







SUMMARY OF ANALYSIS (SAMPLE ID: SA36184)

Order ID: OR10642 **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 M00065C13229911297 Mass: 10pcs Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 10/09/2023 P20231003MAN16 License: ADH 113 License: 00065C **Date Received:** 10/09/2023

Cultivar (Strain) or Sample Description: AR-Wana Gummies Mango Sativa **Date Completed:** 10/11/2023

*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

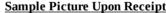
> **Moisture Content (%)** PASS/FAIL Water Activity (aw) **PASS** Not Tested Not Tested

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	ds (Top 3)	<u>(%)</u>	mg/g
Δ9-Τ	HC	0.192	1.92
CB	D	0.0159	0.159
CB	G	0.00636	0.0636
TOTAL	CBD	0.0159	0.159
TOTAL	THC	0.192	1.92
TOTAL CANN	NABINOIDS	0.214	2.14
Terpenes	(Top 5)	<u>(%)</u>	µg/g
α-Pin	ene	0.000945	9.45
α-Bisa	polol		
Campl	nene		
δ-3-Ca	rene		
β-Caryop	hyllene		
TOTAL TE	RPENES	0.000945	9.45
<u>Contaminants</u>	Contaminants PASS/FAIL		e Upon Receipt

Contaminants	PASS/FAIL
Heavy Metals:	PASS
Microbiology:	PASS
Pesticides:	PASS
Residual Solvents	DASS







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.

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The statements and results herein have not been approved and/or endorsed by the FDA.







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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36184)

Testing Location:Customer ID: 2168Order ID: OR10642ArkansasRiver Valley Relief MIPSLot Number:232 S. Broadview St.5601 Old Greenwood Rd Suite CM00065C13229911297Greenbrier, AR 72058Fort Smith, AR 72903Production Run:

Fort Smith, AR 72903 **Production Run:**License: 00065C P20231003MAN16

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

Date Completed: 10/11/2023

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

Analysis Date/Time: 10/10/2023 1815 **Method:** HPLC/DAD **Analyst:** PW **Instrument:** Agilent 1100

Cannabinoid Result Per Per Result LOD LOQ Result **Serving Unit** (mg/ (%) (mg/g)(mg/g)(mg/g)<u>mL)</u> <u>(mg)</u> (mg) **CBC** ND ND 0.00215 0.00503 **CBCA** ND ND 0.00669 0.0156 0.159 **CBD** 0.0159 0.0152 0.0355 0.770 7.70 **CBDA** ND ND 0.00560 0.0131 **CBDV** ND ND 0.00244 0.00569 **CBDVA** ND ND 0.00650 0.0152 3.08 **CBG** 0.00636 0.0636 0.00987 0.0230 0.308 **CBGA** ND ND 0.0140 0.0165 CBL ND ND 0.0114 0.0266 **CBN** ND ND 0.00524 0.0122 ND ND 0.00565 **CBNA** 0.0132 Δ9-ΤΗС 0.192 1.92 0.00628 0.0146 9.30 93.0 Δ8-ΤΗС ND ND 0.00979 0.0229 ND ND **THCA** 0.00340 0.00796 THCV ND ND 0.00817 0.0190 **THCVA** ND ND 0.00261 0.00607 TOTAL 0.214 2.14 10.4 104 **TOTAL CBC TOTAL CBD** 0.0159 0.159 0.770 7.70 TOTAL CBDV **TOTAL CBG** 0.00636 0.0636 0.308 3.08 **TOTAL CBN** 0.192 TOTAL THC 1.92 9.30 93.0 **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): -



Sample Type: Primary

Date Collected: 10/09/2023

Date Received: 10/09/2023

Matrix: Edible

Mass: 10pcs

SERVING MASS (g): 4.84 SERVINGS/UNIT: 10

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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10/11/2023

[&]quot;-" Not detected above LOD.









CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36184)

Customer ID: 2168 Order ID: OR10642 Sample Type: Primary **Testing Location:** Arkansas River Valley Relief MIPS Lot Number: Matrix: Edible 232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13229911297 Mass: 10pcs

Greenbrier, AR 72058 Fort Smith, AR 72903 **Date Collected:** 10/09/2023 **Production Run:** License: 00065C P20231003MAN16 **Date Received:** 10/09/2023 License: ADH 113

Cultivar (Strain) or Sample Description: AR-Wana Gummies Mango Sativa **Date Completed:** 10/11/2023

TERPENOID PROFILE

Analysis Date/Time:10/10/2023 2031 Method: GC/MS **Deviations from SOP:**

Analyst: KF Instrument: Agilent 7890/5975 None

<u>Terpene</u>	Result	Result	
α-Bisabolol	<u>(µg/g)</u> ND	<u>(%)</u> -	
Camphene	ND	_	
δ-3-Carene	ND	_	
β-Caryophyllene	<loq< td=""><td>_</td><td></td></loq<>	_	
Caryophyllene oxide	ND	-	SOUR GUMMIES SOUR GUMMIES
p-Cymene	ND	-	Parker and the second s
Eucalyptol	ND	-	AM NEW TISMICAL
Geraniol	ND	-	
Guaiol	ND	-	
α-Humulene	<loq< td=""><td>-</td><td>Abbreviations: GC - Gas Chromatography, MS - Mass</td></loq<>	-	Abbreviations: GC - Gas Chromatography, MS - Mass
Isopulegol	ND	-	Spectrometry, RL - Reporting Limit
d-Limonene	<loq< td=""><td>-</td><td>Abbreviations: ND - Not Detected, , LOD - Limit of Detection, LOQ - Limit</td></loq<>	-	Abbreviations: ND - Not Detected, , LOD - Limit of Detection, LOQ - Limit
Linalool	ND	-	of Quantitation
β-Myrcene	ND	-	This information is provided as a service 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
α-Pinene	9.45	0.000945	only and should not be used to diagnose, treat, or prevent any
β-Pinene	<loq< td=""><td>-</td><td>medical-related symptoms.</td></loq<>	-	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	9.45	0.000945	Reporting Limit (μg/g):

Reporting Limit (µg/g): 7.88

"-" Not detected above LOD.







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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36184)

Testing Location:Customer ID: 2168Order ID: OR10642Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13229911297 Mass: 10pcs
Greenbrier, AR 72058 Fort Smith, AR 72903 Production Run: Date Collected: 10/09/2023

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

Date Completed: 10/11/2023

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

P20231003MAN16

Analysis Date/Time: 10/10/2023 0333 Method: HS/GC/MS Deviations from SOP:

Analyst: KF Instrument: Agilent 7890/5975 None

License: 00065C

Solvent	Result (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> <u>Level</u> (μg/g)	Solvent	Result (μg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	Action Level (µg/g)
Acetone (67-64-1)	-	129	258	5000	n-Heptane (142-82-5)	-	129	258	5000
Acetonitrile (75-5-8)	-	129	258	410	n-Hexane (110-54-3)	-	45.1	90.2	290
Benzene (71-43-2)	-	1.29	2.58	2	Isobutane (75-28-5)	-	129	258	5000
n-Butane (106-97-2)	-	129	258	5000	Isopropanol (67-63-0)	-	129	258	5000
1-Butanol (71-36-3)	-	129	258	5000	Isopropyl acetate	_	129	258	5000
2-Butanol (78-92-2)	-	129	258	5000	(108-21-4)		120	200	5000
2-Butanone (78-93-3)	-	129	258	5000	Isopropyl benzene (98-82-8)	-	12.9	25.8	70
Cyclohexane (110-82-7)	-	129	258	3880	Methanol (67-56-1)	-	129	258	3000
1,2-Dimethoxyethane (110-71-4)	-	12.9	25.8	100	2-Methylbutane (78-78-4)	-	129	258	5000
N,N-Dimethylacetamide (127-19-5)	-	129	258	1090	Methylene chloride (75-9-2)	-	129	258	600
2,2-Dimethylbutane (75-83-2)	-	45.1	90.2	290	2-Methylpentane (107-83-5)	-	45.1	90.2	290
2,3-Dimethylbutane (79-29-8)	-	45.1	90.2	290	3-Methylpentane (96-10-0) n-Pentane (109-66-0)	-	45.1 129	90.2 258	290 5000
N,N-Dimethylformamide (68-12-2)	-	129	258	880	1-Pentanol (71-41-0) n-Propane (74-98-6)	-	129 129	258 258	5000 5000
Dimethylsulfoxide (67-68-5)	-	129	258	5000	1-Propanol (71-23-8)	-	129	258	5000
1,4-Dioxane (123-91-1)	_	129	258	380	Pyridine (110-86-1)	-	45.1	90.2	200
Ethanol (64-17-5)	_	129	258	5000	Tetrahydrofuran (109-99-9)	-	129	258	720
2-Ethoxyethanol (110-80-5)	-	45.1	90.2	160	Tetramethylene sulfone (126-33-0)	-	45.1	90.2	160
Ethyl ether (60-29-7)	-	129	258	5000	Toluene (108-88-3)	_	129	258	890
Ethyl acetate (141-78-6)	-	129	258	5000	o-Xylene (95-47-6)	_	129	258	2170
Ethyl benzene (100-41-4)	-	129	258	2170	m,p-Xylene (108-38-3 or				
Ethylene glycol (107-21-1)	-	129	258	620	106-42-3)	-	129	258	2170
Ethylene oxide (75-21-8)	-	12.9	25.8	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



