

License: ADH 113





E20240718RAWFSO20



Date Received: 07/18/2024

SUMMARY OF ANALYSIS (SAMPLE ID: SA38428)

Order ID: OR11107 **Testing Location:** Customer ID: 2168 Sample Type: Primary Arkansas River Valley Relief MIPS Lot Number: Matrix: Concentrate

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13355971595 Mass: 4ea Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 07/18/2024

Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil 1g Applicator **Date Completed:** 07/22/2024

*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 (%)** PASS/FAIL Water Activity (aw) **PASS** Not Tested Not Tested

Moisture content/water activity action levels are referenced from the State of Arkansas MMJ testing guidelines.

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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>	mg/g
Δ9-7	ГНС	74.6	746
CI	3G	3.03	30.3
CI	BC	0.568	5.68
TOTA	L CBD	0.208	2.08
TOTA	LTHC	74.6	746
TOTAL CAN	NABINOIDS	79.4	794
<u>Terpene</u>	s (Top 5)	(%)	ħā∖ā
β-Caryo	phyllene	1.80	18000
β-Му	rcene	0.971	9710
Lina	lool	0.826	8260
d-Lim	onene	0.824	8240
α-Hun	nulene	0.593	5930
TOTAL T	ERPENES	6.54	65400
Contaminants	PASS/FAIL	Sample Pictur	e Upon Receipt

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: SA38428)

Testing Location:Customer ID: 2168Order ID: OR11107Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Concentrate232 S. Broadview St.5601 Old Greenwood Rd Suite CM00065C13355971595Mass: 4ea

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 07/18/2024

License: ADH 113 License: 00065C E20240718RAWFSO20 **Date Received:** 07/18/2024 **Cultivar (Strain) or Sample Description:** ArkanRAW Full Spectrum Raw Cannabis Oil 1g Applicator **Date Completed:** 07/22/2024

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

Analysis Date/Time: 07/19/2024 1406 Method: HPLC/DAD
Analyst: PW Instrument: Agilent 1100

<u>Cannabinoid</u>	Result (%)	Result (mg/g)	LOD (mg/g)	LOQ (mg/g)	Result (mg/ mL)	<u>Per</u> Serving (mg)	Per Unit (mg)
CBC	0.568	5.68	0.108	0.253	-	5.68	5.68
CBCA	ND	ND	0.337	0.786	-	-	-
CBD	0.208	2.08	0.765	1.79	-	2.08	2.08
CBDA	ND	ND	0.282	0.658	-	-	-
CBDV	ND	ND	0.123	0.286	-	-	-
CBDVA	ND	ND	0.327	0.764	-	-	-
CBG	3.03	30.3	0.497	1.16	-	30.3	30.3
CBGA	0.198	1.98	0.704	0.831	-	1.98	1.98
CBL	ND	ND	0.574	1.34	-	-	-
CBN	0.245	2.45	0.264	0.616	-	2.45	2.45
CBNA	ND	ND	0.285	0.663	-	-	-
Δ9-ΤΗС	74.6	746	0.316	0.737	-	746	746
Δ8-ΤΗС	ND	ND	0.493	1.15	-	-	-
THCA	ND	ND	0.171	0.401	-	-	-
THCV	0.550	5.50	0.411	0.958	-	5.50	5.50
THCVA	ND	ND	0.131	0.306	-	-	-
TOTAL	79.4	794			-	794	794
TOTAL CBC	0.568	5.68		,,	-	5.68	5.68
TOTAL CBD	0.208	2.08			-	2.08	2.08
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	3.20	32.0			-	32.0	32.0
TOTAL CBN	0.245	2.45			-	2.45	2.45
TOTAL THC	74.6	746			-	746	746
TOTAL THCV	0.550	5.50			-	5.50	5.50

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.







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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA38428)

Customer ID: 2168 Order ID: OR11107 Sample Type: Primary **Testing Location:** Arkansas River Valley Relief MIPS Lot Number: 232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13355971595 Mass: 4ea

Fort Smith, AR 72903 Greenbrier, AR 72058 **Production Run:** License: 00065C E20240718RAWFSO20

Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil 1g Applicator

Matrix: Concentrate

Date Collected: 07/18/2024 **Date Received:** 07/18/2024

Date Completed: 07/22/2024

TERPENOID PROFILE

Analysis Date/Time:07/19/2024 1442 Method: GC/MS **Deviations from SOP: Analyst:** KF Instrument: Agilent 7890/5975 None

