







SUMMARY OF ANALYSIS (SAMPLE ID: SA35290)

Order ID: OR10467 **Testing Location:** Customer ID: 2161 Sample Type: Primary Arkansas River Valley Relief Cultivation Lot Number: Matrix: Flower 232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13130697173 Mass: 22g

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 05/23/2023 5033H20230424RSR-2 License: ADH 113 License: 00065C **Date Received:** 05/23/2023

Cultivar (Strain) or Sample Description: Ryleigh's Reserve **Date Completed:** 05/25/2023

*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

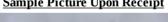
> PASS/FAIL **Moisture Content (%)** Water Activity (aw) 0.465 @ 24.5°C **PASS** 13.7

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.

Cannabinoi	ds (Top 3)	<u>(%)</u>	mg/g
THCA		26.6	266
CBC	GA	0.780	7.80
Δ9-Τ	НС	0.324	3.24
TOTAL	CBD	-	-
TOTAL	THC	23.7	237
TOTAL CAN	NABINOIDS	28.0	280
<u>Terpenes</u>	(Top 5)	<u>(%)</u>	µg/g
d-Limo	onene	0.469	4690
β-Myr	cene	0.240	2400
β-Ocimene		0.217	2170
β-Caryophyllene		0.204	2040
Linalool		0.188	1880
TOTAL TERPENES		1.68	16800
Contaminants PASS/FAIL		Sample Pictur	re Upon Receipt

Contaminants	PASS/FAIL
Heavy Metals:	PASS
Microbiology:	PASS
Pesticides:	PASS







Scan the QR code to verify results.

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The statements and results herein have not been approved and/or endorsed by the FDA.







License: ADH 113







CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35290)

Testing Location: Customer ID: 2161 Arkansas River Valley Relief Cultivation 232 S. Broadview St. 5601 Old Greenwood Rd Suite C Greenbrier, AR 72058 Fort Smith, AR 72903

Cultivar (Strain) or Sample Description: Ryleigh's Reserve

License: 00065C

Order ID: OR10467 Lot Number: M00065C13130697173

Production Run: 5033H20230424RSR-2 Sample Type: Primary Matrix: Flower

Mass: 22g

Date Collected: 05/23/2023 **Date Received:** 05/23/2023

Date Completed: 05/25/2023

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

Analysis Date/Time: 05/23/2023 1705 Method: HPLC/DAD Moisture Content (%): 13.7

Analyst: PW Instrument: Agilent 1100 **Water Activity (aw):** 0.465 @ 24.5°C

Cannabinoid	Result (%)	Result (mg/g)	LOD (mg/g)	LOQ (mg/g)	Result (mg/ mL)	Per Serving (mg)	Per Unit (mg)
CBC	ND	ND	0.0390	0.0910	-	-	-
CBCA	ND	ND	0.121	0.283	-	-	-
CBD	ND	ND	0.275	0.642	-	-	-
CBDA	ND	ND	0.101	0.236	-	-	-
CBDV	ND	ND	0.0441	0.103	-	-	-
CBDVA	ND	ND	0.118	0.275	-	-	-
CBG	0.145	1.45	0.179	0.417	-	1.45	1.45
CBGA	0.780	7.80	0.253	0.299	-	7.80	7.80
CBL	ND	ND	0.206	0.482	-	-	-
CBN	ND	ND	0.0947	0.221	-	-	-
CBNA	ND	ND	0.102	0.238	-	-	-
Δ9-ΤΗС	0.324	3.24	0.114	0.265	-	3.24	3.24
Δ8-ΤΗС	ND	ND	0.177	0.414	-	-	-
THCA	26.6	266	0.0616	0.144	-	266	266
THCV	ND	ND	0.148	0.344	-	-	-
THCVA	0.156	1.56	0.0472	0.110	-	1.56	1.56
TOTAL	28.0	280				280	280
TOTAL CBC	-	-		•	-	-	-
TOTAL CBD	-	-			-	-	-
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	0.830	8.30			-	8.30	8.30
TOTAL CBN	-	-			-	-	-
TOTAL THC	23.7	237			-	237	237
TOTAL THCV	0.135	1.35			-	1.35	1.35

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

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 = $(CBCA \times 0.877) + CBC$ Total CBD = $(CBDA \times 0.877) + CBD$ Total CBDV = (CBDVA \times 0.867) + CBDV Total CBG = $(CBGA \times 0.878) + CBG$ Total CBN = $(CBNA \times 0.876) + CBN$ Total THC = (THCA x 0.877) + $\Delta 9$ -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.





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[&]quot;-" Not detected above LOD.



