



SUMMARY OF ANALYSIS (SAMPLE ID: SA35791)

Testing Location:	Customer ID: 2168	Order ID: OR10562	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13245106176	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 08/07/2023
License: ADH 113	License: 00065C	E20230731LCBLR01	Date Received: 08/08/2023
Cultivar (Strain) or Sample D	Date Completed: 08/10/2023		

*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.

Cannabinoids (Top 3)	<u>(%)</u>	mg/g
THCA	34.4	344
Δ9-ΤΗΟ	27.2	272
CBGA	1.90	19.0
TOTAL CBD	0.129	1.29
TOTAL THC	57.4	574
TOTAL CANNABINOIDS	66.6	666
<u>Terpenes (Top 5)</u>	<u>(%)</u>	<u>µg∕g</u>
<u>Terpenes (Top 5)</u> β-Myrcene	(%) 6.42	<mark>µg/g</mark> 64200
β-Myrcene	6.42	64200
β-Myrcene d-Limonene	6.42 1.99	64200 19900
β-Myrcene d-Limonene β-Caryophyllene	6.42 1.99 1.67	64200 19900 16700

PASS/FAIL
PASS
PASS
PASS
PASS





Scan the QR code to verify results.

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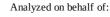
REPORT OF LABORATORY ANALYSIS

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Cultivar (Strain) or Sample I	Date Completed: 08/10/2023		

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

Analysis Date/Time: 08/08/2023 1429 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> (mg/ <u>mL)</u>	<u>Per</u> <u>Serving</u> (mg)	<u>Per</u> g <u>Unit</u> (mg)
CBC	0.332	3.32	0.110	0.258	-	3.32	3.32
CBCA	1.26	12.6	0.343	0.800	-	12.6	12.6
CBD	ND	ND	0.779	1.82	-	-	-
CBDA	0.148	1.48	0.287	0.670	-	1.48	1.48
CBDV	ND	ND	0.125	0.292	-	-	-
CBDVA	ND	ND	0.333	0.778	-	-	-
CBG	0.903	9.03	0.506	1.18	-	9.03	9.03
CBGA	1.90	19.0	0.717	0.846	-	19.0	19.0
CBL	ND	ND	0.584	1.36	-	-	-
CBN	ND	ND	0.268	0.627	-	-	-
CBNA	ND	ND	0.290	0.675	-	-	-
Δ9-ΤΗC	27.2	272	0.322	0.750	-	272	272
$\Delta 8$ -THC	ND	ND	0.502	1.17	-	-	-
THCA	34.4	344	0.174	0.408	-	344	344
THCV	ND	ND	0.419	0.976	-	-	-
THCVA	0.377	3.77	0.134	0.311	-	3.77	3.77
TOTAL	66.6	666			-	666	666
TOTAL CBC	1.43	14.3			-	14.3	14.3
TOTAL CBD	0.129	1.29			-	1.29	1.29
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	2.57	25.7			-	25.7	25.7
TOTAL CBN	-	-			-	-	-
TOTAL THC	57.4	574			-	574	574
TOTAL THCV	0.327	3.27			-	3.27	3.27

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.

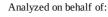
mg/g results are reported as mass component per mass material.













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License: ADH 113	License: 00065C	E20230731LCBLR01	Date Received: 08/08/2023
Cultivar (Strain) or Sample	Date Completed: 08/10/2023		

Analysis Date/Time:08/08/2023 1521 Analyst: KF

Result

 $(\mu g/g)$

465

543

ND

16700

310

<LOQ

ND

ND

426

6380

58.1

19900

1100

64200

ND

2330

1300

1880

3450

3800

96.9

155

756

124000

Terpene

α-Bisabolol

Camphene

δ-3-Carene

Eucalyptol

α-Humulene

Isopulegol

Linalool

β-Myrcene

cis-Nerolidol

α-Ocimene

β-Ocimene

α-Pinene

β-Pinene

α-Terpinene

y-Terpinene

Terpinolene TOTAL

trans-Nerolidol

d-Limonene

Geraniol

Guaiol

oxide p-Cymene

β-Caryophyllene

Caryophyllene

TERPENOID PROFILE

Method: GC/MS Instrument: Agilent 7890/5975

<u>Result</u>

(%)

0.0465

0.0543

_

1.67

0.0310

0.0426

0.638

0.00581

1.99

0.110

6.42

_

0.233

0.130

0.188

0.345

0.380

0.00969

0.0155

0.0756

12.4

Deviations from SOP
None



Spectrometry, RL - Reporting Limit Abbreviations: ND - Not Detected. . LOD - Limit of Detection, LOQ - Limit

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Reporting Limit (µg/g): 48.5

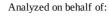
"-" Not detected above LOD.













