

SUMMARY OF ANALYSIS (SAMPLE ID: SA41456)

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: OR11657 Lot Number: M00065C13242434714 Production Run: E20250619SPARKS01	Sample Type: Primary Matrix: Concentrate Mass: 4ea Date Collected: 06/19/2025 Date Received: 06/19/2025
Cultivar (Strain) or Sample Description: Sour Papaya Full Spectrum AR 1g Cart			Date Completed: 06/23/2025

*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.

*Where provided, statements of conformity (e.g. Pass/Fail) are made in accordance with ILAC G8, Binary Statement for Simple Acceptance Rule (w=0, AL=TL). PASS: when the result is within the acceptance interval. FAIL: when the result is outside the acceptance interval

Moisture Content (%)

Not Tested

Water Activity (aw)

Not Tested

PASS/FAIL

PASS

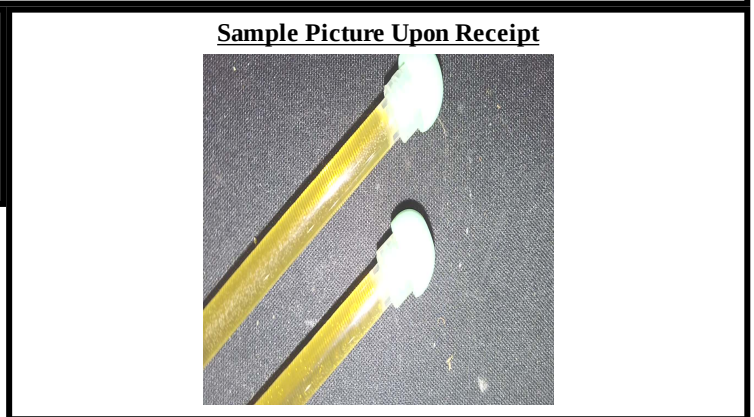
Moisture content/water activity action levels are referenced from the State of Arkansas MMJ testing guidelines.

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
Δ9-THC	68.8	688
THCA	15.2	152
CBGA	1.60	16.0
TOTAL CBD	0.249	2.49
TOTAL THC	82.1	821
TOTAL CANNABINOIDS	86.1	861

Terpenes (Top 5)	(%)	ppm
β-Caryophyllene	3.08	30800
β-Myrcene	2.08	20800
α-Humulene	1.92	19200
d-Limonene	0.637	6370
trans-Nerolidol	0.353	3530
TOTAL TERPENES	8.83	88300

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



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



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA41456)

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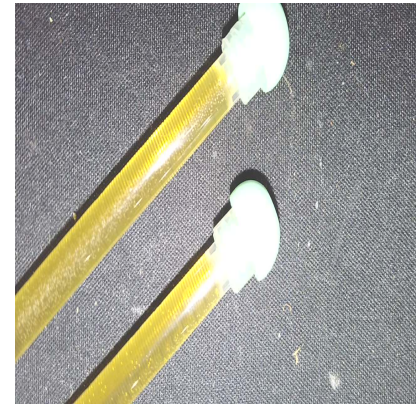
CANNABINOID (POTENCY) PROFILE (SOP: SOP-CANN-001)

Analysis Date/Time: 06/20/2025 1535
Analyst: PW

Method: HPLC/DAD
Instrument: Agilent 1100

Moisture Content (%): -
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	0.0970	0.226	-	0.000	0.000
CBCA	ND	ND	0.301	0.703	-	0.000	0.000
CBD	0.160	1.60	0.684	1.60	-	1.60	1.60
CBDA	0.102	1.02	0.252	0.588	-	1.02	1.02
CBDV	ND	ND	0.110	0.256	-	0.000	0.000
CBDVA	ND	ND	0.293	0.683	-	0.000	0.000
CBG	DET	DET	0.444	1.04	-	0.000	0.000
CBGA	1.60	16.0	0.629	0.743	-	16.0	16.0
CBL	ND	ND	0.513	1.20	-	0.000	0.000
CBN	0.126	1.26	0.236	0.550	-	1.26	1.26
CBNA	ND	ND	0.254	0.593	-	0.000	0.000
Δ9-THC	68.8	688	0.282	0.658	-	688	688
Δ8-THC	ND	ND	0.441	1.03	-	0.000	0.000
THCA	15.2	152	0.153	0.358	-	152	152
THCV	0.124	1.24	0.367	0.857	-	1.24	1.24
THCVA	ND	ND	0.117	0.273	-	0.000	0.000
TOTAL	86.1	861				861	861
TOTAL CBC	0.000	0.000				0.000	0.000
TOTAL CBD	0.249	2.49				2.49	2.49
TOTAL CBDV	0.000	0.000				0.000	0.000
TOTAL CBG	1.40	14.0				14.0	14.0
TOTAL CBN	0.126	1.26				1.26	1.26
TOTAL THC	82.1	821				821	821
TOTAL THCV	0.124	1.24				1.24	1.24



