

**Rainbow Sherbet D8**

Sample ID: SA-260622-82927  
 Batch: RS8-5.18.26  
 Type: Finished Product - Inhalable  
 Matrix: Concentrate - Vape  
 Unit Size (g):  
 Unit Volume (mL): , Density (g/mL):

Received: 06/29/2026  
 Completed: 07/10/2026

**Client**  
 TribeTokes  
 242 W 38th St  
 New York, NY 10018  
 USA


**Summary**

**Test**  
 Cannabinoids

**Date**  
 07/10/2026

**Status**  
 Tested

|              |               |                    |                   |                   |                                 |
|--------------|---------------|--------------------|-------------------|-------------------|---------------------------------|
| <b>ND</b>    | <b>83.7 %</b> | <b>91.0 %</b>      | <b>Not Tested</b> | <b>Not Tested</b> | <b>Yes</b>                      |
| Total Δ9-THC | Δ8-THC        | Total Cannabinoids | Moisture Content  | Foreign Matter    | Internal Standard Normalization |

**Cannabinoids by HPLC-PDA and GC-MS/MS**

| Analyte      | LOD (%) | LOQ (%) | Result (%) | Result (mg/g) |
|--------------|---------|---------|------------|---------------|
| CBC          | 0.0095  | 0.0284  | ND         | ND            |
| CBCA         | 0.0181  | 0.0543  | ND         | ND            |
| CBCV         | 0.006   | 0.018   | ND         | ND            |
| CBD          | 0.0081  | 0.0242  | ND         | ND            |
| CBD A        | 0.0043  | 0.013   | ND         | ND            |
| CBD B        | 0.0133  | 0.04    | ND         | ND            |
| CBD-C8       | 0.0133  | 0.04    | ND         | ND            |
| CBDH         | 0.0133  | 0.04    | ND         | ND            |
| CBDP         | 0.0133  | 0.04    | ND         | ND            |
| CBDV         | 0.0061  | 0.0182  | ND         | ND            |
| CBDVA        | 0.0021  | 0.0063  | ND         | ND            |
| CBG          | 0.0057  | 0.0172  | ND         | ND            |
| CBGA         | 0.0049  | 0.0147  | ND         | ND            |
| CBL          | 0.0112  | 0.0335  | 0.105      | 1.05          |
| CBLA         | 0.0124  | 0.0371  | ND         | ND            |
| CBN          | 0.0056  | 0.0169  | 0.958      | 9.58          |
| CBNA         | 0.006   | 0.0181  | ND         | ND            |
| CBP          | 0.0133  | 0.04    | ND         | ND            |
| CBT          | 0.018   | 0.054   | 0.234      | 2.34          |
| Δ4,8-iso-THC | 0.0133  | 0.04    | 2.18       | 21.8          |
| Δ6a,10a-THC  | 0.0133  | 0.04    | ND         | ND            |
| Δ8-iso-THC   | 0.0133  | 0.04    | 2.45       | 24.5          |
| Δ8-THC       | 0.0104  | 0.0312  | 83.7       | 837           |
| Δ8-THCB      | 0.0133  | 0.04    | ND         | ND            |
| Δ8-THC-C8    | 0.0133  | 0.04    | ND         | ND            |

Cannabinoid results continued on next page...



Generated By: Christopher Kyzar  
 QA/QC Manager  
 Date: 07/10/2026





**KCA Laboratories**  
1509 Bull Lea Rd, Ste 100  
Lexington, KY 40511

+1-833-KCA-LABS  
<https://kcalabs.com>  
KDA Lic.# P\_0058

## Certificate of Analysis

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### Rainbow Sherbet D8

Sample ID: SA-260622-82927  
Batch: RS8-5.18.26  
Type: Finished Product - Inhalable  
Matrix: Concentrate - Vape  
Unit Size (g):  
Unit Volume (mL): , Density (g/mL):

