

**SUMMARY OF ANALYSIS (SAMPLE ID: SA35951)**

<b>Testing Location:</b> Arkansas 232 S. Broadview St. Greenbrier, AR 72058 License: ADA 05_H273	<b>Customer ID:</b> 2116 Beehive Blends 7825 S Highland Dr. Cottonwood Heights, UT 84121 License: Not Entered or N/A	<b>Order ID:</b> OR10595 <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> 90ml <b>Date Collected:</b> 08/31/2023 <b>Date Received:</b> 08/31/2023
<b>Cultivar (Strain) or Sample Description:</b> Relief Roll-On			<b>Date Completed:</b> 09/05/2023

\*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 Utah 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 Utah 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>
CBD	1.72	17.2
CBG	0.161	1.61
CBN	0.103	1.03
TOTAL CBD	1.72	17.2
TOTAL THC	-	-
TOTAL CANNABINOIDS	1.99	19.9

<b>Terpenes (Top 5)</b>	<b>(%)</b>	<b>µg/g</b>
α-Bisabolol		
Camphene		
δ-3-Carene		
β-Caryophyllene		
Caryophyllene oxide		
TOTAL TERPENES	-	-

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

**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: SA35951)**

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**Cultivar (Strain) or Sample Description:** Relief Roll-On **Date Completed:** 09/05/2023

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

**Analysis Date/Time:** 08/31/2023 2000  
**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	0.00230	0.00537	-	-	-
CBCA	ND	ND	0.00714	0.0167	-	-	-
CBD	1.72	17.2	0.0162	0.0379	-	17.2	1550
CBDV	0.00308	0.0308	0.00260	0.00608	-	0.0308	2.77
CBDVA	ND	ND	0.00694	0.0162	-	-	-
CBG	0.161	1.61	0.0105	0.0246	-	1.61	145
CBGA	ND	ND	0.0149	0.0176	-	-	-
CBL	ND	ND	0.0122	0.0284	-	-	-
CBN	0.103	1.03	0.00559	0.0131	-	1.03	93.1
CBNA	ND	ND	0.00603	0.0141	-	-	-
Δ9-THC	ND	ND	0.00670	0.0156	-	-	-
Δ8-THC	ND	ND	0.0105	0.0244	-	-	-
THCA	ND	ND	0.00363	0.00850	-	-	-
THCV	ND	ND	0.00872	0.0203	-	-	-
THCVA	ND	ND	0.00279	0.00648	-	-	-
<b>TOTAL</b>	1.99	19.9			-	19.9	1790
<b>TOTAL CBC</b>	-	-			-	-	-
<b>TOTAL CBD</b>	1.72	17.2			-	17.2	1550
<b>TOTAL CBDV</b>	0.00308	0.0308			-	0.0308	2.77
<b>TOTAL CBG</b>	0.161	1.61			-	1.61	145
<b>TOTAL CBN</b>	0.103	1.03			-	1.03	93.1
<b>TOTAL THC</b>	-	-			-	-	-
<b>TOTAL THC V</b>	-	-			-	-	-



**SERVING MASS (g):** 1.00  
**SERVINGS/UNIT:** 90

"-" 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 THC V = (THCVA x 0.867) + THC V

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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**Cultivar (Strain) or Sample Description:** Relief Roll-On **Date Completed:** 09/05/2023

**TERPENOID PROFILE**

**Analysis Date/Time:** 08/31/2023 2339  
**Analyst:** KF

**Method:** GC/MS  
**Instrument:** Agilent 7890/5975

**Deviations from SOP:**  
None

<u>Terpene</u>	<u>Result (µg/g)</u>	<u>Result (%)</u>
α-Bisabolol	ND	-
Camphene	ND	-
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	ND	-
Linalool	ND	-
β-Myrcene	ND	-
cis-Nerolidol	ND	-
trans-Nerolidol	ND	-
α-Ocimene	ND	-
β-Ocimene	ND	-
α-Pinene	ND	-
β-Pinene	ND	-
α-Terpinene	ND	-
γ-Terpinene	ND	-
Terpinolene	ND	-
<b>TOTAL</b>	0.000	0.000



*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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**Reporting Limit (µg/g): 8.41**

"-" Not detected above LOD.

