

**SUMMARY OF ANALYSIS (SAMPLE ID: SA37427)**

<b>Testing Location:</b> Arkansas 232 S. Broadview St. Greenbrier, AR 72058 License: ADA 05_H273	<b>Customer ID:</b> 2240 MaxOrizonz LLC PO Box 3373 Pearland, TX 77588 License: Not Entered or N/A	<b>Order ID:</b> OR10917 <b>Lot Number:</b> Not Entered <b>Batch Number:</b> Not Entered	<b>Sample Type:</b> Primary <b>Matrix:</b> Lotion/Salve <b>Mass:</b> 1fl oz <b>Date Collected:</b> 03/07/2024 <b>Date Received:</b> 03/20/2024
<b>Cultivar (Strain) or Sample Description:</b> Peppermint Cinnamon Salve			<b>Date Completed:</b> 03/22/2024

\*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 (%)**

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.

<b>Cannabinoids (Top 3)</b>	<b>(%)</b>	<b>mg/g</b>
Δ8-THC	1.82	18.2
Δ9-THC	0.171	1.71
CBD	0.00639	0.0639
TOTAL CBD	0.00639	0.0639
TOTAL THC	0.171	1.71
TOTAL CANNABINOIDS	2.00	20.0

<b>Contaminants</b>	<b>PASS/FAIL</b>
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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**REPORT OF LABORATORY ANALYSIS**

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

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**CERTIFICATE OF ANALYSIS (SAMPLE ID: SA37427)**

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**Cultivar (Strain) or Sample Description:** Peppermint Cinnamon Salve **Date Completed:** 03/22/2024

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

**Analysis Date/Time:** 03/20/2024 1247  
**Analyst:** PW

**Method:** HPLC/DAD  
**Instrument:** Agilent 1100

**Moisture Content (%):** -  
**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	ND	ND	0.00222	0.00517	-	-	-
CBCA	ND	ND	0.00688	0.0161	-	-	-
CBD	0.00639	0.0639	0.0156	0.0365	-	0.0639	1.92
CBDa	ND	ND	0.00575	0.0134	-	-	-
CBDV	0.00297	0.0297	0.00251	0.00585	-	0.0297	0.891
CBDVA	ND	ND	0.00669	0.0156	-	-	-
CBG	ND	ND	0.0101	0.0237	-	-	-
CBGA	ND	ND	0.0144	0.0170	-	-	-
CBL	ND	ND	0.0117	0.0274	-	-	-
CBN	ND	ND	0.00538	0.0126	-	-	-
CBNA	ND	ND	0.00581	0.0135	-	-	-
Δ9-THC	0.171	1.71	0.00645	0.0150	-	1.71	51.4
Δ8-THC	1.82	18.2	0.0101	0.0235	-	18.2	547
THCA	ND	ND	0.00350	0.00818	-	-	-
THCV	ND	ND	0.00840	0.0196	-	-	-
THCVA	ND	ND	0.00268	0.00624	-	-	-
<b>TOTAL</b>	<b>2.00</b>	<b>20.0</b>				<b>20.0</b>	<b>601</b>
<b>TOTAL CBC</b>	-	-				-	-
<b>TOTAL CBD</b>	0.00639	0.0639				0.0639	1.92
<b>TOTAL CBDV</b>	0.00297	0.0297				0.0297	0.891
<b>TOTAL CBG</b>	-	-				-	-
<b>TOTAL CBN</b>	-	-				-	-
<b>TOTAL THC</b>	0.171	1.71				1.71	51.4
<b>TOTAL THCv</b>	-	-				-	-



**SERVING MASS (g):** 1.00  
**SERVINGS/UNIT:** 30

"-" 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) + 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, UM - Measurement Uncertainty

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



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**Cultivar (Strain) or Sample Description:** Peppermint Cinnamon Salve **Date Completed:** 03/22/2024

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

**Analysis Date/Time:** 03/21/2024 1442 **Method:** HS/GC/MS **Deviations from SOP:**  
**Analyst:** KF **Instrument:** Agilent 7890/5975 None

