







## SUMMARY OF ANALYSIS (SAMPLE ID: SA35049)

Testing Location:	Customer ID: 2168	Order ID: OR10416	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13246623704	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	<b>Date Collected:</b> 04/24/2023
License: ADH 113	License: 00065C	E20230417ICMLS01	Date Received: 04/24/2023
Cultivar (Strain) or Sample De	Date Completed: 04/27/2023		

\*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.

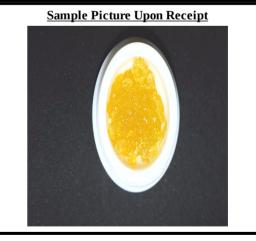
Moisture Content (%)	Water Activity (aw)	PASS/FAIL	
Not Tested	Not Tested	PASS	

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.

Cannabinoids (Top 3)	<u>(%)</u>	mala
		mg/g
THCA	89.6	896
Δ9-ΤΗC	4.77	47.7
CBGA	1.18	11.8
TOTAL CBD	-	-
TOTAL THC	83.4	834
TOTAL CANNABINOIDS	95.9	959
<u>Terpenes (Top 5)</u>	<u>(%)</u>	<u>µg/g</u>
<u>Terpenes (Top 5)</u> Terpinolene	<u>(%)</u> 1.56	<mark>µg/g</mark> 15600
Terpinolene	1.56	15600
Terpinolene β-Myrcene	1.56 1.02	15600 10200
Terpinolene β-Myrcene d-Limonene	1.56 1.02 0.449	15600 10200 4490

<u>Contaminants</u>	PASS/FAIL
Heavy Metals:	PASS
Microbiology:	PASS
Pesticides:	PASS
<b>Residual Solvents:</b>	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: SA35049)**

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Cultivar (Strain) or Sample	<b>Date Completed:</b> 04/27/2023		

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

Analysis Date/Time: 04/25/2023 1645 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> g <u>Unit</u> (mg)
CBC	ND	ND	0.533	1.24	-	-	-
CBCA	ND	ND	1.66	3.86	-	-	-
CBD	ND	ND	3.76	8.78	-	-	-
CBDA	ND	ND	1.38	3.23	-	-	-
CBDV	ND	ND	0.604	1.41	-	-	-
CBDVA	ND	ND	1.61	3.76	-	-	-
CBG	ND	ND	2.44	5.70	-	-	-
CBGA	1.18	11.8	1.75	4.08	-	11.8	11.8
CBL	ND	ND	2.82	6.59	-	-	-
CBN	<loq< td=""><td><loq< td=""><td>1.30</td><td>3.03</td><td>-</td><td>-</td><td>-</td></loq<></td></loq<>	<loq< td=""><td>1.30</td><td>3.03</td><td>-</td><td>-</td><td>-</td></loq<>	1.30	3.03	-	-	-
CBNA	ND	ND	1.40	3.26	-	-	-
Δ9-ТНС	4.77	47.7	1.55	3.62	-	47.7	47.7
$\Delta 8$ -THC	ND	ND	2.42	5.66	-	-	-
THCA	89.6	896	0.842	1.97	-	896	896
THCV	ND	ND	2.02	4.71	-	-	-
THCVA	0.300	3.00	0.646	1.50	-	3.00	3.00
TOTAL	95.9	959			-	959	959
TOTAL CBC	-	-				-	-
TOTAL CBD	-	-			-	-	-
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	1.04	10.4			-	10.4	10.4
TOTAL CBN	-	-			-	-	-
TOTAL THC	83.4	834			-	834	834
TOTAL THCV	0.260	2.60			-	2.60	2.60

\* CBCA, CBDVA, CBL, CBNA, and THCVA are in process of being added to our accreditation scope.

*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): 1.00 SERVINGS/UNIT: 1

"-" 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) +  $\Delta$ 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.

Kyle W. Felling, Ph.D. atory Dire











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License: ADH 113	License: 00065C	E20230417ICMLS01	Date Received: 04/24/2023
Cultivar (Strain) or Sample	<b>Date Completed:</b> 04/27/2023		

