







Date Received: 11/13/2023

SUMMARY OF ANALYSIS (SAMPLE ID: SA36425)

Testing Location:Customer ID: 2168Order ID: OR10701Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Concentrate232 S. Broadview St.5601 Old Greenwood Rd Suite CM00065C13244936466Mass: 4grams

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 11/13/2023

Cultivar (Strain) or Sample Description: Truffle Shuffle Solventless Live Hash Rosin **Date Completed:** 11/15/2023

*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/FAILNot TestedNot TestedPASS

E20231110TFSLHR01

Moisture content/water activity action levels are referenced from the State of Arkansas MMJ testing guidelines.

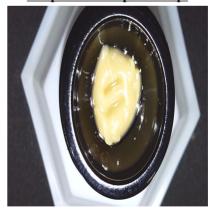
License: 00065C

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

Cannabing	oids (Top 3)	<u>(%)</u>	mg/g
TH	CA	75.7	757
СВ	GA	1.26	12.6
Δ9-7	ГНС	0.862	8.62
TOTA	L CBD	-	-
TOTA	LTHC	67.3	673
TOTAL CAN	NABINOIDS	78.8	788
<u>Terpene</u>	s (Top 5)	<u>(%)</u>	<u>µg/g</u>
d-Lim	onene	3.09	30900
β-Caryo	phyllene	0.996	9960
β-Pi	nene	0.673	6730
α-Pi	nene	0.658	6580
α-Hun	nulene	0.373	3730
TOTAL T	ERPENES	6.65	66500
Contaminants	PASS/FAIL	Sample Pictur	re Upon Receipt

Contaminants	PASS/FAIL
Heavy Metals:	PASS
Microbiology:	PASS
Pesticides:	PASS
Pesidual Solvents:	DASS







Scan the QR code to verify results.

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.

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











E20231110TFSLHR01



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36425)

Testing Location:Customer ID: 2168Order ID: OR10701Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Concentrate232 S. Broadview St.5601 Old Greenwood Rd Suite CM00065C13244936466Mass: 4grams

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run:** Date Collected: 11/13/2023

Cultivar (Strain) or Sample Description: Truffle Shuffle Solventless Live Hash Rosin

Date Completed: 11/15/2023

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

Analysis Date/Time: 11/14/2023 1605 Method: HPLC/DAD
Analyst: PW Instrument: Agilent 1100

License: 00065C

<u>Cannabinoid</u>	Result (%)	Result (mg/g)	LOD (mg/g)	LOQ (mg/g)	Result (mg/ mL)	<u>Per</u> Serving (mg)	Per Unit (mg)
CBC	ND	ND	0.0996	0.232	-	-	-
CBCA	ND	ND	0.309	0.722	-	-	-
CBD	ND	ND	0.702	1.64	-	-	-
CBDA	ND	ND	0.259	0.604	-	-	-
CBDV	ND	ND	0.113	0.263	-	-	-
CBDVA	ND	ND	0.301	0.701	-	-	-
CBG	0.339	3.39	0.456	1.06	-	3.39	3.39
CBGA	1.26	12.6	0.646	0.763	-	12.6	12.6
CBL	ND	ND	0.527	1.23	-	-	-
CBN	ND	ND	0.242	0.565	-	-	-
CBNA	ND	ND	0.261	0.609	-	-	-
Δ9-ΤΗС	0.862	8.62	0.290	0.676	-	8.62	8.62
Δ8-ΤΗС	ND	ND	0.453	1.06	-	-	-
THCA	75.7	757	0.157	0.368	-	757	757
THCV	ND	ND	0.377	0.880	-	-	-
THCVA	0.610	6.10	0.121	0.280	-	6.10	6.10
TOTAL	78.8	788				788	788
TOTAL CBC	-					-	
TOTAL CBD	-	-			-	-	-
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	1.45	14.5			-	14.5	14.5
TOTAL CBN	-	-			-	-	-
TOTAL THC	67.3	673			-	673	673
TOTAL THCV	0.529	5.29			-	5.29	5.29

