



SUMMARY OF ANALYSIS (SAMPLE ID: SA37348)

Testing Location:	Customer ID: 2168	Order ID: OR10900	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13244202149	Mass: 4G
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 03/11/2024
License: ADH 113	License: 00065C	E20240308BZLD01	Date Received: 03/11/2024
Cultivar (Strain) or Sample D	Date Completed: 03/13/2024		

*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.

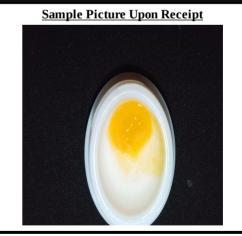
Moisture Content (%)	Water Activity (aw)	PASS/FAIL
Not Tested	Not Tested	PASS

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.

	(0/)	,
<u>Cannabinoids (Top 3)</u>	<u>(%)</u>	<u>mg/g</u>
THCA	78.9	789
CBGA	5.96	59.6
THCVA	0.602	6.02
TOTAL CBD	-	-
TOTAL THC	69.2	692
TOTAL CANNABINOIDS	85.4	854
<u>Terpenes (Top 5)</u>	<u>(%)</u>	µg∕g
<u>Terpenes (Top 5)</u> β-Caryophyllene	<u>(%)</u> 1.93	<mark>µg/g</mark> 19300
β-Caryophyllene	1.93	19300
β-Caryophyllene d-Limonene	1.93 1.31	19300 13100
β-Caryophyllene d-Limonene β-Myrcene	1.93 1.31 0.909	19300 13100 9090

PASS/FAIL
PASS
PASS
PASS
PASS





Scan the QR code to verify results.

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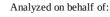
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Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 03/11/2024			
License: ADH 113	License: 00065C	E20240308BZLD01	Date Received: 03/11/2024			
Cultivar (Strain) or Sample Description: Banana Zoap Indica Hybrid Live Badder Date Completed: 03/13/202						

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

Analysis Date/Time: 03/12/2024 1733 Analyst: PW Method: HPLC/DAD Instrument: Agilent 1100

<u>Cannabinoid</u>	<u>Result</u> (<u>%</u>)	<u>Result</u> (mg/g)	<u>LOD</u> (<u>mg/g)</u>	LOQ (mg/g)	<u>Result</u> (mg/ mL)	<u>Per</u> <u>Serving</u> (mg)	<u>Per</u> 5 <u>Unit</u> (<u>mg)</u>
CBC	ND	ND	0.113	0.263	-	-	-
CBCA	ND	ND	0.351	0.818	-	-	-
CBD	ND	ND	0.796	1.86	-	-	-
CBDA	ND	ND	0.293	0.684	-	-	-
CBDV	ND	ND	0.128	0.298	-	-	-
CBDVA	ND	ND	0.341	0.795	-	-	-
CBG	ND	ND	0.517	1.21	-	-	-
CBGA	5.96	59.6	0.733	0.865	-	59.6	59.6
CBL	ND	ND	0.597	1.39	-	-	-
CBN	ND	ND	0.274	0.641	-	-	-
CBNA	ND	ND	0.296	0.690	-	-	-
Δ9-ΤΗC	ND	ND	0.329	0.767	-	-	-
$\Delta 8$ -THC	ND	ND	0.513	1.20	-	-	-
THCA	78.9	789	0.178	0.417	-	789	789
THCV	ND	ND	0.428	0.997	-	-	-
THCVA	0.602	6.02	0.137	0.318	-	6.02	6.02
TOTAL	85.4	854				854	854
TOTAL CBC	-	-			-	-	-
TOTAL CBD	-	-			-	-	-
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	5.23	52.3			-	52.3	52.3
TOTAL CBN	-	-			-	-	-
TOTAL THC	69.2	692			-	692	692
TOTAL THCV	0.522	5.22			-	5.22	5.22

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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|-001) 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.

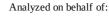
mg/g results are reported as mass component per mass material.













