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Sample: 2312TLL0069.0196

## **Genesis Bioceuticals, LLC**

1120 W Watkins St Phoenix, AZ 85007 shonae.j@genbioaz.com (847) 682-4899 Lic. #00000058DCQU00115543 Harvest Dates: 11/22/2023

Strain: Tropic Banana Parent Batch #:; Batch#: TRBA-2179-20231122; Batch Size: 11 g Sample Received: 12/15/2023; Report Created: 12/21/2023; Expires: 12/21/2024 Manufacturing Date: Sampling: ; Environment:



## **Tropic Banana Flower/Pre-ROII**

Plant, Flower - Cured

Dispensary License #: ; Manufacturing License #: ; Cultivation License #:



### Cannabinoids

TPL_Potency_01				
24.80%	<loq< th=""><th></th><th>NT</th><th></th></loq<>		NT	
Total THC	Total CBD			
Analyte	LOQ	Mass	Mass Q	ualifier
	%	%	mg/g	
THCa	0.20	28.00	280.0	
Δ9-THC	0.20	0.24	2.4	
∆8-THC	0.20	ND	ND	
THCV	0.20	ND	ND	
CBDa	0.20	< LOQ	<loq< td=""><td></td></loq<>	
CBD	0.20	ND	ND	
CBDV	0.20	ND	ND	
CBN	0.20	ND	ND	
CBGa	0.20	1.02	10.2	
CBG	0.20	< LOQ	<loq< td=""><td></td></loq<>	
CBC	0.20	ND	ND	
Total	:	29.26	292.6	

Total THC = THCa \* 0.877 +  $\Delta$ 9-THC Total CBD = CBDa \* 0.877 + CBD Instrument: HPLC-DAD: ; Method: TPL\_Potency\_01



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**Brian DiMarco** 

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Laboratory Director

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Safety

1	Jaicty		
	<b>Pass</b> Pesticides	<b>Pass</b> Microbials	<b>Pass</b> Mycotoxins
ļ	resticides	I I I CI ODIAIS	IVIYCOLOXIIIS
	Complete	Pass	Not Tested
I	Solvents	Metals	Foreign Matter

### Ternenes

β-Caryophyllene1.294012.940Q2Linalool0.39503.950Q2δ-Limonene0.35003.500Q2Ocimene0.31403.140Q2Guaiol0.27002.700Q2β-Myrcene0.20202.020Q2	TPL Terpenes 01				
Analyte         Log         Mass         Mass         Qualifie $\alpha$ -Humulene         %         %         mg/g         g $\alpha$ -Humulene         1.5360         15.360         Q3 $\beta$ -Caryophyllene         1.2940         12.940         Q3           Linalool         0.3950         3.950         Q3 $\delta$ -Limonene         0.3140         3.140         Q3           Ocimene         0.2700         2.700         Q3           Guaiol         0.2200         2.020         Q3	<b>*</b>	Ŷ			
α-Humulene         %         mg/g $α$ -Humulene         1.5360         15.360         Q3 $β$ -Caryophyllene         1.2940         12.940         Q3           Linalool         0.3950         3.950         Q3           δ-Limonene         0.3500         3.500         Q3           Ocimene         0.3140         3.140         Q3           Guaiol         0.2700         2.700         Q3 $β$ -Myrcene         0.2020         2.020         Q3	Hops	Cinnamon		Lavender	
α-Humulene $1.5360$ $15.360$ Q3           β-Caryophyllene $1.2940$ $12.940$ Q3           Linalool $0.3950$ $3.950$ Q3           δ-Limonene $0.3500$ $3.500$ Q3           Ocimene $0.3140$ $3.140$ Q3           Guaiol $0.2700$ $2.700$ Q3           β-Myrcene $0.2020$ $2.020$ Q3	Analyte				ualifier
$\begin{array}{cccccccccccccccccccccccccccccccccccc$	$\begin{array}{l} \beta\text{-Caryophyllene}\\ \text{Linalool}\\ \delta\text{-Limonene}\\ \text{Ocimene}\\ \text{Guaiol}\\ \beta\text{-Myrcene}\\ \alpha\text{-Bisabolol}\\ \beta\text{-Pinene}\\ \text{Terpinolene}\\ \text{y-Terpinene}\\ \text{Eucalyptol}\\ \text{Nerolidol}\\ \text{trans-Nerolidol}\\ \alpha\text{-Pinene}\\ \text{Caryophyllene Oxide}\\ \text{Camphene}\\ 3\text{-Carene}\\ \alpha\text{-Terpinene}\\ \text{cis-Nerolidol}\\ \text{Geraniol}\\ \text{Isopulegol}\\ \text{p-Cymene}\\ \end{array}$		5360 15 2940 12 3950 3 3500 3 3140 3 2700 2 2020 2 2020 2 2020 2 2020 2 1380 1 0760 0 0670 0 0670 0 0480 0 0480 0 0470 0 04470 0 0470 0 0000000000	.360 .940 .950 .5500 .140 .700 .380 .760 .7700 .380 .760 .480 .470 .4470.4470	3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3
Total         5.0580         50.580           Instrument: GCMS; Method: TPL_Terp_01         Notes:         Notes:         Notes:	Instrument: GCMS; Method:	TPL_Terp_01 5.0	<u>)580 50</u>	.580	



