

SUMMARY OF ANALYSIS (SAMPLE ID: SA43683)

Testing Location:	Customer ID: 2168	Order ID: OR12098	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13247396652	Mass: 4ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 04/09/2026
License: ADH 113	License: 00065C	E20260408RVARKS01	Date Received: 04/09/2026
Cultivar (Strain) or Sample Description: Red Velvet Full Spectrum AR 1g Cart			Date Completed: 04/12/2026

*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.

*Where provided, statements of conformity (e.g. Pass/Fail) are made in accordance with ILAC G8, Binary Statement for Simple Acceptance Rule (w=0, AL=TL).
PASS: when the result is within the acceptance interval. FAIL: when the result is outside the acceptance interval

Moisture Content (%)

Not Tested

Water Activity (aw)

Not Tested

PASS/FAIL

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>Cannabinoids (Top 3)</u>	<u>(%)</u>	<u>mg/g</u>
Δ9-THC	65.1	651
THCA	23.6	236
CBGA	1.47	14.7
TOTAL CBD	0.000	0.000
TOTAL THC	85.8	858
TOTAL CANNABINOIDS	92.3	923

<u>Terpenes (Top 5)</u>	<u>(%)</u>	<u>ppm</u>
β-Caryophyllene	1.91	19100
d-Limonene	1.74	17400
β-Myrcene	1.35	13500
α-Humulene	0.567	5670
trans-β-Farnesene	0.305	3050
TOTAL TERPENES	7.23	72300

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

Sample Picture Upon Receipt



Scan the QR code to verify results.

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Kyle W. Felling
Kyle W. Felling, Ph.D.
Laboratory Director



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA43683)

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License: ADH 113	License: 00065C	E0260408RVARKS01	Date Received: 04/09/2026

Cultivar (Strain) or Sample Description: Red Velvet Full Spectrum AR 1g Cart **Date Completed:** 04/12/2026

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

Analysis Date/Time: 04/10/2026 1351

Method: HPLC/DAD

Moisture Content (%): -

Analyst: PW

Instrument: Agilent 1100

Water Activity (aw): -

<u>Cannabinoid</u>	<u>Result (%)</u>	<u>Result (mg/g)</u>	<u>LOD (mg/g)</u>	<u>LOQ (mg/g)</u>	<u>Result (mg/mL)</u>	<u>Per Serving (mg)</u>	<u>Per Unit (mg)</u>
CBC	0.120	1.20	0.100	0.234	-	1.20	1.20
CBCA	1.02	10.2	0.312	0.727	-	10.2	10.2
CBD	ND	ND	0.708	1.65	-	0.000	0.000
CBDV	DET	DET	0.261	0.608	-	0.000	0.000
CBDVA	ND	ND	0.114	0.265	-	0.000	0.000
CBDVA	ND	ND	0.303	0.707	-	0.000	0.000
CBG	0.363	3.63	0.459	1.07	-	3.63	3.63
CBGA	1.47	14.7	0.651	0.768	-	14.7	14.7
CBL	ND	ND	0.531	1.24	-	0.000	0.000
CBN	ND	ND	0.244	0.569	-	0.000	0.000
CBNA	ND	ND	0.263	0.614	-	0.000	0.000
Δ9-THC	65.1	651	0.292	0.681	-	651	651
Δ8-THC	ND	ND	0.456	1.06	-	0.000	0.000
THCA	23.6	236	0.158	0.371	-	236	236
THCV	0.468	4.68	0.380	0.886	-	4.68	4.68
THCVA	0.145	1.45	0.121	0.283	-	1.45	1.45
TOTAL	92.3	923			-	923	923
TOTAL CBC	1.01	10.1			-	10.1	10.1
TOTAL CBD	0.000	0.000			-	0.000	0.000
TOTAL CBDV	0.000	0.000			-	0.000	0.000
TOTAL CBG	1.65	16.5			-	16.5	16.5
TOTAL CBN	0.000	0.000			-	0.000	0.000
TOTAL THC	85.8	858			-	858	858
TOTAL THCV	0.594	5.94			-	5.94	5.94



SERVING MASS (g): 1.00
SERVINGS/UNIT: 1

"-" Not reported for this sample.

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.
Dry percent = Wet percent / (1 - (Moisture Content / 100))

Total CBC = (CBCA x 0.877) + CBC
Total CBD = (CBDV 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) + Δ9-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.

