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Genesis Bioceuticals, LLC

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Lic. #00000058DCQU00115543

Harvest Dates: 10/02/2023

Tropic Banana Shatter

Concentrates & Extracts, Shatter, Extraction Method: Butane Dispensary License #:; Manufacturing License #:; Cultivation License #:



Strain: Tropic Banana Shatter

Parent Batch #: ; Batch #: TRBA-2214-20231207-s; Batch Size: 11 g Sample Received: 12/27/2023; Report Created: 12/31/2023; Expires: 12/31/2024

Manufacturing Date: 12/07/2023

Sampling: ; Environment:





Safety

Pass **Pass Pass Pesticides** Microbials **Mycotoxins Pass Pass Not Tested** Solvents Metals Foreign Matter

Cannabinoids

| alyte | LOQ | Mass | Mass Qualifi |
|-----------|-----------|------|--------------|
| Total THC | Total CBD | | Moisture |
| 77.83% | 0.14% | | NT |
| | | | |

| | | | | 40 |
|---------|------|--|--------------------------------|-----------|
| Analyte | LOQ | Mass | Mass | Qualifier |
| | % | % | mg/g | |
| THCa | 0.10 | 87.29 | 872.9 | M2 |
| Δ9-THC | 0.10 | 1.27 | 12.7 | M2 |
| Δ8-THC | 0.10 | ND | ND | M2 |
| THCV | 0.10 | <loq< th=""><th><loq< th=""><th>M2</th></loq<></th></loq<> | <loq< th=""><th>M2</th></loq<> | M2 |
| CBDa | 0.10 | 0.17 | 1.7 | M2 |
| CBD | 0.10 | ND | ND | M2 |
| CBDV | 0.10 | ND | ND | M2 |
| CBN | 0.10 | ND | ND | M2 |
| CBGa | 0.10 | 3.51 | 35.1 | M2 |
| CBG | 0.10 | 0.40 | 4.0 | M2 |
| CBC | 0.10 | ND | ND | M2 |
| Total | | 92.64 | 926.4 | |

Total THC = THCa * 0.877 + Δ 9-THC Total CBD = CBDa * 0.877 + CBD

Instrument: HPLC-DAD: ; Method: TPL_Potency_01

Terpenes TPL_Terpenes_01







| Analyte LOC | Q Mass | Mass | Qualifier |
|---------------------------------------|--------|--------|-----------|
| | % | mg/g | |
| α-Humulene | 1.5916 | 15.916 | Q3 |
| β-Caryophyllene | 1.3400 | 13.400 | Q3 |
| δ-Limonene | 0.7035 | 7.035 | Q3 |
| Linalool | 0.6535 | 6.535 | Q3 |
| Ocimene | 0.6060 | 6.060 | Q3 |
| β-Myrcene | 0.4056 | 4.056 | Q3 |
| Guaiol | 0.2737 | 2.737 | Q3 |
| β-Pinene | 0.2059 | 2.059 | Q3 |
| α-Bisabolol | 0.1909 | 1.909 | Q3 |
| Nerolidol | 0.1735 | 1.735 | Q3 |
| Caryophyllene Oxide | 0.1464 | 1.464 | Q3 |
| trans-Nerolidol | 0.1085 | 1.085 | Q3 |
| α-Pinene | 0.0927 | 0.927 | Q3 |
| Terpinolene | 0.0887 | 0.887 | Q3 |
| y-Terpinene | 0.0700 | 0.700 | Q3 |
| cis-Nerolidol | 0.0650 | 0.650 | Q3 |
| Eucalyptol | 0.0487 | 0.487 | Q3 |
| Camphene | 0.0288 | 0.288 | Q3 |
| Isopulegol | 0.0068 | 0.068 | Q3 |
| 3-Carene | < | < | Q3 |
| α-Terpinene | < | < | Q3 |
| Geraniol | < | < | Q3 |
| p-Cymene | < | < | Q3 |
| Total | 6.7998 | 67.998 | |
| Instrument: GCMS; Method: TPL Terp 01 | · | | |



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Brian DiMarco **Laboratory Director**

