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: SA40274)

Testing Location:	**Customer ID:** 2168	**Order ID:** OR11449	**Sample Type:** Primary
Arkansas	River Valley Relief MIPS	**Lot Number:**	**Matrix:** Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13357938987	**Mass:** 4ea
Greenbrier, AR 72058	Fort Smith, AR 72903	**Production Run:**	**Date Collected:** 02/17/2025
License: ADH 113	License: 00065C	E20250213RAW26	**Date Received:** 02/17/2025

Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil 1g Applicator **Date Completed:**02/20/2025

*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

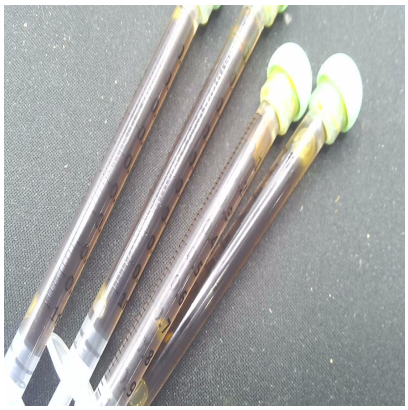
Moisture content/water activity action levels are referenced from the State of Arkansas MMJ testing guidelines.
Moisture content levels less than 15% are recommended but the sample does not fail. Water activity levels must be less than 0.65aw.

Cannabinoids (Top 3)	(%)	mg/g
Δ9-THC	74.1	741
CBG	2.51	25.1
CBC	0.859	8.59
TOTAL CBD	0.000	0.000
TOTAL THC	74.1	741
TOTAL CANNABINOIDS	78.7	787

Terpenes (Top 5)	(%)	µg/g
β-Caryophyllene	2.09	20900
β-Myrcene	1.71	17100
d-Limonene	1.45	14500
α-Humulene	0.702	7020
Linalool	0.509	5090
TOTAL TERPENES	7.26	72600

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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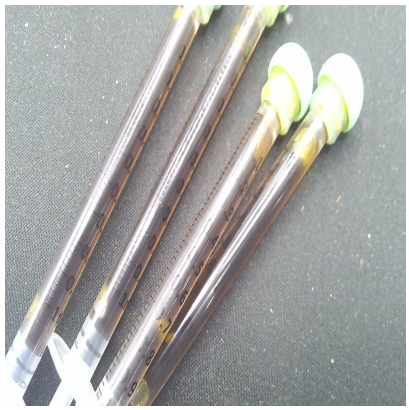
License: ADH 113	License: 00065C	E20250213RAW26	**Date Received:** 02/17/2025

Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil 1g Applicator **Date Completed:**02/20/2025

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

Analysis Date/Time: 02/18/2025 1147
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	0.859	8.59	0.103	0.241	-	8.59	8.59
CBCA	ND	ND	0.320	0.748	-	0.000	0.000
CBD	DET	DET	0.728	1.70	-	0.000	0.000
CBDA	ND	ND	0.268	0.626	-	0.000	0.000
CBDV	ND	ND	0.117	0.272	-	0.000	0.000
CBDVA	ND	ND	0.311	0.727	-	0.000	0.000
CBG	2.51	25.1	0.473	1.10	-	25.1	25.1
CBGA	ND	ND	0.670	0.790	-	0.000	0.000
CBL	ND	ND	0.546	1.27	-	0.000	0.000
CBN	0.796	7.96	0.251	0.586	-	7.96	7.96
CBNA	ND	ND	0.271	0.631	-	0.000	0.000
CBT	ND	ND	0.393	0.916	-	0.000	0.000
Δ9-THC	74.1	741	0.301	0.701	-	741	741
Δ8-THC	ND	ND	0.469	1.09	-	0.000	0.000
THCA	ND	ND	0.163	0.381	-	0.000	0.000
THCV	0.421	4.21	0.391	0.912	-	4.21	4.21
THCVA	ND	ND	0.125	0.291	-	0.000	0.000
TOTAL	78.7	787			-	787	787
TOTAL CBC	0.859	8.59			-	8.59	8.59
TOTAL CBD	0.000	0.000			-	0.000	0.000
TOTAL CBDV	0.000	0.000			-	0.000	0.000
TOTAL CBG	2.51	25.1			-	25.1	25.1
TOTAL CBN	0.796	7.96			-	7.96	7.96
TOTAL THC	74.1	741			-	741	741
TOTAL THCV	0.421	4.21			-	4.21	4.21



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

"-" Not reported for this sample.

