







SUMMARY OF ANALYSIS (SAMPLE ID: SA35054)

Testing Location:	Customer ID: 2168	Order ID: OR10416	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13241421739	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 04/24/2023
License: ADH 113	License: 00065C	E20230418O43HTR01	Date Received: 04/24/2023
Cultivar (Strain) or Sample D	Date Completed: 04/27/2023		

*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.

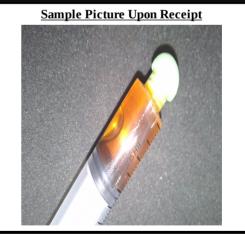
Moisture Content (%)	Water Activity (aw)	PASS/FAIL	
Not Tested	Not Tested	PASS	

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.

<u>Cannabinoids (Top 3)</u>	<u>(%)</u>	<u>mg/g</u>
THCA	33.6	336
Δ9-ТНС	29.4	294
CBD	5.17	51.7
TOTAL CBD	5.17	51.7
TOTAL THC	58.9	589
TOTAL CANNABINOIDS	76.2	762
<u>Terpenes (Top 5)</u>	<u>(%)</u>	µg∕g
<u>Terpenes (Top 5)</u> Terpinolene	<u>(%)</u> 1.78	µg/g 17800
Terpinolene	1.78	17800
Terpinolene β-Myrcene	1.78 0.814	17800 8140
Terpinolene β-Myrcene β-Caryophyllene	1.78 0.814 0.683	17800 8140 6830

PASS/FAIL
PASS
PASS
PASS
PASS





Scan the QR code to verify results.

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.



REPORT OF LABORATORY ANALYSIS

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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35054)

Testing Location:	Customer ID: 2168	Order ID: OR10416	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
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Cultivar (Strain) or Sample	Date Completed: 04/27/2023		

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

Analysis Date/Time: 04/25/2023 1739 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/ <u>mL)</u>	<u>Per</u> <u>Serving</u> (mg)	<u>Per</u> g <u>Unit</u> (mg)
CBC	0.176	1.76	0.558	1.30	-	1.76	1.76
CBCA	ND	ND	1.73	4.04	-	-	-
CBD	5.17	51.7	3.93	9.18	-	51.7	51.7
CBDA	<loq< td=""><td><loq< td=""><td>1.45</td><td>3.38</td><td>-</td><td>-</td><td>-</td></loq<></td></loq<>	<loq< td=""><td>1.45</td><td>3.38</td><td>-</td><td>-</td><td>-</td></loq<>	1.45	3.38	-	-	-
CBDV	ND	ND	0.631	1.47	-	-	-
CBDVA	ND	ND	1.68	3.93	-	-	-
CBG	0.953	9.53	2.55	5.96	-	9.53	9.53
CBGA	3.37	33.7	1.83	4.27	-	33.7	33.7
CBL	2.75	27.5	2.95	6.89	-	27.5	27.5
CBN	0.529	5.29	1.36	3.17	-	5.29	5.29
CBNA	ND	ND	1.46	3.41	-	-	-
Δ9-ΤΗC	29.4	294	1.62	3.79	-	294	294
Δ8-THC	ND	ND	2.53	5.92	-	-	-
THCA	33.6	336	0.881	2.06	-	336	336
THCV	<loq< td=""><td><loq< td=""><td>2.11</td><td>4.93</td><td>-</td><td>-</td><td>-</td></loq<></td></loq<>	<loq< td=""><td>2.11</td><td>4.93</td><td>-</td><td>-</td><td>-</td></loq<>	2.11	4.93	-	-	-
THCVA	0.273	2.73	0.675	1.57	-	2.73	2.73
TOTAL	76.2	762			-	762	762
TOTAL CBC	0.176	1.76			-	1.76	1.76
TOTAL CBD	5.17	51.7			-	51.7	51.7
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	3.91	39.1			-	39.1	39.1
TOTAL CBN	0.529	5.29			-	5.29	5.29
TOTAL THC	58.9	589			-	589	589
TOTAL THCV	0.237	2.37			-	2.37	2.37

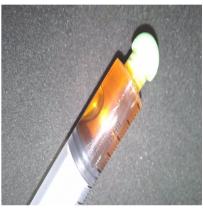
* CBCA, CBDVA, CBL, CBNA, and THCVA are in process of being added to our accreditation scope.

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

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.

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) + 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.

