



SUMMARY OF ANALYSIS (SAMPLE ID: SA36048)

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	M00065C13228154652	Mass: 1ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 09/18/2023
License: ADH 113	License: 00065C	P20230911straw19	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
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>Cannabino</u>	ids (Top 3)	<u>(%)</u>	mg/g
Δ9-7	ГНС	1.58	15.8
CE	BD	0.133	1.33
CE	3G	0.0572	0.572
TOTA	L CBD	0.133	1.33
TOTA	L THC	1.58	15.8
TOTAL CAN	NABINOIDS	1.80	18.0
Terpene	<u>s (Top 5)</u>	<u>(%)</u>	ћã∖ã
cis-Ne	rolidol	0.00435	43.5
Caryophyl	lene oxide	0.00254	25.4
α-Hum	nulene	0.00181	18.1
α-Oci	mene	0.00181	18.1
Gua	aiol	0.00109	10.9
		0.0110	110
TOTAL TI	ERPENES	0.0116	116

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





Scan the QR code to verify results.

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Kyle W. Felling, Ph.D. aboratory Directo

REPORT OF LABORATORY ANALYSIS

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License: ADH 113	License: 00065C	P20230911straw19	Date Received: 09/19/2023
Cultivar (Strain) or Sample	Date Completed: 09/20/2023		

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

Analysis Date/Time: 09/19/2023 1243 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> Unit (mg)
CBC	ND	ND	0.00248	0.00579	-	-	-
CBCA	ND	ND	0.00770	0.0180	-	-	-
CBD	0.133	1.33	0.0175	0.0408	-	0.834	25.0
CBDA	ND	ND	0.00644	0.0150	-	-	-
CBDV	ND	ND	0.00281	0.00655	-	-	-
CBDVA	ND	ND	0.00748	0.0175	-	-	-
CBG	0.0572	0.572	0.0114	0.0265	-	0.360	10.8
CBGA	ND	ND	0.0161	0.0190	-	-	-
CBL	ND	ND	0.0131	0.0306	-	-	-
CBN	0.0180	0.180	0.00602	0.0141	-	0.113	3.40
CBNA	ND	ND	0.00650	0.0152	-	-	-
Δ9-ΤΗC	1.58	15.8	0.00722	0.0168	-	9.91	297
$\Delta 8$ -THC	ND	ND	0.0113	0.0263	-	-	-
THCA	ND	ND	0.00392	0.00916	-	-	-
THCV	0.0140	0.140	0.00940	0.0219	-	0.0877	2.63
THCVA	ND	ND	0.00300	0.00698	-	-	-
TOTAL	1.80	18.0			-	11.3	339
TOTAL CBC	-	-			-	-	-
TOTAL CBD	0.133	1.33			-	0.834	25.0
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	0.0572	0.572			-	0.360	10.8
TOTAL CBN	0.0180	0.180			-	0.113	3.40
TOTAL THC	1.58	15.8			-	9.91	297
TOTAL THCV	0.0140	0.140			-	0.0877	2.63

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): 0.628 SERVINGS/UNIT: 30

"-" 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.

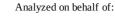
mg/g results are reported as mass component per mass material.

W. Felling, Ph.D. ory Dire











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License: ADH 113	License: 00065C	P20230911straw19	Date Received: 09/19/2023
Cultivar (Strain) or Sample	Date Completed: 09/20/2023		

Analysis Date/Time:09/20/2023 0835 Analyst: KF

TERPENOID PROFILE

Method: GC/MS

Analyst: KF		Instrument: Agilent 7890/597			
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)			
α-Bisabolol	<loq< td=""><td>-</td><td></td></loq<>	-			
Camphene	ND	-			
δ-3-Carene	ND	-			
β-Caryophyllene	ND	-			
Caryophyllene oxide	25.4	0.00254			
p-Cymene	ND	-			
Eucalyptol	ND	-			
Geraniol	ND	-			
Guaiol	10.9	0.00109			
α-Humulene	18.1	0.00181			
Isopulegol	ND	-			
d-Limonene	ND	-			
Linalool	ND	-			
β-Myrcene	ND	-			
cis-Nerolidol	43.5	0.00435			
trans-Nerolidol	ND	-			
α-Ocimene	18.1	0.00181			
β-Ocimene	ND	-			
α-Pinene	ND	-			
β-Pinene	ND	-			
α-Terpinene	ND	-			
γ-Terpinene	ND	-			
Terpinolene	ND	-			
TOTAL	116	0.0116			



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): 9.06

"-" Not detected above LOD.