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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Moisture Content (%): -Water Activity (aw): -



Date Received: 11/13/2023

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) + Δ 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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11/15/2023









CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36425)

Customer ID: 2168 Order ID: OR10701 Sample Type: Primary **Testing Location:** Arkansas River Valley Relief MIPS Lot Number: Matrix: Concentrate 232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13244936466 Mass: 4grams

Fort Smith, AR 72903 **Date Collected:** 11/13/2023 Greenbrier, AR 72058 **Production Run:** License: 00065C E20231110TFSLHR01 **Date Received:** 11/13/2023

Cultivar (Strain) or Sample Description: Truffle Shuffle Solventless Live Hash Rosin **Date Completed:** 11/15/2023

TERPENOID PROFILE

Analysis Date/Time:11/15/2023 0207 Method: GC/MS **Deviations from SOP: Analyst:** KF Instrument: Agilent 7890/5975 None

Terpene Result (μg/g) Result (½) α-Bisabolol 961 0.0961 Camphene 903 0.0903 δ-3-Carene ND - β-Caryophyllene 9960 0.996 Caryophyllene oxide ND - p-Cymene 146 0.0146 Eucalyptol ND - Geraniol ND - Guaiol 2560 0.256 α-Humulene 3730 0.373 Isopulegol ND - d-Limonene 30900 3.09 Linalool 349 0.0349 β-Myrcene 757 0.0757 cis-Nerolidol 670 0.0670 trans-Nerolidol 1310 0.131 α-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - γ-Terpinene	•		
α-Bisabolol 961 0.0961 Camphene 903 0.0903 δ-3-Carene ND - β-Caryophyllene oxide 9960 0.996 Caryophyllene oxide ND - p-Cymene 146 0.0146 Eucalyptol ND - Geraniol ND - Guaiol 2560 0.256 α-Humulene 3730 0.373 Isopulegol ND - d-Limonene 30900 3.09 Linalool 349 0.0349 β-Myrcene 757 0.0757 cis-Nerolidol 670 0.0670 trans-Nerolidol 1310 0.131 α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524	<u>Terpene</u>		
Camphene 903 0.0903 δ-3-Carene ND - β-Caryophyllene oxide 9960 0.996 Caryophyllene oxide ND - p-Cymene 146 0.0146 Eucalyptol ND - Geraniol ND - Guaiol 2560 0.256 α-Humulene 3730 0.373 Isopulegol ND - d-Limonene 30900 3.09 Linalool 349 0.0349 β-Myrcene 757 0.0757 cis-Nerolidol 670 0.0670 trans-Nerolidol 1310 0.131 α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524	α-Bisabolol		
δ-3-Carene ND - β-Caryophyllene oxide 9960 0.996 Caryophyllene oxide ND - p-Cymene 146 0.0146 Eucalyptol ND - Geraniol ND - Guaiol 2560 0.256 α-Humulene 3730 0.373 Isopulegol ND - d-Limonene 30900 3.09 Linalool 349 0.0349 β-Myrcene 757 0.0757 cis-Nerolidol 670 0.0670 trans-Nerolidol 1310 0.131 α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524			
β-Caryophyllene oxide ND - p-Cymene 146 0.0146 Eucalyptol ND - Geraniol ND - Guaiol 2560 0.256 α-Humulene 3730 0.373 Isopulegol ND - d-Limonene 30900 3.09 Linalool 349 0.0349 β-Myrcene 757 0.0757 cis-Nerolidol 670 0.0670 trans-Nerolidol 1310 0.131 α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524	•		0.0303
Caryophyllene oxide ND - p-Cymene 146 0.0146 Eucalyptol ND - Geraniol ND - Guaiol 2560 0.256 α-Humulene 3730 0.373 Isopulegol ND - d-Limonene 30900 3.09 Linalool 349 0.0349 β-Myrcene 757 0.0757 cis-Nerolidol 670 0.0670 trans-Nerolidol 1310 0.131 α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524			-