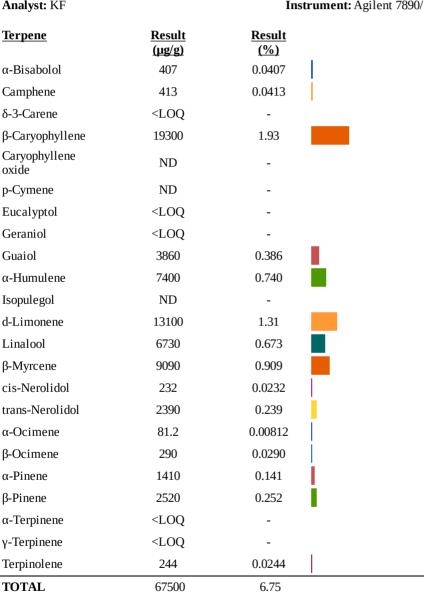
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Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 03/11/2024
License: ADH 113	License: 00065C	E20240308BZLD01	Date Received: 03/11/2024
Cultivar (Strain) or Sample	Date Completed: 03/13/2024		

Analysis Date/Time:03/12/2024 1745 Analyst: KF

TERPENOID PROFILE

Method: GC/MS Instrument: Agilent 7890/5975

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): 49.5

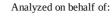
"-" Not detected above LOD.













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Cultivar (Strain) or Sample	Date Completed: 03/13/2024		

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

Analysis Date/Time:	03/12/20	024 0046		ľ	Method: HS/GC/MS			D	eviation	s from SOP:	
Analyst: KF				I	nstrument: Agilent 78	90/5975	5 None				
<u>Solvent</u>	<u>Result</u> (µg/g)	LOD (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)	-	87.1	174	5000	n-Heptane (142-82-5)	-	87.1	174	5000		
Acetonitrile (75-5-8)	-	87.1	174	410	n-Hexane (110-54-3)	-	30.5	61.0	290		
Benzene (71-43-2)	-	0.871	1.74	2	Isobutane (75-28-5)	-	87.1	174	5000		
n-Butane (106-97-2)	-	87.1	174	5000	Isopropanol (67-63-0)	-	87.1	174	5000		
1-Butanol (71-36-3)	-	87.1	174	5000	Isopropyl acetate	_	87.1	174	5000		
2-Butanol (78-92-2)	-	87.1	174	5000	(108-21-4)						
2-Butanone (78-93-3)	-	87.1	174	5000	Isopropyl benzene (98-82-8)	-	8.71	17.4	70		
Cyclohexane (110-82-7)	-	87.1	174	3880	Methanol (67-56-1)		87.1	174	3000	<u>Color Key</u>	
1,2-Dimethoxyethane	_	8.71	17.4	100	2-Methylbutane (78-78-4)	_	87.1	174	5000		
(110-71-4) N,N-Dimethylacetamide (127-19-5)	_	87.1	174	1090	Methylene chloride (75-9-2)	-	87.1	174	600	RESULT < AL RESULT > AL	
(127-19-5) 2,2-Dimethylbutane (75-83-2)	_	30.5	61.0	290	2-Methylpentane (107-83-5)	-	30.5	61.0	290	"DET" detected less than LOQ	
2,3-Dimethylbutane					3-Methylpentane (96-10-0)	-	30.5	61.0	290	"-" not detected above	
(79-29-8)	-	30.5	61.0	290	n-Pentane (109-66-0)	-	87.1	174	5000	LOD	
N,N-Dimethylformamide		07.1	174	880	1-Pentanol (71-41-0)	-	87.1	174	5000	"*" - o,m,p-Xylene and	
(68-12-2)	-	87.1	174	880	n-Propane (74-98-6)	-	87.1	174	5000	Ethylbenzene	
Dimethylsulfoxide	_	87.1	174	5000	1-Propanol (71-23-8)	-	87.1	174	5000	- 9	
(67-68-5)		071	174	380	Pyridine (110-86-1)	-	30.5	61.0	200	Action levels are	
1,4-Dioxane (123-91-1)	-	87.1		5000	Tetrahydrofuran (109-99-9)	-	87.1	174	720	referenced from the State of Arkansas	
Ethanol (64-17-5) 2-Ethoxyethanol (110-80-5)	-	87.1 30.5	174 61.0	160	Tetramethylene sulfone		30.5	61.0	160	MMJ testing	
Ethyl ether (60-29-7)	-	30.3 87.1	174	5000	(126-33-0)					guidelines.	
Ethyl acetate (141-78-6)	-	87.1	174 174	5000	Toluene (108-88-3)	-	87.1	174	890	A value of "-"	
Ethyl benzene (100-41-4)	-	87.1	174 174	2170	o-Xylene (95-47-6)	-	87.1	174	2170	for the action level	
Ethylene glycol (107-21-1)		87.1	174	620	m,p-Xylene (108-38-3 or 106-42-3)	-	87.1	174	2170	means that analyte	
Ethylene oxide (75-21-8)	-	8.71	174	50	Xylenes* (1330-20-7)		43.3	86.7	2170	is not currently	
2 ,	-	•		50	,	-	-	00.7	2170	regulated by the regulations referenced above.	
Solvent		Synonym(s			<u>Solvent</u>	_	Synonym(s)				
Acetonitrile		Methyl Cya			Ethylene glycol		,2-Ethanediol				
1-Butanol		n-Butanol,	5	hol	Isobutane		-Methylpropa				
2-Butanol		sec-Butyl a		E K	Isopropanol		-Propanol, IP/				
2-Butanone		Methyl eth		1EK	Isopropyl Acetate		Acetic acid iso		ľ		
1,2-Dimethoxyethane		Monoglym			Methanol		Aethyl alcoho	1			
2,3-Dimethylbutane		Neohexane			2-Methylbutane		sopentane				
2,3-Dimethylbutane		Diisopropy	/1		Methylene chloride		Dichlorometha	ine			
N,N-Dimethylformamide		DMF			2-Methylpentane		sohexane	,			
Dimethysufoxide		DMSO		1	1-Pentanol		-Amyl alcoho				
2-Ethoxyethanol		Cellosolve		01	1-Propanol		ropyl alcohol				
Ethyl ether		Diethyl eth	ier, Ether		Tetrahydrofuran		THF				
Ethyl acetate		EtOAc			Tetramethylene sulfone		Sulfolane				
Ethyl benzene		Phenyletha	ne		Xylene	Γ	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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License: ADH 113	License: 00065C	E20240308BZLD01	Date Received: 03/11/2024
Cultivar (Strain) or Sample	Date Completed: 03/13/2024		

