



SUMMARY OF ANALYSIS (SAMPLE ID: SA36045)

Testing Location:	Customer ID: 2168	Order ID: OR10613	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13221531121	Mass: 10ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 09/18/2023
License: ADH 113	License: 00065C	P20230912BLU16	Date Received: 09/19/2023
Cultivar (Strain) or Sample D	Date Completed: 09/20/2023		

*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
0.000	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.

Cannabin	<u>oids (Top 3)</u>	<u>(%)</u>	<u>mg/g</u>		
С	BD	0.000	0.000		
CE	DA	0.000	0.000		
CE	DV	0.000	0.000		
TOTA	L CBD	0.000	0.000		
TOTA	L THC	0.000	0.000		
TOTAL CAN	INABINOIDS	0.000	0.000		
Terpen	<u>es (Top 5)</u>	<u>(%)</u>	µg∕g		
cis-Ne	erolidol	0.0498	498		
Gu	aiol	0.00374	37.4		
β-My	rcene	0.00218	21.8		
α-Hu	nulene	0.00156	15.6		
α-Bis	abolol				
TOTAL T	ERPENES	0.0573	573		
Contaminants	PASS/FAIL	Sample Picture Upon Receipt			

1100/11111
PASS
PASS
PASS
PASS





Scan the QR code to verify results.

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Kyle W. Felling, Ph.D. aboratory Director

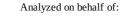
REPORT OF LABORATORY ANALYSIS

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Cultivar (Strain) or Sample D	Date Completed: 09/20/2023		

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

Analysis Date/Time: 09/19/2023 1151 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> (<u>mg/</u> <u>mL)</u>	<u>Per</u> Serving (mg)	<u>Per</u> 5 <u>Unit</u> (mg)
CBC	0.000	0.000	0.00213	0.00497	-	-	-
CBCA	0.000	0.000	0.00662	0.0154	-	-	-
CBD	0.000	0.000	0.0150	0.0351	-	0.242	2.42
CBDA	0.000	0.000	0.00553	0.0129	-	-	-
CBDV	0.000	0.000	0.00241	0.00563	-	-	-
CBDVA	0.000	0.000	0.00643	0.0150	-	-	-
CBG	0.000	0.000	0.00976	0.0228	-	0.266	2.66
CBGA	0.000	0.000	0.0138	0.0163	-	-	-
CBL	0.000	0.000	0.0113	0.0263	-	-	-
CBN	0.000	0.000	0.00518	0.0121	-	-	-
CBNA	0.000	0.000	0.00559	0.0130	-	-	-
Δ9-ΤΗC	0.000	0.000	0.00621	0.0145	-	9.23	92.3
$\Delta 8$ -THC	0.000	0.000	0.00968	0.0226	-	-	-
THCA	0.000	0.000	0.00336	0.00787	-	-	-
THCV	0.000	0.000	0.00807	0.0188	-	-	-
THCVA	0.000	0.000	0.00258	0.00600	-	-	-
TOTAL	0.000	0.000				9.74	97.4
TOTAL CBC	0.000	0.000	· · · · · · · · · · · · · · · · · · ·	,		-	
TOTAL CBD	0.000	0.000			-	0.242	2.42
TOTAL CBDV	0.000	0.000			-	-	-
TOTAL CBG	0.000	0.000			-	0.266	2.66
TOTAL CBN	0.000	0.000			-	-	-
TOTAL THC	0.000	0.000			-	9.23	92.3
TOTAL THCV	0.000	0.000			-	-	-

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 (%): 0.000 Water Activity (aw): -



SERVING MASS (g): 4.60 SERVINGS/UNIT: 10

"-" 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.

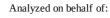
mg/g results are reported as mass component per mass material.

Kyle W. Felling, Ph.D. tory Directo











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Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 09/18/2023
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Cultivar (Strain) or Sample De	Date Completed: 09/20/2023		

