







# **SUMMARY OF ANALYSIS (SAMPLE ID: SA36046)**

Testing Location:Customer ID: 2168Order ID: OR10613Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible232 S. Broadview St.5601 Old Greenwood Rd Suite CM00065C13224222797Mass: 10ea

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run:** Date Collected: 09/18/2023 License: ADH 113 License: 00065C P20230913BLU17 Date Received: 09/19/2023

 Cultivar (Strain) or Sample Description: AR-Wana Gummies Blueberry
 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/FAIL0.000Not TestedPASS

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>	ds (Top 3)	<u>(%)</u>	mg/g
СВ	D	0.000	0.000
CBI	DA .	0.000	0.000
CBI	OV	0.000	0.000
TOTAL	CBD	0.000	0.000
TOTAL	THC	0.000	0.000
TOTAL CAN	NABINOIDS	0.000	0.000
Terpenes	(Top 5)	(%)	µg/g
cis-Ner	olidol	0.0549	549
β-Myr	cene	0.00666	66.6
d-Limo	onene	0.000998	9.98
α-Bisa	bolol		
Camp	hene		
TOTAL TE	RPENES	0.0626	626
<u>Contaminants</u>	PASS/FAIL	Sample Picture Upon Receipt	

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





Scan the QR code to verify results.

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Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 09/18/2023 License: ADH 113 License: 00065C P20230913BLU17 **Date Received:** 09/19/2023 **Cultivar (Strain) or Sample Description:** AR-Wana Gummies Blueberry **Date Completed:** 09/20/2023

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

Analysis Date/Time: 09/19/2023 1204 Method: HPLC/DAD Moisture Content (%): 0.000
Analyst: PW Instrument: Agilent 1100 Water Activity (aw): -

<u>Cannabinoid</u>	Result (%)	Result (mg/g)	LOD (mg/g)	LOQ (mg/g)	Result (mg/ mL)	Per Serving (mg)	Per Unit (mg)
CBC	0.000	0.000	0.00228	0.00531	-	-	-
CBCA	0.000	0.000	0.00707	0.0165	-	-	-
CBD	0.000	0.000	0.0161	0.0375	-	-	-
CBDA	0.000	0.000	0.00591	0.0138	-	-	-
CBDV	0.000	0.000	0.00258	0.00601	-	-	-
CBDVA	0.000	0.000	0.00687	0.0160	-	-	-
CBG	0.000	0.000	0.0104	0.0243	-	0.240	2.40
CBGA	0.000	0.000	0.0148	0.0174	-	-	-
CBL	0.000	0.000	0.0120	0.0281	-	-	-
CBN	0.000	0.000	0.00553	0.0129	-	-	-
CBNA	0.000	0.000	0.00597	0.0139	-	-	-
Δ9-ΤΗС	0.000	0.000	0.00663	0.0155	-	9.55	95.5
Δ8-ΤΗС	0.000	0.000	0.0103	0.0241	-	-	-
THCA	0.000	0.000	0.00359	0.00841	-	-	-
THCV	0.000	0.000	0.00863	0.0201	-	-	-
THCVA	0.000	0.000	0.00276	0.00641	-	-	-
TOTAL	0.000	0.000				9.79	97.9
TOTAL CBC	0.000	0.000				_	
TOTAL CBD	0.000	0.000			-	-	-
TOTAL CBDV	0.000	0.000			-	-	-
TOTAL CBG	0.000	0.000			-	0.240	2.40
TOTAL CBN	0.000	0.000			-	-	-
TOTAL THC	0.000	0.000			-	9.55	95.5
TOTAL THCV	0.000	0.000			-	-	-

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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SERVING MASS (g): 4.73 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.













Order ID: OR10613 Sample Type: Primary **Testing Location:** Customer ID: 2168 Arkansas River Valley Relief MIPS Lot Number: Matrix: Edible 232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13224222797 Mass: 10ea

Fort Smith, AR 72903 Greenbrier, AR 72058 **Production Run: Date Collected:** 09/18/2023 P20230913BLU17 License: ADH 113 License: 00065C **Date Received:** 09/19/2023

Cultivar (Strain) or Sample Description: AR-Wana Gummies Blueberry **Date Completed:** 09/20/2023

#### TERPENOID PROFILE

Analysis Date/Time:09/20/2023 0729 Method: GC/MS **Deviations from SOP: Analyst:** KF Instrument: Agilent 7890/5975 None

