







### SUMMARY OF ANALYSIS (SAMPLE ID: SA34713)

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	M00065C13240855667	Mass: 4g				
Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	<b>Date Collected:</b> 03/15/2023				
License: ADH 113	License: 00065C	E20230313BBFLS01	Date Received: 03/16/2023				
Cultivar (Strain) or Sample Description: Ben's Buffalo Indica Hybrid Live Sugar 1g JarDate 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.

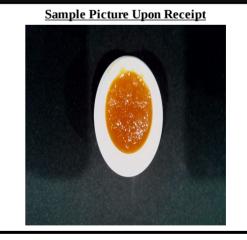
Moisture Content (%)	Water Activity (aw)	PASS/FAIL
Not Tested	Not Tested	PASS

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</u> /g
THCA	73.7	737
Δ9-ТНС	3.62	36.2
THCVA	1.92	19.2
TOTAL CBD	-	-
TOTAL THC	68.3	683
TOTAL CANNABINOIDS	81.7	817
<u>Terpenes (Top 5)</u>	<u>(%)</u>	µg∕g
<u>Terpenes (Top 5)</u> Terpinolene	<u>(%)</u> 2.04	<mark>µg/g</mark> 20400
Terpinolene	2.04	20400
Terpinolene β-Myrcene	2.04 0.934	20400 9340
Terpinolene β-Myrcene d-Limonene	2.04 0.934 0.636	20400 9340 6360

PASS/FAIL
PASS
PASS
PASS
PASS





Scan the QR code to verify results.

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## **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA34713)**

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License: ADH 113	License: 00065C	E20230313BBFLS01	Date Received: 03/16/2023
Cultivar (Strain) or Sample Des	Date Completed: 03/22/2023		

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

Analysis Date/Time: 3/17/2023 1553 Analyst: PW Method: HPLC/DAD Instrument: Agilent 1100

<u>Cannabinoid</u>	<u>Result</u> (%)	<u>Result</u> (mg/g)	LOD (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	0.514	1.20	-	-	-
CBCA	ND	ND	1.60	3.72	-	-	-
CBD	ND	ND	3.62	8.46	-	-	-
CBDA	ND	ND	1.33	3.11	-	-	-
CBDV	ND	ND	0.581	1.36	-	-	-
CBDVA	ND	ND	1.55	3.62	-	-	-
CBG	ND	ND	2.35	5.49	-	-	-
CBGA	1.89	18.9	1.69	3.93	-	18.9	18.9
CBL	ND	ND	2.72	6.34	-	-	-
CBN	0.495	4.95	1.25	2.92	-	4.95	4.95
CBNA	ND	ND	1.35	3.14	-	-	-
Δ9-ΤΗC	3.62	36.2	1.50	3.49	-	36.2	36.2
$\Delta 8$ -THC	ND	ND	2.33	5.45	-	-	-
THCA	73.7	737	0.811	1.90	-	737	737
THCV	ND	ND	1.95	4.54	-	-	-
THCVA	1.92	19.2	0.622	1.45	-	19.2	19.2
TOTAL	81.7	817				817	817
TOTAL CBC	-	-			-	-	-
TOTAL CBD	-	-			-	-	-
TOTAL CBDV	-	-			-	-	-
TOTAL CBG	1.66	16.6			-	16.6	16.6
TOTAL CBN	0.495	4.95			-	4.95	4.95
TOTAL THC	68.3	683			-	683	683
TOTAL THCV	1.67	16.7			-	16.7	16.7

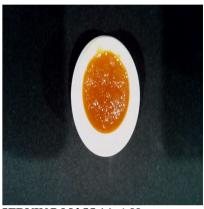
\* CBCA, CBDVA, CBL, CBNA, and THCVA are in process of being added to our accreditation scope.

*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 Dire

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Greenbrier, AR 72058	Fort Smith, AR 72903	Production Run:	Date Collected: 03/15/2023
License: ADH 113	License: 00065C	E20230313BBFLS01	Date Received: 03/16/2023
Cultivar (Strain) or Sample D	Date Completed: 03/22/2023		

Analysis Date/Time:03/22/2023 0109 Analyst: KF

#### **TERPENOID PROFILE**

Method: GC/MS Instrument: Agilent 7890/5975

#### **Deviations from SOP:** None

Abbreviations: GC - Gas

safety of this product.

analysis conducted.

treat, or prevent any medical-related symptoms.

"-" Not detected above RL

the FDA.