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

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

Analysis Date/Time: Analyst: KF	09/19/20	023 1923			Method: HS/GC/MS instrument: Agilent 78	90/5975			eviation	s from SOP:
Solvent	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	Solvent	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	RIVER VALLEY PROVISIONS
Acetone (67-64-1)	_	143	286	5000	n-Heptane (142-82-5)	-	143	286	5000	
Acetonitrile (75-5-8)	-	143	286	410	n-Hexane (110-54-3)	-	50.0	100	290	STRAWBERKIFS T
Benzene (71-43-2)	-	1.43	2.86	2	Isobutane (75-28-5)	-	143	286	5000	
n-Butane (106-97-2)	-	143	286	5000	Isopropanol (67-63-0)	-	143	286	5000	- FREEZE DRIED SLICES
1-Butanol (71-36-3)	-	143	286	5000	Isopropyl acetate		142	200	5000	
2-Butanol (78-92-2)	-	143	286	5000	(108-21-4)	-	143	286	5000	
2-Butanone (78-93-3)	-	143	286	5000	Isopropyl benzene	_	14.3	28.6	70	
Cyclohexane (110-82-7)	-	143	286	3880	(98-82-8)					Color Key
1,2-Dimethoxyethane		140	20.0	100	Methanol (67-56-1)	925	143	286	3000	<u>Color Rey</u>
(110-71-4)	-	14.3	28.6	100	2-Methylbutane (78-78-4)	-	143	286	5000	RESULT < AL
N,N-Dimethylacetamide (127-19-5)	-	143	286	1090	Methylene chloride (75-9-2)	-	143	286	600	RESULT > AL
2,2-Dimethylbutane (75-83-2)	-	50.0	100	290	2-Methylpentane (107-83-5)	-	50.0	100	290	"DET" detected less than LOQ
2,3-Dimethylbutane		50.0	100	290	3-Methylpentane (96-10-0)	-	50.0	100	290	"-" not detected above
(79-29-8)		30.0	100	290	n-Pentane (109-66-0)	-	143	286	5000	LOD
N,N-Dimethylformamide	_	143	286	880	1-Pentanol (71-41-0)	-	143	286	5000	"*" - o,m,p-Xylene and
(68-12-2)		-			n-Propane (74-98-6)	-	143	286	5000	Ethylbenzene
Dimethylsulfoxide (67-68-5)	-	143	286	5000	1-Propanol (71-23-8)	-	143	286	5000	
1,4-Dioxane (123-91-1)	_	143	286	380	Pyridine (110-86-1)	-	50.0	100	200	Action levels are referenced from the State of
Ethanol (64-17-5)	1620	143	286	5000	Tetrahydrofuran (109-99-9)	-	143	286	720	Arkansas
2-Ethoxyethanol (110-80-5)	-	50.0	100	160	Tetramethylene sulfone	_	50.0	100	160	MMJ testing
Ethyl ether (60-29-7)	_	143	286	5000	(126-33-0)			200	000	guidelines.
Ethyl acetate (141-78-6)	_	143	286	5000	Toluene (108-88-3)	-	143	286	890	A value of "-"
Ethyl benzene (100-41-4)	_	143	286	2170	o-Xylene (95-47-6)	-	143	286	2170	for the action level
Ethylene glycol (107-21-1)	_	143	286	620	m,p-Xylene (108-38-3 or 106-42-3)	-	143	286	2170	means that analyte
Ethylene oxide (75-21-8)	_	14.3	28.6	50	Xylenes* (1330-20-7)	_	43.3	86.7	2170	is not currently
Solvent		Synonym(s			Solvent	s	ynonym(s)			regulated by the regulations referenced above.
Acetonitrile		Methyl Cya			Ethylene glycol		,2-Ethanedio			
1-Butanol		n-Butanol,	,		Isobutane		-Methylpropa			
2-Butanol		sec-Butyl a		1101	Isopropanol		-Propanol, IP			
2-Butanone		Methyl ethy		NEK	Isopropyl Acetate		cetic acid iso			
1,2-Dimethoxyethane		Monoglym		ILK	Methanol		fethyl alcoho		Ĺ	
2,3-Dimethylbutane		Neohexane			2-Methylbutane		sopentane	L		
2,3-Dimethylbutane		Diisopropy			Methylene chloride		opentane Dichlorometha			
		DIISOPIOPY	/1		5			lile		
N,N-Dimethylformamide Dimethysufoxide		DMF			2-Methylpentane 1-Pentanol		sohexane	1		
5			Ethyl alr-				-Amyl alcoho			
2-Ethoxyethanol		Cellosolve,		.01	1-Propanol Tetrahydrofyran		ropyl alcohol `HF			
Ethyl ether		Diethyl eth	ier, Ether		Tetrahydrofuran					
Ethyl acetate		EtOAc			Tetramethylene sulfone		ulfolane			
Ethyl benzene		Phenyletha	ne		Xylene	D	Dimethylbenz	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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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36048)