# Analysis Date/Time+04/25/2023 1920

# **TERPENOID PROFILE**

Analysis Date/Time:04/25/2023 1920 Analyst: KF			ethod: GC/MS strument: Agilent 7890/5975	<b>Deviations from SOP:</b> None			
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)					
α-Bisabolol	-	-					
Camphene	583	0.0583					
δ-3-Carene	872	0.0872					
β-Caryophyllene	1580	0.158					
Caryophyllene oxide	-	-					
p-Cymene	523	0.0523					
Eucalyptol	327	0.0327					
Geraniol	-	-					
Guaiol	-	-		Abbrevistioner CC Con			
α-Humulene	963	0.0963		<i>Abbreviations:</i> GC - Gas Chromatography, MS - Mass			
Isopulegol	-	-		Spectrometry, RL - Reporting Limit			
d-Limonene	4490	0.449		This information is provided as a service and makes no claims of efficacy and/or			
Linalool	975	0.0975		safety of this product.			
β-Myrcene	10200	1.02		Results are applicable only for the sample(s) analyzed and for the specific			
cis-Nerolidol	-	-		analysis conducted.			
trans-Nerolidol α-Ocimene	491	0.0491		This report is for informational purposes only and should not be used to diagnose, treat or provot any			
β-Ocimene	2020	0.202		treat, or prevent any medical-related symptoms.			
α-Pinene	1370	0.202	1	The statements and results herein have			
β-Pinene	2010	0.137		not been approved and/or endorsed by the FDA.			
α-Terpinene	995	0.0995	1				
γ-Terpinene	785	0.0785					
Terpinolene	15600	1.56					
				"-" Not detected above RL.			
TOTAL	43800	4.38		<b>Reporting Limit (µg/g):</b> 46.8			

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











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License: ADH 113	License: 00065C	E20230417ICMLS01	Date Received: 04/24/2023
Cultivar (Strain) or Sample	Date Completed: 04/27/2023		

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

Analysis Date/Time: 04/26/2023 0934			Method: HS/GC/MS			<b>Deviations from SOP:</b>				
Analyst: KF				]	nstrument: Agilent 78	90/5975		N	lone	
<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)	-	37.2	74.4	5000	n-Heptane (142-82-5)	-	37.2	74.4	5000	
Acetonitrile (75-5-8)	-	37.2	74.4	410	n-Hexane (110-54-3)	-	13.0	26.0	290	
Benzene (71-43-2)	-	0.372	0.744	2	Isobutane (75-28-5)	-	37.2	74.4	5000	
n-Butane (106-97-2)	414	37.2	74.4	5000	Isopropanol (67-63-0)	-	37.2	74.4	5000	
1-Butanol (71-36-3)	-	37.2	74.4	5000	Isopropyl acetate		37.2	74.4	5000	
2-Butanol (78-92-2)	-	37.2	74.4	5000	(108-21-4)		57.2	/ 4.4	5000	
2-Butanone (78-93-3)	-	37.2	74.4	5000	Isopropyl benzene (98-82-8)	-	3.72	7.44	70	
Cyclohexane (110-82-7)	-	37.2	74.4	3880	(90-02-0) Methanol (67-56-1)		37.2	74.4	3000	Color Key
