







SUMMARY OF ANALYSIS (SAMPLE ID: SA31889)

Testing Location:	Customer ID: 2168	Order ID: OR9915	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13241641866	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Batch Number:	Date Collected: 07/07/2022
License: ADH 113	License: 00065C	Not Entered	Date Received: 07/08/2022
Cultivar (Strain) or Sample Des	Date Completed: 07/10/2022		

*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.

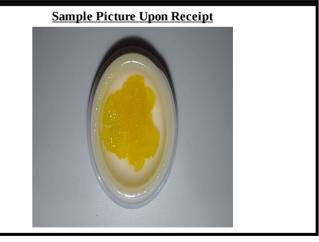
Moisture Content (%)	Water Activity (aw)	PASS/FAIL
Not Tested	Not Tested	N/A

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	80.4	804
CBGA	2.07	20.7
THCVA	0.687	6.87
TOTAL CBD	-	-
TOTAL THC	70.5	705
TOTAL CANNABINOIDS	83.2	832
<u>Terpenes (Top 5)</u>	<u>(%)</u>	<u>µg∕</u> g
<u>Terpenes (Top 5)</u> Terpinolene	<u>(%)</u> 1.14	<mark>µg/g</mark> 11400
Terpinolene	1.14	11400
Terpinolene β-Myrcene	1.14 0.834	11400 8340
Terpinolene β-Myrcene β-Ocimene	1.14 0.834 0.395	11400 8340 3950

<u>Contaminants</u>	PASS/FAIL
Heavy Metals:	PASS
Microbiology:	PASS
Pesticides:	PASS
Residual Solvents:	PASS





Scan the QR code to verify results.

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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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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA31889)

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Cultivar (Strain) or Sample D	Date Completed: 07/10/2022		

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

Analysis Date/Time: 7/8/2022 1826 Analyst: PW Method: HPLC/DAD Instrument: Agilent 1100

<u>Cannabinoid</u>	<u>Result</u> (%)	<u>UM</u> (+/-%)	<u>Result</u> (mg/g)	LOQ (mg/g)	<u>Result</u> (mg/ mL)	<u>Per</u> Serving (mg)	<u>Per</u> g <u>Unit</u> (mg)
CBC	-	-	-	4.79	-	-	-
CBCA	DET	-	DET	4.79	-	-	-
CBD	-	-	-	4.79	-	-	-
CBDA	DET	-	DET	4.79	-	-	-
CBDV	DET	-	DET	4.79	-	-	-
CBDVA	-	-	-	4.79	-	-	-
CBG	-	-	-	4.79	-	-	-
CBGA	2.07	0.228	20.7	4.79	-	20.7	20.7
CBL	DET	-	DET	4.79	-	-	-
CBN	DET	-	DET	4.79	-	-	-
CBNA	-	-	-	4.79	-	-	-
Δ9-ΤΗϹ	DET	-	DET	4.79	-	-	-
$\Delta 8$ -THC	DET	-	DET	4.79	-	-	-
THCA	80.4	8.04	804	4.79	-	804	804
THCV	DET	-	DET	4.79	-	-	-
THCVA	0.687	0.0343	6.87	4.79	-	6.87	6.87
TOTAL	83.2	8.31	832		-	832	832
TOTAL CBC	-	-	-		-	-	-
TOTAL CBD	-	-	-		-	-	-
TOTAL CBDV	-	-	-		-	-	-
TOTAL CBG	1.82	0.200	18.2		-	18.2	18.2
TOTAL CBN	-	-	-		-	-	-
TOTAL THC	70.5	7.05	705		-	705	705
TOTAL THCV	0.596	0.0298	5.96		-	5.96	5.96

* 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) + 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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Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
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Greenbrier, AR 72058	Fort Smith, AR 72903	Batch Number:	Date Collected: 07/07/2022
License: ADH 113	License: 00065C	Not Entered	Date Received: 07/08/2022
Cultivar (Strain) or Sample De	Date Completed: 07/10/2022		

