

SUMMARY OF ANALYSIS (SAMPLE ID: SA37206)

Testing Location: Arkansas 232 S. Broadview St. Greenbrier, AR 72058 License: ADH 113	Customer ID: 2168 River Valley Relief MIPS 5601 Old Greenwood Rd Suite C Fort Smith, AR 72903 License: 00065C	Order ID: OR10884 Lot Number: M00065C13244397329 Production Run: E20240226KPLPLRCD01	Sample Type: Primary Matrix: Concentrate Mass: 8each Date Collected: 02/26/2024 Date Received: 02/26/2024
Cultivar (Strain) or Sample Description: Key Lime Pie Hybrid Live Rosin			Date Completed: 02/28/2024

*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.

Cannabinoids (Top 3)	(%)	mg/g
Δ9-THC	79.4	794
CBG	2.40	24.0
CBC	1.13	11.3
TOTAL CBD	0.520	5.20
TOTAL THC	79.4	794
TOTAL CANNABINOIDS	84.6	846

Terpenes (Top 5)	(%)	µg/g
β-Myrcene	1.36	13600
d-Limonene	1.03	10300
β-Caryophyllene	0.744	7440
Linalool	0.420	4200
α-Humulene	0.399	3990
TOTAL TERPENES	4.84	48400

Contaminants	PASS/FAIL
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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Cultivar (Strain) or Sample Description: Key Lime Pie Hybrid Live Rosin			Date Completed: 02/28/2024

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

Analysis Date/Time: 02/27/2024 1454 **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	1.13	11.3	0.106	0.248	-	11.3	11.3
CBCA	ND	ND	0.329	0.769	-	-	-
CBD	0.520	5.20	0.748	1.75	-	5.20	5.20
CBDV	ND	ND	0.275	0.643	-	-	-
CBDVA	ND	ND	0.320	0.747	-	-	-
CBG	2.40	24.0	0.486	1.13	-	24.0	24.0
CBGA	0.688	6.88	0.689	0.812	-	6.88	6.88
CBL	ND	ND	0.561	1.31	-	-	-
CBN	ND	ND	0.258	0.602	-	-	-
CBNA	ND	ND	0.278	0.649	-	-	-
Δ9-THC	79.4	794	0.309	0.720	-	794	794
Δ8-THC	ND	ND	0.482	1.13	-	-	-
THCA	ND	ND	0.168	0.392	-	-	-
THCV	0.525	5.25	0.402	0.937	-	5.25	5.25
THCVA	ND	ND	0.128	0.299	-	-	-
TOTAL	84.6	846				846	846
TOTAL CBC	1.13	11.3				11.3	11.3
TOTAL CBD	0.520	5.20				5.20	5.20
TOTAL CBDV	-	-				-	-
TOTAL CBG	3.00	30.0				30.0	30.0
TOTAL CBN	-	-				-	-
TOTAL THC	79.4	794				794	794
TOTAL THCV	0.525	5.25				5.25	5.25



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

"-" 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 = (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) + THCVA

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: Key Lime Pie Hybrid Live Rosin			Date Completed: 02/28/2024

TERPENOID PROFILE

Analysis Date/Time: 02/27/2024 1538
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	1320	0.132	█
Camphene	984	0.0984	█
δ-3-Carene	66.1	0.00661	█
β-Caryophyllene	7440	0.744	█
Caryophyllene oxide	1190	0.119	█
p-Cymene	<LOQ	-	
Eucalyptol	65.1	0.00651	█
Geraniol	ND	-	
Guaiol	239	0.0239	█
α-Humulene	3990	0.399	█
Isopulegol	ND	-	
d-Limonene	10300	1.03	█
Linalool	4200	0.420	█
β-Myrcene	13600	1.36	█
cis-Nerolidol	ND	-	
trans-Nerolidol	869	0.0869	█
α-Ocimene	55.8	0.00558	█
β-Ocimene	<LOQ	-	
α-Pinene	3290	0.329	█
β-Pinene	322	0.0322	█
α-Terpinene	110	0.0110	█
γ-Terpinene	77.2	0.00772	█
Terpinolene	262	0.0262	█
TOTAL	48400	4.84	



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): 46.5

"-" Not detected above LOD.

