F.A.S.T. LABORATORIES/RESEARCH

ACCREDITED CERTIFICATE #4473.01

ASA PFC Laboratory

SUMMARY OF ANALYSIS (SAMPLE ID: SA35960)

Testing Location:	**Customer ID:** 2097	**Order ID:** OR10598	**Sample Type:** Primary
Arkansas	Pocono Pharmaceutical	**Lot Number:**	**Matrix:** Patch
232 S. Broadview St.	100 Sweetree St	2111025	**Mass:** 2cnt
Greenbrier, AR 72058	Cherryville , NC 28021	**Batch Number:**	**Date Collected:** 08/31/2023
License: ADA 05_H273	License: Not Entered or N/A	Not Entered	**Date Received:** 09/01/2023
Cultivar (Strain) or Sample Description: Arm Patch			**Date Completed:** 09/11/2023

*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.

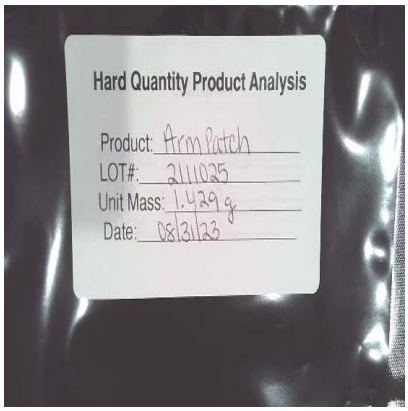
Moisture Content (%)	Water Activity (aw)	PASS/FAIL
Not Tested	Not Tested	PASS

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
CBD	1.90	19.0
CBDV	0.00185	0.0185
CBDA		-
TOTAL CBD	1.90	19.0
TOTAL THC	-	-
TOTAL CANNABINOIDS	1.90	19.0

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, Ph.D.
Laboratory Director

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F.A.S.T. LABORATORIES/RESEARCH

A2LA ACCREDITED CERTIFICATE #4473.01

ASA PFC Laboratory

CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35960)

Testing Location:
Arkansas
232 S. Broadview St.
Greenbrier, AR 72058
License: ADA 05_H273

Customer ID: 2097
Pocono Pharmaceutical
100 Sweetree St
Cherryville , NC 28021
License: Not Entered or N/A

Order ID: OR10598
Lot Number:
2111025
Batch Number:
Not Entered

Sample Type: Primary
Matrix: Patch
Mass: 2cnt
Date Collected: 08/31/2023
Date Received: 09/01/2023

Cultivar (Strain) or Sample Description: Arm Patch

Date Completed: 09/11/2023

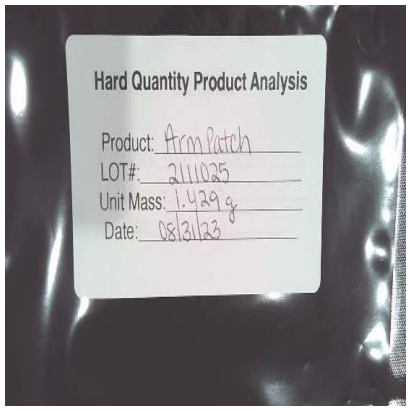
CANNABINOID (POTENCY) PROFILE (SOP: SOP-CANN-001)

