



SUMMARY OF ANALYSIS (SAMPLE ID: SA37616)

Testing Location:	Customer ID: 2168	Order ID: OR10953	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13245972560	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 04/11/2024
License: ADH 113	License: 00065C	E20240410BNZLB03	Date Received: 04/11/2024
Cultivar (Strain) or Sample De	Date Completed: 04/15/2024		

*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.

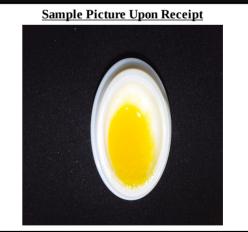
Moisture Content (%)	Water Activity (aw)	PASS/FAIL
Not Tested	Not Tested	PASS

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.

<u>Cannabinoids (Top 3)</u>	<u>(%)</u>	<u>mg/g</u>
THCA	76.6	766
CBGA	7.10	71.0
Δ9-ΤΗC	3.20	32.0
TOTAL CBD	-	_
TOTAL THC	70.4	704
TOTAL CANNABINOIDS	88.0	880
<u>Terpenes (Top 5)</u>	<u>(%)</u>	µg∕g
<u>Terpenes (Top 5)</u> β-Caryophyllene	<u>(%)</u> 2.02	<mark>µg/g</mark> 20200
β-Caryophyllene	2.02	20200
β-Caryophyllene α-Humulene	2.02 0.801	20200 8010
β-Caryophyllene α-Humulene Linalool	2.02 0.801 0.600	20200 8010 6000

PASS/FAIL
PASS
PASS
PASS
PASS





Scan the QR code to verify results.

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REPORT OF LABORATORY ANALYSIS

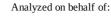
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Cultivar (Strain) or Sample	Date Completed: 04/15/2024		

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

Analysis Date/Time: 04/12/2024 1450 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> (<u>mg)</u>
CBC	ND	ND	0.110	0.256	-	-	-
CBCA	ND	ND	0.341	0.796	-	-	-
CBD	ND	ND	0.775	1.81	-	-	-
CBDA	ND	ND	0.285	0.666	-	-	-
CBDV	ND	ND	0.124	0.290	-	-	-
CBDVA	ND	ND	0.331	0.774	-	-	-
CBG	0.423	4.23	0.503	1.17	-	4.23	4.23
CBGA	7.10	71.0	0.713	0.841	-	71.0	71.0
CBL	ND	ND	0.581	1.36	-	-	-
CBN	ND	ND	0.267	0.623	-	-	-
CBNA	ND	ND	0.288	0.672	-	-	-
Δ9-ΤΗC	3.20	32.0	0.320	0.746	-	32.0	32.0
$\Delta 8$ -THC	ND	ND	0.499	1.16	-	-	-
THCA	76.6	766	0.173	0.406	-	766	766
THCV	ND	ND	0.416	0.970	-	-	-
THCVA	0.605	6.05	0.133	0.309	-	6.05	6.05
TOTAL	88.0	880			-	880	880
TOTAL CBC	-	-			-	-	-
TOTAL CBD	-	-			-	-	-
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	6.66	66.6			-	66.6	66.6
TOTAL CBN	-	-			-	-	-
TOTAL THC	70.4	704			-	704	704
TOTAL THCV	0.525	5.25			-	5.25	5.25

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.

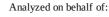
mg/g results are reported as mass component per mass material.













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License: ADH 113	License: 00065C	E20240410BNZLB03	Date Received: 04/11/2024
Cultivar (Strain) or Sample	Date Completed: 04/15/2024		

Analysis Date/Time:04/14/2024 1657 Analys

TERPENOID PROFILE
Method: GC/MS

Analyst: KF		Ins	trument: Agilent 7890/5975
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)	
α-Bisabolol	ND	-	
Camphene	356	0.0356	
δ-3-Carene	ND	-	
β-Caryophyllene	20200	2.02	
Caryophyllene oxide	234	0.0234	
p-Cymene	ND	-	
Eucalyptol	ND	-	
Geraniol	231	0.0231	
Guaiol	2740	0.274	1
α-Humulene	8010	0.801	
Isopulegol	ND	-	
d-Limonene	5590	0.559	
Linalool	6000	0.600	
β-Myrcene	4820	0.482	
cis-Nerolidol	ND	-	
trans-Nerolidol	2730	0.273	
α-Ocimene	283	0.0283	
β-Ocimene	75.7	0.00757	
α-Pinene	299	0.0299	
β-Pinene	863	0.0863	
α-Terpinene	ND	-	
γ-Terpinene	ND	-	
Terpinolene	<loq< td=""><td>-</td><td></td></loq<>	-	
TOTAL	FO 100	= 0.4	

