







### SUMMARY OF ANALYSIS (SAMPLE ID: SA35050)

Testing Location:	Customer ID: 2168	Order ID: OR10416	Sample Type: Primary
Arkansas	<b>River Valley Relief MIPS</b>	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13245725149	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 04/24/2023
License: ADH 113	License: 00065C	E2023KMTHTR01	Date Received: 04/24/2023
Cultivar (Strain) or Sample D	Date Completed: 05/02/2023		

\*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)	<u>(%)</u>	<u>mg/g</u>
THCA	41.4	414
Δ9-THC	21.7	217
CBD	5.02	50.2
TOTAL CBD	5.35	53.5
TOTAL THC	58.0	580
TOTAL CANNABINOIDS	75.5	755
<u>Terpenes (Top 5)</u>	<u>(%)</u>	µg∕g
<u>Terpenes (Top 5)</u> d-Limonene	<u>(%)</u> 3.24	<b>µg/g</b> 32400
d-Limonene	3.24	32400
d-Limonene β-Caryophyllene	3.24 1.81	32400 18100
d-Limonene β-Caryophyllene β-Myrcene	3.24 1.81 1.23	32400 18100 12300

PASS/FAIL
PASS
PASS
PASS
PASS





Scan the QR code to verify results.

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Kyle W. Felling, Ph.D. aboratory Director

#### REPORT OF LABORATORY ANALYSIS

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## **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35050)**

Testing Location:	Customer ID: 2168	Order ID: OR10416	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
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License: ADH 113	License: 00065C	E2023KMTHTR01	Date Received: 04/24/2023
Cultivar (Strain) or Sample I	Date Completed: 05/02/2023		

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

Analysis Date/Time: 04/25/2023 1656 Analyst: PW Method: HPLC/DAD Instrument: Agilent 1100

<u>Cannabinoid</u>	<u>Result</u> (%)	<u>Result</u> (mg/g)	<u>LOD</u> ( <u>mg/g)</u>	LOQ (mg/g)	<u>Result</u> (mg/ S <u>mL)</u>	<u>Per</u> Serving (mg)	<u>Per</u> 5 <u>Unit</u> ( <u>mg)</u>
CBC	ND	ND	0.505	1.18	-	-	-
CBCA	1.14	11.4	1.57	3.66	-	11.4	11.4
CBD	5.02	50.2	3.56	8.31	-	50.2	50.2
CBDA	0.370	3.70	1.31	3.06	-	3.70	3.70
CBDV	ND	ND	0.571	1.33	-	-	-
CBDVA	ND	ND	1.52	3.56	-	-	-
CBG	0.605	6.05	2.31	5.40	-	6.05	6.05
CBGA	2.01	20.1	1.66	3.87	-	20.1	20.1
CBL	2.98	29.8	2.67	6.24	-	29.8	29.8
CBN	ND	ND	1.23	2.87	-	-	-
CBNA	ND	ND	1.32	3.09	-	-	-
Δ9-ΤΗC	21.7	217	1.47	3.43	-	217	217
$\Delta 8$ -THC	ND	ND	2.29	5.36	-	-	-
THCA	41.4	414	0.797	1.86	-	414	414
THCV	<loq< td=""><td><loq< td=""><td>1.91</td><td>4.46</td><td>-</td><td>-</td><td>-</td></loq<></td></loq<>	<loq< td=""><td>1.91</td><td>4.46</td><td>-</td><td>-</td><td>-</td></loq<>	1.91	4.46	-	-	-
THCVA	0.185	1.85	0.611	1.42	-	1.85	1.85
TOTAL	75.5	755			-	755	755
TOTAL CBC	1.00	10.0				10.0	10.0
TOTAL CBD	5.35	53.5			-	53.5	53.5
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	2.37	23.7			-	23.7	23.7
TOTAL CBN	-	-			-	-	-
TOTAL THC	58.0	580			-	580	580
TOTAL THCV	0.160	1.60			-	1.60	1.60

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

*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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Moisture Content (%): -Water Activity (aw): -



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) +  $\Delta$ 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.