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

Analysis Date/Time: 02/26/2024 2117	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)	-	50.7	101	5000	n-Heptane (142-82-5)	-	50.7	101	5000
Acetonitrile (75-5-8)	-	50.7	101	410	n-Hexane (110-54-3)	-	17.8	35.5	290
Benzene (71-43-2)	-	0.507	1.01	2	Isobutane (75-28-5)	-	50.7	101	5000
n-Butane (106-97-2)	-	50.7	101	5000	Isopropanol (67-63-0)	-	50.7	101	5000
1-Butanol (71-36-3)	-	50.7	101	5000	Isopropyl acetate (108-21-4)	-	50.7	101	5000
2-Butanol (78-92-2)	-	50.7	101	5000	Isopropyl benzene (98-82-8)	-	5.07	10.1	70
2-Butanone (78-93-3)	-	50.7	101	5000	Methanol (67-56-1)	-	50.7	101	3000
Cyclohexane (110-82-7)	-	50.7	101	3880	2-Methylbutane (78-78-4)	-	50.7	101	5000
1,2-Dimethoxyethane (110-71-4)	-	5.07	10.1	100	Methylene chloride (75-9-2)	-	50.7	101	600
N,N-Dimethylacetamide (127-19-5)	-	50.7	101	1090	2-Methylpentane (107-83-5)	-	17.8	35.5	290
2,2-Dimethylbutane (75-83-2)	-	17.8	35.5	290	3-Methylpentane (96-10-0)	-	17.8	35.5	290
2,3-Dimethylbutane (79-29-8)	-	17.8	35.5	290	n-Pentane (109-66-0)	-	50.7	101	5000
N,N-Dimethylformamide (68-12-2)	-	50.7	101	880	1-Pentanol (71-41-0)	-	50.7	101	5000
Dimethylsulfoxide (67-68-5)	-	50.7	101	5000	n-Propane (74-98-6)	-	50.7	101	5000
1,4-Dioxane (123-91-1)	-	50.7	101	380	1-Propanol (71-23-8)	-	50.7	101	5000
Ethanol (64-17-5)	-	50.7	101	5000	Pyridine (110-86-1)	-	17.8	35.5	200
2-Ethoxyethanol (110-80-5)	-	17.8	35.5	160	Tetrahydrofuran (109-99-9)	-	50.7	101	720
Ethyl ether (60-29-7)	-	50.7	101	5000	Tetramethylene sulfone (126-33-0)	-	17.8	35.5	160
Ethyl acetate (141-78-6)	-	50.7	101	5000	Toluene (108-88-3)	-	50.7	101	890
Ethyl benzene (100-41-4)	-	50.7	101	2170	o-Xylene (95-47-6)	-	50.7	101	2170
Ethylene glycol (107-21-1)	-	50.7	101	620	m,p-Xylene (108-38-3 or 106-42-3)	-	50.7	101	2170
Ethylene oxide (75-21-8)	-	5.07	10.1	50	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 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: 02/27/2024 1508
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)	-	0.0464	0.371	0.5	Kresoxim-methyl (143390-89-0)	-	0.0464	0.371	0.4
Acephate (30560-19-1)	-	0.0464	0.371	0.4	Malathion (121-75-5)	-	0.0464	0.371	0.2
Acetamiprid (135410-20-7)	-	0.0464	0.371	0.2	Metalaxyl (57837-19-1)	-	0.0464	0.371	0.2
Aldicarb (116-06-3)	-	0.0464	0.371	0.4	Methiocarb (2032-65-7)	-	0.0464	0.371	0.2
Azoxystrobin (131860-33-8)	-	0.0464	0.371	0.2	Methomyl (16752-77-5)	-	0.0464	0.371	0.4
Bifenazate (149877-41-8)	-	0.0464	0.371	0.2	Methyl parathion (298-0-0)	-	0.0464	0.371	0.2
Bifenthrin (82657-04-3)	-	0.0464	0.371	0.2	MGK 264 (113-48-4)	-	0.0464	0.371	0.2
Boscalid (188425-85-6)	-	0.0464	0.371	0.4	Myclobutanil (88671-89-0)	-	0.0464	0.371	0.2
Carbaryl (63-25-2)	-	0.0464	0.371	0.2	Naled (300-76-5)	-	0.0464	0.371	0.5
Carbofuran (1563-66-2)	-	0.0464	0.371	0.2	Oxamyl (23135-22-0)	-	0.0464	0.371	1
Chlorantraniliprole (800008-45-7)	-	0.0464	0.371	0.2	Pacllobutrazol (76738-62-0)	-	0.0464	0.371	0.4
Chlorfenapyr (122453-73-0)	-	0.0464	0.371	1	Permethrins (52645-53-1)	-	0.0464	0.371	0.2
Chlorpyrifos (2921-88-2)	-	0.0464	0.371	0.2	Phosmet (732-11-6)	-	0.0464	0.371	0.2