Analysis Date/Time:07/09/2022 1657

TERPENOID PROFILE

Analysis Date/Time:07/ Analyst: KF	09/2022 1657		ethod: GC/MS strument: Agilent 7890/5975	Deviations from SOP: None			
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)					
α-Bisabolol	-	-					
Camphene	-	-					
δ-3-Carene	-	-					
β-Caryophyllene	1240	0.124					
Caryophyllene oxide	-	-					
p-Cymene	-	-					
Eucalyptol	-	-					
Geraniol	-	-					
Guaiol	131	0.0131		Abbreviations: GC - Gas			
α-Humulene	88.0	0.00880		Chromatography, MS - Mass			
Isopulegol	-	-		Spectrometry, RL - Reporting Limit			
d-Limonene	2840	0.284		This information is provided as a service and makes no claims of efficacy and/or			
Linalool	247	0.0247		safety of this product.			
β-Myrcene	8340	0.834		Results are applicable only for the sample(s) analyzed and for the specific			
cis-Nerolidol	460	0.0460		analysis conducted.			
trans-Nerolidol	272	0.0272		This report is for informational purposes only and should not be used to diagnose,			
α-Ocimene	-	-		treat, or prevent any			
β-Ocimene	3950	0.395		medical-related symptoms.			
α-Pinene	1120	0.112	1	The statements and results herein have not been approved and/or endorsed by			
β-Pinene	1280	0.128		the FDA.			
α-Terpinene	60.7	0.00607					
γ-Terpinene	-	-					
Terpinolene	11400	1.14		"-" Not detected above RL.			
TOTAL	31400	3.14		Reporting Limit (µg/g): 19.6			













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Cultivar (Strain) or Sample Des	Date Completed: 07/10/2022		

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

Analysis Date/Time: Analyst: KF	7/8/2022 2245 Method: HS/GC/MS Instrument: Agilent 7890/5975			Deviations from SOP: 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> <u>Level</u> (µg/g)	
Acetone (67-64-1)	-	36.4	72.9	5000	n-Heptane (142-82-5)	-	36.4	72.9	5000	
Acetonitrile (75-5-8)	-	36.4	72.9	410	n-Hexane (110-54-3)	-	12.8	25.5	290	
Benzene (71-43-2)	-	0.364	0.729	2	Isobutane (75-28-5)	-	36.4	72.9	5000	
n-Butane (106-97-2)	-	36.4	72.9	5000	Isopropanol (67-63-0)	-	36.4	72.9	5000	
1-Butanol (71-36-3)	-	36.4	72.9	5000	Isopropyl acetate		36.4	72.9	5000	
2-Butanol (78-92-2)	-	36.4	72.9	5000	(108-21-4)		50.4	72.9	3000	
2-Butanone (78-93-3)	-	36.4	72.9	5000	Isopropyl benzene	_	3.64	7.29	70	
Cyclohexane (110-82-7)	-	36.4	72.9	3880	(98-82-8)				2000	Color Key
1,2-Dimethoxyethane		3.64	7.29	100	Methanol (67-56-1)	-	36.4	72.9	3000	<u>eolor ney</u>
(110-71-4)	-	3.64	7.29	100	2-Methylbutane (78-78-4)	-	36.4	72.9	5000	RESULT < AL
N,N-Dimethylacetamide (127-19-5)	-	36.4	72.9	1090	Methylene chloride (75-9-2)	-	36.4	72.9	600	RESULT > AL
2,2-Dimethylbutane (75-83-2)	-	12.8	25.5	290	2-Methylpentane (107-83-5)	-	12.8	25.5	290	"DET" detected less than LOQ
2,3-Dimethylbutane		12.8	25.5	290	3-Methylpentane (96-10-0)	-	12.8	25.5	290	"-" not detected above
(79-29-8)		12.0	23.5	250	n-Pentane (109-66-0)	-	36.4	72.9	5000	LOD
N,N-Dimethylformamide	_	36.4	72.9	880	1-Pentanol (71-41-0)	-	36.4	72.9	5000	"*" - o,m,p-Xylene and
(68-12-2)					n-Propane (74-98-6)	-	36.4	72.9	5000	Ethylbenzene
Dimethylsulfoxide (67-68-5)	-	36.4	72.9	5000	1-Propanol (71-23-8)	-	36.4	72.9	5000	Action levels are
1,4-Dioxane (123-91-1)	_	36.4	72.9	380	Pyridine (110-86-1)	-	12.8	25.5	200	referenced from the State of
Ethanol (64-17-5)	_	36.4	72.9	5000	Tetrahydrofuran (109-99-9)	-	36.4	72.9	720	Arkansas
2-Ethoxyethanol (110-80-5)	_	12.8	25.5	160	Tetramethylene sulfone	_	12.8	25.5	160	MMJ testing
Ethyl ether (60-29-7)	-	36.4	72.9	5000	(126-33-0) Toluene (108-88-3)		36.4	72.9	890	guidelines.
Ethyl acetate (141-78-6)	_	36.4	72.9	5000	· · · ·	-				A value of "-"
Ethyl benzene (100-41-4)	_	36.4	72.9	2170	o-Xylene (95-47-6) m,p-Xylene (108-38-3 or	-	36.4	72.9	2170	for the action level
Ethylene glycol (107-21-1)	-	36.4	72.9	620	106-42-3)	-	36.4	72.9	2170	means that analyte
Ethylene oxide (75-21-8)	_	3.64	7.29	50	Xylenes* (1330-20-7)	_	43.3	86.7	2170	is not currently
.,,					j ()					regulated by the regulations referenced above.
Solvent		Synonym(s	5)		Solvent	S	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	A		
2-Butanone		Methyl ethy	yl ketone, N	/IEK	Isopropyl Acetate	А	cetic acid iso	propyl este	r	
1,2-Dimethoxyethane		Monoglym	e		Methanol	Μ	fethyl alcoho	1		
2,3-Dimethylbutane		Neohexane			2-Methylbutane	Is	opentane			
2,3-Dimethylbutane		Diisopropy	71		Methylene chloride	D	ichlorometha	ine		
N,N-Dimethylformamide		DMF			2-Methylpentane	Is	ohexane			
Dimethysufoxide		DMSO			1-Pentanol		-Amyl alcoho	ol		
2-Ethoxyethanol		Cellosolve,	, Ethyl glvc	ol	1-Propanol		ropyl alcohol			
Ethyl ether		Diethyl eth			Tetrahydrofuran		HF			
Ethyl acetate		EtOAc			Tetramethylene sulfone		ulfolane			
Ethyl benzene		Phenyletha	ne		Xylene		imethylbenz	ene		
		- meny reenu				D				

