

SUMMARY OF ANALYSIS (SAMPLE ID: SA36125)

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: OR10628 Lot Number: M00065C13249911251 Production Run: E20230929GBBLHR01	Sample Type: Primary Matrix: Concentrate Mass: 4g Date Collected: 10/02/2023 Date Received: 10/03/2023
Cultivar (Strain) or Sample Description: Glitter Bomb Indica Hybrid Solventless Live Hash Rosin			Date Completed: 10/04/2023

*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 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.

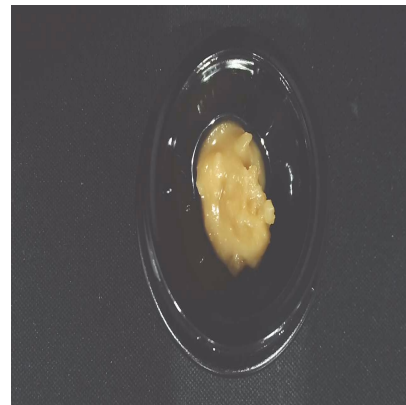
Moisture content levels less than 15% are recommended but the sample does not fail. Water activity levels must be less than 0.65aw.

Cannabinoids (Top 3)	(%)	mg/g
THCA	76.5	765
CBGA	2.76	27.6
Δ9-THC	2.47	24.7
TOTAL CBD	0.0995	0.995
TOTAL THC	69.6	696
TOTAL CANNABINOIDS	83.6	836

Terpenes (Top 5)	(%)	µg/g
β-Myrcene	5.40	54000
β-Caryophyllene	1.08	10800
β-Ocimene	0.400	4000
d-Limonene	0.368	3680
α-Humulene	0.316	3160
TOTAL TERPENES	8.53	85300

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

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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36125)

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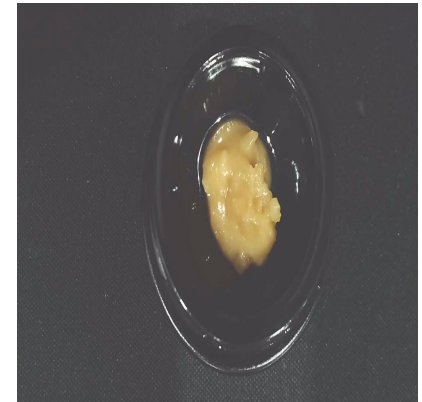
CANNABINOID (POTENCY) PROFILE (SOP: SOP-CANN-001)

Analysis Date/Time: 10/03/2023 1533
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	0.110	0.258	-	-	-
BCA	0.767	7.67	0.343	0.800	-	7.67	7.67
CBD	ND	ND	0.779	1.82	-	-	-
CBDA	0.113	1.13	0.287	0.669	-	1.13	1.13
CBDV	ND	ND	0.125	0.292	-	-	-
CBDVA	ND	ND	0.333	0.778	-	-	-
CBG	0.294	2.94	0.506	1.18	-	2.94	2.94
CBGA	2.76	27.6	0.717	0.846	-	27.6	27.6
CBL	ND	ND	0.584	1.36	-	-	-
CBN	ND	ND	0.268	0.627	-	-	-
CBNA	ND	ND	0.290	0.675	-	-	-
Δ9-THC	2.47	24.7	0.322	0.750	-	24.7	24.7
Δ8-THC	ND	ND	0.502	1.17	-	-	-
THCA	76.5	765	0.174	0.408	-	765	765
THCV	ND	ND	0.418	0.975	-	-	-
THCVA	0.731	7.31	0.134	0.311	-	7.31	7.31
TOTAL	83.6	836				836	836
TOTAL CBC	0.673	6.73				6.73	6.73
TOTAL CBD	0.0995	0.995				0.995	0.995
TOTAL CBDV	-	-				-	-
TOTAL CBG	2.72	27.2				27.2	27.2
TOTAL CBN	-	-				-	-
TOTAL THC	69.6	696				696	696
TOTAL THC V	0.634	6.34				6.34	6.34



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 = (BCA 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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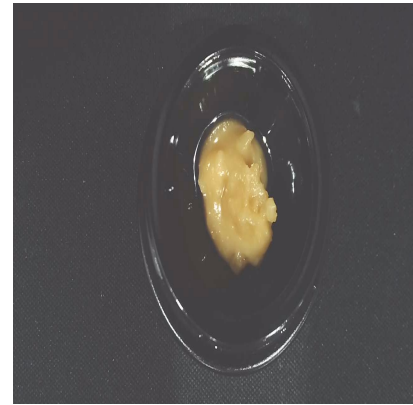
TERPENOID PROFILE

Analysis Date/Time: 10/04/2023 0202
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	807	0.0807
Camphene	194	0.0194
δ-3-Carene	ND	-
β-Caryophyllene	10800	1.08
Caryophyllene oxide	ND	-
p-Cymene	ND	-
Eucalyptol	226	0.0226
Geraniol	ND	-
Guaiol	ND	-
α-Humulene	3160	0.316
Isopulegol	ND	-
d-Limonene	3680	0.368
Linalool	2870	0.287
β-Myrcene	54000	5.40
cis-Nerolidol	646	0.0646
trans-Nerolidol	1420	0.142
α-Ocimene	ND	-
β-Ocimene	4000	0.400
α-Pinene	646	0.0646
β-Pinene	2420	0.242
α-Terpinene	161	0.0161
γ-Terpinene	129	0.0129
Terpinolene	129	0.0129
TOTAL	85300	8.53



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): 80.7

"-" Not detected above LOD.

