

SUMMARY OF ANALYSIS (SAMPLE ID: SA35451)

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: OR10497 Lot Number: M00065C13243006276 Production Run: E20230612FSOLBu02	Sample Type: Primary Matrix: Concentrate Mass: 4g Date Collected: 06/19/2023 Date Received: 06/20/2023
Cultivar (Strain) or Sample Description: Fatso Indica Live Badder 1g Jar			Date Completed: 06/21/2023

*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.

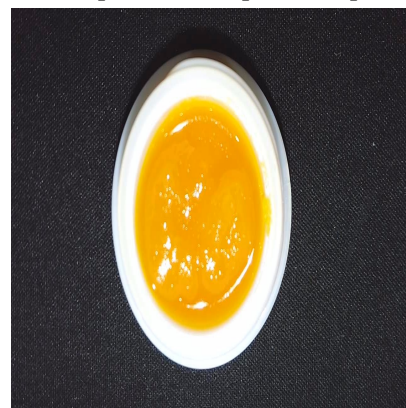
Moisture content levels less than 15% are recommended but the sample does not fail. Water activity levels must be less than 0.65aw.

Cannabinoids (Top 3)	(%)	mg/g
THCA	71.2	712
CBL	1.74	17.4
CBGA	1.73	17.3
TOTAL CBD	0.0908	0.908
TOTAL THC	63.3	633
TOTAL CANNABINOIDS	76.9	769

Terpenes (Top 5)	(%)	µg/g
β-Myrcene	2.55	25500
d-Limonene	0.842	8420
β-Caryophyllene	0.377	3770
Linalool	0.114	1140
β-Ocimene	0.106	1060
TOTAL TERPENES	4.29	42900

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

www.FASTLaboratories.com



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

Analysis Date/Time: 06/20/2023 1436
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	0.107	0.249	-	-	-
CBCA	0.798	7.98	0.331	0.772	-	7.98	7.98
CBD	ND	ND	0.752	1.75	-	-	-
CBDA	0.104	1.04	0.277	0.646	-	1.04	1.04
CBDV	ND	ND	0.121	0.281	-	-	-
CBDVA	ND	ND	0.322	0.751	-	-	-
CBG	<LOQ	<LOQ	0.488	1.14	-	-	-
CBGA	1.73	17.3	0.692	0.816	-	17.3	17.3
CBL	1.74	17.4	0.564	1.32	-	17.4	17.4
CBN	ND	ND	0.259	0.605	-	-	-
CBNA	ND	ND	0.280	0.652	-	-	-
Δ9-THC	0.891	8.91	0.310	0.724	-	8.91	8.91
Δ8-THC	ND	ND	0.484	1.13	-	-	-
THCA	71.2	712	0.168	0.394	-	712	712
THCV	ND	ND	0.404	0.941	-	-	-
THCVA	0.413	4.13	0.129	0.300	-	4.13	4.13
TOTAL	76.9	769			-	769	769
TOTAL CBC	0.700	7.00			-	7.00	7.00
TOTAL CBD	0.0908	0.908			-	0.908	0.908
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	1.52	15.2			-	15.2	15.2
TOTAL CBN	-	-			-	-	-
TOTAL THC	63.3	633			-	633	633
TOTAL THCV	0.358	3.58			-	3.58	3.58



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 = (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) + 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: Fatso Indica Live Badder 1g Jar			Date Completed: 06/21/2023

TERPENOID PROFILE

Analysis Date/Time: 06/21/2023 0143
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	29.4	0.00294	
Camphene	249	0.0249	
δ-3-Carene	20.3	0.00203	
β-Caryophyllene	3770	0.377	█
Caryophyllene oxide	-	-	
p-Cymene	-	-	
Eucalyptol	-	-	
Geraniol	-	-	
Guaiol	17.2	0.00172	
α-Humulene	859	0.0859	█
Isopulegol	-	-	
d-Limonene	8420	0.842	█
Linalool	1140	0.114	█
β-Myrcene	25500	2.55	█
cis-Nerolidol	-	-	
trans-Nerolidol	52.5	0.00525	
α-Ocimene	-	-	
β-Ocimene	1060	0.106	█
α-Pinene	949	0.0949	█
β-Pinene	318	0.0318	█
α-Terpinene	34.7	0.00347	
γ-Terpinene	37.2	0.00372	
Terpinolene	368	0.0368	█
TOTAL	42900	4.29	



Abbreviations: GC - Gas Chromatography, MS - Mass Spectrometry, RL - Reporting Limit

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"-" Not detected above RL.