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Harvest Dates: 10/02/2023

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Sample: 2312TLL0080.0251

Strain: Tropic Banana Shatter

Parent Batch #:; Batch #: TRBA-2214-20231207-s; Batch Size: 11 g

Sample Received: 12/27/2023; Report Created: 12/31/2023; Expires: 12/31/2024 Manufacturing Date: 12/07/2023

Sampling: ; Environment:



Pass

Residual Solvents Pass

| Analyte | LOQ | Limit | Mass StatusQualifie |
|----------------------|--------|--------|----------------------------|
| | PPM | PPM | PPM |
| Acetone | 200.0 | 1000.0 | ND Pass |
| Acetonitrile | 82.0 | 410.0 | ND Pass |
| Benzene | 0.4 | 2.0 | ND Pass |
| n-Butane | | | ND Tested |
| Isobutane | | | ND Tested |
| Butanes | 500.0 | 5000.0 | <loq pass<="" th=""></loq> |
| Chloroform | 12.0 | 60.0 | ND Pass |
| Dichloromethane | 120.0 | 600.0 | ND Pass |
| Ethanol | 1000.0 | 5000.0 | ND Pass |
| Ethyl-Acetate | 1000.0 | 5000.0 | ND Pass |
| Ethyl-Ether | 1000.0 | 5000.0 | ND Pass |
| Heptane | 1000.0 | 5000.0 | ND Pass |
| n-Hexane | | | ND Tested |
| 2-Methyl-Pentane | | | ND Tested |
| 3-Methyl-Pentane | | | ND Tested |
| 2,2-Dimethyl-Butane | | | ND Tested |
| 2,3-Dimethyl-Butane | | | ND Tested |
| Hexanes | 145.0 | 290.0 | ND Pass |
| Isopropyl-Acetate | 1000.0 | 5000.0 | ND Pass |
| Methanol | 600.0 | 3000.0 | ND Pass |
| n-Pentane | | | ND Tested |
| Isopentane | | | ND Tested |
| Neopentane | | | ND Tested |
| Pentanes | 1000.0 | 5000.0 | ND Pass |
| 2-Propanol | 1000.0 | 5000.0 | <loq pass<="" th=""></loq> |
| Propane | | 5000.0 | ND Pass |
| Toluene | 178.0 | 890.0 | ND Pass |
| 1,2-Dimethyl-Benzene | | | ND Tested |
| 1,3-Dimethyl-Benzene | | | ND Tested |
| 1,4-Dimethyl-Benzene | | | ND Tested |
| Ethyl-Benzene | | | ND Tested |
| Xylenes | 868.0 | 2170.0 | ND Pass |

| Heavy Metals | | | | | Pass |
|--------------|-----|-------|------|--------|-----------|
| Analyte | LOQ | Limit | Mass | Status | Qualifier |
| | PPB | PPB | PPB | | |

| | PPB | PPB | PPB | | |
|---------|-------|--------|---|------|------------|
| Arsenic | 200.0 | 400.0 | ND | Pass | |
| Cadmium | 200.0 | 400.0 | ND | Pass | |
| Lead | 500.0 | 1000.0 | ND | Pass | L1 |
| Mercury | 100.0 | 200.0 | <loq< td=""><td>Pass</td><td>B2, L1, M2</td></loq<> | Pass | B2, L1, M2 |
| - | | | | | |

Mycotoxins **Pass** Analyte Mass StatusQualifier PPB В1 ND 11. M2 10 Pass 11, M2 B2 10 0 ND Pass G1 10 0 ND Pass I1, M2 Pass R1 G2 10 ND Ochratoxin A 10 20 ND Pass I1, V1 I1, M2 9 **Total Aflatoxins** 20 ND Pass R1

Instrument: HS-GCMS



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Sample Receiv

Sample: 2312TLL0080.0251

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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: qPCR/Plating



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Pesticides TPL_Pesticides_01