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, ND - Not Detected (less than LOD)

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Laboratory Director



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License: ADH 113	License: 00065C	E0260408RVARKS01	Date Received: 04/09/2026

Cultivar (Strain) or Sample Description: Red Velvet Full Spectrum AR 1g Cart **Date Completed:** 04/12/2026

TERPENOID PROFILE

Analysis Date/Time: 04/10/2026 1623

Method: GC/MS

Deviations from SOP:

Analyst: KF

Instrument: Agilent 7890/5975

None

<u>Terpene</u>	<u>Result (ppm)</u>	<u>Result (%)</u>	<u>Terpene</u>	<u>Result (ppm)</u>	<u>Result (%)</u>
α-Bisabolol	1050	0.105	d-Limonene	17400	1.74
L-Borneol	ND	-	Linalool	1420	0.142
D-Borneol	138	0.0138	l-Menthone	ND	-
Camphene	650	0.0650	Menthol	138	0.0138
Camphor	ND	-	β-Myrcene	13500	1.35
δ-3-Carene	ND	-	Nerol	ND	-
Carvacrol	ND	-	cis-Nerolidol	ND	-
Carvone	ND	-	trans-Nerolidol	386	0.0386
β-Caryophyllene	19100	1.91	Nootkatone	ND	-
Caryophyllene oxide	155	0.0155	cis-β-Ocimene	161	0.0161
Cedrene	103	0.0103	trans-β-Ocimene	866	0.0866
Cedrol	55.6	0.00556	Octyl acetate	ND	-
cis-Citral	ND	-	α-Phellandrene	ND	-
trans-Citral	ND	-	Phytane	ND	-
Citronellol	66.3	0.00663	α-Pinene	2670	0.267
o-Cymene	ND	-	β-Pinene	1340	0.134
m-Cymene	ND	-	Piperitone	ND	-
p-Cymene	ND	-	Pulegone	ND	-
Eucalyptol	ND	-	Sabinene	809	0.0809
trans-β-Farnesene	3050	0.305	Sabinene hydrate	ND	-
Farnesol 1	ND	-	Safranal	ND	-
Farnesol 2	173	0.0173	Squalene	ND	-
Fenchone	239	0.0239	Terpinen-4-ol	ND	-
Fenchyl alcohol	953	0.0953	α-Terpinene	ND	-
Geraniol	ND	-	γ-Terpinene	ND	-
Geranyl acetate	1230	0.123	α-Terpineol	104	0.0104
Guaiol	80.2	0.00802	Terpinolene	250	0.0250
α-Humulene	5670	0.567	α-Thujone	ND	-
Isoborneol	ND	-	Thymol	ND	-
Isobornyl acetate	ND	-	Valencene	523	0.0523
d-Isomenthone	ND	-	d-Valerolactam	ND	-
Isopulegol	ND	-	Verbenone	ND	-



Abbreviations: GC - Gas Chromatography, MS - Mass Spectrometry, RL - Reporting Limit

Abbreviations: ND - Not Detected, LOD - Limit of Detection, LOQ - Limit of Quantitation

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"-" Not detected above LOD.

TOTAL (ppm) 72300 (%) 7.23 **Reporting Limit (ppm):** 44.0

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License: ADH 113	License: 00065C	E0260408RVARKS01	Date Received: 04/09/2026

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

Analysis Date/Time: 04/10/2026 1447	Method: HS/GC/MS	Deviations from SOP:
Analyst: KF	Instrument: Agilent 7890/5975	None

