



### SUMMARY OF ANALYSIS (SAMPLE ID: SA35819)

Testing Location:	Customer ID: 2168	<b>Order ID:</b> OR10570	Sample Type: Primary		
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible		
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13228222330	Mass: 10pieces		
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 08/14/2023		
License: ADH 113	License: 00065C	P20230809MAN14	Date Received: 08/14/2023		
Cultivar (Strain) or Sample Description: AR-Wana Gummies Mango SativaDate Completed: 08/17/2023					

\*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.

Cannabinoi Δ9-Τ CB CB TOTAL TOTAL	HC G D CBD	(%) 0.196 0.00573 0.00518 0.00518 0.196	<u>mg/g</u> 1.96 0.0573 0.0518 0.0518 1.96	
TOTAL CAN		0.207	2.07	
<u>Terpenes</u> β-Caryop α-Bisa Camp δ-3-Ca Caryophyll TOTAL TE	hyllene bolol hene rene ene oxide	(%) µg/g 0.0154 154		
<u>Contaminants</u> Heavy Metals: Microbiology: Pesticides: Residual Solvents:	PASS/FAIL PASS PASS PASS PASS	Sample Picture	Upon Receipt	



Scan the QR code to verify results.

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REPORT OF LABORATORY ANALYSIS

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License: ADH 113	License: 00065C	P20230809MAN14	Date Received: 08/14/2023
Cultivar (Strain) or Sample D	Date Completed: 08/17/2023		

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

Analysis Date/Time: 8/15/2023 1345 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/ <u>mL)</u>	<u>Per</u> Serving (mg)	<u>Per</u> <u>Unit</u> (mg)
CBC	ND	ND	0.00209	0.00489	-	-	-
CBCA	ND	ND	0.00650	0.0152	-	-	-
CBD	0.00518	0.0518	0.0148	0.0345	-	0.247	2.47
CBDA	ND	ND	0.00544	0.0127	-	-	-
CBDV	ND	ND	0.00237	0.00553	-	-	-
CBDVA	ND	ND	0.00632	0.0147	-	-	-
CBG	0.00573	0.0573	0.00959	0.0224	-	0.273	2.73
CBGA	ND	ND	0.0136	0.0160	-	-	-
CBL	ND	ND	0.0111	0.0259	-	-	-
CBN	ND	ND	0.00509	0.0119	-	-	-
CBNA	ND	ND	0.00549	0.0128	-	-	-
Δ9-ΤΗC	0.196	1.96	0.00610	0.0142	-	9.34	93.4
$\Delta 8$ -THC	ND	ND	0.00951	0.0222	-	-	-
THCA	ND	ND	0.00331	0.00773	-	-	-
THCV	ND	ND	0.00793	0.0185	-	-	-
THCVA	ND	ND	0.00253	0.00590	-	-	-
TOTAL	0.207	2.07				9.86	98.6
TOTAL CBC	-	-				-	-
TOTAL CBD	0.00518	0.0518			-	0.247	2.47
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	0.00573	0.0573			-	0.273	2.73
TOTAL CBN	-	-			-	-	-
TOTAL THC	0.196	1.96			-	9.34	93.4
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): 4.77 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) +  $\Delta$ 9-THC Total THCV = (THCVA x 0.867) + THCV

Percentage results are reported by mass.

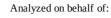
mg/g results are reported as mass component per mass material.

Kyle W. Felling, Ph.D. atory Directo











Testing Location:	Customer ID: 2168	Order ID: OR10570	Sample Type: Primary
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License: ADH 113	License: 00065C	P20230809MAN14	Date Received: 08/14/2023
Cultivar (Strain) or Sample	Date Completed: 08/17/2023		