oxide ND - p-Cymene 146 0.0146 Eucalyptol ND - Geraniol ND - Guaiol 2560 0.256 α-Humulene 3730 0.373 Isopulegol ND - d-Limonene 30900 3.09 Linalool 349 0.0349 β-Myrcene 757 0.0757 cis-Nerolidol 670 0.0670 trans-Nerolidol 1310 0.131 α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524		9960	0.996
Eucalyptol ND - Geraniol ND - Guaiol 2560 0.256 α-Humulene 3730 0.373 Isopulegol ND - d-Limonene 30900 3.09 Linalool 349 0.0349 β-Myrcene 757 0.0757 cis-Nerolidol 670 0.0670 trans-Nerolidol 1310 0.131 α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524		ND	-
Geraniol ND - Guaiol 2560 0.256 α-Humulene 3730 0.373 Isopulegol ND - d-Limonene 30900 3.09 Linalool 349 0.0349 β-Myrcene 757 0.0757 cis-Nerolidol 670 0.0670 trans-Nerolidol 1310 0.131 α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524	p-Cymene	146	0.0146
Guaiol 2560 0.256 α-Humulene 3730 0.373 Isopulegol ND - d-Limonene 30900 3.09 Linalool 349 0.0349 β-Myrcene 757 0.0757 cis-Nerolidol 670 0.0670 trans-Nerolidol 1310 0.131 α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524	Eucalyptol	ND	-
α-Humulene 3730 0.373 Isopulegol ND - d-Limonene 30900 3.09 Linalool 349 0.0349 β-Myrcene 757 0.0757 cis-Nerolidol 670 0.0670 trans-Nerolidol 1310 0.131 α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524	Geraniol	ND	-
Isopulegol ND - d-Limonene 30900 3.09 Linalool 349 0.0349 β-Myrcene 757 0.0757 cis-Nerolidol 670 0.0670 trans-Nerolidol 1310 0.131 α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524	Guaiol	2560	0.256
d-Limonene 30900 3.09 Linalool 349 0.0349 β-Myrcene 757 0.0757 cis-Nerolidol 670 0.0670 trans-Nerolidol 1310 0.131 α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524	α-Humulene	3730	0.373
Linalool 349 0.0349 β-Myrcene 757 0.0757 cis-Nerolidol 670 0.0670 trans-Nerolidol 1310 0.131 α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524	Isopulegol	ND	-
β-Myrcene 757 0.0757 cis-Nerolidol 670 0.0670 trans-Nerolidol 1310 0.131 α -Ocimene ND - β -Ocimene 408 0.0408 α -Pinene 6580 0.658 β -Pinene 6730 0.673 α -Terpinene ND - γ -Terpinene ND - Terpinolene 524 0.0524	d-Limonene	30900	3.09
cis-Nerolidol 670 0.0670 trans-Nerolidol 1310 0.131 α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524	Linalool	349	0.0349
trans-Nerolidol 1310 0.131 α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524	β-Myrcene	757	0.0757
α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524	cis-Nerolidol	670	0.0670
α-Ocimene ND - β-Ocimene 408 0.0408 α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524	trans-Nerolidol	1310	0.131
α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524	α-Ocimene	ND	-
α-Pinene 6580 0.658 β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524	β-Ocimene	408	0.0408
β-Pinene 6730 0.673 α-Terpinene ND - γ-Terpinene ND - Terpinolene 524 0.0524	•	6580	0.658
α-TerpineneND-γ-TerpineneND-Terpinolene5240.0524	β-Pinene		0.673
γ-Terpinene ND - Terpinolene 524 0.0524	•		<u>-</u>
Terpinolene 524 0.0524	•		_
	-		0.0524
	TOTAL	66500	6.65

Reporting Limit (µg/g): 72.8

"-" Not detected above LOD.











