F.A.S.T. LABORATORIES/RESEARCH

ACCREDITED CERTIFICATE #4473.01

ASA PFC Laboratory

Analyzed on behalf of:

RVR RIVER VALLEY RELIEF

SUMMARY OF ANALYSIS (SAMPLE ID: SA35790)

Testing Location:	**Customer ID:** 2168	**Order ID:** OR10562	**Sample Type:** Primary
Arkansas	River Valley Relief MIPS	**Lot Number:**	**Matrix:** Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13351500364	**Mass:** 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	**Production Run:**	**Date Collected:** 08/07/2023
License: ADH 113	License: 00065C	E20230803AKR16	**Date Received:** 08/08/2023

Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil — **Date Completed:** 08/10/2023

*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

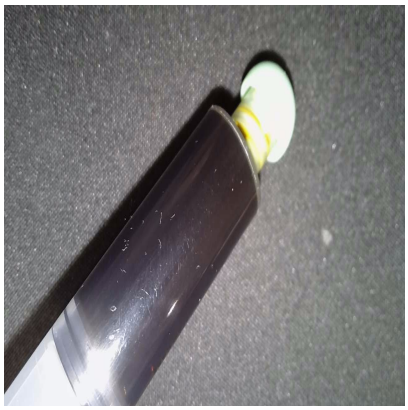
Moisture content/water activity action levels are referenced from the State of Arkansas MMJ testing guidelines.
Moisture content levels less than 15% are recommended but the sample does not fail. Water activity levels must be less than 0.65aw.

Cannabinoids (Top 3)	(%)	mg/g
Δ9-THC	78.8	788
CBG	2.94	29.4
CBC	0.870	8.70
TOTAL CBD	0.251	2.51
TOTAL THC	78.8	788
TOTAL CANNABINOIDS	83.8	838

Terpenes (Top 5)	(%)	µg/g
β-Caryophyllene	1.54	15400
α-Humulene	0.560	5600
d-Limonene	0.312	3120
Linalool	0.122	1220
trans-Nerolidol	0.119	1190
TOTAL TERPENES	3.14	31400

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, Ph.D.
Laboratory Director

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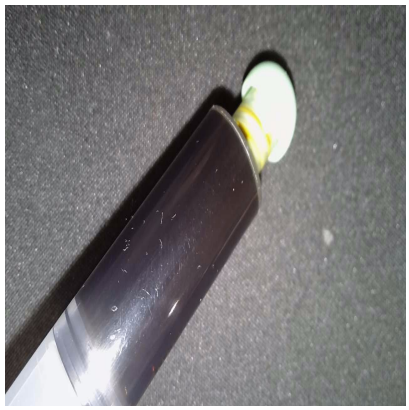
Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil — **Date Completed:** 08/10/2023

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

Analysis Date/Time: 08/08/2023 1416
Analyst: PW
Method: HPLC/DAD
Instrument: Agilent 1100
Moisture Content (%): -
Water Activity (aw): -

Cannabinoid	Result (%)	Result (mg/g)	LOD (mg/g)	LOQ (mg/g)	Result (mg/mL)	Per Serving (mg)	Per Unit (mg)
CBC	0.870	8.70	0.0995	0.232	-	8.70	8.70
CBCA	ND	ND	0.309	0.721	-	-	-
CBD	0.251	2.51	0.702	1.64	-	2.51	2.51
CBDA	ND	ND	0.258	0.603	-	-	-
CBDV	ND	ND	0.113	0.263	-	-	-
CBDVA	ND	ND	0.300	0.701	-	-	-
CBG	2.94	29.4	0.455	1.06	-	29.4	29.4
CBGA	0.116	1.16	0.646	0.762	-	1.16	1.16
CBL	ND	ND	0.526	1.23	-	-	-
CBN	0.410	4.10	0.242	0.565	-	4.10	4.10
CBNA	ND	ND	0.261	0.608	-	-	-
Δ9-THC	78.8	788	0.290	0.675	-	788	788
Δ8-THC	ND	ND	0.452	1.05	-	-	-
THCA	ND	ND	0.157	0.367	-	-	-
THCV	0.390	3.90	0.377	0.879	-	3.90	3.90
THCVA	ND	ND	0.120	0.280	-	-	-
TOTAL	83.8	838			-	838	838
TOTAL CBC	0.870	8.70			-	8.70	8.70
TOTAL CBD	0.251	2.51			-	2.51	2.51
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	3.05	30.5			-	30.5	30.5
TOTAL CBN	0.410	4.10			-	4.10	4.10
TOTAL THC	78.8	788			-	788	788
TOTAL THCV	0.390	3.90			-	3.90	3.90



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) + Δ9-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.

