



SUMMARY OF ANALYSIS (SAMPLE ID: SA36685)

Testing Location:	Customer ID: 2168	Order ID: OR10768	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13244795806	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 12/18/2023
License: ADH 113	License: 00065C	E20231215LRPPA01	Date Received: 12/18/2023
Cultivar (Strain) or Sample D	Date Completed: 12/22/2023		

*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>Cannabinoids (Top 3)</u>	<u>(%)</u>	<u>mg/g</u>
THCA	79.2	792
CBGA	2.20	22.0
Δ9-ТНС	1.76	17.6
TOTAL CBD	-	-
TOTAL THC	71.2	712
TOTAL CANNABINOIDS	83.8	838
<u>Terpenes (Top 5)</u>	<u>(%)</u>	µg∕g
<u>Terpenes (Top 5)</u> d-Limonene	<u>(%)</u> 2.19	<mark>µg/g</mark> 21900
d-Limonene	2.19	21900
d-Limonene β-Myrcene	2.19 0.706	21900 7060
d-Limonene β-Myrcene β-Caryophyllene	2.19 0.706 0.551	21900 7060 5510

<u>Contaminants</u>	PASS/FAIL				
Heavy Metals:	PASS				
Microbiology:	PASS				
Pesticides:	PASS				
Residual Solvents:	PASS				





Scan the QR code to verify results.

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Cultivar (Strain) or Sample	Date Completed: 12/22/2023		

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

Analysis Date/Time: 12/19/2023 2110 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> g <u>Unit</u> (mg)
CBC	ND	ND	0.111	0.258	-	-	-
CBCA	ND	ND	0.344	0.803	-	-	-
CBD	ND	ND	0.781	1.82	-	-	-
CBDA	ND	ND	0.288	0.671	-	-	-
CBDV	ND	ND	0.125	0.292	-	-	-
CBDVA	ND	ND	0.334	0.780	-	-	-
CBG	0.445	4.45	0.507	1.18	-	4.45	4.45
CBGA	2.20	22.0	0.719	0.848	-	22.0	22.0
CBL	ND	ND	0.586	1.37	-	-	-
CBN	ND	ND	0.269	0.629	-	-	-
CBNA	ND	ND	0.291	0.677	-	-	-
Δ9-ΤΗC	1.76	17.6	0.323	0.752	-	17.6	17.6
$\Delta 8$ -THC	ND	ND	0.503	1.17	-	-	-
THCA	79.2	792	0.175	0.409	-	792	792
THCV	ND	ND	0.420	0.978	-	-	-
THCVA	0.137	1.37	0.134	0.312	-	1.37	1.37
TOTAL	83.8	838			-	838	838
TOTAL CBC	-	-			-	-	-
TOTAL CBD	-	-			-	-	-
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	2.38	23.8			-	23.8	23.8
TOTAL CBN	-	-			-	-	-
TOTAL THC	71.2	712			-	712	712
TOTAL THCV	0.119	1.19			-	1.19	1.19

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.

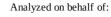
mg/g results are reported as mass component per mass material.













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License: ADH 113	License: 00065C	E20231215LRPPA01	Date Received: 12/18/2023
Cultivar (Strain) or Sample	Date Completed: 12/22/2023		

Analysis Date/Time:12/21/2023 2019 Analyst: KF

FERPENOID PROFIL	E

Method: GC/MS

Analyst: KF		In	strument: Agilent 7890/5975
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)	
α-Bisabolol	ND	-	1.44
Camphene	518	0.0518	
δ-3-Carene	ND	-	
β-Caryophyllene	5510	0.551	
Caryophyllene oxide	ND	-	
p-Cymene	130	0.0130	
Eucalyptol	ND	-	
Geraniol	ND	-	
Guaiol	4280	0.428	
α-Humulene	648	0.0648	Abbre Chron
Isopulegol	1720	0.172	Spectr
d-Limonene	21900	2.19	Abbre LOD
Linalool	1390	0.139	of Qu
β-Myrcene	7060	0.706	This in and m
cis-Nerolidol	ND	-	safety
trans-Nerolidol	ND	-	Result
α-Ocimene	ND	-	sample analys
β-Ocimene	ND	-	This r
α-Pinene	1620	0.162	only a treat, c
β-Pinene	3820	0.382	medic
α-Terpinene	ND	-	The st not be
γ-Terpinene	ND	-	the FL
Terpinolene	194	0.0194	
TOTAL	48800	4.88	