Received: 06/29/2026  
Completed: 07/10/2026

**Client**  
TribeTokes  
242 W 38th St  
New York, NY 10018  
USA

| Analyte             | LOD (%) | LOQ (%) | Result (%)  | Result (mg/g) |
|---------------------|---------|---------|-------------|---------------|
| Δ8-THCH             | 0.0133  | 0.04    | ND          | ND            |
| Δ8-THCP             | 0.0133  | 0.04    | ND          | ND            |
| Δ8-THCV             | 0.0133  | 0.04    | 1.43        | 14.3          |
| Δ9-THC              | 0.0076  | 0.0227  | ND          | ND            |
| Δ9-THCA             | 0.0084  | 0.0251  | ND          | ND            |
| Δ9-THCB             | 0.0133  | 0.04    | ND          | ND            |
| Δ9-THC-C8           | 0.0133  | 0.04    | ND          | ND            |
| Δ9-THCH             | 0.0133  | 0.04    | ND          | ND            |
| Δ9-THCP             | 0.0133  | 0.04    | ND          | ND            |
| Δ9-THCV             | 0.0069  | 0.0206  | ND          | ND            |
| Δ9-THCVA            | 0.0062  | 0.0186  | ND          | ND            |
| (6aR,9R)-Δ10-THC    | 0.0133  | 0.04    | ND          | ND            |
| (6aR,9S)-Δ10-THC    | 0.0133  | 0.04    | ND          | ND            |
| exo-THC             | 0.0133  | 0.04    | ND          | ND            |
| (6aR,9R,10aR)-HHC   | 0.0133  | 0.04    | ND          | ND            |
| (6aR,9S,10aR)-HHC   | 0.0133  | 0.04    | ND          | ND            |
| <b>Total Δ9-THC</b> |         |         | <b>ND</b>   | <b>ND</b>     |
| <b>Total</b>        |         |         | <b>91.0</b> | <b>910</b>    |

ND = Not Detected; NR = (Spike) Not Recoverable, sample matrix interference present which may affect accuracy of results; NT = Not Tested; UA = Unsuitable for Analysis; LOD = Limit of Detection; LOQ = Limit of Quantitation; RL = Reporting Limit; Δ = Delta; Total Δ9-THC = Δ9-THCA \* 0.877 + Δ9-THC; Total CBD = CBDA \* 0.877 + CBD;

Generated By: Christopher Kyzar  
QA/QC Manager  
Date: 07/10/2026

Tested By: Kelsey Rogers  
Senior Scientist  
Date: 07/10/2026



ISO/IEC 17025:2017 Accredited  
Accreditation #108651



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# Gobi Hemp

## Amended Report For: Analytical Report 2605120002-V1 - CDPHE Certified Certificate of Analysis



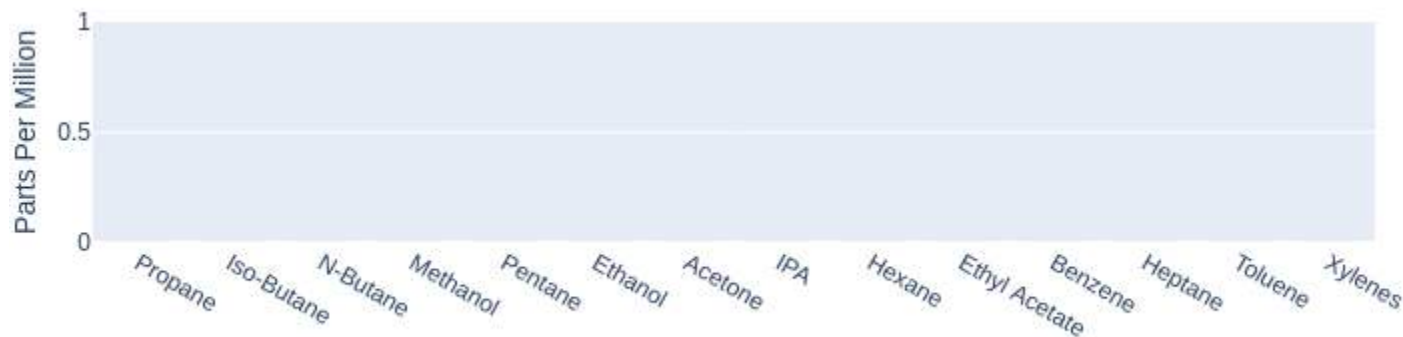
**Manifest:** 2605120002  
**Sample ID:** 1A-GHEMP-2605120002-0001  
**Sample Name:** 344026031632C - FM-942-811  
**Sample Type:** D8 - Distillate  
**Client ID:** CID-00207  
**Client:** FM Labs  
**Address:** 1285 Factory Circle, Fort Lupton, CO 80621

