

License: ADH 113







Date Received: 01/14/2025

SUMMARY OF ANALYSIS (SAMPLE ID: SA39847)

Testing Location:Customer ID: 2168Order ID: OR11375Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13220923443 Mass: 1ea
Greenbrier, AR 72058 Fort Smith, AR 72903 Production Run: Date Collected: 01/13/2025

Cultivar (Strain) or Sample Description: ARV-Sea Salt Caramel Soft Chews 200mg 20pk

Date Completed:01/15/2025

*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/FAILNot TestedNot TestedPASS

P20250106CAR01

Moisture content/water activity action levels are referenced from the State of Arkansas MMJ testing guidelines.

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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>	oids (Top 3)	(%)	<u>mg</u> /g
Δ9-7	ГНС	0.223	2.23
Δ8-	ГНС	0.0154	0.154
Cl	BL	0.00993	0.0993
TOTA	L CBD	0.000	0.000
TOTA	LTHC	0.223	2.23
TOTAL CAN	NABINOIDS	0.250	2.50
<u>Terpene</u>	s (Top 5)	<u>(%)</u>	µg/g
β-Му	rcene	0.00308	30.8
Lina	lool	0.00226	22.6
α-Pi	nene	0.00215	21.5
Cam	bhene	0.00212	21.2
d-Lim	onene	0.00138	13.8
TOTAL T	ERPENES	0.0137	137
Contaminants	PASS/FAIL	Sample Dicture	e Unon Receipt

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







Scan the QR code to verify results.

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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA39847)

Testing Location:Customer ID: 2168Order ID: OR11375Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible232 S. Broadview St5601 Old Greenwood Rd Suite CM00065C13220923443Mass: 1ea

 232 S. Broadview St.
 5601 Old Greenwood Rd Suite C
 M00065C13220923443
 Mass: 1ea

 Greenbrier, AR 72058
 Fort Smith, AR 72903
 Production Run:
 Date Collected: 01/13/2025

 License: ADH 113
 License: 00065C
 P20250106CAR01
 Date Received: 01/14/2025

Cultivar (Strain) or Sample Description: ARV-Sea Salt Caramel Soft Chews 200mg 20pk

Date Completed: 01/15/2025

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

Analysis Date/Time: 01/14/2025 1437 Method: HPLC/DAD
Analyst: PW Instrument: Agilent 1100

<u>Cannabinoid</u>	Result (%)	Result (mg/g)	LOD (mg/g)	LOQ (mg/g)	Result (mg/ mL)	Per Serving (mg)	Per Unit (mg)
CBC	ND	ND	0.00214	0.00500	-	0.000	0.000
CBCA	ND	ND	0.00666	0.0155	-	0.000	0.000
CBD	ND	ND	0.0151	0.0353	-	0.000	0.000
CBDA	ND	ND	0.00557	0.0130	-	0.000	0.000
CBDV	ND	ND	0.00243	0.00566	-	0.000	0.000
CBDVA	ND	ND	0.00647	0.0151	-	0.000	0.000
CBG	ND	ND	0.00982	0.0229	-	0.000	0.000
CBGA	ND	ND	0.0139	0.0164	-	0.000	0.000
CBL	0.00993	0.0993	0.0113	0.0265	-	0.417	8.34
CBN	0.00136	0.0136	0.00521	0.0122	-	0.0571	1.14
CBNA	ND	ND	0.00562	0.0131	-	0.000	0.000
CBT	ND	ND	0.00816	0.0190	-	0.000	0.000
Δ9-ΤΗС	0.223	2.23	0.00625	0.0146	-	9.37	187
Δ8-ΤΗС	0.0154	0.154	0.00974	0.0227	-	0.647	12.9
THCA	ND	ND	0.00339	0.00792	-	0.000	0.000
THCV	ND	ND	0.00813	0.0189	-	0.000	0.000
THCVA	ND	ND	0.00260	0.00604	-	0.000	0.000
TOTAL	0.250	2.50	,			10.5	210
TOTAL CBC	0.000	0.000				0.000	0.000
TOTAL CBD	0.000	0.000			-	0.000	0.000
TOTAL CBDV	0.000	0.000			-	0.000	0.000
TOTAL CBG	0.000	0.000			-	0.000	0.000
TOTAL CBN	0.00136	0.0136			-	0.0571	1.14
TOTAL THC	0.223	2.23			-	9.37	187
TOTAL THCV	0.000	0.000			-	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, ND - Not Detected (less than LOD)

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Moisture Content (%): -Water Activity (aw): -



SERVING MASS (g): 4.20 SERVINGS/UNIT: 20

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.





