







## SUMMARY OF ANALYSIS (SAMPLE ID: SA34719)

Testing Location:	Customer ID: 2168	Order ID: OR10354	Sample Type: Primary
Arkansas	<b>River Valley Relief MIPS</b>	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13247510882	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 03/15/2023
License: ADH 113	License: 00065C	E20230314GVALS01	Date Received: 03/16/2023
Cultivar (Strain) or Sample D	Date Completed: 03/22/2023		

\*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.

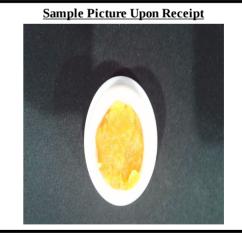
Moisture Content (%)	Water Activity (aw)	PASS/FAIL		
Not Tested	Not Tested	PASS		

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.

<u>Cannabinoids (Top 3)</u>	<u>(%)</u>	<u>mg/g</u>
THCA	75.6	756
Δ9-ΤΗC	4.74	47.4
CBGA	3.25	32.5
TOTAL CBD	-	-
TOTAL THC	71.1	711
TOTAL CANNABINOIDS	84.0	840
<u>Terpenes (Top 5)</u>	<u>(%)</u>	µg∕g
<u>Terpenes (Top 5)</u> d-Limonene	<u>(%)</u> 0.987	<b>µg/g</b> 9870
d-Limonene	0.987	9870
d-Limonene β-Myrcene	0.987 0.755	9870 7550
d-Limonene β-Myrcene β-Caryophyllene	0.987 0.755 0.426	9870 7550 4260

<u>Contaminants</u>	PASS/FAIL				
Heavy Metals:	PASS				
Microbiology:	PASS				
Pesticides:	PASS				
Residual Solvents:	PASS				





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## **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA34719)**

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Cultivar (Strain) or Sample	Date Completed: 03/22/2023		

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

Analysis Date/Time: 3/17/2023 1700 Analyst: PW Method: HPLC/DAD Instrument: Agilent 1100

<u>Cannabinoid</u>	<u>Result</u> (%)	<u>Result</u> (mg/g)	<u>LOD</u> (mg/g)	LOQ (mg/g)	<u>Result</u> ( <u>mg/</u> <u>mL</u> )	<u>Per</u> Serving (mg)	<u>Per</u> g <u>Unit</u> (mg)
CBC	ND	ND	0.474	1.11	-	-	-
CBCA	ND	ND	1.47	3.44	-	-	-
CBD	ND	ND	3.34	7.81	-	-	-
CBDA	ND	ND	1.23	2.87	-	-	-
CBDV	ND	ND	0.537	1.25	-	-	-
CBDVA	ND	ND	1.43	3.34	-	-	-
CBG	<loq< td=""><td><loq< td=""><td>2.17</td><td>5.07</td><td>-</td><td>-</td><td>-</td></loq<></td></loq<>	<loq< td=""><td>2.17</td><td>5.07</td><td>-</td><td>-</td><td>-</td></loq<>	2.17	5.07	-	-	-
CBGA	3.25	32.5	1.56	3.63	-	32.5	32.5
CBL	ND	ND	2.51	5.86	-	-	-
CBN	ND	ND	1.15	2.69	-	-	-
CBNA	ND	ND	1.24	2.90	-	-	-
Δ9-ТНС	4.74	47.4	1.38	3.22	-	47.4	47.4
$\Delta 8$ -THC	ND	ND	2.15	5.03	-	-	-
THCA	75.6	756	0.749	1.75	-	756	756
THCV	ND	ND	1.80	4.19	-	-	-
THCVA	0.384	3.84	0.574	1.34	-	3.84	3.84
TOTAL	84.0	840			-	840	840
TOTAL CBC	-	-			-	-	-
TOTAL CBD	-	-			-	-	-
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	2.85	28.5			-	28.5	28.5
TOTAL CBN	-	-			-	-	-
TOTAL THC	71.1	711			-	711	711
TOTAL THCV	0.333	3.33			-	3.33	3.33

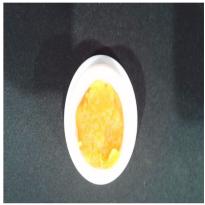
\* CBCA, CBDVA, CBL, CBNA, and THCVA are in process of being added to our accreditation scope.

*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): -



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) +  $\Delta$ 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.