**Test Performed:** RSA  
**Report No:** A-R-2605120002-V1  
**Receive Date:** 2026-05-12  
**Test Date:** 2026-05-14  
**Report Date:** 2026-05-28  
**Sample Condition:** Good  
**Method Reference:** GH-OP-08

**Scope:** The content of fifteen residual solvents was determined by an in-house developed method for Headspace-Gas Chromatography with Flame Ionization Detection.

| Solvents      | LOD (ppm) | LOQ (ppm) | Parts Per Million (ppm) |
|---------------|-----------|-----------|-------------------------|
| Propane       | 47.0      | 142.3     | ND                      |
| Iso-Butane    | 55.5      | 168.0     | ND                      |
| N-Butane      | 68.1      | 206.4     | ND                      |
| Methanol      | 34.8      | 105.4     | ND                      |
| Pentane       | 64.8      | 196.4     | ND                      |
| Ethanol       | 87.8      | 266.1     | ND                      |
| Acetone       | 71.4      | 216.4     | ND                      |
| IPA           | 86.3      | 261.5     | ND                      |
| Hexane        | 0.6       | 35.0      | ND                      |
| Ethyl Acetate | 71.6      | 217.0     | ND                      |
| Benzene       | 0.3       | 1.0       | ND                      |
| Heptane       | 58.8      | 178.2     | ND                      |
| Toluene       | 6.5       | 94.3      | ND                      |
| Xylenes       | 7.8       | 185.9     | ND                      |

ND - not detected; LOD - limit of detection; LOQ - limit of quantitation; ULOQ - upper limit of quantitation;  
 \*Estimated result, greater than the upper limit of quantitation (>ULOQ)



### Lab Comments:

*Walter Marsh*

Walter Marsh Lead Research Lab Analyst

2026-05-28

Date



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**Manifest:** 2605120002  
**Sample ID:** 1A-GHEMP-2605120002-0001  
**Sample Name:** 344026031632C - FM-942-811  
**Sample Type:** D8 - Distillate  
**Client ID:** CID-00207  
**Client:** FM Labs  
**Facility Address:** 1285 Factory Circle, Fort Lupton, CO 80621

**Test Performed:** Pesticide  
**Report No:** A-PE-2605120002-V1  
**Receive Date:** 2026-05-12  
**Test Date:** 2026-05-14  
**Report Date:** 2026-05-28  
**Sample Condition:** Good  
**Method Reference:** GA-OP-11

### Executive Summary:

Sample 1A-GHEMP-2605120002-0001 has **passed** pesticide testing.

The following pesticides were detected in the sample:

### Scope:

The content of the reported pesticide residues were quantified using LC-MS-MS and GC-TQMS. Identification was based on the retention time of each compound and the product mass spectra generated using Single Reaction Monitoring (SRM) or Dramatic Multiple Reaction Monitoring, and quantitation was determined using external standard calibration.