**TERPENOID PROFILE** 

Analysis Date/Time:08/17/2023 0435 Analyst: KF		<b>Method:</b> GC/MS <b>Instrument:</b> Agilent 78	Deviations from SOP: 190/5975 None
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)	
α-Bisabolol	ND	-	
Camphene	ND	-	
δ-3-Carene	ND	-	
β-Caryophyllene	154	0.0154	Wana
Caryophyllene oxide	ND	-	SOUR GUMMES Inter
p-Cymene	ND	-	
Eucalyptol	ND	-	AM NETWELSON (COV
Geraniol	ND	-	
Guaiol	ND	-	
α-Humulene	ND	-	Abbreviations: 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 cis-Nerolidol	ND ND	-	and makes no claims of efficacy and/or safety of this product.
trans-Nerolidol	ND	-	Results are applicable only for the
α-Ocimene	ND	-	sample(s) analyzed and for the specific analysis conducted.
β-Ocimene	ND	-	This report is for informational purposes
α-Pinene	ND	-	only and should not be used to diagnose, treat, or prevent any
β-Pinene	ND	-	medical-related symptoms.
α-Terpinene	ND	-	The statements and results herein have
γ-Terpinene	ND	-	not been approved and/or endorsed by the FDA.
Terpinolene	ND	-	
TOTAL	154	0.0154	Reporting Limit (µg/g): 4

4.59

"-" Not detected above LOD.













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Cultivar (Strain) or Sample	Date Completed: 08/17/2023		

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

Analysis Date/Time:	Analysis Date/Time: 08/16/2023 1132 Method: HS/GC/MS				<b>Deviations from SOP:</b>					
Analyst: KF	Analyst: KF Instrument: Agilent 7890/5		90/5975	75 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)	LOD (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> <u>Level</u> (μg/g)	
Acetone (67-64-1)	-	124	249	5000	n-Heptane (142-82-5)	-	124	249	5000	
Acetonitrile (75-5-8)	-	124	249	410	n-Hexane (110-54-3)	-	43.5	87.1	290	
Benzene (71-43-2)	-	1.24	2.49	2	Isobutane (75-28-5)	-	124	249	5000	
n-Butane (106-97-2)	-	124	249	5000	Isopropanol (67-63-0)	-	124	249	5000	Uana sour gummies
1-Butanol (71-36-3)	-	124	249	5000	Isopropyl acetate		124	249	5000	And
2-Butanol (78-92-2)	-	124	249	5000	(108-21-4)		124	245	5000	AM server istanting
2-Butanone (78-93-3)	-	124	249	5000	Isopropyl benzene (98-82-8)	-	12.4	24.9	70	
Cyclohexane (110-82-7)	-	124	249	3880	. ,		124	249	3000	Color Key
1,2-Dimethoxyethane		12.4	24.9	100	Methanol (67-56-1)		124			
(110-71-4)	-	12.4	24.9	100	2-Methylbutane (78-78-4)	-	124	249	5000	<b>RESULT &lt; AL</b>
N,N-Dimethylacetamide (127-19-5)	-	124	249	1090	Methylene chloride (75-9-2)	-	124	249	600	<b>RESULT &gt; AL</b>
2,2-Dimethylbutane (75-83-2)	-	43.5	87.1	290	2-Methylpentane (107-83-5)	-	43.5	87.1	290	"DET" detected less than LOC
2,3-Dimethylbutane		43.5	87.1	290	3-Methylpentane (96-10-0)	-	43.5	87.1	290	"-" not detected above
(79-29-8)	-	43.5	07.1	290	n-Pentane (109-66-0)	-	124	249	5000	LOD
N,N-Dimethylformamide	-	124	249	880	1-Pentanol (71-41-0)	-	124	249	5000	"*" - o,m,p-Xylene and
(68-12-2)					n-Propane (74-98-6)	-	124	249	5000	Ethylbenzene
Dimethylsulfoxide (67-68-5)	-	124	249	5000	1-Propanol (71-23-8)	-	124	249	5000	
1,4-Dioxane (123-91-1)	_	124	249	380	Pyridine (110-86-1)	-	43.5	87.1	200	Action levels are referenced from the State of
Ethanol (64-17-5)	_	124	249	5000	Tetrahydrofuran (109-99-9)	-	124	249	720	Arkansas
2-Ethoxyethanol (110-80-5)	_	43.5	87.1	160	Tetramethylene sulfone	_	43.5	87.1	160	MMJ testing
Ethyl ether (60-29-7)	_	124	249	5000	(126-33-0)					guidelines.
Ethyl acetate (141-78-6)	_	124	249	5000	Toluene (108-88-3)	-	124	249	890	A value of "-"
Ethyl benzene (100-41-4)	_	124	249	2170	o-Xylene (95-47-6)	-	124	249	2170	for the action level
Ethylene glycol (107-21-1)	_	124	249	620	m,p-Xylene (108-38-3 or 106-42-3)	-	124	249	2170	means that analyte
Ethylene oxide (75-21-8)		124	24.9	50	Xylenes* (1330-20-7)		43.3	86.7	2170	is not currently
,		-		50		-	•	00.7	2170	regulated by the regulations referenced above.
<u>Solvent</u>		Synonym(s			<u>Solvent</u>		vnonym(s)			
Acetonitrile		Methyl Cya	,		Ethylene glycol	,	2-Ethanedio			
1-Butanol			Butyl Alco	hol	Isobutane		Methylpropa			
2-Butanol		sec-Butyl a			Isopropanol		Propanol, IP			
2-Butanone			yl ketone, N	1EK	Isopropyl Acetate		cetic acid iso		r	
1,2-Dimethoxyethane		Monoglym			Methanol		ethyl alcoho	1		
2,3-Dimethylbutane		Neohexane			2-Methylbutane		opentane			
2,3-Dimethylbutane		Diisopropy	1		Methylene chloride		ichlorometha	ane		
N,N-Dimethylformamide		DMF			2-Methylpentane		ohexane			
Dimethysufoxide		DMSO			1-Pentanol		Amyl alcoho			
2-Ethoxyethanol		Cellosolve	Ethyl glyc	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: SA35819)**

