



## SUMMARY OF ANALYSIS (SAMPLE ID: SA39216)

Testing Location:	Customer ID: 2168	<b>Order ID:</b> OR11261	Sample Type: Primary
Arkansas	<b>River Valley Relief MIPS</b>	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13351352162	Mass: 4ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 10/28/2024
License: ADH 113	License: 00065C	E20241025RAW23	Date Received: 10/28/2024
Cultivar (Strain) or Sample De	<b>Date Completed:</b> 10/30/2024		

\*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.

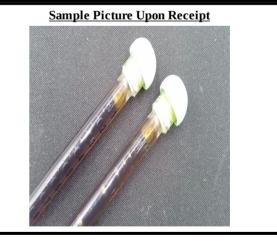
Moisture Content (%)	Water Activity (aw)	PASS/FAIL		
Not Tested	Not Tested	PASS		

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.

	(0/)	,
<u>Cannabinoids (Top 3)</u>	(%)	<u>mg/g</u>
Δ9-ТНС	82.6	826
CBG	2.92	29.2
CBC	0.794	7.94
TOTAL CBD	-	-
TOTAL THC	82.6	826
TOTAL CANNABINOIDS	87.2	872
<u>Terpenes (Top 5)</u>	<u>(%)</u>	<u>µg∕g</u>
<u>Terpenes (Top 5)</u> β-Caryophyllene	<u>(%)</u> 1.09	<mark>µg/g</mark> 10900
β-Caryophyllene	1.09	10900
β-Caryophyllene α-Humulene	1.09 0.435	10900 4350
β-Caryophyllene α-Humulene Guaiol	1.09 0.435 0.417	10900 4350 4170

PASS/FAIL
PASS
PASS
PASS
PASS





Scan the QR code to verify results.

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

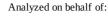
REPORT OF LABORATORY ANALYSIS

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License: ADH 113	License: 00065C	E20241025RAW23	Date Received: 10/28/2024
Cultivar (Strain) or Sample I	<b>Date Completed:</b> 10/30/2024		

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

Analysis Date/Time: 10/29/2024 1450 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	0.794	7.94	0.111	0.260	-	7.94	7.94	
CBCA	ND	ND	0.346	0.808	-	-	-	
CBD	ND	ND	0.786	1.84	-	-	-	
CBDA	ND	ND	0.289	0.676	-	-	-	
CBDV	ND	ND	0.126	0.294	-	-	-	
CBDVA	ND	ND	0.336	0.785	-	-	-	
CBG	2.92	29.2	0.510	1.19	-	29.2	29.2	
CBGA	ND	ND	0.723	0.854	-	-	-	
CBL	ND	ND	0.590	1.38	-	-	-	
CBN	0.485	4.85	0.271	0.633	-	4.85	4.85	
CBNA	ND	ND	0.292	0.681	-	-	-	
Δ9-ΤΗC	82.6	826	0.325	0.757	-	826	826	
$\Delta 8$ -THC	ND	ND	0.506	1.18	-	-	-	
THCA	ND	ND	0.176	0.412	-	-	-	I
THCV	0.341	3.41	0.422	0.985	-	3.41	3.41	ľ
THCVA	ND	ND	0.135	0.314	-	-	-	I
TOTAL	87.2	872				872	872	I
TOTAL CBC	0.794	7.94				7.94	7.94	5
TOTAL CBD	-	-			-	-	-	
TOTAL CBDV	-	-			-	-	-	V
TOTAL CBG	2.92	29.2			-	29.2	29.2	C
TOTAL CBN	0.485	4.85			-	4.85	4.85	
TOTAL THC	82.6	826			-	826	826	]
TOTAL THCV	0.341	3.41			-	3.41	3.41	] 7

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.

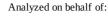
mg/g results are reported as mass component per mass material.











**Deviations from SOP:** 



# **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA39216)**

Testing Location:	Customer ID: 2168	Order ID: OR11261	Sample Type: Primary
Arkansas	<b>River Valley Relief MIPS</b>	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13351352162	Mass: 4ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 10/28/2024
License: ADH 113	License: 00065C	E20241025RAW23	Date Received: 10/28/2024
Cultivar (Strain) or Sample	Date Completed: 10/30/2024		