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): 81.0

"-" Not detected above LOD.













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Cultivar (Strain) or Sample	Date Completed: 12/22/2023		

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

Analysis Date/Time:	12/19/20)23 2114			Method: HS/GC/MS					s 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)	-	141	282	5000	n-Heptane (142-82-5)	-	141	282	5000	
Acetonitrile (75-5-8)	-	141	282	410	n-Hexane (110-54-3)	-	49.3	98.6	290	
Benzene (71-43-2)	-	1.41	2.82	2	Isobutane (75-28-5)	-	141	282	5000	
n-Butane (106-97-2)	-	141	282	5000	Isopropanol (67-63-0)	-	141	282	5000	
1-Butanol (71-36-3)	-	141	282	5000	Isopropyl acetate	_	141	282	5000	
2-Butanol (78-92-2)	-	141	282	5000	(108-21-4)		141	202	5000	
2-Butanone (78-93-3)	-	141	282	5000	Isopropyl benzene (98-82-8)	-	14.1	28.2	70	
Cyclohexane (110-82-7)	-	141	282	3880	(38-62-8) Methanol (67-56-1)		141	282	3000	Color Key
1,2-Dimethoxyethane	_	14.1	28.2	100	2-Methylbutane (78-78-4)		141	282	5000	
(110-71-4) N,N-Dimethylacetamide	_	141	282	1090	Methylene chloride (75-9-2)		141	282	600	RESULT < AL RESULT > AL
(127-19-5) 2,2-Dimethylbutane (75-83-2)	-	49.3	98.6	290	2-Methylpentane (107-83-5)	-	49.3	98.6	290	"DET" detected less than LOQ
2,3-Dimethylbutane					3-Methylpentane (96-10-0)	-	49.3	98.6	290	"-" not detected above
(79-29-8)	-	49.3	98.6	290	n-Pentane (109-66-0)	-	141	282	5000	LOD
N,N-Dimethylformamide (68-12-2)	-	141	282	880	1-Pentanol (71-41-0)	-	141 141	282 282	5000 5000	"*" - o,m,p-Xylene and
Dimethylsulfoxide			202		n-Propane (74-98-6)	-	141 141	282	5000	Ethylbenzene
(67-68-5)	-	141	282	5000	1-Propanol (71-23-8) Pyridine (110-86-1)	-	49.3	282 98.6	200	Action levels are
1,4-Dioxane (123-91-1)	-	141	282	380	Tetrahydrofuran (109-99-9)	-	49.3 141	282	200 720	referenced from the State of
Ethanol (64-17-5)	-	141	282	5000	Tetramethylene sulfone	-	141	202	720	Arkansas
2-Ethoxyethanol (110-80-5)	-	49.3	98.6	160	(126-33-0)	-	49.3	98.6	160	MMJ testing guidelines.
Ethyl ether (60-29-7)	-	141	282	5000	Toluene (108-88-3)	-	141	282	890	guidennes.
Ethyl acetate (141-78-6)	-	141	282	5000	o-Xylene (95-47-6)	-	141	282	2170	A value of "-"
Ethyl benzene (100-41-4)	-	141	282	2170	m,p-Xylene (108-38-3 or		141	282	2170	for the action level means that analyte
Ethylene glycol (107-21-1)	-	141	282	620	106-42-3)	-	141	202	2170	is not currently
Ethylene oxide (75-21-8)	-	14.1	28.2	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170	regulated by the regulations referenced above.
Solvent		Synonym(s			<u>Solvent</u>		nonym(s)			
Acetonitrile		Methyl Cya	nide, ACN		Ethylene glycol	1,2	2-Ethanedio			
1-Butanol		n-Butanol,	Butyl Alco	hol	Isobutane	2-	Methylpropa	ine		
2-Butanol		sec-Butyl a	lcohol		Isopropanol	2-	Propanol, IP	£		
2-Butanone		Methyl ethy	yl ketone, N	1EK	Isopropyl Acetate	A	cetic acid iso	propyl este	r	
1,2-Dimethoxyethane		Monoglym	e		Methanol	Μ	ethyl alcoho	1		
2,3-Dimethylbutane		Neohexane			2-Methylbutane	Ise	opentane			
2,3-Dimethylbutane		Diisopropy	rl		Methylene chloride	Di	ichlorometha	ine		
N,N-Dimethylformamide		DMF			2-Methylpentane	Ise	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	Di	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: SA36685)