Analysis Date/Time: 03/12/2024 2118 Analyst: KF

PESTICIDES PROFILE (SOP: SOP-PEST-001)

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

Analyst: KF	Instrument: Shimadzu LC-8050				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)	Pesticide	<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)	-	0.0492	0.394	0.5	Kresoxim-methyl (143390-89-0)	_	0.0492	0.394	0.4	
Acephate (30560-19-1)	-	0.0492	0.394	0.4	· · · · · ·		0.0492	0.394	0.2	
Acequinocyl (57960-19-7)	-	0.0492	0.394	2	Malathion (121-75-5)	-	0.0492	0.394	0.2	
Acetamiprid (135410-20-7)	-	0.0492	0.394	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.0492	0.394	0.2	
Aldicarb (116-06-3)	-	0.0492	0.394	0.4	Methomyl (16752-77-5)	-	0.0492	0.394	0.4	
Azoxystrobin (131860-33-8)	-	0.0492	0.394	0.2	Methyl parathion (298-0-0)	-	0.0492	0.394	0.2	
Bifenazate (149877-41-8)	-	0.0492	0.394	0.2	MGK 264 (113-48-4)	-	0.0492	0.394	0.2	"DET" d
Bifenthrin (82657-04-3)	-	0.0492	0.394	0.2	Myclobutanil		0.0492	0.394	0.2	"-" T
Boscalid (188425-85-6)	-	0.0492	0.394	0.4	(88671-89-0)					- 1
Carbaryl (63-25-2)	-	0.0492	0.394	0.2	Naled (300-76-5)	-	0.0492	0.394	0.5	D (
Carbofuran (1563-66-2)	-	0.0492	0.394	0.2	Oxamyl (23135-22-0)	-	0.0492	0.394	1	Permet
Chlorantraniliprole (800008-45-7)	-	0.0492	0.394	0.2	Paclobutrazol (76738-62-0)	-	0.0492	0.394	0.4	trans-
Chlorfenapyr		0.0492	0.394	1	Permethrins (52645-53-1)	-	0.0492	0.394	0.2	Pyreth cumu
(122453-73-0)	-	0.0492	0.394	1	Phosmet (732-11-6)	-	0.0492	0.394	0.2	pyrethrin
Chlorpyrifos (2921-88-2)	-	0.0492	0.394	0.2	Piperonyl butoxide	_	0.0492	0.394	2	
Clofentezine (74115-24-5)	-	0.0492	0.394	0.2	(51-03-6)		0.0400	0.004		Action le
Cyfluthrin (68359-37-5)	-	0.0492	0.394	1	Prallethrins (2331-36-9)	-	0.0492	0.394	0.2	recton re
Cypermethrin (52315-07-8)	-	0.0492	0.394	1	Propiconazole (60207-90-1))	-	0.0492	0.394	0.4	State of
Daminozide (1596-84-5)	-	0.0492	0.394	1	Propoxur (114-26-1)	-	0.0492	0.394	0.2	A value o
DDVP (62-73-7)	-	0.0492	0.394	0.1	Pyrethrins (8003-34-7)	-	0.0492	0.394	1	mean
Diazinon (333-41-5)	-	0.0492	0.394	0.2	Pyridaben (96489-71-3)	-	0.0492	0.394	0.2	curren
Dimethoate (60-51-5)	-	0.0492	0.394	0.2	Spinosad (168316-95-8)	-	0.0492	0.394	0.2	regulati
Ethoprophos (13194-48-4)	-	0.0492	0.394	0.2	Spiromesifen	-	0.0492	0.394	0.2	Disclaim
Etofenprox (80844-07-1)	-	0.0492	0.394	0.4	(283594-90-1)					provided no claims
Etoxazole (153233-91-1)	-	0.0492	0.394	0.2	Spirotetramat (203313-25-1)	-	0.0492	0.394	0.2	of this
Fenoxycarb (72490-01-8)	-	0.0492	0.394	0.2	Spiroxamine		0.0400	0.004		applicabl
(E)-Fenpyroximate (134098-61-6)	-	0.0492	0.394	0.4	(118134-30-8) Tebuconazole	-	0.0492	0.394	0.4	analyze analysis c for inforr
