

SUMMARY OF ANALYSIS (SAMPLE ID: SA36126)

Testing Location: Arkansas 232 S. Broadview St. Greenbrier, AR 72058 License: ADH 113	Customer ID: 2168 River Valley Relief MIPS 5601 Old Greenwood Rd Suite C Fort Smith, AR 72903 License: 00065C	Order ID: OR10628 Lot Number: M00065C13247672701 Production Run: E2023USKLHR01	Sample Type: Primary Matrix: Concentrate Mass: 4g Date Collected: 10/02/2023 Date Received: 10/03/2023
Cultivar (Strain) or Sample Description: Unkle Sink Hybrid Solventless Live Hash Rosin			Date Completed: 10/04/2023

*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 (%)

Not Tested

Water Activity (aw)

Not Tested

PASS/FAIL

PASS

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

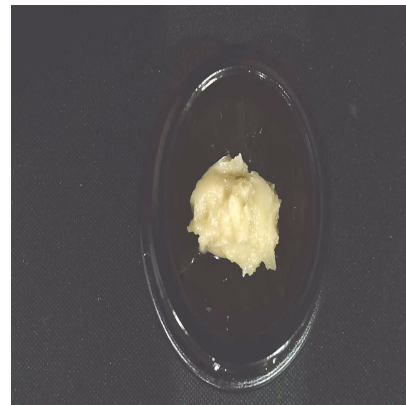
Moisture content levels less than 15% are recommended but the sample does not fail. Water activity levels must be less than 0.65aw.

Cannabinoids (Top 3)	(%)	mg/g
THCA	81.2	812
CBGA	2.53	25.3
Δ9-THC	0.841	8.41
TOTAL CBD	-	-
TOTAL THC	72.1	721
TOTAL CANNABINOIDS	85.6	856

Terpenes (Top 5)	(%)	µg/g
d-Limonene	1.36	13600
β-Caryophyllene	1.22	12200
β-Myrcene	0.625	6250
α-Humulene	0.468	4680
Linalool	0.367	3670
TOTAL TERPENES	5.08	50800

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

Sample Picture Upon Receipt



Scan the QR code to verify results.

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

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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36126)

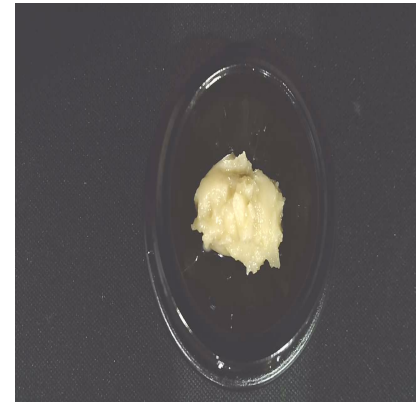
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Cultivar (Strain) or Sample Description: Unkle Sink Hybrid Solventless Live Hash Rosin **Date Completed:** 10/04/2023

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

Analysis Date/Time: 10/03/2023 1546 **Method:** HPLC/DAD **Moisture Content (%):** -
Analyst: PW **Instrument:** Agilent 1100 **Water Activity (aw):** -

<u>Cannabinoid</u>	<u>Result (%)</u>	<u>Result (mg/g)</u>	<u>LOD (mg/g)</u>	<u>LOQ (mg/g)</u>	<u>Result (mg/mL)</u>	<u>Per Serving (mg)</u>	<u>Per Unit (mg)</u>
CBC	ND	ND	0.107	0.250	-	-	-
CBCA	0.440	4.40	0.333	0.778	-	4.40	4.40
CBD	ND	ND	0.757	1.77	-	-	-
CBDV	ND	ND	0.279	0.651	-	-	-
CBDVA	ND	ND	0.121	0.283	-	-	-
CBG	ND	ND	0.324	0.756	-	-	-
CBGA	0.133	1.33	0.491	1.15	-	1.33	1.33
CBL	2.53	25.3	0.697	0.822	-	25.3	25.3
CBN	ND	ND	0.568	1.33	-	-	-
CBNA	ND	ND	0.261	0.609	-	-	-
Δ9-THC	ND	ND	0.282	0.656	-	-	-
Δ8-THC	0.841	8.41	0.313	0.729	-	8.41	8.41
THCA	ND	ND	0.488	1.14	-	-	-
THCV	81.2	812	0.169	0.396	-	812	812
THCVA	ND	ND	0.407	0.948	-	-	-
THCVA	0.405	4.05	0.130	0.302	-	4.05	4.05
TOTAL	85.6	856				856	856
TOTAL CBC	0.386	3.86				3.86	3.86
TOTAL CBD	-	-				-	-
TOTAL CBDV	-	-				-	-
TOTAL CBG	2.35	23.5				23.5	23.5
TOTAL CBN	-	-				-	-
TOTAL THC	72.1	721				721	721
TOTAL THCV	0.351	3.51				3.51	3.51



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 = (CBDV 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) + Δ9-THC
Total THCV = (THCVA x 0.867) + THCVA

Percentage results are reported by mass.