License: ADH 113







CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35290)

Testing Location: Customer ID: 2161
Arkansas River Valley Relief Cultivation
232 S. Broadview St. 5601 Old Greenwood Rd Suite C
Greenbrier, AR 72058 Fort Smith, AR 72903

Cultivar (Strain) or Sample Description: Ryleigh's Reserve

License: 00065C

Order ID: OR10467 **Lot Number:** M00065C13130697173

Production Run: 5033H20230424RSR-2

Sample Type: Primary **Matrix:** Flower

Mass: 22g

Date Collected: 05/23/2023 **Date Received:** 05/23/2023

Date Completed: 05/25/2023

TERPENOID PROFILE

Analysis Date/Time:05/24/2023 0138Method: GC/MSDeviations from SOP:Analyst: KFInstrument: Agilent 7890/5975None

-			
<u>Terpene</u>	<u>Result</u> (μg/g)	Result (%)	
α-Bisabolol	59.1	0.00591	
Camphene	192	0.0192	
δ-3-Carene	11.2	0.00112	
β-Caryophyllene	2040	0.204	
Caryophyllene oxide	-	-	
p-Cymene	13.9	0.00139	
Eucalyptol	-	-	
Geraniol	-	-	
Guaiol	-	-	
α-Humulene	552	0.0552	
Isopulegol	-	-	
d-Limonene	4690	0.469	
Linalool	1880	0.188	
β-Myrcene	2400	0.240	
cis-Nerolidol	125	0.0125	
trans-Nerolidol	947	0.0947	
α-Ocimene	-	-	
β-Ocimene	2170	0.217	
α-Pinene	659	0.0659	
β-Pinene	936	0.0936	
α-Terpinene	9.89	0.000989	
γ-Terpinene	4.13	0.000413	
Terpinolene	122	0.0122	
TOTAL	16800	1.68	



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

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"-" Not detected above RL

TOTAL 16800 1.68 Reporting Limit (μg/g): 49.2





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05/25/2023









CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35290)

Testing Location: Arkansas

232 S. Broadview St. Greenbrier, AR 72058

License: ADH 113 Cultivar (Strain) or Sample Description: Ryleigh's Reserve

Customer ID: 2161

River Valley Relief Cultivation 5601 Old Greenwood Rd Suite C

Fort Smith, AR 72903 License: 00065C

Lot Number: M00065C13130697173

Order ID: OR10467

Production Run: 5033H20230424RSR-2 Sample Type: Primary

Matrix: Flower Mass: 22g

Date Collected: 05/23/2023 **Date Received:** 05/23/2023

Date Completed: 05/25/2023

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 05/24/2023 1032

Analyst: KF

Method: LC/MS/MS

Instrument: Shimadzu LC-8050

Deviations from SOP:

None

Acephate (30560-19-1)	on (121-75-5) 0.0427 or (121-75-5) 0.0427 or (2032-65-7) 0.0427 or (16752-77-5) 0.0427 or (113-48-4) 0.0427	0.341 0.341 0.341 0.341 0.341 0.341 0.341 0.341 0.341 0.341	0.4 0.2 0.2 0.2 0.4 0.2 0.2 0.5 1 0.4
Acequinocyl (57960-19-7)	on (121-75-5) 0.0427 vl (57837-19-1) 0.0427 arb (2032-65-7) 0.0427 vl (16752-77-5) 0.0427 varathion 0.0427 ttanil 39-0) 0.0427 varathion 0.0427 ctanil 0.0427	0.341 0.341 0.341 0.341 0.341 0.341 0.341 0.341 0.341 0.341	0.2 0.2 0.4 0.2 0.2 0.2 0.5 1
Acetamicyl (3730-137) Acetamicyl (135410-20-7) Aldicarb (116-06-3) Azoxystrobin (131860-33-8) Bifenazate (149877-41-8) Bifenthrin (82657-04-3) Boscalid (188425-85-6) Carbaryl (63-25-2) Chlorentraniliprole (80008-45-7) Chlorenapyr (122453-73-0) Chlorpyrifos (2921-88-2) Clofentezine (74115-24-5) Cypermethrin (52315-07-8) Daminozide (1596-84-5) DDVP (62-73-7) Diazinon (333-41-5) Dimethoate (60-51-5) Metalax: Metalax: Methom Metalax: Methoic Met	At (57837-19-1) 0.0427 arb (2032-65-7) 0.0427 yl (16752-77-5) 0.0427 vl (113-48-4) 0.0427 attanil 0.0427 (23135-22-0) 0.0427 razol 0.0427 ins (52645-53-1) 0.0427	0.341 0.341 0.341 0.341 0.341 0.341 0.341 0.341 0.341	0.2 0.2 0.4 0.2 0.2 0.2 0.5 1
Acteaning Transport (135410-20-7) Aldicarb (116-06-3) Azoxystrobin (131860-33-8) Bifenazate (149877-41-8) Bifenazate (149877-41-8) Boscalid (188425-85-6) Carbaryl (63-25-2) Chlorantraniliprole (80008-45-7) Chlorapyr (122453-73-0) Chloryprifos (2921-88-2) Clofentezine (74115-24-5) Cypermethrin (52315-07-8) Daminozide (1596-84-5) DDVP (62-73-7) Diazinon (333-41-5) Dimethoate (60-51-5) Methiod O.0427 O.341 O.2 Methyl (228-0-C Methyl (228-0-C Methor O.0427 O.341 O.2 Methyl (228-0-C Methyl (228-0-C Methyl (228-0-C Methyl (228-0-C Methor O.0427 O.341 O.2 Methor Methom O.0427 O.341 O.2 Methyl (298-0-C O.0427 O.341 O.2 Methyl Methom O.0427 O.341 O.2 Methyl Methor Methor O.0427 O.341 O.2 Methyl Methyl O.2 Methyl O.0427 O.341 O.2 Methyl O.2 Methyl O.2 Methyl O.2 Methyl O.2 Methyl O.2 Methyl O.2 Me	arb (2032-65-7)	0.341 0.341 0.341 0.341 0.341 0.341 0.341 0.341	0.2 0.4 0.2 0.2 0.2 0.5 1
Azoxystrobin (131860-33-8) Bifenazate (149877-41-8)	Arrathion 0.0427 4 (113-48-4) 0.0427 Atanil 0.0427 00-76-5) 0.0427 (23135-22-0) 0.0427 razol 0.0427 rins (52645-53-1) 0.0427	0.341 0.341 0.341 0.341 0.341 0.341	0.2 0.2 0.2 0.5 1
Carborura (1563-66-2) Carbofuran (1563-66-2) Carborura (152453-73-0) Chlorapyri (122453-73-0) Chlorapyri (63291-88-2) Cultura (74115-24-5) Cultura (68359-37-5) Cultura (1563-68-4) Cultura (1563-68-2) Cultura (1563-67-1) Chlorapyri (1563-67-1) Chlo	0.0427 4 (113-48-4) 0.0427 101111 0.0427 00-76-5) 0.0427 (23135-22-0) 0.0427 10201 0.0427 10311 0.0427 10311 0.0427 10311 0.0427 10311 0.0427	0.341 0.341 0.341 0.341 0.341	0.2 0.2 0.5 1 0.4
Bifenthrin (82657-04-3)	tanil 39-0) - 0.0427 00-76-5) - 0.0427 (23135-22-0) - 0.0427 razol 52-0) - 0.0427 ins (52645-53-1) - 0.0427	0.341 0.341 0.341 0.341 0.341	0.2 0.5 1 0.4
Boscalid (188425-85-6) - 0.0427 0.341 0.4 (88671-4	39-0) 0.0427 00-76-5) 0.0427 (23135-22-0) 0.0427 razol 0.0427 ins (52645-53-1) 0.0427	0.341 0.341 0.341 0.341	0.5 1 0.4
Carbaryl (63-25-2) - 0.0427 0.341 0.2 Naled (3 Carbofuran (1563-66-2) - 0.0427 0.341 0.2 Oxamyl Paclobut (800008-45-7) - 0.0427 0.341 0.2 (76738-160008-45-7) - 0.0427 0.341 0.2 Piperony (122453-73-0) - 0.0427 0.341 0.2 Piperony (122453-73-0) - 0.0427 0.341 0.2 Piperony (12453-73-6) - 0.0427 0.341 0.2 Piperony (12453-73-6) - 0.0427 0.341 1 Prallethr (68359-37-5) - 0.0427 0.341 1 Prallethr (52315-07-8) - 0.0427 0.341 1 Propoxu (60207-4600000000000000000000000000000000000	(23135-22-0) 0.0427 (23135-22-0) 0.0427 razol 0.0427 rins (52645-53-1) 0.0427	0.341 0.341 0.341 0.341	0.5 1 0.4
Carbofuran (1563-66-2) - 0.0427 0.341 0.2 Oxamyl Paclobut (800008-45-7) - 0.0427 0.341 0.2 (76738-1800008-45-7) - 0.0427 0.341 0.2 Piperony (122453-73-0) - 0.0427 0.341 0.2 Piperony Clofentezine (74115-24-5) - 0.0427 0.341 0.2 Piperony Clofentezine (74115-24-5) - 0.0427 0.341 1 Prallethrin (68359-37-5) - 0.0427 0.341 1 Prallethrin (52315-07-8) - 0.0427 0.341 1 Propoxu (60207-5) DDVP (62-73-7) - 0.0427 0.341 1 Propoxu DDVP (62-73-7) - 0.0427 0.341 1 Propoxu DDVP (62-73-7) - 0.0427 0.341 0.1 Pyrethrin Diazinon (333-41-5) - 0.0427 0.341 0.2 Spinosac	(23135-22-0) 0.0427 razol 0.0427 62-0) 0.0427 ins (52645-53-1) 0.0427	0.341 0.341 0.341	1 0.4
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Cypermethrin (52315-07-8) - 0.0427 0.341 1 Propicor (60207-9	ins (2331-36-9) 0.0427	0.341	0.2
0.0427 0.341 1 (60207-160207-	` '	0.541	
DDVP (62-73-7) - 0.0427 0.341 0.1 Pyrethring Diazinon (333-41-5) - 0.0427 0.341 0.2 Pyridabe Dimethoate (60-51-5) - 0.0427 0.341 0.2 Spinosac	90-1))		0.4
Diazinon (333-41-5) Dimethoate (60-51-5) Dimethoate (60-51-5) Dimethoate (60-51-5) Dimethoate (60-51-5) Dimethoate (60-51-5)	r (114-26-1) - 0.0427		0.2
Dimethoate (60-51-5) - 0.0427 0.341 0.2 Spinosac	ns (8003-34-7) 0.0427		1
0.0427 0.341 0.2	en (96489-71-3) 0.0427		0.2
	1 (168316-95-8)	0.341	0.2
Ethoprophos (13194-48-4) 0.0427 0.341 0.2 Spirome (283594		0.341	0.2
Etofenprox (80844-07-1) - 0.0427 0.341 0.4 Spiroteti	· ·		
Etoxazole (153233-91-1) 0.0427 0.341 0.2 (203313	0.0427	0.341	0.2
Fenoxycarb (72490-01-8) 0.0427 0.341 0.2 Spiroxar (118134		0.341	0.4
(134098-61-6) 0.0427 0.341 0.4 Tebucon	azole 0.0427	0.341	0.4
Fipronil (120068-37-3) 0.0427 0.341 0.4 (80443-4	1 1-0)	0.541	0.4
Flonicamid (158062-67-0) 0.0427 0.341 1 Thiaclog		0.341	0.2
Fludioxinil (131341-86-1) 0.0427 0.341 0.4 (111988	*		
Hexythiazox (78587-05-0) 0.0427 0.341 1 Thiamet (153719	- 0.0427	0.341	0.2
Imazalil (35554-44-0) 0.0427 0.341 0.2 Trifloxy	-23-4)		
Imidacloprid (138261-41-3) - 0.0427 0.341 0.4 (141517	-23-4)	0.341	0.2



