







SUMMARY OF ANALYSIS (SAMPLE ID: SA33963)

Testing Location:Customer ID: 2168Order ID: OR10230Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Concentrate

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13249703028 Mass: 4g

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 12/21/2022 License: ADH 113 License: 00065C E20221215TGKC01 **Date Received:** 12/21/2022 **Cultivar (Strain) or Sample Description:** Triangle Kush Indica Live Resin .5g Cart **Date Completed:** 12/28/2022

*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/FAILNot TestedNot TestedPASS

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.

Cannabino	oids (Top 3)	<u>(%)</u>	mg/g
TH	CA	33.0	330
Δ9-7	ГНС	29.4	294
CB	GA	2.97	29.7
TOTA	L CBD	0.213	2.13
TOTA	LTHC	58.4	584
TOTAL CAN	NABINOIDS	69.6	696
<u>Terpene</u>	s (Top 5)	<u>(%)</u>	µg/g
β-Му	rcene	1.89	18900
d-Lim	onene	1.00	10000
β-Caryo	phyllene	0.465	4650
Lina	alool	0.314	3140
α-Hun	nulene	0.190	1900
TOTAL T	ERPENES	4.47	44700
Contaminants	PASS/FAIL	Sample Pictur	re 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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License: ADH 113





E20221215TGKC01



Date Received: 12/21/2022

CERTIFICATE OF ANALYSIS (SAMPLE ID: SA33963)

Order ID: OR10230 **Testing Location:** Customer ID: 2168 Sample Type: Primary River Valley Relief MIPS Lot Number: Matrix: Concentrate Arkansas

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13249703028 Mass: 4g Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 12/21/2022

Cultivar (Strain) or Sample Description: Triangle Kush Indica Live Resin .5g Cart **Date Completed:** 12/28/2022

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

Analysis Date/Time: 12/27/2022 1612 Method: HPLC/DAD Moisture Content (%): -**Analyst: PW Instrument:** Agilent 1100 Water Activity (aw): -

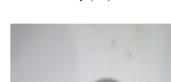
<u>Cannabinoid</u>	Result (%)	<u>UM</u> (+/-%)	Result (mg/g)	LOQ (mg/g)	Result (mg/ mL)	Per Serving (mg)	Per Unit (mg)
CBC	-	-	-	0.377	-	-	-
CBCA	1.37	0.123	13.7	0.377	-	13.7	13.7
CBD	-	-	-	0.377	-	-	-
CBDA	0.243	0.0219	2.43	0.377	-	2.43	2.43
CBDV	-	-	-	0.377	-	-	-
CBDVA	-	-	-	0.377	-	-	-
CBG	1.29	0.116	12.9	0.377	-	12.9	12.9
CBGA	2.97	0.268	29.7	0.377	-	29.7	29.7
CBL	-	-	-	0.377	-	-	-
CBN	0.149	0.0135	1.49	0.377	-	1.49	1.49
CBNA	-	-	-	0.377	-	-	-
Δ9-ΤΗС	29.4	2.65	294	0.377	-	294	294
Δ8-ΤΗС	-	-	-	0.377	-	-	-
THCA	33.0	2.97	330	0.377	-	330	330
THCV	0.592	0.0533	5.92	0.377	-	5.92	5.92
THCVA	0.525	0.0473	5.25	0.377	-	5.25	5.25
TOTAL	69.6	6.26	696		_	696	696
TOTAL CBC	1.20	0.108	12.0		· -	12.0	12.0
TOTAL CBD	0.213	0.0192	2.13		-	2.13	2.13
TOTAL CBDV	-	-	-		-	-	-
TOTAL CBG	3.90	0.351	39.0		-	39.0	39.0
TOTAL CBN	0.149	0.0135	1.49		-	1.49	1.49
TOTAL THC	58.4	5.25	584		-	584	584
TOTAL THCV	1.05	0.0943	10.5		-	10.5	10.5

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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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SERVING MASS (g): 1.00 **SERVINGS/UNIT: 1**

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 = $(CBCA \times 0.877) + CBC$ Total CBD = $(CBDA \times 0.877) + CBD$ Total CBDV = $(CBDVA \times 0.867) + CBDV$ Total CBG = $(CBGA \times 0.878) + CBG$ Total CBN = $(CBNA \times 0.876) + CBN$ 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.