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.
Dry percent = Wet percent / (1-(Moisture Content/100))

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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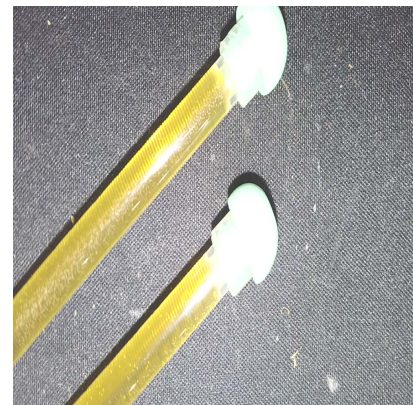
TERPENOID PROFILE

Analysis Date/Time: 06/20/2025 1629
Analyst: KF

Method: GC/MS
Instrument: Agilent 7890/5975

Deviations from SOP:
None

<u>Terpene</u>	<u>Result (ppm)</u>	<u>Result (%)</u>	
α-Bisabolol	506	0.0506	
Camphene	282	0.0282	
δ-3-Carene	<LOQ	-	
β-Caryophyllene	30800	3.08	█
Caryophyllene oxide	302	0.0302	
p-Cymene	ND	-	
Eucalyptol	345	0.0345	
Geraniol	ND	-	
Guaiol	<LOQ	-	
α-Humulene	19200	1.92	█
Isopulegol	ND	-	
d-Limonene	6370	0.637	█
Linalool	1050	0.105	
β-Myrcene	20800	2.08	█
cis-Nerolidol	ND	-	
trans-Nerolidol	3530	0.353	█
α-Ocimene	64.4	0.00644	
β-Ocimene	1660	0.166	
α-Pinene	1420	0.142	
β-Pinene	1730	0.173	
α-Terpinene	<LOQ	-	
γ-Terpinene	64.5	0.00645	
Terpinolene	190	0.0190	
TOTAL	88300	8.83	



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 (ppm): 42.5

"-" Not detected above LOD.

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

Analysis Date/Time: 06/19/2025 1744 **Method:** HS/GC/MS **Deviations from SOP:**
Analyst: KF **Instrument:** Agilent 7890/5975 **None**

Solvent	Result (ppm)	LOD (ppm)	LOQ (ppm)	Action Level (ppm)	Solvent	Result (ppm)	LOD (ppm)	LOQ (ppm)	Action Level (ppm)
Acetone (67-64-1)	-	32.2	64.3	5000	n-Heptane (142-82-5)	-	32.2	64.3	5000
Acetonitrile (75-5-8)	-	32.2	64.3	410	n-Hexane (110-54-3)	-	11.3	22.5	290
Benzene (71-43-2)	-	0.322	0.643	2	Isobutane (75-28-5)	-	32.2	64.3	5000
n-Butane (106-97-2)	-	32.2	64.3	5000	Isopropanol (67-63-0)	-	32.2	64.3	5000
1-Butanol (71-36-3)	-	32.2	64.3	5000	Isopropyl acetate (108-21-4)	-	32.2	64.3	5000
2-Butanol (78-92-2)	-	32.2	64.3	5000	Isopropyl benzene (98-82-8)	-	3.22	6.43	70
2-Butanone (78-93-3)	-	32.2	64.3	5000	Methanol (67-56-1)	-	32.2	64.3	3000
Cyclohexane (110-82-7)	-	32.2	64.3	3880	2-Methylbutane (78-78-4)	-	32.2	64.3	5000
1,2-Dimethoxyethane (110-71-4)	-	3.22	6.43	100	Methylene chloride (75-9-2)	-	32.2	64.3	600
N,N-Dimethylacetamide (127-19-5)	-	32.2	64.3	1090	2-Methylpentane (107-83-5)	-	11.3	22.5	290
2,2-Dimethylbutane (75-83-2)	-	11.3	22.5	290	3-Methylpentane (96-10-0)	-	11.3	22.5	290
2,3-Dimethylbutane (79-29-8)	-	11.3	22.5	290	n-Pentane (109-66-0)	-	32.2	64.3	5000
N,N-Dimethylformamide (68-12-2)	-	32.2	64.3	880	1-Pentanol (71-41-0)	-	32.2	64.3	5000
Dimethylsulfoxide (67-68-5)	-	32.2	64.3	5000	n-Propane (74-98-6)	-	32.2	64.3	5000
1,4-Dioxane (123-91-1)	-	32.2	64.3	380	1-Propanol (71-23-8)	-	32.2	64.3	5000
Ethanol (64-17-5)	-	32.2	64.3	5000	Pyridine (110-86-1)	-	11.3	22.5	200
2-Ethoxyethanol (110-80-5)	-	11.3	22.5	160	Tetrahydrofuran (109-99-9)	-	32.2	64.3	720
Ethyl ether (60-29-7)	-	32.2	64.3	5000	Tetramethylene sulfone (126-33-0)	-	11.3	22.5	160
Ethyl acetate (141-78-6)	-	32.2	64.3	5000	Toluene (108-88-3)	-	32.2	64.3	890
Ethyl benzene (100-41-4)	-	32.2	64.3	2170	o-Xylene (95-47-6)	-	32.2	64.3	2170
Ethylene glycol (107-21-1)	-	32.2	64.3	620	m,p-Xylene (108-38-3 or 106-42-3)	-	32.2	64.3	2170
Ethylene oxide (75-21-8)	-	3.22	6.43	50	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
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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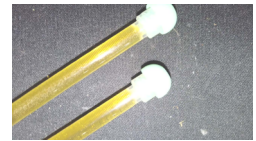
PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 06/20/2025 1603
Analyst: KF