Kyle W. Felling, Ph.D. atory Dire

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Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 04/24/2023
License: ADH 113	License: 00065C	E2023KMTHTR01	Date Received: 04/24/2023
Cultivar (Strain) or Sample	Date Completed: 05/02/2023		

#### **TERPENOID PROFILE**

Analysis Date/Time:04/25/2023 1942 Analyst: KF			ethod: GC/MS strument: Agilent 7890/5975	<b>Deviations from SOP:</b> None			
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)					
α-Bisabolol	2100	0.210					
Camphene	1390	0.139		Tank the second			
δ-3-Carene	-	-					
β-Caryophyllene	18100	1.81					
Caryophyllene oxide	-	-					
p-Cymene	622	0.0622		· 93			
Eucalyptol	286	0.0286					
Geraniol	-	-					
Guaiol	1160	0.116					
α-Humulene	6600	0.660		Abbreviations: GC - Gas Chromatography, MS - Mass			
Isopulegol	-	-		Spectrometry, RL - Reporting Limit			
d-Limonene	32400	3.24		This information is provided as a service and makes no claims of efficacy and/or			
Linalool	8540	0.854		safety of this product.			
β-Myrcene cis-Nerolidol	12300 -	1.23 -	•	Results are applicable only for the sample(s) analyzed and for the specific analysis conducted.			
trans-Nerolidol	2510	0.251	1	This report is for informational purposes			
α-Ocimene	239	0.0239		only and should not be used to diagnose, treat, or prevent any			
β-Ocimene	2270	0.227		medical-related symptoms.			
α-Pinene	4580	0.458	1	The statements and results herein have			
β-Pinene	5720	0.572	i	not been approved and/or endorsed by the FDA.			
' α-Terpinene	-	-	-				
γ-Terpinene	-	-					
Terpinolene	1600	0.160	1	"-" Not detected above RL.			
TOTAL	100000	10.0	-	Reporting Limit (µg/g): 4			













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License: ADH 113	License: 00065C	E2023KMTHTR01	Date Received: 04/24/2023
Cultivar (Strain) or Sample	Date Completed: 05/02/2023		