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, ND - Not Detected (less than LOD)

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Laboratory Director

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RVR RIVER VALLEY RELIEF

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License: ADH 113	License: 00065C	E20250213RAW26	**Date Received:** 02/17/2025

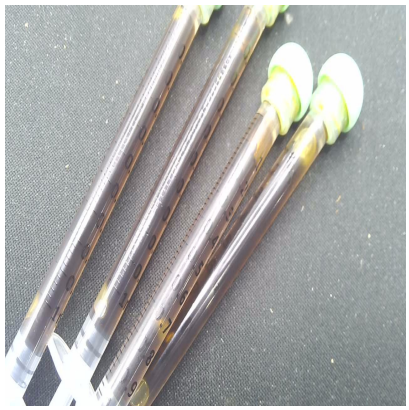
Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil 1g Applicator **Date Completed:**02/20/2025

TERPENOID PROFILE

Analysis Date/Time:02/19/2025 2032 **Method:** GC/MS **Deviations from SOP:**

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

Terpene	Result (µg/g)	Result (%)
α-Bisabolol	3040	0.304
Camphene	ND	-
δ-3-Carene	ND	-
β-Caryophyllene	20900	2.09
Caryophyllene oxide	338	0.0338
p-Cymene	ND	-
Eucalyptol	ND	-
Geraniol	ND	-
Guaiol	<LOQ	-
α-Humulene	7020	0.702
Isopulegol	ND	-
d-Limonene	14500	1.45
Linalool	5090	0.509
β-Myrcene	17100	1.71
cis-Nerolidol	ND	-
trans-Nerolidol	1790	0.179
α-Ocimene	ND	-
β-Ocimene	191	0.0191
α-Pinene	1220	0.122
β-Pinene	1320	0.132
α-Terpinene	ND	-
γ-Terpinene	ND	-
Terpinolene	145	0.0145
TOTAL	72600	7.26



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): 45.3

"-" Not detected above LOD.

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License: ADH 113	License: 00065C	E20250213RAW26	**Date Received:** 02/17/2025

Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil 1g Applicator **Date Completed:**02/20/2025

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

Analysis Date/Time: 02/17/2025 1628 **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)	-	40.0	80.0	5000
Acetonitrile (75-5-8)	-	40.0	80.0	410
Benzene (71-43-2)	-	0.400	0.800	2
n-Butane (106-97-2)	-	40.0	80.0	5000
1-Butanol (71-36-3)	-	40.0	80.0	5000
2-Butanol (78-92-2)	-	40.0	80.0	5000
2-Butanone (78-93-3)	-	40.0	80.0	5000
Cyclohexane (110-82-7)	-	40.0	80.0	3880
1,2-Dimethoxyethane (110-71-4)	-	4.00	8.00	100
N,N-Dimethylacetamide (127-19-5)	-	40.0	80.0	1090
2,2-Dimethylbutane (75-83-2)	-	14.0	28.0	290
2,3-Dimethylbutane (79-29-8)	-	14.0	28.0	290
N,N-Dimethylformamide (68-12-2)	-	40.0	80.0	880
Dimethylsulfoxide (67-68-5)	-	40.0	80.0	5000
1,4-Dioxane (123-91-1)	-	40.0	80.0	380
Ethanol (64-17-5)	-	40.0	80.0	5000
2-Ethoxyethanol (110-80-5)	-	14.0	28.0	160
Ethyl ether (60-29-7)	-	40.0	80.0	5000
Ethyl acetate (141-78-6)	-	40.0	80.0	5000
Ethyl benzene (100-41-4)	-	40.0	80.0	2170
Ethylene glycol (107-21-1)	-	40.0	80.0	620
Ethylene oxide (75-21-8)	-	4.00	8.00	50

