







SUMMARY OF ANALYSIS (SAMPLE ID: SA34487)

Testing Location:	Customer ID: 2168	Order ID: OR10317	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13242531430	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 02/16/2023
License: ADH 113	License: 00065C	E20230213RFLLS02	Date Received: 02/17/2023
Cultivar (Strain) or Sample D	Date Completed: 02/23/2023		

*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>	<u>mg</u> /g
THCA	80.3	803
CBGA	2.98	29.8
Δ9-THC	1.33	13.3
TOTAL CBD	-	-
TOTAL THC	71.8	718
TOTAL CANNABINOIDS	85.5	855
<u>Terpenes (Top 5)</u>	<u>(%)</u>	<u>µg/g</u>
<u>Terpenes (Top 5)</u> d-Limonene	<u>(%)</u> 0.832	<u>µg/g</u> 8320
d-Limonene	0.832	8320
d-Limonene β-Caryophyllene	0.832 0.550	8320 5500
d-Limonene β-Caryophyllene β-Myrcene	0.832 0.550 0.492	8320 5500 4920

<u>Contaminants</u>	PASS/FAIL
Heavy Metals:	PASS
Microbiology:	PASS
Pesticides:	PASS
Residual Solvents:	PASS





Scan the QR code to verify results.

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.

Kyle W. Felling, Ph.D. aboratory Director

REPORT OF LABORATORY ANALYSIS

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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA34487)

Testing Location:	Customer ID: 2168	Order ID: OR10317	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13242531430	Mass: 4g
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Cultivar (Strain) or Sample	Date Completed: 02/23/2023		

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

Analysis Date/Time: 2/17/2023 1449 Analyst: PW Method: HPLC/DAD Instrument: Agilent 1100

<u>Cannabinoid</u>	<u>Result</u> (%)	<u>Result</u> (mg/g)	LOD (mg/g)	<u>LOQ</u> (mg/g)	<u>Result</u> (mg/ mL)	<u>Per</u> Serving (mg)	<u>Per</u> g <u>Unit</u> (mg)
CBC	ND	ND	0.492	1.15	-	-	-
CBCA	0.477	4.77	1.53	3.56	-	4.77	4.77
CBD	ND	ND	3.47	8.10	-	-	-
CBDA	ND	ND	1.28	2.98	-	-	-
CBDV	ND	ND	0.557	1.30	-	-	-
CBDVA	ND	ND	1.48	3.47	-	-	-
CBG	<loq< td=""><td><loq< td=""><td>2.25</td><td>5.26</td><td>-</td><td>-</td><td>-</td></loq<></td></loq<>	<loq< td=""><td>2.25</td><td>5.26</td><td>-</td><td>-</td><td>-</td></loq<>	2.25	5.26	-	-	-
CBGA	2.98	29.8	1.61	3.77	-	29.8	29.8
CBL	ND	ND	2.60	6.08	-	-	-
CBN	ND	ND	1.20	2.79	-	-	-
CBNA	ND	ND	1.29	3.01	-	-	-
Δ9-ΤΗC	1.33	13.3	1.43	3.34	-	13.3	13.3
$\Delta 8$ -THC	ND	ND	2.24	5.22	-	-	-
THCA	80.3	803	0.777	1.82	-	803	803
THCV	ND	ND	1.86	4.35	-	-	-
THCVA	0.368	3.68	0.596	1.39	-	3.68	3.68
TOTAL	85.5	855				855	855
TOTAL CBC	0.419	4.19			-	4.19	4.19
TOTAL CBD	-	-			-	-	-
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	2.62	26.2			-	26.2	26.2
TOTAL CBN	-	-			-	-	-
TOTAL THC	71.8	718			-	718	718
TOTAL THCV	0.319	3.19			-	3.19	3.19

* 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) + 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. atory Directo

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Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 02/16/2023
License: ADH 113	License: 00065C	E20230213RFLLS02	Date Received: 02/17/2023
Cultivar (Strain) or Sample	Date Completed: 02/23/2023		

