







SUMMARY OF ANALYSIS (SAMPLE ID: SA33699)

Testing Location:	Customer ID: 2168	Order ID: OR10217	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13241008157	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 12/14/2022
License: ADH 113	License: 00065C	E20221128O43HLR01	Date Received: 12/14/2022
Cultivar (Strain) or Sample De	Date Completed: 12/21/2022		

*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.

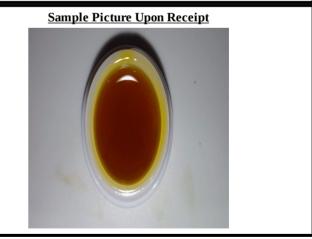
Moisture Content (%)	Water Activity (aw)	PASS/FAIL	
Not Tested	Not Tested	PASS	

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.

<u>Cannabinoids (Top 3)</u>	<u>(%)</u>	<u>mg/g</u>
Δ9-ТНС	29.9	299
THCA	28.4	284
CBGA	3.69	36.9
TOTAL CBD	0.342	3.42
TOTAL THC	54.8	548
TOTAL CANNABINOIDS	64.0	640
<u>Terpenes (Top 5)</u>	<u>(%)</u>	<u>µg/g</u>
<u>Terpenes (Top 5)</u> Terpinolene	<u>(%)</u> 2.92	<u>µg/g</u> 29200
Terpinolene	2.92	29200
Terpinolene β-Myrcene	2.92 2.15	29200 21500
Terpinolene β-Myrcene d-Limonene	2.92 2.15 0.925	29200 21500 9250

Contaminants	PASS/FAIL
Heavy Metals:	PASS
Microbiology:	PASS
Pesticides:	PASS
Residual Solvents:	PASS





Scan the QR code to verify results.

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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA33699)

Testing Location:	Customer ID: 2168	Order ID: OR10217	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
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License: ADH 113	License: 00065C	E20221128O43HLR01	Date Received: 12/14/2022
Cultivar (Strain) or Sample I	Date Completed: 12/21/2022		

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

Analysis Date/Time: 12/15/2022 2053 Analyst: PW Method: HPLC/DAD Instrument: Agilent 1100

<u>Cannabinoid</u>	<u>Result</u> (%)	<u>UM</u> (+/-%)	<u>Result</u> (mg/g)	LOQ (mg/g)	<u>Result</u> (mg/ <u>mL</u>)	<u>Per</u> Serving (mg)	<u>Per</u> g <u>Unit</u> (mg)
CBC	-	-	-	2.30	-	-	-
CBCA	-	-	-	2.30	-	-	-
CBD	-	-	-	2.30	-	-	-
CBDA	0.390	0.0351	3.90	2.30	-	3.90	3.90
CBDV	-	-	-	2.30	-	-	-
CBDVA	-	-	-	2.30	-	-	-
CBG	1.55	0.139	15.5	2.30	-	15.5	15.5
CBGA	3.69	0.332	36.9	2.30	-	36.9	36.9
CBL	-	-	-	2.30	-	-	-
CBN	-	-	-	2.30	-	-	-
CBNA	-	-	-	2.30	-	-	-
Δ9-ΤΗC	29.9	2.69	299	2.30	-	299	299
$\Delta 8$ -THC	-	-	-	2.30	-	-	-
THCA	28.4	2.56	284	2.30	-	284	284
THCV	-	-	-	2.30	-	-	-
THCVA	-	-	-	2.30	-	-	-
TOTAL	64.0	5.76	640		-	640	640
TOTAL CBC	-	-	-		-	-	-
TOTAL CBD	0.342	0.0308	3.42		-	3.42	3.42
TOTAL CBDV	-	-	-		-	-	-
TOTAL CBG	4.78	0.431	47.8		-	47.8	47.8
TOTAL CBN	-	-	-		-	-	-
TOTAL THC	54.8	4.94	548		-	548	548
TOTAL THCV	-	-	-		-	-	-

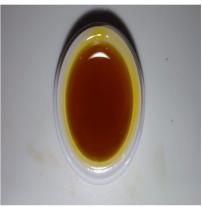
* CBCA, CBDVA, CBL, CBNA, and THCVA are in process of being added to our accreditation scope.