Analysis Date/Time:10/30/2024 0136 Analyst: KF

**Result** 

<u>(µg/g)</u> 2580

138

ND

10900

639

<LOQ

<LOQ

159

4170

4350

ND

4060

3510

3050

ND

1790

ND

1010

707

805

<LOQ

<LOQ

158

38000

Terpene

α-Bisabolol Camphene

δ-3-Carene

Eucalyptol

α-Humulene

Isopulegol

Linalool

β-Myrcene

cis-Nerolidol

α-Ocimene

β-Ocimene

α-Pinene

**B**-Pinene

α-Terpinene

y-Terpinene

Terpinolene

TOTAL

trans-Nerolidol

d-Limonene

Geraniol

Guaiol

oxide p-Cymene

β-Caryophyllene

Caryophyllene

#### **TERPENOID PROFILE**

**Method:** GC/MS **Instrument:** Agilent 7890/5975

<u>Result</u>

(%)

0.258

0.0138

\_

1.09

0.0639

0.0159

0.417

0.435

\_

0.406

0.351

0.305

0.179

0.101

0.0707

0.0805

0.0158

3.80

975	None
	Abbreviations: GC - Gas Chromatography, MS - Mass

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): 48.9

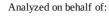
"-" Not detected above LOD.













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License: ADH 113	License: 00065C	E20241025RAW23	Date Received: 10/28/2024
Cultivar (Strain) or Sample	<b>Date Completed:</b> 10/30/2024		

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

Analysis Date/Time:	nalysis Date/Time: 10/29/2024 0118 Method: HS/GC/MS		<b>Deviations from SOP:</b>								
Analyst: KF				I	<b>Instrument:</b> Agilent 78	90/5975		None			
<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)	-	33.7	67.4	5000	n-Heptane (142-82-5)	-	33.7	67.4	5000		
Acetonitrile (75-5-8)	-	33.7	67.4	410	n-Hexane (110-54-3)	-	11.8	23.6	290		
Benzene (71-43-2)	-	0.337	0.674	2	Isobutane (75-28-5)	-	33.7	67.4	5000		
n-Butane (106-97-2)	-	33.7	67.4	5000	Isopropanol (67-63-0)	-	33.7	67.4	5000		
1-Butanol (71-36-3)	-	33.7	67.4	5000	Isopropyl acetate	_	33.7	67.4	5000		
2-Butanol (78-92-2)	-	33.7	67.4	5000	(108-21-4)		55.7	07.4	5000		
2-Butanone (78-93-3)	-	33.7	67.4	5000	Isopropyl benzene (98-82-8)	-	3.37	6.74	70		
Cyclohexane (110-82-7)	-	33.7	67.4	3880	(38-82-8) Methanol (67-56-1)		33.7	67.4	3000	Color Key	
1,2-Dimethoxyethane (110-71-4)	-	3.37	6.74	100	2-Methylbutane (78-78-4)	-	33.7	67.4	5000	RESULT < AL	
N,N-Dimethylacetamide (127-19-5)	-	33.7	67.4	1090	Methylene chloride (75-9-2)	-	33.7	67.4	600	<b>RESULT &gt; AL</b>	
2,2-Dimethylbutane (75-83-2)	-	11.8	23.6	290	2-Methylpentane (107-83-5)	-	11.8	23.6	290	"-" not detected above 1/2 Action Level	
2,3-Dimethylbutane		11.0	22.6	200	3-Methylpentane (96-10-0)	-	11.8	23.6	290		
(79-29-8)	-	11.8	23.6	290	n-Pentane (109-66-0)	-	33.7	67.4	5000	"*" - o,m,p-Xylene and	
N,N-Dimethylformamide (68-12-2)	-	33.7	67.4	880	1-Pentanol (71-41-0) n-Propane (74-98-6)	-	33.7 33.7	67.4 67.4	5000 5000	Ethylbenzene Action levels are	
Dimethylsulfoxide		33.7	67.4	5000	1-Propanol (71-23-8)	_	33.7	67.4	5000	referenced from the State of	
(67-68-5)					Pyridine (110-86-1)	_	11.8	23.6	200	Arkansas	
1,4-Dioxane (123-91-1)	-	33.7	67.4	380	Tetrahydrofuran (109-99-9)	_	33.7	67.4	720	MMJ testing	
Ethanol (64-17-5)	-	33.7	67.4	5000	Tetramethylene sulfone					guidelines.	
2-Ethoxyethanol (110-80-5)	-	11.8	23.6	160	(126-33-0)	-	11.8	23.6	160	A value of "-"	
Ethyl ether (60-29-7)	-	33.7	67.4	5000	Toluene (108-88-3)	-	33.7	67.4	890	for the action level	
Ethyl acetate (141-78-6)	-	33.7	67.4	5000	o-Xylene (95-47-6)	-	33.7	67.4	2170	means that analyte	
Ethyl benzene (100-41-4)	-	33.7	67.4	2170	m,p-Xylene (108-38-3 or	_	33.7	67.4	2170	is not currently regulated by the	
Ethylene glycol (107-21-1)	-	33.7	67.4	620	106-42-3)					regulations referenced above.	
Ethylene oxide (75-21-8)	-	3.37	6.74	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170		
Solvent		Synonym(s	5)		Solvent	Sy	nonym(s)				
Acetonitrile		Methyl Cya	anide, ACN		Ethylene glycol	1,2	2-Ethanediol				
1-Butanol		n-Butanol,	Butyl Alco	hol	Isobutane	2-	Methylpropa	ne			
2-Butanol		sec-Butyl a	lcohol		Isopropanol	2-	Propanol, IPA	1			
2-Butanone		Methyl ethy	yl ketone, N	/IEK	Isopropyl Acetate	A	cetic acid iso	propyl ester			
1,2-Dimethoxyethane		Monoglym	e		Methanol	Μ	ethyl alcohol				
2,3-Dimethylbutane		Neohexane			2-Methylbutane	Ise	opentane				
2,3-Dimethylbutane		Diisopropy	rl		Methylene chloride	Di	ichlorometha	ne			
N,N-Dimethylformamide		DMF			2-Methylpentane	Ise	ohexane				
Dimethysufoxide		DMSO			1-Pentanol	n-	Amyl alcoho	1			
2-Ethoxyethanol		Cellosolve,	Ethyl glyc	ol	1-Propanol	Pr	opyl alcohol				
Ethyl ether		Diethyl eth	er, Ether		Tetrahydrofuran	TI	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