E20231110TFSLHR01



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36425)

Testing Location:Customer ID: 2168Order ID: OR10701Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Concentrate232 S. Broadview St.5601 Old Greenwood Rd Suite CM00065C13244936466Mass: 4grams

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run:** Date Collected: 11/13/2023

Cultivar (Strain) or Sample Description: Truffle Shuffle Solventless Live Hash Rosin

Date Completed: 11/15/2023

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

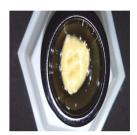
Analysis Date/Time: 11/14/2023 2228 Method: HS/GC/MS Deviations from SOP:

Analyst: KF Instrument: Agilent 7890/5975 None

License: 00065C

Solvent	Result (μg/g)	<u>LOD</u> (μg/g)	LOQ (µg/g)	<u>Action</u> <u>Level</u> (μg/g)	Solvent	Result (μg/g)	<u>LOD</u> (µg/g)	LOQ (µg/g)	<u>Action</u> <u>Level</u> (μg/g)
Acetone (67-64-1)	-	137	274	5000	n-Heptane (142-82-5)	-	137	274	5000
Acetonitrile (75-5-8)	-	137	274	410	n-Hexane (110-54-3)	-	47.9	95.9	290
Benzene (71-43-2)	-	1.37	2.74	2	Isobutane (75-28-5)	-	137	274	5000
n-Butane (106-97-2)	-	137	274	5000	Isopropanol (67-63-0)	-	137	274	5000
1-Butanol (71-36-3)	-	137	274	5000	Isopropyl acetate	_	137	274	5000
2-Butanol (78-92-2)	-	137	274	5000	(108-21-4)		157	2/4	5000
2-Butanone (78-93-3)	-	137	274	5000	Isopropyl benzene (98-82-8)	-	13.7	27.4	70
Cyclohexane (110-82-7)	-	137	274	3880	Methanol (67-56-1)	-	137	274	3000
1,2-Dimethoxyethane (110-71-4)	-	13.7	27.4	100	2-Methylbutane (78-78-4)	-	137	274	5000
N,N-Dimethylacetamide (127-19-5)	-	137	274	1090	Methylene chloride (75-9-2)	-	137	274	600
2,2-Dimethylbutane (75-83-2)	-	47.9	95.9	290	2-Methylpentane (107-83-5)	-	47.9	95.9	290
2,3-Dimethylbutane (79-29-8)	-	47.9	95.9	290	3-Methylpentane (96-10-0) n-Pentane (109-66-0)	-	47.9 137	95.9 274	290 5000
N,N-Dimethylformamide (68-12-2)	-	137	274	880	1-Pentanol (71-41-0) n-Propane (74-98-6)	-	137 137	274 274	5000 5000
Dimethylsulfoxide (67-68-5)	-	137	274	5000	1-Propanol (71-23-8)	-	137	274	5000
1,4-Dioxane (123-91-1)	_	137	274	380	Pyridine (110-86-1)	-	47.9	95.9	200
Ethanol (64-17-5)	_	137	274	5000	Tetrahydrofuran (109-99-9)	-	137	274	720
2-Ethoxyethanol (110-80-5)	-	47.9	95.9	160	Tetramethylene sulfone (126-33-0)	-	47.9	95.9	160
Ethyl ether (60-29-7)	-	137	274	5000	Toluene (108-88-3)	-	137	274	890
Ethyl acetate (141-78-6)	-	137	274	5000	o-Xylene (95-47-6)	_	137	274	2170
Ethyl benzene (100-41-4)	-	137	274	2170	m,p-Xylene (108-38-3 or				
Ethylene glycol (107-21-1)	-	137	274	620	106-42-3)	-	137	274	2170
Ethylene oxide (75-21-8)	-	13.7	27.4	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170

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



Date Received: 11/13/2023

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.