Date Received: 10/09/2023

Color Key

RESULT < AL RESULT > AL

"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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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36184)

Testing Location:Customer ID: 2168Order ID: OR10642Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13229911297 Mass: 10pcs
Greenbrier, AR 72058 Fort Smith, AR 72903 Production Run: Date Collected: 10/09/2023

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

Date Completed: 10/11/2023

PESTICIDES PROFILE (SOP: SOP-PEST-001)

P20231003MAN16

Analysis Date/Time: 10/10/2023 1722 Method: LC/MS/MS Deviations from SOP:

Analyst: KF Instrument: Shimadzu LC-8050 None

License: 00065C

<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)	-	0.00940	0.0752	0.5	Kresoxim-methyl	_	0.00940	0.0752	0.4
Acephate (30560-19-1)	-	0.00940	0.0752	0.4	(143390-89-0)				
Acequinocyl (57960-19-7)	-	0.00940	0.0752	2	Malathion (121-75-5)	-	0.00940	0.0752	0.2
Acetamiprid (135410-20-7)	-	0.00940	0.0752	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.00940 0.00940	0.0752 0.0752	0.2 0.2
Aldicarb (116-06-3)	-	0.00940	0.0752	0.4	Methomyl (16752-77-5)	-	0.00940	0.0752	0.4
Azoxystrobin (131860-33-8)	-	0.00940	0.0752	0.2	Methyl parathion (298-0-0)	-	0.00940	0.0752	0.2
Bifenazate (149877-41-8)	-	0.00940	0.0752	0.2	MGK 264 (113-48-4)	-	0.00940	0.0752	0.2
Bifenthrin (82657-04-3)	-	0.00940	0.0752	0.2	Myclobutanil		0.00940	0.0752	0.2
Boscalid (188425-85-6)	-	0.00940	0.0752	0.4	(88671-89-0)		0.00540	0.0732	
Carbaryl (63-25-2)	-	0.00940	0.0752	0.2	Naled (300-76-5)	-	0.00940	0.0752	0.5
Carbofuran (1563-66-2)	-	0.00940	0.0752	0.2	Oxamyl (23135-22-0)	-	0.00940	0.0752	1
Chlorantraniliprole (800008-45-7)	-	0.00940	0.0752	0.2	Paclobutrazol (76738-62-0)	-	0.00940	0.0752	0.4
Chlorfenapyr (122453-73-0)	-	0.00940	0.0752	1	Permethrins (52645-53-1)	-	0.00940	0.0752	0.2
Chlorpyrifos (2921-88-2)	-	0.00940	0.0752	0.2	Phosmet (732-11-6)	-	0.00940	0.0752	0.2
Clofentezine (74115-24-5)	-	0.00940	0.0752	0.2	Piperonyl butoxide	_	0.00940	0.0752	2
Cyfluthrin (68359-37-5)	-	0.00940	0.0752	1	(51-03-6) Prallethrins (2331-36-9)		0.00940	0.0752	0.2
Cypermethrin (52315-07-8)	-	0.00940	0.0752	1	Propiconazole (60207-90-1))	-	0.00940	0.0752	0.4
Daminozide (1596-84-5)	-	0.00940	0.0752	1	Propoxur (114-26-1)		0.00940	0.0752	0.2
DDVP (62-73-7)	-	0.00940	0.0752	0.1	Pyrethrins (8003-34-7)		0.00940	0.0752	1
Diazinon (333-41-5)	-	0.00940	0.0752	0.2	Pyridaben (96489-71-3)	-	0.00940	0.0752	0.2
Dimethoate (60-51-5)	-	0.00940	0.0752	0.2	Spinosad (168316-95-8)		0.00940	0.0752	0.2
Ethoprophos (13194-48-4)	-	0.00940	0.0752	0.2	Spiromesifen		0.00940	0.0752	0.2
Etofenprox (80844-07-1)	-	0.00940	0.0752	0.4	(283594-90-1)				