<u>Terpene</u>	Result (µg/g)	Result (%)
α-Bisabolol	ND	<u>(70)</u> -
Camphene	<loq< td=""><td>-</td></loq<>	-
δ-3-Carene	ND	-
β-Caryophyllene	ND	-
Caryophyllene oxide	ND	-
p-Cymene	ND	-
Eucalyptol	ND	-
Geraniol	ND	-
Guaiol	ND	-
α-Humulene	ND	-
Isopulegol	ND	-
d-Limonene	9.98	0.000998
Linalool	ND	-
β-Myrcene	66.6	0.00666
cis-Nerolidol	549	0.0549
trans-Nerolidol	ND	-
α-Ocimene	ND	-
β-Ocimene	ND	-
α-Pinene	<loq< td=""><td>-</td></loq<>	-
β-Pinene	ND	-
α-Terpinene	ND	-
γ-Terpinene	ND	-
Terpinolene	ND	-
TOTAL	626	0.0626

Reporting Limit (µg/g): 8.32

"-" Not detected above LOD.













**Order ID:** OR10613 **Testing Location:** Customer ID: 2168 Sample Type: Primary Arkansas River Valley Relief MIPS Lot Number: Matrix: Edible

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13224222797 Mass: 10ea Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 09/18/2023

P20230913BLU17 License: ADH 113 License: 00065C **Date Received:** 09/19/2023 **Date Completed:** 09/20/2023 Cultivar (Strain) or Sample Description: AR-Wana Gummies Blueberry

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

**Analysis Date/Time:** 09/19/2023 1819 Method: HS/GC/MS **Deviations from SOP:** 

**Analyst: KF Instrument:** Agilent 7890/5975 None

Solvent	Result (µg/g)	<u>LOD</u> (µg/g)	LOQ (µg/g)	<u>Action</u> <u>Level</u> (μg/g)	Solvent	Result (μg/g)	<u>LOD</u> (µg/g)	LOQ (μg/g)	<u>Action</u> <u>Level</u> (μg/g)
Acetone (67-64-1)	-	134	268	5000	n-Heptane (142-82-5)	-	134	268	5000
Acetonitrile (75-5-8)	-	134	268	410	n-Hexane (110-54-3)	-	46.8	93.6	290
Benzene (71-43-2)	-	1.34	2.68	2	Isobutane (75-28-5)	-	134	268	5000
n-Butane (106-97-2)	-	134	268	5000	Isopropanol (67-63-0)	-	134	268	5000
1-Butanol (71-36-3)	-	134	268	5000	Isopropyl acetate	_	134	268	5000
2-Butanol (78-92-2)	-	134	268	5000	(108-21-4)		154	200	5000
2-Butanone (78-93-3)	-	134	268	5000	Isopropyl benzene (98-82-8)	-	13.4	26.8	70
Cyclohexane (110-82-7)	-	134	268	3880	Methanol (67-56-1)	_	134	268	3000
1,2-Dimethoxyethane (110-71-4)	-	13.4	26.8	100	2-Methylbutane (78-78-4)	-	134	268	5000
N,N-Dimethylacetamide (127-19-5)	-	134	268	1090	Methylene chloride (75-9-2)	-	134	268	600
2,2-Dimethylbutane (75-83-2)	-	46.8	93.6	290	2-Methylpentane (107-83-5)	-	46.8	93.6	290
2,3-Dimethylbutane		46.8	93.6	290	3-Methylpentane (96-10-0)	-	46.8	93.6	290
(79-29-8)	_	40.0	93.0	290	n-Pentane (109-66-0)	-	134	268	5000
N,N-Dimethylformamide	_	134	268	880	1-Pentanol (71-41-0)	-	134	268	5000
(68-12-2)					n-Propane (74-98-6)	-	134	268	5000
Dimethylsulfoxide (67-68-5)	-	134	268	5000	1-Propanol (71-23-8)	-	134	268	5000
1,4-Dioxane (123-91-1)	_	134	268	380	Pyridine (110-86-1)	-	46.8	93.6	200
Ethanol (64-17-5)	_	134	268	5000	Tetrahydrofuran (109-99-9)	-	134	268	720
2-Ethoxyethanol (110-80-5)	-	46.8	93.6	160	Tetramethylene sulfone (126-33-0)	-	46.8	93.6	160
Ethyl ether (60-29-7)	-	134	268	5000	Toluene (108-88-3)	_	134	268	890
Ethyl acetate (141-78-6)	-	134	268	5000	o-Xylene (95-47-6)	_	134	268	2170
Ethyl benzene (100-41-4)	-	134	268	2170	m,p-Xylene (108-38-3 or		_		
Ethylene glycol (107-21-1)	-	134	268	620	106-42-3)	-	134	268	2170
Ethylene oxide (75-21-8)	-	13.4	26.8	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170