Complete

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#### **Residual Solvents**

Residual Solvents	•		Complete
Analyte	LOQ	Limit	Mass StatusQualifier
	PPM	PPM	PPM
Acetone	400.0	1000.0	NR NT Q3
Acetonitrile	164.0	410.0	NR NT Q3
Benzene	0.8	2.0	NR NT Q3
n-Butane			NR NT Q3
Isobutane			NR NT Q3
Butanes	2000.0	5000.0	NR NT Q3
Chloroform	24.0	60.0	NR NT Q3
Dichloromethane	240.0	600.0	NR NT Q3
Ethanol	2000.0	5000.0	NR NT Q3
Ethyl-Acetate	2000.0	5000.0	NR NT Q3
Ethyl-Ether	2000.0	5000.0	NR NT Q3
Heptane	2000.0	5000.0	NR NT Q3
n-Hexane			NR NT Q3
2-Methyl-Pentane			NR NT Q3
3-Methyl-Pentane			NR NT Q3
2,2-Dimethyl-Butane			NR NT Q3
2,3-Dimethyl-Butane			NR NT Q3
Hexanes	116.0	290.0	NR NT Q3
Isopropyl-Acetate	2000.0	5000.0	NR NT Q3
Methanol	1200.0	3000.0	NR NT Q3
n-Pentane			NR NT Q3
Isopentane			NR NT Q3
Neopentane			NR NT Q3
Pentanes	2000.0	5000.0	NR NT Q3
2-Propanol	2000.0	5000.0	NR NT Q3
Propane	2000.0	5000.0	NR NT Q3
Toluene	356.0	890.0	NR NT Q3
1,2-Dimethyl-Benzene			NR NT Q3
1,3-Dimethyl-Benzene			NR NT Q3
1,4-Dimethyl-Benzene			NR NT Q3
Ethyl-Benzene			NR NT Q3
Xylenes	868.0	2170.0	NR NT Q3

Heavy Met	als				Pass
Analyte	LOQ	Limit	Mass	Status	Qualifier
	PPB	PPB	PPB		
Arsenic	200.0	400.0	ND	Pass	V1
Cadmium	200.0	400.0	<loq< th=""><th>Pass</th><th></th></loq<>	Pass	
Lead	500.0	1000.0	ND	Pass	V1
Mercury	100.0	200.0	<loq< th=""><th>Pass</th><th></th></loq<>	Pass	

Mycotoxins					Pass
Analyte	LOQ	Limit	Mass S	StatusC	) ualifier
	PPB	PPB	PPB		
B1	9	0	ND	Pass	
B2	9	0	ND	Pass	
G1	9	0	ND	Pass	11
G2	5	0	ND	Pass	
Ochratoxin A	9	20	ND	Pass	I1,R1
Total Aflatoxins	1	20	ND	Pass	

Instrument: HS-GCMS ; Method: TPL\_ResSolv\_01



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### **Microbials**

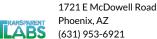


Pass

Analyte	LOQ	Limit	Result	Status	Qualifier
	CFU/g	CFU/g	CFU/g		
E. Coli	10	100	<10	Pass	

Analyte	Limit	Result	Status	Qualifier
Salmonella	Detectable in 1g	Not Detected	Pass	
Aspergillus	Detectable in 1g	Not Detected	Pass	
Aspergillus fumigatus	Detectable in 1g	Not Detected	Pass	
Aspergillus niger	Detectable in 1g	Not Detected	Pass	
Aspergillus flavus	Detectable in 1g	Not Detected	Pass	
Aspergillus terreus	Detectable in 1g	Not Detected	Pass	

Instrument: : Method:



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Pass

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### Pesticides TPL\_Pesticides\_01