Chromatography, MS - Mass Spectrometry, RL - Reporting Limit

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sample(s) analyzed and for the specific

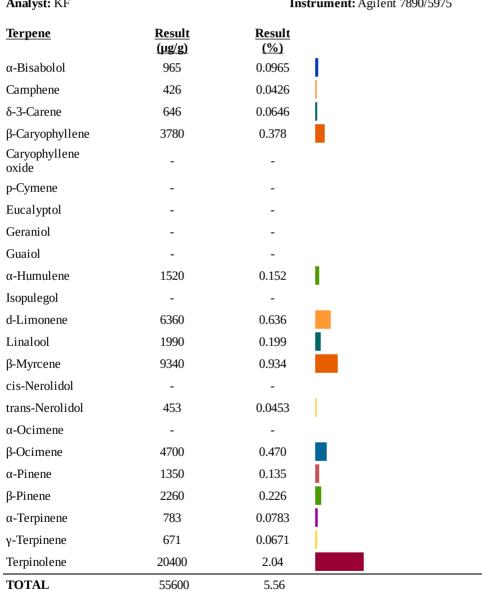
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5.56

55600

Reporting Limit (µg/g): 49.2













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

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

Analysis Date/Time:	3/18/202	3 0226		Ν	lethod: HS/GC/MS			Ď	eviations	from SOP:
Analyst: KF				Iı	nstrument: Agilent 789	0/5975		Ν	one	
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)	LOD (µg/g)	<u>LOQ</u> (µg/g)	<u>Action</u> Level (µg/g)	
Acetone (67-64-1)	-	39.0	78.0	5000	n-Heptane (142-82-5)	-	39.0	78.0	5000	
Acetonitrile (75-5-8)	-	39.0	78.0	410	n-Hexane (110-54-3)	-	13.6	27.3	290	
Benzene (71-43-2)	-	0.390	0.780	2	Isobutane (75-28-5)	-	39.0	78.0	5000	
n-Butane (106-97-2)	-	39.0	78.0	5000	Isopropanol (67-63-0)	-	39.0	78.0	5000	
1-Butanol (71-36-3)	-	39.0	78.0	5000	Isopropyl acetate	_	39.0	78.0	5000	(manager)
2-Butanol (78-92-2)	-	39.0	78.0	5000	(108-21-4)					
2-Butanone (78-93-3)	-	39.0	78.0	5000	Isopropyl benzene (98-82-8)	-	3.90	7.80	70	
Cyclohexane (110-82-7)	-	39.0	78.0	3880	Methanol (67-56-1)	_	39.0	78.0	3000	<u>Color Key</u>
1,2-Dimethoxyethane	_	3.90	7.80	100	2-Methylbutane (78-78-4)	_	39.0	78.0	5000	
(110-71-4) N,N-Dimethylacetamide	_	39.0	78.0	1090	Methylene chloride (75-9-2)	-	39.0	78.0	600	RESULT < AL RESULT > AL
(127-19-5) 2,2-Dimethylbutane (75-83-2)	-	13.6	27.3	290	2-Methylpentane (107-83-5)	-	13.6	27.3	290	"DET" detected less than LOQ
2,3-Dimethylbutane					3-Methylpentane (96-10-0)	-	13.6	27.3	290	"-" not detected above
(79-29-8)	-	13.6	27.3	290	n-Pentane (109-66-0)	-	39.0	78.0	5000	LOD
N,N-Dimethylformamide		39.0	78.0	880	1-Pentanol (71-41-0)	-	39.0	78.0	5000	"*" - o,m,p-Xylene and
(68-12-2)		39.0	/0.0	000	n-Propane (74-98-6)	-	39.0	78.0	5000	Ethylbenzene
Dimethylsulfoxide (67-68-5)		39.0	78.0	5000	1-Propanol (71-23-8)	-	39.0	78.0	5000	5
(07-00-3) 1,4-Dioxane (123-91-1)		39.0	78.0	380	Pyridine (110-86-1)	-	13.6	27.3	200	Action levels are referenced from the State of
Ethanol (64-17-5)	-	39.0	78.0	5000	Tetrahydrofuran (109-99-9)	-	39.0	78.0	720	Arkansas
2-Ethoxyethanol (110-80-5)	-	13.6	27.3	160	Tetramethylene sulfone	_	13.6	27.3	160	MMJ testing
Ethyl ether (60-29-7)		39.0	78.0	5000	(126-33-0)					guidelines.
Ethyl acetate (141-78-6)		39.0	78.0	5000	Toluene (108-88-3)	-	39.0	78.0	890	A value of "-"
Ethyl benzene (100-41-4)		39.0	78.0	2170	o-Xylene (95-47-6)	-	39.0	78.0	2170	for the action level
Ethylene glycol (107-21-1)		39.0	78.0	620	m,p-Xylene (108-38-3 or 106-42-3)	-	39.0	78.0	2170	means that analyte
Ethylene oxide (75-21-8)		3.90	7.80	50	Xylenes* (1330-20-7)		43.3	86.7	2170	is not currently
Solvent		Synonym(s		50	Solvent		-	00.7	2170	regulated by the regulations referenced above.
<u>Sorvent</u> Acetonitrile		Methyl Cya			<u>Sorvent</u> Ethylene glycol		<b>Synonym(s)</b> 1,2-Ethanediol			
1-Butanol		n-Butanol,		hal	Isobutane					
		,	5	101			2-Methylpropane			
2-Butanol 2-Butanone		sec-Butyl a		Œν	Isopropanol		2-Propanol, IPA			
1,2-Dimethoxyethane		Monoglym			Acetic acid isopropyl ester					
		0.0			Methyl alcohol					
2,3-Dimethylbutane 2,3-Dimethylbutane		Neohexane	1		2-Methylbutane Methylene chloride		sopentane Dichlorometha	20		
N,N-Dimethylformamide		Diisopropy DMF	1		2-Methylpentane		sohexane	ne		
-		DMF			1-Pentanol			1		
Dimethysufoxide 2-Ethoxyethanol		Cellosolve,	Ethyl alve	ol			n-Amyl alcoho Propyl alcohol	1		
5				01	1-Propanol Totrahydrofuran		Propyl alcohol FHF			
Ethyl ether		Diethyl eth EtOAc	ei, Etilei		Tetrahydrofuran Tetramethylona culfono		Sulfolane			
Ethyl acetate					Tetramethylene sulfone			200		
Ethyl benzene		Phenyletha	ne		Xylene	1	Dimethylbenze	ne		