TERPENOID PROFILE

Analysis Date/Time:2/20/2023 1916 Analyst: KF			e thod: GC/MS s trument: Agilent 7890/5975	Deviations from SOP: None			
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)		e l'an			
α-Bisabolol	1490	0.149					
Camphene	404	0.0404					
δ-3-Carene	-	-					
β-Caryophyllene	5500	0.550					
Caryophyllene oxide	-	-					
p-Cymene	208	0.0208					
Eucalyptol	-	-					
Geraniol	-	-		a share a share a			
Guaiol	-	-					
α-Humulene	2130	0.213		<i>Abbreviations:</i> GC - Gas Chromatography, MS - Mass			
Isopulegol	-	-		Spectrometry, RL - Reporting Limit			
d-Limonene	8320	0.832		This information is provided as a service and makes no claims of efficacy and/or			
Linalool	3660	0.366		safety of this product.			
β-Myrcene cis-Nerolidol	4920	0.492		Results are applicable only for the sample(s) analyzed and for the specific analysis conducted.			
trans-Nerolidol	-	-		This report is for informational purposes			
α-Ocimene	-	-		only and should not be used to diagnose, treat, or prevent any			
β-Ocimene	-	-		medical-related symptoms.			
α-Pinene	747	0.0747	1	The statements and results herein have			
β-Pinene	1260	0.126	i	not been approved and/or endorsed by the FDA.			
α-Terpinene	-	-	-				
γ-Terpinene	-	-					
Terpinolene	424	0.0424	1	"-" Not detected above RL.			
TOTAL	29100	2.91		Reporting Limit (µg/g): 49.			













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Cultivar (Strain) or Sample	Date Completed: 02/23/2023		

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

Analysis Date/Time:	2/19/202	23 1232	100	Method: HS/GC/MS				Deviations from SOP:			
Analyst: KF				I	nstrument: Agilent 789	0/5975	None				
Solvent	<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)	LOQ (µg/g)	<u>Action</u> Level (µg/g)		
Acetone (67-64-1)	-	38.2	76.4	5000	n-Heptane (142-82-5)	-	38.2	76.4	5000		
Acetonitrile (75-5-8)	-	38.2	76.4	410	n-Hexane (110-54-3)	-	13.4	26.7	290		
Benzene (71-43-2)	-	0.382	0.764	2	Isobutane (75-28-5)	-	38.2	76.4	5000		
n-Butane (106-97-2)	-	38.2	76.4	5000	Isopropanol (67-63-0)	-	38.2	76.4	5000		
1-Butanol (71-36-3)	-	38.2	76.4	5000	Isopropyl acetate	2	38.2	76.4	5000		
2-Butanol (78-92-2)	-	38.2	76.4	5000	(108-21-4)						
2-Butanone (78-93-3)	-	38.2	76.4	5000	Isopropyl benzene (98-82-8)	-	3.82	7.64	70	a sufficient of the second	
Cyclohexane (110-82-7)	-	38.2	76.4	3880	Methanol (67-56-1)	_	38.2	76.4	3000	<u>Color Key</u>	
1,2-Dimethoxyethane	-	3.82	7.64	100	2-Methylbutane (78-78-4)	-	38.2	76.4	5000	DECHITZAI	
(110-71-4) N,N-Dimethylacetamide (127-19-5)	_	38.2	76.4	1090	Methylene chloride (75-9-2)	-	38.2	76.4	600	RESULT < AL RESULT > AL	
(127-13-3) 2,2-Dimethylbutane (75-83-2)	-	13.4	26.7	290	2-Methylpentane (107-83-5)	-	13.4	26.7	290	"DET" detected less than LOQ	
2,3-Dimethylbutane					3-Methylpentane (96-10-0)	-	13.4	26.7	290	"-" not detected above	
(79-29-8)	-	13.4	26.7	290	n-Pentane (109-66-0)	-	38.2	76.4	5000	LOD	
N,N-Dimethylformamide		38.2	76.4	880	1-Pentanol (71-41-0)	-	38.2	76.4	5000	"*" - o,m,p-Xylene and	
(68-12-2)		50.2	70.4	000	n-Propane (74-98-6)	-	38.2	76.4	5000	Ethylbenzene	
Dimethylsulfoxide (67-68-5)	-	38.2	76.4	5000	1-Propanol (71-23-8)	-	38.2	76.4	5000		
(07-00-3) 1,4-Dioxane (123-91-1)		38.2	76.4	380	Pyridine (110-86-1)	-	13.4	26.7	200	Action levels are referenced from the State of	
Ethanol (64-17-5)		38.2	76.4	5000	Tetrahydrofuran (109-99-9)	-	38.2	76.4	720	Arkansas	
2-Ethoxyethanol (110-80-5)	_	13.4	26.7	160	Tetramethylene sulfone	_	13.4	26.7	160	MMJ testing	
Ethyl ether (60-29-7)	_	38.2	76.4	5000	(126-33-0)					guidelines.	
Ethyl acetate (141-78-6)	_	38.2	76.4	5000	Toluene (108-88-3)	-	38.2	76.4	890	A value of "-"	
Ethyl benzene (100-41-4)	-	38.2	76.4	2170	o-Xylene (95-47-6)	-	38.2	76.4	2170	for the action level	
Ethylene glycol (107-21-1)	-	38.2	76.4	620	m,p-Xylene (108-38-3 or 106-42-3)	-	38.2	76.4	2170	means that analyte	
Ethylene oxide (75-21-8)	-	3.82	7.64	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170	is not currently	
Solvent		Synonym(s			Solvent	s	Synonym(s)			regulated by the regulations referenced above.	
Acetonitrile		Methyl Cya	-		Ethylene glycol	_	,2-Ethanediol				
1-Butanol		n-Butanol,		hol	Isobutane		-Methylpropa				
2-Butanol		sec-Butyl a		1101	Isopropanol		-Propanol, IP/				
2-Butanone		Methyl eth		/EK	Isopropyl Acetate		Acetic acid iso		r		
1,2-Dimethoxyethane		Monoglym			Methanol		Aethyl alcohol				
2,3-Dimethylbutane		Neohexane			2-Methylbutane		sopentane				
2,3-Dimethylbutane		Diisopropy			Methylene chloride		Dichlorometha	ne			
N,N-Dimethylformamide		DMF			2-Methylpentane		sohexane				
Dimethysufoxide		DMSO			1-Pentanol		-Amyl alcoho	1			
2-Ethoxyethanol		Cellosolve	, Ethyl glvc	ol	1-Propanol		ropyl alcohol				
Ethyl ether		Diethyl eth			Tetrahydrofuran		THF				
Ethyl acetate		EtOAc			Tetramethylene sulfone	S	Sulfolane				
Ethyl benzene		Phenyletha	ne		Xylene		Dimethylbenze	ene			
-		2					-				

