



SUMMARY OF ANALYSIS (SAMPLE ID: SA35377)

Testing Location:	Customer ID: 2168	Order ID: OR10484	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13220972768	Mass: 10pieces
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 06/05/2023
License: ADH 113	License: 00065C	P20230601PEA012	Date Received: 06/05/2023
Cultivar (Strain) or Sample Do	Date Completed: 06/07/2023		

*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.

Cannabin	<u>pids (Top 3)</u>	<u>(%)</u>	<u>mg/g</u>
Δ9-'	ГНС	0.198	1.98
C	BC	0.0133	0.133
C	BG	0.00677	0.0677
TOTA	L CBD	-	-
TOTA	L THC	0.198	1.98
TOTAL CAN	INABINOIDS	0.222	2.22
Terpene	e <u>s (Top 5)</u>	<u>(%)</u>	ħā\ā
d-Lin	nonene	0.000419	4.19
ar Di			
u-Pl	nene	0.000419	4.19
	abolol	0.000419 0.000	4.19 0.000
α-Bis			
α-Bis Cam	abolol	0.000	0.000
α-Bis Cam δ-3-C	abolol phene	0.000 0.000	0.000 0.000

Containinanto	11100/11112
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

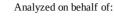
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Cultivar (Strain) or Sample	Date Completed: 06/07/2023		

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

Analysis Date/Time: 06/06/2023 1626 Analyst: PW Method: HPLC/DAD Instrument: Agilent 1100

<u>Cannabinoid</u>	<u>Result</u> (%)	<u>Result</u> (mg/g)	<u>LOD</u> (mg/g)	<u>LOQ</u> (mg/g)	<u>Result</u> (<u>mg/</u> <u>mL)</u>	<u>Per</u> <u>Serving</u> (mg)	<u>Per</u> 5 <u>Unit</u> (mg)
CBC	0.0133	0.133	0.00214	0.00499	-	0.669	6.69
CBCA	ND	ND	0.00664	0.0155	-	-	-
CBD	ND	ND	0.0151	0.0352	-	-	-
CBDA	ND	ND	0.00555	0.0130	-	-	-
CBDV	ND	ND	0.00242	0.00565	-	-	-
CBDVA	ND	ND	0.00645	0.0151	-	-	-
CBG	0.00677	0.0677	0.00980	0.0229	-	0.341	3.41
CBGA	ND	ND	0.0139	0.0164	-	-	-
CBL	ND	ND	0.0113	0.0264	-	-	-
CBN	0.00383	0.0383	0.00520	0.0121	-	0.193	1.93
CBNA	ND	ND	0.00561	0.0131	-	-	-
Δ9-ΤΗC	0.198	1.98	0.00623	0.0145	-	9.97	99.7
$\Delta 8$ -THC	ND	ND	0.00972	0.0227	-	-	-
THCA	ND	ND	0.00338	0.00790	-	-	-
THCV	ND	ND	0.00811	0.0189	-	-	-
THCVA	ND	ND	0.00259	0.00602	-	-	-
TOTAL	0.222	2.22				11.2	112
TOTAL CBC	0.0133	0.133				0.669	6.69
TOTAL CBD	-	-			-	-	-
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	0.00677	0.0677			-	0.341	3.41
TOTAL CBN	0.00383	0.0383			-	0.193	1.93
TOTAL THC	0.198	1.98			-	9.97	99.7
TOTAL THCV	-	-			-	-	-

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): 5.04 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.

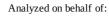
mg/g results are reported as mass component per mass material.

Kyle W. Felling, Ph.D. ory Dire











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Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 06/05/2023
License: ADH 113	License: 00065C	P20230601PEA012	Date Received: 06/05/2023
Cultivar (Strain) or Sample	Date Completed: 06/07/2023		

