







## SUMMARY OF ANALYSIS (SAMPLE ID: SA35325)

Testing Location:	Customer ID: 2168	<b>Order ID:</b> OR10474	Sample Type: Primary		
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible		
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13224782109	Mass: 10pcs		
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 05/30/2023		
License: ADH 113	License: 00065C	P20230525PEA011	Date Received: 05/30/2023		
Cultivar (Strain) or Sample De	Date Completed: 06/01/2023				

\*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>		
Δ9-ΤΗϹ	0.189	1.89		
CBL	0.0291	0.291		
CBG	0.00533	0.0533		
TOTAL CBD	-	-		
TOTAL THC	0.189	1.89		
TOTAL CANNABINOIDS	0.227	2.27		
<u>Terpenes (Top 5)</u>	<u>(%)</u>	<u>µg/g</u>		
d-Limonene	0.000818	8.18		
α-Pinene	0.000409	4.09		
β-Myrcene	0.000124	1.24		
Terpinolene	0.000111	1.11		
α-Bisabolol	0.000	0.000		
TOTAL TERPENES	0.00146	14.6		
Contaminants PASS/FAIL	<u>Sample Picture Upon Receipt</u>			

Heavy Metals:	PASS
Microbiology:	PASS
Pesticides:	PASS
Residual Solvents:	PASS





Scan the QR code to verify results.

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## **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35325)**

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Arkansas	<b>River Valley Relief MIPS</b>	Lot Number:	Matrix: Edible
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Cultivar (Strain) or Sample De	Date Completed: 06/01/2023		

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

Analysis Date/Time: 05/31/2023 1521 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> (mg/ mL)	<u>Per</u> Serving (mg)	<u>Per</u> g <u>Unit</u> (mg)
CBC	ND	ND	0.00213	0.00497	-	-	-
CBCA	ND	ND	0.00662	0.0154	-	-	-
CBD	ND	ND	0.0150	0.0351	-	-	-
CBDA	ND	ND	0.00553	0.0129	-	-	-
CBDV	ND	ND	0.00241	0.00563	-	-	-
CBDVA	ND	ND	0.00643	0.0150	-	-	-
CBG	0.00533	0.0533	0.00976	0.0228	-	0.261	2.61
CBGA	ND	ND	0.0138	0.0163	-	-	-
CBL	0.0291	0.291	0.0113	0.0263	-	1.43	14.3
CBN	0.00318	0.0318	0.00518	0.0121	-	0.156	1.56
CBNA	ND	ND	0.00559	0.0130	-	-	-
Δ9-ΤΗC	0.189	1.89	0.00621	0.0145	-	9.27	92.7
$\Delta 8$ -THC	ND	ND	0.00968	0.0226	-	-	-
THCA	ND	ND	0.00336	0.00787	-	-	-
THCV	ND	ND	0.00807	0.0188	-	-	-
THCVA	ND	ND	0.00258	0.00600	-	-	-
TOTAL	0.227	2.27			-	11.1	111
TOTAL CBC	-	-			-	-	-
TOTAL CBD	-	-			-	-	-
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	0.00533	0.0533			-	0.261	2.61
TOTAL CBN	0.00318	0.0318			-	0.156	1.56
TOTAL THC	0.189	1.89			-	9.27	92.7
TOTAL THCV	-	-			-	-	-

*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): 4.91 SERVINGS/UNIT: 10

"-" 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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Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 05/30/2023
License: ADH 113	License: 00065C	P20230525PEA011	Date Received: 05/30/2023
Cultivar (Strain) or Sample	Date Completed: 06/01/2023		

#### **TERPENOID PROFILE**

Analysis Date/Time:06/01/2023 0713		Ν	fethod: GC/MS	<b>Deviations from SOP:</b>			
Analyst: KF		I	nstrument: Agilent 7890/5975	None			
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> <u>(%)</u>					
α-Bisabolol	-	-					
Camphene	-	-					
δ-3-Carene	-	-					
β-Caryophyllene	-	-		PEACH BELLINI			
Caryophyllene oxide	-	-					
p-Cymene	-	-		M NET WE 158m (45g)			
Eucalyptol	-	-					
Geraniol	-	-					
Guaiol	-	-		Abbreviations: GC - Gas			
α-Humulene	-	-		Chromatography, MS - Mass			
Isopulegol	-	-		Spectrometry, RL - Reporting Limit			
d-Limonene	8.18	0.000818		This information is provided as a service and makes no claims of efficacy and/or safety of this product.			
Linalool	-	-		Results are applicable only for the			
β-Myrcene cis-Nerolidol	1.24	0.000124 -	•	sample(s) analyzed and for the specific analysis conducted.			
trans-Nerolidol	-	-		This report is for informational purposes			
α-Ocimene	-	-		only and should not be used to diagnose, treat, or prevent any			
β-Ocimene	-	-		medical-related symptoms.			
α-Pinene	4.09	0.000409		The statements and results herein have not been approved and/or endorsed by			
β-Pinene	-	-		the FDA.			
α-Terpinene	-	-					
γ-Terpinene	-	-					
Terpinolene	1.11	0.000111		"-" Not detected above RL.			
TOTAL	14.6	0.00146		Reporting Limit (µg/g): 44.4			