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

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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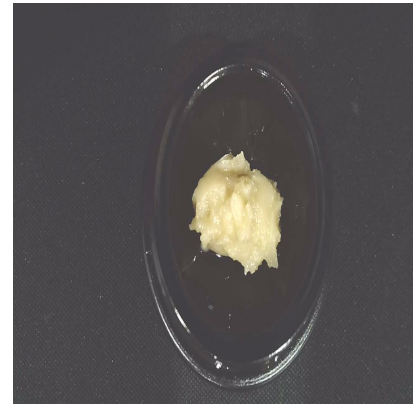
TERPENOID PROFILE

Analysis Date/Time: 10/04/2023 0225
Analyst: KF

Method: GC/MS
Instrument: Agilent 7890/5975

Deviations from SOP:
None

<u>Terpene</u>	<u>Result (µg/g)</u>	<u>Result (%)</u>	
α-Bisabolol	659	0.0659	
Camphene	502	0.0502	
δ-3-Carene	ND	-	
β-Caryophyllene	12200	1.22	█
Caryophyllene oxide	ND	-	
p-Cymene	126	0.0126	
Eucalyptol	ND	-	
Geraniol	ND	-	
Guaiol	690	0.0690	
α-Humulene	4680	0.468	█
Isopulegol	ND	-	
d-Limonene	13600	1.36	█
Linalool	3670	0.367	█
β-Myrcene	6250	0.625	█
cis-Nerolidol	534	0.0534	
trans-Nerolidol	1190	0.119	
α-Ocimene	ND	-	
β-Ocimene	753	0.0753	
α-Pinene	2450	0.245	█
β-Pinene	3200	0.320	█
α-Terpinene	<LOQ	-	
γ-Terpinene	ND	-	
Terpinolene	282	0.0282	
TOTAL	50800	5.08	



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): 78.5

"-" Not detected above LOD.

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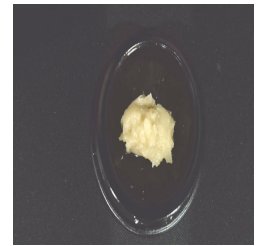
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Cultivar (Strain) or Sample Description: Unkle Sink Hybrid Solventless Live Hash Rosin **Date Completed:** 10/04/2023

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

Analysis Date/Time: 10/03/2023 2144 **Method:** HS/GC/MS **Deviations from SOP:**
Analyst: KF **Instrument:** Agilent 7890/5975 **None**

Solvent	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)	Solvent	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
Acetone (67-64-1)	-	150	300	5000	n-Heptane (142-82-5)	-	150	300	5000
Acetonitrile (75-5-8)	-	150	300	410	n-Hexane (110-54-3)	-	52.5	105	290
Benzene (71-43-2)	-	1.50	3.00	2	Isobutane (75-28-5)	-	150	300	5000
n-Butane (106-97-2)	-	150	300	5000	Isopropanol (67-63-0)	-	150	300	5000
1-Butanol (71-36-3)	-	150	300	5000	Isopropyl acetate (108-21-4)	-	150	300	5000
2-Butanol (78-92-2)	-	150	300	5000	Isopropyl benzene (98-82-8)	-	15.0	30.0	70
2-Butanone (78-93-3)	-	150	300	5000	Methanol (67-56-1)	-	150	300	3000
Cyclohexane (110-82-7)	-	150	300	3880	2-Methylbutane (78-78-4)	-	150	300	5000
1,2-Dimethoxyethane (110-71-4)	-	15.0	30.0	100	Methylene chloride (75-9-2)	-	150	300	600
N,N-Dimethylacetamide (127-19-5)	-	150	300	1090	2-Methylpentane (107-83-5)	-	52.5	105	290
2,2-Dimethylbutane (75-83-2)	-	52.5	105	290	3-Methylpentane (96-10-0)	-	52.5	105	290
2,3-Dimethylbutane (79-29-8)	-	52.5	105	290	n-Pentane (109-66-0)	-	150	300	5000
N,N-Dimethylformamide (68-12-2)	-	150	300	880	1-Pentanol (71-41-0)	-	150	300	5000
Dimethylsulfoxide (67-68-5)	-	150	300	5000	n-Propane (74-98-6)	-	150	300	5000
1,4-Dioxane (123-91-1)	-	150	300	380	1-Propanol (71-23-8)	-	150	300	5000
Ethanol (64-17-5)	-	150	300	5000	Pyridine (110-86-1)	-	52.5	105	200
2-Ethoxyethanol (110-80-5)	-	52.5	105	160	Tetrahydrofuran (109-99-9)	-	150	300	720
Ethyl ether (60-29-7)	-	150	300	5000	Tetramethylene sulfone (126-33-0)	-	52.5	105	160
Ethyl acetate (141-78-6)	-	150	300	5000	Toluene (108-88-3)	-	150	300	890
Ethyl benzene (100-41-4)	-	150	300	2170	o-Xylene (95-47-6)	-	150	300	2170
Ethylene glycol (107-21-1)	-	150	300	620	m,p-Xylene (108-38-3 or 106-42-3)	-	150	300	2170
Ethylene oxide (75-21-8)	-	15.0	30.0	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170