Kyle W. Felling, Ph.D. atory Dire



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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35054)

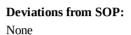
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License: ADH 113	License: 00065C	E20230418O43HTR01	Date Received: 04/24/2023
Cultivar (Strain) or Sample	Date Completed: 04/27/2023		

Analysis Date/Time:04/25/2023 2109 Analyst: KF

TERPENOID PROFILE

Method: GC/MS Instrument: Agilent 7890/5975

Analyst. M		111	strument. Agricut 7030/3373	INOL		
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)				
α-Bisabolol	2530	0.253				
Camphene	712	0.0712				
δ-3-Carene	1020	0.102				
β-Caryophyllene	6830	0.683				
Caryophyllene oxide	-	-				
p-Cymene	632	0.0632				
Eucalyptol	511	0.0511				
Geraniol	-	-				
Guaiol	999	0.0999		Abburictions		
α-Humulene	2600	0.260		Abbreviations: G Chromatography		
Isopulegol	-	-		Spectrometry, RI		
d-Limonene	5840	0.584		This information and makes no cla		
Linalool	2540	0.254	T	safety of this pro		
β-Myrcene	8140	0.814		Results are applies sample(s) analyz		
cis-Nerolidol	-	-		analysis conducte		
trans-Nerolidol	1060	0.106	1	This report is for		
α-Ocimene	-	-		only and should treat, or prevent a		
β-Ocimene	2590	0.259	1	medical-related s		
α-Pinene	2070	0.207	1	The statements an not been approve		
β-Pinene	2850	0.285		the FDA.		
α-Terpinene	1350	0.135	Ī			
γ-Terpinene	1250	0.125				
Terpinolene	17800	1.78		"-" Not detected		
TOTAL	61400	6.14				





GC - Gas ny, MS - Mass RL - Reporting Limit

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or informational purposes d not be used to diagnose, any

symptoms.

and results herein have ved and/or endorsed by

> d above RL Reporting Limit (µg/g): 49.0

4) Kyle W. Felling, Ph.D. aboratory Director











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License: ADH 113	License: 00065C	E20230418O43HTR01	Date Received: 04/24/2023
Cultivar (Strain) or Sample	Date Completed: 04/27/2023		

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

Analysis Date/Time: 04/26/2023 1210			Ι	Method: HS/GC/MS			Deviations from SOP:			
Analyst: KF				1	nstrument: Agilent 78	90/5975	5 None			
<u>Solvent</u>	<u>Result</u> (µg/g)	<u>LOD</u> (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)	\land
Acetone (67-64-1)	-	36.2	72.5	5000	n-Heptane (142-82-5)	-	36.2	72.5	5000	\checkmark
Acetonitrile (75-5-8)	-	36.2	72.5	410	n-Hexane (110-54-3)	-	12.7	25.4	290	
Benzene (71-43-2)	-	0.362	0.725	2	Isobutane (75-28-5)	-	36.2	72.5	5000	
n-Butane (106-97-2)	1950	36.2	72.5	5000	Isopropanol (67-63-0)	-	36.2	72.5	5000	
1-Butanol (71-36-3)	-	36.2	72.5	5000	Isopropyl acetate	_	36.2	72.5	5000	
2-Butanol (78-92-2)	-	36.2	72.5	5000	(108-21-4)		50.2	/2.0	5000	