### Lab Comments:

Walter Marsh Lead Research Lab Analyst

2026-05-28

Date



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# Gobi Hemp

## Amended Report For: Pesticide Residues

### Report 2605120002-V1



| Pesticide           | Limits (ppm) |            | Result (ppm) |      | Pesticide          | Limits (ppm) |            | Result (ppm) |      | Pesticide          | Limits (ppm) |            | Result (ppm) |      |
|---------------------|--------------|------------|--------------|------|--------------------|--------------|------------|--------------|------|--------------------|--------------|------------|--------------|------|
|                     | Regulatory   | Reporting* |              |      |                    | Regulatory   | Reporting* |              |      |                    | Regulatory   | Reporting* |              |      |
| Abamectin           |              | 0.10000    | ND           | LCMS | Dodemorph          |              | 0.10000    | ND           | LCMS | Oxamyl             |              | 1.50000    | ND           | LCMS |
| Acephate            |              | 0.10000    | ND           | LCMS | Endosulfan sulfate |              | 0.10000    | ND           | GCMS | Paclobutrazol      |              | 0.10000    | ND           | LCMS |
| Acequinocyl         |              | 0.10000    | ND           | LCMS | Endosulfan-alpha   |              | 0.20000    | ND           | GCMS | Parathion-methyl   |              | 0.10000    | ND           | GCMS |
| Acetamiprid         |              | 0.10000    | ND           | LCMS | Endosulfan-beta    |              | 0.10000    | ND           | GCMS | Permethrins        |              | 0.50000    | ND           | LCMS |
| Aldicarb            |              | 0.10000    | ND           | LCMS | Ethoprophos        |              | 0.10000    | ND           | LCMS | Phenothrin         |              | 0.10000    | ND           | LCMS |
| Allethrin           |              | 0.10000    | ND           | LCMS | Etofenprox         |              | 0.10000    | ND           | LCMS | Phosmet            |              | 0.10000    | ND           | LCMS |
| Atrazine            |              | 0.10000    | ND           | LCMS | Etoazole           |              | 0.10000    | ND           | LCMS | Piperonyl butoxide |              | 1.00000    | ND           | LCMS |
| Azadirachtin        |              | 0.50000    | ND           | LCMS | Etridiazole        |              | 0.10000    | ND           | GCMS | Pirimicarb         |              | 0.10000    | ND           | LCMS |
| Azoxystrobin        |              | 0.10000    | ND           | LCMS | Fenhexamid         |              | 0.12500    | ND           | LCMS | Prallethrin        |              | 0.10000    | ND           | LCMS |
| Benzovindiflupyr    |              | 0.10000    | ND           | LCMS | Fenoxycarb         |              | 0.10000    | ND           | LCMS | Propiconazole      |              | 0.10000    | ND           | LCMS |
| Bifenazate          |              | 0.10000    | ND           | LCMS | Fenpyroximate      |              | 0.10000    | ND           | LCMS | Propoxur           |              | 0.10000    | ND           | LCMS |
| Bifenthrin          |              | 1.00000    | ND           | LCMS | Fensulfothion      |              | 0.10000    | ND           | LCMS | Pyraclostrobin     |              | 0.10000    | ND           | LCMS |
| Boscalid            |              | 0.10000    | ND           | LCMS | Fenthion           |              | 0.10000    | ND           | GCMS | Pyrethrins         |              | 0.10000    | ND           | LCMS |
| Buprofezin          |              | 0.10000    | ND           | LCMS | Fenvalerate        |              | 0.10000    | ND           | GCMS | Pyridaben          |              | 0.10000    | ND           | LCMS |
| Carbaryl            |              | 0.10000    | ND           | LCMS | Fipronil           |              | 0.10000    | ND           | LCMS | Pyriproxyfen       |              | 0.10000    | ND           | LCMS |
