

SUMMARY OF ANALYSIS (SAMPLE ID: SA43390)

Testing Location:	Customer ID: 2168	Order ID: OR12036	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13249906850	Mass: 3ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 03/02/2026
License: ADH 113	License: 00065C	E20260302HLBH01	Date Received: 03/02/2026
Cultivar (Strain) or Sample Description: Hoop Loops Solventless Bubble Hash 1.5g Jar			Date Completed: 03/04/2026

*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.

*Where provided, statements of conformity (e.g. Pass/Fail) are made in accordance with ILAC G8, Binary Statement for Simple Acceptance Rule (w=0, AL=TL).
PASS: when the result is within the acceptance interval. FAIL: when the result is outside the acceptance interval

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	61.6	616
CBGA	2.15	21.5
Δ9-THC	1.45	14.5
TOTAL CBD	0.390	3.90
TOTAL THC	55.5	555
TOTAL CANNABINOIDS	67.5	675
Terpenes (Top 5)	(%)	ppm
β-Caryophyllene	0.697	6970
Linalool	0.581	5810
d-Limonene	0.510	5100
β-Myrcene	0.239	2390
α-Humulene	0.205	2050
TOTAL TERPENES	3.20	32000

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

Analysis Date/Time: 03/03/2026 1520

Method: HPLC/DAD

Moisture Content (%): -

Analyst: PW

Instrument: Agilent 1100

Water Activity (aw): -

Cannabinoid	Result (%)	Result (mg/g)	LOD (mg/g)	LOQ (mg/g)	Result (mg/mL)	Per Serving (mg)	Per Unit (mg)
CBC	ND	ND	0.109	0.254	-	0.000	0.000
CBCA	1.07	10.7	0.338	0.789	-	10.7	10.7
CBD	0.285	2.85	0.768	1.79	-	2.85	2.85
CBDV	0.120	1.20	0.283	0.660	-	1.20	1.20
CBDVA	ND	ND	0.123	0.287	-	0.000	0.000
CBDVA	ND	ND	0.329	0.767	-	0.000	0.000
CBG	0.327	3.27	0.499	1.16	-	3.27	3.27
CBGA	2.15	21.5	0.707	0.834	-	21.5	21.5
CBL	ND	ND	0.576	1.34	-	0.000	0.000
CBN	0.157	1.57	0.265	0.618	-	1.57	1.57
CBNA	ND	ND	0.286	0.666	-	0.000	0.000
Δ9-THC	1.45	14.5	0.317	0.739	-	14.5	14.5
Δ8-THC	ND	ND	0.495	1.15	-	0.000	0.000
THCA	61.6	616	0.172	0.402	-	616	616
THCV	ND	ND	0.413	0.962	-	0.000	0.000
THCVA	0.303	3.03	0.132	0.307	-	3.03	3.03
TOTAL	67.5	675			-	675	675
TOTAL CBC	0.938	9.38			-	9.38	9.38
TOTAL CBD	0.390	3.90			-	3.90	3.90
TOTAL CBDV	0.000	0.000			-	0.000	0.000
TOTAL CBG	2.21	22.1			-	22.1	22.1
TOTAL CBN	0.157	1.57			-	1.57	1.57
TOTAL THC	55.5	555			-	555	555
TOTAL THCV	0.263	2.63			-	2.63	2.63



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

"-" Not reported for this sample.

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.
Dry percent = Wet percent / (1-(Moisture Content/100))

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) + 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, ND - Not Detected (less than LOD)

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Laboratory Director



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Cultivar (Strain) or Sample Description: Hoop Loops Solventless Bubble Hash 1.5g Jar			Date Completed: 03/04/2026

TERPENOID PROFILE

Analysis Date/Time: 03/03/2026 2227

Method: GC/MS

Deviations from SOP:

Analyst: KF

Instrument: Agilent 7890/5975

None

<u>Terpene</u>	<u>Result (ppm)</u>	<u>Result (%)</u>	<u>Terpene</u>	<u>Result (ppm)</u>	<u>Result (%)</u>
α-Bisabolol	342	0.0342	d-Limonene	5100	0.510
L-Borneol	ND	-	Linalool	5810	0.581
D-Borneol	146	0.0146	l-Menthone	ND	-
Camphene	275	0.0275	Menthol	134	0.0134
Camphor	ND	-	β-Myrcene	2390	0.239
δ-3-Carene	<LOQ	-	Nerol	ND	-
Carvacrol	ND	-	cis-Nerolidol	ND	-
Carvone	ND	-	trans-Nerolidol	679	0.0679
β-Caryophyllene	6970	0.697	Nootkatone	ND	-
Caryophyllene oxide	188	0.0188	cis-β-Ocimene	51.5	0.00515
Cedrene	57.6	0.00576	trans-β-Ocimene	<LOQ	-
Cedrol	ND	-	Octyl acetate	<LOQ	-
cis-Citral	ND	-	α-Phellandrene	<LOQ	-
trans-Citral	ND	-	Phytane	ND	-
Citronellol	ND	-	α-Pinene	811	0.0811
o-Cymene	ND	-	β-Pinene	1050	0.105
m-Cymene	ND	-	Piperitone	ND	-
p-Cymene	ND	-	Pulegone	ND	-
Eucalyptol	<LOQ	-	Sabinene	624	0.0624
trans-β-Farnesene	1260	0.126	Sabinene hydrate	<LOQ	-
Farnesol 1	132	0.0132	Safranal	ND	-
Farnesol 2	362	0.0362	Squalene	125	0.0125
Fenchone	<LOQ	-	Terpinen-4-ol	<LOQ	-
Fenchyl alcohol	1030	0.103	α-Terpinene	<LOQ	-
Geraniol	88.7	0.00887	γ-Terpinene	<LOQ	-
Geranyl acetate	439	0.0439	α-Terpineol	919	0.0919
Guaiol	510	0.0510	Terpinolene	98.5	0.00985
α-Humulene	2050	0.205	α-Thujone	ND	-
Isoborneol	<LOQ	-	Thymol	69.7	0.00697
Isobornyl acetate	ND	-	Valencene	166	0.0166
d-Isomenthone	ND	-	d-Valerolactam	ND	-
Isopulegol	117	0.0117	Verbenone	ND	-



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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"-" Not detected above LOD.

TOTAL (ppm) 32000 (%) 3.20 **Reporting Limit (ppm):** 47.8

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

Analysis Date/Time: 03/03/2026 2106	Method: HS/GC/MS	Deviations from SOP:
Analyst: KF	Instrument: Agilent 7890/5975	None

<u>Solvent</u>	<u>Result</u> (ppm)	<u>LOD</u> (ppm)	<u>LOQ</u> (ppm)	<u>Action</u> <u>Level</u> (ppm)	<u>Solvent</u>	<u>Result</u> (ppm)	<u>LOD</u> (ppm)	<u>LOQ</u> (ppm)	<u>Action</u> <u>Level</u> (ppm)
Acetone (67-64-1)	-	39.0	78.0	5000	n-Heptane (142-82-5)	-	39.0	78.0	5000
Acetonitrile (75-5-8)	-	39.0	78.0	410	n-Hexane (110-54-3)	-	13.7	27.3	290
Benzene (71-43-2)	-	0.390	0.780	2	Isobutane (75-28-5)	-	39.0	78.0	5000
n-Butane (106-97-2)	-	39.0	78.0	5000	Isopropanol (67-63-0)	-	39.0	78.0	5000
1-Butanol (71-36-3)	-	39.0	78.0	5000	Isopropyl acetate (108-21-4)	-	39.0	78.0	5000
2-Butanol (78-92-2)	-	39.0	78.0	5000	Isopropyl benzene (98-82-8)	-	3.90	7.80	70
2-Butanone (78-93-3)	-	39.0	78.0	5000	Methanol (67-56-1)	-	39.0	78.0	3000
Cyclohexane (110-82-7)	-	39.0	78.0	3880	2-Methylbutane (78-78-4)	-	39.0	78.0	5000
1,2-Dimethoxyethane (110-71-4)	-	3.90	7.80	100	Methylene chloride (75-9-2)	-	39.0	78.0	600
N,N-Dimethylacetamide (127-19-5)	-	39.0	78.0	1090	2-Methylpentane (107-83-5)	-	13.7	27.3	290
2,2-Dimethylbutane (75-83-2)	-	13.7	27.3	290	3-Methylpentane (96-10-0)	-	13.7	27.3	290
2,3-Dimethylbutane (79-29-8)	-	13.7	27.3	290	n-Pentane (109-66-0)	-	39.0	78.0	5000
N,N-Dimethylformamide (68-12-2)	-	39.0	78.0	880	1-Pentanol (71-41-0)	-	39.0	78.0	5000
Dimethylsulfoxide (67-68-5)	-	39.0	78.0	5000	n-Propane (74-98-6)	-	39.0	78.0	5000
1,4-Dioxane (123-91-1)	-	39.0	78.0	380	1-Propanol (71-23-8)	-	39.0	78.0	5000
Ethanol (64-17-5)	-	39.0	78.0	5000	Pyridine (110-86-1)	-	13.7	27.3	200
2-Ethoxyethanol (110-80-5)	-	13.7	27.3	160	Tetrahydrofuran (109-99-9)	-	39.0	78.0	720
Ethyl ether (60-29-7)	-	39.0	78.0	5000	Tetramethylene sulfone (126-33-0)	-	13.7	27.3	160
Ethyl acetate (141-78-6)	-	39.0	78.0	5000	Toluene (108-88-3)	-	39.0	78.0	890
Ethyl benzene (100-41-4)	-	39.0	78.0	2170	o-Xylene (95-47-6)	-	39.0	78.0	2170
Ethylene glycol (107-21-1)	-	39.0	78.0	620	m,p-Xylene (108-38-3 or 106-42-3)	-	39.0	78.0	2170
Ethylene oxide (75-21-8)	-	3.90	7.80	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170