Reporting Limit (µg/g): 278

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Cultivar (Strain) or Sample Description: Fatso Indica Live Badder 1g Jar **Date Completed:** 06/21/2023

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

Analysis Date/Time: 06/20/2023 1607	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)	-	143	287	5000	n-Heptane (142-82-5)	-	143	287	5000
Acetonitrile (75-5-8)	-	143	287	410	n-Hexane (110-54-3)	-	50.2	100	290
Benzene (71-43-2)	-	1.43	2.87	2	Isobutane (75-28-5)	-	143	287	5000
n-Butane (106-97-2)	-	143	287	5000	Isopropanol (67-63-0)	-	143	287	5000
1-Butanol (71-36-3)	-	143	287	5000	Isopropyl acetate (108-21-4)	-	143	287	5000
2-Butanol (78-92-2)	-	143	287	5000	Isopropyl benzene (98-82-8)	-	14.3	28.7	70
2-Butanone (78-93-3)	-	143	287	5000	Methanol (67-56-1)	-	143	287	3000
Cyclohexane (110-82-7)	-	143	287	3880	2-Methylbutane (78-78-4)	-	143	287	5000
1,2-Dimethoxyethane (110-71-4)	-	14.3	28.7	100	Methylene chloride (75-9-2)	-	143	287	600
N,N-Dimethylacetamide (127-19-5)	-	143	287	1090	2-Methylpentane (107-83-5)	-	50.2	100	290
2,2-Dimethylbutane (75-83-2)	-	50.2	100	290	3-Methylpentane (96-10-0)	-	50.2	100	290
2,3-Dimethylbutane (79-29-8)	-	50.2	100	290	n-Pentane (109-66-0)	-	143	287	5000
N,N-Dimethylformamide (68-12-2)	-	143	287	880	1-Pentanol (71-41-0)	-	143	287	5000
Dimethylsulfoxide (67-68-5)	-	143	287	5000	n-Propane (74-98-6)	-	143	287	5000
1,4-Dioxane (123-91-1)	-	143	287	380	1-Propanol (71-23-8)	-	143	287	5000
Ethanol (64-17-5)	-	143	287	5000	Pyridine (110-86-1)	-	50.2	100	200
2-Ethoxyethanol (110-80-5)	-	50.2	100	160	Tetrahydrofuran (109-99-9)	-	143	287	720
Ethyl ether (60-29-7)	-	143	287	5000	Tetramethylene sulfone (126-33-0)	-	50.2	100	160
Ethyl acetate (141-78-6)	-	143	287	5000	Toluene (108-88-3)	-	143	287	890
Ethyl benzene (100-41-4)	-	143	287	2170	o-Xylene (95-47-6)	-	143	287	2170
Ethylene glycol (107-21-1)	-	143	287	620	m,p-Xylene (108-38-3 or 106-42-3)	-	143	287	2170
Ethylene oxide (75-21-8)	-	14.3	28.7	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170



Color Key

RESULT < AL (Green bar)
RESULT > AL (Red bar)

"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: Fatso Indica Live Badder 1g Jar **Date Completed:** 06/21/2023