Color Key

"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

<u>Pesticide</u>	Synonym(s)	<u>Pesticide</u>	Synonym(s)	<u>Pesticide</u>	Synonym(s)
Cyfluthrin	Baythroid	Myclobutanil	Systhane	Propiconazole	Tilt
DDVP	Dichlorvos	Naled	Dibrom	Propoxur	Baygon
Ethoprophos	Prophos	Phosmet	Imidan		











M00065C13130697173



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35290)

Order ID: OR10467 **Testing Location:** Customer ID: 2161 Sample Type: Primary Arkansas River Valley Relief Cultivation Lot Number: Matrix: Flower

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 05/23/2023 5033H20230424RSR-2 License: ADH 113 License: 00065C **Date Received:** 05/23/2023 **Date Completed:** 05/25/2023

Cultivar (Strain) or Sample Description: Ryleigh's Reserve

5601 Old Greenwood Rd Suite C

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

Analysis Date/Time: 05/23/2023 1712 (ICP/OES) Method: ICP/MS **Deviations from SOP:**

Analysis Date/Time: - (DMA) **Instrument:** Agilent 7500ce None

Analyst: KF

232 S. Broadview St.

Heavy Metal	<u>Result</u> (μg/kg)	<u>LOD</u> (µg/kg)	<u>LOQ</u> (μg/kg)	<u>Action Level</u> (μg/kg)
Arsenic (As)	-	56.9	90.0	200
Cadmium (Cd)	-	56.9	90.0	200
Lead (Pb)	-	56.9	90.0	500
Mercury (Hg)	-	56.9	90.0	100



Mass: 22g

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 < AI RESULT > AI

"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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CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35290)

Testing Location:Customer ID: 2161Sample ID: SA35290Sample Type: PrimaryArkansasRiver Valley Relief CultivationLot Number:Matrix: Flower232 S. Broadview St.5601 Old Greenwood Rd Suite CM00065C13130697173Mass: 22g

 Greenbrier, AR 72058
 Fort Smith, AR 72903
 Production Run:
 Date Collected: 05/23/2023

 License: ADA 05_H273
 License: 00065C
 5033H20230424RSR-2
 Date Received: 05/23/2023

Cultivar (Strain) or Sample Description: Ryleigh's Reserve **Date Completed:** 05/25/2023

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 05/24/2023 1130 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)

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.