Solvent	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
n-Heptane (142-82-5)	-	40.0	80.0	5000
n-Hexane (110-54-3)	-	14.0	28.0	290
Isobutane (75-28-5)	-	40.0	80.0	5000
Isopropanol (67-63-0)	-	40.0	80.0	5000
Isopropyl acetate (108-21-4)	-	40.0	80.0	5000
Isopropyl benzene (98-82-8)	-	4.00	8.00	70
Methanol (67-56-1)	-	40.0	80.0	3000
2-Methylbutane (78-78-4)	-	40.0	80.0	5000
Methylene chloride (75-9-2)	-	40.0	80.0	600
2-Methylpentane (107-83-5)	-	14.0	28.0	290
3-Methylpentane (96-10-0)	-	14.0	28.0	290
n-Pentane (109-66-0)	-	40.0	80.0	5000
1-Pentanol (71-41-0)	-	40.0	80.0	5000
n-Propane (74-98-6)	-	40.0	80.0	5000
1-Propanol (71-23-8)	-	40.0	80.0	5000
Pyridine (110-86-1)	-	14.0	28.0	200
Tetrahydrofuran (109-99-9)	-	40.0	80.0	720
Tetramethylene sulfone (126-33-0)	-	14.0	28.0	160
Toluene (108-88-3)	-	40.0	80.0	890
o-Xylene (95-47-6)	-	40.0	80.0	2170
m,p-Xylene (108-38-3 or 106-42-3)	-	40.0	80.0	2170
Xylenes* (1330-20-7)	-	43.3	86.7	2170

Color Key

RESULT < AL

RESULT > AL

"-" not detected above 1/2 Action Level

"*" - 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
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: 02/18/2025 1211 **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)	-	0.0450	0.360	0.5
Acephate (30560-19-1)	-	0.0450	0.360	0.4
Acequinocyl (57960-19-7)	DET	0.0450	0.360	2
Acetamiprid (135410-20-7)	-	0.0450	0.360	0.2
Aldicarb (116-06-3)	-	0.0450	0.360	0.4
Azoxystrobin (131860-33-8)	-	0.0450	0.360	0.2
Bifenazate (149877-41-8)	-	0.0450	0.360	0.2
Bifenthrin (82657-04-3)	-	0.0450	0.360	0.2
Boscalid (188425-85-6)	-	0.0450	0.360	0.4
Carbaryl (63-25-2)	-	0.0450	0.360	0.2
Carbofuran (1563-66-2)	-	0.0450	0.360	0.2
Chlorantraniliprole (800008-45-7)	-	0.0450	0.360	0.2
Chlorfenapyr (122453-73-0)	-	0.0450	0.360	1
Chlorpyrifos (2921-88-2)	-	0.0450	0.360	0.2
Clofentezine (74115-24-5)	-	0.0450	0.360	0.2
Cyfluthrin (68359-37-5)	DET	0.0450	0.360	1
Cypermethrin (52315-07-8)	-	0.0450	0.360	1
Daminozide (1596-84-5)	-	0.0450	0.360	1
DDVP (62-73-7)	-	0.0450	0.360	0.1
Diazinon (333-41-5)	-	0.0450	0.360	0.2
Dimethoate (60-51-5)	-	0.0450	0.360	0.2
Ethoprophos (13194-48-4)	-	0.0450	0.360	0.2
Etofenprox (80844-07-1)	-	0.0450	0.360	0.4
Etoxazole (153233-91-1)	-	0.0450	0.360	0.2
Fenoxycarb (72490-01-8)	-	0.0450	0.360	0.2
(E)-Fenpyroximate (134098-61-6)	-	0.0450	0.360	0.4
Fipronil (120068-37-3)	-	0.0450	0.360	0.4
Flonicamid (158062-67-0)	-	0.0450	0.360	1
Fludioxinil (131341-86-1)	-	0.0450	0.360	0.4
Hexythiazox (78587-05-0)	-	0.0450	0.360	1
Imazalil (35554-44-0)	-	0.0450	0.360	0.2
Imidacloprid (138261-41-3)	-	0.0450	0.360	0.4

