



SUMMARY OF ANALYSIS (SAMPLE ID: SA39547)

Testing Location:	Customer ID: 2168	Order ID: OR11313	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13222392749	Mass: 1ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 12/02/2024
License: ADH 113	License: 00065C	P20241127GHEE12	Date Received: 12/02/2024
Cultivar (Strain) or Sample De	Date Completed:12/04/2024		

*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.

Cannabinoid	ts (Tan 3)	<u>(%)</u>	<u>mg</u> /g
<u>Δ</u> 9-TI		0.662	6.62
CBI		0.002	0.119
THC		0.0113	0.112
TOTAL		0.0112	0.112
TOTAL		0.662	6.62
TOTAL CANN		0.703	7.03
<u>Terpenes</u> α-Bisab		<u>(%)</u>	hā\ā
Camph			
δ-3-Cai			
β-Caryoph			
Caryophylle			
TOTAL TE			-
Contaminants	PASS/FAIL	Sample Picture	<u>Upon Receipt</u>
Heavy Metals:	PASS		
Microbiology:	PASS		
Pesticides:	PASS		
Residual Solvents:	PASS		



Scan the QR code to verify results.

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Kyle W. Felling, Ph.D. Laboratory Director

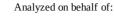
REPORT OF LABORATORY ANALYSIS

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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA39547)

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Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13222392749	Mass: 1ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 12/02/2024
License: ADH 113	Date Received: 12/02/2024		
Cultivar (Strain) or Sample I	Date Completed:12/04/2024		

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

Analysis Date/Time: 12/03/2024 1313 Analyst: PW Method: HPLC/DAD Instrument: Agilent 1100

<u>Cannabinoid</u>	<u>Result</u> (%)	<u>Result</u> (mg/g)	<u>LOD</u> (mg/g)	LOQ (mg/g)	<u>Result</u> (mg/ mL)	<u>Per</u> Serving (mg)	<u>Per</u> <u>Unit</u> (mg)
CBC	ND	ND	0.00226	0.00526	-	0.000	0.000
CBCA	ND	ND	0.00700	0.0163	-	0.000	0.000
CBD	0.0119	0.119	0.0159	0.0371	-	0.178	10.7
CBDA	ND	ND	0.00586	0.0137	-	0.000	0.000
CBDV	ND	ND	0.00255	0.00596	-	0.000	0.000
CBDVA	ND	ND	0.00681	0.0159	-	0.000	0.000
CBG	0.00909	0.0909	0.0103	0.0241	-	0.136	8.16
CBGA	ND	ND	0.0146	0.0173	-	0.000	0.000
CBL	ND	ND	0.0119	0.0279	-	0.000	0.000
CBN	0.00890	0.0890	0.00548	0.0128	-	0.134	8.04
CBNA	ND	ND	0.00592	0.0138	-	0.000	0.000
CBT	ND	ND	0.00859	0.0200	-	0.000	0.000
Δ9-ΤΗC	0.662	6.62	0.00657	0.0153	-	9.93	596
$\Delta 8$ -THC	ND	ND	0.0103	0.0239	-	0.000	0.000
THCA	ND	ND	0.00356	0.00833	-	0.000	0.000
THCV	0.0112	0.112	0.00855	0.0199	-	0.168	10.1
THCVA	ND	ND	0.00273	0.00635	-	0.000	0.000
TOTAL	0.703	7.03			-	10.5	630
TOTAL CBC	0.000	0.000			-	0.000	0.000
TOTAL CBD	0.0119	0.119			-	0.178	10.7
TOTAL CBDV	0.000	0.000			-	0.000	0.000
TOTAL CBG	0.00909	0.0909			-	0.136	8.16
TOTAL CBN	0.00890	0.0890			-	0.134	8.04
TOTAL THC	0.662	6.62			-	9.93	596
TOTAL THCV	0.0112	0.112			-	0.168	10.1

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



SERVING MASS (g): 1.50 SERVINGS/UNIT: 60

"-" 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) + Δ 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.



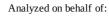
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Deviations from SOP:



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Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13222392749	Mass: 1ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 12/02/2024
License: ADH 113	License: 00065C	P20241127GHEE12	Date Received: 12/02/2024
Cultivar (Strain) or Sample De	Date Completed:12/04/2024		

