

Product Name:	Harmony Focus Gummy		
Product Batch:	WB21343		
Certificate ID Number:	CB220114008		
Date Tested:	1/14/2022		

Cannabinoid Profile & Potency Liquid Tincture:					
D9-THC:	1.18mg/unit				
CBD:	23.40mg/unit				
CBDV:	ND				
CBG:	9.39mg/unit				
CBC:	0.68mg/unit				
CBN:	ND				
Total Count:					
Total THC:	1.18mg/unit				
Total CBD:	23.40mg/unit				
Manufactured By: Palmetto Sy	nergistic Research				

Manufacturer Date: 1/1/2022

Elemental Analysis:	Pass
Microbiological Contaminants:	Pass
Pathogenic Bacterial Contaminants:	Pass
Mycotoxin Testing:	Pass
Pesticide Analysis:	Pass
Terpene Profile:	Please see the full lab for multiple terpene profiles.
Analysis of Volatile Organic Compounds:	Pass

This product has been reviewed by CannaBusiness Laboratories, LLC. The product contains less than 0.3% THC per the Farm Bill of 2018. This product is not intended to diagnose, treat, cure or prevent any disease. The FDA has not evaluated this



CannaBusiness Laboratories, LLC

2554 Palumbo Dr. Lexington, KY 40509

Certificate of Analysis

Customer: Palmetto Synergistic Research 8856 Pee Dee Hwy Conway, SC 29527 / 843-331-1246

Collected Date: Received Date: 1/14/2022 COA Released: 1/19/2022

Comments:

CANNABINOID PROFILE (Product Size = 3 g) Analyte LOQ (%) % weight mg/g mg/unit CBC 0.01 0.022 0.225 0.68 CBD 0.780 7.799 0.01 23.40 CBDa 0.01 ND ND ND 0.01 ND ND CBDV ND CBG 0.01 0.313 3.130 9.39 0.01 ND ND ND CBGa 0.01 ND ND ND CBN 0.01 ND ND ND d8-THC 0.01 0.039 0.392 d9-THC 1.18 THCa 0.01 ND ND ND **Total Cannabinoids** 1.154 11.54 34.62 **Total Potential THC** 1.18 0.039 0.392 **Total Potential CBD** 0.780 7.799 23.40 **Total Potential CBG** 0.313 3.130 9.39 Ratio of Total Potential CBD to Total Potential THC 20.00:1 Ratio of Total Potential CBG to Total Potential THC 8.03:1



Cannabinoids (% weight)

Sample ID: 220114022

External Sample ID:

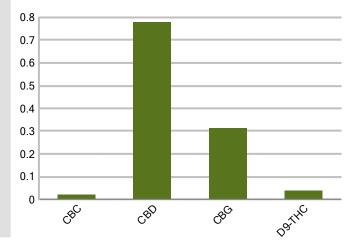
Product Type: Edible

Sample Type: Edible

Batch Number: WB21343

Order Number: CB220114008

Sample Name: HAR Focus Gummy



*Total Cannabinoids refers to the sum of all cannabinoids detected.

*Total Potential CBD = (0.877 x CBDa) + CBD. *Total Potential THC = (0.877 x THCa) + THC. *Total Potential CBG = (0.877 x CBGa) + CBG.

*Total Potential THC/CBD are calculated to take into account the loss of an acid group during decarboxylation.