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

Analysis Date/Time: 05/01/2023 2235 Metho				Method: HS/GC/MS			D	eviation	s from SOP:	
Analyst: KF				1	Instrument: Agilent 78	90/5975	5	Ν	one	
<u>Solvent</u>	<u>Result</u> (µg/g)	LOD (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	<u>Solvent</u>	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	
Acetone (67-64-1)	-	34.5	69.1	5000	n-Heptane (142-82-5)	-	34.5	69.1	5000	
Acetonitrile (75-5-8)	-	34.5	69.1	410	n-Hexane (110-54-3)	-	12.1	24.2	290	
Benzene (71-43-2)	-	0.345	0.691	2	Isobutane (75-28-5)	-	34.5	69.1	5000	0 9
n-Butane (106-97-2)	2330	34.5	69.1	5000	Isopropanol (67-63-0)	-	34.5	69.1	5000	0.00
1-Butanol (71-36-3)	-	34.5	69.1	5000	Isopropyl acetate	_	34.5	69.1	5000	03
2-Butanol (78-92-2)	-	34.5	69.1	5000	(108-21-4)					
2-Butanone (78-93-3)	-	34.5	69.1	5000	Isopropyl benzene (98-82-8)	-	3.45	6.91	70	
Cyclohexane (110-82-7)	-	34.5	69.1	3880	Methanol (67-56-1)	_	34.5	69.1	3000	<u>Color Key</u>
1,2-Dimethoxyethane	_	3.45	6.91	100	2-Methylbutane (78-78-4)	_	34.5	69.1	5000	
(110-71-4) N,N-Dimethylacetamide (127-19-5)	_	34.5	69.1	1090	Methylene chloride (75-9-2)	-	34.5	69.1	600	RESULT < AL RESULT > AL
(127-13-3) 2,2-Dimethylbutane (75-83-2)	-	12.1	24.2	290	2-Methylpentane (107-83-5)	-	12.1	24.2	290	"DET" detected less than LOQ
2,3-Dimethylbutane					3-Methylpentane (96-10-0)	-	12.1	24.2	290	"-" not detected above
(79-29-8)	-	12.1	24.2	290	n-Pentane (109-66-0)	-	34.5	69.1	5000	LOD
N,N-Dimethylformamide		34.5	69.1	880	1-Pentanol (71-41-0)	-	34.5	69.1	5000	"*" - o,m,p-Xylene and
(68-12-2)		54.5	09.1	000	n-Propane (74-98-6)	-	34.5	69.1	5000	Ethylbenzene
Dimethylsulfoxide (67-68-5)	-	34.5	69.1	5000	1-Propanol (71-23-8)	-	34.5	69.1	5000	U U
(67-66-5) 1,4-Dioxane (123-91-1)		34.5	69.1	380	Pyridine (110-86-1)	-	12.1	24.2	200	Action levels are referenced from the State of
Ethanol (64-17-5)	-	34.5	69.1	5000	Tetrahydrofuran (109-99-9)	-	34.5	69.1	720	Arkansas
2-Ethoxyethanol (110-80-5)	-	12.1	24.2	160	Tetramethylene sulfone	_	12.1	24.2	160	MMJ testing
Ethyl ether (60-29-7)		34.5	69.1	5000	(126-33-0)					guidelines.
Ethyl acetate (141-78-6)	-	34.5	69.1	5000	Toluene (108-88-3)	-	34.5	69.1	890	A value of "-"
Ethyl benzene (100-41-4)	-	34.5	69.1	2170	o-Xylene (95-47-6)	-	34.5	69.1	2170	for the action level
Ethylene glycol (107-21-1)		34.5	69.1	620	m,p-Xylene (108-38-3 or 106-42-3)	-	34.5	69.1	2170	means that analyte
Ethylene oxide (75-21-8)		3.45	6.91	50	Xylenes* (1330-20-7)		43.3	86.7	2170	is not currently
		-		50				00.7	2170	regulated by the regulations referenced above.
<u>Solvent</u> Acetonitrile		Synonym(s Methyl Cya	-		<u>Solvent</u> Ethylene glycol		<b>Synonym(s)</b> 1,2-Ethanediol			
1-Butanol		n-Butanol,		hol	Isobutane		2-Methylpropa	20		
2-Butanol		sec-Butyl a	5	1101			2-Methylplopa 2-Propanol, IPA			
2-Butanone		Methyl ethy		1 F K	Isopropanol Isopropyl Acetate		Acetic acid iso			
1,2-Dimethoxyethane		Monoglym		iLIX	Methanol		Methyl alcohol			
2,3-Dimethylbutane		Neohexane			2-Methylbutane		lsopentane			
2,3-Dimethylbutane		Diisopropy			Methylene chloride		Dichlorometha	ne		
N,N-Dimethylformamide		DIISOPIOPY			2-Methylpentane		Isohexane	iic.		
Dimethysufoxide		DMF			1-Pentanol		n-Amyl alcoho	1		
2-Ethoxyethanol		Cellosolve,	Ethyl alve	ol	1-Propanol		Propyl alcohol	1		
Ethyl ether		Diethyl eth		01	Tetrahydrofuran		гюруг асоног ГНF			
Ethyl acetate		EtOAc	ei, Eulei		Tetramethylene sulfone		Sulfolane			
Ethyl benzene		Phenyletha	no		Xylene		Dimethylbenze	no		
Ethyl Delizene		rnenyietha	ne		лутене	1	onneuryibenze	ne		

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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**Deviations from SOP:** 

None

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License: ADH 113	License: 00065C	E2023KMTHTR01	Date Received: 04/24/2023
Cultivar (Strain) or Sample	<b>Date Completed:</b> 05/02/2023		

