



SUMMARY OF ANALYSIS (SAMPLE ID: SA36239)

Testing Location:	Customer ID: 2168	Order ID: OR10656	Sample Type: Primary			
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible			
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13224603833	Mass: 10pcs			
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 10/16/2023			
License: ADH 113	License: 00065C	P20231011BLU19	Date Received: 10/16/2023			
Cultivar (Strain) or Sample Description: AR-Wana Gummies Blueberry IndicaDate Completed: 10/18/2023						

*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>	<u>mg</u> /g			
Δ9-Τ	ТНС	0.200	2.00			
CB	3D		-			
CBI	DA		-			
TOTAI	CBD	-	-			
TOTAI	L THC	0.200	2.00			
TOTAL CAN	NABINOIDS	0.200	2.00			
<u>Terpenes</u>	<u>s (Top 5)</u>	<u>(%)</u>	µg∕g			
α-Pir	nene	0.000943	9.43			
α-Bisa	bolol					
Camp	hene					
δ-3-Ca	arene					
β-Caryop	hyllene					
TOTAL TH	ERPENES	0.000943	9.43			
Contaminants	PASS/FAIL	Sample Picture	Upon Receipt			
Heavy Metals:	PASS					
Microbiology:	PASS					
Pesticides:	PASS					
Residual Solvents:	PASS					
		UDDA SOUR GUMMIES	Wana			



Scan the QR code to verify results.

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.

Kyle W. Felling, Ph.D. Laboratory Director

REPORT OF LABORATORY ANALYSIS

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Cultivar (Strain) or Sample	Date Completed: 10/18/2023		

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

Analysis Date/Time: 10/17/2023 1527 **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	0.00215	0.00502	-	-	-
CBCA	ND	ND	0.00668	0.0156	-	-	-
CBD	ND	ND	0.0152	0.0354	-	-	-
CBDA	ND	ND	0.00558	0.0130	-	-	-
CBDV	ND	ND	0.00243	0.00568	-	-	-
CBDVA	ND	ND	0.00649	0.0152	-	-	-
CBG	ND	ND	0.00985	0.0230	-	-	-
CBGA	ND	ND	0.0140	0.0165	-	-	-
CBL	ND	ND	0.0114	0.0266	-	-	-
CBN	ND	ND	0.00523	0.0122	-	-	-
CBNA	ND	ND	0.00564	0.0132	-	-	-
Δ9-ΤΗC	0.200	2.00	0.00626	0.0146	-	9.31	93.1
$\Delta 8$ -THC	ND	ND	0.00977	0.0228	-	-	-
THCA	ND	ND	0.00340	0.00794	-	-	-
THCV	ND	ND	0.00815	0.0190	-	-	-
THCVA	ND	ND	0.00260	0.00606	-	-	-
TOTAL	0.200	2.00		,	-	9.31	93.1
TOTAL CBC	-	-		,	-	-	-
TOTAL CBD	-	-			-	-	-
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	-	-			-	-	-
TOTAL CBN	-	-			-	-	-
TOTAL THC	0.200	2.00			-	9.31	93.1
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): 4.64 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.

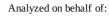
mg/g results are reported as mass component per mass material.

Kyle W. Felling, Ph.D. ory Directo











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License: ADH 113	License: 00065C	P20231011BLU19	Date Received: 10/16/2023
Cultivar (Strain) or Sample	Date Completed: 10/18/2023		

TERPENOID PROFILE

Analysis Date/Time:10, Analyst: KF	/17/2023 1615	Method: GC/MS Instrument: Agilent 7890/5975	Deviations from SOP: None		
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)			
α-Bisabolol	ND	-			
Camphene	<loq< td=""><td>-</td><td></td></loq<>	-			
δ-3-Carene	ND	-			
β-Caryophyllene	ND	-	Mr. F. M.		
Caryophyllene oxide	ND	-	SOUR GUMMIES BUERUMMES BUERUMMES		
p-Cymene	ND	-			
Eucalyptol	ND	-	AM NETWEI SAM (64)		
Geraniol	ND	-			
Guaiol	ND	-			
α-Humulene	ND	-	<i>Abbreviations:</i> GC - Gas Chromatography, MS - Mass		
Isopulegol	ND	-	Spectrometry, RL - Reporting Limit		
d-Limonene	ND	-	Abbreviations: ND - Not Detected, , LOD - Limit of Detection, LOQ - Limit of Quantitation		
Linalool	ND	-	This information is provided as a service		
β-Myrcene	ND	-	and makes no claims of efficacy and/or		
cis-Nerolidol	ND	-	safety of this product.		
trans-Nerolidol	ND	-	Results are applicable only for the sample(s) analyzed and for the specific		
α-Ocimene	ND	-	analysis conducted.		
β-Ocimene	ND	-	This report is for informational purposes only and should not be used to diagnose,		
α-Pinene	9.43	0.000943	treat, or prevent any		
β-Pinene	ND	-	medical-related symptoms.		
α-Terpinene	ND	-	The statements and results herein have not been approved and/or endorsed by		
γ-Terpinene	ND	-	the FDA.		
Terpinolene	ND	-			
TOTAL	9.43	0.000943	Reporting Limit (µg/g):		

"-" Not detected above LOD.