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Cultivar (Strain) or Sample	Date Completed: 08/10/2023		

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

Analysis Date/Time:	08/08/20	023 2156		ľ	Method: HS/GC/MS		Deviations from SOP:		s from SOP:	
Analyst: KF				I	nstrument: Agilent 78	90/5975		N	one	
<u>Solvent</u>	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	LOQ (µg/g)	<u>Action</u> Level (µg/g)	<u>Solvent</u>	<u>Result</u> (µg/g)	LOD (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	
Acetone (67-64-1)	-	118	235	5000	n-Heptane (142-82-5)	-	118	235	5000	
Acetonitrile (75-5-8)	-	118	235	410	n-Hexane (110-54-3)	-	41.2	82.4	290	and the second
Benzene (71-43-2)	-	1.18	2.35	2	Isobutane (75-28-5)	-	118	235	5000	
n-Butane (106-97-2)	1570	118	235	5000	Isopropanol (67-63-0)	-	118	235	5000	
1-Butanol (71-36-3)	-	118	235	5000	Isopropyl acetate	_	118	235	5000	
2-Butanol (78-92-2)	-	118	235	5000	(108-21-4)		110	255	3000	
2-Butanone (78-93-3)	-	118	235	5000	Isopropyl benzene (98-82-8)	-	11.8	23.5	70	
Cyclohexane (110-82-7)	-	118	235	3880	(98-82-8) Methanol (67-56-1)		118	235	3000	Color Key
1,2-Dimethoxyethane		11.8	23.5	100	. ,	-				
(110-71-4)		11.0	23.5	100	2-Methylbutane (78-78-4)	-	118	235	5000	RESULT < AL
N,N-Dimethylacetamide (127-19-5)	-	118	235	1090	Methylene chloride (75-9-2)	-	118	235	600	RESULT > AL
2,2-Dimethylbutane (75-83-2)	-	41.2	82.4	290	2-Methylpentane (107-83-5)	-	41.2	82.4	290	"DET" detected less than LOQ
2,3-Dimethylbutane		41.2	82.4	290	3-Methylpentane (96-10-0)	-	41.2	82.4	290	"-" not detected above
(79-29-8)		41.2	02.4	290	n-Pentane (109-66-0)	-	118	235	5000	LOD
N,N-Dimethylformamide	_	118	235	880	1-Pentanol (71-41-0)	-	118	235	5000	"*" - o,m,p-Xylene and
(68-12-2)					n-Propane (74-98-6)	-	118	235	5000	Ethylbenzene
Dimethylsulfoxide (67-68-5)	-	118	235	5000	1-Propanol (71-23-8)	-	118	235	5000	A stion levels are
1,4-Dioxane (123-91-1)	_	118	235	380	Pyridine (110-86-1)	-	41.2	82.4	200	Action levels are referenced from the State of
Ethanol (64-17-5)	669	118	235	5000	Tetrahydrofuran (109-99-9)	-	118	235	720	Arkansas
2-Ethoxyethanol (110-80-5)	-	41.2	82.4	160	Tetramethylene sulfone	_	41.2	82.4	160	MMJ testing
Ethyl ether (60-29-7)	_	118	235	5000	(126-33-0)			225	000	guidelines.
Ethyl acetate (141-78-6)	_	118	235	5000	Toluene (108-88-3)	-	118	235	890	A value of "-"
Ethyl benzene (100-41-4)	_	118	235	2170	o-Xylene (95-47-6)	-	118	235	2170	for the action level
Ethylene glycol (107-21-1)	_	118	235	620	m,p-Xylene (108-38-3 or 106-42-3)	-	118	235	2170	means that analyte
Ethylene oxide (75-21-8)	_	11.8	23.5	50	Xylenes* (1330-20-7)	_	43.3	86.7	2170	is not currently
		1110	2010	50	11516165 (1555 257)		1010	001/	21/0	regulated by the regulations referenced above.
Solvent		Synonym(s	5)		Solvent	S	ynonym(s)			regulations referenced above.
Acetonitrile		Methyl Cya	anide, ACN		Ethylene glycol	1	,2-Ethanediol			
1-Butanol		n-Butanol,	Butyl Alco	hol	Isobutane	2	-Methylpropa	ine		
2-Butanol		sec-Butyl a	lcohol		Isopropanol	2	-Propanol, IP/	Ą		
2-Butanone		Methyl ethy	vl ketone, N	/IEK	Isopropyl Acetate		Acetic acid iso		r	
1,2-Dimethoxyethane		Monoglym			Methanol		/lethyl alcoho			
2,3-Dimethylbutane		Neohexane			2-Methylbutane		sopentane			
2,3-Dimethylbutane		Diisopropy			Methylene chloride		Dichlorometha	ine		
N,N-Dimethylformamide		DMF			2-Methylpentane		sohexane	-		
Dimethysufoxide		DMSO			1-Pentanol		-Amyl alcoho	1		
2-Ethoxyethanol		Cellosolve,	Ethyl glyc	nl	1-Propanol		ropyl alcohol			
Ethyl ether		Diethyl eth			Tetrahydrofuran		TOP YI LICONOI THF			
Ethyl acetate		EtOAc	,		Tetramethylene sulfone		ulfolane			
Ethyl benzene		Phenyletha	no		Xylene		Dimethylbenze	000		
Ethyl benzene		rnenyietha	ne		лутене	L	memyibenze	ine		