Analyte	LOQ	Limit	Mass	Status Q	ualifier	Analyte	LOQ	Limit	Mass	Status Q	ualifier
	PPM	PPM	PPM				PPM	PPM	PPM		
Abamectin	0.23	0.50	ND	Pass		Hexythiazox	0.46	1.00	ND	Pass	M2
Acephate	0.19	0.40	ND	Pass		Imazalil	0.09	0.20	ND	Pass	
Acetamiprid	0.09	0.20	ND	Pass		Imidacloprid	0.19	0.40	ND	Pass	
Aldicarb	0.19	0.40	ND	Pass		Kresoxim	0.19	0.40	ND	Pass	
Azoxystrobin	0.09	0.20	ND	Pass		Methyl	0.17	0.40	ND	F 855	
Bifenazate	0.09	0.20	ND	Pass		Malathion	0.09	0.20	ND	Pass	M1
Bifenthrin	0.09	0.20	ND	Pass		Metalaxyl	0.09	0.20	ND	Pass	
Boscalid	0.19	0.40	ND	Pass	M1	Methiocarb	0.09	0.20	ND	Pass	M1
Carbaryl	0.09	0.20	ND	Pass		Methomyl	0.19	0.40	ND	Pass	
Carbofuran	0.09	0.20	ND	Pass		Myclobutanil	0.09	0.20	ND	Pass	M1
Chlorantraniliprole	0.09	0.20	ND	Pass	M1	Naled	0.23	0.50	ND	Pass	
Chlorfenapyr	0.46	1.00	ND	Pass	I1,M2	Oxamyl	0.46	1.00	ND	Pass	
Chlorpyrifos	0.09	0.20	ND	Pass	M2	Paclobutrazol	0.19	0.40	ND	Pass	
Clofentezine	0.09	0.20	ND	Pass		Permethrin	0.09	0.20	ND	Pass	
Cyfluthrin	0.46	1.00	ND	Pass		Phosmet	0.09	0.20	ND	Pass	M1
Cypermethrin	0.46	1.00	ND	Pass		Piperonyl	0.93	2.00	<loq< td=""><td>Pass</td><td>M2</td></loq<>	Pass	M2
Daminozide	0.46	1.00	ND	Pass		Butoxide	0.75	2.00	LUQ	газэ	1112
Diazinon	0.09	0.20	ND	Pass		Prallethrin	0.09	0.20	ND	Pass	
Dichlorvos	0.05	0.10	ND	Pass	11	Propiconazole	0.19	0.40	ND	Pass	
Dimethoate	0.09	0.20	ND	Pass		Propoxur	0.09	0.20	ND	Pass	M1
Ethoprophos	0.09	0.20	ND	Pass	M1	Pyrethrins	0.46	1.00	ND	Pass	M1
Etofenprox	0.19	0.40	ND	Pass		Pyridaben	0.09	0.20	ND	Pass	M2
Etoxazole	0.09	0.20	ND	Pass		Spinosad	0.09	0.20	ND	Pass	
Fenoxycarb	0.09	0.20	ND	Pass		Spiromesifen	0.09	0.20	ND	Pass	
Fenpyroximate	0.19	0.40	ND	Pass		Spirotetramat	0.09	0.20	ND	Pass	
Fipronil	0.19	0.40	ND	Pass	M1	Spiroxamine	0.19	0.40	ND	Pass	
Flonicamid	0.46	1.00	ND	Pass		Tebuconazole	0.19	0.40	ND	Pass	
Fludioxonil	0.19	0.40	ND	Pass	M1	Thiacloprid	0.09	0.20	ND	Pass	
						Thiamethoxam	0.09	0.20	ND	Pass	
						Trifloxystrobin	0.09	0.20	ND	Pass	

Instrument: LC-QQQ



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B1 = Target analyte detected in calibration blank was above LOQ but the concentration of cannabinoid was blow LOQ,

B2 = Target analyte detected in calibration blank was above LOQ but was below the maximum allowable concentration.

D1 = The limit of quantitation and the sample results were adjusted to reflect sample dilution,

I1 = The relative intensity of a characteristic ion in a sample analyte exceeded the acceptance criteria with respect to the reference spectra, indicating interference,

L1 = The percent recovery of a laboratory control sample is greater than the acceptance limits in A.A.C 17 R9-17-404.03(K)(2)(C), but the sample's target analytes were not detected above the maximum allowed concentration,

M1 = The recovery from the matrix spike was high, but the recovery from the laboratory control sample was within acceptance criteria,

M2 = The recovery from the matrix spike was low, but the recovery from the laboratory control sample was within acceptance criteria,

M3 = The recovery from the matrix spike was unusable because the analyte concentration was disproportionate to the spike level, but the recovery from the laboratory control sample was within acceptance criteria,

M4 = The analysis of a spiked sample required a dilution such that the spike recovery calculation does not provide useful information, but the recovery from the associated laboratory control sample was within acceptance criteria,

M5 = The analyte concentration was determined by the method of standard addition, in which the standard is added directly to the aliquots of the analyzed sample,

N1 - A description of the variance is described in the final report of testing,

R1 = The relative percent difference for the laboratory control sample and duplicate exceeded the limit in A.A.C 17 R9-17-404.03(K)(3), but the recover in subsection A.A.C 17 R9-17-404.03 (K)(2) was within accepted criteria,

R2 = The relative percent difference for a sample and duplicated exceeded the limit in subsection A.A.C 17 R9-17-404.03 (0)

Q1 = Sample integrity was not maintained,

Q2 = The sample is heterogenous and sample homogeneity could not be readily achieved using routine laboratory practices

Q3 = Testing result is for informational purposes only and cannot be used to satisfy dispensary testing requirements in R9-17-317.01(A) or labeling requirements in R9-17-317

V1 = The recovery from continuing calibration verification standards exceeded the acceptance limits denoted in A.C.C 17 R9-17-403.03(J)(1)(b), but the sample's target analytes were not detected above the maximum allowable concentrations for the analytes in the sample.

TLABS

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