TERPENOID PROFILE

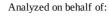
Analysis Date/Time:06/	06/2023 1932	M	ethod: GC/MS	Deviations from SOP:
Analyst: KF		Ins	strument: Agilent 7890/5975	None
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)		
α-Bisabolol	-	-		
Camphene	-	-		
δ-3-Carene	-	-		
β-Caryophyllene	-	-		PEACH BELLINI
Caryophyllene oxide	-	-		
p-Cymene	-	-		NT WE USE ((5))
Eucalyptol	-	-		
Geraniol	-	-		
Guaiol	-	-		Abbreviations: GC - Gas
α-Humulene	-	-		Chromatography, MS - Mass
Isopulegol	-	-		Spectrometry, RL - Reporting Limit
d-Limonene Linalool	4.19	0.000419		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
β-Myrcene cis-Nerolidol	-	-		sample(s) analyzed and for the specific analysis conducted.
trans-Nerolidol	-	-		This report is for informational purposes
α-Ocimene	-	-		only and should not be used to diagnose, treat, or prevent any
β-Ocimene	-	-		medical-related symptoms.
α-Pinene	4.19	0.000419		The statements and results herein have not been approved and/or endorsed by
β-Pinene	-	-		the FDA.
α-Terpinene	-	-		
γ-Terpinene	-	-		
Terpinolene	-	-		"-" Not detected above RL.
TOTAL	8.39	0.000839		Reporting Limit (µg/g): 12.6













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License: ADH 113	License: 00065C	P20230601PEA012	Date Received: 06/05/2023
Cultivar (Strain) or Sample	Date Completed: 06/07/2023		

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

Analysis Date/Time:	rsis Date/Time: 06/06/2023 1632 Method: HS/GC/MS				Deviations from SOP:						
Analyst: KF				Ι	nstrument: Agilent 78	90/5975		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> Level (µg/g)	_	
Acetone (67-64-1)	-	37.9	75.7	5000	n-Heptane (142-82-5)	-	37.9	75.7	5000		
Acetonitrile (75-5-8)	-	37.9	75.7	410	n-Hexane (110-54-3)	-	13.3	26.5	290		
Benzene (71-43-2)	-	0.379	0.757	2	Isobutane (75-28-5)	-	37.9	75.7	5000	DEACH BELLINI	
n-Butane (106-97-2)	-	37.9	75.7	5000	Isopropanol (67-63-0)	-	37.9	75.7	5000		
1-Butanol (71-36-3)	-	37.9	75.7	5000	Isopropyl acetate	_	37.9	75.7	5000	NET WE ISSUE (Gg)	
2-Butanol (78-92-2)	-	37.9	75.7	5000	(108-21-4)		57.5	/ 3./	5000		
2-Butanone (78-93-3)	-	37.9	75.7	5000	Isopropyl benzene (98-82-8)	-	3.79	7.57	70		
Cyclohexane (110-82-7)	-	37.9	75.7	3880	(98-82-8) Methanol (67-56-1)		37.9	75.7	3000	Color Key	
1,2-Dimethoxyethane		3.79	7.57	100	. ,	-					
(110-71-4) N,N-Dimethylacetamide					2-Methylbutane (78-78-4) Methylene chloride	-	37.9 37.9	75.7 75.7	5000 600	RESULT < AL RESULT > AL	
(127-19-5) 2,2-Dimethylbutane	-	37.9 13.3	75.7 26.5	1090 290	(75-9-2) 2-Methylpentane (107-83-5)	-	13.3	26.5	290	"DET" detected less than LOQ	
(75-83-2)		15.5	20.5	250	(107-83-3) 3-Methylpentane (96-10-0)		13.3	26.5	290	"-" not detected above	
2,3-Dimethylbutane (79-29-8)	-	13.3	26.5	290	n-Pentane (109-66-0)	-	37.9	20.3 75.7	5000	LOD	
(75-25-6) N,N-Dimethylformamide					1-Pentanol (71-41-0)	-	37.9	75.7	5000	-	
(68-12-2)	-	37.9	75.7	880	n-Propane (74-98-6)	-	37.9	75.7	5000	"*" - o,m,p-Xylene and	
Dimethylsulfoxide (67-68-5)	-	37.9	75.7	5000	1-Propanol (71-23-8)	-	37.9	75.7	5000	Ethylbenzene	
1,4-Dioxane (123-91-1)		37.9	75.7	380	Pyridine (110-86-1)	-	13.3	26.5	200	Action levels are referenced from the State of	
Ethanol (64-17-5)		37.9	75.7	5000	Tetrahydrofuran (109-99-9)	-	37.9	75.7	720	Arkansas	
2-Ethoxyethanol (110-80-5)		13.3	26.5	160	Tetramethylene sulfone	_	13.3	26.5	160	MMJ testing	
Ethyl ether (60-29-7)	_	37.9	75.7	5000	(126-33-0)					guidelines.	
Ethyl acetate (141-78-6)		37.9	75.7	5000	Toluene (108-88-3)	-	37.9	75.7	890	A value of "-"	
Ethyl benzene (100-41-4)		37.9	75.7	2170	o-Xylene (95-47-6)	-	37.9	75.7	2170	for the action level	
Ethylene glycol (107-21-1)	_	37.9	75.7	620	m,p-Xylene (108-38-3 or 106-42-3)	-	37.9	75.7	2170	means that analyte	
Ethylene oxide (75-21-8)		3.79	7.57	50	Xylenes* (1330-20-7)		43.3	86.7	2170	is not currently	
		-		50	· · · /		_	00.7	2170	regulated by the regulations referenced above.	
<u>Solvent</u>		Synonym(s			<u>Solvent</u>		ynonym(s)				
Acetonitrile		Methyl Cya			Ethylene glycol		,2-Ethanediol				
1-Butanol		n-Butanol,	5	hol	Isobutane		-Methylpropa				
2-Butanol		sec-Butyl a		mi	Isopropanol		-Propanol, IP/				
2-Butanone		Methyl ethy		1EK	Isopropyl Acetate		cetic acid iso		r		
1,2-Dimethoxyethane		Monoglym			Methanol		fethyl alcoho	l			
2,3-Dimethylbutane		Neohexane			2-Methylbutane		sopentane				
2,3-Dimethylbutane		Diisopropy	1		Methylene chloride		ichlorometha	ne			
N,N-Dimethylformamide		DMF			2-Methylpentane		sohexane				
Dimethysufoxide		DMSO			1-Pentanol		-Amyl alcoho				
2-Ethoxyethanol		Cellosolve,		10	1-Propanol		ropyl alcohol				
Ethyl ether		Diethyl eth	er, Ether		Tetrahydrofuran		ΉF				
Ethyl acetate		EtOAc			Tetramethylene sulfone		ulfolane				
Ethyl benzene		Phenyletha	ne		Xylene	D	imethylbenze	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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PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 06/06/2023 1447 Analyst: KF