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.

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
Dimethylsulfoxide	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

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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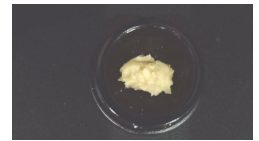
PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 10/03/2023 1626
Analyst: KF

Method: LC/MS/MS
Instrument: Shimadzu LC-8050

Deviations from SOP:
None

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



Color Key

RESULT < AL
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.
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Pesticide	Synonym(s)	Pesticide	Synonym(s)	Pesticide	Synonym(s)
Cyfluthrin	Baythroid	Myclobutanil	Systhane	Propiconazole	Tilt
DDVP	Dichlorvos	Naled	Dibrom	Propoxur	Baygon
Ethoprophos	Prophos	Phosmet	Imidan		

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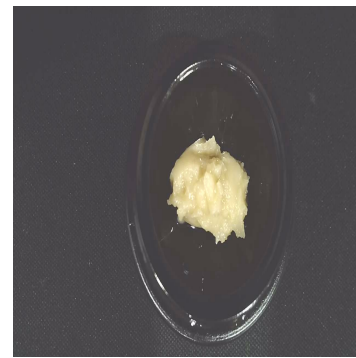
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HEAVY METAL PROFILE (SOP: SOP-ICP-200.6)

Analysis Date/Time: 10/03/2023 1847 (ICP/OES) **Method:** ICP/MS **Deviations from SOP:**
Analysis Date/Time: - (DMA) **Instrument:** Agilent 7500ce 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)	-	58.7	93.0	200
Cadmium (Cd)	-	58.7	93.0	200
Lead (Pb)	-	58.7	93.0	500
Mercury (Hg)	-	58.7	93.0	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



"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.

REPORT OF LABORATORY ANALYSIS

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 Felling Analytical Services and Technology (F.A.S.T.), LLC

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



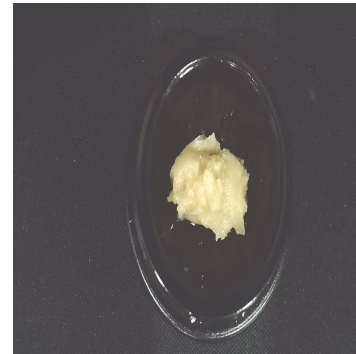
CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36126)

Testing Location: Arkansas 232 S. Broadview St. Greenbrier, AR 72058 License: ADA 05_H273	Customer ID: 2168 River Valley Relief MIPS 5601 Old Greenwood Rd Suite C Fort Smith, AR 72903 License: 00065C	Sample ID: SA36126 Lot Number: M00065C13247672701 Production Run: E2023USKLHR01	Sample Type: Primary Matrix: Concentrate Mass: 4g Date Collected: 10/02/2023 Date Received: 10/03/2023
Cultivar (Strain) or Sample Description: Unkle Sink Hybrid Solventless Live Hash Rosin			Date Completed: 10/04/2023

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 10/04/2023 0838 **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 (Green background)

RESULT > AL (Red background)

Reporting Limit (CFU/g)
1

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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