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CannaBusiness Laboratories, LLC 2554 Palumbo Dr. Lexington, KY 40509



Certificate of Analysis

Sample ID:	22011
Sample Name:	HAR F
Sample Type:	Edible

220114022 Focus Gummy

Customer Palmetto Synergistic Research 8856 Pee Dee Hwy Conway, SC 29527 /



		Sample N
Overall Bat	tch Results	Sample ID
Pesticide	Moisture Content	Product T
Potency	Water Activity	Sample Ty Collected
Mycotoxins	Heavy Metals	Received
Microbial Screen	Residual Solvents	Batch Nui Batch Size
Terpenoids		Sample Si COA relea

ample Name:	HAR Focus (Gummy
ample ID:	220114022	
oduct Type:	Edible	
ample Type:	Edible	
ollected Date:		
eceived Date:	01/14/2022	
atch Number:	WB21343	
atch Size:		
ample Size:		
OA released:	01/19/2022	2:50 PM

Potency (mg/g)							
Date Tested: 01/15/20	CB-SOP-028	3					
Instrument:							
0.039 %	, D	1.154 %		11.	11.54 mg/g		
Total THC	Total CBD		Total Ca	annabinoids	Total Cannabinoids		
Analyte	Result	Units	LOQ	Result	Units		
CBC (Cannabichromer	0.022	%	0.010	0.225	mg/g		
CBD (Cannabidiol)	0.780	%	0.010	7.799	mg/g		
CBDa (Cannabidiolic A	ND	%	0.010	ND	mg/g		
CBDV (Cannabidivarin)	ND	%	0.010	ND	mg/g	
CBG (Cannabigerol)		0.313	%	0.010	3.130	mg/g	
CBGa (Cannabigerolic	ND	%	0.010	ND	mg/g		
CBN (Cannabinol)	ND	%	0.010	ND	mg/g		
D8-THC (D8-Tetrahydr	ND	%	0.010	ND	mg/g		
D9-THC (D9-Tetrahydr	0.039	%	0.010	0.392	mg/g		
THCa (Tetrahydrocann	abinolic Acid)	ND	%	0.010	ND	mg/g	

Terpenoids					
Date Tested: 01/15/2022		Method: C	B-SOP-02	26	
Instrument:					
Analyte	Result	Unit	LOQ	Result	Unit
alpha-Bisabolol	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
alpha-humulene	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
alpha-pinene	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
alpha-terpinene	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
beta-caryophyllene	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
Beta-myrcene	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
Beta-pinene	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
cis-Nerolidol	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
Camphene	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
d-Limonene	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
delta-3-Carene	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
Eucalyptol	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
gamma-Terpinene	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
Geraniol	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
Guaiol	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
Isopulegol	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
Linalool	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
Ocimene (mixture of isomers)	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
p-Isopropyltoluene (p-Cymene)	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
trans-beta-Ocimene	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
trans-Nerolidol	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%
Terpinolene	<loq< td=""><td>mg/g</td><td>0.100</td><td><loq< td=""><td>%</td></loq<></td></loq<>	mg/g	0.100	<loq< td=""><td>%</td></loq<>	%

Pesticides

Date Tested: 01/14/2022	Method: CB-SOP-025	Instrume	nt:				
Analyte	Result Units	LOQ	Result	Analyte	Result Units	LOQ	Result
Acephate	ND ppm	0.010		Acetamiprid	ND ppm	0.010	
Aldicarb	ND ppm	0.010		Azoxystrobin	ND ppm	0.010	
Bifenazate	ND ppm	0.010		Bifenthrin	ND ppm	0.100	
Boscalid	ND ppm	0.010		Carbaryl	ND ppm	0.010	
Carbofuran	ND ppm	0.010		Chlorantraniliprole	ND ppm	0.010	
Chlorpyrifos	ND ppm	0.010		Clofentezine	ND ppm	0.010	
Coumaphos	ND ppm	0.010		Daminozide	ND ppm	0.010	
Diazinon	ND ppm	0.010		Dichlorvos	ND ppm	0.010	
Dimethoate	ND ppm	0.010		Etofenprox	ND ppm	0.010	
Etoxazole	ND ppm	0.010		Fenhexamid	ND ppm	0.010	
Fenoxycarb	ND ppm	0.010		Fenpyroximate	ND ppm	0.010	
Fipronil	ND ppm	0.010		Flonicamid	ND ppm	0.100	