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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Laboratory Director

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License: ADH 113	License: 00065C	E20230803AKR16	**Date Received:** 08/08/2023

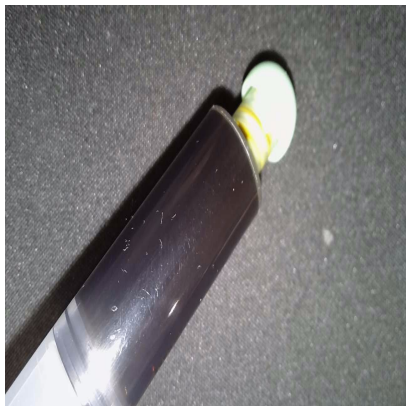
Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil — **Date Completed:** 08/10/2023

TERPENOID PROFILE

Analysis Date/Time:08/08/2023 1455 — **Method:** GC/MS — **Deviations from SOP:**

Analyst: KF — **Instrument:** Agilent 7890/5975 — None

Terpene	Result (µg/g)	Result (%)
α-Bisabolol	175	0.0175
Camphene	122	0.0122
δ-3-Carene	<LOQ	-
β-Caryophyllene	15400	1.54
Caryophyllene oxide	332	0.0332
p-Cymene	87.3	0.00873
Eucalyptol	52.4	0.00524
Geraniol	ND	-
Guaiol	314	0.0314
α-Humulene	5600	0.560
Isopulegol	52.4	0.00524
d-Limonene	3120	0.312
Linalool	1220	0.122
β-Myrcene	925	0.0925
cis-Nerolidol	69.8	0.00698
trans-Nerolidol	1190	0.119
α-Ocimene	87.3	0.00873
β-Ocimene	384	0.0384
α-Pinene	524	0.0524
β-Pinene	419	0.0419
α-Terpinene	105	0.0105
γ-Terpinene	157	0.0157
Terpinolene	1060	0.106
TOTAL	31400	3.14



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): 43.6

"-" Not detected above LOD.

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License: ADH 113	License: 00065C	E20230803AKR16	**Date Received:** 08/08/2023

Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil — **Date Completed:** 08/10/2023

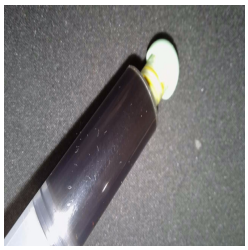
RESIDUAL SOLVENT PROFILE (SOP: SOP-RS-001)

Analysis Date/Time: 08/08/2023 2123 — **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)
Acetone (67-64-1)	-	143	286	5000
Acetonitrile (75-5-8)	-	143	286	410
Benzene (71-43-2)	-	1.43	2.86	2
n-Butane (106-97-2)	-	143	286	5000
1-Butanol (71-36-3)	-	143	286	5000
2-Butanol (78-92-2)	-	143	286	5000
2-Butanone (78-93-3)	-	143	286	5000
Cyclohexane (110-82-7)	-	143	286	3880
1,2-Dimethoxyethane (110-71-4)	-	14.3	28.6	100
N,N-Dimethylacetamide (127-19-5)	-	143	286	1090
2,2-Dimethylbutane (75-83-2)	-	50.0	100	290
2,3-Dimethylbutane (79-29-8)	-	50.0	100	290
N,N-Dimethylformamide (68-12-2)	-	143	286	880
Dimethylsulfoxide (67-68-5)	-	143	286	5000
1,4-Dioxane (123-91-1)	-	143	286	380
Ethanol (64-17-5)	-	143	286	5000
2-Ethoxyethanol (110-80-5)	-	50.0	100	160
Ethyl ether (60-29-7)	-	143	286	5000
Ethyl acetate (141-78-6)	-	143	286	5000
Ethyl benzene (100-41-4)	-	143	286	2170
Ethylene glycol (107-21-1)	-	143	286	620
Ethylene oxide (75-21-8)	-	14.3	28.6	50