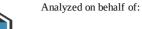
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**Deviations from SOP:** 

None



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Cultivar (Strain) or Sample De	Date Completed: 10/30/2024		

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

Analysis Date/Time: 10/29/2024 1558 Analyst: KF

#### 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> (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)	-	0.0486	0.389	0.5	Kresoxim-methyl		0.0406	0.200		
Acephate (30560-19-1)	-	0.0486	0.389	0.4	(143390-89-0)	-	0.0486	0.389	0.4	
Acequinocyl (57960-19-7)	-	0.0486	0.389	2	Malathion (121-75-5)	-	0.0486	0.389	0.2	
Acetamiprid		0.0400	0.200	0.2	Metalaxyl (57837-19-1)	-	0.0486	0.389	0.2	
(135410-20-7)		0.0486	0.389	0.2	Methiocarb (2032-65-7)	-	0.0486	0.389	0.2	<u>Color Key</u>
Aldicarb (116-06-3)	-	0.0486	0.389	0.4	Methomyl (16752-77-5)	-	0.0486	0.389	0.4	<b>RESULT &lt; AL</b>
Azoxystrobin (131860-33-8)	-	0.0486	0.389	0.2	Methyl parathion (298-0-0)	-	0.0486	0.389	0.2	RESULT > AL
Bifenazate (149877-41-8)	-	0.0486	0.389	0.2	MGK 264 (113-48-4)	-	0.0486	0.389	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3)	-	0.0486	0.389	0.2	Myclobutanil	_	0.0486	0.389	0.2	"-" not detected above
Boscalid (188425-85-6)	-	0.0486	0.389	0.4	(88671-89-0)			0.000		LOD
Carbaryl (63-25-2)	-	0.0486	0.389	0.2	Naled (300-76-5)	-	0.0486	0.389	0.5	Permethrins measured as the
Carbofuran (1563-66-2)	-	0.0486	0.389	0.2	Oxamyl (23135-22-0)	-	0.0486	0.389	1	cumulative residue of the <i>cis</i> - and
Chlorantraniliprole (800008-45-7)	-	0.0486	0.389	0.2	Paclobutrazol (76738-62-0)	-	0.0486	0.389	0.4	trans- permethrin isomers.
Chlorfenapyr		0.0486	0.389	1	Permethrins (52645-53-1)	-	0.0486	0.389	0.2	Pyrethrins measured as the cumulative residue of the
(122453-73-0)		0.0400	0.369	1	Phosmet (732-11-6)	-	0.0486	0.389	0.2	pyrethrin I, cinerin I, and jasmolin
Chlorpyrifos (2921-88-2)	-	0.0486	0.389	0.2	Piperonyl butoxide	-	0.0486	0.389	2	I isomers.
Clofentezine (74115-24-5)	-	0.0486	0.389	0.2	(51-03-6)		0.0486	0.389	0.2	Action levels are
Cyfluthrin (68359-37-5)	-	0.0486	0.389	1	Prallethrins (2331-36-9)	-	0.0486	0.389	0.2	referenced from the State of
Cypermethrin (52315-07-8)	-	0.0486	0.389	1	Propiconazole (60207-90-1))	-	0.0486	0.389	0.4	Arkansas MMJ testing
Daminozide (1596-84-5)	-	0.0486	0.389	1	Propoxur (114-26-1)	-	0.0486	0.389	0.2	guidelines.
DDVP (62-73-7)	-	0.0486	0.389	0.1	Pyrethrins (8003-34-7)	-	0.0486	0.389	1	A value of "-" for the action level
Diazinon (333-41-5)	-	0.0486	0.389	0.2	Pyridaben (96489-71-3)	-	0.0486	0.389	0.2	means that analyte is not currently regulated by the
Dimethoate (60-51-5)	-	0.0486	0.389	0.2	Spinosad (168316-95-8)	-	0.0486	0.389	0.2	regulations referenced above.
Ethoprophos (13194-48-4)	-	0.0486	0.389	0.2	Spiromesifen (283594-90-1)	-	0.0486	0.389	0.2	Disclaimer: This information is
Etofenprox (80844-07-1)	-	0.0486	0.389	0.4	Spirotetramat					provided as a service and makes