Result	Result (%)
	(%) 0.219
	0.0382
	0.0362
18000	1.80
ND	-
ND	-
162	0.0162
137	0.0137
483	0.0483
5930	0.593
<loq< td=""><td>-</td></loq<>	-
8240	0.824
8260	0.826
9710	0.971
444	0.0444
3330	0.333
140	0.0140
1720	0.172
2310	0.231
1380	0.138
187	0.0187
220	0.0220
2180	0.218
65400	6.54
	(µg/g) 2190 382 55.4 18000 ND ND 162 137 483 5930 <loq 1380="" 140="" 1720="" 187="" 2180<="" 220="" 2310="" 3330="" 444="" 8240="" 8260="" 9710="" td=""></loq>

Reporting Limit (µg/g): 47.6

"-" Not detected above LOD.













CERTIFICATE OF ANALYSIS (SAMPLE ID: SA38428)

Testing Location:Customer ID: 2168Order ID: OR11107Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Concentrate

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13355971595 **Mass:** 4ea

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 07/18/2024

License: ADH 113 License: 00065C E20240718RAWFSO20 **Date Received:** 07/18/2024 **Cultivar (Strain) or Sample Description:** ArkanRAW Full Spectrum Raw Cannabis Oil 1g Applicator **Date Completed:** 07/22/2024

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

Analysis Date/Time: 07/19/2024 0406 Method: HS/GC/MS Deviations from SOP:

Analyst: KF Instrument: Agilent 7890/5975 None

Solvent	Result (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> <u>Level</u> (μg/g)	Solvent	Result (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> <u>Level</u> (μg/g)
Acetone (67-64-1)	-	39.5	79.1	5000	n-Heptane (142-82-5)	-	39.5	79.1	5000
Acetonitrile (75-5-8)	-	39.5	79.1	410	n-Hexane (110-54-3)	-	13.8	27.7	290
Benzene (71-43-2)	-	0.395	0.791	2	Isobutane (75-28-5)	-	39.5	79.1	5000
n-Butane (106-97-2)	1810	39.5	79.1	5000	Isopropanol (67-63-0)	-	39.5	79.1	5000
1-Butanol (71-36-3)	-	39.5	79.1	5000	Isopropyl acetate	_	39.5	79.1	5000
2-Butanol (78-92-2)	-	39.5	79.1	5000	(108-21-4)		55.5	7511	5000
2-Butanone (78-93-3)	-	39.5	79.1	5000	Isopropyl benzene (98-82-8)	-	3.95	7.91	70
Cyclohexane (110-82-7)	-	39.5	79.1	3880	Methanol (67-56-1)	_	39.5	79.1	3000
1,2-Dimethoxyethane (110-71-4)	-	3.95	7.91	100	2-Methylbutane (78-78-4)	-	39.5	79.1	5000
N,N-Dimethylacetamide (127-19-5)	-	39.5	79.1	1090	Methylene chloride (75-9-2)	-	39.5	79.1	600
2,2-Dimethylbutane (75-83-2)	-	13.8	27.7	290	2-Methylpentane (107-83-5)	-	13.8	27.7	290
2,3-Dimethylbutane (79-29-8)	-	13.8	27.7	290	3-Methylpentane (96-10-0) n-Pentane (109-66-0)	-	13.8 39.5	27.7 79.1	290 5000
N,N-Dimethylformamide (68-12-2)	-	39.5	79.1	880	1-Pentanol (71-41-0) n-Propane (74-98-6)	-	39.5 39.5	79.1 79.1	5000 5000
Dimethylsulfoxide (67-68-5)	-	39.5	79.1	5000	1-Propanol (71-23-8)	-	39.5	79.1	5000
1,4-Dioxane (123-91-1)	_	39.5	79.1	380	Pyridine (110-86-1)	-	13.8	27.7	200
Ethanol (64-17-5)	DET	39.5	79.1	5000	Tetrahydrofuran (109-99-9)	-	39.5	79.1	720
2-Ethoxyethanol (110-80-5)	-	13.8	27.7	160	Tetramethylene sulfone (126-33-0)	-	13.8	27.7	160
Ethyl ether (60-29-7)	-	39.5	79.1	5000	Toluene (108-88-3)	_	39.5	79.1	890
Ethyl acetate (141-78-6)	-	39.5	79.1	5000	o-Xylene (95-47-6)	_	39.5	79.1	2170
Ethyl benzene (100-41-4)	-	39.5	79.1	2170	m,p-Xylene (108-38-3 or				
Ethylene glycol (107-21-1)	-	39.5	79.1	620	106-42-3)	-	39.5	79.1	2170
Ethylene oxide (75-21-8)	-	3.95	7.91	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170