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Cultivar (Strain) or Sample	Date Completed: 10/18/2023		

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

Analysis Date/Time:	10/18/20)23 0139		Γ	Method: HS/GC/MS			D	eviation	s from SOP:
Analyst: KF				1	nstrument: Agilent 78	90/5975	5 None			
Solvent	<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)	-	88.1	176	5000	n-Heptane (142-82-5)	-	88.1	176	5000	
Acetonitrile (75-5-8)	-	88.1	176	410	n-Hexane (110-54-3)	-	30.8	61.7	290	C III
Benzene (71-43-2)	-	0.881	1.76	2	Isobutane (75-28-5)	-	88.1	176	5000	iuana M.
n-Butane (106-97-2)	-	88.1	176	5000	Isopropanol (67-63-0)	-	88.1	176	5000	SOUR GUMMIES
1-Butanol (71-36-3)	-	88.1	176	5000	Isopropyl acetate	_	88.1	176	5000	Sector Sector
2-Butanol (78-92-2)	-	88.1	176	5000	(108-21-4)					AM services
2-Butanone (78-93-3)	-	88.1	176	5000	Isopropyl benzene (98-82-8)	-	8.81	17.6	70	
Cyclohexane (110-82-7)	-	88.1	176	3880	Methanol (67-56-1)	213	88.1	176	3000	<u>Color Key</u>
1,2-Dimethoxyethane (110-71-4)	-	8.81	17.6	100	2-Methylbutane (78-78-4)	-	88.1	176	5000	RESULT < AL
(110-71-4) N,N-Dimethylacetamide (127-19-5)	-	88.1	176	1090	Methylene chloride (75-9-2)	-	88.1	176	600	RESULT > AL
(12, 10, 0) 2,2-Dimethylbutane (75-83-2)	-	30.8	61.7	290	2-Methylpentane (107-83-5)	-	30.8	61.7	290	"DET" detected less than LOQ
2,3-Dimethylbutane		20.0	o	200	3-Methylpentane (96-10-0)	-	30.8	61.7	290	"-" not detected above
(79-29-8)	-	30.8	61.7	290	n-Pentane (109-66-0)	-	88.1	176	5000	LOD
N,N-Dimethylformamide		88.1	176	880	1-Pentanol (71-41-0)	-	88.1	176	5000	"*" - o,m,p-Xylene and
(68-12-2)		00.1	170	000	n-Propane (74-98-6)	-	88.1	176	5000	Ethylbenzene
Dimethylsulfoxide (67-68-5)	-	88.1	176	5000	1-Propanol (71-23-8)	-	88.1	176	5000	Action levels are
1,4-Dioxane (123-91-1)	_	88.1	176	380	Pyridine (110-86-1)	-	30.8	61.7	200	referenced from the State of
Ethanol (64-17-5)	_	88.1	176	5000	Tetrahydrofuran (109-99-9)	-	88.1	176	720	Arkansas
2-Ethoxyethanol (110-80-5)	-	30.8	61.7	160	Tetramethylene sulfone (126-33-0)	-	30.8	61.7	160	MMJ testing
Ethyl ether (60-29-7)	-	88.1	176	5000	Toluene (108-88-3)	_	88.1	176	890	guidelines.
Ethyl acetate (141-78-6)	-	88.1	176	5000	o-Xylene (95-47-6)		88.1	176	2170	A value of "-"
Ethyl benzene (100-41-4)	-	88.1	176	2170	m,p-Xylene (108-38-3 or					for the action level
Ethylene glycol (107-21-1)	-	88.1	176	620	106-42-3)	-	88.1	176	2170	means that analyte
Ethylene oxide (75-21-8)	-	8.81	17.6	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170	is not currently regulated by the regulations referenced above.
Solvent		Synonym(s	5)		Solvent	<u>S</u>	Synonym(s)			
Acetonitrile		Methyl Cya	nide, ACN		Ethylene glycol	1	,2-Ethanediol			
1-Butanol		n-Butanol,	Butyl Alcol	hol	Isobutane	2	-Methylpropa	ine		
2-Butanol		sec-Butyl a	lcohol		Isopropanol	2	Propanol, IP	A		
2-Butanone		Methyl ethy	yl ketone, N	1EK	Isopropyl Acetate	A	Acetic acid iso	propyl este	r	
1,2-Dimethoxyethane		Monoglym	e		Methanol	Ν	Methyl alcoho	1		
2,3-Dimethylbutane		Neohexane			2-Methylbutane	I	sopentane			
2,3-Dimethylbutane		Diisopropy	'l		Methylene chloride	Γ	Dichlorometha	ine		
N,N-Dimethylformamide		DMF			2-Methylpentane	I	sohexane			
Dimethysufoxide		DMSO			1-Pentanol	n	-Amyl alcoho	ol		
2-Ethoxyethanol		Cellosolve,	Ethyl glyc	ol	1-Propanol	P	ropyl alcohol			
Ethyl ether		Diethyl eth	er, Ether		Tetrahydrofuran	Г	THF			
Ethyl acetate		EtOAc			Tetramethylene sulfone	S	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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Deviations from SOP:

None



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36239)

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Analysis Date/Time: 10/17/2023 1556 Analyst: KF

PESTICIDES PROFILE (SOP: SOP-PEST-001)

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

Instrument: Shimadzu LC-8050

<u>Pesticide</u>	<u>Result</u>	LOD	LOQ	<u>Action</u> Level	<u>Pesticide</u>	<u>Result</u>	LOD	LOQ	<u>Action</u> Level	
	<u>(µg/g)</u>	<u>(µg/g)</u>	<u>(µg/g)</u>	(µg/g)		<u>(µg/g)</u>	<u>(µg/g)</u>	(µg/g)	(µg/g)	
Abamectin (71751-41-2)	-	0.00944	0.0756	0.5	Kresoxim-methyl	-	0.00944	0.0756	0.4	
Acephate (30560-19-1)	-	0.00944	0.0756	0.4	(143390-89-0)		0.000.4.4	0.0756		
Acequinocyl (57960-19-7)	-	0.00944	0.0756	2	Malathion (121-75-5)	-	0.00944	0.0756	0.2	wana M
Acetamiprid	-	0.00944	0.0756	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.00944 0.00944	0.0756 0.0756	0.2 0.2	SOUR CUMMIES INTEREST
(135410-20-7) Aldicarb (116-06-3)		0.00944	0.0756	0.4	Methomyl (16752-77-5)	-	0.00944	0.0756	0.2	
Alucaib (110-00-3)	-				Methyl parathion	-				ASM verveissant
(131860-33-8)	-	0.00944	0.0756	0.2	(298-0-0)	-	0.00944	0.0756	0.2	
Bifenazate (149877-41-8)	-	0.00944	0.0756	0.2	MGK 264 (113-48-4)	-	0.00944	0.0756	0.2	<u>Color Key</u>
Bifenthrin (82657-04-3)	-	0.00944	0.0756	0.2	Myclobutanil	-	0.00944	0.0756	0.2	RESULT < AL
Boscalid (188425-85-6)	-	0.00944	0.0756	0.4	(88671-89-0) Naled (300-76-5)		0.00944	0.0756	0.5	RESULT > AL
Carbaryl (63-25-2)	-	0.00944	0.0756	0.2	Oxamyl (23135-22-0)	-	0.00944	0.0756	0.5	"DET" detected less than LOQ
Carbofuran (1563-66-2)	-	0.00944	0.0756	0.2	Paclobutrazol	-		0.0730	1	· ·
Chlorantraniliprole (800008-45-7)	-	0.00944	0.0756	0.2	(76738-62-0)	-	0.00944	0.0756	0.4	"-" not detected above LOD
Chlorfenapyr (122453-73-0)	-	0.00944	0.0756	1	Permethrins (52645-53-1)	-	0.00944	0.0756	0.2	Permethrins measured as the
Chlorpyrifos (2921-88-2)	-	0.00944	0.0756	0.2	Phosmet (732-11-6)	-	0.00944	0.0756	0.2	cumulative residue of the <i>cis</i> - and <i>trans</i> - permethrin isomers.
Clofentezine (74115-24-5)	-	0.00944	0.0756	0.2	Piperonyl butoxide	_	0.00944	0.0756	2	
Cyfluthrin (68359-37-5)	-	0.00944	0.0756	1	(51-03-6)				0.2	Pyrethrins measured as the cumulative residue of the
Cypermethrin		0.00944	0.0756	1	Prallethrins (2331-36-9)	-	0.00944	0.0756	0.2	pyrethrin I, cinerin I, and jasmolin
(52315-07-8)					Propiconazole (60207-90-1))	-	0.00944	0.0756	0.4	I isomers.
Daminozide (1596-84-5)	-	0.00944	0.0756	1	Propoxur (114-26-1)	-	0.00944	0.0756	0.2	Action levels are referenced from
DDVP (62-73-7)	-	0.00944	0.0756	0.1	Pyrethrins (8003-34-7)	-	0.00944	0.0756	1	the DBU is it is
Diazinon (333-41-5)	-	0.00944	0.0756	0.2	Pyridaben (96489-71-3)	-	0.00944	0.0756	0.2	State of Arkansas MMJ testing guidelines.
Dimethoate (60-51-5)	-	0.00944	0.0756	0.2	Spinosad (168316-95-8)	-	0.00944	0.0756	0.2	5
Ethoprophos (13194-48-4)	-	0.00944	0.0756	0.2	Spiromesifen (283594-90-1)	-	0.00944	0.0756	0.2	A value of "-" for the action level means that analyte is not
Etofenprox (80844-07-1)	-	0.00944	0.0756	0.4	Spirotetramat					currently regulated by the regulations referenced above.
Etoxazole (153233-91-1)	-	0.00944	0.0756	0.2	(203313-25-1)	-	0.00944	0.0756	0.2	0
Fenoxycarb (72490-01-8)	-	0.00944	0.0756	0.2	Spiroxamine	_	0.00944	0.0756	0.4	Disclaimer: This information is provided as a service and makes
(E)-Fenpyroximate (134098-61-6)	-	0.00944	0.0756	0.4	(118134-30-8) Tebuconazole					no claims of efficacy and/or safety of this product. Results are
Fipronil (120068-37-3)	-	0.00944	0.0756	0.4	(80443-41-0)	-	0.00944	0.0756	0.4	applicable only for the sample(s)
Flonicamid (158062-67-0)	-	0.00944	0.0756	1	Thiacloprid	_	0.00944	0.0756	0.2	analyzed and for the specific analysis conducted. This report is
Fludioxinil (131341-86-1)	-	0.00944	0.0756	0.4	(111988-49-9)					for informational purposes only
Hexythiazox (78587-05-0)	-	0.00944	0.0756	1	Thiamethoxam (153719-23-4)	-	0.00944	0.0756	0.2	and should not be used to
Imazalil (35554-44-0)	-	0.00944	0.0756	0.2	Trifloxystrobin		0.0004.5	0.0750	0.2	diagnose, treat, or prevent any medical-related symptoms. The
Imidacloprid (138261-41-3)	-	0.00944	0.0756	0.4	(141517-21-7)		0.00944	0.0756	0.2	statements and results herein have not been approved and/or endorsed by the FDA.