| Carbofuran          |              | 0.10000    | ND           | LCMS | Flonicamid         |              | 0.10000    | ND           | LCMS | Quintozene         |              | 0.10000    | ND           | GCMS |
| Chlorantraniliprole |              | 0.10000    | ND           | LCMS | Fludioxonil        |              | 0.10000    | ND           | LCMS | Resmethrin         |              | 0.10000    | ND           | LCMS |
| Chlorphenapyr       |              | 0.10000    | ND           | GCMS | Fluopyram          |              | 0.10000    | ND           | LCMS | Spinetoram         |              | 0.10000    | ND           | LCMS |
| Chlorpyrifos        |              | 0.10000    | ND           | LCMS | Hexythiazox        |              | 0.10000    | ND           | LCMS | Spinosad           |              | 0.10000    | ND           | LCMS |
| Clofentezine        |              | 0.10000    | ND           | LCMS | Imazalil           |              | 0.10000    | ND           | LCMS | Spirodiclofen      |              | 0.25000    | ND           | LCMS |
| Clothianidin        |              | 0.10000    | ND           | LCMS | Imidacloprid       |              | 0.10000    | ND           | LCMS | Spiromesifen       |              | 3.00000    | ND           | LCMS |
| Coumaphos           |              | 0.10000    | ND           | LCMS | Iprodione          |              | 0.50000    | ND           | LCMS | Spirotetramat      |              | 0.10000    | ND           | LCMS |
| Cyantraniliprole    |              | 0.10000    | ND           | LCMS | Kinoprene          |              | 0.10000    | ND           | GCMS | Spiroxamine        |              | 0.10000    | ND           | LCMS |
| Cyfluthrin          |              | 0.20000    | ND           | GCMS | Kresoxim-methyl    |              | 0.10000    | ND           | LCMS | Tebuconazole       |              | 0.10000    | ND           | LCMS |
| Cypermethrin        |              | 0.25000    | ND           | GCMS | MGK-264            |              | 0.10000    | ND           | GCMS | Tebufenozide       |              | 0.10000    | ND           | LCMS |
| Cyprodinil          |              | 0.10000    | ND           | LCMS | Malathion          |              | 0.10000    | ND           | LCMS | Teflubenzuron      |              | 0.10000    | ND           | LCMS |
| Daminozide          |              | 0.10000    | ND           | LCMS | Metalaxyl          |              | 0.10000    | ND           | LCMS | Tetrachlorvinphos  |              | 0.10000    | ND           | LCMS |
| Deltamethrin        |              | 0.50000    | ND           | LCMS | Methiocarb         |              | 0.10000    | ND           | LCMS | Tetramethrin       |              | 0.10000    | ND           | LCMS |
| Diazinon            |              | 0.10000    | ND           | LCMS | Methomyl           |              | 0.10000    | ND           | LCMS | Thiabendazole      |              | 0.10000    | ND           | LCMS |
| Dichlorvos          |              | 0.10000    | ND           | GCMS | Methoprene         |              | 2.00000    | ND           | LCMS | Thiacloprid        |              | 0.10000    | ND           | LCMS |
| Dimethoate          |              | 0.10000    | ND           | LCMS | Mevinphos          |              | 0.10000    | ND           | LCMS | Thiamethoxam       |              | 0.10000    | ND           | LCMS |
| Dimethomorph        |              | 0.10000    | ND           | LCMS | Myclobutanil       |              | 0.10000    | ND           | LCMS | Thiophanate-methyl |              | 0.10000    | ND           | LCMS |
| Dinotefuran         |              | 0.10000    | ND           | LCMS | Naled              |              | 0.10000    | ND           | LCMS | Trifloxystrobin    |              | 0.10000    | ND           | LCMS |
| Diuron              |              | 0.10000    | ND           | LCMS | Novaluron          |              | 0.10000    | ND           | LCMS | lambda-Cyhalothrin |              | 0.20000    | ND           | GCMS |