Color Key

RESULT < AL

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>	<u>Synonym(s)</u>	<u>Solvent</u>	<u>Synonym(s)</u>
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: 03/03/2026 1546

Analyst: KF

Method: LC/MS/MS

Instrument: Shimadzu LC-8050

Deviations from SOP:

None

Pesticide	Result (ppm)	LOD (ppm)	LOQ (ppm)	Action Level (ppm)	Pesticide	Result (ppm)	LOD (ppm)	LOQ (ppm)	Action Level (ppm)
Abamectin (71751-41-2)	-	0.0477	0.382	0.5	Kresoxim-methyl (143390-89-0)	-	0.0477	0.382	0.4
Acephate (30560-19-1)	-	0.0477	0.382	0.4	Malathion (121-75-5)	-	0.0477	0.382	0.2
Acetaminocyl (57960-19-7)	-	0.0477	0.382	2	Metalaxyl (57837-19-1)	-	0.0477	0.382	0.2
Acetamiprid (135410-20-7)	-	0.0477	0.382	0.2	Methiocarb (2032-65-7)	-	0.0477	0.382	0.2
Aldicarb (116-06-3)	-	0.0477	0.382	0.4	Methomyl (16752-77-5)	-	0.0477	0.382	0.4
Azoxystrobin (131860-33-8)	-	0.0477	0.382	0.2	Methyl parathion (298-0-0)	-	0.0477	0.382	0.2
Bifenazate (149877-41-8)	-	0.0477	0.382	0.2	MGK 264 (113-48-4)	-	0.0477	0.382	0.2
Bifenthrin (82657-04-3)	-	0.0477	0.382	0.2	Myclobutanil (88671-89-0)	-	0.0477	0.382	0.2
Boscalid (188425-85-6)	-	0.0477	0.382	0.4	Naled (300-76-5)	-	0.0477	0.382	0.5
Carbaryl (63-25-2)	-	0.0477	0.382	0.2	Oxamyl (23135-22-0)	-	0.0477	0.382	1
Carbofuran (1563-66-2)	-	0.0477	0.382	0.2	Paclobutrazol (76738-62-0)	-	0.0477	0.382	0.4
Chlorantraniliprole (800008-45-7)	-	0.0477	0.382	0.2	Permethrins (52645-53-1)	-	0.0477	0.382	0.2
Chlorfenapyr (122453-73-0)	-	0.0477	0.382	1	Phosmet (732-11-6)	-	0.0477	0.382	0.2
Chlorpyrifos (2921-88-2)	-	0.0477	0.382	0.2	Piperonyl butoxide (51-03-6)	-	0.0477	0.382	2
Clofentezine (74115-24-5)	-	0.0477	0.382	0.2	Prallethrin (2331-36-9)	-	0.0477	0.382	0.2
Cyfluthrin (68359-37-5)	-	0.0477	0.382	1	Propiconazole (60207-90-1)	-	0.0477	0.382	0.4
Cypermethrin (52315-07-8)	-	0.0477	0.382	1	Propoxur (114-26-1)	-	0.0477	0.382	0.2
Daminozide (1596-84-5)	-	0.0477	0.382	1	Pyrethrins (8003-34-7)	-	0.0477	0.382	1
DDVP (62-73-7)	-	0.0477	0.382	0.1	Pyridaben (96489-71-3)	-	0.0477	0.382	0.2
Diazinon (333-41-5)	-	0.0477	0.382	0.2	Spinosad (168316-95-8)	-	0.0477	0.382	0.2
Dimethoate (60-51-5)	-	0.0477	0.382	0.2	Spiromesifen (283594-90-1)	-	0.0477	0.382	0.2
Ethoprophos (13194-48-4)	-	0.0477	0.382	0.2	Spirotetramat (203313-25-1)	-	0.0477	0.382	0.2
Etofenprox (80844-07-1)	-	0.0477	0.382	0.4	Spiroxamine (118134-30-8)	-	0.0477	0.382	0.4
Etoxazole (153233-91-1)	-	0.0477	0.382	0.2	Tebuconazole (80443-41-0)	-	0.0477	0.382	0.4
Fenoxycarb (72490-01-8)	-	0.0477	0.382	0.2	Thiacloprid (111988-49-9)	-	0.0477	0.382	0.2
(E)-Fenpyroximate (134098-61-6)	-	0.0477	0.382	0.4	Thiamethoxam (153719-23-4)	-	0.0477	0.382	0.2
Fipronil (120068-37-3)	-	0.0477	0.382	0.4	Trifloxystrobin (141517-21-7)	-	0.0477	0.382	0.2
Flonicamid (158062-67-0)	-	0.0477	0.382	1					
Fludioxinil (131341-86-1)	-	0.0477	0.382	0.4					
Hexythiazox (78587-05-0)	-	0.0477	0.382	1					
Imazalil (35554-44-0)	-	0.0477	0.382	0.2					
Imidacloprid (138261-41-3)	-	0.0477	0.382	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	Systhane	Propiconazole	Tilt
DDVP	Dichlorvos	Naled	Dibrom	Propoxur	Baygon
Ethoprophos	Prophos	Phosmet	Imidan		