1,2-Dimethoxyethane	_	3.72	7.44	100	2-Methylbutane (78-78-4)		37.2	74.4	5000	
(110-71-4) N,N-Dimethylacetamide	_	37.2	74.4	1090	Methylene chloride (75-9-2)	-	37.2	74.4	600	RESULT < AL RESULT > AL
(127-19-5) 2,2-Dimethylbutane (75-83-2)	_	13.0	26.0	290	2-Methylpentane (107-83-5)	-	13.0	26.0	290	"DET" detected less than LOQ
2,3-Dimethylbutane					3-Methylpentane (96-10-0)	-	13.0	26.0	290	"-" not detected above
(79-29-8)	-	13.0	26.0	290	n-Pentane (109-66-0)	-	37.2	74.4	5000	LOD
N,N-Dimethylformamide		27.2	74.4	880	1-Pentanol (71-41-0)	-	37.2	74.4	5000	"*" - o,m,p-Xylene and
(68-12-2)		37.2	/4.4	000	n-Propane (74-98-6)	-	37.2	74.4	5000	Ethylbenzene
Dimethylsulfoxide (67-68-5)	-	37.2	74.4	5000	1-Propanol (71-23-8)	-	37.2	74.4	5000	
(07-00-3) 1,4-Dioxane (123-91-1)		37.2	74.4	380	Pyridine (110-86-1)	-	13.0	26.0	200	Action levels are referenced from the State of
Ethanol (64-17-5)	-	37.2	74.4	5000	Tetrahydrofuran (109-99-9)	-	37.2	74.4	720	Arkansas
2-Ethoxyethanol (110-80-5)	-	13.0	26.0	160	Tetramethylene sulfone	_	13.0	26.0	160	MMJ testing
Ethyl ether (60-29-7)		37.2	20.0 74.4	5000	(126-33-0)					guidelines.
Ethyl acetate (141-78-6)		37.2	74.4	5000	Toluene (108-88-3)	-	37.2	74.4	890	A value of "-"
Ethyl benzene (100-41-4)		37.2	74.4	2170	o-Xylene (95-47-6)	-	37.2	74.4	2170	for the action level
Ethylene glycol (107-21-1)	_	37.2	74.4	620	m,p-Xylene (108-38-3 or 106-42-3)	-	37.2	74.4	2170	means that analyte
Ethylene oxide (75-21-8)	-	3.72	7.44	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170	is not currently regulated by the regulations referenced above.
Solvent		Synonym(s	5)		Solvent	<u>S</u>	ynonym(s)			regulations referenced above.
Acetonitrile		Methyl Cya	anide, ACN		Ethylene glycol	1	,2-Ethanedio	I		
1-Butanol		n-Butanol,	Butyl Alco	hol	Isobutane	2	-Methylpropa	ane		
2-Butanol		sec-Butyl a	lcohol		Isopropanol	2	-Propanol, IP.	A		
2-Butanone		Methyl eth	yl ketone, N	/IEK	Isopropyl Acetate	А	cetic acid iso	propyl este	r	
1,2-Dimethoxyethane		Monoglym	e		Methanol	N	lethyl alcoho	1		
2,3-Dimethylbutane		Neohexane	!		2-Methylbutane	Is	sopentane			
2,3-Dimethylbutane		Diisopropy	/l		Methylene chloride	D	Dichlorometha	ane		
N,N-Dimethylformamide		DMF			2-Methylpentane	Is	sohexane			
Dimethysufoxide		DMSO			1-Pentanol	n	-Amyl alcoho	ol		
2-Ethoxyethanol		Cellosolve	, Ethyl glyc	ol	1-Propanol	P	ropyl alcohol	l		
Ethyl ether		Diethyl eth	ier, Ether		Tetrahydrofuran	Т	ΉF			
Ethyl acetate		EtOAc			Tetramethylene sulfone	S	ulfolane			
Ethyl benzene		Phenyletha	ne		Xylene	D	Dimethylbenz	ene		