NT = Not tested, ND = Not detected; LOQ = Limit of Quantitation; <LOQ = Detected; >ULOL = Above upper limit of linearity; CFU/g = Colony forming units per 1 gram; TNTC = Too numerous to count

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Sample ID: Sample Name: Sample Type:

220114022 e: HAR Focus Gummy Edible

Certificate of Analysis

Date Tested: 01/14/2022	Method: CB-SOP-025	Instrume	nt:				
Analyte	Result Units	LOQ	Result	Analyte	Result Units	LOQ	Resul
Fludioxonil	ND ppm	0.010		Hexythiazox	ND ppm	0.010	
Imazalil	ND ppm	0.010		Imidacloprid	ND ppm	0.010	
Malathion	ND ppm	0.010		Metalaxyl	ND ppm	0.010	
Methiocarb	ND ppm	0.010		Methomyl	ND ppm	0.010	
Myclobutanil	ND ppm	0.010		Naled	ND ppm	0.010	
Oxamyl	ND ppm	0.010		Paclobutrazol	ND ppm	0.010	
Phosmet	ND ppm	0.010		Prallethrin	ND ppm	0.010	
Propiconazole	ND ppm	0.010		Propoxur	ND ppm	0.010	
Pyrethrin I	ND ppm	0.010		Pyrethrin II	ND ppm	0.010	
Pyridaben	ND ppm	0.010		Spinetoram	ND ppm	0.010	
Spiromesifen	ND ppm	0.010		Spirotetramat	ND ppm	0.010	
Tebuconazole	ND ppm	0.010		Thiacloprid	ND ppm	0.010	
Thiamethoxam	ND ppm	0.010		Trifloxystrobin	ND ppm	0.010	
Ethoprophos	ND ppm	0.010		Kresoxym-methyl	ND ppm	0.010	
Permethrins	ND ppm	0.010		Piperonyl Butoxide	ND ppm	0.010	
Spinosyn A	ND ppm	0.010		Spiroxamine-1	ND ppm	0.010	
AbamectinB1a	ND ppm	0.010		Spinosyn D	ND ppm	0.010	
Mycotoxins							
Date Tested: 01/14/2022	Method: CB-SOP-025	Instrume	nt:				
Analyte	Result Units	LOQ	Result	Analyte	Result Units	LOQ	Resul
Ochratoxin A	ND ppm	0.010		Aflatoxin B1	ND ppm	0.010	
Aflatoxin G2	ND ppm	0.010		Aflatoxin B2	ND ppm	0.010	
Aflatoxin G1	ND ppm	0.010					
Metals							
Date Tested: 01/15/2022	Method: CB-SOP-027	Instrume	nt:				
Analyte	Result Units	LOQ	Result	Analyte	Result Units	LOQ	Resul
Arsenic	<loq ppm<="" td=""><td>0.200</td><td></td><td>Cadmium</td><td><loq ppm<="" td=""><td>0.200</td><td></td></loq></td></loq>	0.200		Cadmium	<loq ppm<="" td=""><td>0.200</td><td></td></loq>	0.200	
Lead	<loq ppm<="" td=""><td>0.200</td><td></td><td>Mercury</td><td><loq ppm<="" td=""><td>0.200</td><td></td></loq></td></loq>	0.200		Mercury	<loq ppm<="" td=""><td>0.200</td><td></td></loq>	0.200	
Microbial							
Date Tested: 01/19/2022	Method:	Instrume	nt:				
Analyte	Result Units	LOQ	Result	Analyte	Result Units	LOQ	Resu
STEC (E. coli)	Negative			Salmonella	Negative		
L. monocytogenes	Negative			Yeast/Mold (qPCR)	0 CFUs		
Residual Solvent							
Date Tested: 01/15/2022	Method: CB-SOP-032	Instrume	nt:				
Analyte	Result Units	LOQ	Result	Analyte	Result Units	LOQ	Resu
1-4 Dioxane	<loq ppm<="" td=""><td>29</td><td></td><td>2-Butanol</td><td><loq ppm<="" td=""><td>175</td><td></td></loq></td></loq>	29		2-Butanol	<loq ppm<="" td=""><td>175</td><td></td></loq>	175	
2-Ethoxyethanol	<loq ppm<="" td=""><td>24</td><td></td><td>2-Methylpentane</td><td><loq ppm<="" td=""><td>87</td><td></td></loq></td></loq>	24		2-Methylpentane	<loq ppm<="" td=""><td>87</td><td></td></loq>	87	
3-Methylpentane	<loq ppm<="" td=""><td>87</td><td></td><td>2-Propanol</td><td><loq ppm<="" td=""><td>350</td><td></td></loq></td></loq>	87		2-Propanol	<loq ppm<="" td=""><td>350</td><td></td></loq>	350	
Cyclohexane	<loq ppm<="" td=""><td>146</td><td></td><td>Ether</td><td><loq ppm<="" td=""><td>350</td><td></td></loq></td></loq>	146		Ether	<loq ppm<="" td=""><td>350</td><td></td></loq>	350	
Ethylbenzene	<loq ppm<="" td=""><td>81</td><td></td><td>Acetone</td><td><loq ppm<="" td=""><td>350</td><td></td></loq></td></loq>	81		Acetone	<loq ppm<="" td=""><td>350</td><td></td></loq>	350	
Isopropyl Acetate	<loq ppm<="" td=""><td>175</td><td></td><td>Methylbutane</td><td><loq ppm<="" td=""><td>350</td><td></td></loq></td></loq>	175		Methylbutane	<loq ppm<="" td=""><td>350</td><td></td></loq>	350	
n-Heptane	<loq ppm<="" td=""><td>350</td><td></td><td>n-Hexane</td><td><loq ppm<="" td=""><td>87</td><td></td></loq></td></loq>	350		n-Hexane	<loq ppm<="" td=""><td>87</td><td></td></loq>	87	
n-Pentane	<loq ppm<="" td=""><td>350</td><td></td><td>Tetrahydrofuran</td><td><loq ppm<="" td=""><td>54</td><td></td></loq></td></loq>	350		Tetrahydrofuran	<loq ppm<="" td=""><td>54</td><td></td></loq>	54	
Acetonitrile	<loq ppm<="" td=""><td>123</td><td></td><td>Ethanol</td><td><loq ppm<="" td=""><td>350</td><td></td></loq></td></loq>	123		Ethanol	<loq ppm<="" td=""><td>350</td><td></td></loq>	350	