(138261-41-3)

Method: LC/MS/MS

Instrument: Shimadzu LC-8050 And an Developed

<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)	-	0.00944	0.0755	0.5	Kresoxim-methyl	_	0.00944	0.0755	0.4	
Acephate (30560-19-1)	-	0.00944	0.0755	0.4	(143390-89-0)		0.000.4.4	0.0755	0.2	
Acequinocyl (57960-19-7)	-	0.00944	0.0755	2	Malathion (121-75-5)	-	0.00944	0.0755	0.2	
Acetamiprid (135410-20-7)	-	0.00944	0.0755	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.00944 0.00944	0.0755 0.0755	0.2 0.2	
Aldicarb (116-06-3)	-	0.00944	0.0755	0.4	Methomyl (16752-77-5)	-	0.00944	0.0755	0.4	
Azoxystrobin (131860-33-8)	-	0.00944	0.0755	0.2	Methyl parathion (298-0-0)	-	0.00944	0.0755	0.2	
Bifenazate (149877-41-8)	-	0.00944	0.0755	0.2	MGK 264 (113-48-4)	-	0.00944	0.0755	0.2	
Bifenthrin (82657-04-3)	-	0.00944	0.0755	0.2	Myclobutanil		0.00944	0.0755	0.2	
Boscalid (188425-85-6)	-	0.00944	0.0755	0.4	(88671-89-0)					
Carbaryl (63-25-2)	-	0.00944	0.0755	0.2	Naled (300-76-5)	-	0.00944	0.0755	0.5	
Carbofuran (1563-66-2)	-	0.00944	0.0755	0.2	Oxamyl (23135-22-0)	-	0.00944	0.0755	1	
Chlorantraniliprole (800008-45-7)	-	0.00944	0.0755	0.2	Paclobutrazol (76738-62-0)	-	0.00944	0.0755	0.4	
Chlorfenapyr (122453-73-0)	-	0.00944	0.0755	1	Permethrins (52645-53-1)	-	0.00944	0.0755	0.2	
Chlorpyrifos (2921-88-2)	-	0.00944	0.0755	0.2	Phosmet (732-11-6)	-	0.00944	0.0755	0.2	CL
Clofentezine (74115-24-5)	-	0.00944	0.0755	0.2	Piperonyl butoxide	_	0.00944	0.0755	2	
Cyfluthrin (68359-37-5)	-	0.00944	0.0755	1	(51-03-6) Prallethrins (2331-36-9)		0.00944	0.0755	0.2	
Cypermethrin (52315-07-8)	-	0.00944	0.0755	1	Propiconazole (60207-90-1))	-	0.00944	0.0755	0.2	ру
Daminozide (1596-84-5)	-	0.00944	0.0755	1	(00207-50-1)) Propoxur (114-26-1)		0.00944	0.0755	0.2	А
DDVP (62-73-7)	-	0.00944	0.0755	0.1	Pyrethrins (8003-34-7)	-	0.00944	0.0755	1	Г
Diazinon (333-41-5)	-	0.00944	0.0755	0.2	Pyridaben (96489-71-3)	-	0.00944	0.0755	0.2	
Dimethoate (60-51-5)	-	0.00944	0.0755	0.2	Spinosad (168316-95-8)		0.00944	0.0755	0.2	
Ethoprophos (13194-48-4)	-	0.00944	0.0755	0.2	Spiromesifen (283594-90-1)	-	0.00944	0.0755	0.2	A
Etofenprox (80844-07-1)	-	0.00944	0.0755	0.4	(283394-90-1) Spirotetramat					
Etoxazole (153233-91-1)	-	0.00944	0.0755	0.2	(203313-25-1)	-	0.00944	0.0755	0.2	
Fenoxycarb (72490-01-8)	-	0.00944	0.0755	0.2	Spiroxamine		0.00044	0.0755	0.4	Ι
(E)-Fenpyroximate (134098-61-6)	-	0.00944	0.0755	0.4	(118134-30-8) Tebuconazole	-	0.00944	0.0755	0.4	r no
Fipronil (120068-37-3)	-	0.00944	0.0755	0.4	(80443-41-0)	-	0.00944	0.0755	0.4	a
Flonicamid (158062-67-0)	-	0.00944	0.0755	1	Thiacloprid	_	0.00944	0.0755	0.2	
Fludioxinil (131341-86-1)	-	0.00944	0.0755	0.4	(111988-49-9)		2.000.4	5.67.55	0.2	ar f
Hexythiazox (78587-05-0)	-	0.00944	0.0755	1	Thiamethoxam (153719-23-4)	-	0.00944	0.0755	0.2	
Imazalil (35554-44-0)	-	0.00944	0.0755	0.2	Trifloxystrobin				A -	
Imidacloprid	-	0.00944	0.0755	0.4	(141517-21-7)	-	0.00944	0.0755	0.2	st