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<b>Cultivar (Strain) or Sample Description:</b> Relief Roll-On			<b>Date Completed:</b> 09/05/2023

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

<b>Analysis Date/Time:</b> 09/01/2023 0938	<b>Method:</b> HS/GC/MS	<b>Deviations from SOP:</b>
<b>Analyst:</b> KF	<b>Instrument:</b> 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)	-	159	319	5000	n-Heptane (142-82-5)	-	159	319	5000
Acetonitrile (75-5-8)	-	159	319	410	n-Hexane (110-54-3)	-	55.8	112	290
Benzene (71-43-2)	-	1.59	3.19	2	Isobutane (75-28-5)	-	159	319	5000
n-Butane (106-97-2)	-	159	319	5000	Isopropanol (67-63-0)	1360000	159	319	5000
1-Butanol (71-36-3)	-	159	319	5000	Isopropyl acetate (108-21-4)	-	159	319	5000
2-Butanol (78-92-2)	-	159	319	5000	Isopropyl benzene (98-82-8)	-	15.9	31.9	70
2-Butanone (78-93-3)	-	159	319	5000	Methanol (67-56-1)	-	159	319	3000
Cyclohexane (110-82-7)	-	159	319	3880	2-Methylbutane (78-78-4)	-	159	319	5000
1,2-Dimethoxyethane (110-71-4)	-	15.9	31.9	100	Methylene chloride (75-9-2)	-	159	319	600
N,N-Dimethylacetamide (127-19-5)	-	159	319	1090	2-Methylpentane (107-83-5)	-	55.8	112	290
2,2-Dimethylbutane (75-83-2)	-	55.8	112	290	3-Methylpentane (96-10-0)	-	55.8	112	290
2,3-Dimethylbutane (79-29-8)	-	55.8	112	290	n-Pentane (109-66-0)	-	159	319	5000
N,N-Dimethylformamide (68-12-2)	-	159	319	880	1-Pentanol (71-41-0)	-	159	319	5000
Dimethylsulfoxide (67-68-5)	-	159	319	5000	n-Propane (74-98-6)	-	159	319	5000
1,4-Dioxane (123-91-1)	-	159	319	380	1-Propanol (71-23-8)	-	159	319	5000
Ethanol (64-17-5)	-	159	319	5000	Pyridine (110-86-1)	-	55.8	112	100
2-Ethoxyethanol (110-80-5)	-	55.8	112	160	Tetrahydrofuran (109-99-9)	-	159	319	720
Ethyl ether (60-29-7)	-	159	319	5000	Tetramethylene sulfone (126-33-0)	-	55.8	112	160
Ethyl acetate (141-78-6)	-	159	319	5000	Toluene (108-88-3)	-	159	319	890
Ethyl benzene (100-41-4)	-	159	319	2170	o-Xylene (95-47-6)	-	159	319	2170
Ethylene glycol (107-21-1)	-	159	319	620	m,p-Xylene (108-38-3 or 106-42-3)	-	159	319	2170
Ethylene oxide (75-21-8)	-	15.9	31.9	-	Xylenes* (1330-20-7)	-	43.3	86.7	2170



**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 Utah 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:** Relief Roll-On **Date Completed:** 09/05/2023

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

**Analysis Date/Time:** 08/31/2023 2020  
**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)	-	0.0101	0.0809	0.5	Kresoxim-methyl (143390-89-0)	-	0.0101	0.0809	0.4
Acephate (30560-19-1)	-	0.0101	0.0809	0.4	Malathion (121-75-5)	-	0.0101	0.0809	0.2
Acetaminocyl (57960-19-7)	-	0.0101	0.0809	2	Metalaxyl (57837-19-1)	-	0.0101	0.0809	0.2
Acetamiprid (135410-20-7)	-	0.0101	0.0809	0.2	Methiocarb (2032-65-7)	-	0.0101	0.0809	0.2
Aldicarb (116-06-3)	-	0.0101	0.0809	0.4	Methomyl (16752-77-5)	-	0.0101	0.0809	0.4
Azoxystrobin (131860-33-8)	-	0.0101	0.0809	0.2	Methyl parathion (298-0-0)	-	0.0101	0.0809	0.2
Bifenazate (149877-41-8)	-	0.0101	0.0809	0.2	MGK 264 (113-48-4)	-	0.0101	0.0809	0.2
Bifenthrin (82657-04-3)	-	0.0101	0.0809	0.2	Myclobutanil (88671-89-0)	-	0.0101	0.0809	0.2
Boscalid (188425-85-6)	-	0.0101	0.0809	0.4	Naled (300-76-5)	-	0.0101	0.0809	0.5
Carbaryl (63-25-2)	-	0.0101	0.0809	0.2	Oxamyl (23135-22-0)	-	0.0101	0.0809	1
Carbofuran (1563-66-2)	-	0.0101	0.0809	0.2	Pacllobutrazol (76738-62-0)	-	0.0101	0.0809	0.4
Chlorantraniliprole (800008-45-7)	-	0.0101	0.0809	0.2	Permethrins (52645-53-1)	-	0.0101	0.0809	0.2
Chlorfenapyr (122453-73-0)	-	0.0101	0.0809	1	Phosmet (732-11-6)	-	0.0101	0.0809	0.2
Chlorpyrifos (2921-88-2)	-	0.0101	0.0809	0.2	Piperonyl butoxide (51-03-6)	-	0.0101	0.0809	2
Clofentezine (74115-24-5)	-	0.0101	0.0809	0.2	Prallethrins (2331-36-9)	-	0.0101	0.0809	0.2
Cyfluthrin (68359-37-5)	-	0.0101	0.0809	1	Propiconazole (60207-90-1)	-	0.0101	0.0809	0.4
Cypermethrin (52315-07-8)	-	0.0101	0.0809	1	Propoxur (114-26-1)	-	0.0101	0.0809	0.2
Daminozide (1596-84-5)	-	0.0101	0.0809	1	Pyrethrins (8003-34-7)	-	0.0101	0.0809	1
DDVP (62-73-7)	-	0.0101	0.0809	0.1	Pyridaben (96489-71-3)	-	0.0101	0.0809	0.2
Diazinon (333-41-5)	-	0.0101	0.0809	0.2	Spinosad (168316-95-8)	-	0.0101	0.0809	0.2
Dimethoate (60-51-5)	-	0.0101	0.0809	0.2	Spiromesifen (283594-90-1)	-	0.0101	0.0809	0.2
Ethoprophos (13194-48-4)	-	0.0101	0.0809	0.2	Spirotetramat (203313-25-1)	-	0.0101	0.0809	0.2
Etofenprox (80844-07-1)	-	0.0101	0.0809	0.4	Spiroxamine (118134-30-8)	-	0.0101	0.0809	0.4
Etoxazole (153233-91-1)	-	0.0101	0.0809	0.2	Tebuconazole (80443-41-0)	-	0.0101	0.0809	0.4
Fenoxycarb (72490-01-8)	-	0.0101	0.0809	0.2	Thiacloprid (111988-49-9)	-	0.0101	0.0809	0.2
(E)-Fenpyroximate (134098-61-6)	-	0.0101	0.0809	0.4	Thiamethoxam (153719-23-4)	-	0.0101	0.0809	0.2
Fipronil (120068-37-3)	-	0.0101	0.0809	0.4	Trifloxystrobin (141517-21-7)	-	0.0101	0.0809	0.2
Fonicamid (158062-67-0)	-	0.0101	0.0809	1					
Fludioxinil (131341-86-1)	-	0.0101	0.0809	0.4					
Hexythiazox (78587-05-0)	-	0.0101	0.0809	1					
Imazalil (35554-44-0)	-	0.0101	0.0809	0.2					
Imidacloprid (138261-41-3)	-	0.0101	0.0809	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 Utah MMJ testing guidelines.  
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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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**Cultivar (Strain) or Sample Description:** Relief Roll-On **Date Completed:** 09/05/2023