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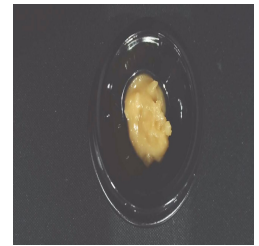
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RESIDUAL SOLVENT PROFILE (SOP: SOP-RS-001)

Analysis Date/Time: 10/03/2023 2114	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)	-	159	317	5000	n-Heptane (142-82-5)	-	159	317	5000
Acetonitrile (75-5-8)	-	159	317	410	n-Hexane (110-54-3)	-	55.6	111	290
Benzene (71-43-2)	-	1.59	3.17	2	Isobutane (75-28-5)	-	159	317	5000
n-Butane (106-97-2)	-	159	317	5000	Isopropanol (67-63-0)	-	159	317	5000
1-Butanol (71-36-3)	-	159	317	5000	Isopropyl acetate (108-21-4)	-	159	317	5000
2-Butanol (78-92-2)	-	159	317	5000	Isopropyl benzene (98-82-8)	-	15.9	31.7	70
2-Butanone (78-93-3)	-	159	317	5000	Methanol (67-56-1)	-	159	317	3000
Cyclohexane (110-82-7)	-	159	317	3880	2-Methylbutane (78-78-4)	-	159	317	5000
1,2-Dimethoxyethane (110-71-4)	-	15.9	31.7	100	Methylene chloride (75-9-2)	-	159	317	600
N,N-Dimethylacetamide (127-19-5)	-	159	317	1090	2-Methylpentane (107-83-5)	-	55.6	111	290
2,2-Dimethylbutane (75-83-2)	-	55.6	111	290	3-Methylpentane (96-10-0)	-	55.6	111	290
2,3-Dimethylbutane (79-29-8)	-	55.6	111	290	n-Pentane (109-66-0)	-	159	317	5000
N,N-Dimethylformamide (68-12-2)	-	159	317	880	1-Pentanol (71-41-0)	-	159	317	5000
Dimethylsulfoxide (67-68-5)	-	159	317	5000	n-Propane (74-98-6)	-	159	317	5000
1,4-Dioxane (123-91-1)	-	159	317	380	1-Propanol (71-23-8)	-	159	317	5000
Ethanol (64-17-5)	-	159	317	5000	Pyridine (110-86-1)	-	55.6	111	200
2-Ethoxyethanol (110-80-5)	-	55.6	111	160	Tetrahydrofuran (109-99-9)	-	159	317	720
Ethyl ether (60-29-7)	-	159	317	5000	Tetramethylene sulfone (126-33-0)	-	55.6	111	160
Ethyl acetate (141-78-6)	-	159	317	5000	Toluene (108-88-3)	-	159	317	890
Ethyl benzene (100-41-4)	-	159	317	2170	o-Xylene (95-47-6)	-	159	317	2170
Ethylene glycol (107-21-1)	-	159	317	620	m,p-Xylene (108-38-3 or 106-42-3)	-	159	317	2170
Ethylene oxide (75-21-8)	-	15.9	31.7	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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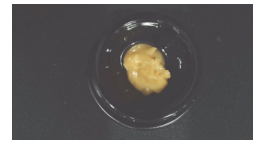
PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 10/03/2023 1611
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)	-	0.0482	0.385	0.5	Kresoxim-methyl (143390-89-0)	-	0.0482	0.385	0.4
Acephate (30560-19-1)	-	0.0482	0.385	0.4	Malathion (121-75-5)	-	0.0482	0.385	0.2
Acequinocyl (57960-19-7)	-	0.0482	0.385	2	Metalaxyl (57837-19-1)	-	0.0482	0.385	0.2
Acetamiprid (135410-20-7)	-	0.0482	0.385	0.2	Methiocarb (2032-65-7)	-	0.0482	0.385	0.2
Aldicarb (116-06-3)	-	0.0482	0.385	0.4	Methomyl (16752-77-5)	-	0.0482	0.385	0.4
Azoxystrobin (131860-33-8)	-	0.0482	0.385	0.2	Methyl parathion (298-0-0)	-	0.0482	0.385	0.2
Bifenazate (149877-41-8)	-	0.0482	0.385	0.2	MGK 264 (113-48-4)	-	0.0482	0.385	0.2
Bifenthrin (82657-04-3)	-	0.0482	0.385	0.2	Myclobutanil (88671-89-0)	-	0.0482	0.385	0.2
Boscalid (188425-85-6)	-	0.0482	0.385	0.4	Naled (300-76-5)	-	0.0482	0.385	0.5
Carbaryl (63-25-2)	-	0.0482	0.385	0.2	Oxamyl (23135-22-0)	-	0.0482	0.385	1
Carbofuran (1563-66-2)	-	0.0482	0.385	0.2	Pacllobutrazol (76738-62-0)	-	0.0482	0.385	0.4
Chlorantraniliprole (800008-45-7)	-	0.0482	0.385	0.2	Permethrins (52645-53-1)	-	0.0482	0.385	0.2
Chlorfenapyr (122453-73-0)	-	0.0482	0.385	1	Phosmet (732-11-6)	-	0.0482	0.385	0.2
Chlorpyrifos (2921-88-2)	-	0.0482	0.385	0.2	Piperonyl butoxide (51-03-6)	-	0.0482	0.385	2
Clofentezine (74115-24-5)	-	0.0482	0.385	0.2	Prallethrins (2331-36-9)	-	0.0482	0.385	0.2
Cyfluthrin (68359-37-5)	-	0.0482	0.385	1	Propiconazole (60207-90-1))	-	0.0482	0.385	0.4
Cypermethrin (52315-07-8)	-	0.0482	0.385	1	Propoxur (114-26-1)	-	0.0482	0.385	0.2
Daminozide (1596-84-5)	-	0.0482	0.385	1	Pyrethrins (8003-34-7)	-	0.0482	0.385	1
DDVP (62-73-7)	-	0.0482	0.385	0.1	Pyridaben (96489-71-3)	-	0.0482	0.385	0.2
Diazinon (333-41-5)	-	0.0482	0.385	0.2	Spinosad (168316-95-8)	-	0.0482	0.385	0.2
Dimethoate (60-51-5)	-	0.0482	0.385	0.2	Spiromesifen (283594-90-1)	-	0.0482	0.385	0.2
Ethoprophos (13194-48-4)	-	0.0482	0.385	0.2	Spirotetramat (203313-25-1)	-	0.0482	0.385	0.2
Etofenprox (80844-07-1)	-	0.0482	0.385	0.4	Spiroxamine (118134-30-8)	-	0.0482	0.385	0.4
Etoxazole (153233-91-1)	-	0.0482	0.385	0.2	Tebuconazole (80443-41-0)	-	0.0482	0.385	0.4
Fenoxycarb (72490-01-8)	-	0.0482	0.385	0.2	Thiacloprid (111988-49-9)	-	0.0482	0.385	0.2
(E)-Fenpyroximate (134098-61-6)	-	0.0482	0.385	0.4	Thiamethoxam (153719-23-4)	-	0.0482	0.385	0.2
Fipronil (120068-37-3)	-	0.0482	0.385	0.4	Trifloxystrobin (141517-21-7)	-	0.0482	0.385	0.2
Fonicamid (158062-67-0)	-	0.0482	0.385	1					
Fludioxinil (131341-86-1)	-	0.0482	0.385	0.4					
Hexythiazox (78587-05-0)	-	0.0482	0.385	1					
Imazalil (35554-44-0)	-	0.0482	0.385	0.2					
Imidacloprid (138261-41-3)	-	0.0482	0.385	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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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		