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Cultivar (Strain) or Sample	Date Completed: 06/01/2023		

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

Analysis Date/Time:	05/31/20	)23 1835		ľ	Method: HS/GC/MS			, D	eviation	s from SOP:	
Analyst: KF				I	nstrument: Agilent 78	90/5975	None				
Solvent	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	LOQ (µg/g)	<u>Action</u> Level (µg/g)	<u>Solvent</u>	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)		
Acetone (67-64-1)	-	35.7	71.4	5000	n-Heptane (142-82-5)	-	35.7	71.4	5000		
Acetonitrile (75-5-8)	-	35.7	71.4	410	n-Hexane (110-54-3)	-	12.5	25.0	290	Wana Sarra	
Benzene (71-43-2)	-	0.357	0.714	2	Isobutane (75-28-5)	-	35.7	71.4	5000	PEACH BELLINI	
n-Butane (106-97-2)	-	35.7	71.4	5000	Isopropanol (67-63-0)	-	35.7	71.4	5000		
1-Butanol (71-36-3)	-	35.7	71.4	5000	Isopropyl acetate	_	35.7	71.4	5000	NET WIT 1584 (165)	
2-Butanol (78-92-2)	-	35.7	71.4	5000	(108-21-4)						
2-Butanone (78-93-3)	-	35.7	71.4	5000	Isopropyl benzene (98-82-8)	-	3.57	7.14	70		
Cyclohexane (110-82-7)	-	35.7	71.4	3880	Methanol (67-56-1)	414	35.7	71.4	3000	<u>Color Key</u>	
1,2-Dimethoxyethane	_	3.57	7.14	100	2-Methylbutane (78-78-4)		35.7	71.4	5000		
(110-71-4) N,N-Dimethylacetamide	_	35.7	71.4	1090	Methylene chloride (75-9-2)	-	35.7	71.4	600	RESULT < AL RESULT > AL	
(127-19-5) 2,2-Dimethylbutane (75-83-2)	_	12.5	25.0	290	2-Methylpentane (107-83-5)	-	12.5	25.0	290	"DET" detected less than LOQ	
2,3-Dimethylbutane					3-Methylpentane (96-10-0)	-	12.5	25.0	290	"-" not detected above	
(79-29-8)	-	12.5	25.0	290	n-Pentane (109-66-0)	-	35.7	71.4	5000	LOD	
N,N-Dimethylformamide		35.7	71.4	880	1-Pentanol (71-41-0)	-	35.7	71.4	5000	"*" - o,m,p-Xylene and	
(68-12-2)	-	35./	/1.4	000	n-Propane (74-98-6)	-	35.7	71.4	5000	Ethylbenzene	
Dimethylsulfoxide	_	35.7	71.4	5000	1-Propanol (71-23-8)	-	35.7	71.4	5000	0	
(67-68-5)		35.7	71.4	380	Pyridine (110-86-1)	-	12.5	25.0	200	Action levels are	
1,4-Dioxane (123-91-1)			71.4 71.4	5000	Tetrahydrofuran (109-99-9)	-	35.7	71.4	720	referenced from the State of Arkansas	
Ethanol (64-17-5) 2-Ethoxyethanol (110-80-5)	-	35.7 12.5	25.0	160	Tetramethylene sulfone	_	12.5	25.0	160	MMJ testing	
Ethyl ether (60-29-7)	-	35.7	23.0 71.4	5000	(126-33-0)					guidelines.	
Ethyl acetate (141-78-6)	-	35.7	71.4	5000	Toluene (108-88-3)	-	35.7	71.4	890	A value of "-"	
Ethyl benzene (100-41-4)	-	35.7	71.4	2170	o-Xylene (95-47-6)	-	35.7	71.4	2170	for the action level	
Ethylene glycol (107-21-1)		35.7	71.4	620	m,p-Xylene (108-38-3 or 106-42-3)	-	35.7	71.4	2170	means that analyte	
Ethylene oxide (75-21-8)		3.57	7.14	50	Xylenes* (1330-20-7)		43.3	86.7	2170	is not currently	
Solvent		Synonym(s		50	Solvent		Synonym(s)	00.7	2170	regulated by the regulations referenced above.	
Acetonitrile		Methyl Cya	-		Ethylene glycol	_	,2-Ethanediol				
1-Butanol		n-Butanol, 1		hol	Isobutane		-Methylpropa	20			
2-Butanol		sec-Butyl a	5	101	Isopropanol		2-Propanol, IPA				
2-Butanone		Methyl ethy		1FK	Isopropyl Acetate		Acetic acid iso				
1,2-Dimethoxyethane		Monoglym		ШК	Methanol		Methyl alcohol				
2,3-Dimethylbutane		Neohexane	e		2-Methylbutane		sopentane	L			
2,3-Dimethylbutane		Diisopropy	1		Methylene chloride		Dichlorometha	no			
N,N-Dimethylformamide		DIISOPIOPY			2-Methylpentane		sohexane	iic.			
Dimethysufoxide		DMF			1-Pentanol		i-Amyl alcoho	1			
2-Ethoxyethanol		Cellosolve,	Ethyl glyc	ol	1-Propanol		ropyl alcohol				
Ethyl ether		Diethyl eth		01	Tetrahydrofuran		THF				
Ethyl acetate		EtOAc	с, шис		Tetramethylene sulfone		Sulfolane				
Ethyl benzene		Phenyletha	ne		Xylene		Dimethylbenze	ne			
Luiyi Denzene		rnenyreuld	uc		Ay ICHC	L	Junetity i DeilZe	.11C			