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PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 12/19/2023 2240 Analyst: KF

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> Level (µg/g)	
Abamectin (71751-41-2)	-	0.0483	0.386	0.5	Kresoxim-methyl		0.0483	0.386	0.4	
Acephate (30560-19-1)	-	0.0483	0.386	0.4	(143390-89-0)					
Acequinocyl (57960-19-7)	-	0.0483	0.386	2	Malathion (121-75-5)	-	0.0483	0.386	0.2	—
Acetamiprid (135410-20-7)	-	0.0483	0.386	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.0483 0.0483	0.386 0.386	0.2 0.2	<u>Color Key</u>
Aldicarb (116-06-3)	-	0.0483	0.386	0.4	Methomyl (16752-77-5)	-	0.0483	0.386	0.4	RESULT < AL
Azoxystrobin (131860-33-8)	-	0.0483	0.386	0.2	Methyl parathion (298-0-0)	-	0.0483	0.386	0.2	RESULT > AL
Bifenazate (149877-41-8)	-	0.0483	0.386	0.2	MGK 264 (113-48-4)	-	0.0483	0.386	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3)	-	0.0483	0.386	0.2	Myclobutanil	_	0.0483	0.386	0.2	"-" not detected above
Boscalid (188425-85-6)	-	0.0483	0.386	0.4	(88671-89-0)					LOD
Carbaryl (63-25-2)	-	0.0483	0.386	0.2	Naled (300-76-5)	-	0.0483	0.386	0.5	Permethrins measured as the
Carbofuran (1563-66-2)	-	0.0483	0.386	0.2	Oxamyl (23135-22-0)	-	0.0483	0.386	1	cumulative residue of the <i>cis</i> - and
Chlorantraniliprole (800008-45-7)	-	0.0483	0.386	0.2	Paclobutrazol (76738-62-0)	-	0.0483	0.386	0.4	trans- permethrin isomers.
Chlorfenapyr	_	0.0483	0.386	1	Permethrins (52645-53-1)	-	0.0483	0.386	0.2	Pyrethrins measured as the cumulative residue of the
(122453-73-0)					¹ Phosmet (732-11-6)	-	0.0483	0.386	0.2	pyrethrin I, cinerin I, and jasmolin
Chlorpyrifos (2921-88-2)	-	0.0483	0.386	0.2	Piperonyl butoxide (51-03-6)	-	0.0483	0.386	2	I isomers.
Clofentezine (74115-24-5)	-	0.0483	0.386	0.2	(31-03-0) Prallethrins (2331-36-9)		0.0483	0.386	0.2	Action levels are referenced from
Cyfluthrin (68359-37-5)	-	0.0483	0.386	1	Propiconazole					the
Cypermethrin (52315-07-8)	-	0.0483	0.386	1	(60207-90-1))	-	0.0483	0.386	0.4	State of Arkansas MMJ testing guidelines.
Daminozide (1596-84-5)	-	0.0483	0.386	1	Propoxur (114-26-1)	-	0.0483	0.386	0.2	A value of "-" for the action level
DDVP (62-73-7)	-	0.0483	0.386	0.1	Pyrethrins (8003-34-7)	-	0.0483	0.386	1	means that analyte is not
Diazinon (333-41-5)	-	0.0483	0.386	0.2	Pyridaben (96489-71-3)	-	0.0483	0.386	0.2	currently regulated by the regulations referenced above.
Dimethoate (60-51-5)	-	0.0483	0.386	0.2	Spinosad (168316-95-8)	-	0.0483	0.386	0.2	5
Ethoprophos (13194-48-4)	-	0.0483	0.386	0.2	Spiromesifen (283594-90-1)	-	0.0483	0.386	0.2	<i>Disclaimer</i> : This information is provided as a service and makes
Etofenprox (80844-07-1)	-	0.0483	0.386	0.4	Spirotetramat					no claims of efficacy and/or safety