Deviations from SOP:

None

A -+-



Color Key

"DET" detected less than LOQ "-" not detected above LOD Permethrins measured as the cumulative residue of the cis- and trans-permethrin isomers. Pyrethrins measured as the

cumulative residue of the pyrethrin I, cinerin I, and jasmolin I isomers.

Action levels 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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Abbreviations: LC - Liquid Chromatography, MS - Mass Spectrometry, RL - Reporting Limit, AL - Action Level, CAS - Chemical Abstract Services, LOD - Limit of Detection, LOQ - Limit of Ouantification

Quantinearion					
<u>Pesticide</u>	<u>Synonym(s)</u>	Pesticide	<u>Synonym(s)</u>	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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232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13220972768	Mass: 10pieces
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 06/05/2023
License: ADH 113	License: 00065C	P20230601PEA012	Date Received: 06/05/2023
Cultivar (Strain) or Sample	Date Completed: 06/07/2023		

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

Analysis Date/Time: 06/06/2023 1501 (ICP/OES) Analysis Date/Time: - (DMA) Analyst: KF			Method: ICP/MS Instrument: 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)	-	59.4	94.0	200	
Cadmium (Cd)	-	59.4	94.0	200	
Lead (Pb)	-	59.4	94.0	500	
Mercury (Hg)	-	59.4	94.0	100	PEACH BELLINI
					AN REFERENCES

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



"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.











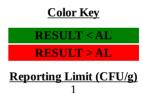


Testing Location:	Customer ID: 2168	Sample ID: SA35377	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13220972768	Mass: 10pieces
Greenbrier, AR 72058 Fort Smith, AR 72903 Production Run:		Production Run:	Date Collected: 06/05/2023
License: ADA 05_H273	License: 00065C	P20230601PEA012	Date Received: 06/05/2023
Cultivar (Strain) or Sample I	Date Completed: 06/07/2023		

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 06/07/20 Analyst: PW		l: Hardy Diagnostics CompactDry nent: 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	-	
Pseudomonas aeruginosa	NT	-	SE DE
Salmonella spp.	NT	-	
Staphylococcus aureus	NT	-	er m is 1304 (45g)

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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