2-Butanone (78-93-3)	-	36.2	72.5	5000	Isopropyl benzene (98-82-8)	-	3.62	7.25	70	
Cyclohexane (110-82-7)	-	36.2	72.5	3880	(56-62-6) Methanol (67-56-1)		36.2	72.5	3000	Color Key
1,2-Dimethoxyethane (110-71-4)	-	3.62	7.25	100	2-Methylbutane (78-78-4)	-	36.2	72.5	5000	RESULT < AL
N,N-Dimethylacetamide (127-19-5)	-	36.2	72.5	1090	Methylene chloride (75-9-2)	-	36.2	72.5	600	RESULT > AL
(2,2-Dimethylbutane (75-83-2)	-	12.7	25.4	290	2-Methylpentane (107-83-5)	-	12.7	25.4	290	"DET" detected less than LOQ
2,3-Dimethylbutane		10 5		200	3-Methylpentane (96-10-0)	-	12.7	25.4	290	"-" not detected above
(79-29-8)	-	12.7	25.4	290	n-Pentane (109-66-0)	-	36.2	72.5	5000	LOD
N,N-Dimethylformamide		36.2	72.5	880	1-Pentanol (71-41-0)	-	36.2	72.5	5000	"*" - o,m,p-Xylene and
(68-12-2)		50.2	/2.0	000	n-Propane (74-98-6)	-	36.2	72.5	5000	Ethylbenzene
Dimethylsulfoxide (67-68-5)	-	36.2	72.5	5000	1-Propanol (71-23-8)	-	36.2	72.5	5000	
1,4-Dioxane (123-91-1)		36.2	72.5	380	Pyridine (110-86-1)	-	12.7	25.4	200	Action levels are referenced from the State of
Ethanol (64-17-5)		36.2	72.5	5000	Tetrahydrofuran (109-99-9)	-	36.2	72.5	720	Arkansas
2-Ethoxyethanol (110-80-5)		12.7	25.4	160	Tetramethylene sulfone	_	12.7	25.4	160	MMJ testing
Ethyl ether (60-29-7)	_	36.2	72.5	5000	(126-33-0)					guidelines.
Ethyl acetate (141-78-6)	_	36.2	72.5	5000	Toluene (108-88-3)	-	36.2	72.5	890	A value of "-"
Ethyl benzene (100-41-4)	_	36.2	72.5	2170	o-Xylene (95-47-6)		36.2	72.5	2170	for the action level
Ethylene glycol (107-21-1)	_	36.2	72.5	620	m,p-Xylene (108-38-3 or 106-42-3)	-	36.2	72.5	2170	means that analyte
Ethylene oxide (75-21-8)	-	3.62	7.25	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170	is not currently regulated by the regulations referenced above.
Solvent		Synonym(s	5)		Solvent	<u>s</u>	ynonym(s)			regulations referenced above.
Acetonitrile		Methyl Cya	anide, ACN		Ethylene glycol	1	,2-Ethanedio	l		
1-Butanol		n-Butanol,	Butyl Alco	hol	Isobutane	2	-Methylpropa	ane		
2-Butanol		sec-Butyl a	lcohol		Isopropanol	2	-Propanol, IP.	4		
2-Butanone		Methyl ethy	yl ketone, N	⁄IEK	Isopropyl Acetate	A	Acetic acid iso	propyl este	r	
1,2-Dimethoxyethane		Monoglym	e		Methanol	Ν	Aethyl alcoho	1		
2,3-Dimethylbutane		Neohexane			2-Methylbutane	I	sopentane			
2,3-Dimethylbutane		Diisopropy	7l		Methylene chloride	Γ	Dichlorometha	ane		
N,N-Dimethylformamide		DMF			2-Methylpentane	I	sohexane			
Dimethysufoxide		DMSO			1-Pentanol	n	-Amyl alcoho	ol		
2-Ethoxyethanol		Cellosolve,	, Ethyl glyc	ol	1-Propanol	P	ropyl alcohol			
Ethyl ether		Diethyl eth	er, Ether		Tetrahydrofuran	Г	THF			
Ethyl acetate		EtOAc			Tetramethylene sulfone	S	Sulfolane			
Ethyl benzene		Phenyletha	ne		Xylene	Γ	Dimethylbenz	ene		