TOTAL

52400

5.24

Deviations from SOP:

None

Abbreviations: GC - Gas Chromatography, MS - Mass Spectrometry, RL - Reporting Limit Abbreviations: ND - Not Detected. . LOD - Limit of Detection, LOQ - Limit of Quantitation

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Reporting Limit (µg/g): 48.2

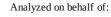
"-" Not detected above LOD.













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Cultivar (Strain) or Sample	Date Completed: 04/15/2024		

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

Analysis Date/Time:	04/11/20	04/11/2024 2147 Method: HS/GC/MS		Deviations 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)	-	126	252	5000	n-Heptane (142-82-5)	-	126	252	5000	
Acetonitrile (75-5-8)	-	126	252	410	n-Hexane (110-54-3)	-	44.0	88.1	290	
Benzene (71-43-2)	-	1.26	2.52	2	Isobutane (75-28-5)	-	126	252	5000	
n-Butane (106-97-2)	-	126	252	5000	Isopropanol (67-63-0)	-	126	252	5000	
1-Butanol (71-36-3)	-	126	252	5000	Isopropyl acetate	_	126	252	5000	
2-Butanol (78-92-2)	-	126	252	5000	(108-21-4)		120	202	5000	
2-Butanone (78-93-3)	-	126	252	5000	Isopropyl benzene (98-82-8)	-	12.6	25.2	70	
Cyclohexane (110-82-7)	-	126	252	3880	(98-62-6) Methanol (67-56-1)		126	252	3000	<u>Color Key</u>
1,2-Dimethoxyethane	_	12.6	25.2	100	2-Methylbutane (78-78-4)		126	252	5000	
(110-71-4) N,N-Dimethylacetamide	_	126	252	1090	Methylene chloride (75-9-2)	-	126	252	600	RESULT < AL RESULT > AL
(127-19-5) 2,2-Dimethylbutane (75-83-2)	_	44.0	88.1	290	2-Methylpentane (107-83-5)	-	44.0	88.1	290	"DET" detected less than LOQ
2,3-Dimethylbutane					3-Methylpentane (96-10-0)	-	44.0	88.1	290	"-" not detected above
(79-29-8)	-	44.0	88.1	290	n-Pentane (109-66-0)	-	126	252	5000	LOD
N,N-Dimethylformamide		100	252	880	1-Pentanol (71-41-0)	-	126	252	5000	"*" - o,m,p-Xylene and
(68-12-2)		126	252	000	n-Propane (74-98-6)	-	126	252	5000	Ethylbenzene
Dimethylsulfoxide (67-68-5)	-	126	252	5000	1-Propanol (71-23-8)	-	126	252	5000	Action levels are
1,4-Dioxane (123-91-1)	-	126	252	380	Pyridine (110-86-1)	-	44.0	88.1	200	referenced from the State of
Ethanol (64-17-5)	-	126	252	5000	Tetrahydrofuran (109-99-9)	-	126	252	720	Arkansas
2-Ethoxyethanol (110-80-5)	-	44.0	88.1	160	Tetramethylene sulfone (126-33-0)	-	44.0	88.1	160	MMJ testing
Ethyl ether (60-29-7)	-	126	252	5000	Toluene (108-88-3)	-	126	252	890	guidelines.
Ethyl acetate (141-78-6)	-	126	252	5000	o-Xylene (95-47-6)	-	126	252	2170	A value of "-"
Ethyl benzene (100-41-4)	-	126	252	2170	m,p-Xylene (108-38-3 or					for the action level
Ethylene glycol (107-21-1)	-	126	252	620	106-42-3)	-	126	252	2170	means that analyte is not currently
Ethylene oxide (75-21-8)	-	12.6	25.2	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170	regulated by the regulations referenced above.
Solvent		Synonym(s	<u>s)</u>		Solvent	<u>s</u>	ynonym(s)			0
Acetonitrile		Methyl Cya	anide, ACN		Ethylene glycol	1	,2-Ethanedio	1		
1-Butanol		n-Butanol,	Butyl Alco	hol	Isobutane	2	-Methylprop	ane		
2-Butanol		sec-Butyl a	lcohol		Isopropanol	2	-Propanol, IP	A		
2-Butanone		Methyl eth	yl ketone, N	/IEK	Isopropyl Acetate	А	cetic acid iso	propyl este	r	
1,2-Dimethoxyethane		Monoglym	e		Methanol	Ν	lethyl alcoho	1		
2,3-Dimethylbutane		Neohexane	1		2-Methylbutane	Is	sopentane			
2,3-Dimethylbutane		Diisopropy	/l		Methylene chloride	Γ	ichlorometha	ane		
N,N-Dimethylformamide		DMF			2-Methylpentane	Is	sohexane			
Dimethysufoxide		DMSO			1-Pentanol		-Amyl alcoho			
2-Ethoxyethanol		Cellosolve	, Ethyl glyc	ol	1-Propanol		ropyl alcoho	l		
Ethyl ether		Diethyl eth	ier, Ether		Tetrahydrofuran	Т	ΉF			
Ethyl acetate		EtOAc			Tetramethylene sulfone	S	ulfolane			
Ethyl benzene		Phenyletha	ne		Xylene	E	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