Solvent	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)	Solvent	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
Acetone (67-64-1)	-	137	273	5000	n-Heptane (142-82-5)	-	137	273	5000
Acetonitrile (75-5-8)	-	137	273	410	n-Hexane (110-54-3)	-	47.8	95.6	290
Benzene (71-43-2)	-	1.37	2.73	2	Isobutane (75-28-5)	-	137	273	5000
n-Butane (106-97-2)	-	137	273	5000	Isopropanol (67-63-0)	-	137	273	5000
1-Butanol (71-36-3)	-	137	273	5000	Isopropyl acetate (108-21-4)	-	137	273	5000
2-Butanol (78-92-2)	-	137	273	5000	Isopropyl benzene (98-82-8)	-	13.7	27.3	70
2-Butanone (78-93-3)	-	137	273	5000	Methanol (67-56-1)	-	137	273	3000
Cyclohexane (110-82-7)	-	137	273	3880	2-Methylbutane (78-78-4)	-	137	273	5000
1,2-Dimethoxyethane (110-71-4)	-	13.7	27.3	100	Methylene chloride (75-9-2)	-	137	273	600
N,N-Dimethylacetamide (127-19-5)	-	137	273	1090	2-Methylpentane (107-83-5)	-	47.8	95.6	290
2,2-Dimethylbutane (75-83-2)	-	47.8	95.6	290	3-Methylpentane (96-10-0)	-	47.8	95.6	290
2,3-Dimethylbutane (79-29-8)	-	47.8	95.6	290	n-Pentane (109-66-0)	-	137	273	5000
N,N-Dimethylformamide (68-12-2)	-	137	273	880	1-Pentanol (71-41-0)	-	137	273	5000
Dimethylsulfoxide (67-68-5)	-	137	273	5000	n-Propane (74-98-6)	-	137	273	5000
1,4-Dioxane (123-91-1)	-	137	273	380	1-Propanol (71-23-8)	-	137	273	5000
Ethanol (64-17-5)	-	137	273	5000	Pyridine (110-86-1)	-	47.8	95.6	200
2-Ethoxyethanol (110-80-5)	-	47.8	95.6	160	Tetrahydrofuran (109-99-9)	-	137	273	720
Ethyl ether (60-29-7)	-	137	273	5000	Tetramethylene sulfone (126-33-0)	-	47.8	95.6	160
Ethyl acetate (141-78-6)	-	137	273	5000	Toluene (108-88-3)	-	137	273	890
Ethyl benzene (100-41-4)	-	137	273	2170	o-Xylene (95-47-6)	-	137	273	2170
Ethylene glycol (107-21-1)	-	137	273	620	m,p-Xylene (108-38-3 or 106-42-3)	-	137	273	2170
Ethylene oxide (75-21-8)	-	13.7	27.3	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170



**Color Key**

**RESULT < AL** (Green background)  
**RESULT > AL** (Red background)

"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.

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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**Cultivar (Strain) or Sample Description:** Peppermint Cinnamon Salve **Date Completed:** 03/22/2024