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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Analysis Date/Time: 7/8/2022 1922 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)	-	0.0196	0.157	0.5	Kresoxim-methyl	_	0.0196	0.157	0.4	
Acephate (30560-19-1)	-	0.0196	0.157	0.4	(143390-89-0)					
Acequinocyl (57960-19-7)	-	0.0196	0.157	2	Malathion (121-75-5)	-	0.0196	0.157	0.2	
Acetamiprid (135410-20-7)	-	0.0196	0.157	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.0196 0.0196	0.157 0.157	0.2 0.2	<u>Color Key</u>
Aldicarb (116-06-3)	-	0.0196	0.157	0.4	Methomyl (16752-77-5)	-	0.0196	0.157	0.4	RESULT < AL
Azoxystrobin (131860-33-8)	-	0.0196	0.157	0.2	Methyl parathion (298-0-0)	-	0.0196	0.157	0.2	RESULT > AL
Bifenazate (149877-41-8)	-	0.0196	0.157	0.2	MGK 264 (113-48-4)	-	0.0196	0.157	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3)	-	0.0196	0.157	0.2	Myclobutanil	_	0.0196	0.157	0.2	"-" not detected above
Boscalid (188425-85-6)	-	0.0196	0.157	0.4	(88671-89-0)					LOD
Carbaryl (63-25-2)	-	0.0196	0.157	0.2	Naled (300-76-5)	-	0.0196	0.157	0.5	Permethrins measured as the
Carbofuran (1563-66-2)	-	0.0196	0.157	0.2	Oxamyl (23135-22-0)	-	0.0196	0.157	1	cumulative residue of the <i>cis</i> - and
Chlorantraniliprole (800008-45-7)	-	0.0196	0.157	0.2	Paclobutrazol (76738-62-0)	-	0.0196	0.157	0.4	trans- permethrin isomers.
Chlorfenapyr	_	0.0196	0.157	1	Permethrins (52645-53-1)	-	0.0196	0.157	0.2	Pyrethrins measured as the cumulative residue of the
(122453-73-0)	-	0.0190	0.137	1	Phosmet (732-11-6)	-	0.0196	0.157	0.2	pyrethrin I, cinerin I, and jasmolin
Chlorpyrifos (2921-88-2)	-	0.0196	0.157	0.2	Piperonyl butoxide	_	0.0196	0.157	2	I isomers.
Clofentezine (74115-24-5)	-	0.0196	0.157	0.2	(51-03-6) Prallethrins (2331-36-9)		0.0196	0.157	0.2	Action levels are referenced from
Cyfluthrin (68359-37-5)	-	0.0196	0.157	1	Propiconazole	-	0.0196	0.157	0.2	the
Cypermethrin (52315-07-8)	-	0.0196	0.157	1	(60207-90-1))	-	0.0196	0.157	0.4	State of Arkansas MMJ testing guidelines.
Daminozide (1596-84-5)	-	0.0196	0.157	1	Propoxur (114-26-1)	-	0.0196	0.157	0.2	A value of "-" for the action level
DDVP (62-73-7)	-	0.0196	0.157	0.1	Pyrethrins (8003-34-7)	-	0.0196	0.157	1	means that analyte is not
Diazinon (333-41-5)	-	0.0196	0.157	0.2	Pyridaben (96489-71-3)	-	0.0196	0.157	0.2	currently regulated by the regulations referenced above.
Dimethoate (60-51-5)	-	0.0196	0.157	0.2	Spinosad (168316-95-8)	-	0.0196	0.157	0.2	
Ethoprophos (13194-48-4)	-	0.0196	0.157	0.2	Spiromesifen (283594-90-1)	-	0.0196	0.157	0.2	Disclaimer: This information is
Etofenprox (80844-07-1)	-	0.0196	0.157	0.4	Spirotetramat					provided as a service and makes no claims of efficacy and/or safety
Etoxazole (153233-91-1)	-	0.0196	0.157	0.2	(203313-25-1)	-	0.0196	0.157	0.2	of this product. Results are
Fenoxycarb (72490-01-8)	-	0.0196	0.157	0.2	Spiroxamine	_	0.0196	0.157	0.4	applicable only for the sample(s) analyzed and for the specific
(E)-Fenpyroximate (134098-61-6)	-	0.0196	0.157	0.4	(118134-30-8) Tebuconazole		0.0196	0.157	0.4	analysis conducted. This report is for informational purposes only
Fipronil (120068-37-3)	-	0.0196	0.157	0.4	(80443-41-0)	-	0.0196	0.157	0.4	and should not be used to
Flonicamid (158062-67-0)	-	0.0196	0.157	1	Thiacloprid	_	0.0196	0.157	0.2	diagnose, treat, or prevent any medical-related symptoms. The
Fludioxinil (131341-86-1)	-	0.0196	0.157	0.4	(111988-49-9)					statements and results herein have
Hexythiazox (78587-05-0)	-	0.0196	0.157	1	Thiamethoxam (153719-23-4)	-	0.0196	0.157	0.2	not been approved and/or
Imazalil (35554-44-0)	-	0.0196	0.157	0.2	Trifloxystrobin		0.0100	0.155	0.5	endorsed by the FDA.
Imidacloprid (138261-41-3)	-	0.0196	0.157	0.4	(141517-21-7)	-	0.0196	0.157	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	<u>Synonym(s)</u>
Cyfluthrin	Baythroid	Myclobutanil	Systhane	Propiconazole	Tilt
DDVP	Dichlorvos	Naled	Dibrom	Propoxur	Baygon
Ethoprophos	Prophos	Phosmet	Imidan		