Etoxazole (153233-91-1)	-	0.0486	0.389	0.2	(203313-25-1)	-	0.0486	0.389	0.2	no claims of efficacy and/or safety
Fenoxycarb (72490-01-8)	-	0.0486	0.389	0.2	Spiroxamine	_	0.0486	0.389	0.4	of this product. Results are applicable only for the sample(s)
(E)-Fenpyroximate (134098-61-6)	-	0.0486	0.389	0.4	(118134-30-8) Tebuconazole					analyzed and for the specific analysis conducted. This report is
Fipronil (120068-37-3)	-	0.0486	0.389	0.4	(80443-41-0)	-	0.0486	0.389	0.4	for informational purposes only
Flonicamid (158062-67-0)	-	0.0486	0.389	1	Thiacloprid	_	0.0486	0.389	0.2	and should not be used to diagnose, treat, or prevent any
Fludioxinil (131341-86-1)	-	0.0486	0.389	0.4	(111988-49-9)					medical-related symptoms. The
Hexythiazox (78587-05-0)	-	0.0486	0.389	1	Thiamethoxam (153719-23-4)	-	0.0486	0.389	0.2	statements and results herein have
Imazalil (35554-44-0)	-	0.0486	0.389	0.2	Trifloxystrobin					not been approved and/or endorsed by the FDA.
Imidacloprid (138261-41-3)	-	0.0486	0.389	0.4	(141517-21-7)		0.0486	0.389	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	<u>Synonym(s)</u>	Pesticide	Synonym(s)	Pesticide	<u>Synonym(s)</u>
Cyfluthrin	Baythroid	Myclobutanil	Systhane	Propiconazole	Tilt
DDVP	Dichlorvos	Naled	Dibrom	Propoxur	Baygon
Ethoprophos	Prophos	Phosmet	Imidan		



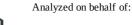
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Testing Location:	Customer ID: 2168	Order ID: OR11261	Sample Type: Primary
Arkansas	<b>River Valley Relief MIPS</b>	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13351352162	Mass: 4ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 10/28/2024
License: ADH 113	License: 00065C	E20241025RAW23	Date Received: 10/28/2024
Cultivar (Strain) or Sample	<b>Date Completed:</b> 10/30/2024		

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

Analysis Date/T	<b>`ime:</b> 10/30/2024 (	0033	<b>Method:</b> ICP/MS <b>Instrument:</b> Agilent 7500ce		<b>Deviations from SOP:</b> None	
Analyst: KF				-		
<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)	-	59.7	94.6	200		
Cadmium (Cd)	-	59.7	94.6	200		
Lead (Pb)	-	59.7	94.6	500		
Mercury (Hg)	-	59.7	94.6	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





"DET" detected less than LOQ

"-" not detected above LOD

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

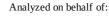
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Testing Location:	Customer ID: 2168	Sample ID: SA39216	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13351352162	Mass: 4ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 10/28/2024
License: ADA 05_H273	Date Received: 10/28/2024		
Cultivar (Strain) or Sample I	Date Completed: 10/30/2024		

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

Analysis Date/Time: 10/29/20 Analyst: PW		Hardy Diagnostics Compact ent: Thermo Incubator	Dry <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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