Analysis Date/Time: 09/06/2023 1418
Analyst: PW

Method: HPLC/DAD
Instrument: Agilent 1100

Moisture Content (%): -
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	0.00247	0.00576	-	-	-
CBCA	ND	ND	0.00766	0.0179	-	-	-
CBD	1.90	19.0	0.0174	0.0406	-	36.1	36.1
CBDA	ND	ND	0.00641	0.0150	-	-	-
CBDV	0.00185	0.0185	0.00279	0.00652	-	0.0352	0.0352
CBDVA	ND	ND	0.00745	0.0174	-	-	-
CBG	ND	ND	0.0113	0.0264	-	-	-
CBGA	ND	ND	0.0160	0.0189	-	-	-
CBL	ND	ND	0.0131	0.0305	-	-	-
CBN	ND	ND	0.00600	0.0140	-	-	-
CBNA	ND	ND	0.00647	0.0151	-	-	-
Δ9-THC	ND	ND	0.00719	0.0168	-	-	-
Δ8-THC	ND	ND	0.0112	0.0262	-	-	-
THCA	ND	ND	0.00390	0.00911	-	-	-
THCV	ND	ND	0.00935	0.0218	-	-	-
THCVA	ND	ND	0.00299	0.00695	-	-	-
TOTAL	1.90	19.0			-	36.1	36.1
TOTAL CBC	-	-			-	-	-
TOTAL CBD	1.90	19.0			-	36.1	36.1
TOTAL CBDV	0.00185	0.0185			-	0.0352	0.0352
TOTAL CBG	-	-			-	-	-
TOTAL CBN	-	-			-	-	-
TOTAL THC	-	-			-	-	-
TOTAL THCV	-	-			-	-	-



SERVING MASS (g): 1.90
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) + 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, UM - Measurement Uncertainty

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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35960)

Testing Location:	**Customer ID:** 2097	**Order ID:** OR10598	**Sample Type:** Primary
Arkansas	Pocono Pharmaceutical	**Lot Number:**	**Matrix:** Patch
232 S. Broadview St.	100 Sweetree St	2111025	**Mass:** 2cnt
Greenbrier, AR 72058	Cherryville , NC 28021	**Batch Number:**	**Date Collected:** 08/31/2023
License: ADA 05_H273	License: Not Entered or N/A	Not Entered	**Date Received:** 09/01/2023
Cultivar (Strain) or Sample Description: Arm Patch			**Date Completed:** 09/11/2023

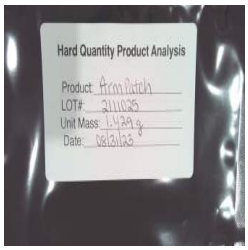
RESIDUAL SOLVENT PROFILE (SOP: SOP-RS-001)

Analysis Date/Time: 09/07/2023 0004 **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)
Acetone (67-64-1)	-	139	278	5000
Acetonitrile (75-5-8)	-	139	278	410
Benzene (71-43-2)	-	1.39	2.78	2
n-Butane (106-97-2)	-	139	278	5000
1-Butanol (71-36-3)	-	139	278	5000
2-Butanol (78-92-2)	-	139	278	5000
2-Butanone (78-93-3)	-	139	278	5000
Cyclohexane (110-82-7)	-	139	278	3880
1,2-Dimethoxyethane (110-71-4)	-	13.9	27.8	100
N,N-Dimethylacetamide (127-19-5)	-	139	278	1090
2,2-Dimethylbutane (75-83-2)	-	48.6	97.2	290
2,3-Dimethylbutane (79-29-8)	-	48.6	97.2	290
N,N-Dimethylformamide (68-12-2)	-	139	278	880
Dimethylsulfoxide (67-68-5)	-	139	278	5000
1,4-Dioxane (123-91-1)	-	139	278	380
Ethanol (64-17-5)	-	139	278	5000
2-Ethoxyethanol (110-80-5)	-	48.6	97.2	160
Ethyl ether (60-29-7)	-	139	278	5000
Ethyl acetate (141-78-6)	-	139	278	5000
Ethyl benzene (100-41-4)	-	139	278	2170
Ethylene glycol (107-21-1)	-	139	278	620
Ethylene oxide (75-21-8)	-	13.9	27.8	50

