







SUMMARY OF ANALYSIS (SAMPLE ID: SA31757)

Testing Location:	Customer ID: 2168	Order ID: OR9887	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13245542023	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Batch Number:	Date Collected: 06/22/2022
License: ADH 113	License: 00065C	M00065C13246232520	Date Received: 06/22/2022
Cultivar (Strain) or Sample D	Date Completed: 06/24/2022		

*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.

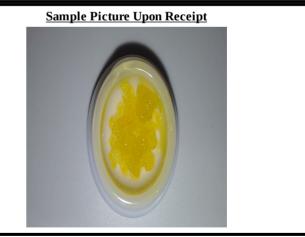
Moisture Content (%)	Water Activity (aw)	PASS/FAIL
Not Tested	Not Tested	N/A

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	88.6	886
THCVA	1.02	10.2
CBGA	0.693	6.93
TOTAL CBD	-	
TOTAL THC	77.7	777
TOTAL CANNABINOIDS	90.3	903
<u>Terpenes (Top 5)</u>	<u>(%)</u>	<u>µg∕g</u>
<u>Terpenes (Top 5)</u> β-Myrcene	<u>(%)</u> 3.14	µg/g 31400
β-Myrcene	3.14	31400
β-Myrcene d-Limonene	3.14 1.65	31400 16500
β-Myrcene d-Limonene β-Caryophyllene	3.14 1.65 0.934	31400 16500 9340

PASS/FAIL
PASS
PASS
PASS
PASS





Scan the QR code to verify results.

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cbd tincture non-mandatory Gummies mandatory concentrates mandatory



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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA31757)

Testing Location:	Customer ID: 2168	Order ID: OR9887	Sample Type: Primary
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232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13245542023	Mass: 4g
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Cultivar (Strain) or Sample	Date Completed: 06/24/2022		

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

Analysis Date/Time: 06/24/2022 0142 Analyst: PW Method: HPLC/DAD Instrument: Agilent 1100

<u>Cannabinoid</u>	<u>Result</u> (%)	<u>UM</u> (+/-%)	<u>Result</u> (mg/g)	<u>LOQ</u> (mg/g)	<u>Result</u> (mg/ mL)	<u>Per</u> <u>Serving</u> (mg)	<u>Per</u> g <u>Unit</u> (mg)
CBC	-	-	-	6.74	-	-	-
CBCA	-	-	-	6.74	-	-	-
CBD	DET	-	DET	6.74	-	-	-
CBDA	DET	-	DET	6.74	-	-	-
CBDV	-	-	-	6.74	-	-	-
CBDVA	DET	-	DET	6.74	-	-	-
CBG	-	-	-	6.74	-	-	-
CBGA	0.693	0.0763	6.93	6.74	-	6.93	6.93
CBL	DET	-	DET	6.74	-	-	-
CBN	-	-	-	6.74	-	-	-
CBNA	-	-	-	6.74	-	-	-
Δ9-ΤΗC	-	-	-	6.74	-	-	-
Δ8- THC	DET	-	DET	6.74	-	-	-
THCA	88.6	8.86	886	6.74	-	886	886
THCV	DET	-	DET	6.74	-	-	-
THCVA	1.02	0.0509	10.2	6.74	-	10.2	10.2
TOTAL	90.3	8.99	903		-	903	903
TOTAL CBC	-	-	-		-	-	-
TOTAL CBD	-	-	-		-	-	-
TOTAL CBDV	-	-	-		-	-	-
TOTAL CBG	0.609	0.0670	6.09		-	6.09	6.09
TOTAL CBN	-	-	-		-	-	-
TOTAL THC	77.7	7.77	777		-	777	777
TOTAL THCV	0.882	0.0441	8.82		-	8.82	8.82

* 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. tory Dire



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Greenbrier, AR 72058	Fort Smith, AR 72903	Batch Number:	Date Collected: 06/22/2022
License: ADH 113	License: 00065C	M00065C13246232520	Date Received: 06/22/2022
Cultivar (Strain) or Sample D	Date Completed: 06/24/2022		