Kyle W. Felling, Ph.D. atory Directo

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Cultivar (Strain) or Sample	<b>Date Completed:</b> 03/22/2023		

Analysis Date/Time:03/22/2023 0315 Analyst: KF

# TERPENOID PROFILE

Method: GC/MS

Analyst: KF		Instrument: Agilent 7890/597			
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)			
α-Bisabolol	1030	0.103			
Camphene	431	0.0431			
δ-3-Carene	-	-			
β-Caryophyllene	4260	0.426			
Caryophyllene oxide	3100	0.310	•		
p-Cymene	-	-			
Eucalyptol	-	-			
Geraniol	-	-			
Guaiol	1520	0.152	I		
α-Humulene	1950	0.195			
Isopulegol	-	-			
d-Limonene	9870	0.987			
Linalool	3530	0.353			
β-Myrcene	7550	0.755			
cis-Nerolidol	-	-			
trans-Nerolidol	756	0.0756			
α-Ocimene	-	-			
β-Ocimene	-	-			
α-Pinene	968	0.0968			
β-Pinene	1500	0.150	1		
α-Terpinene	-	-			
γ-Terpinene	-	-			
Terpinolene	915	0.0915			
TOTAL	37400	3.74			



**Deviations from SOP:** 

None

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

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"-" Not detected above RL

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Kyle W. Felling, Ph.D. Laboratory Director



Reporting Limit (µg/g): 46.1









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Cultivar (Strain) or Sample	Date Completed: 03/22/2023		

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

Analysis Date/Time: 3/18/2023 0529			Method: HS/GC/MS				<b>Deviations from SOP:</b>			
Analyst: KF				Iı	nstrument: Agilent 789	0/5975		Ν	one	
<u>Solvent</u>	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	<u>Solvent</u>	<u>Result</u> (µg/g)		<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	
Acetone (67-64-1)	-	36.4	72.9	5000	n-Heptane (142-82-5)	-	36.4	72.9	5000	
Acetonitrile (75-5-8)	-	36.4	72.9	410	n-Hexane (110-54-3)	-	12.8	25.5	290	
Benzene (71-43-2)	-	0.364	0.729	2	Isobutane (75-28-5)	-	36.4	72.9	5000	
n-Butane (106-97-2)	322	36.4	72.9	5000	Isopropanol (67-63-0)	-	36.4	72.9	5000	
1-Butanol (71-36-3)	-	36.4	72.9	5000	Isopropyl acetate	_	36.4	72.9	5000	
2-Butanol (78-92-2)	-	36.4	72.9	5000	(108-21-4)		50.4	/2.5	5000	
2-Butanone (78-93-3)	-	36.4	72.9	5000	Isopropyl benzene (98-82-8)	-	3.64	7.29	70	
Cyclohexane (110-82-7)	-	36.4	72.9	3880	(56-62-6) Methanol (67-56-1)		36.4	72.9	3000	Color Key
1,2-Dimethoxyethane	_	3.64	7.29	100	2-Methylbutane (78-78-4)		36.4	72.9	5000	
(110-71-4) N,N-Dimethylacetamide	_	36.4	72.9	1090	Methylene chloride (75-9-2)		36.4	72.9	600	RESULT < AL RESULT > AL
(127-19-5) 2,2-Dimethylbutane	_	12.8	25.5	290	2-Methylpentane (107-83-5)	-	12.8	25.5	290	"DET" detected less than LOQ
(75-83-2)					3-Methylpentane (96-10-0)		12.8	25.5	290	"-" not detected above
2,3-Dimethylbutane (79-29-8)	-	12.8	25.5	290	n-Pentane (109-66-0)		36.4	72.9	5000	LOD
N.N-Dimethylformamide		<b>0</b> 0 4			1-Pentanol (71-41-0)	-	36.4	72.9	5000	"*" V
(68-12-2)	-	36.4	72.9	880	n-Propane (74-98-6)	-	36.4	72.9	5000	"*" - o,m,p-Xylene and Ethylbenzene
Dimethylsulfoxide (67-68-5)	-	36.4	72.9	5000	1-Propanol (71-23-8)	-	36.4	72.9	5000	Action levels are
1,4-Dioxane (123-91-1)	-	36.4	72.9	380	Pyridine (110-86-1)	-	12.8	25.5	200	referenced from the State of
Ethanol (64-17-5)	-	36.4	72.9	5000	Tetrahydrofuran (109-99-9)	-	36.4	72.9	720	Arkansas
2-Ethoxyethanol (110-80-5)	-	12.8	25.5	160	Tetramethylene sulfone (126-33-0)	-	12.8	25.5	160	MMJ testing
Ethyl ether (60-29-7)	-	36.4	72.9	5000	Toluene (108-88-3)	_	36.4	72.9	890	guidelines.
Ethyl acetate (141-78-6)	-	36.4	72.9	5000	o-Xylene (95-47-6)	_	36.4	72.9	2170	A value of "-"
Ethyl benzene (100-41-4)	-	36.4	72.9	2170	m,p-Xylene (108-38-3 or					for the action level
Ethylene glycol (107-21-1)	-	36.4	72.9	620	106-42-3)	-	36.4	72.9	2170	means that analyte
Ethylene oxide (75-21-8)	-	3.64	7.29	50	Xylenes* (1330-20-7)	-	43.3	86.7	2170	is not currently regulated by the regulations referenced above.
Solvent		Synonym(s	5)		Solvent	5	Synonym(s)			-0
Acetonitrile		Methyl Cya	nide, ACN		Ethylene glycol	1	1,2-Ethanedio	l		
1-Butanol		n-Butanol,	Butyl Alcol	hol	Isobutane	2	2-Methylpropa	ane		
2-Butanol		sec-Butyl a	lcohol		Isopropanol	2	2-Propanol, IP	4		
2-Butanone		Methyl ethy	yl ketone, N	1EK	Isopropyl Acetate	1	Acetic acid iso	propyl este	r	
1,2-Dimethoxyethane		Monoglym	e		Methanol	I	Methyl alcoho	1		
2,3-Dimethylbutane		Neohexane			2-Methylbutane	1	Isopentane			
2,3-Dimethylbutane		Diisopropy	'l		Methylene chloride	1	Dichlorometha	ane		
N,N-Dimethylformamide		DMF			2-Methylpentane	1	Isohexane			
Dimethysufoxide		DMSO			1-Pentanol	1	n-Amyl alcoho	ol		
2-Ethoxyethanol		Cellosolve,	Ethyl glyc	ol	1-Propanol	1	Propyl alcohol			
Ethyl ether		Diethyl eth	er, Ether		Tetrahydrofuran	5	THF			
Ethyl acetate		EtOAc			Tetramethylene sulfone	5	Sulfolane			
Ethyl benzene		Phenyletha	ne		Xylene	I	Dimethylbenz	ene		