[&]quot;-" Not reported for this sample.



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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA39847)

Customer ID: 2168 **Testing Location:** Arkansas River Valley Relief MIPS 232 S. Broadview St. 5601 Old Greenwood Rd Suite C Fort Smith, AR 72903 Greenbrier, AR 72058

License: 00065C

Cultivar (Strain) or Sample Description: ARV-Sea Salt Caramel Soft Chews 200mg 20pk

Order ID: OR11375 Lot Number: M00065C13220923443 **Production Run:**

P20250106CAR01

Matrix: Edible Mass: 1ea **Date Collected:** 01/13/2025

Sample Type: Primary

Date Received: 01/14/2025 **Date Completed:**01/15/2025

TERPENOID PROFILE

Analysis Date/Time:01/14/2025 2011 Method: GC/MS **Deviations from SOP: Analyst: KF** Instrument: Agilent 7890/5975 None

<u>Terpene</u>	<u>Result</u> (μg/g)	Result (%)	
α-Bisabolol	ND	<u>(70)</u> -	
Camphene	21.2	0.00212	
δ-3-Carene	ND	-	
3-Caryophyllene	ND	-	
Caryophyllene oxide	ND	-	
p-Cymene	11.2	0.00112	
Eucalyptol	7.80	0.000780	
Geraniol	ND	-	
Guaiol	ND	-	Albert interes CC Core
x-Humulene	ND	-	Abbreviations: GC - Gas Chromatography, MS - Mass
sopulegol	ND	-	Spectrometry, RL - Reporting Limit
d-Limonene	13.8	0.00138	Abbreviations: ND - Not Detected, LOD - Limit of Detection, LOQ - L
Linalool	22.6	0.00226	of Quantitation
β-Myrcene	30.8	0.00308	This information is provided as a se and makes no claims of efficacy and
cis-Nerolidol	ND	-	safety of this product.
rans-Nerolidol	ND	-	Results are applicable only for the sample(s) analyzed and for the spec
x-Ocimene	ND	-	analysis conducted.
3-Ocimene	ND	-	This report is for informational purp
x-Pinene	21.5	0.00215	only and should not be used to diag treat, or prevent any
3-Pinene	ND	-	medical-related symptoms.
x-Terpinene	ND	-	The statements and results herein ha not been approved and/or endorsed
γ-Terpinene	ND	-	the FDA.
Terpinolene	7.53	0.000753	
TOTAL	137	0.0137	Reporting Limit

Reporting Limit (µg/g): 4.70

"-" Not detected above LOD.













CERTIFICATE OF ANALYSIS (SAMPLE ID: SA39847)

Testing Location:Customer ID: 2168Order ID: OR11375Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13220923443 **Mass:** 1ea

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run:** Date Collected: 01/13/2025 License: ADH 113 License: 00065C P20250106CAR01 Date Received: 01/14/2025

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

Analysis Date/Time: 01/14/2025 2022 Method: HS/GC/MS Deviations from SOP:

Analyst: KF Instrument: Agilent 7890/5975 None

Cultivar (Strain) or Sample Description: ARV-Sea Salt Caramel Soft Chews 200mg 20pk