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 Dire

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Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 12/14/2022
License: ADH 113	License: 00065C	E20221128O43HLR01	Date Received: 12/14/2022
Cultivar (Strain) or Sample	Date Completed: 12/21/2022		

Analysis Date/Time:12/18/2022 1803 Analyst: KF

TERPENOID PROFILE

Method: GC/MS Instrument: Agilent 7890/5975

Deviations from SOP: None

Analyst. Iti		1112	di diffetit. Agricia 7050/5575	5 None			
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)					
α-Bisabolol	680	0.0680					
Camphene	553	0.0553					
δ-3-Carene	1500	0.150					
β-Caryophyllene	4850	0.485					
Caryophyllene oxide	249	0.0249					
p-Cymene	-	-					
Eucalyptol	-	-					
Geraniol	674	0.0674					
Guaiol	-	-		Abbreviations: GC - Gas			
α-Humulene	1460	0.146	1	Chromatography, MS - Mass			
Isopulegol	-	-	_	Spectrometry, RL - Reporting Limit			
d-Limonene Linalool	9250 2020	0.925 0.202	•	This information is provided as a service and makes no claims of efficacy and/or safety of this product.			
β-Myrcene	21500	2.15		Results are applicable only for the sample(s) analyzed and for the specific			
cis-Nerolidol	-	-		analysis conducted. This report is for informational purposes			
trans-Nerolidol	352	0.0352		only and should not be used to diagnose,			
α-Ocimene	-	-		treat, or prevent any medical-related symptoms.			
β-Ocimene	4040	0.404		The statements and results herein have			
α-Pinene	2540	0.254	<u> </u>	not been approved and/or endorsed by			
β-Pinene	4180	0.418		the FDA.			
α-Terpinene	1650	0.165					
γ-Terpinene	1220	0.122					
Terpinolene	29200	2.92		"-" Not detected above RL.			

Reporting Limit (µg/g): 85.9



TOTAL

85900

8.59











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Cultivar (Strain) or Sample	Date Completed: 12/21/2022		

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

Analysis Date/Time: 12/16/2022 0247 Method: HS/G		Method: HS/GC/MS			D	Deviations from SOP:				
Analyst: KF				I	nstrument: Agilent 78	90/5975	5 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)	LOQ (µg/g)	<u>Action</u> <u>Level</u> (µg/g)	
Acetone (67-64-1)	-	39.6	79.3	5000	n-Heptane (142-82-5)	-	39.6	79.3	5000	
Acetonitrile (75-5-8)	-	39.6	79.3	410	n-Hexane (110-54-3)	-	13.9	27.8	290	
Benzene (71-43-2)	-	0.396	0.793	2	Isobutane (75-28-5)	-	39.6	79.3	5000	
n-Butane (106-97-2)	1530	39.6	79.3	5000	Isopropanol (67-63-0)	-	39.6	79.3	5000	