Pesticide	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
Kresoxim-methyl (143390-89-0)	-	0.0450	0.360	0.4
Malathion (121-75-5)	-	0.0450	0.360	0.2
Metalaxyl (57837-19-1)	-	0.0450	0.360	0.2
Methiocarb (2032-65-7)	-	0.0450	0.360	0.2
Methomyl (16752-77-5)	-	0.0450	0.360	0.4
Methyl parathion (298-0-0)	-	0.0450	0.360	0.2
MGK 264 (113-48-4)	-	0.0450	0.360	0.2
Myclobutanil (88671-89-0)	-	0.0450	0.360	0.2
Naled (300-76-5)	-	0.0450	0.360	0.5
Oxamyl (23135-22-0)	-	0.0450	0.360	1
Paclobutrazol (76738-62-0)	-	0.0450	0.360	0.4
Permethrins (52645-53-1)	-	0.0450	0.360	0.2
Phosmet (732-11-6)	-	0.0450	0.360	0.2
Piperonyl butoxide (51-03-6)	-	0.0450	0.360	2
Prallethrins (2331-36-9)	-	0.0450	0.360	0.2
Propiconazole (60207-90-1))	-	0.0450	0.360	0.4
Propoxur (114-26-1)	-	0.0450	0.360	0.2
Pyrethrins (8003-34-7)	-	0.0450	0.360	1
Pyridaben (96489-71-3)	-	0.0450	0.360	0.2
Spinosad (168316-95-8)	-	0.0450	0.360	0.2
Spiromesifen (283594-90-1)	-	0.0450	0.360	0.2
Spirotetramat (203313-25-1)	-	0.0450	0.360	0.2
Spiroxamine (118134-30-8)	-	0.0450	0.360	0.4
Tebuconazole (80443-41-0)	-	0.0450	0.360	0.4
Thiacloprid (111988-49-9)	-	0.0450	0.360	0.2
Thiamethoxam (153719-23-4)	-	0.0450	0.360	0.2
Trifloxystrobin (141517-21-7)	-	0.0450	0.360	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		

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: SA40274)

Testing Location:	**Customer ID:** 2168	**Order ID:** OR11449	**Sample Type:** Primary
Arkansas	River Valley Relief MIPS	**Lot Number:**	**Matrix:** Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13357938987	**Mass:** 4ea
Greenbrier, AR 72058	Fort Smith, AR 72903	**Production Run:**	**Date Collected:** 02/17/2025
License: ADH 113	License: 00065C	E20250213RAW26	**Date Received:** 02/17/2025

Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil 1g Applicator **Date Completed:**02/20/2025

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

Analysis Date/Time: 02/19/2025 1220

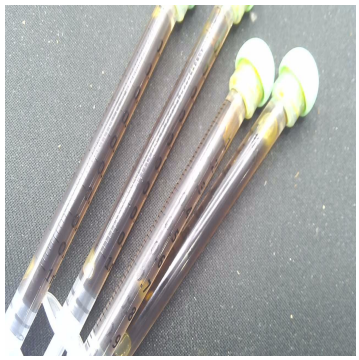
Method: ICP/MS

Instrument: Agilent 7700x

Deviations from SOP: None

Analyst: KF

Heavy Metal	Result (µg/kg)	LOD (µg/kg)	LOQ (µg/kg)	Action Level (µg/kg)
Arsenic (As)	-	59.6	94.4	200
Cadmium (Cd)	-	59.6	94.4	200
Lead (Pb)	-	59.6	94.4	500
Mercury (Hg)	-	59.6	94.4	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.

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ASA PFC Laboratory

Analyzed on behalf of:

RVR RIVER VALLEY RELIEF

CERTIFICATE OF ANALYSIS (SAMPLE ID: SA40274)

Testing Location:	**Customer ID:** 2168	**Sample ID:** SA40274	**Sample Type:** Primary
Arkansas	River Valley Relief MIPS	**Lot Number:**	**Matrix:** Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13357938987	**Mass:** 4ea
Greenbrier, AR 72058	Fort Smith, AR 72903	**Production Run:**	**Date Collected:** 02/17/2025
License: ADA 05_H273	License: 00065C	E20250213RAW26	**Date Received:** 02/17/2025

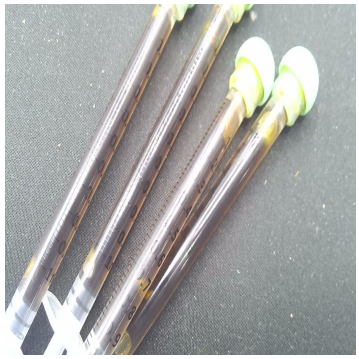
Cultivar (Strain) or Sample Description: ArkanRAW Full Spectrum Raw Cannabis Oil 1g Applicator **Date Completed:**02/20/2025

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 02/13/2025 1308 **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)
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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