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

Analysis Date/Time: 09/19/2023 1322 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> <u>Level</u>	<u>Pesticide</u>	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> <u>Level</u>	RIVER VALLEY PROVISIO
	(µs/s)			<u>(µg/g)</u>		(µg/g)	(µg/g)	(µg/g)	<u>(µg/g)</u>	STRAWBERRIES
Abamectin (71751-41-2)	-	0.0108	0.0867	0.5	Kresoxim-methyl (143390-89-0)	-	0.0108	0.0867	0.4	- FREEZE DRIED SLICES
Acephate (30560-19-1)	-	0.0108	0.0867	0.4	Malathion (121-75-5)		0.0108	0.0867	0.2	
Acequinocyl (57960-19-7)	-	0.0108	0.0867	2	Metalaxyl (57837-19-1)	_	0.0108	0.0867	0.2	
Acetamiprid (135410-20-7)	-	0.0108	0.0867	0.2	Methiocarb (2032-65-7)	-	0.0108	0.0867	0.2	<u>Color Key</u>
Aldicarb (116-06-3)	-	0.0108	0.0867	0.4	Methomyl (16752-77-5)	-	0.0108	0.0867	0.4	RESULT < AL
Azoxystrobin (131860-33-8)	-	0.0108	0.0867	0.2	Methyl parathion (298-0-0)	-	0.0108	0.0867	0.2	RESULT > AL
Bifenazate (149877-41-8)	-	0.0108	0.0867	0.2	MGK 264 (113-48-4)	-	0.0108	0.0867	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3)	-	0.0108	0.0867	0.2	Myclobutanil	_	0.0108	0.0867	0.2	"-" not detected above
Boscalid (188425-85-6)	-	0.0108	0.0867	0.4	(88671-89-0)					LOD
Carbaryl (63-25-2)	-	0.0108	0.0867	0.2	Naled (300-76-5)	-	0.0108	0.0867	0.5	Permethrins measured as the
Carbofuran (1563-66-2)	-	0.0108	0.0867	0.2	Oxamyl (23135-22-0)	-	0.0108	0.0867	1	cumulative residue of the <i>cis</i> - and
Chlorantraniliprole (800008-45-7)	-	0.0108	0.0867	0.2	Paclobutrazol (76738-62-0)	-	0.0108	0.0867	0.4	<i>trans</i> - permethrin isomers.
Chlorfenapyr		0.0108	0.0867	1	Permethrins (52645-53-1)	-	0.0108	0.0867	0.2	Pyrethrins measured as the cumulative residue of the
(122453-73-0)					Phosmet (732-11-6)	-	0.0108	0.0867	0.2	pyrethrin I, cinerin I, and jasmolin
Chlorpyrifos (2921-88-2)	-	0.0108	0.0867	0.2	Piperonyl butoxide (51-03-6)	-	0.0108	0.0867	2	I isomers.
Clofentezine (74115-24-5)	-	0.0108	0.0867	0.2	(31-03-0) Prallethrins (2331-36-9)		0.0108	0.0867	0.2	Action levels are referenced from
Cyfluthrin (68359-37-5)	-	0.0108	0.0867	1	Propiconazole					the
Cypermethrin (52315-07-8)	-	0.0108	0.0867	1	(60207-90-1))	-	0.0108	0.0867	0.4	State of Arkansas MMJ testing guidelines.
Daminozide (1596-84-5)	-	0.0108	0.0867	1	Propoxur (114-26-1)	-	0.0108	0.0867	0.2	A value of "-" for the action level
DDVP (62-73-7)	-	0.0108	0.0867	0.1	Pyrethrins (8003-34-7)	-	0.0108	0.0867	1	means that analyte is not
Diazinon (333-41-5)	-	0.0108	0.0867	0.2	Pyridaben (96489-71-3)	-	0.0108	0.0867	0.2	currently regulated by the regulations referenced above.
Dimethoate (60-51-5)	-	0.0108	0.0867	0.2	Spinosad (168316-95-8)	-	0.0108	0.0867	0.2	0
Ethoprophos (13194-48-4)	-	0.0108	0.0867	0.2	Spiromesifen (283594-90-1)	-	0.0108	0.0867	0.2	Disclaimer: This information is provided as a service and makes
Etofenprox (80844-07-1)	-	0.0108	0.0867	0.4	Spirotetramat		0.0100	0.0007		no claims of efficacy and/or safety
Etoxazole (153233-91-1)	-	0.0108	0.0867	0.2	(203313-25-1)	-	0.0108	0.0867	0.2	of this product. Results are
Fenoxycarb (72490-01-8)	-	0.0108	0.0867	0.2	Spiroxamine	_	0.0108	0.0867	0.4	applicable only for the sample(s) analyzed and for the specific
(E)-Fenpyroximate (134098-61-6)	-	0.0108	0.0867	0.4	(118134-30-8) Tebuconazole	_	0.0108	0.0867	0.4	analysis conducted. This report is for informational purposes only
Fipronil (120068-37-3)	-	0.0108	0.0867	0.4	(80443-41-0)		0.0100	0.0007	0.4	and should not be used to diagnose, treat, or prevent any
Flonicamid (158062-67-0)	-	0.0108	0.0867	1	Thiacloprid (111988-49-9)	-	0.0108	0.0867	0.2	medical-related symptoms. The
Fludioxinil (131341-86-1)	DET	0.0108	0.0867	0.4	Thiamethoxam					statements and results herein have
Hexythiazox (78587-05-0)	-	0.0108	0.0867	1	(153719-23-4)	-	0.0108	0.0867	0.2	not been approved and/or endorsed by the FDA.
Imazalil (35554-44-0)	-	0.0108	0.0867	0.2	Trifloxystrobin		0.0108	0.0867	0.2	chaoisca by the FDA.
Imidacloprid (138261-41-3)	-	0.0108	0.0867	0.4	(141517-21-7)		0.0108	0.000/	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 Pesticide Pesticide Synonym(s) Synonym(s) Synonym(s) Cyfluthrin Baythroid Tilt Myclobutanil Systhane Propiconazole DDVP Baygon Dichlorvos Naled Dibrom Propoxur Ethoprophos Prophos Phosmet Imidan