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:

- -

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Cultivar (Strain) or Sample I	Date Completed: 02/23/2023		

Analysis Date/Time: 2/20/2023 1842

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Method: LC/MS/MS <u>.</u>...

Analyst: KF					Instrument: Shimadzu LC-8050			None		
<u>Pesticide</u>	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	Pesticide	<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)	-	0.0492	0.394	0.5	Kresoxim-methyl	_	0.0492	0.394	0.4	
Acephate (30560-19-1)	-	0.0492	0.394	0.4	(143390-89-0)					
Acequinocyl (57960-19-7)	-	0.0492	0.394	2	Malathion (121-75-5)	-	0.0492	0.394	0.2	
Acetamiprid (135410-20-7)	-	0.0492	0.394	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.0492 0.0492	0.394 0.394	0.2 0.2	
Aldicarb (116-06-3)	-	0.0492	0.394	0.4	Methomyl (16752-77-5)	-	0.0492	0.394	0.4	
Azoxystrobin (131860-33-8)	-	0.0492	0.394	0.2	Methyl parathion (298-0-0)	-	0.0492	0.394	0.2	<u>Color Key</u>
Bifenazate (149877-41-8)	-	0.0492	0.394	0.2	MGK 264 (113-48-4)	-	0.0492	0.394	0.2	RESULT < AL
Bifenthrin (82657-04-3)	-	0.0492	0.394	0.2	Myclobutanil		0.0492	0.394	0.2	RESULT > AL
Boscalid (188425-85-6)	-	0.0492	0.394	0.4	(88671-89-0)	-	0.0492	0.394	0.2	"DET" detected less than LOQ
Carbaryl (63-25-2)	-	0.0492	0.394	0.2	Naled (300-76-5)	-	0.0492	0.394	0.5	" "
Carbofuran (1563-66-2)	-	0.0492	0.394	0.2	Oxamyl (23135-22-0)	-	0.0492	0.394	1	"-" not detected above LOD
Chlorantraniliprole (800008-45-7)	-	0.0492	0.394	0.2	Paclobutrazol (76738-62-0)	-	0.0492	0.394	0.4	Permethrins measured as the
Chlorfenapyr		0.0400	0.004		Permethrins (52645-53-1)	-	0.0492	0.394	0.2	cumulative residue of the <i>cis</i> - and <i>trans</i> - permethrin isomers.
(122453-73-0)	-	0.0492	0.394	1	Phosmet (732-11-6)	-	0.0492	0.394	0.2	trans-permetititi isomers.
Chlorpyrifos (2921-88-2)	-	0.0492	0.394	0.2	Piperonyl butoxide	_	0.0492	0.394	2	Pyrethrins measured as the
Clofentezine (74115-24-5)	-	0.0492	0.394	0.2	(51-03-6)					cumulative residue of the pyrethrin I, cinerin I, and jasmolin
Cyfluthrin (68359-37-5)	-	0.0492	0.394	1	Prallethrins (2331-36-9)	-	0.0492	0.394	0.2	I isomers.
Cypermethrin (52315-07-8)	-	0.0492	0.394	1	Propiconazole (60207-90-1))	-	0.0492	0.394	0.4	Action levels are referenced from
Daminozide (1596-84-5)	-	0.0492	0.394	1	Propoxur (114-26-1)	-	0.0492	0.394	0.2	the State of Arkansas MMJ testing
DDVP (62-73-7)	-	0.0492	0.394	0.1	Pyrethrins (8003-34-7)	-	0.0492	0.394	1	guidelines.
Diazinon (333-41-5)	-	0.0492	0.394	0.2	Pyridaben (96489-71-3)	-	0.0492	0.394	0.2	A value of "-" for the action level
Dimethoate (60-51-5)	-	0.0492	0.394	0.2	Spinosad (168316-95-8)	-	0.0492	0.394	0.2	means that analyte is not
Ethoprophos (13194-48-4)	-	0.0492	0.394	0.2	Spiromesifen	_	0.0492	0.394	0.2	currently regulated by the
Etofenprox (80844-07-1)	-	0.0492	0.394	0.4	(283594-90-1)					regulations referenced above.