<u>Solvent</u>	Synonym(s)	Solvent	Synonym(s)
Acetonitrile	Methyl Cyanide, ACN	Ethylene glycol	1,2-Ethanediol
1-Butanol	n-Butanol, Butyl Alcohol	Isobutane	2-Methylpropane
2-Butanol	sec-Butyl alcohol	Isopropanol	2-Propanol, IPA
2-Butanone	Methyl ethyl ketone, MEK	Isopropyl Acetate	Acetic acid isopropyl ester
1,2-Dimethoxyethane	Monoglyme	Methanol	Methyl alcohol
2,3-Dimethylbutane	Neohexane	2-Methylbutane	Isopentane
2,3-Dimethylbutane	Diisopropyl	Methylene chloride	Dichloromethane
N,N-Dimethylformamide	DMF	2-Methylpentane	Isohexane
Dimethysufoxide	DMSO	1-Pentanol	n-Amyl alcohol
2-Ethoxyethanol	Cellosolve, Ethyl glycol	1-Propanol	Propyl alcohol
Ethyl ether	Diethyl ether, Ether	Tetrahydrofuran	THF
Ethyl acetate	EtOAc	Tetramethylene sulfone	Sulfolane
Ethyl benzene	Phenylethane	Xylene	Dimethylbenzene



**Color Key** 

"DET" detected less than LOQ

"-" not detected above LOD

"\*" - o,m,p-Xylene and Ethylbenzene

Action levels 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.

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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232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13224222797 Mass: 10ea
Greenbrier, AR 72058 Fort Smith, AR 72903 Production Run: Date Collected: 09/18/2023

License: ADH 113 License: 00065C P20230913BLU17 **Date Received:** 09/19/2023 **Cultivar (Strain) or Sample Description:** AR-Wana Gummies Blueberry **Date Completed:** 09/20/2023

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

Analysis Date/Time: 09/19/2023 1236 Method: LC/MS/MS Deviations from SOP:

Analyst: KF Instrument: Shimadzu LC-8050 None

<u>Pesticide</u>	Result (μg/g)	<u>LOD</u> (μg/g)	<u>LOQ</u> (µg/g)	Action Level (µg/g)	<u>Pesticide</u>	Result (μg/g)	<u>LOD</u> (µg/g)	LOQ (µg/g)	<u>Action</u> <u>Level</u> (μg/g)
Abamectin (71751-41-2)	-	0.00998	0.0799	0.5	Kresoxim-methyl	_	0.00998	0.0799	0.4
Acephate (30560-19-1)	-	0.00998	0.0799	0.4	(143390-89-0)				
Acequinocyl (57960-19-7)	-	0.00998	0.0799	2	Malathion (121-75-5)	-	0.00998	0.0799	0.2
Acetamiprid (135410-20-7)	-	0.00998	0.0799	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.00998 0.00998	0.0799 0.0799	0.2 0.2
Aldicarb (116-06-3)	-	0.00998	0.0799	0.4	Methomyl (16752-77-5)	-	0.00998	0.0799	0.4
Azoxystrobin (131860-33-8)	-	0.00998	0.0799	0.2	Methyl parathion (298-0-0)	-	0.00998	0.0799	0.2
Bifenazate (149877-41-8)	-	0.00998	0.0799	0.2	MGK 264 (113-48-4)	_	0.00998	0.0799	0.2
Bifenthrin (82657-04-3)	-	0.00998	0.0799	0.2	Myclobutanil		0.00998	0.0799	0.2
Boscalid (188425-85-6)	-	0.00998	0.0799	0.4	(88671-89-0)	_	0.00998	0.0799	0.2
Carbaryl (63-25-2)	-	0.00998	0.0799	0.2	Naled (300-76-5)	-	0.00998	0.0799	0.5
Carbofuran (1563-66-2)	-	0.00998	0.0799	0.2	Oxamyl (23135-22-0)	-	0.00998	0.0799	1
Chlorantraniliprole (800008-45-7)	-	0.00998	0.0799	0.2	Paclobutrazol (76738-62-0)	-	0.00998	0.0799	0.4
Chlorfenapyr (122453-73-0)	-	0.00998	0.0799	1	Permethrins (52645-53-1)	-	0.00998	0.0799	0.2
Chlorpyrifos (2921-88-2)	-	0.00998	0.0799	0.2	Phosmet (732-11-6)	-	0.00998	0.0799	0.2
Clofentezine (74115-24-5)	-	0.00998	0.0799	0.2	Piperonyl butoxide	_	0.00998	0.0799	2
Cyfluthrin (68359-37-5)	-	0.00998	0.0799	1	(51-03-6) Prallethrins (2331-36-9)		0.00998	0.0799	0.2
Cypermethrin (52315-07-8)	-	0.00998	0.0799	1	Propiconazole	-	0.00998	0.0799	0.2
Daminozide (1596-84-5)	-	0.00998	0.0799	1	(60207-90-1)) Propoxur (114-26-1)		0.00998	0.0799	0.2
DDVP (62-73-7)	-	0.00998	0.0799	0.1	Pyrethrins (8003-34-7)		0.00998	0.0799	1
Diazinon (333-41-5)	-	0.00998	0.0799	0.2	Pyridaben (96489-71-3)	-	0.00998	0.0799	0.2
Dimethoate (60-51-5)	-	0.00998	0.0799	0.2	Spinosad (168316-95-8)		0.00998	0.0799	0.2
Ethoprophos (13194-48-4)	-	0.00998	0.0799	0.2	Spiromesifen (283594-90-1)		0.00998	0.0799	0.2
Etofenprox (80844-07-1)	-	0.00998	0.0799	0.4	` /				
Etoxazole (153233-91-1)	-	0.00998	0.0799	0.2	Spirotetramat (203313-25-1)	-	0.00998	0.0799	0.2
Fenoxycarb (72490-01-8)	-	0.00998	0.0799	0.2	Spiroxamine		0.00000	0.0700	0.4
(E)-Fenpyroximate (134098-61-6)	-	0.00998	0.0799	0.4	(118134-30-8) Tebuconazole	-	0.00998	0.0799	0.4
Fipronil (120068-37-3)	-	0.00998	0.0799	0.4	(80443-41-0)	-	0.00998	0.0799	0.4
Flonicamid (158062-67-0)	-	0.00998	0.0799	1	Thiacloprid	_	0.00998	0.0799	0.2
Fludioxinil (131341-86-1)	-	0.00998	0.0799	0.4	(111988-49-9)		0.00000	0.07.55	0.2
Hexythiazox (78587-05-0)	-	0.00998	0.0799	1	Thiamethoxam (153719-23-4)	-	0.00998	0.0799	0.2
Imazalil (35554-44-0)	-	0.00998	0.0799	0.2	Trifloxystrobin				
Imidacloprid (138261-41-3)	-	0.00998	0.0799	0.4	(141517-21-7)	-	0.00998	0.0799	0.2