Solvent	Result (ppm)	LOD (ppm)	LOQ (ppm)	Action Level (ppm)	Solvent	Result (ppm)	LOD (ppm)	LOQ (ppm)	Action Level (ppm)
Acetone (67-64-1)	-	35.0	69.9	5000	n-Heptane (142-82-5)	-	35.0	69.9	5000
Acetonitrile (75-5-8)	-	35.0	69.9	410	n-Hexane (110-54-3)	-	12.2	24.5	290
Benzene (71-43-2)	-	0.350	0.699	2	Isobutane (75-28-5)	-	35.0	69.9	5000
n-Butane (106-97-2)	-	35.0	69.9	5000	Isopropanol (67-63-0)	-	35.0	69.9	5000
1-Butanol (71-36-3)	-	35.0	69.9	5000	Isopropyl acetate (108-21-4)	-	35.0	69.9	5000
2-Butanol (78-92-2)	-	35.0	69.9	5000	Isopropyl benzene (98-82-8)	-	3.50	6.99	70
2-Butanone (78-93-3)	-	35.0	69.9	5000	Methanol (67-56-1)	-	35.0	69.9	3000
Cyclohexane (110-82-7)	-	35.0	69.9	3880	2-Methylbutane (78-78-4)	-	35.0	69.9	5000
1,2-Dimethoxyethane (110-71-4)	-	3.50	6.99	100	Methylene chloride (75-9-2)	-	35.0	69.9	600
N,N-Dimethylacetamide (127-19-5)	-	35.0	69.9	1090	2-Methylpentane (107-83-5)	-	12.2	24.5	290
2,2-Dimethylbutane (75-83-2)	-	12.2	24.5	290	3-Methylpentane (96-10-0)	-	12.2	24.5	290
2,3-Dimethylbutane (79-29-8)	-	12.2	24.5	290	n-Pentane (109-66-0)	-	35.0	69.9	5000
N,N-Dimethylformamide (68-12-2)	-	35.0	69.9	880	1-Pentanol (71-41-0)	-	35.0	69.9	5000
Dimethylsulfoxide (67-68-5)	-	35.0	69.9	5000	n-Propane (74-98-6)	-	35.0	69.9	5000
1,4-Dioxane (123-91-1)	-	35.0	69.9	380	1-Propanol (71-23-8)	-	35.0	69.9	5000
Ethanol (64-17-5)	-	35.0	69.9	5000	Pyridine (110-86-1)	-	12.2	24.5	200
2-Ethoxyethanol (110-80-5)	-	12.2	24.5	160	Tetrahydrofuran (109-99-9)	-	35.0	69.9	720
Ethyl ether (60-29-7)	-	35.0	69.9	5000	Tetramethylene sulfone (126-33-0)	-	12.2	24.5	160
Ethyl acetate (141-78-6)	-	35.0	69.9	5000	Toluene (108-88-3)	-	35.0	69.9	890
Ethyl benzene (100-41-4)	-	35.0	69.9	2170	o-Xylene (95-47-6)	-	35.0	69.9	2170
Ethylene glycol (107-21-1)	-	35.0	69.9	620	m,p-Xylene (108-38-3 or 106-42-3)	-	35.0	69.9	2170
Ethylene oxide (75-21-8)	-	3.50	6.99	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170



Color Key

RESULT < AL

RESULT > AL

"-" not detected above 1/2 Action Level

"*" - 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.

Solvent	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
Dimethylsulfoxide	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

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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PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 04/10/2026 1321

Method: LC/MS/MS

Deviations from SOP:

Analyst: KF

Instrument: Shimadzu LC-8050

None

Pesticide	Result (ppm)	LOD (ppm)	LOQ (ppm)	Action Level (ppm)	Pesticide	Result (ppm)	LOD (ppm)	LOQ (ppm)	Action Level (ppm)
Abamectin (71751-41-2)	-	0.0439	0.351	0.5	Kresoxim-methyl (143390-89-0)	-	0.0439	0.351	0.4
Acephate (30560-19-1)	-	0.0439	0.351	0.4	Malathion (121-75-5)	-	0.0439	0.351	0.2
Acetaminocyl (57960-19-7)	-	0.0439	0.351	2	Metalaxyl (57837-19-1)	-	0.0439	0.351	0.2
Acetamiprid (135410-20-7)	-	0.0439	0.351	0.2	Methiocarb (2032-65-7)	-	0.0439	0.351	0.2
Aldicarb (116-06-3)	-	0.0439	0.351	0.4	Methomyl (16752-77-5)	-	0.0439	0.351	0.4
Azoxystrobin (131860-33-8)	-	0.0439	0.351	0.2	Methyl parathion (298-0-0)	-	0.0439	0.351	0.2
Bifenazate (149877-41-8)	-	0.0439	0.351	0.2	MGK 264 (113-48-4)	-	0.0439	0.351	0.2
Bifenthrin (82657-04-3)	-	0.0439	0.351	0.2	Myclobutanil (88671-89-0)	-	0.0439	0.351	0.2
Boscalid (188425-85-6)	-	0.0439	0.351	0.4	Naled (300-76-5)	-	0.0439	0.351	0.5
Carbaryl (63-25-2)	-	0.0439	0.351	0.2	Oxamyl (23135-22-0)	-	0.0439	0.351	1
Carbofuran (1563-66-2)	-	0.0439	0.351	0.2	Paclobutrazol (76738-62-0)	-	0.0439	0.351	0.4
Chlorantraniliprole (800008-45-7)	-	0.0439	0.351	0.2	Permethrins (52645-53-1)	-	0.0439	0.351	0.2
Chlorfenapyr (122453-73-0)	-	0.0439	0.351	1	Phosmet (732-11-6)	-	0.0439	0.351	0.2
Chlorpyrifos (2921-88-2)	-	0.0439	0.351	0.2	Piperonyl butoxide (51-03-6)	-	0.0439	0.351	2
Clofentezine (74115-24-5)	-	0.0439	0.351	0.2	Prallethrin (2331-36-9)	-	0.0439	0.351	0.2
Cyfluthrin (68359-37-5)	-	0.0439	0.351	1	Propiconazole (60207-90-1)	-	0.0439	0.351	0.4
Cypermethrin (52315-07-8)	-	0.0439	0.351	1	Propoxur (114-26-1)	-	0.0439	0.351	0.2
Daminozide (1596-84-5)	-	0.0439	0.351	1	Pyrethrins (8003-34-7)	-	0.0439	0.351	1
DDVP (62-73-7)	-	0.0439	0.351	0.1	Pyridaben (96489-71-3)	-	0.0439	0.351	0.2
Diazinon (333-41-5)	-	0.0439	0.351	0.2	Spinosad (168316-95-8)	-	0.0439	0.351	0.2
Dimethoate (60-51-5)	-	0.0439	0.351	0.2	Spiromesifen (283594-90-1)	-	0.0439	0.351	0.2
Ethoprophos (13194-48-4)	-	0.0439	0.351	0.2	Spirotetramat (203313-25-1)	-	0.0439	0.351	0.2
Etofenprox (80844-07-1)	-	0.0439	0.351	0.4	Spiroxamine (118134-30-8)	-	0.0439	0.351	0.4
Etoxazole (153233-91-1)	-	0.0439	0.351	0.2	Tebuconazole (80443-41-0)	-	0.0439	0.351	0.4
Fenoxycarb (72490-01-8)	-	0.0439	0.351	0.2	Thiacloprid (111988-49-9)	-	0.0439	0.351	0.2
(E)-Fenpyroximate (134098-61-6)	-	0.0439	0.351	0.4	Thiamethoxam (153719-23-4)	-	0.0439	0.351	0.2
Fipronil (120068-37-3)	-	0.0439	0.351	0.4	Trifloxystrobin (141517-21-7)	-	0.0439	0.351	0.2
Flonicamid (158062-67-0)	-	0.0439	0.351	1					
Fludioxinil (131341-86-1)	-	0.0439	0.351	0.4					
Hexythiazox (78587-05-0)	-	0.0439	0.351	1					
Imazalil (35554-44-0)	-	0.0439	0.351	0.2					
Imidacloprid (138261-41-3)	-	0.0439	0.351	0.4					



Color Key

RESULT < AL

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.

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

REPORT OF LABORATORY ANALYSIS

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Kyle W. Felling, Ph.D.
Laboratory Director



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA43683)

Testing Location:	Customer ID: 2168	Order ID: OR12098	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13247396652	Mass: 4ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 04/09/2026
License: ADH 113	License: 00065C	E20260408RVARKS01	Date Received: 04/09/2026
Cultivar (Strain) or Sample Description: Red Velvet Full Spectrum AR 1g Cart			Date Completed: 04/12/2026

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

Analysis Date/Time: 04/11/2026 1221

Method: ICP/MS

Deviations from SOP:

Instrument: Agilent 7700x

None

Analyst: KF

Heavy Metal	Result (ppb)	LOD (ppb)	LOQ (ppb)	Action Level (ppb)
Arsenic (As)	ND	57.4	90.9	200
Cadmium (Cd)	ND	57.4	90.9	200
Lead (Pb)	ND	57.4	90.9	500
Mercury (Hg)	ND	57.4	90.9	100



Abbreviations: ICP - Inductively-Coupled Plasma, MS - Mass Spectrometry,
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

"ND" not detected above LOD

Reporting limit is 1/2 the action level of each analyte.

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.

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Kyle W. Felling
Kyle W. Felling, Ph.D.
Laboratory Director



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA43683)

Testing Location:	Customer ID: 2168	Sample ID: SA43683	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13247396652	Mass: 4ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 04/09/2026
-	License: 00065C	E20260408RVARKS01	Date Received: 04/09/2026
Cultivar (Strain) or Sample Description: Red Velvet Full Spectrum AR 1g Cart			Date Completed: 04/12/2026

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 04/11/2026 1543	Method: Hardy Diagnostics CompactDry	Deviations from SOP:
Analyst: PW	Instrument: Thermo Incubator	None

Bacteria/Microbe	Result (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	-



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)

1

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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