**PharmLabs San Diego** Certificate of Analysis

Sample **THC Buzzed Gummies - Delta 8 Boosted**

| Delta9 THC | THCa | Total THC (THCa * 0.877 + THC) | Delta8 THC |
|---|---|---|---|
| 0.28% | <LOQ | 0.28% | 0.56% |

| | | | |
|---|---|---|---|
| Sample ID **SD260518-022 (139189)** | | | Matrix **Edible** |
| Tested for **TribeTokes** | | | |
| Received **May 18, 2026** | | Reported **May 28, 2026** | |
| Analyses executed **CANX, D9C** | Unit Mass (g) **75.172** | Num. of Servings **20** | Serving Size (g) **3.76** |

Summary D9C: **The total Δ9-THC content in this sample is 0.28%. For the most accurate Δ9-THC concentration, refer to the GC MS/MS section of this COA.**This sample was tested using HPLC and GC MS/MS. HPLC analysis can yield inconsistent results for **Δ8-THC** and **Δ9-THC** due to isomer interference: GC MS/MS was employed to avoid this issue.

## D9C - D9 Confirmation

Analyzed **May 19, 2026** | Instrument **GC MS/MS** | Method **SOP-041 D9C**
The expanded Uncertainty of the D9 Confirmation analysis is approximately **±7.81%** at the 95% Confidence Level

| Analyte | LOD ppb | LOQ ppb | Result % | Result mg/g | Result mg/Serving | Result mg/Unit |
|---|---|---|---|---|---|---|
| Δ9-Tetrahydrocannabinol (Δ9-THC) | 1.462 | 4.432 | 0.28 | 2.83 | 10.64 | 212.74 |
| **Total Cannabinoids Analyzed** | - | - | 0.28 | 2.83 | 10.64 | 212.74 |

## CANx - Cannabinoids

Analyzed **May 18, 2026** | Instrument **HPLC-VWD** | Method **SOP-001**
The expanded Uncertainty of the Cannabinoids analysis is approximately **±7.81%** at the 95% Confidence Level