REPORT OF LABORATORY ANALYSIS

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Kyle W. Felling, Ph.D.
Laboratory Director



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA43390)

Testing Location:	Customer ID: 2168	Order ID: OR12036	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13249906850	Mass: 3ea
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 03/02/2026
License: ADH 113	License: 00065C	E20260302HLBH01	Date Received: 03/02/2026
Cultivar (Strain) or Sample Description: Hoop Loops Solventless Bubble Hash 1.5g Jar			Date Completed: 03/04/2026

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

Analysis Date/Time: 03/03/2026 1614

Method: ICP/MS

Deviations from SOP:

Instrument: Agilent 7700x

None

Analyst: KF

<u>Heavy Metal</u>	<u>Result</u> <u>(ppb)</u>	<u>LOD</u> <u>(ppb)</u>	<u>LOQ</u> <u>(ppb)</u>	<u>Action</u> <u>Level</u> <u>(ppb)</u>
Arsenic (As)	ND	57.7	91.4	200
Cadmium (Cd)	ND	57.7	91.4	200
Lead (Pb)	ND	57.7	91.4	500
Mercury (Hg)	ND	57.7	91.4	100



Abbreviations: ICP - Inductively-Coupled Plasma, MS - Mass Spectrometry,
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

"ND" not detected above LOD

Reporting limit is 1/2 the action level of each analyte.

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

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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA43390)

Testing Location:	Customer ID: 2168	Sample ID: SA43390	Sample Type: Primary
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-	License: 00065C	E20260302HLBH01	Date Received: 03/02/2026
Cultivar (Strain) or Sample Description: Hoop Loops Solventless Bubble Hash 1.5g Jar			Date Completed: 03/04/2026

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 03/04/2026 0547	Method: Hardy Diagnostics CompactDry	Deviations from SOP:
Analyst: PW	Instrument: Thermo Incubator	None

Bacteria/Microbe	Result (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	-



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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