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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**Deviations from SOP:** 

None

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## **PESTICIDES PROFILE (SOP: SOP-PEST-001)**

Analysis Date/Time: 03/17/2023 2209 Analyst: KF

# Method: LC/MS/MS

#### Instrument: Shimadzu LC-8050

<u>Pesticide</u>	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	<u>Pesticide</u>	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	
Abamectin (71751-41-2)	-	0.0461	0.369	0.5	Kresoxim-methyl		0.0464	0.000		
Acephate (30560-19-1)	-	0.0461	0.369	0.4	(143390-89-0)	-	0.0461	0.369	0.4	
Acequinocyl (57960-19-7)	-	0.0461	0.369	2	Malathion (121-75-5)	-	0.0461	0.369	0.2	
Acetamiprid		0.0461	0.369	0.2	Metalaxyl (57837-19-1)	-	0.0461	0.369	0.2	
(135410-20-7)	-	0.0461	0.369	0.2	Methiocarb (2032-65-7)	-	0.0461	0.369	0.2	<u>Color Key</u>
Aldicarb (116-06-3)	-	0.0461	0.369	0.4	Methomyl (16752-77-5)	-	0.0461	0.369	0.4	<b>RESULT &lt; AL</b>
Azoxystrobin (131860-33-8)	-	0.0461	0.369	0.2	Methyl parathion (298-0-0)	-	0.0461	0.369	0.2	<b>RESULT &gt; AL</b>
Bifenazate (149877-41-8)	-	0.0461	0.369	0.2	MGK 264 (113-48-4)	-	0.0461	0.369	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3)	-	0.0461	0.369	0.2	Myclobutanil	_	0.0461	0.369	0.2	"-" not detected above
Boscalid (188425-85-6)	-	0.0461	0.369	0.4	(88671-89-0)			0.260	0.5	LOD
Carbaryl (63-25-2)	-	0.0461	0.369	0.2	Naled (300-76-5)	-	0.0461	0.369	0.5	Permethrins measured as the
Carbofuran (1563-66-2)	-	0.0461	0.369	0.2	Oxamyl (23135-22-0)	-	0.0461	0.369	1	cumulative residue of the cis- and
Chlorantraniliprole (800008-45-7)	-	0.0461	0.369	0.2	Paclobutrazol (76738-62-0)	-	0.0461	0.369	0.4	trans- permethrin isomers.
Chlorfenapyr	_	0.0461	0.369	1	Permethrins (52645-53-1)	-	0.0461	0.369	0.2	Pyrethrins measured as the cumulative residue of the
(122453-73-0)					Phosmet (732-11-6)	-	0.0461	0.369	0.2	pyrethrin I, cinerin I, and jasmolin
Chlorpyrifos (2921-88-2)	-	0.0461	0.369	0.2	Piperonyl butoxide (51-03-6)	-	0.0461	0.369	2	I isomers.
Clofentezine (74115-24-5)	-	0.0461	0.369	0.2	(31-03-0) Prallethrins (2331-36-9)		0.0461	0.369	0.2	Action levels are referenced from
Cyfluthrin (68359-37-5)	-	0.0461	0.369	1	Propiconazole					the
Cypermethrin (52315-07-8)	-	0.0461	0.369	1	(60207-90-1))	-	0.0461	0.369	0.4	State of Arkansas MMJ testing guidelines.
Daminozide (1596-84-5)	-	0.0461	0.369	1	Propoxur (114-26-1)	-	0.0461	0.369	0.2	A value of "-" for the action level
DDVP (62-73-7)	-	0.0461	0.369	0.1	Pyrethrins (8003-34-7)	-	0.0461	0.369	1	means that analyte is not
Diazinon (333-41-5)	-	0.0461	0.369	0.2	Pyridaben (96489-71-3)	-	0.0461	0.369	0.2	currently regulated by the regulations referenced above.
Dimethoate (60-51-5)	-	0.0461	0.369	0.2	Spinosad (168316-95-8)	-	0.0461	0.369	0.2	
Ethoprophos (13194-48-4)	-	0.0461	0.369	0.2	Spiromesifen (283594-90-1)	-	0.0461	0.369	0.2	Disclaimer: This information is provided as a service and makes
Etofenprox (80844-07-1)	-	0.0461	0.369	0.4	Spirotetramat					no claims of efficacy and/or safety
Etoxazole (153233-91-1)	-	0.0461	0.369	0.2	(203313-25-1)	-	0.0461	0.369	0.2	of this product. Results are
Fenoxycarb (72490-01-8)	-	0.0461	0.369	0.2	Spiroxamine	_	0.0461	0.369	0.4	applicable only for the sample(s) analyzed and for the specific
(E)-Fenpyroximate (134098-61-6)	-	0.0461	0.369	0.4	(118134-30-8) Tebuconazole		0.0461	0.369	0.4	analysis conducted. This report is for informational purposes only
Fipronil (120068-37-3)	-	0.0461	0.369	0.4	(80443-41-0)	-	0.0401	0.309	0.4	and should not be used to
Flonicamid (158062-67-0)	-	0.0461	0.369	1	Thiacloprid	_	0.0461	0.369	0.2	diagnose, treat, or prevent any medical-related symptoms. The
Fludioxinil (131341-86-1)	-	0.0461	0.369	0.4	(111988-49-9)					statements and results herein have
Hexythiazox (78587-05-0)	-	0.0461	0.369	1	Thiamethoxam (153719-23-4)	-	0.0461	0.369	0.2	not been approved and/or
Imazalil (35554-44-0)	-	0.0461	0.369	0.2	Trifloxystrobin		0.0464	0.000	0.0	endorsed by the FDA.
Imidacloprid (138261-41-3)	-	0.0461	0.369	0.4	(141517-21-7)	-	0.0461	0.369	0.2	