Analysis Date/Time:06/24/2022 1332 Analyst: KF

Result

 $(\mu g/g)$ 480

519

_

9340

59.5

3680

50.4

16500

1260

31400

904

49.6

98.6

_

1510

2750

_

_

0.151

0.275

Terpene

α-Bisabolol Camphene

δ-3-Carene

β-Caryophyllene

Caryophyllene

α-Humulene

Isopulegol

Linalool

β-Myrcene

cis-Nerolidol

α-Ocimene

β-Ocimene

α-Pinene

β-Pinene

α-Terpinene y-Terpinene

trans-Nerolidol

d-Limonene

oxide p-Cymene Eucalyptol Geraniol Guaiol

TERPENOID PROFILE

Method: GC/MS Instrument: Agilent 7890/5975

Deviations from SOP: None

<u>Result</u> (%) 0.0480 0.0519 0.934 0.00595 Abbreviations: GC - Gas 0.368 Chromatography, MS - Mass Spectrometry, RL - Reporting Limit 0.00504 This information is provided as a service 1.65 and makes no claims of efficacy and/or safety of this product. 0.126 Results are applicable only for the 3.14 sample(s) analyzed and for the specific 0.0904 analysis conducted. This report is for informational purposes 0.00496 only and should not be used to diagnose, 0.00986 treat, or prevent any

medical-related symptoms.

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Terpinolene	283	0.0283	"-" Not detected above RL.
TOTAL	68800	6.88	Reporting Limit (μg/g): 19.5













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Cultivar (Strain) or Sample De	Date Completed: 06/24/2022		

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

Analysis Date/Time: 6/23/2022 2021 Analyst: KF			Method: HS/GC/MS Instrument: Agilent 7890/5975				Deviations from SOP: None			
Solvent	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	Action Level (µg/g)	Solvent	<u>Result</u> (µg/g)		LOQ (µg/g)	<u>Action</u> Level (µg/g)	
Acetone (67-64-1)	_	36.0	72.0	5000	n-Heptane (142-82-5)	-	36.0	72.0	5000	
Acetonitrile (75-5-8)	-	36.0	72.0	410	n-Hexane (110-54-3)	_	12.6	25.2	290	
Benzene (71-43-2)	_	0.360	0.720	2	Isobutane (75-28-5)		36.0	72.0	5000	
n-Butane (106-97-2)	319	36.0	72.0	5000	Isopropanol (67-63-0)	_	36.0	72.0	5000	
1-Butanol (71-36-3)	-	36.0	72.0	5000	Isopropyl acetate					
2-Butanol (78-92-2)	-	36.0	72.0	5000	(108-21-4)	-	36.0	72.0	5000	
2-Butanone (78-93-3)	_	36.0	72.0	5000	Isopropyl benzene	_	3.60	7.20	70	
Cyclohexane (110-82-7)	-	36.0	72.0	3880	(98-82-8)					Color Key
1,2-Dimethoxyethane					Methanol (67-56-1)	-	36.0	72.0	3000	<u>Color Rey</u>
(110-71-4)	-	3.60	7.20	100	2-Methylbutane (78-78-4)	-	36.0	72.0	5000	RESULT < AL