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

## Analysis Date/Time: 06/01/2023 0723 Analyst: KF

#### Method: LC/MS/MS Instrument: Shimadzu I C-8050

Analyst: KF					<b>Instrument:</b> Shimadzu LC-8050				None	
<u>Pesticide</u>	<u>Result</u> (µg/g)	<u>LOD</u> (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	Pesticide	<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)	-	0.00940	0.0752	0.5	Kresoxim-methyl (143390-89-0)	-	0.00940	0.0752	0.4	
Acephate (30560-19-1)	-	0.00940	0.0752	0.4	(143350-85-0) Malathion (121-75-5)		0.00940	0.0752	0.2	
Acequinocyl (57960-19-7)	-	0.00940	0.0752	2	Metalaxyl (57837-19-1)		0.00940	0.0752	0.2	
Acetamiprid (135410-20-7)	-	0.00940	0.0752	0.2	Methiocarb (2032-65-7)	-	0.00940	0.0752	0.2	
Aldicarb (116-06-3)	-	0.00940	0.0752	0.4	Methomyl (16752-77-5)	-	0.00940	0.0752	0.4	
Azoxystrobin (131860-33-8)	-	0.00940	0.0752	0.2	Methyl parathion (298-0-0)	-	0.00940	0.0752	0.2	
Bifenazate (149877-41-8)	-	0.00940	0.0752	0.2	MGK 264 (113-48-4)	-	0.00940	0.0752	0.2	
Bifenthrin (82657-04-3)	-	0.00940	0.0752	0.2	Myclobutanil	_	0.00940	0.0752	0.2	
Boscalid (188425-85-6)	-	0.00940	0.0752	0.4	(88671-89-0)					
Carbaryl (63-25-2)	-	0.00940	0.0752	0.2	Naled (300-76-5)	-	0.00940	0.0752	0.5	"D
Carbofuran (1563-66-2)	-	0.00940	0.0752	0.2	Oxamyl (23135-22-0)	-	0.00940	0.0752	1	D
Chlorantraniliprole (800008-45-7)	-	0.00940	0.0752	0.2	Paclobutrazol (76738-62-0)	-	0.00940	0.0752	0.4	
Chlorfenapyr (122453-73-0)	-	0.00940	0.0752	1	Permethrins (52645-53-1)	-	0.00940	0.0752	0.2	P cum
Chlorpyrifos (2921-88-2)	-	0.00940	0.0752	0.2	Phosmet (732-11-6)	-	0.00940	0.0752	0.2	t
Clofentezine (74115-24-5)	-	0.00940	0.0752	0.2	Piperonyl butoxide (51-03-6)	-	0.00940	0.0752	2	,
Cyfluthrin (68359-37-5)	-	0.00940	0.0752	1	Prallethrins (2331-36-9)	-	0.00940	0.0752	0.2	1
Cypermethrin (52315-07-8)	-	0.00940	0.0752	1	Propiconazole (60207-90-1))	-	0.00940	0.0752	0.4	pyre
Daminozide (1596-84-5)	-	0.00940	0.0752	1	Propoxur (114-26-1)	_	0.00940	0.0752	0.2	Act
DDVP (62-73-7)	-	0.00940	0.0752	0.1	Pyrethrins (8003-34-7)	-	0.00940	0.0752	1	
Diazinon (333-41-5)	-	0.00940	0.0752	0.2	Pyridaben (96489-71-3)	_	0.00940	0.0752	0.2	Sta
Dimethoate (60-51-5)	-	0.00940	0.0752	0.2	Spinosad (168316-95-8)	-	0.00940	0.0752	0.2	
Ethoprophos (13194-48-4)	-	0.00940	0.0752	0.2	Spiromesifen (283594-90-1)	-	0.00940	0.0752	0.2	Av
Etofenprox (80844-07-1)	-	0.00940	0.0752	0.4	Spirotetramat					re
Etoxazole (153233-91-1)	-	0.00940	0.0752	0.2	(203313-25-1)	-	0.00940	0.0752	0.2	
Fenoxycarb (72490-01-8)	-	0.00940	0.0752	0.2	Spiroxamine	_	0.00940	0.0752	0.4	Dis pro
(E)-Fenpyroximate (134098-61-6)	-	0.00940	0.0752	0.4	(118134-30-8) Tebuconazole		0.00940	0.0752	0.4	no c
Fipronil (120068-37-3)	-	0.00940	0.0752	0.4	(80443-41-0)		0.00540	0.0732	0.4	app
Flonicamid (158062-67-0)	-	0.00940	0.0752	1	Thiacloprid (111988-49-9)	-	0.00940	0.0752	0.2	anal
Fludioxinil (131341-86-1)	-	0.00940	0.0752	0.4	(111988-49-9) Thiamethoxam					for
Hexythiazox (78587-05-0)	-	0.00940	0.0752	1	(153719-23-4)	-	0.00940	0.0752	0.2	di
Imazalil (35554-44-0)	-	0.00940	0.0752	0.2	Trifloxystrobin		0.00940	0.0752	0.2	me
Imidacloprid (138261-41-3)	-	0.00940	0.0752	0.4	(141517-21-7)	_	0.00940	0.0752	0.2	state