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA37616)

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Analysis Date/Time: 04/12/2024 1547 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> <u>Level</u> (µg/g)	
Abamectin (71751-41-2)	-	0.0479	0.383	0.5	Kresoxim-methyl	_	0.0479	0.383	0.4	
Acephate (30560-19-1)	-	0.0479	0.383	0.4	(143390-89-0)		0.0470	0.000	0.0	
Acequinocyl (57960-19-7)	-	0.0479	0.383	2	Malathion (121-75-5)	-	0.0479	0.383	0.2	
Acetamiprid (135410-20-7)	_	0.0479	0.383	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.0479 0.0479	0.383 0.383	0.2 0.2	Color Key
(135410-20-7) Aldicarb (116-06-3)		0.0479	0.383	0.4	Methomyl (16752-77-5)	-	0.0479	0.383	0.2	
Azoxystrobin					Methyl parathion	-				RESULT < AL
(131860-33-8)	-	0.0479	0.383	0.2	(298-0-0)	-	0.0479	0.383	0.2	RESULT > AL
Bifenazate (149877-41-8)	-	0.0479	0.383	0.2	MGK 264 (113-48-4)	-	0.0479	0.383	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3)	-	0.0479	0.383	0.2	Myclobutanil	_	0.0479	0.383	0.2	"-" not detected above
Boscalid (188425-85-6)	-	0.0479	0.383	0.4	(88671-89-0)					LOD
Carbaryl (63-25-2)	-	0.0479	0.383	0.2	Naled (300-76-5)	-	0.0479	0.383	0.5	Permethrins measured as the
Carbofuran (1563-66-2)	-	0.0479	0.383	0.2	Oxamyl (23135-22-0)	-	0.0479	0.383	1	cumulative residue of the cis- and
Chlorantraniliprole (800008-45-7)	-	0.0479	0.383	0.2	Paclobutrazol (76738-62-0)	-	0.0479	0.383	0.4	trans- permethrin isomers.
Chlorfenapyr		0.0479	0.383	1	Permethrins (52645-53-1)	-	0.0479	0.383	0.2	Pyrethrins measured as the cumulative residue of the
(122453-73-0)					Phosmet (732-11-6)	-	0.0479	0.383	0.2	pyrethrin I, cinerin I, and jasmolin
Chlorpyrifos (2921-88-2)	-	0.0479	0.383	0.2	Piperonyl butoxide (51-03-6)	-	0.0479	0.383	2	I isomers.
Clofentezine (74115-24-5)	-	0.0479	0.383	0.2	(31-03-0) Prallethrins (2331-36-9)		0.0479	0.383	0.2	Action levels are referenced from
Cyfluthrin (68359-37-5)	-	0.0479	0.383	1	Propiconazole					the
Cypermethrin (52315-07-8)	-	0.0479	0.383	1	(60207-90-1))	-	0.0479	0.383	0.4	State of Arkansas MMJ testing guidelines.
Daminozide (1596-84-5)	-	0.0479	0.383	1	Propoxur (114-26-1)	-	0.0479	0.383	0.2	A value of "-" for the action level
DDVP (62-73-7)	-	0.0479	0.383	0.1	Pyrethrins (8003-34-7)	-	0.0479	0.383	1	means that analyte is not
Diazinon (333-41-5)	-	0.0479	0.383	0.2	Pyridaben (96489-71-3)	-	0.0479	0.383	0.2	currently regulated by the regulations referenced above.
Dimethoate (60-51-5)	-	0.0479	0.383	0.2	Spinosad (168316-95-8)	-	0.0479	0.383	0.2	5
Ethoprophos (13194-48-4)	-	0.0479	0.383	0.2	Spiromesifen (283594-90-1)	-	0.0479	0.383	0.2	<i>Disclaimer:</i> This information is provided as a service and makes
Etofenprox (80844-07-1)	-	0.0479	0.383	0.4	Spirotetramat					no claims of efficacy and/or safety
Etoxazole (153233-91-1)	-	0.0479	0.383	0.2	(203313-25-1)	-	0.0479	0.383	0.2	of this product. Results are