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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**Deviations from SOP:** 

None

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Cultivar (Strain) or Sample D	Date Completed: 04/27/2023		

## Analysis Date/Time: 04/25/2023 1807 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> Level (µg/g)	<u>Pesticide</u>	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	
Abamectin (71751-41-2)	-	0.0468	0.375	0.5	Kresoxim-methyl	_	0.0468	0.375	0.4	
Acephate (30560-19-1)	-	0.0468	0.375	0.4	(143390-89-0)					
Acequinocyl (57960-19-7)	-	0.0468	0.375	2	Malathion (121-75-5)	-	0.0468	0.375	0.2	
Acetamiprid (135410-20-7)	-	0.0468	0.375	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.0468 0.0468	0.375 0.375	0.2 0.2	<u>Color Key</u>
Aldicarb (116-06-3)	-	0.0468	0.375	0.4	Methomyl (16752-77-5)	-	0.0468	0.375	0.4	<b>RESULT &lt; AL</b>
Azoxystrobin (131860-33-8)	-	0.0468	0.375	0.2	Methyl parathion (298-0-0)	-	0.0468	0.375	0.2	<b>RESULT &gt; AL</b>
Bifenazate (149877-41-8)	-	0.0468	0.375	0.2	MGK 264 (113-48-4)	-	0.0468	0.375	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3)	-	0.0468	0.375	0.2	Myclobutanil	_	0.0468	0.375	0.2	"-" not detected above
Boscalid (188425-85-6)	-	0.0468	0.375	0.4	(88671-89-0)					LOD
Carbaryl (63-25-2)	-	0.0468	0.375	0.2	Naled (300-76-5)	-	0.0468	0.375	0.5	Permethrins measured as the
Carbofuran (1563-66-2)	-	0.0468	0.375	0.2	Oxamyl (23135-22-0)	-	0.0468	0.375	1	cumulative residue of the <i>cis</i> - and
Chlorantraniliprole (800008-45-7)	-	0.0468	0.375	0.2	Paclobutrazol (76738-62-0)	-	0.0468	0.375	0.4	trans- permethrin isomers.
Chlorfenapyr	_	0.0468	0.375	1	Permethrins (52645-53-1)	-	0.0468	0.375	0.2	Pyrethrins measured as the cumulative residue of the
(122453-73-0)	-			1	Phosmet (732-11-6)	-	0.0468	0.375	0.2	pyrethrin I, cinerin I, and jasmolin
Chlorpyrifos (2921-88-2)	-	0.0468	0.375	0.2	Piperonyl butoxide (51-03-6)	-	0.0468	0.375	2	I isomers.
Clofentezine (74115-24-5)	-	0.0468	0.375	0.2	(51-03-0) Prallethrins (2331-36-9)		0.0468	0.375	0.2	Action levels are referenced from
Cyfluthrin (68359-37-5)	-	0.0468	0.375	1	Propiconazole					the
Cypermethrin (52315-07-8)	-	0.0468	0.375	1	(60207-90-1))	-	0.0468	0.375	0.4	State of Arkansas MMJ testing guidelines.
Daminozide (1596-84-5)	-	0.0468	0.375	1	Propoxur (114-26-1)	-	0.0468	0.375	0.2	A value of "-" for the action level
DDVP (62-73-7)	-	0.0468	0.375	0.1	Pyrethrins (8003-34-7)	-	0.0468	0.375	1	means that analyte is not
Diazinon (333-41-5)	-	0.0468	0.375	0.2	Pyridaben (96489-71-3)	-	0.0468	0.375	0.2	currently regulated by the regulations referenced above.
Dimethoate (60-51-5)	-	0.0468	0.375	0.2	Spinosad (168316-95-8)	-	0.0468	0.375	0.2	5
Ethoprophos (13194-48-4)	-	0.0468	0.375	0.2	Spiromesifen (283594-90-1)	-	0.0468	0.375	0.2	Disclaimer: This information is
Etofenprox (80844-07-1)	-	0.0468	0.375	0.4	Spirotetramat					provided as a service and makes no claims of efficacy and/or safety
Etoxazole (153233-91-1)	-	0.0468	0.375	0.2	(203313-25-1)	-	0.0468	0.375	0.2	of this product. Results are
Fenoxycarb (72490-01-8)	-	0.0468	0.375	0.2	Spiroxamine	_	0.0468	0.375	0.4	applicable only for the sample(s) analyzed and for the specific
(E)-Fenpyroximate (134098-61-6)	-	0.0468	0.375	0.4	(118134-30-8) Tebuconazole		0.0468	0.375	0.4	analysis conducted. This report is for informational purposes only
Fipronil (120068-37-3)	-	0.0468	0.375	0.4	(80443-41-0)		0.0400	0.373	0.4	and should not be used to
Flonicamid (158062-67-0)	-	0.0468	0.375	1	Thiacloprid (111988-49-9)	-	0.0468	0.375	0.2	diagnose, treat, or prevent any medical-related symptoms. The
Fludioxinil (131341-86-1)	-	0.0468	0.375	0.4	· · · · · ·					statements and results herein have
Hexythiazox (78587-05-0)	-	0.0468	0.375	1	Thiamethoxam (153719-23-4)	-	0.0468	0.375	0.2	not been approved and/or endorsed by the FDA.
Imazalil (35554-44-0)	-	0.0468	0.375	0.2	Trifloxystrobin		0.0460	0.275	0.2	enuoised by the FDA.
Imidacloprid (138261-41-3)	-	0.0468	0.375	0.4	(141517-21-7)	-	0.0468	0.375	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

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



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Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13246623704	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	<b>Production Run:</b>	Date Collected: 04/24/2023
License: ADH 113	License: 00065C	E20230417ICMLS01	Date Received: 04/24/2023
Cultivar (Strain) or Sample	Date Completed: 04/27/2023		

#### HEAVY METAL PROFILE (SOP: SOP-ICP-200.7)

Analysis Date/Time: 04/25/2023 1743 (ICP/OES) Analysis Date/Time: - (DMA) Analyst: KF		<b>Method:</b> ICP/OES <b>Instrument:</b> Agilent 720-ES		<b>Deviations from SOP:</b> ES None	
<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)	-	57.8	91.5	200	
Cadmium (Cd)	-	57.8	91.5	200	
Lead (Pb)	-	57.8	91.5	500	
Mercury (Hg)	-	57.8	91.5	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



"DET" detected less than LOQ

"-" not detected above LOD

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.













# **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35049)**

Testing Location:	Customer ID: 2168	Sample ID: SA35049	Sample Type: Primary
Arkansas	<b>River Valley Relief MIPS</b>	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13246623704	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 04/24/2023
License: ADA 05_H273	License: 00065C	E20230417ICMLS01	Date Received: 04/24/2023
Cultivar (Strain) or Sample	<b>Date Completed:</b> 04/27/2023		

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

Analysis Date/Time: 2023042 Analyst: PW		ardy Diagnostics Compac : 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	1	
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





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.

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