Etoxazole (153233-91-1)	-	0.00940	0.0752	0.2	Spirotetramat (203313-25-1)	-	0.00940	0.0752	0.2
Fenoxycarb (72490-01-8)	-	0.00940	0.0752	0.2	Spiroxamine		0.000.40	0.0750	0.4
(E)-Fenpyroximate (134098-61-6)	-	0.00940	0.0752	0.4	(118134-30-8) Tebuconazole	-	0.00940	0.0752	0.4
Fipronil (120068-37-3)	-	0.00940	0.0752	0.4	(80443-41-0)	-	0.00940	0.0752	0.4
Flonicamid (158062-67-0)	-	0.00940	0.0752	1	Thiacloprid	_	0.00940	0.0752	0.2
Fludioxinil (131341-86-1)	-	0.00940	0.0752	0.4	(111988-49-9)		0.000.0	0.07.52	0.2
Hexythiazox (78587-05-0)	-	0.00940	0.0752	1	Thiamethoxam (153719-23-4)	-	0.00940	0.0752	0.2
Imazalil (35554-44-0)	-	0.00940	0.0752	0.2	Trifloxystrobin				
Imidacloprid (138261-41-3)	-	0.00940	0.0752	0.4	(141517-21-7)	-	0.00940	0.0752	0.2



Color Key

RESULT < AL

Date Received: 10/09/2023

"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.

A value of "-" for the action level means that analyte is not currently regulated by the regulations referenced above.

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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















CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36184)

 Testing Location:
 Customer ID: 2168
 Order ID: OR10642
 Sample Type: Primary

 Arkansas
 River Valley Relief MIPS
 Lot Number:
 Matrix: Edible

 232 S. Broadview St.
 5601 Old Greenwood Rd Suite C
 M00065C13229911297
 Mass: 10pcs

 Greenbrier, AR 72058
 Fort Smith, AR 72903
 Production Run:
 Date Collected: 10/09/2023

 License: ADH 113
 License: 00065C
 P20231003MAN16
 Date Received: 10/09/2023

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

Date Completed: 10/11/2023

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

Analysis Date/Time: 10/10/2023 1814 (ICP/OES) Method: ICP/MS Deviations from SOP:

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

Analyst: KF

<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)	-	58.5	92.7	200
Cadmium (Cd)	-	58.5	92.7	200
Lead (Pb)	-	58.5	92.7	500
Mercury (Hg)	-	58.5	92.7	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

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: 10/09/2023

CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36184)

Testing Location:Customer ID: 2168Sample ID: SA36184Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13229911297 Mass: 10pcs
Greenbrier, AR 72058 Fort Smith, AR 72903 Production Run: Date Collected: 10/09/2023

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

Date Completed: 10/11/2023

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

P20231003MAN16

Analysis Date/Time: 10/10/2023 1355 **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)

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.