Solvent	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
n-Heptane (142-82-5)	-	143	286	5000
n-Hexane (110-54-3)	-	50.0	100	290
Isobutane (75-28-5)	-	143	286	5000
Isopropanol (67-63-0)	-	143	286	5000
Isopropyl acetate (108-21-4)	-	143	286	5000
Isopropyl benzene (98-82-8)	-	14.3	28.6	70
Methanol (67-56-1)	-	143	286	3000
2-Methylbutane (78-78-4)	-	143	286	5000
Methylene chloride (75-9-2)	-	143	286	600
2-Methylpentane (107-83-5)	-	50.0	100	290
3-Methylpentane (96-10-0)	-	50.0	100	290
n-Pentane (109-66-0)	-	143	286	5000
1-Pentanol (71-41-0)	-	143	286	5000
n-Propane (74-98-6)	-	143	286	5000
1-Propanol (71-23-8)	-	143	286	5000
Pyridine (110-86-1)	-	50.0	100	200
Tetrahydrofuran (109-99-9)	-	143	286	720
Tetramethylene sulfone (126-33-0)	-	50.0	100	160
Toluene (108-88-3)	-	143	286	890
o-Xylene (95-47-6)	-	143	286	2170
m,p-Xylene (108-38-3 or 106-42-3)	-	143	286	2170
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.

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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

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: 08/09/2023 0550 — **Method:** LC/MS/MS — **Deviations from SOP:**

Analyst: KF — **Instrument:** Shimadzu LC-8050 — None

Pesticide	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
Abamectin (71751-41-2)	-	0.0436	0.348	0.5
Acephate (30560-19-1)	-	0.0436	0.348	0.4
Acequinocyl (57960-19-7)	-	0.0436	0.348	2
Acetamiprid (135410-20-7)	-	0.0436	0.348	0.2
Aldicarb (116-06-3)	-	0.0436	0.348	0.4
Azoxystrobin (131860-33-8)	-	0.0436	0.348	0.2
Bifenazate (149877-41-8)	-	0.0436	0.348	0.2
Bifenthrin (82657-04-3)	-	0.0436	0.348	0.2
Boscalid (188425-85-6)	-	0.0436	0.348	0.4
Carbaryl (63-25-2)	-	0.0436	0.348	0.2
Carbofuran (1563-66-2)	-	0.0436	0.348	0.2
Chlorantraniliprole (800008-45-7)	-	0.0436	0.348	0.2
Chlorfenapyr (122453-73-0)	-	0.0436	0.348	1
Chlorpyrifos (2921-88-2)	-	0.0436	0.348	0.2
Clofentezine (74115-24-5)	-	0.0436	0.348	0.2
Cyfluthrin (68359-37-5)	-	0.0436	0.348	1
Cypermethrin (52315-07-8)	-	0.0436	0.348	1
Daminozide (1596-84-5)	-	0.0436	0.348	1
DDVP (62-73-7)	-	0.0436	0.348	0.1
Diazinon (333-41-5)	-	0.0436	0.348	0.2
Dimethoate (60-51-5)	-	0.0436	0.348	0.2
Ethoprophos (13194-48-4)	-	0.0436	0.348	0.2
Etofenprox (80844-07-1)	-	0.0436	0.348	0.4
Etoxazole (153233-91-1)	-	0.0436	0.348	0.2
Fenoxycarb (72490-01-8)	-	0.0436	0.348	0.2
(E)-Fenpyroximate (134098-61-6)	-	0.0436	0.348	0.4
Fipronil (120068-37-3)	-	0.0436	0.348	0.4
Flonicamid (158062-67-0)	-	0.0436	0.348	1
Fludioxinil (131341-86-1)	-	0.0436	0.348	0.4
Hexythiazox (78587-05-0)	-	0.0436	0.348	1
Imazalil (35554-44-0)	-	0.0436	0.348	0.2
Imidacloprid (138261-41-3)	-	0.0436	0.348	0.4
Kresoxim-methyl (143390-89-0)	-	0.0436	0.348	0.4
Malathion (121-75-5)	-	0.0436	0.348	0.2
Metalaxyl (57837-19-1)	-	0.0436	0.348	0.2
Methiocarb (2032-65-7)	-	0.0436	0.348	0.2
Methomyl (16752-77-5)	-	0.0436	0.348	0.4
Methyl parathion (298-0-0)	-	0.0436	0.348	0.2
MGK 264 (113-48-4)	-	0.0436	0.348	0.2
Myclobutanil (88671-89-0)	-	0.0436	0.348	0.2
Naled (300-76-5)	-	0.0436	0.348	0.5
Oxamyl (23135-22-0)	-	0.0436	0.348	1
Paclobutrazol (76738-62-0)	-	0.0436	0.348	0.4
Permethrins (52645-53-1)	-	0.0436	0.348	0.2
Phosmet (732-11-6)	-	0.0436	0.348	0.2
Piperonyl butoxide (51-03-6)	-	0.0436	0.348	2
Prallethrins (2331-36-9)	-	0.0436	0.348	0.2
Propiconazole (60207-90-1))	-	0.0436	0.348	0.4
Propoxur (114-26-1)	-	0.0436	0.348	0.2
Pyrethrins (8003-34-7)	-	0.0436	0.348	1
Pyridaben (96489-71-3)	-	0.0436	0.348	0.2
Spinosad (168316-95-8)	-	0.0436	0.348	0.2
Spiromesifen (283594-90-1)	-	0.0436	0.348	0.2
Spirotetramat (203313-25-1)	-	0.0436	0.348	0.2
Spiroxamine (118134-30-8)	-	0.0436	0.348	0.4
Tebuconazole (80443-41-0)	-	0.0436	0.348	0.4
Thiacloprid (111988-49-9)	-	0.0436	0.348	0.2
Thiamethoxam (153719-23-4)	-	0.0436	0.348	0.2
Trifloxystrobin (141517-21-7)	-	0.0436	0.348	0.2