\*or Lower Limit of Quantitation (LLOQ).  
 ND (Not Detected) = sample result is below MDL.  
 >HLOQ = sample result is above Higher LOQ.  
 \*\*

*Walter Marsh*

Walter Marsh Lead Research Lab Analyst

2026-05-28

Date



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## Amended Report For: Analytical Report 2605120002-V1 - CDPHE Certified Certificate of Analysis



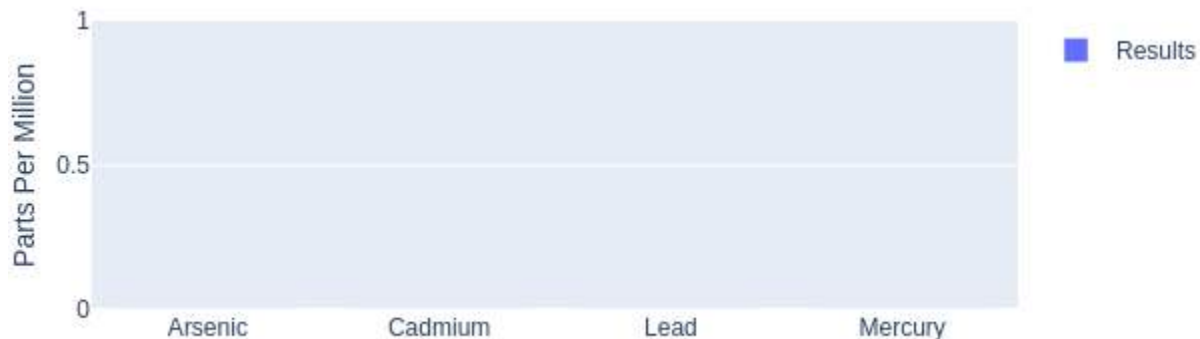
**Manifest:** 2605120002  
**Sample ID:** 1A-GHEMP-2605120002-0001  
**Sample Name:** 344026031632C - FM-942-811  
**Sample Type:** D8 - Distillate  
**Client ID:** CID-00207  
**Client:** FM Labs  
**Address:** 1285 Factory Circle, Fort Lupton, CO 80621

**Test Performed:** Elemental impurities  
**Intended Use:** Inhaled or Audited Product  
**Report No:** A-MT-2605120002-V1  
**Receive Date:** 2026-05-12  
**Test Date:** 2026-05-15  
**Report Date:** 2026-05-28  
**Sample Condition:** Good  
**Method Reference:** GH-OP-17

**Scope:** Arsenic, Cadmium, Lead and Mercury were determined by an Inductively Coupled Plasma Mass Spectrometer (ICP-MS) using an in-house developed method.

| Elemental Impurities | LOD (ppm) | LOQ (ppm) | Parts Per Million (ppm) |
|----------------------|-----------|-----------|-------------------------|
| Arsenic              | 0.007     | 0.025     | ND                      |
| Cadmium              | 0.003     | 0.01      | ND                      |
| Lead                 | 0.003     | 0.01      | ND                      |
| Mercury              | 0.0009    | 0.003     | ND                      |

ND - not detected; ULOQ - upper limit of quantitation; LOD - limit of detection; LOQ - limit of quantitation



**Lab Comments:**

*Walter Marsh*

Walter Marsh Lead Research Lab Analyst

2026-05-28

Date



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## Amended Report For: Analytical Report 2605120002-V1 - CDPHE Certified Certificate of Analysis



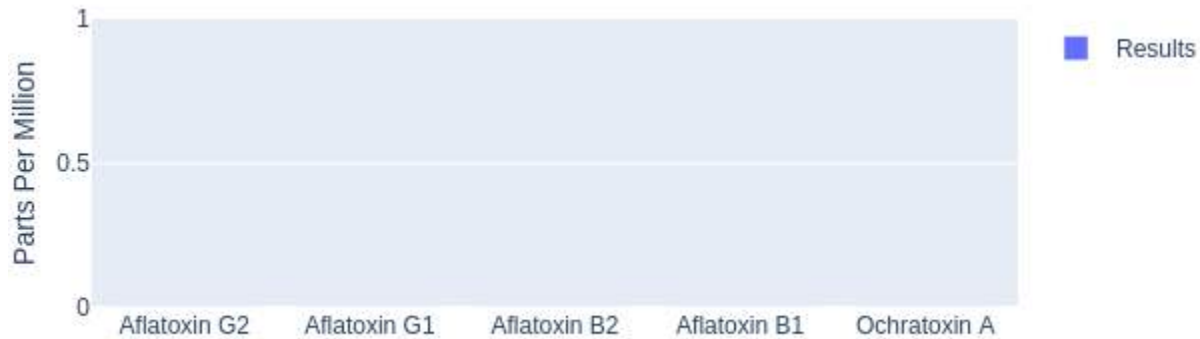
**Manifest:** 2605120002  
**Sample ID:** 1A-GHEMP-2605120002-0001  
**Sample Name:** 344026031632C - FM-942-811  
**Sample Type:** D8 - Distillate  
**Client ID:** CID-00207  
**Client:** FM Labs  
**Address:** 1285 Factory Circle, Fort Lupton, CO 80621