Fipronil (120068-37-3)	-	0.0492	0.394	0.4	(80443-41-0)	-	0.0492	0.394	0.4	and sh
Flonicamid (158062-67-0)	_	0.0492	0.394	1	Thiacloprid		0.0402	0.204	0.2	diagnos
Fludioxinil (131341-86-1)	_	0.0492	0.394	0.4	(111988-49-9)		0.0492	0.394	0.2	medical- statements
Hexythiazox (78587-05-0)	_	0.0492	0.394	1	Thiamethoxam	_	0.0492	0.394	0.2	not be
Imazalil (35554-44-0)	-	0.0492	0.394	0.2	(153719-23-4)					end
Imidacloprid (138261-41-3)	-	0.0492	0.394	0.4	Trifloxystrobin (141517-21-7)	-	0.0492	0.394	0.2	
·										

Deviations from SOP:

None



<u>Color Key</u>

RESULT < AL RESULT > AL ET" detected less than LOQ "-" not detected above LOD ermethrins measured as the ulative residue of the *cis*- and *rans*- permethrin isomers. Pyrethrins measured as the cumulative residue of the thrin I, cinerin I, and jasmolin I isomers.

ction 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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Synonym(s)

Tilt

Baygon

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** Cyfluthrin Myclobutanil Baythroid Systhane Propiconazole DDVP Dichlorvos Naled Dibrom Propoxur Prophos Phosmet Imidan Ethoprophos



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

Analysis Date/T Analysis Date/T Analyst: KF	`ime: 03/12/2024 2 `ime: - (DMA)	2158 (ICP/OES)	_	t hod: ICP/MS rument: Agilent 7500	Deviations from SOP: ce 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)	-	57.7	91.4	200	
Cadmium (Cd)	-	57.7	91.4	200	
Lead (Pb)	-	57.7	91.4	500	
Mercury (Hg)	-	57.7	91.4	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.

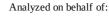
Disclaimer: This information is provided as a service and makes no claims of efficacy and/or safety of this product. Results are applicable only for the sample(s) analyzed and for the specific analysis conducted. This report is for informational purposes only and should not be used to diagnose, treat, 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: SA37348	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13244202149	Mass: 4G
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 03/11/2024
License: ADA 05_H273	License: 00065C	E20240308BZLD01	Date Received: 03/11/2024
Cultivar (Strain) or Sample	Date Completed: 03/13/2024		

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 03/12/20 Analyst: PW		lardy Diagnostics Compac at: Thermo Incubator	tDry 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





1

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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