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

[&]quot;-" Not detected above LOD.









CERTIFICATE OF ANALYSIS (SAMPLE ID: SA33963)

Testing Location:Customer ID: 2168Order ID: OR10230Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Concentrate232 S. Broadview St.5601 Old Greenwood Rd Suite CM00065C13249703028Mass: 4g

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 12/21/2022 License: ADH 113 License: 00065C E20221215TGKC01 **Date Received:** 12/21/2022

Cultivar (Strain) or Sample Description: Triangle Kush Indica Live Resin .5g Cart **Date Completed:** 12/28/2022

TERPENOID PROFILE

Analysis Date/Time:12/28/2022 0903 Method: GC/MS Deviations from SOP:
Analyst: KF Instrument: Agilent 7890/5975 None

Analyst: KF		Inst	Instrument: Agilent 7890/5975			
<u>Terpene</u>	<u>Result</u> (μg/g)	Result (%)				
α-Bisabolol	1110	0.111				
Camphene	375	0.0375				
δ-3-Carene	-	-				
β-Caryophyllene	4650	0.465				
Caryophyllene oxide	327	0.0327	1			
p-Cymene	-	-				
Eucalyptol	-	-				
Geraniol	313	0.0313				
Guaiol	-	-				
α-Humulene	1900	0.190				

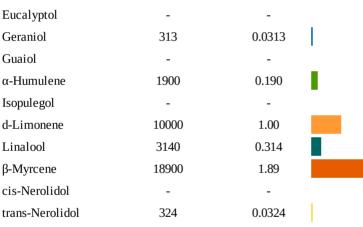
0.0407

0.0878

0.171

0.0246

0.0448



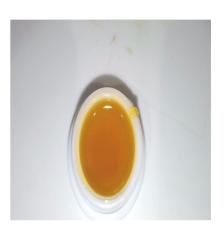
407

878

1710

246

448



Abbreviations: GC - Gas Chromatography, MS - Mass Spectrometry, RL - Reporting Limit

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"-" Not detected above RL

TOTAL 44700 4.47 Reporting Limit (μg/g): 34.1



α-Ocimene

β-Ocimene

α-Pinene

B-Pinene

α-Terpinene γ-Terpinene

Terpinolene





Greenbrier, AR 72058





Production Run:



Date Collected: 12/21/2022

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License: ADH 113 License: 00065C E20221215TGKC01 **Date Received:** 12/21/2022

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

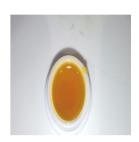
Analysis Date/Time: 12/27/2022 1643 Method: HS/GC/MS Deviations from SOP:

Analyst: KF Instrument: Agilent 7890/5975 None

Fort Smith, AR 72903

Solvent	<u>Result</u> (μg/g)	<u>LOD</u> (µg/g)	LOQ (µg/g)	Action Level (µg/g)	Solvent	Result (µg/g)	<u>LOD</u> (µg/g)	LOQ (µg/g)	<u>Action</u> <u>Level</u> (μg/g)
Acetone (67-64-1)	-	38.1	76.2	5000	n-Heptane (142-82-5)	-	38.1	76.2	5000
Acetonitrile (75-5-8)	-	38.1	76.2	410	n-Hexane (110-54-3)	-	13.3	26.7	290
Benzene (71-43-2)	-	0.381	0.762	2	Isobutane (75-28-5)	-	38.1	76.2	5000
n-Butane (106-97-2)	747	38.1	76.2	5000	Isopropanol (67-63-0)	-	38.1	76.2	5000
1-Butanol (71-36-3)	-	38.1	76.2	5000	Isopropyl acetate		38.1	76.2	5000
2-Butanol (78-92-2)	-	38.1	76.2	5000	(108-21-4)		30.1	70.2	3000
2-Butanone (78-93-3)	-	38.1	76.2	5000	Isopropyl benzene (98-82-8)	_	3.81	7.62	70
Cyclohexane (110-82-7)	-	38.1	76.2	3880	Methanol (67-56-1)	_	38.1	76.2	3000
1,2-Dimethoxyethane (110-71-4)	-	3.81	7.62	100	2-Methylbutane (78-78-4)	-	38.1	76.2	5000
N,N-Dimethylacetamide (127-19-5)	-	38.1	76.2	1090	Methylene chloride (75-9-2)	-	38.1	76.2	600
2,2-Dimethylbutane (75-83-2)	-	13.3	26.7	290	2-Methylpentane (107-83-5)	-	13.3	26.7	290
2,3-Dimethylbutane		13.3	26.7	290	3-Methylpentane (96-10-0)	-	13.3	26.7	290
(79-29-8)	_	13.3	20./	290	n-Pentane (109-66-0)	-	38.1	76.2	5000
N,N-Dimethylformamide	_	38.1	76.2	880	1-Pentanol (71-41-0)	-	38.1	76.2	5000
(68-12-2)					n-Propane (74-98-6)	-	38.1	76.2	5000
Dimethylsulfoxide (67-68-5)	-	38.1	76.2	5000	1-Propanol (71-23-8)	-	38.1	76.2	5000
1,4-Dioxane (123-91-1)	_	38.1	76.2	380	Pyridine (110-86-1)	-	13.3	26.7	200
Ethanol (64-17-5)	_	38.1	76.2	5000	Tetrahydrofuran (109-99-9)	-	38.1	76.2	720
2-Ethoxyethanol (110-80-5)	-	13.3	26.7	160	Tetramethylene sulfone (126-33-0)	-	13.3	26.7	160
Ethyl ether (60-29-7)	-	38.1	76.2	5000	Toluene (108-88-3)	_	38.1	76.2	890
Ethyl acetate (141-78-6)	-	38.1	76.2	5000	o-Xylene (95-47-6)	_	38.1	76.2	2170
Ethyl benzene (100-41-4)	-	38.1	76.2	2170	m,p-Xylene (108-38-3 or				
Ethylene glycol (107-21-1)	-	38.1	76.2	620	106-42-3)	-	38.1	76.2	2170
Ethylene oxide (75-21-8)	-	3.81	7.62	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170

Solvent	Synonym(s)	Solvent	Synonym(s)
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

"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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E20221215TGKC01



Date Received: 12/21/2022

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Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 12/21/2022

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PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 12/27/2022 1857 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)	Action Level (µg/g)	<u>Pesticide</u>	Result (μg/g)	<u>LOD</u> (µg/g)	LOQ (µg/g)	Action Level (µg/g)
Abamectin (71751-41-2)	-	0.0341	0.273	0.5	Kresoxim-methyl	2	0.0341	0.273	0.4
Acephate (30560-19-1)	-	0.0341	0.273	0.4	(143390-89-0)		0.0241	0.272	0.2
Acequinocyl (57960-19-7)	-	0.0341	0.273	2	Malathion (121-75-5)	-	0.0341	0.273	0.2
Acetamiprid (135410-20-7)	-	0.0341	0.273	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.0341 0.0341	0.273 0.273	0.2 0.2
Aldicarb (116-06-3)	-	0.0341	0.273	0.4	Methomyl (16752-77-5)	-	0.0341	0.273	0.4
Azoxystrobin (131860-33-8)	-	0.0341	0.273	0.2	Methyl parathion (298-0-0)	-	0.0341	0.273	0.2
Bifenazate (149877-41-8)	-	0.0341	0.273	0.2	MGK 264 (113-48-4)	-	0.0341	0.273	0.2
Bifenthrin (82657-04-3)	-	0.0341	0.273	0.2	Myclobutanil	_	0.0341	0.273	0.2
Boscalid (188425-85-6)	-	0.0341	0.273	0.4	(88671-89-0)				
Carbaryl (63-25-2)	-	0.0341	0.273	0.2	Naled (300-76-5)	-	0.0341	0.273	0.5
Carbofuran (1563-66-2)	-	0.0341	0.273	0.2	Oxamyl (23135-22-0)	-	0.0341	0.273	1
Chlorantraniliprole (800008-45-7)	-	0.0341	0.273	0.2	Paclobutrazol (76738-62-0)	-	0.0341	0.273	0.4
Chlorfenapyr (122453-73-0)	-	0.0341	0.273	1	Permethrins (52645-53-1) Phosmet (732-11-6)	-	0.0341 0.0341	0.273 0.273	0.2 0.2
Chlorpyrifos (2921-88-2)	-	0.0341	0.273	0.2	Piperonyl butoxide		0.0241	0.272	2
Clofentezine (74115-24-5)	_	0.0341	0.273	0.2	(51-03-6)	_	0.0341	0.273	2
Cyfluthrin (68359-37-5)	-	0.0341	0.273	1	Prallethrins (2331-36-9)	-	0.0341	0.273	0.2
Cypermethrin (52315-07-8)	-	0.0341	0.273	1	Propiconazole (60207-90-1))	-	0.0341	0.273	0.4
Daminozide (1596-84-5)	_	0.0341	0.273	1	Propoxur (114-26-1)	-	0.0341	0.273	0.2
DDVP (62-73-7)	_	0.0341	0.273	0.1	Pyrethrins (8003-34-7)	-	0.0341	0.273	1
Diazinon (333-41-5)	_	0.0341	0.273	0.2	Pyridaben (96489-71-3)	-	0.0341	0.273	0.2
Dimethoate (60-51-5)	-	0.0341	0.273	0.2	Spinosad (168316-95-8)	-	0.0341	0.273	0.2
Ethoprophos (13194-48-4)	-	0.0341	0.273	0.2	Spiromesifen		0.0341	0.273	0.2
Etofenprox (80844-07-1)	-	0.0341	0.273	0.4	(283594-90-1)				
Etoxazole (153233-91-1)	-	0.0341	0.273	0.2	Spirotetramat (203313-25-1)	-	0.0341	0.273	0.2
Fenoxycarb (72490-01-8)	-	0.0341	0.273	0.2	Spiroxamine				
(E)-Fenpyroximate (134098-61-6)	-	0.0341	0.273	0.4	(118134-30-8) Tebuconazole	-	0.0341	0.273	0.4
Fipronil (120068-37-3)	_	0.0341	0.273	0.4	(80443-41-0)	-	0.0341	0.273	0.4
Flonicamid (158062-67-0)	-	0.0341	0.273	1	Thiacloprid		0.0341	0.273	0.2
Fludioxinil (131341-86-1)	_	0.0341	0.273	0.4	(111988-49-9)	-	0.0341	0.273	0.2
Hexythiazox (78587-05-0)	-	0.0341	0.273	1	Thiamethoxam		0.0341	0.273	0.2
Imazalil (35554-44-0)	-	0.0341	0.273	0.2	(153719-23-4) Trifloxystrobin				
Imidacloprid (138261-41-3)	-	0.0341	0.273	0.4	(141517-21-7)	-	0.0341	0.273	0.2