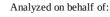
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HEAVY METAL PROFILE (SOP: SOP-ICP-200.6)

Analysis Date/Time: 09/19/2023 1849 (ICP/OES) Analysis Date/Time: - (DMA) Analyst: KF				thod: ICP/MS trument: Agilent 7500cc	Deviations from SOP: e None
Heavy Metal Arsenic (As) Cadmium (Cd) Lead (Pb) Mercury (Hg)	<u>Result</u> (µg/kg) - - - -	LOD (μg/kg) 55.6 55.6 55.6 55.6	LOQ (µg/kg) 88.1 88.1 88.1 88.1	Action Level (µg/kg) 200 200 500 100	RIVER VALLEY PROVISIONS STRAWBERRIES FREEZE DRIED SLICES

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.

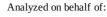
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Testing Location:	Customer ID: 2168	Sample ID: SA36048	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13228154652	Mass: 1ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 09/18/2023
License: ADA 05_H273	License: 00065C	P20230911straw19	Date Received: 09/19/2023
Cultivar (Strain) or Sample	Date Completed: 09/20/2023		

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 9/20/202 Analyst: PW		Hardy Diagnostics CompactDry nt: Thermo Incubator	Deviations from SOP: None
<u>Bacteria/Microbe</u>	<u>Result</u> (CFU/g)	<u>Action Level</u> (CFU/g)	RIVER VALLEY PROVISIONS
Aerobic Plate Count	NT	-	
Coliforms, Total	Absent	1	ATD AWIDED DUE
Escherichia Coli (E. Coli)	Absent	100	SI KAWBEKKIPS 📈
Mold/Yeast	NT	-	VIII CIIII CIIII CIIII
Pseudomonas aeruginosa	NT	-	FREEZE DRIED SLICES
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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