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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E20231110TFSLHR01



CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36425)

Testing Location:Customer ID: 2168Order ID: OR10701Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Concentrate232 S. Broadview St.5601 Old Greenwood Rd Suite CM00065C13244936466Mass: 4grams

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run:** Date Collected: 11/13/2023

Cultivar (Strain) or Sample Description: Truffle Shuffle Solventless Live Hash Rosin

Date Completed: 11/15/2023

PESTICIDES PROFILE (SOP: SOP-PEST-001)

Analysis Date/Time: 11/14/2023 1709 Method: LC/MS/MS Deviations from SOP:

Analyst: KF Instrument: Shimadzu LC-8050 None

License: 00065C

<u>Pesticide</u>	<u>Result</u> (μg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> <u>Level</u> (μg/g)	<u>Pesticide</u>	<u>Result</u> (μg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	Action Level (µg/g)
Abamectin (71751-41-2)	-	0.0435	0.348	0.5	Kresoxim-methyl	_	0.0435	0.348	0.4
Acephate (30560-19-1)	-	0.0435	0.348	0.4	(143390-89-0)				
Acequinocyl (57960-19-7)	-	0.0435	0.348	2	Malathion (121-75-5)	-	0.0435	0.348	0.2
Acetamiprid (135410-20-7)	-	0.0435	0.348	0.2	Metalaxyl (57837-19-1) Methiocarb (2032-65-7)	-	0.0435 0.0435	0.348 0.348	0.2 0.2
Aldicarb (116-06-3)	-	0.0435	0.348	0.4	Methomyl (16752-77-5)	-	0.0435	0.348	0.4
Azoxystrobin (131860-33-8)	-	0.0435	0.348	0.2	Methyl parathion (298-0-0)	-	0.0435	0.348	0.2
Bifenazate (149877-41-8)	-	0.0435	0.348	0.2	MGK 264 (113-48-4)	-	0.0435	0.348	0.2
Bifenthrin (82657-04-3)	-	0.0435	0.348	0.2	Myclobutanil		0.0435	0.348	0.2
Boscalid (188425-85-6)	-	0.0435	0.348	0.4	(88671-89-0)				
Carbaryl (63-25-2)	-	0.0435	0.348	0.2	Naled (300-76-5)	-	0.0435	0.348	0.5
Carbofuran (1563-66-2)	-	0.0435	0.348	0.2	Oxamyl (23135-22-0)	-	0.0435	0.348	1
Chlorantraniliprole (800008-45-7)	-	0.0435	0.348	0.2	Paclobutrazol (76738-62-0)	-	0.0435	0.348	0.4
Chlorfenapyr		0.0435	0.348	1	Permethrins (52645-53-1)	-	0.0435	0.348	0.2
(122453-73-0)	_	0.0435	0.346	1	Phosmet (732-11-6)	-	0.0435	0.348	0.2
Chlorpyrifos (2921-88-2)	-	0.0435	0.348	0.2	Piperonyl butoxide	_	0.0435	0.348	2
Clofentezine (74115-24-5)	-	0.0435	0.348	0.2	(51-03-6)		0.0425	0.240	0.0
Cyfluthrin (68359-37-5)	-	0.0435	0.348	1	Prallethrins (2331-36-9)	-	0.0435	0.348	0.2
Cypermethrin (52315-07-8)	-	0.0435	0.348	1	Propiconazole (60207-90-1))	-	0.0435	0.348	0.4
Daminozide (1596-84-5)	-	0.0435	0.348	1	Propoxur (114-26-1)	-	0.0435	0.348	0.2
DDVP (62-73-7)	-	0.0435	0.348	0.1	Pyrethrins (8003-34-7)	-	0.0435	0.348	1
Diazinon (333-41-5)	-	0.0435	0.348	0.2	Pyridaben (96489-71-3)	-	0.0435	0.348	0.2
Dimethoate (60-51-5)	-	0.0435	0.348	0.2	Spinosad (168316-95-8)	-	0.0435	0.348	0.2
Ethoprophos (13194-48-4)	-	0.0435	0.348	0.2	Spiromesifen (283594-90-1)	-	0.0435	0.348	0.2
Etofenprox (80844-07-1)	-	0.0435	0.348	0.4	Spirotetramat				
Etoxazole (153233-91-1)	-	0.0435	0.348	0.2	(203313-25-1)	-	0.0435	0.348	0.2
Fenoxycarb (72490-01-8)	-	0.0435	0.348	0.2	Spiroxamine		0.0435	0.348	0.4
(E)-Fenpyroximate (134098-61-6)	-	0.0435	0.348	0.4	(118134-30-8) Tebuconazole	-			
Fipronil (120068-37-3)	-	0.0435	0.348	0.4	(80443-41-0)	-	0.0435	0.348	0.4
Flonicamid (158062-67-0)	-	0.0435	0.348	1	Thiacloprid	_	0.0435	0.348	0.2
Fludioxinil (131341-86-1)	-	0.0435	0.348	0.4	(111988-49-9)				
Hexythiazox (78587-05-0)	-	0.0435	0.348	1	Thiamethoxam (153719-23-4)	-	0.0435	0.348	0.2
Imazalil (35554-44-0)	-	0.0435	0.348	0.2	Trifloxystrobin				
Imidacloprid (138261-41-3)	-	0.0435	0.348	0.4	(141517-21-7)		0.0435	0.348	0.2