N,N-Dimethylacetamide (127-19-5)	-	36.0	72.0	1090	Methylene chloride (75-9-2)	-	36.0	72.0	600	RESULT > AL
2,2-Dimethylbutane (75-83-2)	-	12.6	25.2	290	2-Methylpentane (107-83-5)	-	12.6	25.2	290	"DET" detected less than LOQ
2,3-Dimethylbutane		12.6	25.2	290	3-Methylpentane (96-10-0)	-	12.6	25.2	290	"-" not detected above
(79-29-8)	-	12.6	25.2	290	n-Pentane (109-66-0)	-	36.0	72.0	5000	LOD
N,N-Dimethylformamide	_	36.0	72.0	880	1-Pentanol (71-41-0)	-	36.0	72.0	5000	"*" - o,m,p-Xylene and
(68-12-2)					n-Propane (74-98-6)	-	36.0	72.0	5000	Ethylbenzene
Dimethylsulfoxide (67-68-5)	-	36.0	72.0	5000	1-Propanol (71-23-8)	-	36.0	72.0	5000	
1,4-Dioxane (123-91-1)	_	36.0	72.0	380	Pyridine (110-86-1)	-	12.6	25.2	200	Action levels are referenced from the State of
Ethanol (64-17-5)		36.0	72.0	5000	Tetrahydrofuran (109-99-9)	-	36.0	72.0	720	Arkansas
2-Ethoxyethanol (110-80-5)	_	12.6	25.2	160	Tetramethylene sulfone	_	12.6	25.2	160	MMJ testing
Ethyl ether (60-29-7)	_	36.0	72.0	5000	(126-33-0)					guidelines.
Ethyl acetate (141-78-6)	_	36.0	72.0	5000	Toluene (108-88-3)	-	36.0	72.0	890	A value of "-"
Ethyl benzene (100-41-4)	_	36.0	72.0	2170	o-Xylene (95-47-6)	-	36.0	72.0	2170	for the action level
Ethylene glycol (107-21-1)	_	36.0	72.0	620	m,p-Xylene (108-38-3 or 106-42-3)	-	36.0	72.0	2170	means that analyte
Ethylene oxide (75-21-8)	_	3.60	7.20	50	Xylenes* (1330-20-7)	_	43.3	86.7	2170	is not currently
,		-		50			_	000	21/0	regulated by the regulations referenced above.
Solvent		Synonym(s			<u>Solvent</u>		Synonym(s)			
Acetonitrile		Methyl Cya			Ethylene glycol		1,2-Ethanediol			
1-Butanol		n-Butanol,	5	hol	Isobutane		2-Methylpropane			
2-Butanol		sec-Butyl a		m v	Isopropanol		2-Propanol, IP/			
2-Butanone		Methyl ethy	-	1EK	Isopropyl Acetate		Acetic acid iso		r	
1,2-Dimethoxyethane		Monoglym			Methanol		Methyl alcohol -			
2,3-Dimethylbutane		Neohexane			2-Methylbutane		Isopentane			
2,3-Dimethylbutane		Diisopropy	71		Methylene chloride		Dichlorometha	ne		
N,N-Dimethylformamide		DMF			2-Methylpentane		Isohexane	,		
Dimethysufoxide		DMSO		,	1-Pentanol		n-Amyl alcoho	1		
2-Ethoxyethanol			, Ethyl glyc	101	1-Propanol		Propyl alcohol			
Ethyl ether		Diethyl eth	er, Ether		Tetrahydrofuran		THF			
Ethyl acetate		EtOAc			Tetramethylene sulfone		Sulfolane			
Ethyl benzene		Phenyletha	ne		Xylene	1	Dimethylbenze	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:

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Cultivar (Strain) or Sample	Date Completed: 06/24/2022		

Analysis Date/Time: 06/24/2022 0658

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Method: LC/MS/MS

Analyst: KF					Instrument: Shimadzu LC-8050			None		
<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> <u>Level</u> (µg/g)	
Abamectin (71751-41-2)	-	0.0195	0.156	0.5	Kresoxim-methyl	_	0.0195	0.156	0.4	
Acephate (30560-19-1)	-	0.0195	0.156	0.4	(143390-89-0)					
Acequinocyl (57960-19-7)	-	0.0195	0.156	2	Malathion (121-75-5)	-	0.0195	0.156	0.2	
Acetamiprid (135410-20-7)	-	0.0195	0.156	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.0195 0.0195	0.156 0.156	0.2 0.2	<u>Color Key</u>
Aldicarb (116-06-3)	-	0.0195	0.156	0.4	Methomyl (16752-77-5)	-	0.0195	0.156	0.4	RESULT < AL
Azoxystrobin (131860-33-8)	-	0.0195	0.156	0.2	Methyl parathion (298-0-0)	-	0.0195	0.156	0.2	RESULT > AL
Bifenazate (149877-41-8)	-	0.0195	0.156	0.2	MGK 264 (113-48-4)	-	0.0195	0.156	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3) Boscalid (188425-85-6)	-	0.0195 0.0195	0.156 0.156	0.2 0.4	Myclobutanil (88671-89-0)	-	0.0195	0.156	0.2	"-" not detected above LOD
Carbaryl (63-25-2)	_	0.0195	0.156	0.2	Naled (300-76-5)	-	0.0195	0.156	0.5	
Carbofuran (1563-66-2)	_	0.0195	0.156	0.2	Oxamyl (23135-22-0)	-	0.0195	0.156	1	Permethrins measured as the cumulative residue of the <i>cis</i> - and
Chlorantraniliprole (800008-45-7)	-	0.0195	0.156	0.2	Paclobutrazol (76738-62-0)	-	0.0195	0.156	0.4	<i>trans</i> - permethrin isomers.
Chlorfenapyr		0.0105	0.150		Permethrins (52645-53-1)	-	0.0195	0.156	0.2	Pyrethrins measured as the
(122453-73-0)	-	0.0195	0.156	1	Phosmet (732-11-6)	-	0.0195	0.156	0.2	cumulative residue of the pyrethrin I, cinerin I, and jasmolin
Chlorpyrifos (2921-88-2)	-	0.0195	0.156	0.2	Piperonyl butoxide	_	0.0195	0.156	2	I isomers.
Clofentezine (74115-24-5)	-	0.0195	0.156	0.2	(51-03-6) Prallethrins (2331-36-9)		0.0195	0.156	0.2	Action levels are referenced from
Cyfluthrin (68359-37-5)	-	0.0195	0.156	1	Propiconazole	-				the
Cypermethrin (52315-07-8)	-	0.0195	0.156	1	(60207-90-1))	-	0.0195	0.156	0.4	State of Arkansas MMJ testing guidelines.
Daminozide (1596-84-5)	-	0.0195	0.156	1	Propoxur (114-26-1)	-	0.0195	0.156	0.2	A value of "-" for the action level
DDVP (62-73-7)	-	0.0195	0.156	0.1	Pyrethrins (8003-34-7)	-	0.0195	0.156	1	means that analyte is not
Diazinon (333-41-5)	-	0.0195	0.156	0.2	Pyridaben (96489-71-3)	-	0.0195	0.156	0.2	currently regulated by the regulations referenced above.
Dimethoate (60-51-5)	-	0.0195	0.156	0.2	Spinosad (168316-95-8)	-	0.0195	0.156	0.2	5
Ethoprophos (13194-48-4)	-	0.0195	0.156	0.2	Spiromesifen (283594-90-1)	-	0.0195	0.156	0.2	Disclaimer: This information is provided as a service and makes
Etofenprox (80844-07-1)	-	0.0195	0.156	0.4	Spirotetramat		0.0105	0.450	0.0	no claims of efficacy and/or safety
Etoxazole (153233-91-1)	-	0.0195	0.156	0.2	(203313-25-1)	-	0.0195	0.156	0.2	of this product. Results are
Fenoxycarb (72490-01-8)	-	0.0195	0.156	0.2	Spiroxamine	_	0.0195	0.156	0.4	applicable only for the sample(s) analyzed and for the specific
(E)-Fenpyroximate (134098-61-6)	-	0.0195	0.156	0.4	(118134-30-8) Tebuconazole	_	0.0195	0.156	0.4	analysis conducted. This report is for informational purposes only
Fipronil (120068-37-3)	-	0.0195	0.156	0.4	(80443-41-0)		010100	0.1200		and should not be used to
Flonicamid (158062-67-0)	-	0.0195	0.156	1	Thiacloprid (111988-49-9)	-	0.0195	0.156	0.2	diagnose, treat, or prevent any medical-related symptoms. The
Fludioxinil (131341-86-1)	-	0.0195	0.156	0.4	Thiamethoxam					statements and results herein have
Hexythiazox (78587-05-0)	-	0.0195	0.156	1	(153719-23-4)	-	0.0195	0.156	0.2	not been approved and/or endorsed by the FDA.
Imazalil (35554-44-0)	-	0.0195	0.156	0.2	Trifloxystrobin	_	0.0195	0.156	0.2	chuoiscu by the i DA.
Imidacloprid (138261-41-3)	-	0.0195	0.156	0.4	(141517-21-7)					

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 Myclobutanil Baythroid Systhane Propiconazole Tilt DDVP Dichlorvos Naled Dibrom Propoxur Baygon Prophos Phosmet Imidan Ethoprophos



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

Analysis Date/Time: 06/24/2022 1010 (ICP/OES) Analysis Date/Time: - (DMA) Analyst: KF		Method: ICP/OES Instrument: Agilent 720-ES		Deviations from SOP:.SNone	
<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)	-	49.2	93.6	200	
Cadmium (Cd)	-	49.2	93.6	200	
Lead (Pb)	-	49.2	93.6	500	
Mercury (Hg)	-	49.2	93.6	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: SA31757)

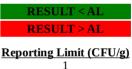
Testing Location:	Customer ID: 2168	Sample ID: SA31757	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13245542023	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Batch Number:	Date Collected: 06/22/2022
License: ADA 05_H273	License: 00065C	M00065C13246232520	Date Received: 06/22/2022
Cultivar (Strain) or Sample	Date Completed: 06/24/2022		

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 06/23/2022 1410 Analyst: PW	Method: Hardy Diagnostics CompactDry Instrument: Thermo Incubator		Deviations from SOP None
<u>Bacteria/Microbe</u>	<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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