Analysis Date/Time:12/03/2024 1811 Analyst: KF

TERPENOID PROFILE
Method: GC/MS

Analyst: KF		Instrument: Agilent 7890/5975	None
<u>Terpene</u>	<u>Result</u> (µg/g)	Result (%)	
α-Bisabolol	<loq< td=""><td>-</td><td></td></loq<>	-	
Camphene	ND	-	
δ-3-Carene	ND	-	
β-Caryophyllene	<loq< td=""><td>-</td><td></td></loq<>	-	
Caryophyllene oxide	ND	-	
p-Cymene	ND	-	
Eucalyptol	ND	-	
Geraniol	ND	-	
Guaiol	ND	-	Abbreviations: GC - Gas
α-Humulene	ND	-	Chromatography, MS - Mass
Isopulegol	ND	-	Spectrometry, RL - Reporting Limit
d-Limonene	<loq< td=""><td>-</td><td><i>Abbreviations:</i> ND - Not Detected, , LOD - Limit of Detection, LOQ - Limit</td></loq<>	-	<i>Abbreviations:</i> ND - Not Detected, , LOD - Limit of Detection, LOQ - Limit
Linalool	<loq< td=""><td>-</td><td>of Quantitation</td></loq<>	-	of Quantitation
β-Myrcene	<loq< td=""><td>-</td><td>This information is provided as a service and makes no claims of efficacy and/or</td></loq<>	-	This information is provided as a service and makes no claims of efficacy and/or
cis-Nerolidol	ND	-	safety of this product.
trans-Nerolidol	ND	-	Results are applicable only for the sample(s) analyzed and for the specific
α-Ocimene	ND	-	analysis conducted.
β-Ocimene	ND	-	This report is for informational purposes only and should not be used to diagnose,
α-Pinene	ND	-	treat, or prevent any
β-Pinene	ND	-	medical-related symptoms.
α-Terpinene	ND	-	The statements and results herein have not been approved and/or endorsed by
γ-Terpinene	ND	-	the FDA.
Terpinolene	ND	-	
TOTAL	0.000	0.000	Reporting Limit (µg/g): 4.

Reporting Limit (µg/g): 4.95

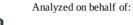
"-" Not detected above LOD.













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Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
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Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 12/02/2024
License: ADH 113	License: 00065C	P20241127GHEE12	Date Received: 12/02/2024
Cultivar (Strain) or Sample	Date Completed:12/04/2024		

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

Analysis Date/Time: Analyst: KF	Analysis Date/Time: 12/02/2024 1555Method: HS/GC/MSAnalyst: KFInstrument: Agilent 7890/5975		Deviations from SOP: 75 None							
Solvent	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	Solvent	<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)	-	29.1	58.1	5000	n-Heptane (142-82-5)	-	29.1	58.1	5000	
Acetonitrile (75-5-8)	_	29.1	58.1	410	n-Hexane (110-54-3)	-	10.2	20.3	290	
Benzene (71-43-2)	-	0.291	0.581	2	Isobutane (75-28-5)	-	29.1	58.1	5000	
n-Butane (106-97-2)	_	29.1	58.1	5000	Isopropanol (67-63-0)	-	29.1	58.1	5000	
