



SUMMARY OF ANALYSIS (SAMPLE ID: SA35453)

Testing Location:	Customer ID: 2168	Order ID: OR10497	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13244975869	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 06/19/2023
License: ADH 113	License: 00065C	E20230612RFLHR02	Date Received: 06/20/2023
Cultivar (Strain) or Sample De	Date Completed: 06/21/2023		

*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.

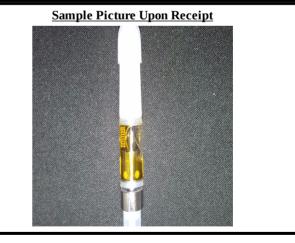
Moisture Content (%)	Water Activity (aw)	PASS/FAIL
Not Tested	Not Tested	PASS

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</u> /g
THCA	39.6	396
Δ9-ΤΗC	26.1	261
CBGA	3.78	37.8
TOTAL CBD	0.172	1.72
TOTAL THC	60.8	608
TOTAL CANNABINOIDS	73.5	735
<u>Terpenes (Top 5)</u>	<u>(%)</u>	µg∕g
<u>Terpenes (Top 5)</u> d-Limonene	<u>(%)</u> 1.46	<mark>µg/g</mark> 14600
d-Limonene	1.46	14600
d-Limonene β-Myrcene	1.46 1.02	14600 10200
d-Limonene β-Myrcene α-Pinene	1.46 1.02 0.142	14600 10200 1420

Contaminants	PASS/FAIL
Heavy Metals:	PASS
Microbiology:	PASS
Pesticides:	PASS
Residual Solvents:	PASS





Scan the QR code to verify results.

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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: SA35453)

Testing Location:	Customer ID: 2168	Order ID: OR10497	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13244975869	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 06/19/2023
License: ADH 113	License: 00065C	E20230612RFLHR02	Date Received: 06/20/2023
Cultivar (Strain) or Sample 1	Date Completed: 06/21/2023		

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

Analysis Date/Time: 06/21/2023 1043 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 Per (mg/ Servin mL) (mg)</u>	<u>Per</u> g <u>Unit</u> (<u>mg)</u>
CBC	0.107	1.07	0.106	0.248	- 1.07	1.07
CBCA	1.35	13.5	0.329	0.769	- 13.5	13.5
CBD	ND	ND	0.748	1.75		-
CBDA	0.196	1.96	0.275	0.643	- 1.96	1.96
CBDV	ND	ND	0.120	0.280		-
CBDVA	ND	ND	0.320	0.747		-
CBG	1.21	12.1	0.486	1.13	- 12.1	12.1
CBGA	3.78	37.8	0.689	0.812	- 37.8	37.8
CBL	1.00	10.0	0.561	1.31	- 10.0	10.0
CBN	ND	ND	0.258	0.602		-
CBNA	ND	ND	0.278	0.649		-
Δ9-ΤΗC	26.1	261	0.309	0.720	- 261	261
$\Delta 8$ -THC	ND	ND	0.482	1.13		-
THCA	39.6	396	0.168	0.392	- 396	396
THCV	ND	ND	0.402	0.937		-
THCVA	0.243	2.43	0.128	0.299	- 2.43	2.43
TOTAL	73.5	735			- 735	735
TOTAL CBC	1.29	12.9			- 12.9	12.9
TOTAL CBD	0.172	1.72			- 1.72	1.72
TOTAL CBDV	-	-				-
TOTAL CBG	4.52	45.2			- 45.2	45.2
TOTAL CBN	-	-				-
TOTAL THC	60.8	608			- 608	608
TOTAL THCV	0.210	2.10			- 2.10	2.10

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 = $(CBCA \times 0.877) + CBC$ Total CBD = (CBDA x 0.877) + CBD Total CBDV = (CBDVA x 0.867) + CBDV Total CBG = $(CBGA \times 0.878) + CBG$ Total CBN = (CBNA x 0.876) + CBN Total THC = (THCA x 0.877) + Δ 9-THC Total THCV = $(THCVA \times 0.867) + THCV$

Percentage results are reported by mass.

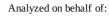
mg/g results are reported as mass component per mass material.