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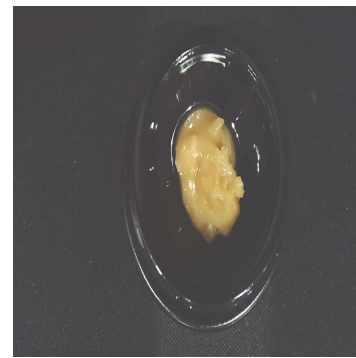
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HEAVY METAL PROFILE (SOP: SOP-ICP-200.6)

Analysis Date/Time: 10/03/2023 1840 (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)	-	59.7	94.6	200
Cadmium (Cd)	-	59.7	94.6	200
Lead (Pb)	-	59.7	94.6	500
Mercury (Hg)	-	59.7	94.6	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

"DET" detected less than LOQ

"-" not detected above LOD

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.

REPORT OF LABORATORY ANALYSIS

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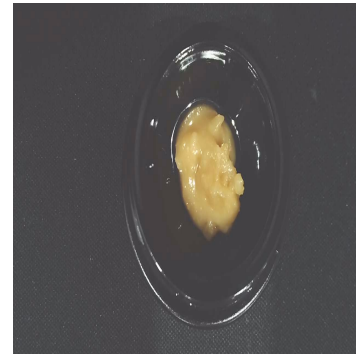
CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36125)

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: SA36125 Lot Number: M00065C13249911251 Production Run: E20230929GBBLHR01	Sample Type: Primary Matrix: Concentrate Mass: 4g Date Collected: 10/02/2023 Date Received: 10/03/2023 Date Completed: 10/04/2023
Cultivar (Strain) or Sample Description: Glitter Bomb Indica Hybrid Solventless Live Hash Rosin			

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 10/04/2023 0838 **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.

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