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

Deviations from SOP:
None

Pesticide	Result (ppm)	LOD (ppm)	LOQ (ppm)	Action Level (ppm)	Pesticide	Result (ppm)	LOD (ppm)	LOQ (ppm)	Action Level (ppm)
Abamectin (71751-41-2)	-	0.0424	0.339	0.5	Kresoxim-methyl (143390-89-0)	-	0.0424	0.339	0.4
Acephate (30560-19-1)	-	0.0424	0.339	0.4	Malathion (121-75-5)	-	0.0424	0.339	0.2
Acequinocyl (57960-19-7)	-	0.0424	0.339	2	Metalaxyl (57837-19-1)	-	0.0424	0.339	0.2
Acetamiprid (135410-20-7)	-	0.0424	0.339	0.2	Methiocarb (2032-65-7)	-	0.0424	0.339	0.2
Aldicarb (116-06-3)	-	0.0424	0.339	0.4	Methomyl (16752-77-5)	-	0.0424	0.339	0.4
Azoxystrobin (131860-33-8)	-	0.0424	0.339	0.2	Methyl parathion (298-0-0)	-	0.0424	0.339	0.2
Bifenazate (149877-41-8)	-	0.0424	0.339	0.2	MGK 264 (113-48-4)	-	0.0424	0.339	0.2
Bifenthrin (82657-04-3)	-	0.0424	0.339	0.2	Myclobutanil (88671-89-0)	-	0.0424	0.339	0.2
Boscalid (188425-85-6)	-	0.0424	0.339	0.4	Naled (300-76-5)	-	0.0424	0.339	0.5
Carbaryl (63-25-2)	-	0.0424	0.339	0.2	Oxamyl (23135-22-0)	-	0.0424	0.339	1
Carbofuran (1563-66-2)	-	0.0424	0.339	0.2	Pacllobutrazol (76738-62-0)	-	0.0424	0.339	0.4
Chlorantraniliprole (800008-45-7)	-	0.0424	0.339	0.2	Permethrins (52645-53-1)	-	0.0424	0.339	0.2
Chlorfenapyr (122453-73-0)	-	0.0424	0.339	1	Phosmet (732-11-6)	-	0.0424	0.339	0.2
Chlorpyrifos (2921-88-2)	-	0.0424	0.339	0.2	Piperonyl butoxide (51-03-6)	-	0.0424	0.339	2
Clofentezine (74115-24-5)	-	0.0424	0.339	0.2	Prallethrins (2331-36-9)	-	0.0424	0.339	0.2
Cyfluthrin (68359-37-5)	-	0.0424	0.339	1	Propiconazole (60207-90-1)	-	0.0424	0.339	0.4
Cypermethrin (52315-07-8)	-	0.0424	0.339	1	Propoxur (114-26-1)	-	0.0424	0.339	0.2
Daminozide (1596-84-5)	-	0.0424	0.339	1	Pyrethrins (8003-34-7)	-	0.0424	0.339	1
DDVP (62-73-7)	-	0.0424	0.339	0.1	Pyridaben (96489-71-3)	-	0.0424	0.339	0.2
Diazinon (333-41-5)	-	0.0424	0.339	0.2	Spinosad (168316-95-8)	-	0.0424	0.339	0.2
Dimethoate (60-51-5)	-	0.0424	0.339	0.2	Spiromesifen (283594-90-1)	-	0.0424	0.339	0.2
Ethoprophos (13194-48-4)	-	0.0424	0.339	0.2	Spirotetramat (203313-25-1)	-	0.0424	0.339	0.2
Etofenprox (80844-07-1)	-	0.0424	0.339	0.4	Spiroxamine (118134-30-8)	-	0.0424	0.339	0.4
Etoxazole (153233-91-1)	-	0.0424	0.339	0.2	Tebuconazole (80443-41-0)	-	0.0424	0.339	0.4
Fenoxycarb (72490-01-8)	-	0.0424	0.339	0.2	Thiacloprid (111988-49-9)	-	0.0424	0.339	0.2
(E)-Fenpyroximate (134098-61-6)	-	0.0424	0.339	0.4	Thiamethoxam (153719-23-4)	-	0.0424	0.339	0.2
Fipronil (120068-37-3)	-	0.0424	0.339	0.4	Trifloxystrobin (141517-21-7)	-	0.0424	0.339	0.2
Fonicamid (158062-67-0)	-	0.0424	0.339	1					
Fludioxinil (131341-86-1)	-	0.0424	0.339	0.4					
Hexythiazox (78587-05-0)	-	0.0424	0.339	1					
Imazalil (35554-44-0)	-	0.0424	0.339	0.2					
Imidacloprid (138261-41-3)	-	0.0424	0.339	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.