**MYCOTOXIN PROFILE (SOP: SOP-MYCO-001)**

<b>Analysis Date/Time:</b> 08/31/2023 2020	<b>Method:</b> LC/MS/MS	<b>Deviations from SOP:</b>
<b>Analyst:</b> KF	<b>Instrument:</b> Shimadzu LC-8050	None

<u>Mycotoxin</u>	<u>Result</u> ( $\mu\text{g}/\text{kg}$ )	<u>Action Level</u> ( $\mu\text{g}/\text{kg}$ )
Aflatoxin B1	-	20
Aflatoxin B2	-	20
Aflatoxin G1	-	20
Aflatoxin G2	-	20
Ochratoxin A	-	20



Abbreviations: LC - Liquid Chromatography, MS - Mass Spectrometry, RL - Reporting Limit, AL - Action Level

**Color Key**



**Reporting Limit ( $\mu\text{g}/\text{kg}$ )**  
Action Level

"-" not detected above reporting limit

Action levels for mycotoxins are referenced from the State of Utah 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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**REPORT OF LABORATORY ANALYSIS**

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



**CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35951)**

<b>Testing Location:</b> Arkansas 232 S. Broadview St. Greenbrier, AR 72058 License: ADA 05_H273	<b>Customer ID:</b> 2116 Beehive Blends 7825 S Highland Dr. Cottonwood Heights, UT 84121 License: Not Entered or N/A	<b>Order ID:</b> OR10595 <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> 90ml <b>Date Collected:</b> 08/31/2023 <b>Date Received:</b> 08/31/2023
<b>Cultivar (Strain) or Sample Description:</b> Relief Roll-On			<b>Date Completed:</b> 09/05/2023

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

<b>Analysis Date/Time:</b> 08/31/2023 2123 (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 (µg/kg)</u>	<u>LOD (µg/kg)</u>	<u>LOQ (µg/kg)</u>	<u>Action Level (µg/kg)</u>
Arsenic (As)	-	54.9	86.9	2000
Cadmium (Cd)	-	54.9	86.9	820
Lead (Pb)	-	54.9	86.9	1200
Mercury (Hg)	-	54.9	86.9	400



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 Utah MMJ testing guidelines.

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<b>Cultivar (Strain) or Sample Description:</b> Relief Roll-On			<b>Date Completed:</b> 09/05/2023

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

**Analysis Date/Time:** 09/01/2023 1025      **Method:** Hardy Diagnostics CompactDry      **Deviations from SOP:**  
**Analyst:** PW      **Instrument:** Thermo Incubator      None

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



**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 Utah 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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