TERPENOID PROFILE

Analysis Date/Time:09/2 Analyst: KF	20/2023 0708		ethod: GC/MS strument: Agilent 7890/5975	Deviations from SOP: None
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> <u>(%)</u>		
α-Bisabolol	ND	-		
Camphene	<loq< td=""><td>-</td><td></td><td></td></loq<>	-		
δ-3-Carene	ND	-		
β-Caryophyllene	ND	-		
Caryophyllene oxide	ND	-		
p-Cymene	ND	-		
Eucalyptol	ND	-		AM NET WEISCHICH
Geraniol	ND	-		
Guaiol	37.4	0.00374		Abburnistianes CC Core
α-Humulene	15.6	0.00156		Abbreviations: GC - Gas Chromatography, MS - Mass
Isopulegol	ND	-		Spectrometry, RL - Reporting Limit
d-Limonene	ND	-		<i>Abbreviations:</i> ND - Not Detected, , LOD - Limit of Detection, LOQ - Limit of Quantitation
Linalool	ND	-		This information is provided as a service
β-Myrcene cis-Nerolidol	21.8 498	0.00218 0.0498		and makes no claims of efficacy and/or safety of this product.
trans-Nerolidol	ND	-		Results are applicable only for the
α-Ocimene	ND	-		sample(s) analyzed and for the specific analysis conducted.
β-Ocimene	ND	-		This report is for informational purposes
α-Pinene	<loq< td=""><td>-</td><td></td><td>only and should not be used to diagnose, treat, or prevent any</td></loq<>	-		only and should not be used to diagnose, treat, or prevent any
β-Pinene	ND	-		medical-related symptoms.
α-Terpinene	ND	-		The statements and results herein have not been approved and/or endorsed by
γ-Terpinene	ND	-		the FDA.
Terpinolene	ND	-		
TOTAL	573	0.0573		Reporting Limit (µg/g): 7.79

"-" Not detected above LOD.













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

Analysis Date/Time:	09/19/20	023 1716		I	Method: HS/GC/MS			D	eviation	s from SOP:
Analyst: KF				I	nstrument: Agilent 78	90/5975		Ν	one	
<u>Solvent</u>	<u>Result</u> (µg/g)	LOD (µg/g)	LOQ (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)	-	145	290	5000	n-Heptane (142-82-5)	-	145	290	5000	08
Acetonitrile (75-5-8)	-	145	290	410	n-Hexane (110-54-3)	-	50.7	101	290	
Benzene (71-43-2)	-	1.45	2.90	2	Isobutane (75-28-5)	-	145	290	5000	Wana M
n-Butane (106-97-2)	-	145	290	5000	Isopropanol (67-63-0)	-	145	290	5000	SOUR QUINHIES HUREKAY
1-Butanol (71-36-3)	-	145	290	5000	Isopropyl acetate	_	145	290	5000	SUCCESSION
2-Butanol (78-92-2)	-	145	290	5000	(108-21-4)		145	250	5000	AM HTWISSHOW
2-Butanone (78-93-3)	-	145	290	5000	Isopropyl benzene (98-82-8)	-	14.5	29.0	70	
Cyclohexane (110-82-7)	-	145	290	3880	(98-62-8) Methanol (67-56-1)		145	290	3000	Color Key
1,2-Dimethoxyethane		14.5	29.0	100	2-Methylbutane (78-78-4)	-	145	290	5000	
(110-71-4) N,N-Dimethylacetamide		145	290	100	Methylene chloride (75-9-2)	-	145	290 290	600	RESULT < AL RESULT > AL
(127-19-5) 2,2-Dimethylbutane	-	50.7	101	290	2-Methylpentane (107-83-5)	-	50.7	101	290	"DET" detected less than LOQ
(75-83-2) 2,3-Dimethylbutane					3-Methylpentane (96-10-0)	-	50.7	101	290	"-" not detected above
(79-29-8)	-	50.7	101	290	n-Pentane (109-66-0)	-	145	290	5000	LOD
N,N-Dimethylformamide	_	145	290	880	1-Pentanol (71-41-0)	-	145	290	5000	"*" - o,m,p-Xylene and
(68-12-2)		145	250	000	n-Propane (74-98-6)	-	145	290	5000	Ethylbenzene
Dimethylsulfoxide (67-68-5)	-	145	290	5000	1-Propanol (71-23-8)	-	145	290	5000	Action lovals are
1,4-Dioxane (123-91-1)	_	145	290	380	Pyridine (110-86-1)	-	50.7	101	200	Action levels are referenced from the State of
Ethanol (64-17-5)	_	145	290	5000	Tetrahydrofuran (109-99-9)	-	145	290	720	Arkansas
2-Ethoxyethanol (110-80-5)	-	50.7	101	160	Tetramethylene sulfone (126-33-0)	-	50.7	101	160	MMJ testing
Ethyl ether (60-29-7)	-	145	290	5000	(120-33-0) Toluene (108-88-3)		145	290	890	guidelines.
Ethyl acetate (141-78-6)	-	145	290	5000	o-Xylene (95-47-6)		145	290	2170	A value of "-"
Ethyl benzene (100-41-4)	-	145	290	2170	m,p-Xylene (108-38-3 or					for the action level
Ethylene glycol (107-21-1)	-	145	290	620	106-42-3)	-	145	290	2170	means that analyte
Ethylene oxide (75-21-8)	-	14.5	29.0	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170	is not currently regulated by the regulations referenced above.
Solvent		Synonym(s	5)		Solvent	Sy	nonym(s)			5
Acetonitrile		Methyl Cya	anide, ACN		Ethylene glycol	1,	2-Ethanedio			
1-Butanol		n-Butanol,	Butyl Alco	hol	Isobutane	2-	Methylpropa	ine		
2-Butanol		sec-Butyl a	lcohol		Isopropanol	2-	Propanol, IP	A		
2-Butanone		Methyl ethy	yl ketone, N	/IEK	Isopropyl Acetate	A	cetic acid iso	propyl este	r	
1,2-Dimethoxyethane		Monoglym	e		Methanol	М	ethyl alcoho	l		
2,3-Dimethylbutane		Neohexane			2-Methylbutane	Is	opentane			
2,3-Dimethylbutane		Diisopropy	'l		Methylene chloride	D	ichlorometha	ine		
N,N-Dimethylformamide		DMF			2-Methylpentane	Is	ohexane			
Dimethysufoxide		DMSO			1-Pentanol		Amyl alcoho			
2-Ethoxyethanol		Cellosolve,		ol	1-Propanol		opyl alcohol			
Ethyl ether		Diethyl eth	er, Ether		Tetrahydrofuran	TI	HF			
Ethyl acetate		EtOAc			Tetramethylene sulfone	Su	ılfolane			
Ethyl benzene		Phenyletha	ne		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