**Test Performed:** Mycotoxin  
**Report No:** A-R-2605120002-V1  
**Receive Date:** 2026-05-12  
**Test Date:** 2026-05-15  
**Report Date:** 2026-05-28  
**Sample Condition:** Good  
**Method Reference:** GH-OP-16

**Scope:** Ochratoxin and Total Aflatoxin were quantified using liquid chromatography coupled to multiple mass spectrometry (LC-MS/MS) equipped with electrospray ionization (ESI) in positive mode after sample extraction. Identification was based on the retention time of each compound and the product mass generated using single reaction monitoring (SRM). Quantitation was determined using external calibration.

| Mycotoxins   | LOD (ppm) | LOQ (ppm) | Reporting Limits (ppm) | Parts Per Million (ppm) |
|--------------|-----------|-----------|------------------------|-------------------------|
| Aflatoxin G2 | 0.0019    | 0.0050    | 0.0050                 | ND                      |
| Aflatoxin G1 | 0.0011    | 0.0050    | 0.0050                 | ND                      |
| Aflatoxin B2 | 0.0017    | 0.0050    | 0.0050                 | ND                      |
| Aflatoxin B1 | 0.0015    | 0.0050    | 0.0050                 | ND                      |
| Ochratoxin A | 0.0033    | 0.0050    | 0.0050                 | ND                      |

ND - not detected; ULOQ - upper limit of quantitation; LOD - limit of detection; LOQ - limit of quantitation



Lab Comments:

*Walter Marsh*

Walter Marsh Lead Research Lab Analyst

2026-05-28

Date



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# Gobi Hemp

## Amended Report For: Microbial Contaminant Report

### 2605120002-V1 - Certificate of Analysis



**Manifest:** 2605120002  
**Sample ID:** 1A-GHEMP-2605120002-0001  
**Sample Name:** 344026031632C - FM-942-811  
**Sample Type:** D8 - Distillate  
**Client ID:** CID-00207  
**Client:** FM Labs  
**Address:** 1285 Factory Circle, Fort Lupton, CO 80621

**Test Performed:** Microbial  
**Report No:** A-M-2605120002-V1  
**Receive Date:** 2026-05-12  
**Test Date:** 2026-05-12  
**Report Date:** 2026-05-28  
**Sample Condition:** Good  
**Method Reference:** MBH-OP-02, MBH-OP-03, MBH-OP-05, MBH-OP-10, MBH-OP-11

**Scope:** Contaminant testing for the identified pathogens *Salmonella spp.* and *Shiga Toxin Virulence Genes, O26, O45, O103, O111, O121, O145 and O157:H7 serogroups of Escherichia coli* (STEC) was performed through Polymerase Chain Reaction (PCR) presumptive experimentation, and confirmed through cultural methodology where applicable. Results for *Salmonella spp.* and STEC are represented as a negative or positive determination, a negative result indicating no detection of the respective contaminant.

Total Yeast and Mold Count (TYMC)/Total Aerobic Count (TAC)/Total Coliform Count (TCC) were determined through 3M™ Petrifilm™ plating technology. The TYMC/TAC/TCC is represented as a count in colony forming units per gram (cfu/g).

| Microbial Contaminants | Results    |
|------------------------|------------|
| <i>Salmonella spp.</i> | ND         |
| STEC                   | ND         |
| Total Yeast and Mold   | <100 CFU/g |
| Total Aerobic          | <100 CFU/g |
| Total Coliform         | <100 CFU/g |

STEC - shiga toxin-producing *Escherichia coli*; TYMC - total yeast and mold count;  
TAC - Total Aerobic Count; TCC - Total Coliform Count; NT - Not Tested;

**Lab Comments:**

Walter Marsh Lead Research Lab Analyst

2026-05-28

Date



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