1-Butanol (71-36-3)	-	39.6	79.3	5000	Isopropyl acetate	_	39.6	79.3	5000	
2-Butanol (78-92-2)	-	39.6	79.3	5000	(108-21-4)		55.0	75.5	5000	
2-Butanone (78-93-3)	-	39.6	79.3	5000	Isopropyl benzene (98-82-8)	-	3.96	7.93	70	
Cyclohexane (110-82-7)	-	39.6	79.3	3880	(90-02-0) Methanol (67-56-1)		39.6	79.3	3000	Color Key
1,2-Dimethoxyethane	_	3.96	7.93	100	· · · ·	-				
(110-71-4) N,N-Dimethylacetamide		39.6	79.3	100	2-Methylbutane (78-78-4) Methylene chloride (75-9-2)		39.6 39.6	79.3 79.3	5000 600	RESULT < AL RESULT > AL
(127-19-5) 2,2-Dimethylbutane		13.9	27.8	290	(73-3-2) 2-Methylpentane (107-83-5)	-	13.9	27.8	290	"DET" detected less than LOQ
(75-83-2)		15.5	27.0	250	(107-03-3) 3-Methylpentane (96-10-0)		13.9	27.8	290	"-" not detected above
2,3-Dimethylbutane	-	13.9	27.8	290	51 ()			27.8 79.3		LOD
(79-29-8) N,N-Dimethylformamide					n-Pentane (109-66-0)		39.6 39.6	79.3 79.3	5000 5000	-
(68-12-2)	-	39.6	79.3	880	1-Pentanol (71-41-0)	-				"*" - o,m,p-Xylene and
Dimethylsulfoxide (67-68-5)	-	39.6	79.3	5000	n-Propane (74-98-6) 1-Propanol (71-23-8)	-	39.6 39.6	79.3 79.3	5000 5000	Ethylbenzene Action levels are
1,4-Dioxane (123-91-1)	-	39.6	79.3	380	Pyridine (110-86-1)	-	13.9	27.8	200	referenced from the State of
Ethanol (64-17-5)	-	39.6	79.3	5000	Tetrahydrofuran (109-99-9)	-	39.6	79.3	720	Arkansas
2-Ethoxyethanol (110-80-5)	-	13.9	27.8	160	Tetramethylene sulfone	_	13.9	27.8	160	MMJ testing
Ethyl ether (60-29-7)	-	39.6	79.3	5000	(126-33-0)		39.6	79.3	890	guidelines.
Ethyl acetate (141-78-6)	-	39.6	79.3	5000	Toluene (108-88-3)		39.6	79.3 79.3	2170	A value of "-"
Ethyl benzene (100-41-4)	_	39.6	79.3	2170	o-Xylene (95-47-6) m,p-Xylene (108-38-3 or		39.0	79.5	2170	for the action level
Ethylene glycol (107-21-1)	_	39.6	79.3	620	106-42-3)	-	39.6	79.3	2170	means that analyte
Ethylene oxide (75-21-8)	-	3.96	7.93	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	nide, ACN		Ethylene glycol	1	,2-Ethanediol			
1-Butanol		n-Butanol,	Butyl Alcol	hol	Isobutane	2	-Methylpropa	ne		
2-Butanol		sec-Butyl a	5		Isopropanol		-Propanol, IP/			
2-Butanone		Methyl ethy	yl ketone, N	1EK	Isopropyl Acetate	А	cetic acid iso	propyl este		
1,2-Dimethoxyethane		Monoglym			Methanol		fethyl alcohol			
2,3-Dimethylbutane		Neohexane			2-Methylbutane		sopentane			
2,3-Dimethylbutane		Diisopropy	1		Methylene chloride)ichlorometha	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		ΉF			
Ethyl acetate		EtOAc	,		Tetramethylene sulfone		ulfolane			
Ethyl benzene		Phenyletha	ne		Xylene		imethylbenze	ne		