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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	Sythane	Propiconazole	Tilt
DDVP	Dichlorvos	Naled	Dibrom	Propoxur	Baygon
Ethoprophos	Prophos	Phosmet	Imidan		

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Laboratory Director



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA41456)

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: OR11657 Lot Number: M00065C13242434714 Production Run: E20250619SPARKS01	Sample Type: Primary Matrix: Concentrate Mass: 4ea Date Collected: 06/19/2025 Date Received: 06/19/2025
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Cultivar (Strain) or Sample Description: Sour Papaya Full Spectrum AR 1g Cart **Date Completed:** 06/23/2025

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

Analysis Date/Time: 06/20/2025 1703

Method: ICP/MS

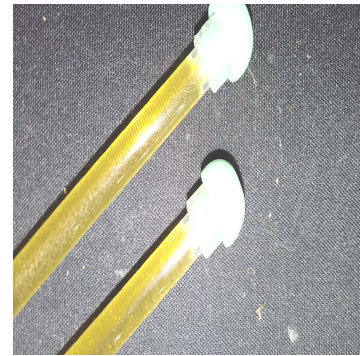
Deviations from SOP:

Instrument: Agilent 7700x

None

Analyst: KF

<u>Heavy Metal</u>	<u>Result (ppb)</u>	<u>LOD (ppb)</u>	<u>LOQ (ppb)</u>	<u>Action Level (ppb)</u>
Arsenic (As)	ND	58.9	93.3	200
Cadmium (Cd)	ND	58.9	93.3	200
Lead (Pb)	ND	58.9	93.3	500
Mercury (Hg)	ND	58.9	93.3	100



Abbreviations: ICP - Inductively-Coupled Plasma, MS - Mass Spectrometry, RL - Reporting Limit, AL - Action Level, LOD - Limit of Detection, LOQ - Limit of Quantitation

Color Key



"DET" detected less than LOQ

"ND" not detected above LOD

Reporting limit is 1/2 the action level of each analyte.

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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F.A.S.T. Personnel perform sampling following the Sampling SOP (SOP-02).

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



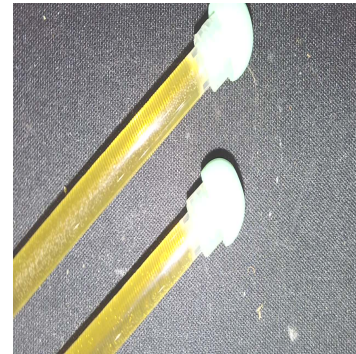
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Cultivar (Strain) or Sample Description: Sour Papaya Full Spectrum AR 1g Cart			Date Completed: 06/23/2025

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 06/20/2025 1210 **Method:** Hardy Diagnostics CompactDry **Deviations from SOP:**
Analyst: PW **Instrument:** Thermo Incubator None

<u>Bacteria/Microbe</u>	<u>Result (CFU/g)</u>	<u>Action Level (CFU/g)</u>
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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