Etoxazole (153233-91-1)	-	0.0492	0.394	0.2	Spirotetramat (203313-25-1)	-	0.0492	0.394	0.2	Disclaimer: This information is
Fenoxycarb (72490-01-8)	-	0.0492	0.394	0.2	Spiroxamine		0.0400	0.004	<u>.</u>	provided as a service and makes no claims of efficacy and/or safety
(E)-Fenpyroximate (134098-61-6)	-	0.0492	0.394	0.4	(118134-30-8) Tebuconazole	-	0.0492	0.394	0.4	of this product. Results are applicable only for the sample(s)
Fipronil (120068-37-3)	-	0.0492	0.394	0.4	(80443-41-0)	-	0.0492	0.394	0.4	analyzed and for the specific
Flonicamid (158062-67-0)	-	0.0492	0.394	1	Thiacloprid	_	0.0492	0.394	0.2	analysis conducted. This report is for informational purposes only
Fludioxinil (131341-86-1)	-	0.0492	0.394	0.4	(111988-49-9)		5.0452	0.334	0.2	and should not be used to
Hexythiazox (78587-05-0)	-	0.0492	0.394	1	Thiamethoxam (153719-23-4)	-	0.0492	0.394	0.2	diagnose, treat, or prevent any
Imazalil (35554-44-0)	-	0.0492	0.394	0.2	(153719-23-4) Trifloxystrobin					medical-related symptoms. The statements and results herein have
Imidacloprid (138261-41-3)	-	0.0492	0.394	0.4	(141517-21-7)	-	0.0492	0.394	0.2	not been approved and/or endorsed by the FDA.

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) Cyfluthrin Baythroid DDVP Dichlorvos Prophos Ethoprophos

Pesticide Myclobutanil Naled Phosmet

Synonym(s) Systhane Dibrom Imidan

Pesticide Propiconazole Propoxur

Synonym(s) Tilt Baygon



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License: ADH 113	License: 00065C	E20230213RFLLS02	Date Received: 02/17/2023
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HEAVY METAL PROFILE (SOP: SOP-ICP-200.7)

Analysis Date/T Analysis Date/T Analyst: KF	ïme: 2/21/2023 11 ïme: - (DMA)	09 (ICP/OES)		nod: ICP/OES rument: Agilent 720-	Deviations from SOP:ESNone
<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)	- I and
Arsenic (As)	-	58.1	91.9	200	
Cadmium (Cd)	-	58.1	91.9	200	
Lead (Pb)	-	58.1	91.9	500	
Mercury (Hg)	-	58.1	91.9	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.

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: SA34487)

Testing Location:	Customer ID: 2168	Sample ID: SA34487	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13242531430	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 02/16/2023
License: ADA 05_H273	License: 00065C	E20230213RFLLS02	Date Received: 02/17/2023
Cultivar (Strain) or Sample Des	Date Completed: 02/23/2023		

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 2023022 Analyst: PW		ardy Diagnostics Compact t: Thermo Incubator	tDry Deviations from SOP: None
Bacteria/Microbe	<u>Result</u> (CFU/g)	<u>Action Level</u> (CFU/g)	e l'an
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





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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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