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36045)

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Analysis Date/Time: 09/19/2023 1221 Analyst: KF

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Method: LC/MS/MS Instrument: Shimadzu LC-8050

Pesticide	<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)	-	0.00935	0.0748	0.5	Kresoxim-methyl	_	0.00935	0.0748	0.4	4
Acephate (30560-19-1)	-	0.00935	0.0748	0.4	(143390-89-0)					
Acequinocyl (57960-19-7)	-	0.00935	0.0748	2	Malathion (121-75-5)	-	0.00935	0.0748	0.2	Wana
Acetamiprid (135410-20-7)	-	0.00935	0.0748	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.00935 0.00935	0.0748 0.0748	0.2 0.2	SOUR GUMMES
Aldicarb (116-06-3)	-	0.00935	0.0748	0.4	Methomyl (16752-77-5)	-	0.00935	0.0748	0.4	
Azoxystrobin (131860-33-8)	-	0.00935	0.0748	0.2	Methyl parathion (298-0-0)	-	0.00935	0.0748	0.2	
Bifenazate (149877-41-8)	-	0.00935	0.0748	0.2	MGK 264 (113-48-4)	-	0.00935	0.0748	0.2	<u>Color Ke</u>
Bifenthrin (82657-04-3)	-	0.00935	0.0748	0.2	Myclobutanil		0.00935	0.0748	0.2	RESULT <
Boscalid (188425-85-6)	-	0.00935	0.0748	0.4	(88671-89-0)					RESULT >
Carbaryl (63-25-2)	-	0.00935	0.0748	0.2	Naled (300-76-5)	-	0.00935	0.0748	0.5	
Carbofuran (1563-66-2)	-	0.00935	0.0748	0.2	Oxamyl (23135-22-0)	-	0.00935	0.0748	1	"DET" detected less
Chlorantraniliprole (800008-45-7)	-	0.00935	0.0748	0.2	Paclobutrazol (76738-62-0)	-	0.00935	0.0748	0.4	"-" not detected LOD
Chlorfenapyr (122453-73-0)	-	0.00935	0.0748	1	Permethrins (52645-53-1)	-	0.00935	0.0748	0.2	Permethrins measu
Chlorpyrifos (2921-88-2)	-	0.00935	0.0748	0.2	Phosmet (732-11-6)	-	0.00935	0.0748	0.2	cumulative residue of trans- permethrin
Clofentezine (74115-24-5)	-	0.00935	0.0748	0.2	Piperonyl butoxide (51-03-6)	-	0.00935	0.0748	2	•
Cyfluthrin (68359-37-5)	-	0.00935	0.0748	1	(51-03-0) Prallethrins (2331-36-9)		0.00935	0.0748	0.2	Pyrethrins measur cumulative residu
Cypermethrin (52315-07-8)	-	0.00935	0.0748	1	Propiconazole (60207-90-1))	-	0.00935	0.0748	0.2	pyrethrin I, cinerin I, a I isomers
Daminozide (1596-84-5)	-	0.00935	0.0748	1	(00207-50-1)) Propoxur (114-26-1)		0.00935	0.0748	0.2	Action levels are refe
DDVP (62-73-7)	-	0.00935	0.0748	0.1	Pyrethrins (8003-34-7)	-	0.00935	0.0748	1	the
Diazinon (333-41-5)	-	0.00935	0.0748	0.2	Pyridaben (96489-71-3)	-	0.00935	0.0748	0.2	State of Arkansas M
Dimethoate (60-51-5)	-	0.00935	0.0748	0.2	Spinosad (168316-95-8)		0.00935	0.0748	0.2	guidelines
Ethoprophos (13194-48-4)	-	0.00935	0.0748	0.2	Spiromesifen (283594-90-1)	_	0.00935	0.0748	0.2	A value of "-" for the means that analy
Etofenprox (80844-07-1)	-	0.00935	0.0748	0.4	Spirotetramat					currently regulate regulations referen
Etoxazole (153233-91-1)	-	0.00935	0.0748	0.2	(203313-25-1)	-	0.00935	0.0748	0.2	regulations reference
Fenoxycarb (72490-01-8)	-	0.00935	0.0748	0.2	Spiroxamine		0.00935	0.0740	0.4	Disclaimer: This inf
(E)-Fenpyroximate (134098-61-6)	-	0.00935	0.0748	0.4	(118134-30-8) Tebuconazole	-		0.0748	0.4	provided as a service no claims of efficacy a of this product. Re
Fipronil (120068-37-3)	-	0.00935	0.0748	0.4	(80443-41-0)	-	0.00935	0.0748	0.4	applicable only for th
Flonicamid (158062-67-0)	-	0.00935	0.0748	1	Thiacloprid	_	0.00935	0.0748	0.2	analyzed and for th analysis conducted. T
Fludioxinil (131341-86-1)	-	0.00935	0.0748	0.4	(111988-49-9)					for informational pu
Hexythiazox (78587-05-0)	-	0.00935	0.0748	1	Thiamethoxam (153719-23-4)	-	0.00935	0.0748	0.2	and should not be
Imazalil (35554-44-0) Imidacloprid	-	0.00935	0.0748	0.2	Trifloxystrobin (141517-21-7)	-	0.00935	0.0748	0.2	diagnose, treat, or p medical-related sym statements and results
(138261-41-3)	-	0.00935	0.0748	0.4	(17101/21-/)					not been approve