Fenoxycarb (72490-01-8)	-	0.0479	0.383	0.2	Spiroxamine	_	0.0479	0.383	0.4	applicable only for the sample(s) analyzed and for the specific
(E)-Fenpyroximate (134098-61-6)	-	0.0479	0.383	0.4	(118134-30-8) Tebuconazole		0.0479	0.383	0.4	analysis conducted. This report is for informational purposes only
Fipronil (120068-37-3)	-	0.0479	0.383	0.4	(80443-41-0)	-	0.0479	0.365	0.4	and should not be used to
Flonicamid (158062-67-0)	-	0.0479	0.383	1	Thiacloprid	-	0.0479	0.383	0.2	diagnose, treat, or prevent any medical-related symptoms. The
Fludioxinil (131341-86-1)	-	0.0479	0.383	0.4	(111988-49-9)					statements and results herein have
Hexythiazox (78587-05-0)	-	0.0479	0.383	1	Thiamethoxam (153719-23-4)	-	0.0479	0.383	0.2	not been approved and/or endorsed by the FDA.
Imazalil (35554-44-0)	-	0.0479	0.383	0.2	Trifloxystrobin		0.0470	0.202	0.0	endorsed by the FDA.
Imidacloprid (138261-41-3)	-	0.0479	0.383	0.4	(141517-21-7)		0.0479	0.383	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** Synonym(s) Cyfluthrin Myclobutanil Baythroid Systhane Propiconazole Tilt DDVP Dichlorvos Naled Dibrom Propoxur Baygon Prophos Phosmet Imidan Ethoprophos



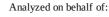
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Cultivar (Strain) or Sample	Date Completed: 04/15/2024		

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

Analysis Date/Time: 04/14/2024 1536 (ICP/OES) Analysis Date/Time: - (DMA) Analyst: KF			-	thod: ICP/MS trument: Agilent 7500ce	Deviations from SOP: 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)	-	58.9	93.2	200	\sim
Cadmium (Cd)	-	58.9	93.2	200	
Lead (Pb)	-	58.9	93.2	500	
Mercury (Hg)	-	58.9	93.2	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

<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.

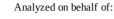
Disclaimer: This information is provided as a service and makes no claims of efficacy and/or safety of this product. Results are applicable only for the sample(s) analyzed and for the specific analysis conducted. This report is for informational purposes only and should not be used to diagnose, treat, or prevent any medical-related symptoms. The statements and results herein have not been approved and/or endorsed by the FDA.













Testing Location:	Customer ID: 2168	Sample ID: SA37616	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13245972560	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 04/11/2024
License: ADA 05_H273	License: 00065C	E20240410BNZLB03	Date Received: 04/11/2024
Cultivar (Strain) or Sample	Date Completed: 04/15/2024		

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 04/12/20 Analyst: PW		Iardy Diagnostics Compact nt: Thermo Incubator	Dry 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	-	
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





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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