Solvent	Synonym(s)	Solvent	Synonym(s)
Solvent	<u>Synonym(s)</u>	Solvent	<u>Symonym(s)</u>
Acetonitrile	Methyl Cyanide, ACN	Ethylene glycol	1,2-Ethanediol
1-Butanol	n-Butanol, Butyl Alcohol	Isobutane	2-Methylpropane
2-Butanol	sec-Butyl alcohol	Isopropanol	2-Propanol, IPA
2-Butanone	Methyl ethyl ketone, MEK	Isopropyl Acetate	Acetic acid isopropyl ester
1,2-Dimethoxyethane	Monoglyme	Methanol	Methyl alcohol
2,3-Dimethylbutane	Neohexane	2-Methylbutane	Isopentane
2,3-Dimethylbutane	Diisopropyl	Methylene chloride	Dichloromethane
N,N-Dimethylformamide	DMF	2-Methylpentane	Isohexane
Dimethysufoxide	DMSO	1-Pentanol	n-Amyl alcohol
2-Ethoxyethanol	Cellosolve, Ethyl glycol	1-Propanol	Propyl alcohol
Ethyl ether	Diethyl ether, Ether	Tetrahydrofuran	THF
Ethyl acetate	EtOAc	Tetramethylene sulfone	Sulfolane
Ethyl benzene	Phenylethane	Xylene	Dimethylbenzene



Color Key

RESULT < AL RESULT > AL

"DET" detected less than LOQ

"-" not detected above LOD

"*" - o,m,p-Xylene and Ethylbenzene

Action levels 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.

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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E20240718RAWFSO20



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA38428)

Testing Location:Customer ID: 2168Order ID: OR11107Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Concentrate

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13355971595 **Mass:** 4ea

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 07/18/2024

Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil 1g Applicator Date Completed: 07/22/2024

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 07/19/2024 1351 Method: LC/MS/MS Deviations from SOP:

Analyst: KF Instrument: Shimadzu LC-8050 None

License: 00065C

<u>Pesticide</u>	Result (µg/g)	<u>LOD</u> (µg/g)	LOQ (µg/g)	<u>Action</u> <u>Level</u> (μg/g)	<u>Pesticide</u>	Result (µg/g)	<u>LOD</u> (µg/g)	LOQ (µg/g)	<u>Action</u> <u>Level</u> (μg/g)
Abamectin (71751-41-2)	-	0.0472	0.378	0.5	Kresoxim-methyl	_	0.0472	0.378	0.4
Acephate (30560-19-1)	-	0.0472	0.378	0.4	(143390-89-0)				
Acequinocyl (57960-19-7)	-	0.0472	0.378	2	Malathion (121-75-5)	-	0.0472	0.378	0.2
Acetamiprid (135410-20-7)	-	0.0472	0.378	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.0472 0.0472	0.378 0.378	0.2 0.2
Aldicarb (116-06-3)	-	0.0472	0.378	0.4	Methomyl (16752-77-5)	-	0.0472	0.378	0.4
Azoxystrobin (131860-33-8)	-	0.0472	0.378	0.2	Methyl parathion (298-0-0)	-	0.0472	0.378	0.2
Bifenazate (149877-41-8)	-	0.0472	0.378	0.2	MGK 264 (113-48-4)	-	0.0472	0.378	0.2
Bifenthrin (82657-04-3)	-	0.0472	0.378	0.2	Myclobutanil		0.0472	0.378	0.2
Boscalid (188425-85-6)	-	0.0472	0.378	0.4	(88671-89-0)		0.0472	0.570	0.2
Carbaryl (63-25-2)	-	0.0472	0.378	0.2	Naled (300-76-5)	-	0.0472	0.378	0.5
Carbofuran (1563-66-2)	-	0.0472	0.378	0.2	Oxamyl (23135-22-0)	-	0.0472	0.378	1
Chlorantraniliprole (800008-45-7)	-	0.0472	0.378	0.2	Paclobutrazol (76738-62-0)	-	0.0472	0.378	0.4
Chlorfenapyr		0.0472	0.378	1	Permethrins (52645-53-1)	-	0.0472	0.378	0.2
(122453-73-0)	_	0.0472	0.376	1	Phosmet (732-11-6)	-	0.0472	0.378	0.2
Chlorpyrifos (2921-88-2)	-	0.0472	0.378	0.2	Piperonyl butoxide	_	0.0472	0.378	2
Clofentezine (74115-24-5)	-	0.0472	0.378	0.2	(51-03-6)		0.0450	0.050	0.0
Cyfluthrin (68359-37-5)	-	0.0472	0.378	1	Prallethrins (2331-36-9)		0.0472	0.378	0.2
Cypermethrin (52315-07-8)	-	0.0472	0.378	1	Propiconazole (60207-90-1))	-	0.0472	0.378	0.4
Daminozide (1596-84-5)	-	0.0472	0.378	1	Propoxur (114-26-1)	-	0.0472	0.378	0.2
DDVP (62-73-7)	-	0.0472	0.378	0.1	Pyrethrins (8003-34-7)	-	0.0472	0.378	1
Diazinon (333-41-5)	-	0.0472	0.378	0.2	Pyridaben (96489-71-3)	-	0.0472	0.378	0.2
Dimethoate (60-51-5)	-	0.0472	0.378	0.2	Spinosad (168316-95-8)	-	0.0472	0.378	0.2
Ethoprophos (13194-48-4)	-	0.0472	0.378	0.2	Spiromesifen	2.0	0.0472	0.378	0.2
Etofenprox (80844-07-1)	-	0.0472	0.378	0.4	(283594-90-1) Spirotetramat				
Etoxazole (153233-91-1)	-	0.0472	0.378	0.2	(203313-25-1)	-	0.0472	0.378	0.2
Fenoxycarb (72490-01-8)	-	0.0472	0.378	0.2	Spiroxamine		0.0450	0.050	
(E)-Fenpyroximate (134098-61-6)	-	0.0472	0.378	0.4	(118134-30-8) Tebuconazole	-	0.0472	0.378	0.4
Fipronil (120068-37-3)	_	0.0472	0.378	0.4	(80443-41-0)	-	0.0472	0.378	0.4
Flonicamid (158062-67-0)	-	0.0472	0.378	1	Thiacloprid		0.0472	0.378	0.2
Fludioxinil (131341-86-1)	-	0.0472	0.378	0.4	(111988-49-9)		0.04/2	0.370	0.2
Hexythiazox (78587-05-0)	-	0.0472	0.378	1	Thiamethoxam	-	0.0472	0.378	0.2
Imazalil (35554-44-0)	-	0.0472	0.378	0.2	(153719-23-4) Trifloxystrobin				
Imidacloprid (138261-41-3)	-	0.0472	0.378	0.4	(141517-21-7)	-	0.0472	0.378	0.2



Color Key

Date Received: 07/18/2024

RESULT < AL

"DET" detected less than LOQ

"-" not detected above LOD

Permethrins measured as the cumulative residue of the *cis*- and *trans*- permethrin isomers.

Pyrethrins measured as the cumulative residue of the pyrethrin I, cinerin I, and jasmolin I isomers.

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











E20240718RAWFSO20



Date Received: 07/18/2024

CERTIFICATE OF ANALYSIS (SAMPLE ID: SA38428)

 Testing Location:
 Customer ID: 2168
 Order ID: OR11107
 Sample Type: Primary

 Arkansas
 River Valley Relief MIPS
 Lot Number:
 Matrix: Concentrate

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13355971595 Mass: 4ea

Greenbrier, AR 72058 Fort Smith, AR 72903 Production Run: Date Collected: 07/18/2024

Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil 1g Applicator Date Completed: 07/22/2024

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

Analysis Date/Time: 07/18/2024 2119 Method: ICP/MS Deviations from SOP:

Instrument: Agilent 7500ce None

Analyst: KF

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<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)	-	54. 5	86.3	200
Cadmium (Cd)	-	54. 5	86.3	200
Lead (Pb)	-	54. 5	86.3	500
Mercury (Hg)	-	54.5	86.3	100

License: 00065C



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

RESULT < AL
RESULT > AL

"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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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA38428)

Testing Location:Customer ID: 2168Sample ID: SA38428Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Concentrate

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13355971595 **Mass:** 4ea

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 07/18/2024

Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil 1g Applicator **Date Completed:** 07/22/2024

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 07/19/2024 1133 Method: Hardy Diagnostics CompactDry Deviations from SOP:

Analyst: PW Instrument: Thermo Incubator None

Bacteria/Microbe	Result (CFU/g)	Action Level (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

Color Key

RESULT < AL

RESULT > AL

Reporting Limit (CFU/g)

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.

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.