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

x					
<u>Pesticide</u>	<u>Synonym(s)</u>	Pesticide	<u>Synonym(s)</u>	<u>Pesticide</u>	<u>Synonym(s)</u>
Cyfluthrin	Baythroid	Myclobutanil	Systhane	Propiconazole	Tilt
DDVP	Dichlorvos	Naled	Dibrom	Propoxur	Baygon
Ethoprophos	Prophos	Phosmet	Imidan		



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eferenced from MMJ testing

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ne action level lyte is not ated by the enced above.

nformation is ice and makes y and/or safety Results are the sample(s) the specific . This report is ourposes only be used to prevent any mptoms. The llts herein have ved and/or









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License: ADH 113	License: 00065C	P20230912BLU16	Date Received: 09/19/2023
Cultivar (Strain) or Sample	Date Completed: 09/20/2023		

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

Analysis Date/T Analysis Date/T Analyst: KF	ime: 09/19/2023 1 ime: - (DMA)	829 (ICP/OES)	_	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)	-	56.4	89.3	200	
Cadmium (Cd)	-	56.4	89.3	200	
Lead (Pb)	-	56.4	89.3	500	
Mercury (Hg)	-	56.4	89.3	100	CONTROL AND A CO

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.

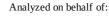
Disclaimer: This information is provided as a service and makes no claims of efficacy and/or safety of this product. Results are applicable only for the sample(s) analyzed and for the specific analysis conducted. This report is for informational purposes only and should not be used to diagnose, treat, 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: SA36045	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13221531121	Mass: 10ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 09/18/2023
License: ADA 05_H273	Date Received: 09/19/2023		
Cultivar (Strain) or Sample	Date Completed: 09/20/2023		

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 9/20/2023 Analyst: PW		Hardy Diagnostics CompactD nt: Thermo Incubator	bry 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	-	M
Pseudomonas aeruginosa	NT	-	
Salmonella spp.	NT	-	SOUR GUMMILS BUDGERNY
Staphylococcus aureus	NT	-	AN NIVELING

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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