Solvent	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
n-Heptane (142-82-5)	-	139	278	5000
n-Hexane (110-54-3)	-	48.6	97.2	290
Isobutane (75-28-5)	-	139	278	5000
Isopropanol (67-63-0)	-	139	278	5000
Isopropyl acetate (108-21-4)	-	139	278	5000
Isopropyl benzene (98-82-8)	-	13.9	27.8	70
Methanol (67-56-1)	-	139	278	3000
2-Methylbutane (78-78-4)	-	139	278	5000
Methylene chloride (75-9-2)	-	139	278	600
2-Methylpentane (107-83-5)	-	48.6	97.2	290
3-Methylpentane (96-10-0)	-	48.6	97.2	290
n-Pentane (109-66-0)	-	139	278	5000
1-Pentanol (71-41-0)	-	139	278	5000
n-Propane (74-98-6)	-	139	278	5000
1-Propanol (71-23-8)	-	139	278	5000
Pyridine (110-86-1)	-	48.6	97.2	200
Tetrahydrofuran (109-99-9)	-	139	278	720
Tetramethylene sulfone (126-33-0)	-	48.6	97.2	160
Toluene (108-88-3)	-	139	278	890
o-Xylene (95-47-6)	-	139	278	2170
m,p-Xylene (108-38-3 or 106-42-3)	-	139	278	2170
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
Dimethysufoxide	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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REPORT OF LABORATORY ANALYSIS

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Arkansas	Pocono Pharmaceutical	**Lot Number:**	**Matrix:** Patch
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Cultivar (Strain) or Sample Description: Arm Patch — **Date Completed:** 09/11/2023

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 09/06/2023 1355 — **Method:** LC/MS/MS — **Deviations from SOP:**

Analyst: KF — **Instrument:** Shimadzu LC-8050 — None

Pesticide	Result (µg/g)	LOD (µg/g)	LOQ (µg/g)	Action Level (µg/g)
Abamectin (71751-41-2)	-	0.0108	0.0865	0.5
Acephate (30560-19-1)	-	0.0108	0.0865	0.4
Acequinocyl (57960-19-7)	-	0.0108	0.0865	2
Acetamiprid (135410-20-7)	-	0.0108	0.0865	0.2
Aldicarb (116-06-3)	-	0.0108	0.0865	0.4
Azoxystrobin (131860-33-8)	-	0.0108	0.0865	0.2
Bifenazate (149877-41-8)	-	0.0108	0.0865	0.2
Bifenthrin (82657-04-3)	-	0.0108	0.0865	0.2
Boscalid (188425-85-6)	-	0.0108	0.0865	0.4
Carbaryl (63-25-2)	-	0.0108	0.0865	0.2
Carbofuran (1563-66-2)	-	0.0108	0.0865	0.2
Chlorantraniliprole (800008-45-7)	-	0.0108	0.0865	0.2
Chlorfenapyr (122453-73-0)	-	0.0108	0.0865	1
Chlorpyrifos (2921-88-2)	-	0.0108	0.0865	0.2
Clofentezine (74115-24-5)	-	0.0108	0.0865	0.2
Cyfluthrin (68359-37-5)	-	0.0108	0.0865	1
Cypermethrin (52315-07-8)	-	0.0108	0.0865	1
Daminozide (1596-84-5)	-	0.0108	0.0865	1
DDVP (62-73-7)	-	0.0108	0.0865	0.1
Diazinon (333-41-5)	-	0.0108	0.0865	0.2
Dimethoate (60-51-5)	-	0.0108	0.0865	0.2
Ethoprophos (13194-48-4)	-	0.0108	0.0865	0.2
Etofenprox (80844-07-1)	-	0.0108	0.0865	0.4
Etoxazole (153233-91-1)	-	0.0108	0.0865	0.2
Fenoxycarb (72490-01-8)	-	0.0108	0.0865	0.2
(E)-Fenpyroximate (134098-61-6)	-	0.0108	0.0865	0.4
Fipronil (120068-37-3)	-	0.0108	0.0865	0.4
Flonicamid (158062-67-0)	-	0.0108	0.0865	1
Fludioxinil (131341-86-1)	-	0.0108	0.0865	0.4
Hexythiazox (78587-05-0)	-	0.0108	0.0865	1
Imazalil (35554-44-0)	-	0.0108	0.0865	0.2
Imidacloprid (138261-41-3)	-	0.0108	0.0865	0.4
Kresoxim-methyl (143390-89-0)	-	0.0108	0.0865	0.4
Malathion (121-75-5)	-	0.0108	0.0865	0.2
Metalaxyl (57837-19-1)	-	0.0108	0.0865	0.2
Methiocarb (2032-65-7)	-	0.0108	0.0865	0.2
Methomyl (16752-77-5)	-	0.0108	0.0865	0.4
Methyl parathion (298-0-0)	-	0.0108	0.0865	0.2
MGK 264 (113-48-4)	-	0.0108	0.0865	0.2
Myclobutanil (88671-89-0)	-	0.0108	0.0865	0.2
Naled (300-76-5)	-	0.0108	0.0865	0.5
Oxamyl (23135-22-0)	-	0.0108	0.0865	1
Paclobutrazol (76738-62-0)	-	0.0108	0.0865	0.4
Permethrins (52645-53-1)	-	0.0108	0.0865	0.2
Phosmet (732-11-6)	-	0.0108	0.0865	0.2
Piperonyl butoxide (51-03-6)	-	0.0108	0.0865	2
Prallethrins (2331-36-9)	-	0.0108	0.0865	0.2
Propiconazole (60207-90-1))	-	0.0108	0.0865	0.4
Propoxur (114-26-1)	-	0.0108	0.0865	0.2
Pyrethrins (8003-34-7)	-	0.0108	0.0865	1
Pyridaben (96489-71-3)	-	0.0108	0.0865	0.2
Spinosad (168316-95-8)	-	0.0108	0.0865	0.2
Spiromesifen (283594-90-1)	-	0.0108	0.0865	0.2
Spirotetramat (203313-25-1)	-	0.0108	0.0865	0.2
Spiroxamine (118134-30-8)	-	0.0108	0.0865	0.4
Tebuconazole (80443-41-0)	-	0.0108	0.0865	0.4
Thiacloprid (111988-49-9)	-	0.0108	0.0865	0.2
Thiamethoxam (153719-23-4)	-	0.0108	0.0865	0.2
Trifloxystrobin (141517-21-7)	-	0.0108	0.0865	0.2