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Sample ID: Sample Name: Sample Type:

D: 220114022 Name: HAR Focus Gummy Type: Edible

Certificate of Analysis

Residual Solvent							
Date Tested: 01/15/2022	Method: CB-SOP-032	Instrume	ent:				
Analyte	Result Units	LOQ	Result	Analyte	Result Units	LOQ	Result
Ethyl acetate	<loq ppm<="" td=""><td>175</td><td></td><td>o-Xylene</td><td><loq ppm<="" td=""><td>81</td><td></td></loq></td></loq>	175		o-Xylene	<loq ppm<="" td=""><td>81</td><td></td></loq>	81	
m+p-Xylene	<loq ppm<="" td=""><td>163</td><td></td><td>Methanol</td><td><loq ppm<="" td=""><td>250</td><td></td></loq></td></loq>	163		Methanol	<loq ppm<="" td=""><td>250</td><td></td></loq>	250	
Methylene Chloride	<loq ppm<="" td=""><td>90</td><td></td><td>Toluene</td><td><loq ppm<="" td=""><td>67</td><td></td></loq></td></loq>	90		Toluene	<loq ppm<="" td=""><td>67</td><td></td></loq>	67	
(A)			ized Signat	ure	Jamie Hobgood	01/19/2022	2:50 PM
PJLA Texting Accreditation V109558		Laborato	ry Manager			Date	Time

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