1-Butanol (71-36-3)	-	29.1	58.1	5000	Isopropyl acetate		20.1	F0 1	5000	
2-Butanol (78-92-2)	-	29.1	58.1	5000	(108-21-4)	-	29.1	58.1	5000	
2-Butanone (78-93-3)	-	29.1	58.1	5000	Isopropyl benzene	_	2.91	5.81	70	
Cyclohexane (110-82-7)	-	29.1	58.1	3880	(98-82-8)					Color Key
1,2-Dimethoxyethane	_	2.91	5.81	100	Methanol (67-56-1) 2-Methylbutane (78-78-4)	-	29.1 29.1	58.1 58.1	3000 5000	
(110-71-4) N,N-Dimethylacetamide	_	29.1	58.1	1090	Methylene chloride (75-9-2)	-	29.1	58.1	600	RESULT < AL RESULT > AL
(127-19-5) 2,2-Dimethylbutane (75-83-2)	-	10.2	20.3	290	2-Methylpentane (107-83-5)	-	10.2	20.3	290	"-" not detected above 1/2 Action Level
2,3-Dimethylbutane					3-Methylpentane (96-10-0)	-	10.2	20.3	290	
(79-29-8)	-	10.2	20.3	290	n-Pentane (109-66-0)	-	29.1	58.1	5000	"*" - o,m,p-Xylene and
N,N-Dimethylformamide	_	29.1	58.1	880	1-Pentanol (71-41-0)	-	29.1	58.1	5000	Ethylbenzene
(68-12-2)		25.1	50.1	000	n-Propane (74-98-6)	-	29.1	58.1	5000	Action levels are
Dimethylsulfoxide (67-68-5)	-	29.1	58.1	5000	1-Propanol (71-23-8)	-	29.1	58.1	5000	referenced from the State of
1,4-Dioxane (123-91-1)	_	29.1	58.1	380	Pyridine (110-86-1)	-	10.2	20.3	200	Arkansas MMJ testing
Ethanol (64-17-5)	_	29.1	58.1	5000	Tetrahydrofuran (109-99-9)	-	29.1	58.1	720	guidelines.
2-Ethoxyethanol (110-80-5)	-	10.2	20.3	160	Tetramethylene sulfone	-	10.2	20.3	160	0
Ethyl ether (60-29-7)	_	29.1	58.1	5000	(126-33-0) Toluene (108-88-3)		29.1	58.1	890	A value of "-" for the action level
Ethyl acetate (141-78-6)	-	29.1	58.1	5000	o-Xylene (95-47-6)	-	29.1	58.1	2170	means that analyte
Ethyl benzene (100-41-4)	-	29.1	58.1	2170	m,p-Xylene (108-38-3 or	-				is not currently
Ethylene glycol (107-21-1)	-	29.1	58.1	620	106-42-3)	-	29.1	58.1	2170	regulated by the
Ethylene oxide (75-21-8)	-	2.91	5.81	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170	regulations referenced above.
Solvent		Synonym(s	<u>5)</u>		Solvent	<u>S</u>	ynonym(s)			
Acetonitrile		Methyl Cya	nide, ACN		Ethylene glycol	1,	,2-Ethanediol			
1-Butanol		n-Butanol,	Butyl Alco	hol	Isobutane	2.	-Methylpropa	ne		
2-Butanol		sec-Butyl a	lcohol		Isopropanol	2.	-Propanol, IPA	A		
2-Butanone		Methyl ethy	yl ketone, N	/IEK	Isopropyl Acetate	А	cetic acid iso	propyl este	r	
1,2-Dimethoxyethane		Monoglym	e		Methanol	Ν	fethyl alcohol			
2,3-Dimethylbutane		Neohexane			2-Methylbutane	Is	opentane			
2,3-Dimethylbutane		Diisopropy	rl -		Methylene chloride	D	ichlorometha	ne		
N,N-Dimethylformamide		DMF			2-Methylpentane		ohexane			
Dimethysufoxide		DMSO			1-Pentanol		-Amyl alcoho			
2-Ethoxyethanol		Cellosolve,		ol	1-Propanol		ropyl alcohol			
Ethyl ether		Diethyl eth	er, Ether		Tetrahydrofuran		HF			
Ethyl acetate		EtOAc			Tetramethylene sulfone	S	ulfolane			