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 06/20/2023 1447
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)	-	0.0466	0.373	0.5	Kresoxim-methyl (143390-89-0)	-	0.0466	0.373	0.4
Acephate (30560-19-1)	-	0.0466	0.373	0.4	Malathion (121-75-5)	-	0.0466	0.373	0.2
Acetamiprid (135410-20-7)	-	0.0466	0.373	0.2	Metalaxyl (57837-19-1)	-	0.0466	0.373	0.2
Aldicarb (116-06-3)	-	0.0466	0.373	0.4	Methiocarb (2032-65-7)	-	0.0466	0.373	0.2
Azoxystrobin (131860-33-8)	-	0.0466	0.373	0.2	Methomyl (16752-77-5)	-	0.0466	0.373	0.4
Bifenazate (149877-41-8)	-	0.0466	0.373	0.2	Methyl parathion (298-0-0)	-	0.0466	0.373	0.2
Bifenthrin (82657-04-3)	-	0.0466	0.373	0.2	MGK 264 (113-48-4)	-	0.0466	0.373	0.2
Boscalid (188425-85-6)	-	0.0466	0.373	0.4	Myclobutanil (88671-89-0)	-	0.0466	0.373	0.2
Carbaryl (63-25-2)	-	0.0466	0.373	0.2	Naled (300-76-5)	-	0.0466	0.373	0.5
Carbofuran (1563-66-2)	-	0.0466	0.373	0.2	Oxamyl (23135-22-0)	-	0.0466	0.373	1
Chlorantraniliprole (800008-45-7)	-	0.0466	0.373	0.2	Pacllobutrazol (76738-62-0)	-	0.0466	0.373	0.4
Chlorfenapyr (122453-73-0)	-	0.0466	0.373	1	Permethrins (52645-53-1)	-	0.0466	0.373	0.2
Chlorpyrifos (2921-88-2)	-	0.0466	0.373	0.2	Phosmet (732-11-6)	-	0.0466	0.373	0.2
Clofentezine (74115-24-5)	-	0.0466	0.373	0.2	Piperonyl butoxide (51-03-6)	-	0.0466	0.373	2
Cyfluthrin (68359-37-5)	-	0.0466	0.373	1	Prallethrins (2331-36-9)	-	0.0466	0.373	0.2
Cypermethrin (52315-07-8)	-	0.0466	0.373	1	Propiconazole (60207-90-1)	-	0.0466	0.373	0.4
Daminozide (1596-84-5)	-	0.0466	0.373	1	Propoxur (114-26-1)	-	0.0466	0.373	0.2
DDVP (62-73-7)	-	0.0466	0.373	0.1	Pyrethrins (8003-34-7)	-	0.0466	0.373	1
Diazinon (333-41-5)	-	0.0466	0.373	0.2	Pyridaben (96489-71-3)	-	0.0466	0.373	0.2
Dimethoate (60-51-5)	-	0.0466	0.373	0.2	Spinosad (168316-95-8)	-	0.0466	0.373	0.2
Ethoprophos (13194-48-4)	-	0.0466	0.373	0.2	Spiromesifen (283594-90-1)	-	0.0466	0.373	0.2
Etofenprox (80844-07-1)	-	0.0466	0.373	0.4	Spirotetramat (203313-25-1)	-	0.0466	0.373	0.2
Etoxazole (153233-91-1)	-	0.0466	0.373	0.2	Spiroxamine (118134-30-8)	-	0.0466	0.373	0.4
Fenoxycarb (72490-01-8)	-	0.0466	0.373	0.2	Tebuconazole (80443-41-0)	-	0.0466	0.373	0.4
(E)-Fenpyroximate (134098-61-6)	-	0.0466	0.373	0.4	Thiacloprid (111988-49-9)	-	0.0466	0.373	0.2
Fipronil (120068-37-3)	-	0.0466	0.373	0.4	Thiamethoxam (153719-23-4)	-	0.0466	0.373	0.2
Fonicamid (158062-67-0)	-	0.0466	0.373	1	Trifloxystrobin (141517-21-7)	-	0.0466	0.373	0.2
Fludioxinil (131341-86-1)	-	0.0466	0.373	0.4					
Hexythiazox (78587-05-0)	-	0.0466	0.373	1					
Imazalil (35554-44-0)	-	0.0466	0.373	0.2					
Imidacloprid (138261-41-3)	-	0.0466	0.373	0.4					



Color Key

RESULT < AL (Green bar)
RESULT > AL (Red bar)

"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: 06/20/2023 1704 (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)	-	59.9	94.8	200
Cadmium (Cd)	-	59.9	94.8	200
Lead (Pb)	-	59.9	94.8	500
Mercury (Hg)	-	59.9	94.8	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: SA35451)

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: SA35451 Lot Number: M00065C13243006276 Production Run: E20230612FSOLBu02	Sample Type: Primary Matrix: Concentrate Mass: 4g Date Collected: 06/19/2023 Date Received: 06/20/2023
Cultivar (Strain) or Sample Description: Fatso Indica Live Badder 1g Jar			Date Completed: 06/21/2023

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 06/21/2023 0953 **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.

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