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	Date Completed: 12/21/2022		

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 12/19/2022 1733 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> 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)	-	0.0190	0.152	0.5	Kresoxim-methyl	_	0.0190	0.152	0.4	
Acephate (30560-19-1)	-	0.0190	0.152	0.4	(143390-89-0)					
Acequinocyl (57960-19-7)	-	0.0190	0.152	2	Malathion (121-75-5)	-	0.0190	0.152	0.2	
Acetamiprid	_	0.0190	0.152	0.2	Metalaxyl (57837-19-1)	-	0.0190	0.152	0.2	Color Key
(135410-20-7)					Methiocarb (2032-65-7)	-	0.0190	0.152	0.2	<u>Color Key</u>
Aldicarb (116-06-3)	-	0.0190	0.152	0.4	Methomyl (16752-77-5)	-	0.0190	0.152	0.4	RESULT < AL
Azoxystrobin (131860-33-8)	-	0.0190	0.152	0.2	Methyl parathion (298-0-0)	-	0.0190	0.152	0.2	RESULT > AL
Bifenazate (149877-41-8)	-	0.0190	0.152	0.2	MGK 264 (113-48-4)	-	0.0190	0.152	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3)	-	0.0190	0.152	0.2	Myclobutanil	_	0.0190	0.152	0.2	"-" not detected above
Boscalid (188425-85-6)	-	0.0190	0.152	0.4	(88671-89-0)		0.0400	0.150	o -	LOD
Carbaryl (63-25-2)	-	0.0190	0.152	0.2	Naled (300-76-5)	-	0.0190	0.152	0.5	Permethrins measured as the
Carbofuran (1563-66-2)	-	0.0190	0.152	0.2	Oxamyl (23135-22-0)	-	0.0190	0.152	1	cumulative residue of the <i>cis</i> - and
Chlorantraniliprole (800008-45-7)	-	0.0190	0.152	0.2	Paclobutrazol (76738-62-0)	-	0.0190	0.152	0.4	trans- permethrin isomers.
Chlorfenapyr		0.0190	0.152	1	Permethrins (52645-53-1)	-	0.0190	0.152	0.2	Pyrethrins measured as the cumulative residue of the
(122453-73-0)		0.0190	0.152	1	¹ Phosmet (732-11-6)	-	0.0190	0.152	0.2	pyrethrin I, cinerin I, and jasmolin
Chlorpyrifos (2921-88-2)	-	0.0190	0.152	0.2	Piperonyl butoxide	_	0.0190	0.152	2	I isomers.
Clofentezine (74115-24-5)	-	0.0190	0.152	0.2	(51-03-6)		0.0100	0.152	0.2	Action levels are referenced from
Cyfluthrin (68359-37-5)	-	0.0190	0.152	1	Prallethrins (2331-36-9)	-	0.0190	0.152	0.2	the
Cypermethrin (52315-07-8)	-	0.0190	0.152	1	Propiconazole (60207-90-1))	-	0.0190	0.152	0.4	State of Arkansas MMJ testing guidelines.
Daminozide (1596-84-5)	-	0.0190	0.152	1	Propoxur (114-26-1)	-	0.0190	0.152	0.2	A value of "-" for the action level
DDVP (62-73-7)	-	0.0190	0.152	0.1	Pyrethrins (8003-34-7)	-	0.0190	0.152	1	means that analyte is not
Diazinon (333-41-5)	-	0.0190	0.152	0.2	Pyridaben (96489-71-3)	-	0.0190	0.152	0.2	currently regulated by the
Dimethoate (60-51-5)	-	0.0190	0.152	0.2	Spinosad (168316-95-8)	-	0.0190	0.152	0.2	regulations referenced above.
Ethoprophos (13194-48-4)	-	0.0190	0.152	0.2	Spiromesifen (283594-90-1)	-	0.0190	0.152	0.2	Disclaimer: This information is
Etofenprox (80844-07-1)	-	0.0190	0.152	0.4	(2033)4-50-1) Spirotetramat					provided as a service and makes no claims of efficacy and/or safety
Etoxazole (153233-91-1)	-	0.0190	0.152	0.2	(203313-25-1)	-	0.0190	0.152	0.2	of this product. Results are
Fenoxycarb (72490-01-8)	-	0.0190	0.152	0.2	Spiroxamine		0.0190	0.152	0.4	applicable only for the sample(s) analyzed and for the specific
(E)-Fenpyroximate (134098-61-6)	-	0.0190	0.152	0.4	(118134-30-8) Tebuconazole	-				analysis conducted. This report is for informational purposes only
Fipronil (120068-37-3)	-	0.0190	0.152	0.4	(80443-41-0)	-	0.0190	0.152	0.4	and should not be used to
Flonicamid (158062-67-0)	-	0.0190	0.152	1	Thiacloprid	-	0.0190	0.152	0.2	diagnose, treat, or prevent any medical-related symptoms. The
Fludioxinil (131341-86-1)	-	0.0190	0.152	0.4	(111988-49-9)					statements and results herein have
Hexythiazox (78587-05-0)	-	0.0190	0.152	1	Thiamethoxam (153719-23-4)	-	0.0190	0.152	0.2	not been approved and/or
Imazalil (35554-44-0)	-	0.0190	0.152	0.2	Trifloxystrobin					endorsed by the FDA.
Imidacloprid (138261-41-3)	-	0.0190	0.152	0.4	(141517-21-7)	-	0.0190	0.152	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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Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 12/14/2022
License: ADH 113	License: 00065C	E20221128O43HLR01	Date Received: 12/14/2022
Cultivar (Strain) or Sample	Date Completed: 12/21/2022		

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

Analysis Date/Time: 12/21/2022 0805 (ICP/OES) Analysis Date/Time: - (DMA) Analyst: KF		Method: ICP/OES Instrument: Agilent 720-ES		Deviations from SOP: ES None	
<u>Heavy Metal</u>	<u>Result</u> (µg/kg)	<u>LOD</u> (µg/kg)	LOQ (µg/kg)	<u>Action Level</u> (µg/kg)	
Arsenic (As)	-	56.4	89.2	200	
Cadmium (Cd)	-	56.4	89.2	200	
Lead (Pb)	-	56.4	89.2	500	
Mercury (Hg)	-	56.4	89.2	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: SA33699)

Testing Location:	Customer ID: 2168	Sample ID: SA33699	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13241008157	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 12/14/2022
License: ADA 05_H273	License: 00065C	E20221128O43HLR01	Date Received: 12/14/2022
Cultivar (Strain) or Sample D	Date Completed: 12/21/2022		

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 2022121 Analyst: PW		ardy Diagnostics CompactDry t: Thermo Incubator	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	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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