#### **PESTICIDES PROFILE (SOP: SOP-PEST-001)**

### Analysis Date/Time: 04/25/2023 1822 Analyst: KF

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

<u>Pesticide</u>	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	<u>Pesticide</u>	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	
Abamectin (71751-41-2)	-	0.0443	0.354	0.5	Kresoxim-methyl	_	0.0443	0.354	0.4	
Acephate (30560-19-1)	-	0.0443	0.354	0.4	(143390-89-0)					10.00
Acequinocyl (57960-19-7)	-	0.0443	0.354	2	Malathion (121-75-5)	-	0.0443	0.354	0.2	
Acetamiprid (135410-20-7)	-	0.0443	0.354	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.0443 0.0443	0.354 0.354	0.2 0.2	<u>Color Key</u>
Aldicarb (116-06-3)	-	0.0443	0.354	0.4	Methomyl (16752-77-5)	-	0.0443	0.354	0.4	<b>RESULT &lt; AL</b>
Azoxystrobin (131860-33-8)	-	0.0443	0.354	0.2	Methyl parathion (298-0-0)	-	0.0443	0.354	0.2	<b>RESULT &gt; AL</b>
Bifenazate (149877-41-8)	-	0.0443	0.354	0.2	MGK 264 (113-48-4)	-	0.0443	0.354	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3)	-	0.0443	0.354	0.2	Myclobutanil	_	0.0443	0.354	0.2	"-" not detected above
Boscalid (188425-85-6)	-	0.0443	0.354	0.4	(88671-89-0)					LOD
Carbaryl (63-25-2)	-	0.0443	0.354	0.2	Naled (300-76-5)	-	0.0443	0.354	0.5	Permethrins measured as the
Carbofuran (1563-66-2)	-	0.0443	0.354	0.2	Oxamyl (23135-22-0)	-	0.0443	0.354	1	cumulative residue of the <i>cis</i> - and
Chlorantraniliprole (800008-45-7)	-	0.0443	0.354	0.2	Paclobutrazol (76738-62-0)	-	0.0443	0.354	0.4	trans- permethrin isomers.
Chlorfenapyr		0.0443	0.354	1	Permethrins (52645-53-1)	-	0.0443	0.354	0.2	Pyrethrins measured as the cumulative residue of the
(122453-73-0)					Phosmet (732-11-6)	-	0.0443	0.354	0.2	pyrethrin I, cinerin I, and jasmolin
Chlorpyrifos (2921-88-2)	-	0.0443	0.354	0.2	Piperonyl butoxide (51-03-6)	-	0.0443	0.354	2	I isomers.
Clofentezine (74115-24-5)	-	0.0443	0.354	0.2	(31-03-0) Prallethrins (2331-36-9)	-	0.0443	0.354	0.2	Action levels are referenced from
Cyfluthrin (68359-37-5)	-	0.0443	0.354	1	Propiconazole					the
Cypermethrin (52315-07-8)	-	0.0443	0.354	1	(60207-90-1))	-	0.0443	0.354	0.4	State of Arkansas MMJ testing guidelines.
Daminozide (1596-84-5)	-	0.0443	0.354	1	Propoxur (114-26-1)	-	0.0443	0.354	0.2	A value of "-" for the action level
DDVP (62-73-7)	-	0.0443	0.354	0.1	Pyrethrins (8003-34-7)	-	0.0443	0.354	1	means that analyte is not
Diazinon (333-41-5)	-	0.0443	0.354	0.2	Pyridaben (96489-71-3)	-	0.0443	0.354	0.2	currently regulated by the regulations referenced above.
Dimethoate (60-51-5)	-	0.0443	0.354	0.2	Spinosad (168316-95-8)	-	0.0443	0.354	0.2	U U
Ethoprophos (13194-48-4)	-	0.0443	0.354	0.2	Spiromesifen (283594-90-1)	-	0.0443	0.354	0.2	<i>Disclaimer</i> : This information is provided as a service and makes
Etofenprox (80844-07-1)	-	0.0443	0.354	0.4	Spirotetramat		0.0440	0.054	0.0	no claims of efficacy and/or safety
Etoxazole (153233-91-1)	-	0.0443	0.354	0.2	(203313-25-1)	-	0.0443	0.354	0.2	of this product. Results are
Fenoxycarb (72490-01-8)	-	0.0443	0.354	0.2	Spiroxamine	_	0.0443	0.354	0.4	applicable only for the sample(s) analyzed and for the specific
(E)-Fenpyroximate (134098-61-6)	-	0.0443	0.354	0.4	(118134-30-8) Tebuconazole	_	0.0443	0.354	0.4	analysis conducted. This report is for informational purposes only
Fipronil (120068-37-3)	-	0.0443	0.354	0.4	(80443-41-0)		0.0445	0.004	0.4	and should not be used to
Flonicamid (158062-67-0)	-	0.0443	0.354	1	Thiacloprid (111988-49-9)	-	0.0443	0.354	0.2	diagnose, treat, or prevent any medical-related symptoms. The
Fludioxinil (131341-86-1)	-	0.0443	0.354	0.4	Thiamethoxam					statements and results herein have
Hexythiazox (78587-05-0)	-	0.0443	0.354	1	(153719-23-4)	-	0.0443	0.354	0.2	not been approved and/or endorsed by the FDA.
Imazalil (35554-44-0)	-	0.0443	0.354	0.2	Trifloxystrobin		0.0443	0.354	0.2	endorsed by the rDA.
Imidacloprid (138261-41-3)	-	0.0443	0.354	0.4	(141517-21-7)		0.0443	0.354	0.2	