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)	-	39.1	78.1	5000	n-Heptane (142-82-5)	-	39.1	78.1	5000
Acetonitrile (75-5-8)	-	39.1	78.1	410	n-Hexane (110-54-3)	-	13.7	27.3	290
Benzene (71-43-2)	-	0.391	0.781	2	Isobutane (75-28-5)	-	39.1	78.1	5000
n-Butane (106-97-2)	-	39.1	78.1	5000	Isopropanol (67-63-0)	-	39.1	78.1	5000
1-Butanol (71-36-3)	-	39.1	78.1	5000	Isopropyl acetate	_	39.1	78.1	5000
2-Butanol (78-92-2)	-	39.1	78.1	5000	(108-21-4)		55.1	70.1	5000
2-Butanone (78-93-3)	-	39.1	78.1	5000	Isopropyl benzene (98-82-8)	-	3.91	7.81	70
Cyclohexane (110-82-7)	-	39.1	78.1	3880	Methanol (67-56-1)	_	39.1	78.1	3000
1,2-Dimethoxyethane (110-71-4)	-	3.91	7.81	100	2-Methylbutane (78-78-4)	-	39.1	78.1	5000
N,N-Dimethylacetamide (127-19-5)	-	39.1	78.1	1090	Methylene chloride (75-9-2)	-	39.1	78.1	600
2,2-Dimethylbutane (75-83-2)	-	13.7	27.3	290	2-Methylpentane (107-83-5)	-	13.7	27.3	290
2,3-Dimethylbutane		13.7	27.3	290	3-Methylpentane (96-10-0)	-	13.7	27.3	290
(79-29-8)	-	13./	27.3	290	n-Pentane (109-66-0)	-	39.1	78.1	5000
N,N-Dimethylformamide	_	39.1	78.1	880	1-Pentanol (71-41-0)	-	39.1	78.1	5000
(68-12-2)					n-Propane (74-98-6)	-	39.1	78.1	5000
Dimethylsulfoxide (67-68-5)	-	39.1	78.1	5000	1-Propanol (71-23-8)	-	39.1	78.1	5000
1,4-Dioxane (123-91-1)	_	39.1	78.1	380	Pyridine (110-86-1)	-	13.7	27.3	200
Ethanol (64-17-5)	_	39.1	78.1	5000	Tetrahydrofuran (109-99-9)	-	39.1	78.1	720
2-Ethoxyethanol (110-80-5)	-	13.7	27.3	160	Tetramethylene sulfone (126-33-0)	-	13.7	27.3	160
Ethyl ether (60-29-7)	-	39.1	78.1	5000	Toluene (108-88-3)	_	39.1	78.1	890
Ethyl acetate (141-78-6)	-	39.1	78.1	5000	o-Xylene (95-47-6)	_	39.1	78.1	2170
Ethyl benzene (100-41-4)	-	39.1	78.1	2170	m,p-Xylene (108-38-3 or				
Ethylene glycol (107-21-1)	-	39.1	78.1	620	106-42-3)	-	39.1	78.1	2170
Ethylene oxide (75-21-8)	-	3.91	7.81	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170



Date Completed:01/15/2025

Color Key

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.

<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

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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Analysis Date/Time: 01/14/2025 1813







CERTIFICATE OF ANALYSIS (SAMPLE ID: SA39847)

Testing Location:Customer ID: 2168Order ID: OR11375Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13220923443 Mass: 1ea

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run:** Date Collected: 01/13/2025 License: ADH 113 License: 00065C P20250106CAR01 Date Received: 01/14/2025

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Cultivar (Strain) or Sample Description: ARV-Sea Salt Caramel Soft Chews 200mg 20pk

Method: LC/MS/MS Deviations from SOP:

Analyst: KF Instrument: Shimadzu LC-8050 None

Docticido				Action	Desti si de				Action
<u>Pesticide</u>	<u>Result</u> (μg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)	<u>Pesticide</u>	<u>Result</u> (μg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
Abamectin (71751-41-2)	-	0.00943	0.0754	0.5	Kresoxim-methyl		0.00943	0.0754	0.4
Acephate (30560-19-1)	-	0.00943	0.0754	0.4	(143390-89-0)				
Acequinocyl (57960-19-7)	-	0.00943	0.0754	2	Malathion (121-75-5)	-	0.00943	0.0754	0.2
Acetamiprid (135410-20-7)	-	0.00943	0.0754	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.00943 0.00943	0.0754 0.0754	0.2 0.2
Aldicarb (116-06-3)	-	0.00943	0.0754	0.4	Methomyl (16752-77-5)	-	0.00943	0.0754	0.4
Azoxystrobin (131860-33-8)	-	0.00943	0.0754	0.2	Methyl parathion (298-0-0)	-	0.00943	0.0754	0.2
Bifenazate (149877-41-8)	-	0.00943	0.0754	0.2	MGK 264 (113-48-4)	-	0.00943	0.0754	0.2
Bifenthrin (82657-04-3)	-	0.00943	0.0754	0.2	Myclobutanil	_	0.00943	0.0754	0.2
Boscalid (188425-85-6)	-	0.00943	0.0754	0.4	(88671-89-0)	_	0.00545	0.0734	0.2
Carbaryl (63-25-2)	-	0.00943	0.0754	0.2	Naled (300-76-5)	-	0.00943	0.0754	0.5
Carbofuran (1563-66-2)	_	0.00943	0.0754	0.2	Oxamyl (23135-22-0)	-	0.00943	0.0754	1
Chlorantraniliprole (800008-45-7)	-	0.00943	0.0754	0.2	Paclobutrazol (76738-62-0)	-	0.00943	0.0754	0.4
Chlorfenapyr (122453-73-0)	-	0.00943	0.0754	1	Permethrins (52645-53-1)	-	0.00943	0.0754	0.2
Chlorpyrifos (2921-88-2)	_	0.00943	0.0754	0.2	Phosmet (732-11-6)	-	0.00943	0.0754	0.2
Clofentezine (74115-24-5)	_	0.00943	0.0754	0.2	Piperonyl butoxide	_	0.00943	0.0754	2
Cyfluthrin (68359-37-5)	-	0.00943	0.0754	1	(51-03-6)				
Cypermethrin (52315-07-8)	-	0.00943	0.0754	1	Prallethrins (2331-36-9) Propiconazole	-	0.00943 0.00943	0.0754 0.0754	0.2
Daminozide (1596-84-5)	_	0.00943	0.0754	1	(60207-90-1))				
DDVP (62-73-7)	_	0.00943	0.0754	0.1	Propoxur (114-26-1)	-	0.00943	0.0754	0.2
Diazinon (333-41-5)	_	0.00943	0.0754	0.2	Pyrethrins (8003-34-7)	-	0.00943	0.0754	1
Dimethoate (60-51-5)	_	0.00943	0.0754	0.2	Pyridaben (96489-71-3)	-	0.00943	0.0754	0.2
Ethoprophos		0.00943	0.0754	0.2	Spinosad (168316-95-8)	-	0.00943	0.0754	0.2
(13194-48-4)					Spiromesifen (283594-90-1)	-	0.00943	0.0754	0.2
Etofenprox (80844-07-1)	-	0.00943	0.0754	0.4	Spirotetramat		0.00943	0.0754	0.2
Etoxazole (153233-91-1)	-	0.00943	0.0754	0.2	(203313-25-1)	_	0.00943	0.0754	0.2
Fenoxycarb (72490-01-8)	-	0.00943	0.0754	0.2	Spiroxamine		0.00943	0.0754	0.4
(E)-Fenpyroximate (134098-61-6)	-	0.00943	0.0754	0.4	(118134-30-8) Tebuconazole		0.00943	0.0754	0.4
Fipronil (120068-37-3)	-	0.00943	0.0754	0.4	(80443-41-0)		0.00545	0.0734	0.4
Flonicamid (158062-67-0)	-	0.00943	0.0754	1	Thiacloprid	_	0.00943	0.0754	0.2
Fludioxinil (131341-86-1)	-	0.00943	0.0754	0.4	(111988-49-9)				
Hexythiazox (78587-05-0)	-	0.00943	0.0754	1	Thiamethoxam (153719-23-4)	-	0.00943	0.0754	0.2
Imazalil (35554-44-0)	-	0.00943	0.0754	0.2	Trifloxystrobin				
Imidacloprid (138261-41-3)	-	0.00943	0.0754	0.4	(141517-21-7)		0.00943	0.0754	0.2



Color Key

Date Completed:01/15/2025

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.

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

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













Date Received: 01/14/2025

CERTIFICATE OF ANALYSIS (SAMPLE ID: SA39847)

Testing Location:Customer ID: 2168Order ID: OR11375Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13220923443 Mass: 1ea

Greenbrier, AR 72058 Fort Smith, AR 72903 Production Run: Date Collected: 01/13/2025

Cultivar (Strain) or Sample Description: ARV-Sea Salt Caramel Soft Chews 200mg 20pk

Date Completed: 01/15/2025

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

P20250106CAR01

Analysis Date/Time: 01/14/2025 1807 Method: ICP/MS Deviations from SOP:

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)	-	56.2	89.0	200
Cadmium (Cd)	-	56.2	89.0	200
Lead (Pb)	-	56.2	89.0	500
Mercury (Hg)	-	56.2	89.0	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: 01/14/2025

CERTIFICATE OF ANALYSIS (SAMPLE ID: SA39847)

Testing Location:Customer ID: 2168Sample ID: SA39847Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Edible

232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13220923443 Mass: 1ea

Greenbrier, AR 72058 Fort Smith, AR 72903 Production Run: Date Collected: 01/13/2025

Cultivar (Strain) or Sample Description: ARV-Sea Salt Caramel Soft Chews 200mg 20pk

Date Completed: 01/15/2025

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

P20250106CAR01

Analysis Date/Time: 01/15/2025 0454 **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.