Pass

| Analyte | LOQ | Limit | Mass | Status C | Qualifier | Analyte | LOQ | Limit | Mass | Status Q | ualifier |
|---------------------|------|-------|------|----------|-----------|-----------------|------|-------|---|----------|----------|
| | PPM | PPM | PPM | | | | PPM | PPM | PPM | | |
| Abamectin | 0.24 | 0.50 | ND | Pass | M2 | Hexythiazox | 0.47 | 1.00 | ND | Pass | M2 |
| Acephate | 0.19 | 0.40 | ND | Pass | | lmazalil | 0.10 | 0.20 | ND | Pass | |
| Acetamiprid | 0.10 | 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 | M2 |
| Azoxystrobin | 0.10 | 0.20 | ND | Pass | | Methyl | 0.17 | 0.40 | ND | F d 5 5 | IVIZ |
| Bifenazate | 0.10 | 0.20 | ND | Pass | | Malathion | 0.10 | 0.20 | ND | Pass | |
| Bifenthrin | 0.10 | 0.20 | ND | Pass | M2 | Metalaxyl | 0.10 | 0.20 | ND | Pass | |
| Boscalid | 0.19 | 0.40 | ND | Pass | M2 | Methiocarb | 0.10 | 0.20 | ND | Pass | |
| Carbaryl | 0.10 | 0.20 | ND | Pass | | Methomyl | 0.19 | 0.40 | ND | Pass | |
| Carbofuran | 0.10 | 0.20 | ND | Pass | | Myclobutanil | 0.10 | 0.20 | ND | Pass | M2 |
| Chlorantraniliprole | 0.10 | 0.20 | ND | Pass | | Naled | 0.24 | 0.50 | ND | Pass | M2 |
| Chlorfenapyr | 0.47 | 1.00 | ND | Pass | I1, M2 | Oxamyl | 0.47 | 1.00 | ND | Pass | |
| Chlorpyrifos | 0.10 | 0.20 | ND | Pass | M2 | Paclobutrazol | 0.19 | 0.40 | ND | Pass | |
| Clofentezine | 0.10 | 0.20 | ND | Pass | M2 | Permethrin | 0.10 | 0.20 | ND | Pass | M2 |
| Cyfluthrin | 0.47 | 1.00 | ND | Pass | M2 | Phosmet | 0.10 | 0.20 | ND | Pass | M2 |
| Cypermethrin | 0.47 | 1.00 | ND | Pass | M2 | Piperonyl | 0.95 | 2.00 | <loq< td=""><td>Pass</td><td>M2</td></loq<> | Pass | M2 |
| Daminozide | 0.47 | 1.00 | ND | Pass | | Butoxide | 0.95 | 2.00 | LOQ | Pass | IVIZ |
| Diazinon | 0.10 | 0.20 | ND | Pass | | Prallethrin | 0.10 | 0.20 | ND | Pass | |
| Dichlorvos | 0.05 | 0.10 | ND | Pass | M2 | Propiconazole | 0.19 | 0.40 | ND | Pass | |
| Dimethoate | 0.10 | 0.20 | ND | Pass | | Propoxur | 0.10 | 0.20 | ND | Pass | |
| Ethoprophos | 0.10 | 0.20 | ND | Pass | M2 | Pyrethrins | 0.40 | 1.00 | <loq< td=""><td>Pass</td><td></td></loq<> | Pass | |
| Etofenprox | 0.19 | 0.40 | ND | Pass | M2 | Pyridaben | 0.10 | 0.20 | ND | Pass | M2 |
| Etoxazole | 0.10 | 0.20 | ND | Pass | | Spinosad | 0.10 | 0.20 | ND | Pass | M2 |
| Fenoxycarb | 0.10 | 0.20 | ND | Pass | M2 | Spiromesifen | 0.10 | 0.20 | ND | Pass | |
| - enpyroximate | 0.19 | 0.40 | ND | Pass | M2 | Spirotetramat | 0.10 | 0.20 | ND | Pass | |
| Fipronil | 0.19 | 0.40 | ND | Pass | I1 | Spiroxamine | 0.19 | 0.40 | ND | Pass | |
| Flonicamid | 0.47 | 1.00 | ND | Pass | | Tebuconazole | 0.19 | 0.40 | ND | Pass | M2 |
| Fludioxonil | 0.19 | 0.40 | ND | Pass | M2 | Thiacloprid | 0.10 | 0.20 | ND | Pass | M2 |
| | | | | | | Thiamethoxam | 0.10 | 0.20 | ND | Pass | |
| | | | | | | Trifloxystrobin | 0.10 | 0.20 | ND | Pass | M2 |

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 (O)
- 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.

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