**Deviations from SOP:** 

Nono



Color Key

DET" detected less than LOQ "-" not detected above LOD Permethrins measured as the mulative residue of the cis- and trans-permethrin isomers. Pyrethrins measured as the cumulative residue of the rethrin I, cinerin I, and jasmolin I isomers. ction levels are referenced from the State of Arkansas MMJ testing guidelines. value of "-" for the action level means that analyte is not currently regulated by the regulations referenced above. isclaimer: This information is rovided as a service and makes claims of efficacy and/or safety of this product. Results are oplicable only for the sample(s) analyzed and for the specific alysis conducted. This report is or informational purposes only and should not be used to liagnose, treat, or prevent any nedical-related symptoms. The atements 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 Ouantification

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



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Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 05/30/2023
License: ADH 113	License: 00065C	P20230525PEA011	Date Received: 05/30/2023
Cultivar (Strain) or Sample	Date Completed: 06/01/2023		

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

Analysis Date/Time: 05/31/2023 1902 (ICP/OES) Analysis Date/Time: - (DMA) Analyst: KF		<b>Method:</b> ICP/MS <b>Instrument:</b> Agilent 7500ce		<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)	-	58.8	93.1	200	
Cadmium (Cd)	-	58.8	93.1	200	
Lead (Pb)	-	58.8	93.1	500	
Mercury (Hg)	-	58.8	93.1	100	PEACH BELLINI
					NET WILSon (Sa)

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



"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.













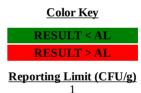
# **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA35325)**

Testing Location:	Customer ID: 2168	Sample ID: SA35325	Sample Type: Primary
Arkansas	River Valley Relief MIPS	Lot Number:	Matrix: Edible
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13224782109	Mass: 10pcs
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 05/30/2023
License: ADA 05_H273	License: 00065C	P20230525PEA011	Date Received: 05/30/2023
Cultivar (Strain) or Sample De	Date Completed: 06/01/2023		

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

Analysis Date/Time: 06/01/20 Analyst: PW		od: Hardy Diagnostics CompactDry ument: Thermo Incubator	<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	FASTACTING GUMMES	
Mold/Yeast	NT	-	OTCA SATIVA	
Pseudomonas aeruginosa	NT	-	A Der	
Salmonella spp.	NT	-	AM	
Staphylococcus aureus	NT	-	NET WE LSSee (45g)	

*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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