Testing Location:	Customer ID: 2168	Order ID: OR10570	Sample Type: Primary
Arkansas	<b>River Valley Relief MIPS</b>	Lot Number:	Matrix: Edible
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Greenbrier, AR 72058	Fort Smith, AR 72903	<b>Production Run:</b>	Date Collected: 08/14/2023
License: ADH 113	License: 00065C	P20230809MAN14	Date Received: 08/14/2023
Cultivar (Strain) or Sample	<b>Date Completed:</b> 08/17/2023		

### Analysis Date/Time: 08/15/2023 1356 Analyst: KF

### **PESTICIDES PROFILE (SOP: SOP-PEST-001)**

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

PesticideResultLODLOQActionPesticideResultLODLO(µg/g)(µg/g)(µg/g)(µg/g)(µg/g)(µg/g)(µg/g)(µg/g)(µg/g)		
Abamectin (71751-41-2) 0.00925 0.0740 0.5 Kresoxim-methyl 0.00925 0.07	740 0.4	
Acephate (30560-19-1) 0.00925 0.0740 0.4 (143390-89-0)		
Acequinocyl (57960-19-7) 0.00925 0.0740 2 Malathion (121-75-5) 0.00925 0.07		
Acetamiprid (125 (112 (112 (112 (112 (112 (112 (112		Wana Sour Gummes
(135410-20-7) 0.00925 0.0740 0.2 Methiocarb (2032-65-7) 0.00925 0.07925   Aldicarb (116-06-3) 0.00925 0.0740 0.4 Methomyl (16752-77-5) 0.00925 0.07		
Azovystehin Methyl parathion		AM server issued
(131860-33-8) 0.00925 0.0740 0.2 (298-0-0) 0.00925 0.07		Color Kay
Bifenazate (149877-41-8) 0.00925 0.0740 0.2 MGK 264 (113-48-4) 0.00925 0.07	740 0.2	<u>Color Key</u>
Bifenthrin (82657-04-3) 0.00925 0.0740 0.2 Myclobutanil 0.00925 0.07 Passelid (188425 85 6) 0.00025 0.0740 0.4 (88671-89-0) 0.00925 0.07	740 0.2	<b>RESULT &lt; AL</b>
Buscalla (100425-05-0) - 0.00925 0.0740 0.4 Noted (200.76.5) 0.00025 0.07	740 0.5	<b>RESULT &gt; AL</b>
Calibary (05-25-2) 0.00925 0.0740 0.2 Overwel (22125-22.0) 0.000925 0.07		"DET" detected less than LOQ
Calobilian (1565-66-2) - 0.00925 0.0740 0.2 Paclobuttazol		
(800008-45-7) 0.00925 0.0740 0.2 (76738-62-0) 0.00925 0.07	740 0.4	"-" not detected above LOD
Chlorfenapyr (122453-73-0) Permethrins 0.00925 Permethrins 0.0740 0.00925 0.07	740 0.2	Permethrins measured as the
Chlorpyrifos (2921-88-2) 0.00925 0.0740 0.2 Phosmet (732-11-6) 0.00925 0.07	740 0.2	cumulative residue of the <i>cis</i> - and <i>trans</i> - permethrin isomers.
Clofentezine (74115-24-5) 0.00925 0.0740 0.2 Piperonyl butoxide 0.00925 0.07	740 2	
Cyfluthrin (68359-37-5) 0.00925 0.0740 1 (51-03-6) 0.00925 0.0740 Prallethrins (2331-36-9) 0.00925 0.07	740 0.2	Pyrethrins measured as the cumulative residue of the
Cypermethrin - 0.00925 0.0740 1 Propiconazole 0.00925 0.07		pyrethrin I, cinerin I, and jasmolin I isomers.
(60207-90-1)) Daminozide (1596-84-5)	740 0.2	Action levels are referenced from
DDVP (62-73-7) 0.00925 0.0740 0.1 Propositi (114-26-1) 0.00925 0.07 Propositi (114-26-1) 0.00925 0.07		the
Diazinon (333-41-5) 0.00925 0.0740 0.2 Pyridaben (96489-71-3) 0.00925 0.07		State of Arkansas MMJ testing
Dimethoate (60-51-5) 0.00925 0.0740 0.2 Spinosad (168316-95-8) 0.00925 0.07		guidelines.
Ethoprophos 0.00925 0.0740 0.2 Spiromesifen 0.00925 0.07		A value of "-" for the action level means that analyte is not