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 2037 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)	-	0.0492	0.394	0.5	Kresoxim-methyl	_	0.0492	0.394	0.4	
Acephate (30560-19-1)	-	0.0492	0.394	0.4	(143390-89-0)					
Acequinocyl (57960-19-7)	-	0.0492	0.394	2	Malathion (121-75-5)	-	0.0492	0.394	0.2	
Acetamiprid	-	0.0492	0.394	0.2	Metalaxyl (57837-19-1)	-	0.0492	0.394	0.2	Color Key
(135410-20-7) Aldicarb (116-06-3)		0.0492	0.394	0.4	Methiocarb (2032-65-7) Methomyl (16752-77-5)	-	0.0492 0.0492	0.394 0.394	0.2 0.4	
Azoxystrobin	-				Methyl parathion	-				RESULT < AL
(131860-33-8)	-	0.0492	0.394	0.2	(298-0-0)	-	0.0492	0.394	0.2	<b>RESULT &gt; AL</b>
Bifenazate (149877-41-8)	-	0.0492	0.394	0.2	MGK 264 (113-48-4)	-	0.0492	0.394	0.2	"DET" detected less than LOQ
Bifenthrin (82657-04-3)	-	0.0492	0.394	0.2	Myclobutanil	_	0.0492	0.394	0.2	"-" not detected above
Boscalid (188425-85-6)	-	0.0492	0.394	0.4	(88671-89-0)		0.0400	0.004	0 -	LOD
Carbaryl (63-25-2)	-	0.0492	0.394	0.2	Naled (300-76-5)	-	0.0492	0.394	0.5	Permethrins measured as the
Carbofuran (1563-66-2)	-	0.0492	0.394	0.2	Oxamyl (23135-22-0)	-	0.0492	0.394	1	cumulative residue of the <i>cis</i> - and
Chlorantraniliprole (800008-45-7)	-	0.0492	0.394	0.2	Paclobutrazol (76738-62-0)	-	0.0492	0.394	0.4	trans- permethrin isomers.
Chlorfenapyr	_	0.0492	0.394	1	Permethrins (52645-53-1)	-	0.0492	0.394	0.2	Pyrethrins measured as the cumulative residue of the
(122453-73-0)	-	0.0492	0.394	1	Phosmet (732-11-6)	-	0.0492	0.394	0.2	pyrethrin I, cinerin I, and jasmolin
Chlorpyrifos (2921-88-2)	-	0.0492	0.394	0.2	Piperonyl butoxide	_	0.0492	0.394	2	I isomers.
Clofentezine (74115-24-5)	-	0.0492	0.394	0.2	(51-03-6) Duall athring (2221-26-0)		0.0492	0.394	0.2	Action levels are referenced from
Cyfluthrin (68359-37-5)	-	0.0492	0.394	1	Prallethrins (2331-36-9)	-	0.0492	0.394	0.2	the
Cypermethrin (52315-07-8)	-	0.0492	0.394	1	Propiconazole (60207-90-1))	-	0.0492	0.394	0.4	State of Arkansas MMJ testing guidelines.
Daminozide (1596-84-5)	-	0.0492	0.394	1	Propoxur (114-26-1)	-	0.0492	0.394	0.2	A value of "-" for the action level
DDVP (62-73-7)	-	0.0492	0.394	0.1	Pyrethrins (8003-34-7)	-	0.0492	0.394	1	means that analyte is not
Diazinon (333-41-5)	-	0.0492	0.394	0.2	Pyridaben (96489-71-3)	-	0.0492	0.394	0.2	currently regulated by the
Dimethoate (60-51-5)	-	0.0492	0.394	0.2	Spinosad (168316-95-8)	-	0.0492	0.394	0.2	regulations referenced above.
Ethoprophos (13194-48-4)	-	0.0492	0.394	0.2	Spiromesifen (283594-90-1)	-	0.0492	0.394	0.2	Disclaimer: This information is
Etofenprox (80844-07-1)	-	0.0492	0.394	0.4	Spirotetramat					provided as a service and makes no claims of efficacy and/or safety
Etoxazole (153233-91-1)	-	0.0492	0.394	0.2	(203313-25-1)	-	0.0492	0.394	0.2	of this product. Results are
Fenoxycarb (72490-01-8)	-	0.0492	0.394	0.2	Spiroxamine	_	0.0492	0.394	0.4	applicable only for the sample(s) analyzed and for the specific
(E)-Fenpyroximate (134098-61-6)	-	0.0492	0.394	0.4	(118134-30-8) Tebuconazole		0.0492	0.394	0.4	analysis conducted. This report is for informational purposes only
Fipronil (120068-37-3)	-	0.0492	0.394	0.4	(80443-41-0)		0.0452	0.554	0.4	and should not be used to
Flonicamid (158062-67-0)	-	0.0492	0.394	1	Thiacloprid	-	0.0492	0.394	0.2	diagnose, treat, or prevent any medical-related symptoms. The
Fludioxinil (131341-86-1)	-	0.0492	0.394	0.4	(111988-49-9)					statements and results herein have
Hexythiazox (78587-05-0)	-	0.0492	0.394	1	Thiamethoxam (153719-23-4)	-	0.0492	0.394	0.2	not been approved and/or
Imazalil (35554-44-0)	-	0.0492	0.394	0.2	Trifloxystrobin		0.0.405	0.204	0.2	endorsed by the FDA.
Imidacloprid (138261-41-3)	-	0.0492	0.394	0.4	(141517-21-7)		0.0492	0.394	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) Cyfluthrin Baythroid Systhane Myclobutanil Propiconazole DDVP Dichlorvos Naled Dibrom Propoxur Ethoprophos Prophos Phosmet Imidan