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	<u>Synonym(s)</u>	Pesticide	<u>Synonym(s)</u>	<u>Pesticide</u>	<u>Synonym(s)</u>
Cyfluthrin	Baythroid	Myclobutanil	Systhane	Propiconazole	Tilt
DDVP	Dichlorvos	Naled	Dibrom	Propoxur	Baygon
Ethoprophos	Prophos	Phosmet	Imidan		



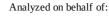
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Cultivar (Strain) or Sample	Date Completed: 10/18/2023		

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

Analysis Date/T Analysis Date/T Analyst: KF	ime: 10/17/2023 1 ime: - (DMA)	806 (ICP/OES)		thod: ICP/MS rument: 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)	-	56.7	89.7	200	
Cadmium (Cd)	-	56.7	89.7	200	
Lead (Pb)	-	56.7	89.7	500	
Mercury (Hg)	-	56.7	89.7	100	CONTROL CONTRO

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.

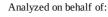
Disclaimer: This information is provided as a service and makes no claims of efficacy and/or safety of this product. Results are applicable only for the sample(s) analyzed and for the specific analysis conducted. This report is for informational purposes only and should not be used to diagnose, treat, 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: SA36239	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13224603833	Mass: 10pcs
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 10/16/2023
License: ADA 05_H273	License: 00065C	P20231011BLU19	Date Received: 10/16/2023
Cultivar (Strain) or Sample	Date Completed: 10/18/2023		

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 10/18/2023 0857 Analyst: PW		hod: Hardy Diagnostics CompactDry r ument: Thermo Incubator	Deviations from SOP: None
Bacteria/Microbe	<u>Result</u> (CFU/g)	Action Level (CFU/g)	
Aerobic Plate Count	Absent	-	
Coliforms, Total	Absent	1	
Escherichia Coli (E. Coli)	NT	100	
Mold/Yeast	NT		Mr. Chill
Pseudomonas aeruginosa	NT		
Salmonella spp.	NT	-	
Staphylococcus aureus	NT		AM nerver use year

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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