W. Felling, Ph.D. ory Dire









Deviations from SOP:



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35453)

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Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 06/19/2023
License: ADH 113	License: 00065C	E20230612RFLHR02	Date Received: 06/20/2023
Cultivar (Strain) or Sample	Date Completed: 06/21/2023		

Analysis Date/Time:06/21/2023 0234 Analyst: KF

TERPENOID PROFILE
Method: GC/MS

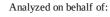
Analyst: KF		Inst	rument: Agilent 7890/5975	None
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)		
α-Bisabolol	43.3	0.00433		
Camphene	1120	0.112		
δ-3-Carene	16.3	0.00163		
β-Caryophyllene	967	0.0967		
Caryophyllene oxide	-	-		
p-Cymene	-	-		
Eucalyptol	-	-		
Geraniol	-	-		
Guaiol	-	-		Abbreviations: GC - Gas
α-Humulene	627	0.0627		Chromatography, MS - Mass
Isopulegol	-	-		Spectrometry, RL - Reporting Limit
d-Limonene	14600	1.46		This information is provided as a service and makes no claims of efficacy and/or
Linalool	407	0.0407		safety of this product.
β-Myrcene	10200	1.02		Results are applicable only for the sample(s) analyzed and for the specific
cis-Nerolidol	-	-		analysis conducted.
trans-Nerolidol	37.6	0.00376		This report is for informational purposes only and should not be used to diagnose,
α-Ocimene	1120	0.112		treat, or prevent any
β-Ocimene	655	0.0655		medical-related symptoms.
α-Pinene	1420	0.142		The statements and results herein have not been approved and/or endorsed by
β-Pinene	41.8	0.00418		the FDA.
α-Terpinene	251	0.0251		
γ-Terpinene	117	0.0117		
Terpinolene	1160	0.116		"-" Not detected above RL.
TOTAL	32800	3.28		Reporting Limit (µg/g): 355

Kyle W. Felling, Ph.D. Liboratory Director











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Cultivar (Strain) or Sample	Date Completed: 06/21/2023		

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

Analysis Date/Time:	06/20/20)23 1854		I	Method: HS/GC/MS			D	eviation	s from SOP:
Analyst: KF				1	nstrument: Agilent 78	90/5975		N	lone	
<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)	
Acetone (67-64-1)	-	151	302	5000	n-Heptane (142-82-5)	-	151	302	5000	
Acetonitrile (75-5-8)	-	151	302	410	n-Hexane (110-54-3)	-	52.9	106	290	Contraction of the second s
Benzene (71-43-2)	-	1.51	3.02	2	Isobutane (75-28-5)	-	151	302	5000	21
n-Butane (106-97-2)	3510	151	302	5000	Isopropanol (67-63-0)	-	151	302	5000	r 2
1-Butanol (71-36-3)	-	151	302	5000	Isopropyl acetate	_	151	302	5000	60
2-Butanol (78-92-2)	-	151	302	5000	(108-21-4)		151	502	5000	
2-Butanone (78-93-3)	-	151	302	5000	Isopropyl benzene	-	15.1	30.2	70	