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.

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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		

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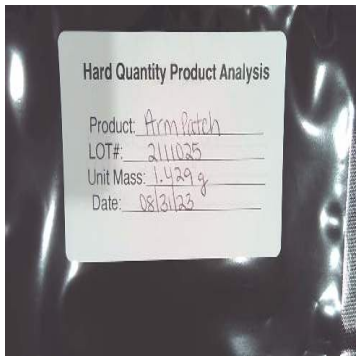
HEAVY METAL PROFILE (SOP: SOP-ICP-200.6)

Analysis Date/Time: 09/12/2023 1948 (ICP/OES)
Analysis Date/Time: - (DMA)
Analyst: KF

Method: ICP/MS
Instrument: Agilent 7500ce

Deviations from SOP:
None

Heavy Metal	Result (µg/kg)	LOD (µg/kg)	LOQ (µg/kg)	Action Level (µg/kg)
Arsenic (As)	-	55.9	88.6	200
Cadmium (Cd)	-	55.9	88.6	200
Lead (Pb)	DET	55.9	88.6	500
Mercury (Hg)	-	55.9	88.6	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.

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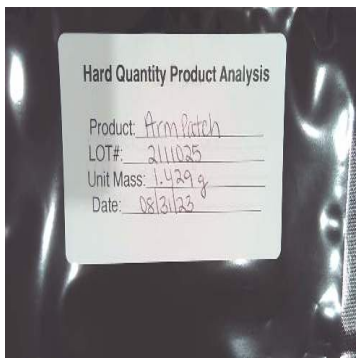
MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 09/07/2023 0935
Analyst: PW

Method: Hardy Diagnostics CompactDry
Instrument: Thermo Incubator

Deviations from SOP:
None

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



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.

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