Color Key

Date Received: 11/13/2023

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.

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.

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		













Date Completed: 11/15/2023

CERTIFICATE OF ANALYSIS (SAMPLE ID: SA36425)

Order ID: OR10701 **Testing Location:** Customer ID: 2168 Sample Type: Primary Arkansas River Valley Relief MIPS Lot Number: Matrix: Concentrate 232 S. Broadview St. 5601 Old Greenwood Rd Suite C M00065C13244936466 Mass: 4grams

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 11/13/2023 E20231110TFSLHR01 License: ADH 113 License: 00065C **Date Received:** 11/13/2023 Cultivar (Strain) or Sample Description: Truffle Shuffle Solventless Live Hash Rosin

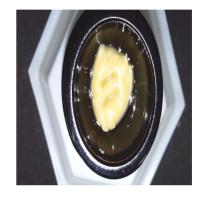
HEAVY METAL PROFILE (SOP: SOP-ICP-200.6)

Analysis Date/Time: 11/14/2023 1803 (ICP/OES) Method: ICP/MS **Deviations from SOP:**

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

Analyst: KF

<u>Heavy Metal</u>	<u>Result</u> (μg/kg)	<u>LOD</u> (µg/kg)	LOQ (µg/kg)	<u>Action Level</u> (μg/kg)
Arsenic (As)	-	57.5	91.0	200
Cadmium (Cd)	-	57.5	91.0	200
Lead (Pb)	-	57.5	91.0	500
Mercury (Hg)	-	57.5	91.0	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 < 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: SA36425)

Testing Location:Customer ID: 2168Sample ID: SA36425Sample Type: PrimaryArkansasRiver Valley Relief MIPSLot Number:Matrix: Concentrate232 S. Broadview St.5601 Old Greenwood Rd Suite CM00065C13244936466Mass: 4grams

Greenbrier, AR 72058 Fort Smith, AR 72903 **Production Run: Date Collected:** 11/13/2023 License: ADA 05_H273 License: 00065C E20231110TFSLHR01 **Date Received:** 11/13/2023

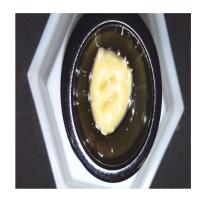
Cultivar (Strain) or Sample Description: Truffle Shuffle Solventless Live Hash Rosin Date Completed: 11/15/2023

MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 11/15/2023 0948 Method: Hardy Diagnostics CompactDry Deviations from SOP:

Analyst: PW Instrument: Thermo Incubator None

Bacteria/Microbe	<u>Result</u> (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

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.

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