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Cultivar (Strain) or Sample Des	Date Completed: 07/10/2022		

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

Analysis Date/Time: 07/09/2022 0956 (ICP/OES) Analysis Date/Time: - (DMA) Analyst: KF		Method: ICP/OES Instrument: Agilent 720-ES		Deviations from SOP:ESNone	
<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)	-	45.7	86.9	200	
Cadmium (Cd)	-	45.7	86.9	200	
Lead (Pb)	-	45.7	86.9	500	
Mercury (Hg)	-	45.7	86.9	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: SA31889)

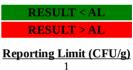
Testing Location:	Customer ID: 2168	Sample ID: SA31889	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13241641866	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Batch Number:	Date Collected: 07/07/2022
License: ADA 05_H273	License: 00065C	Not Entered	Date Received: 07/08/2022
Cultivar (Strain) or Sample	Date Completed: 07/10/2022		

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 7/9/2022 Analyst: PW		ardy Diagnostics CompactDry t: Thermo Incubator	Deviations from SOP: None
Bacteria/Microbe	<u>Result</u> (CFU/g)	<u>Action Level</u> (CFU/g)	
Aerobic Plate Count	NT	-	
Coliforms, Total	Absent	1	
Escherichia Coli (E. Coli)	Absent	100	
Mold/Yeast	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





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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