Color Key

# RESULT > AL

"DET" detected less than LOQ

"-" not detected above LOD

Permethrins measured as the cumulative residue of the *cis*- and *trans*- permethrin isomers.

Pyrethrins measured as the cumulative residue of the pyrethrin I, cinerin I, and jasmolin I isomers.

Action levels are referenced from the

State of Arkansas MMJ testing guidelines.

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 $Abbreviations: LC-Liquid\ Chromatography, MS-Mass\ Spectrometry, RL-Reporting\ Limit, AL-Action\ Level, CAS-Chemical\ Abstract\ Services, LOD-Limit\ of\ Detection, LOQ-Limit\ of\ Quantification$ 

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













**Date Received:** 09/19/2023

# **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36046)**

Testing Location:Customer ID: 2168Order ID: OR10613Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13224222797 Mass: 10ea
Greenbrier, AR 72058 Fort Smith, AR 72903 Production Run: Date Collected: 09/18/2023

Cultivar (Strain) or Sample Description: AR-Wana Gummies Blueberry

Date Completed: 09/20/2023

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

P20230913BLU17

Analysis Date/Time: 09/19/2023 1836 (ICP/OES) Method: ICP/MS Deviations from SOP:

Analysis Date/Time: - (DMA) Instrument: Agilent 7500ce None

**Analyst:** KF

License: ADH 113

<u>Heavy Metal</u>	<u>Result</u> (μg/kg)	<u>LOD</u> (µg/kg)	LOQ (µg/kg)	<u>Action Level</u> (μg/kg)
Arsenic (As)	-	56.0	88.7	200
Cadmium (Cd)	-	56.0	88.7	200
Lead (Pb)	-	56.0	88.7	500
Mercury (Hg)	-	56.0	88.7	100

License: 00065C



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

### Color Key

RESULT < AL
RESULT > AL

"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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License: ADA 05 H273







**Date Received:** 09/19/2023

### **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36046)**

Testing Location:Customer ID: 2168Sample ID: SA36046Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13224222797 Mass: 10ea
Greenbrier, AR 72058 Fort Smith, AR 72903 Production Run: Date Collected: 09/18/2023

Cultivar (Strain) or Sample Description: AR-Wana Gummies Blueberry

Date Completed: 09/20/2023

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

P20230913BLU17

**Analysis Date/Time:** 9/20/2023 0821 **Method:** Hardy Diagnostics CompactDry **Deviations from SOP:** 

Analyst: PW Instrument: Thermo Incubator None

Bacteria/Microbe	<u>Result</u> (CFU/g)	Action Level (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	-

License: 00065C



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

Color Key

RESULT < AL RESULT > AL

Reporting Limit (CFU/g)

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