Etoxazole (153233-91-1)	-	0.0483	0.386	0.2	(203313-25-1)	-	0.0483	0.386	0.2	of this product. Results are
Fenoxycarb (72490-01-8)	-	0.0483	0.386	0.2	Spiroxamine	_	0.0483	0.386	0.4	applicable only for the sample(s) analyzed and for the specific
(E)-Fenpyroximate (134098-61-6)	-	0.0483	0.386	0.4	(118134-30-8) Tebuconazole		0.0483	0.386	0.4	analysis conducted. This report is for informational purposes only
Fipronil (120068-37-3)	-	0.0483	0.386	0.4	(80443-41-0)		0.0405	0.500	0.4	and should not be used to
Flonicamid (158062-67-0)	-	0.0483	0.386	1	Thiacloprid (111988-49-9)	-	0.0483	0.386	0.2	diagnose, treat, or prevent any medical-related symptoms. The
Fludioxinil (131341-86-1)	-	0.0483	0.386	0.4	(111988-49-9) Thiamethoxam					statements and results herein have
Hexythiazox (78587-05-0)	-	0.0483	0.386	1	(153719-23-4)	-	0.0483	0.386	0.2	not been approved and/or endorsed by the FDA.
Imazalil (35554-44-0)	-	0.0483	0.386	0.2	Trifloxystrobin		0.0483	0.386	0.2	endoised by the FDA.
Imidacloprid (138261-41-3)	-	0.0483	0.386	0.4	(141517-21-7)	-	0.0483	0.386	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) Cyfluthrin Baythroid Myclobutanil Systhane Propiconazole DDVP Dichlorvos Naled Dibrom Propoxur Ethoprophos Prophos Phosmet Imidan



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Synonym(s)

Tilt

Baygon

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

Analysis Date/Time: 12/19/2023 2141 (ICP/OES) Analysis Date/Time: - (DMA) Analyst: KF			_	t hod: ICP/MS rument: Agilent 750	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)	-	59.6	94.4	200	
Cadmium (Cd)	-	59.6	94.4	200	
Lead (Pb)	-	59.6	94.4	500	
Mercury (Hg)	-	59.6	94.4	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.

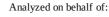
Disclaimer: This information is provided as a service and makes no claims of efficacy and/or safety of this product. Results are applicable only for the sample(s) analyzed and for the specific analysis conducted. This report is for informational purposes only and should not be used to diagnose, treat, 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: SA36685	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13244795806	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 12/18/2023
License: ADA 05_H273	License: 00065C	E20231215LRPPA01	Date Received: 12/18/2023
Cultivar (Strain) or Sample	Date Completed: 12/22/2023		

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 12/20/20 Analyst: PW		Hardy Diagnostics Compac nt: 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





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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