Color Key

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		













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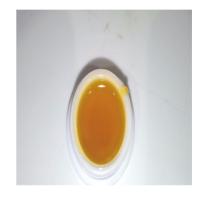
HEAVY METAL PROFILE (SOP: SOP-ICP-200.7)

Analysis Date/Time: 12/27/2022 1115 (ICP/OES) Method: ICP/OES Deviations from SOP:

Analysis Date/Time: - (DMA) Instrument: Agilent 720-ES None

Analyst: KF

<u>Heavy Metal</u>	<u>Result</u> (μg/kg)	<u>LOD</u> (µg/kg)	<u>LOQ</u> (µg/kg)	Action Level (µg/kg)
Arsenic (As)	-	55.8	88.3	200
Cadmium (Cd)	-	55.8	88.3	200
Lead (Pb)	-	55.8	88.3	500
Mercury (Hg)	-	55.8	88.3	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

RESULT < AL RESULT > AL

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"-" 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.







License: ADA 05 H273





E20221215TGKC01



Date Received: 12/21/2022

CERTIFICATE OF ANALYSIS (SAMPLE ID: SA33963)

Testing Location:Customer ID: 2168Sample ID: SA33963Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Concentrate

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13249703028 Mass: 4g
Greenbrier, AR 72058 Fort Smith, AR 72903 Production Run: Date Collected: 12/21/2022

Cultivar (Strain) or Sample Description: Triangle Kush Indica Live Resin .5g Cart **Date Completed:** 12/28/2022

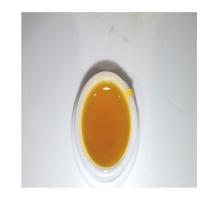
MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 12/23/2022 1030 **Method:** Hardy Diagnostics CompactDry **Deviations from SOP:**

Analyst: PW Instrument: Thermo Incubator None

Bacteria/Microbe	<u>Result</u> (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	1
Salmonella spp.	NT	-
Staphylococcus aureus	NT	-

License: 00065C



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.

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