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

CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35054)

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Analysis Date/Time: 04/25/2023 1924 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)	-	0.0490	0.392	0.5	Kresoxim-methyl	_	0.0490	0.392	0.4	
Acephate (30560-19-1)	-	0.0490	0.392	0.4	(143390-89-0)					
Acequinocyl (57960-19-7)	-	0.0490	0.392	2	Malathion (121-75-5)	-	0.0490	0.392	0.2	
Acetamiprid (135410-20-7)	-	0.0490	0.392	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.0490 0.0490	0.392 0.392	0.2 0.2	<u>Color Key</u>
Aldicarb (116-06-3)	-	0.0490	0.392	0.4	Methomyl (16752-77-5)	-	0.0490	0.392	0.4	RESULT < AL
Azoxystrobin (131860-33-8)	-	0.0490	0.392	0.2	Methyl parathion (298-0-0)	-	0.0490	0.392	0.2	RESULT > AL
Bifenazate (149877-41-8)	-	0.0490	0.392	0.2	MGK 264 (113-48-4)	-	0.0490	0.392	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3)	-	0.0490	0.392	0.2	Myclobutanil	_	0.0490	0.392	0.2	"-" not detected above
Boscalid (188425-85-6)	-	0.0490	0.392	0.4	(88671-89-0)					LOD
Carbaryl (63-25-2)	-	0.0490	0.392	0.2	Naled (300-76-5)	-	0.0490	0.392	0.5	Permethrins measured as the
Carbofuran (1563-66-2)	-	0.0490	0.392	0.2	Oxamyl (23135-22-0)	-	0.0490	0.392	1	cumulative residue of the <i>cis</i> - and
Chlorantraniliprole (800008-45-7)	-	0.0490	0.392	0.2	Paclobutrazol (76738-62-0)	-	0.0490	0.392	0.4	trans- permethrin isomers.
Chlorfenapyr		0.0490	0.392	1	Permethrins (52645-53-1)	-	0.0490	0.392	0.2	Pyrethrins measured as the cumulative residue of the
(122453-73-0)	-	0.0490	0.392	1	Phosmet (732-11-6)	-	0.0490	0.392	0.2	pyrethrin I, cinerin I, and jasmolin
Chlorpyrifos (2921-88-2)	-	0.0490	0.392	0.2	Piperonyl butoxide	_	0.0490	0.392	2	I isomers.
Clofentezine (74115-24-5)	-	0.0490	0.392	0.2	(51-03-6)		0.0490	0.392	0.2	Action levels are referenced from
Cyfluthrin (68359-37-5)	-	0.0490	0.392	1	Prallethrins (2331-36-9)	-	0.0490	0.392	0.2	the
Cypermethrin (52315-07-8)	-	0.0490	0.392	1	Propiconazole (60207-90-1))	-	0.0490	0.392	0.4	State of Arkansas MMJ testing guidelines.
Daminozide (1596-84-5)	-	0.0490	0.392	1	Propoxur (114-26-1)	-	0.0490	0.392	0.2	A value of "-" for the action level
DDVP (62-73-7)	-	0.0490	0.392	0.1	Pyrethrins (8003-34-7)	-	0.0490	0.392	1	means that analyte is not
Diazinon (333-41-5)	-	0.0490	0.392	0.2	Pyridaben (96489-71-3)	-	0.0490	0.392	0.2	currently regulated by the
Dimethoate (60-51-5)	-	0.0490	0.392	0.2	Spinosad (168316-95-8)	-	0.0490	0.392	0.2	regulations referenced above.
Ethoprophos (13194-48-4)	-	0.0490	0.392	0.2	Spiromesifen (283594-90-1)	-	0.0490	0.392	0.2	Disclaimer: This information is
Etofenprox (80844-07-1)	-	0.0490	0.392	0.4	(283394-90-1) Spirotetramat					provided as a service and makes no claims of efficacy and/or safety
Etoxazole (153233-91-1)	-	0.0490	0.392	0.2	(203313-25-1)	-	0.0490	0.392	0.2	of this product. Results are
Fenoxycarb (72490-01-8)	-	0.0490	0.392	0.2	Spiroxamine		0.0490	0.392	0.4	applicable only for the sample(s) analyzed and for the specific
(E)-Fenpyroximate (134098-61-6)	-	0.0490	0.392	0.4	(118134-30-8) Tebuconazole	-	0.0490	0.392		analysis conducted. This report is for informational purposes only
Fipronil (120068-37-3)	-	0.0490	0.392	0.4	(80443-41-0)	-	0.0490	0.392	0.4	and should not be used to
Flonicamid (158062-67-0)	-	0.0490	0.392	1	Thiacloprid	_	0.0490	0.392	0.2	diagnose, treat, or prevent any medical-related symptoms. The
Fludioxinil (131341-86-1)	-	0.0490	0.392	0.4	(111988-49-9)					statements and results herein have
Hexythiazox (78587-05-0)	-	0.0490	0.392	1	Thiamethoxam (153719-23-4)	-	0.0490	0.392	0.2	not been approved and/or
Imazalil (35554-44-0)	-	0.0490	0.392	0.2	Trifloxystrobin					endorsed by the FDA.
Imidacloprid (138261-41-3)	-	0.0490	0.392	0.4	(141517-21-7)	-	0.0490	0.392	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	<u>Synonym(s)</u>	Pesticide	Synonym(s)	Pesticide	Synonym
Cyfluthrin	Baythroid	Myclobutanil	Systhane	Propiconazole	Tilt
DDVP	Dichlorvos	Naled	Dibrom	Propoxur	Baygon
Ethoprophos	Prophos	Phosmet	Imidan		



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/**m(s)**









CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35054)

Testing Location:	Customer ID: 2168	Order ID: OR10416	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13241421739	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 04/24/2023
License: ADH 113	License: 00065C	E20230418O43HTR01	Date Received: 04/24/2023
Cultivar (Strain) or Sample	Date Completed: 04/27/2023		

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

Analysis Date/Time: 04/25/2023 1835 (ICP/OES) Analysis Date/Time: - (DMA) Analyst: KF		Method: ICP/OES Instrument: Agilent 720-ES		-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)	-	58.3	92.3	200	\sim
Cadmium (Cd)	-	58.3	92.3	200	
Lead (Pb)	-	58.3	92.3	500	
Mercury (Hg)	-	58.3	92.3	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.

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: SA35054)

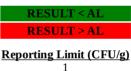
Testing Location:	Customer ID: 2168	Sample ID: SA35054	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13241421739	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 04/24/2023
License: ADA 05_H273	License: 00065C	E20230418O43HTR01	Date Received: 04/24/2023
Cultivar (Strain) or Sample D	Date Completed: 04/27/2023		

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 2023042 Analyst: PW		ardy Diagnostics CompactI t: 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	-	\sim
Coliforms, Total	Absent	1	
Escherichia Coli (E. Coli)	Absent	100	
Mold/Yeast	NT	-	
Pseudomonas aeruginosa	NT	1	R SY
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





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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