Cyclohexane (110-82-7)	-	151	302	3880	(98-82-8)		151	302	3000	Color Key
1,2-Dimethoxyethane		15.1	30.2	100	Methanol (67-56-1)	-	151			
(110-71-4)		15.1	30.2	100	2-Methylbutane (78-78-4)	-	151	302	5000	RESULT < AL
N,N-Dimethylacetamide (127-19-5)	-	151	302	1090	Methylene chloride (75-9-2)	-	151	302	600	RESULT > AL
2,2-Dimethylbutane (75-83-2)	-	52.9	106	290	2-Methylpentane (107-83-5)	-	52.9	106	290	"DET" detected less than LOQ
2,3-Dimethylbutane	_	52.9	106	290	3-Methylpentane (96-10-0)	-	52.9	106	290	"-" not detected above LOD
(79-29-8)		52.5	100	250	n-Pentane (109-66-0)	-	151	302	5000	LOD
N,N-Dimethylformamide (68-12-2)	_	151	302	880	1-Pentanol (71-41-0)	-	151	302	5000	"*" - o,m,p-Xylene and
(66-12-2) Dimethylsulfoxide					n-Propane (74-98-6)	-	151	302	5000	Ethylbenzene
(67-68-5)	-	151	302	5000	1-Propanol (71-23-8)	-	151	302	5000	Action levels are
1,4-Dioxane (123-91-1)	_	151	302	380	Pyridine (110-86-1)	-	52.9	106	200	referenced from the State of
Ethanol (64-17-5)	418	151	302	5000	Tetrahydrofuran (109-99-9)	-	151	302	720	Arkansas
2-Ethoxyethanol (110-80-5)	_	52.9	106	160	Tetramethylene sulfone (126-33-0)	-	52.9	106	160	MMJ testing
Ethyl ether (60-29-7)	-	151	302	5000	(120-33-0) Toluene (108-88-3)		151	302	890	guidelines.
Ethyl acetate (141-78-6)	-	151	302	5000	o-Xylene (95-47-6)	-	151	302	2170	A value of "-"
Ethyl benzene (100-41-4)	-	151	302	2170	m,p-Xylene (108-38-3 or	-	151	302	2170	for the action level
Ethylene glycol (107-21-1)	-	151	302	620	106-42-3)	-	151	302	2170	means that analyte
Ethylene oxide (75-21-8)	-	15.1	30.2	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170	is not currently regulated by the
Solvent		Synonym(s	5)		Solvent	s	ynonym(s)			regulations referenced above.
Acetonitrile		Methyl Cya			Ethylene glycol		,2-Ethanedio			
1-Butanol		n-Butanol,		hol	Isobutane		-Methylpropa			
2-Butanol		sec-Butyl a		1101	Isopropanol		-Propanol, IP			
2-Butanone		Methyl eth		/FK	Isopropyl Acetate		cetic acid iso		r	
1,2-Dimethoxyethane		Monoglym	-		Methanol		fethyl alcoho		L	
2,3-Dimethylbutane		Neohexane			2-Methylbutane		opentane	1		
2,3-Dimethylbutane		Diisopropy			Methylene chloride		ichlorometha	000		
N,N-Dimethylformamide		DIISOPIOPS			2-Methylpentane		ohexane			
Dimethysufoxide		DMF			1-Pentanol		-Amyl alcoho			
2-Ethoxyethanol		Cellosolve	Ethyl glyc	ol	1-Propanol		ropyl alcohol			
Ethyl ether		Diethyl eth		.01	Tetrahydrofuran		HF			
		EtOAc	iei, Euller				нг ulfolane			
Ethyl acetate					Tetramethylene sulfone					
Ethyl benzene		Phenyletha	пе		Xylene	D	imethylbenz	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