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

**Analysis Date/Time:** 03/20/2024 1330  
**Analyst:** KF

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

**Deviations from SOP:**  
None

Pesticide	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)	Pesticide	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
Abamectin (71751-41-2)	-	0.00979	0.0783	0.5	Kresoxim-methyl (143390-89-0)	-	0.00979	0.0783	0.4
Acephate (30560-19-1)	-	0.00979	0.0783	0.4	Malathion (121-75-5)	-	0.00979	0.0783	0.2
Acequinocyl (57960-19-7)	-	0.00979	0.0783	2	Metalaxyl (57837-19-1)	-	0.00979	0.0783	0.2
Acetamiprid (135410-20-7)	-	0.00979	0.0783	0.2	Methiocarb (2032-65-7)	-	0.00979	0.0783	0.2
Aldicarb (116-06-3)	-	0.00979	0.0783	0.4	Methomyl (16752-77-5)	-	0.00979	0.0783	0.4
Azoxystrobin (131860-33-8)	-	0.00979	0.0783	0.2	Methyl parathion (298-0-0)	-	0.00979	0.0783	0.2
Bifenazate (149877-41-8)	-	0.00979	0.0783	0.2	MGK 264 (113-48-4)	-	0.00979	0.0783	0.2
Bifenthrin (82657-04-3)	-	0.00979	0.0783	0.2	Myclobutanil (88671-89-0)	-	0.00979	0.0783	0.2
Boscalid (188425-85-6)	-	0.00979	0.0783	0.4	Naled (300-76-5)	-	0.00979	0.0783	0.5
Carbaryl (63-25-2)	-	0.00979	0.0783	0.2	Oxamyl (23135-22-0)	-	0.00979	0.0783	1
Carbofuran (1563-66-2)	-	0.00979	0.0783	0.2	Paclobotrazol (76738-62-0)	-	0.00979	0.0783	0.4
Chlorantraniliprole (800008-45-7)	-	0.00979	0.0783	0.2	Permethrins (52645-53-1)	-	0.00979	0.0783	0.2
Chlorfenapyr (122453-73-0)	-	0.00979	0.0783	1	Phosmet (732-11-6)	-	0.00979	0.0783	0.2
Chlorpyrifos (2921-88-2)	-	0.00979	0.0783	0.2	Piperonyl butoxide (51-03-6)	-	0.00979	0.0783	2
Clofentezine (74115-24-5)	-	0.00979	0.0783	0.2	Prallethrin (2331-36-9)	-	0.00979	0.0783	0.2
Cyfluthrin (68359-37-5)	-	0.00979	0.0783	1	Propiconazole (60207-90-1)	-	0.00979	0.0783	0.4
Cypermethrin (52315-07-8)	-	0.00979	0.0783	1	Propoxur (114-26-1)	-	0.00979	0.0783	0.2
Daminozide (1596-84-5)	-	0.00979	0.0783	1	Pyrethrins (8003-34-7)	-	0.00979	0.0783	1
DDVP (62-73-7)	-	0.00979	0.0783	0.1	Pyridaben (96489-71-3)	-	0.00979	0.0783	0.2
Diazinon (333-41-5)	-	0.00979	0.0783	0.2	Spinosad (168316-95-8)	-	0.00979	0.0783	0.2
Dimethoate (60-51-5)	-	0.00979	0.0783	0.2	Spiromesifen (283594-90-1)	-	0.00979	0.0783	0.2
Ethoprophos (13194-48-4)	-	0.00979	0.0783	0.2	Spirotetramat (203313-25-1)	-	0.00979	0.0783	0.2
Etofenprox (80844-07-1)	-	0.00979	0.0783	0.4	Spiroxamine (118134-30-8)	-	0.00979	0.0783	0.4
Etoxazole (153233-91-1)	-	0.00979	0.0783	0.2	Tebuconazole (80443-41-0)	-	0.00979	0.0783	0.4
Fenoxycarb (72490-01-8)	-	0.00979	0.0783	0.2	Thiacloprid (111988-49-9)	-	0.00979	0.0783	0.2
(E)-Fenpyroximate (134098-61-6)	-	0.00979	0.0783	0.4	Thiamethoxam (153719-23-4)	-	0.00979	0.0783	0.2
Fipronil (120068-37-3)	-	0.00979	0.0783	0.4	Trifloxystrobin (141517-21-7)	-	0.00979	0.0783	0.2
Flonicamid (158062-67-0)	-	0.00979	0.0783	1					
Fludioxinil (131341-86-1)	-	0.00979	0.0783	0.4					
Hexythiazox (78587-05-0)	-	0.00979	0.0783	1					
Imazalil (35554-44-0)	-	0.00979	0.0783	0.2					
Imidacloprid (138261-41-3)	-	0.00979	0.0783	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.  
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	Synonym(s)	Pesticide	Synonym(s)	Pesticide	Synonym(s)
Cyfluthrin	Baythroid	Myclobutanil	Sythane	Propiconazole	Tilt
DDVP	Dichlorvos	Naled	Dibrom	Propoxur	Baygon
Ethoprophos	Prophos	Phosmet	Imidan		

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<b>Cultivar (Strain) or Sample Description:</b> Peppermint Cinnamon Salve			<b>Date Completed:</b> 03/22/2024

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

<b>Analysis Date/Time:</b> 03/21/2024 1826 (ICP/OES)	<b>Method:</b> ICP/MS	<b>Deviations from SOP:</b>
<b>Analysis Date/Time:</b> - (DMA)	<b>Instrument:</b> Agilent 7500ce	None
<b>Analyst:</b> 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)	-	58.7	93.0	200
Cadmium (Cd)	-	58.7	93.0	200
Lead (Pb)	-	58.7	93.0	500
Mercury (Hg)	-	58.7	93.0	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**

<b>RESULT &lt; AL</b>
<b>RESULT &gt; AL</b>

"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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<b>Cultivar (Strain) or Sample Description:</b> Peppermint Cinnamon Salve			<b>Date Completed:</b> 03/22/2024

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

**Analysis Date/Time:** 03/21/2024 1403      **Method:** Hardy Diagnostics CompactDry      **Deviations from SOP:**  
**Analyst:** PW      **Instrument:** Thermo Incubator      None

<u>Bacteria/Microbe</u>	<u>Result (CFU/g)</u>	<u>Action Level (CFU/g)</u>
Aerobic Plate Count	Absent	-
Coliforms, Total	Absent	1
Escherichia Coli (E. Coli)	Absent	100
Mold/Yeast	Absent	-
Pseudomonas aeruginosa	NT	-
Salmonella spp.	Absent	-
Staphylococcus aureus	Absent	-



**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.

**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.

**REPORT OF LABORATORY ANALYSIS**

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