.2	Oxamyl (23135-22-0)	-	0.00934	0.0748	1	cumulative residue of the <i>cis</i> - and
Chlorantraniliprole (800008-45-7)	-	0.00934	0.0748	0.2	Paclobutrazol (76738-62-0)	-	0.00934	0.0748	0.4	trans- permethrin isomers.
Chlorfenapyr (122453-73-0)	-	0.00934	0.0748	1	Permethrins (52645-53-1)	-	0.00934	0.0748	0.2	Pyrethrins measured as the cumulative residue of the
Chlorpyrifos (2921-88-2)	-	0.00934	0.0748	0.2	Phosmet (732-11-6)	-	0.00934	0.0748	0.2	pyrethrin I, cinerin I, and jasmolin I isomers.
Clofentezine (74115-24-5)	-	0.00934	0.0748	0.2	Piperonyl butoxide	_	0.00934	0.0748	2	
Cyfluthrin (68359-37-5)	-	0.00934	0.0748	1	(51-03-6) Prallethrins (2331-36-9)		0.00934	0.0748	0.2	Action levels are referenced from the State of
Cypermethrin (52315-07-8)	-	0.00934	0.0748	1	Propiconazole (60207-90-1))	-	0.00934	0.0748	0.2	Arkansas MMJ testing
Daminozide (1596-84-5)	-	0.00934	0.0748	1	< <i>//</i>		0.00934	0.0748	0.2	guidelines.
DDVP (62-73-7)	-	0.00934	0.0748	0.1	Propoxur (114-26-1) Pyrethrins (8003-34-7)	-	0.00934	0.0748	0.2	A value of "-" for the action level
Diazinon (333-41-5)	-	0.00934	0.0748	0.2	Pyridaben (96489-71-3)	-	0.00934	0.0748	0.2	means that analyte is not
Dimethoate (60-51-5)	-	0.00934	0.0748	0.2	Spinosad (168316-95-8)	-	0.00934	0.0748	0.2	currently regulated by the regulations referenced above.
Ethoprophos (13194-48-4)	-	0.00934	0.0748	0.2	Spiromesifen	-	0.00934	0.0748	0.2	Disclaimer: This information is
Etofenprox (80844-07-1)	-	0.00934	0.0748	0.4	(283594-90-1)					provided as a service and makes no claims of efficacy and/or safety
Etoxazole (153233-91-1)	-	0.00934	0.0748	0.2	Spirotetramat (203313-25-1)	-	0.00934	0.0748	0.2	of this product. Results are
Fenoxycarb (72490-01-8)	-	0.00934	0.0748	0.2	Spiroxamine		0.0002.4	0.0740	0.4	applicable only for the sample(s)
(E)-Fenpyroximate (134098-61-6)	-	0.00934	0.0748	0.4	(118134-30-8) Tebuconazole		0.00934	0.0748	0.4	analyzed and for the specific analysis conducted. This report is for informational purposes only
Fipronil (120068-37-3)	-	0.00934	0.0748	0.4	(80443-41-0)	-	0.00934	0.0748	0.4	and should not be used to
Flonicamid (158062-67-0)	-	0.00934	0.0748	1	Thiacloprid	_	0.00934	0.0748	0.2	diagnose, treat, or prevent any medical-related symptoms. The
Fludioxinil (131341-86-1)	-	0.00934	0.0748	0.4	(111988-49-9)					statements and results herein have
Hexythiazox (78587-05-0)	-	0.00934	0.0748	1	Thiamethoxam (153719-23-4)	-	0.00934	0.0748	0.2	not been approved and/or
Imazalil (35554-44-0)	-	0.00934	0.0748	0.2	Trifloxystrobin		0.0002.5	0.0740	0.0	endorsed by the FDA.
Imidacloprid (138261-41-3)	-	0.00934	0.0748	0.4	(141517-21-7)	-	0.00934	0.0748	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>Synonym(s)</u> Baythroid Dichlorvos Prophos <u>Pesticide</u> Myclobutanil Naled Phosmet **Synonym(s)** Systhane Dibrom Imidan

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



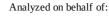
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Testing Location:	Customer ID: 2168	Order ID: OR11215	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13220003538	Mass: 1ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 09/26/2024
License: ADH 113	License: 00065C	P20240925GHEE10	Date Received: 09/27/2024
Cultivar (Strain) or Sample Description: ARV-Infused Ghee Butter 600mg Jar			Date Completed: 09/29/2024

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

Analysis Date/Time: 09/27/2024 1950		Method: ICP/MS Instrument: Agilent 7500ce		Deviations from SOP: None	
Analyst: KF				C	
<u>Heavy Metal</u>	<u>Result</u> (µg/kg)	<u>LOD</u> (µg/kg)	<u>LOQ</u> (µg/kg)	<u>Action Level</u> (µg/kg)	
Arsenic (As)	-	57.7	91.3	200	
Cadmium (Cd)	-	57.7	91.3	200	
Lead (Pb)	-	57.7	91.3	500	
Mercury (Hg)	-	57.7	91.3	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

<u>Color Key</u>



"DET" detected less than LOQ

"-" not detected above LOD

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

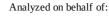
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Testing Location:	Customer ID: 2168	Sample ID: SA38965	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13220003538	Mass: 1ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 09/26/2024
License: ADA 05_H273	License: 00065C	P20240925GHEE10	Date Received: 09/27/2024
Cultivar (Strain) or Sample Description: ARV-Infused Ghee Butter 600mg Jar			Date Completed: 09/29/2024

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 09/29/20 Analyst: PW		Hardy Diagnostics Compac nt: Thermo Incubator	ctDry Deviations from SOP: 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	-	A CONTRACTOR
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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