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Cultivar (Strain) or Sample	Date Completed: 06/21/2023		

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 06/20/2023 1518 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)	-	0.0450	0.360	0.5	Kresoxim-methyl	_	0.0450	0.360	0.4	T
Acephate (30560-19-1)	-	0.0450	0.360	0.4	(143390-89-0)					
Acequinocyl (57960-19-7)	-	0.0450	0.360	2	Malathion (121-75-5)	-	0.0450	0.360	0.2	
Acetamiprid	_	0.0450	0.360	0.2	Metalaxyl (57837-19-1)	-	0.0450	0.360	0.2	Color Key
(135410-20-7)		0.0450	0.260		Methiocarb (2032-65-7)	-	0.0450	0.360	0.2	
Aldicarb (116-06-3)	-	0.0450	0.360	0.4	Methomyl (16752-77-5)	-	0.0450	0.360	0.4	RESULT < AL
Azoxystrobin (131860-33-8)	-	0.0450	0.360	0.2	Methyl parathion (298-0-0)	-	0.0450	0.360	0.2	RESULT > AL
Bifenazate (149877-41-8)	-	0.0450	0.360	0.2	MGK 264 (113-48-4)	-	0.0450	0.360	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3)	-	0.0450	0.360	0.2	Myclobutanil	_	0.0450	0.360	0.2	"-" not detected above
Boscalid (188425-85-6)	-	0.0450	0.360	0.4	(88671-89-0)		0.0450	0.200	0.5	LOD
Carbaryl (63-25-2)	-	0.0450	0.360	0.2	Naled (300-76-5)	-	0.0450	0.360	0.5	Permethrins measured as the
Carbofuran (1563-66-2)	-	0.0450	0.360	0.2	Oxamyl (23135-22-0) Paclobutrazol	-	0.0450	0.360	1	cumulative residue of the cis- and
Chlorantraniliprole (800008-45-7)	-	0.0450	0.360	0.2	(76738-62-0)	-	0.0450	0.360	0.4	trans- permethrin isomers.
Chlorfenapyr		0.0450	0.360	1	Permethrins (52645-53-1)	-	0.0450	0.360	0.2	Pyrethrins measured as the cumulative residue of the
(122453-73-0)					Phosmet (732-11-6)	-	0.0450	0.360	0.2	pyrethrin I, cinerin I, and jasmolin
Chlorpyrifos (2921-88-2)	-	0.0450	0.360	0.2	Piperonyl butoxide (51-03-6)	-	0.0450	0.360	2	I isomers.
Clofentezine (74115-24-5)	-	0.0450	0.360	0.2	(51-03-0) Prallethrins (2331-36-9)	_	0.0450	0.360	0.2	Action levels are referenced from
Cyfluthrin (68359-37-5)	-	0.0450	0.360	1	Propiconazole	-				the
Cypermethrin (52315-07-8)	-	0.0450	0.360	1	(60207-90-1))	-	0.0450	0.360	0.4	State of Arkansas MMJ testing guidelines.
Daminozide (1596-84-5)	-	0.0450	0.360	1	Propoxur (114-26-1)	-	0.0450	0.360	0.2	A value of "-" for the action level
DDVP (62-73-7)	-	0.0450	0.360	0.1	Pyrethrins (8003-34-7)	-	0.0450	0.360	1	means that analyte is not
Diazinon (333-41-5)	-	0.0450	0.360	0.2	Pyridaben (96489-71-3)	-	0.0450	0.360	0.2	currently regulated by the regulations referenced above.
Dimethoate (60-51-5)	-	0.0450	0.360	0.2	Spinosad (168316-95-8)	-	0.0450	0.360	0.2	
Ethoprophos (13194-48-4)	-	0.0450	0.360	0.2	Spiromesifen (283594-90-1)	-	0.0450	0.360	0.2	Disclaimer: This information is
Etofenprox (80844-07-1)	-	0.0450	0.360	0.4	Spirotetramat					provided as a service and makes no claims of efficacy and/or safety
Etoxazole (153233-91-1)	-	0.0450	0.360	0.2	(203313-25-1)	-	0.0450	0.360	0.2	of this product. Results are
Fenoxycarb (72490-01-8)	-	0.0450	0.360	0.2	Spiroxamine	_	0.0450	0.360	0.4	applicable only for the sample(s) analyzed and for the specific
(E)-Fenpyroximate (134098-61-6)	-	0.0450	0.360	0.4	(118134-30-8) Tebuconazole					analysis conducted. This report is for informational purposes only
Fipronil (120068-37-3)	-	0.0450	0.360	0.4	(80443-41-0)	-	0.0450	0.360	0.4	and should not be used to
Flonicamid (158062-67-0)	-	0.0450	0.360	1	Thiacloprid	-	0.0450	0.360	0.2	diagnose, treat, or prevent any medical-related symptoms. The
Fludioxinil (131341-86-1)	-	0.0450	0.360	0.4	(111988-49-9)					statements and results herein have
Hexythiazox (78587-05-0)	-	0.0450	0.360	1	Thiamethoxam (153719-23-4)	-	0.0450	0.360	0.2	not been approved and/or
Imazalil (35554-44-0)	-	0.0450	0.360	0.2	Trifloxystrobin		0.0450	0.000	0.0	endorsed by the FDA.
Imidacloprid (138261-41-3)	-	0.0450	0.360	0.4	(141517-21-7)	-	0.0450	0.360	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	Synonym(s)	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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Analyzed on behalf of:



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License: ADH 113	License: 00065C	E20230612RFLHR02	Date Received: 06/20/2023
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HEAVY METAL PROFILE (SOP: SOP-ICP-200.6)

Analysis Date/Time: 06/20/2023 1716 (ICP/OES) Analysis Date/Time: - (DMA) Analyst: KF				t hod: ICP/MS rument: Agilent 750	Deviations from SOP: Oce 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)	-	58.6	92.8	200	
Cadmium (Cd)	-	58.6	92.8	200	
Lead (Pb)	-	58.6	92.8	500	
Mercury (Hg)	-	58.6	92.8	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.

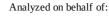
Disclaimer: This information is provided as a service and makes no claims of efficacy and/or safety of this product. Results are applicable only for the sample(s) analyzed and for the specific analysis conducted. This report is for informational purposes only and should not be used to diagnose, treat, 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: SA35453)

Testing Location:	Customer ID: 2168	Sample ID: SA35453	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13244975869	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 06/19/2023
License: ADA 05_H273	License: 00065C	E20230612RFLHR02	Date Received: 06/20/2023
Cultivar (Strain) or Sample D	Date Completed: 06/21/2023		

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 06/21/20 Analyst: PW		Hardy Diagnostics Compac nt: Thermo Incubator	ctDry 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	-	2/1
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





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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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