Clofentezine (74115-24-5)	-	0.0464	0.371	0.2	Piperonyl butoxide (51-03-6)	-	0.0464	0.371	2
Cyfluthrin (68359-37-5)	-	0.0464	0.371	1	Prallethrins (2331-36-9)	-	0.0464	0.371	0.2
Cypermethrin (52315-07-8)	-	0.0464	0.371	1	Propiconazole (60207-90-1))	-	0.0464	0.371	0.4
Daminozide (1596-84-5)	-	0.0464	0.371	1	Propoxur (114-26-1)	-	0.0464	0.371	0.2
DDVP (62-73-7)	-	0.0464	0.371	0.1	Pyrethrins (8003-34-7)	-	0.0464	0.371	1
Diazinon (333-41-5)	-	0.0464	0.371	0.2	Pyridaben (96489-71-3)	-	0.0464	0.371	0.2
Dimethoate (60-51-5)	-	0.0464	0.371	0.2	Spinosad (168316-95-8)	-	0.0464	0.371	0.2
Ethoprophos (13194-48-4)	-	0.0464	0.371	0.2	Spiromesifen (283594-90-1)	-	0.0464	0.371	0.2
Etofenprox (80844-07-1)	-	0.0464	0.371	0.4	Spirotetramat (203313-25-1)	-	0.0464	0.371	0.2
Etoxazole (153233-91-1)	-	0.0464	0.371	0.2	Spiroxamine (118134-30-8)	-	0.0464	0.371	0.4
Fenoxycarb (72490-01-8)	-	0.0464	0.371	0.2	Tebuconazole (80443-41-0)	-	0.0464	0.371	0.4
(E)-Fenpyroximate (134098-61-6)	-	0.0464	0.371	0.4	Thiacloprid (111988-49-9)	-	0.0464	0.371	0.2
Fipronil (120068-37-3)	-	0.0464	0.371	0.4	Thiamethoxam (153719-23-4)	-	0.0464	0.371	0.2
Fonicamid (158062-67-0)	-	0.0464	0.371	1	Trifloxystrobin (141517-21-7)	-	0.0464	0.371	0.2
Fludioxinil (131341-86-1)	-	0.0464	0.371	0.4					
Hexythiazox (78587-05-0)	-	0.0464	0.371	1					
Imazalil (35554-44-0)	-	0.0464	0.371	0.2					
Imidacloprid (138261-41-3)	-	0.0464	0.371	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.
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	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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HEAVY METAL PROFILE (SOP: SOP-ICP-200.6)

Analysis Date/Time: 02/27/2024 1755 (ICP/OES) **Method:** ICP/MS **Deviations from SOP:**
Analysis Date/Time: - (DMA) **Instrument:** Agilent 7500ce None
Analyst: 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)	-	58.5	92.6	200
Cadmium (Cd)	-	58.5	92.6	200
Lead (Pb)	-	58.5	92.6	500
Mercury (Hg)	-	58.5	92.6	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

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 "-" not detected above LOD

Action levels for heavy metals are referenced from the State of Arkansas MMJ testing guidelines.

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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA37206)

Testing Location: Arkansas 232 S. Broadview St. Greenbrier, AR 72058 License: ADA 05_H273	Customer ID: 2168 River Valley Relief MIPS 5601 Old Greenwood Rd Suite C Fort Smith, AR 72903 License: 00065C	Sample ID: SA37206 Lot Number: M00065C13244397329 Production Run: E20240226KPLRCD01	Sample Type: Primary Matrix: Concentrate Mass: 8each Date Collected: 02/26/2024 Date Received: 02/26/2024
Cultivar (Strain) or Sample Description: Key Lime Pie Hybrid Live Rosin			Date Completed: 02/28/2024

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 02/27/2024 1500 **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	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.

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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