(283594-90-1) (283594-90-1)		currently regulated by the
Etoxazole (153233-91-1) 0.00925 0.0740 0.2 (203313-25-1) 0.00925 0.07	740 0.2	regulations referenced above.
Fenoxycarb (72490-01-8) 0.00925 0.0740 0.2 Spiroxymine		Disclaimer: This information is
(E)-Fenpyroximate 0.00925 0.0740 0.4 (118134-30-8) 0.00925 0.07 (134098-61-6) 1.00925 0.0740 0.4 Tebuconazole 0.00925 0.07		provided as a service and makes no claims of efficacy and/or safety of this product. Results are
Fipronil (120068-37-3) 0.00925 0.0740 0.4 (80443-41-0) 0.00925 0.07	740 0.4	applicable only for the sample(s)
Flonicamid (158062-67-0) 0.00925 0.0740 1 Thiacloprid 0.00925 0.07	740 0.2	analyzed and for the specific
Fludioxinil (131341-86-1) - 0.00925 0.0740 0.4 (111988-49-9)	/40 0.2	analysis conducted. This report is for informational purposes only
Hexythiazox (78587-05-0) 0.00925 0.0740 1 Thiamethoxam 0.00925 0.07	740 0.2	and should not be used to
Imazalil (35554-44-0) - 0.00925 0.0740 0.2 Triflowystrobia		diagnose, treat, or prevent any medical-related symptoms. The
Imidacloprid (138261-41-3) 0.00925 0.0740 0.4 (141517-21-7) 0.00925 0.07	740 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 Ouantification

<u>Pesticide</u>	<u>Synonym(s)</u>	<u>Pesticide</u>	<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		



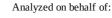
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### HEAVY METAL PROFILE (SOP: SOP-ICP-200.6)

Analysis Date/Time: 08/15/2023 1841 (ICP/OES) Analysis Date/Time: - (DMA) Analyst: KF		<b>Method:</b> ICP/MS <b>Instrument:</b> Agilent 7500ce		<b>Deviations from SOP:</b> 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	89.8	200	
Cadmium (Cd)	-	56.7	89.8	200	
Lead (Pb)	-	56.7	89.8	500	
Mercury (Hg)		56.7	89.8	100	WARDER STATES

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.

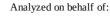
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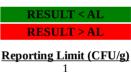
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Arkansas	<b>River Valley Relief MIPS</b>	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13228222330	Mass: 10pieces
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 08/14/2023
License: ADA 05_H273	License: 00065C	P20230809MAN14	Date Received: 08/14/2023
Cultivar (Strain) or Sample Description: AR-Wana Gummies Mango Sativa			<b>Date Completed:</b> 08/17/2023

### MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

<b>Analysis Date/Time:</b> 08/16/20 <b>Analyst:</b> PW		<b>d:</b> Hardy Diagnostics CompactDry <b>ment:</b> Thermo Incubator	<b>Deviations from SOP:</b> 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	-	Wana
Pseudomonas aeruginosa	NT	-	
Salmonella spp.	NT		UIANA SOUR GUMMIES NUM
Staphylococcus aureus	NT		AM HETWITTER

*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.

*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.