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Synonym(s)

Tilt

Baygon

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License: ADH 113	License: 00065C	E20230313BBFLS01	Date Received: 03/16/2023
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#### HEAVY METAL PROFILE (SOP: SOP-ICP-200.7)

Analysis Date/Time: 3/21/2023 2222 (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)	-	54.8	86.8	200	
Cadmium (Cd)	-	54.8	86.8	200	
Lead (Pb)	-	54.8	86.8	500	and and
Mercury (Hg)	-	54.8	86.8	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.













# **CERTIFICATE OF ANALYSIS (SAMPLE ID: SA34713)**

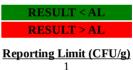
Testing Location:	Customer ID: 2168	Sample ID: SA34713	Sample Type: Primary
Arkansas	<b>River Valley Relief MIPS</b>	Lot Number:	Matrix: Concentrate
232 S. Broadview St.	5601 Old Greenwood Rd Suite C	M00065C13240855667	Mass: 4g
Greenbrier, AR 72058	Fort Smith, AR 72903	<b>Production Run:</b>	Date Collected: 03/15/2023
License: ADA 05_H273	License: 00065C	E20230313BBFLS01	Date Received: 03/16/2023
Cultivar (Strain) or Sample	<b>Date Completed:</b> 03/22/2023		

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

Analysis Date/Time: 20230318 Analyst: PW		ardy Diagnostics Compact <b>t:</b> Thermo Incubator	5	<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		A CONTRACTOR OF THE OWNER OF THE
Escherichia Coli (E. Coli)	Absent	100		Section .
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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