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: SA35791)

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PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 08/09/2023 0606 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)	-	0.0485	0.388	0.5	Kresoxim-methyl	_	0.0485	0.388	0.4	
Acephate (30560-19-1)	-	0.0485	0.388	0.4	(143390-89-0)					1
Acequinocyl (57960-19-7)	-	0.0485	0.388	2	Malathion (121-75-5)	-	0.0485	0.388	0.2	
Acetamiprid (135410-20-7)	-	0.0485	0.388	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.0485 0.0485	0.388 0.388	0.2 0.2	<u>Color Key</u>
Aldicarb (116-06-3)	-	0.0485	0.388	0.4	Methomyl (16752-77-5)	-	0.0485	0.388	0.4	RESULT < AL
Azoxystrobin (131860-33-8)	-	0.0485	0.388	0.2	Methyl parathion (298-0-0)	-	0.0485	0.388	0.2	RESULT > AL
Bifenazate (149877-41-8)	-	0.0485	0.388	0.2	MGK 264 (113-48-4)	-	0.0485	0.388	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3)	-	0.0485	0.388	0.2	Myclobutanil	_	0.0485	0.388	0.2	"-" not detected above
Boscalid (188425-85-6)	-	0.0485	0.388	0.4	(88671-89-0)					LOD
Carbaryl (63-25-2)	-	0.0485	0.388	0.2	Naled (300-76-5)	-	0.0485	0.388	0.5	Permethrins measured as the
Carbofuran (1563-66-2)	-	0.0485	0.388	0.2	Oxamyl (23135-22-0)	-	0.0485	0.388	1	cumulative residue of the <i>cis</i> - and
Chlorantraniliprole (800008-45-7)	-	0.0485	0.388	0.2	Paclobutrazol (76738-62-0)	-	0.0485	0.388	0.4	<i>trans</i> - permethrin isomers.
Chlorfenapyr	_	0.0485	0.388	1	Permethrins (52645-53-1)	-	0.0485	0.388	0.2	Pyrethrins measured as the cumulative residue of the
(122453-73-0)					Phosmet (732-11-6)	-	0.0485	0.388	0.2	pyrethrin I, cinerin I, and jasmolin
Chlorpyrifos (2921-88-2)	-	0.0485	0.388	0.2	Piperonyl butoxide (51-03-6)	-	0.0485	0.388	2	I isomers.
Clofentezine (74115-24-5)	-	0.0485	0.388	0.2	Prallethrins (2331-36-9)		0.0485	0.388	0.2	Action levels are referenced from
Cyfluthrin (68359-37-5)	-	0.0485	0.388	1	Propiconazole					the Drug and the
Cypermethrin (52315-07-8)	-	0.0485	0.388	1	(60207-90-1))	-	0.0485	0.388	0.4	State of Arkansas MMJ testing guidelines.
Daminozide (1596-84-5)	-	0.0485	0.388	1	Propoxur (114-26-1)	-	0.0485	0.388	0.2	A value of "-" for the action level
DDVP (62-73-7)	-	0.0485	0.388	0.1	Pyrethrins (8003-34-7)	-	0.0485	0.388	1	means that analyte is not
Diazinon (333-41-5)	-	0.0485	0.388	0.2	Pyridaben (96489-71-3)	-	0.0485 0.0485	0.388 0.388	0.2 0.2	currently regulated by the regulations referenced above.
Dimethoate (60-51-5)	-	0.0485	0.388	0.2	Spinosad (168316-95-8) Spiromesifen	-	0.0465	0.300	0.2	0
Ethoprophos (13194-48-4)	-	0.0485	0.388	0.2	(283594-90-1)	-	0.0485	0.388	0.2	<i>Disclaimer</i> : This information is provided as a service and makes
Etofenprox (80844-07-1)	-	0.0485	0.388	0.4	Spirotetramat		0.0485	0.388	0.2	no claims of efficacy and/or safety
Etoxazole (153233-91-1)	-	0.0485	0.388	0.2	(203313-25-1)	-	0.0465	0.300	0.2	of this product. Results are applicable only for the sample(s)