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 Pesticide** Synonym(s) Cyfluthrin Baythroid Myclobutanil DDVP Dichlorvos Naled Ethoprophos Prophos Phosmet

Synonym(s) Systhane Dibrom Imidan

**Pesticide** Propiconazole Propoxur

Synonym(s) Tilt Baygon



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Arkansas	<b>River Valley Relief MIPS</b>	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13245725149	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	<b>Production Run:</b>	Date Collected: 04/24/2023
License: ADH 113	License: 00065C	E2023KMTHTR01	Date Received: 04/24/2023
Cultivar (Strain) or Sample Description: Kush Mints Hybrid HTFSE Resin .5g Cart			Date Completed: 05/02/2023

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

Analysis Date/T Analysis Date/T Analyst: KF	<b>ime:</b> 04/25/2023 1 <b>ime: -</b> (DMA)	747 (ICP/OES)	_	thod: ICP/OES trument: Agilent 720	<b>Deviations from SOP:</b> D-ES None
<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)	-	55.7	88.2	200	
Cadmium (Cd)	-	55.7	88.2	200	totte the second s
Lead (Pb)	-	55.7	88.2	500	
Mercury (Hg)	-	55.7	88.2	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

### <u>Color Key</u>



"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.













# **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35050)**

Testing Location:	Customer ID: 2168	Sample ID: SA35050	Sample Type: Primary
Arkansas	<b>River Valley Relief MIPS</b>	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13245725149	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 04/24/2023
License: ADA 05_H273	License: 00065C	E2023KMTHTR01	Date Received: 04/24/2023
Cultivar (Strain) or Sample Description: Kush Mints Hybrid HTFSE Resin .5g Cart			Date Completed: 05/02/2023

#### MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 2023042 Analyst: PW		Hardy Diagnostics CompactD <b>nt:</b> Thermo Incubator	ry <b>Deviations from SOP:</b> None
Bacteria/Microbe	<u>Result</u> (CFU/g)	<u>Action Level</u> (CFU/g)	
Aerobic Plate Count	NT	-	
Coliforms, Total	Absent	1	Latin Charles I and the
Escherichia Coli (E. Coli)	Absent	100	
Mold/Yeast	NT	-	
Pseudomonas aeruginosa	NT	1	
Salmonella spp.	NT	-	00
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





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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