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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Xylene

Phenylethane



Ethyl benzene

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Dimethylbenzene



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Deviations from SOP:

None



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Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 12/02/2024
License: ADH 113	License: 00065C	P20241127GHEE12	Date Received: 12/02/2024
Cultivar (Strain) or Sample D	Date Completed:12/04/2024		

Analysis Date/Time: 12/03/2024 1329 Analyst: KF

PESTICIDES PROFILE (SOP: SOP-PEST-001)

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)	-	0.00993	0.0794	0.5	Kresoxim-methyl	_	0.00993	0.0794	0.4	
Acephate (30560-19-1)	-	0.00993	0.0794	0.4	(143390-89-0)		0.00000	0.0504		
Acequinocyl (57960-19-7)	-	0.00993	0.0794	2	Malathion (121-75-5)	-	0.00993	0.0794	0.2	
Acetamiprid (135410-20-7)	-	0.00993	0.0794	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.00993 0.00993	0.0794 0.0794	0.2 0.2	<u>Color Key</u>
Aldicarb (116-06-3)	-	0.00993	0.0794	0.4	Methomyl (16752-77-5)	-	0.00993	0.0794	0.4	RESULT < AL
Azoxystrobin (131860-33-8)	-	0.00993	0.0794	0.2	Methyl parathion (298-0-0)	-	0.00993	0.0794	0.2	RESULT > AL
Bifenazate (149877-41-8)	-	0.00993	0.0794	0.2	MGK 264 (113-48-4)	-	0.00993	0.0794	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3)	-	0.00993	0.0794	0.2	Myclobutanil		0.00993	0.0794	0.2	"-" not detected above
Boscalid (188425-85-6)	-	0.00993	0.0794	0.4	(88671-89-0)	-	0.00993	0.0794	0.2	- not detected above LOD
Carbaryl (63-25-2)	-	0.00993	0.0794	0.2	Naled (300-76-5)	-	0.00993	0.0794	0.5	-
Carbofuran (1563-66-2)	-	0.00993	0.0794	0.2	Oxamyl (23135-22-0)	-	0.00993	0.0794	1	Permethrins measured as the cumulative residue of the <i>cis</i> - and
Chlorantraniliprole (800008-45-7)	-	0.00993	0.0794	0.2	Paclobutrazol (76738-62-0)	-	0.00993	0.0794	0.4	trans- permethrin isomers.
Chlorfenapyr (122453-73-0)	-	0.00993	0.0794	1	Permethrins (52645-53-1)	-	0.00993	0.0794	0.2	Pyrethrins measured as the cumulative residue of the
(chlorpyrifos (2921-88-2)	_	0.00993	0.0794	0.2	Phosmet (732-11-6)	-	0.00993	0.0794	0.2	pyrethrin I, cinerin I, and jasmolin I isomers.
Clofentezine (74115-24-5)	_	0.00993	0.0794	0.2	Piperonyl butoxide	_	0.00993	0.0794	2	
Cyfluthrin (68359-37-5)	-	0.00993	0.0794	1	(51-03-6)					Action levels are referenced from the State of
Cypermethrin (52315-07-8)	-	0.00993	0.0794	1	Prallethrins (2331-36-9) Propiconazole	-	0.00993 0.00993	0.0794 0.0794	0.2 0.4	Arkansas MMJ testing
Daminozide (1596-84-5)	-	0.00993	0.0794	1	(60207-90-1))					guidelines.
DDVP (62-73-7)	_	0.00993	0.0794	0.1	Propoxur (114-26-1)	-	0.00993	0.0794	0.2	A value of "-" for the action level
Diazinon (333-41-5)	_	0.00993	0.0794	0.2	Pyrethrins (8003-34-7)	-	0.00993	0.0794	1	means that analyte is not
Dimethoate (60-51-5)	_	0.00993	0.0794	0.2	Pyridaben (96489-71-3)	-	0.00993	0.0794	0.2	currently regulated by the
Ethoprophos	-	0.00993	0.0794	0.2	Spinosad (168316-95-8) Spiromesifen	-	0.00993	0.0794	0.2	regulations referenced above.
(13194-48-4)			0.0704		(283594-90-1)	-	0.00993	0.0794	0.2	<i>Disclaimer</i> : This information is provided as a service and makes
Etofenprox (80844-07-1)	-	0.00993	0.0794	0.4	Spirotetramat	_	0.00993	0.0794	0.2	no claims of efficacy and/or safety
Etoxazole (153233-91-1)	-	0.00993	0.0794	0.2	(203313-25-1)		0.00555	0.0734	0.2	of this product. Results are applicable only for the sample(s)
Fenoxycarb (72490-01-8)	-	0.00993	0.0794	0.2	Spiroxamine (118134-30-8)	-	0.00993	0.0794	0.4	analyzed and for the specific
(E)-Fenpyroximate (134098-61-6)	-	0.00993	0.0794	0.4	Tebuconazole	_	0.00993	0.0794	0.4	analysis conducted. This report is for informational purposes only
Fipronil (120068-37-3)	-	0.00993	0.0794	0.4	(80443-41-0)					and should not be used to diagnose, treat, or prevent any
Flonicamid (158062-67-0)	-	0.00993	0.0794	1	Thiacloprid (111988-49-9)	-	0.00993	0.0794	0.2	medical-related symptoms. The
Fludioxinil (131341-86-1)	-	0.00993	0.0794	0.4	Thiamethoxam					statements and results herein have
Hexythiazox (78587-05-0)	-	0.00993	0.0794	1	(153719-23-4)	-	0.00993	0.0794	0.2	not been approved and/or endorsed by the FDA.
Imazalil (35554-44-0)	-	0.00993	0.0794	0.2	Trifloxystrobin		0.00993	0.0794	0.2	chubiscu by the i DA.
Imidacloprid (138261-41-3)	-	0.00993	0.0794	0.4	(141517-21-7)	-	0.00993	0.0794	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 Cyfluthrin DDVP Ethoprophos

Synonym(s) Baythroid Dichlorvos Prophos

Pesticide Myclobutanil Naled Phosmet

Synonym(s) Systhane Dibrom Imidan

Pesticide Propiconazole Propoxur

Synonym(s) Tilt Baygon



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Analyzed on behalf of:



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA39547)

Testing Location:	Customer ID: 2168	Order ID: OR11313	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13222392749	Mass: 1ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 12/02/2024
License: ADH 113	License: 00065C	P20241127GHEE12	Date Received: 12/02/2024
Cultivar (Strain) or Sample Description: ARV-Infused Ghee Butter 600mg Jar			Date Completed:12/04/2024

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

Analysis Date/Time: 12/03/2024 1748		Method: ICP/MS Instrument: Agilent 7500ce		Deviations from SOP: None	
Analyst: KF				0	
<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)	-	59.5	94.2	200	
Cadmium (Cd)	-	59.5	94.2	200	
Lead (Pb)	-	59.5	94.2	500	
Mercury (Hg)	-	59.5	94.2	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





"DET" detected less than LOQ

"-" not detected above LOD

Action levels for heavy metals are referenced from the State of Arkansas MMJ testing guidelines.

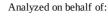
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MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 12/02/20 Analyst: PW		Hardy Diagnostics Compact nt: Thermo Incubator	Dry Deviations from SOP: None
Bacteria/Microbe	<u>Result</u> (CFU/g)	<u>Action Level</u> (<u>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





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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