Fenoxycarb (72490-01-8)	-	0.0485	0.388	0.2	Spiroxamine	-	0.0485	0.388	0.4	analyzed and for the specific
(E)-Fenpyroximate (134098-61-6)	-	0.0485	0.388	0.4	(118134-30-8) Tebuconazole	_	0.0485	0.388	0.4	analysis conducted. This report is for informational purposes only
Fipronil (120068-37-3)	-	0.0485	0.388	0.4	(80443-41-0)					and should not be used to diagnose, treat, or prevent any
Flonicamid (158062-67-0)	-	0.0485	0.388	1	Thiacloprid (111988-49-9)	-	0.0485	0.388	0.2	medical-related symptoms. The
Fludioxinil (131341-86-1)		0.0485	0.388	0.4	Thiamethoxam					statements and results herein have
Hexythiazox (78587-05-0)		0.0485	0.388	1	(153719-23-4)	-	0.0485	0.388	0.2	not been approved and/or endorsed by the FDA.
Imazalil (35554-44-0)	-	0.0485	0.388	0.2	Trifloxystrobin		0.0485	0.388	0.2	
Imidacloprid (138261-41-3)	-	0.0485	0.388	0.4	(141517-21-7)		0.0403	0.500	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	Synonym(s)	Pesticide	<u>Synonym(s)</u>	Pesticide
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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Cultivar (Strain) or Sample Description: Log Cabin Sativa Hybrid Live Resin Cart			Date Completed: 08/10/2023

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

Analysis Date/Time: 08/10/2023 0745 (ICP/OES) Analysis Date/Time: - (DMA) Analyst: KF		Method: ICP/MS Instrument: Agilent 7500ce		Deviations from SOP: Dce 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)	-	57.8	91.5	200	
Cadmium (Cd)	-	57.8	91.5	200	
Lead (Pb)	-	57.8	91.5	500	
Mercury (Hg)	-	57.8	91.5	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





"DET" detected less than LOQ

"-" not detected above LOD

Action levels for heavy metals are referenced from the State of Arkansas MMJ testing guidelines.

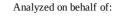
Disclaimer: This information is provided as a service and makes no claims of efficacy and/or safety of this product. Results are applicable only for the sample(s) analyzed and for the specific analysis conducted. This report is for informational purposes only and should not be used to diagnose, treat, 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: SA35791	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13245106176	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 08/07/2023
License: ADA 05_H273	License: 00065C	E20230731LCBLR01	Date Received: 08/08/2023
Cultivar (Strain) or Sample Description: Log Cabin Sativa Hybrid Live Resin Cart			Date Completed: 08/10/2023

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 08/10/20 Analyst: PW		Hardy Diagnostics Compac nt: Thermo Incubator	t Dry 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	-	and the second
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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