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.

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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

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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Testing Location:	**Customer ID:** 2168	**Order ID:** OR10562	**Sample Type:** Primary
Arkansas	River Valley Relief MIPS	**Lot Number:**	**Matrix:** Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13351500364	**Mass:** 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	**Production Run:**	**Date Collected:** 08/07/2023
License: ADH 113	License: 00065C	E20230803AKR16	**Date Received:** 08/08/2023

Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil — **Date Completed:** 08/10/2023

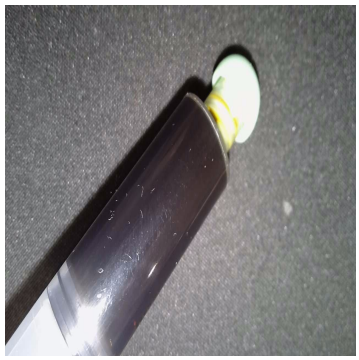
HEAVY METAL PROFILE (SOP: SOP-ICP-200.6)

Analysis Date/Time: 08/10/2023 0738 (ICP/OES)
Analysis Date/Time: - (DMA)
Analyst: KF

Method: ICP/MS
Instrument: Agilent 7500ce

Deviations from SOP:
None

Heavy Metal	Result (µg/kg)	LOD (µg/kg)	LOQ (µg/kg)	Action Level (µg/kg)
Arsenic (As)	-	58.2	92.2	200
Cadmium (Cd)	-	58.2	92.2	200
Lead (Pb)	-	58.2	92.2	500
Mercury (Hg)	-	58.2	92.2	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

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.

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.

Kyle W. Felling, Ph.D.
Laboratory Director

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F.A.S.T. LABORATORIES/RESEARCH

ACCREDITED CERTIFICATE #4473.01

ASA PFC Laboratory

Analyzed on behalf of:

RVR RIVER VALLEY RELIEF

CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35790)

Testing Location:	**Customer ID:** 2168	**Sample ID:** SA35790	**Sample Type:** Primary
Arkansas	River Valley Relief MIPS	**Lot Number:**	**Matrix:** Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13351500364	**Mass:** 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	**Production Run:**	**Date Collected:** 08/07/2023
License: ADA 05_H273	License: 00065C	E20230803AKR16	**Date Received:** 08/08/2023
Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil			**Date Completed:** 08/10/2023

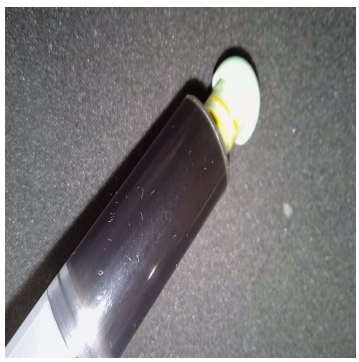
MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 08/10/2023 0832
Analyst: PW

Method: Hardy Diagnostics CompactDry
Instrument: Thermo Incubator

Deviations from SOP:
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)
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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