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



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# **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA34719)**

Testing Location:	Customer ID: 2168	Order ID: OR10354	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13247510882	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 03/15/2023
License: ADH 113	License: 00065C	E20230314GVALS01	Date Received: 03/16/2023
Cultivar (Strain) or Sample	<b>Date Completed:</b> 03/22/2023		

## HEAVY METAL PROFILE (SOP: SOP-ICP-200.7)

Analysis Date/Time: 3/21/2023 2243 (ICP/OES) Analysis Date/Time: - (DMA) Analyst: KF		<b>Method:</b> ICP/OES <b>Instrument:</b> Agilent 720-ES		<b>Deviations from SOP:</b> None	
<u>Heavy Metal</u>	<u>Result</u> (µg/kg)	<u>LOD</u> (µg/kg)	<u>LOQ</u> (µg/kg)	<u>Action Level</u> (μg/kg)	
Arsenic (As)	-	58.5	92.7	200	
Cadmium (Cd)	-	58.5	92.7	200	$\sim$
Lead (Pb)	-	58.5	92.7	500	
Mercury (Hg)	-	58.5	92.7	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

## <u>Color Key</u>



"DET" detected less than LOQ

"-" not detected above LOD

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

*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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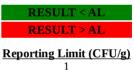
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Arkansas	<b>River Valley Relief MIPS</b>	Lot Number:	Matrix: Concentrate
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License: ADA 05_H273	License: 00065C	E20230314GVALS01	Date Received: 03/16/2023
Cultivar (Strain) or Sample D	Date Completed: 03/22/2023		

## MICROBIOLOGICAL PROFILE (SOP: SOP-Micro-001)

Analysis Date/Time: 2023031 Analyst: PW		Hardy Diagnostics CompactD <b>nt:</b> Thermo Incubator	ry <b>Deviations from SOP:</b> None
Bacteria/Microbe	<u>Result</u> (CFU/g)	<u>Action Level</u> (CFU/g)	
Aerobic Plate Count	NT	-	
Coliforms, Total	Absent	1	
Escherichia Coli (E. Coli)	Absent	100	
Mold/Yeast	NT	-	
Pseudomonas aeruginosa	NT	1	
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





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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