| Analyte | LOD mg/g | LOQ mg/g | Result % | Result mg/g | Result mg/Serving | Result mg/Unit |
|---|---|---|---|---|---|---|
| 11-Hydroxy-Δ8-Tetrahydrocannabivarin (11-Hyd-Δ8-THCV) | 0.013 | 0.041 | ND | ND | ND | ND |
| Cannabidiorcin (CBDO) | 0.006 | 0.02 | ND | ND | ND | ND |
| Abnormal Cannabidiorcin (a-CBDO) | 0.013 | 0.038 | ND | ND | ND | ND |
| (+/-)-9B-hydroxy-Hexahydrocannibinol (9b-HHC) | 0.015 | 0.045 | ND | ND | ND | ND |
| 11-Hydroxy-Δ8-Tetrahydrocannabinol (11-Hyd-Δ8-THC) | 0.015 | 0.045 | ND | ND | ND | ND |
| Cannabidiolic Acid (CBDA) | 0.033 | 0.16 | ND | ND | ND | ND |
| Cannabigerol Acid (CBGA) | 0.033 | 0.16 | ND | ND | ND | ND |
| Cannabigerol (CBG) | 0.048 | 0.16 | 0.02 | 0.19 | 0.71 | 14.28 |
| Cannabidiol (CBD) | 0.069 | 0.229 | <LOQ | <LOQ | <LOQ | <LOQ |
| 1(S)-Tetrahydrocannabidiol (1(S)-H4-CBD) | 0.008 | 0.026 | ND | ND | ND | ND |
| 1(R)-Tetrahydrocannabidiol (1(R)-H4-CBD) | 0.016 | 0.049 | ND | ND | ND | ND |
| Tetrahydrocannabivarin (THCV) | 0.049 | 0.162 | ND | ND | ND | ND |
| Δ8-tetrahydrocannabivarin (Δ8-THCV) | 0.012 | 0.036 | ND | ND | ND | ND |
| Cannabidihexol (CBDH) | 0.014 | 0.042 | ND | ND | ND | ND |
| Tetrahydrocannabutol (Δ9-THCB) | 0.01 | 0.029 | <LOQ | <LOQ | <LOQ | <LOQ |
| Cannabinol (CBN) | 0.047 | 0.16 | <LOQ | <LOQ | <LOQ | <LOQ |
| Cannabidiphorol (CBDP) | 0.016 | 0.049 | ND | ND | ND | ND |
| exo-THC (exo-THC) | 0.016 | 0.8 | ND | ND | ND | ND |
| Tetrahydrocannabinol (Δ9-THC) | 0.092 | 0.307 | D9C | D9C | D9C | D9C |
| Δ8-tetrahydrocannabinol (Δ8-THC) | 0.044 | 0.16 | 0.56 | 5.60 | 21.06 | 420.96 |
| (6aR,9S)-Δ10-Tetrahydrocannabinol ((6aR,9S)-Δ10) | 0.015 | 0.8 | ND | ND | ND | ND |
| Hexahydrocannabinol (S Isomer) (9s-HHC) | 0.017 | 0.8 | ND | ND | ND | ND |
| (6aR,9R)-Δ10-Tetrahydrocannabinol ((6aR,9R)-Δ10) | 0.007 | 0.8 | ND | ND | ND | ND |
| Hexahydrocannabinol (R Isomer) (9r-HHC) | 0.016 | 0.8 | ND | ND | ND | ND |
| Tetrahydrocannabinolic Acid (THCA) | 0.117 | 0.389 | <LOQ | <LOQ | <LOQ | <LOQ |
| Δ9-Tetrahydrocannabihexol (Δ9-THCH) | 0.02 | 0.061 | ND | ND | ND | ND |
| Cannabinol Acetate (CBNO) | 0.009 | 0.027 | ND | ND | ND | ND |
| 9(S)-Hexahydrocannabinolic Acid (9(S)-HHCa) | 0.063 | 0.065 | ND | ND | ND | ND |
| 9(R)-Hexahydrocannabinolic Acid (9(R)-HHCa) | 0.191 | 0.196 | ND | ND | ND | ND |
| Δ9-Tetrahydrocannabiphorol (Δ9-THCP) | 0.017 | 0.8 | ND | ND | ND | ND |
| Δ8-Tetrahydrocannabiphorol (Δ8-THCP) | 0.041 | 0.8 | ND | ND | ND | ND |
| Cannabicitran (CBT) | 0.005 | 0.16 | <LOQ | <LOQ | <LOQ | <LOQ |
| Δ8-THC-O-acetate (Δ8-THCO) | 0.076 | 0.8 | ND | ND | ND | ND |
| 9(S)-HHCP (s-HHCP) | 0.013 | 0.041 | ND | ND | ND | ND |
| Δ9-THC-O-acetate (Δ9-THCO) | 0.066 | 0.8 | ND | ND | ND | ND |
| 9(R)-HHCP (r-HHCP) | 0.015 | 0.045 | ND | ND | ND | ND |
| 9(S)-HHC-O-acetate (s-HHCO) | 0.037 | 0.112 | ND | ND | ND | ND |
| 9(R)-HHC-O-acetate (r-HHCO) | 0.031 | 0.093 | ND | ND | ND | ND |
| 3-octyl-Δ8-Tetrahydrocannabinol (Δ8-THC-C8) | 0.021 | 0.062 | ND | ND | ND | ND |
| **Total THC** ( THCa * 0.877 + Δ9THC ) | | | D9C | D9C | D9C | D9C |
| **Total THC + Δ8THC + Δ10THC** ( THCa * 0.877 + Δ9THC + Δ8THC + Δ10THC ) | | | 0.56 | 5.60 | 21.06 | 420.96 |
| **Total CBD** ( CBDa * 0.877 + CBD ) | | | <LOQ | <LOQ | <LOQ | <LOQ |
| **Total CBG** ( CBGa * 0.877 + CBG ) | | | 0.02 | 0.19 | 0.71 | 14.28 |
| **Total HHC** ( 9r-HHC + 9s-HHC ) | | | ND | ND | ND | ND |
| **Total Cannabinoids Analyzed** | | | 0.58 | 5.79 | 21.77 | 435.25 |


**UI** Unidentified
**ND** Not Detected
**N/A** Not Applicable
**NT** Not Reported
**LOD** Limit of Detection
**LOQ** Limit of Quantification
**<LOQ** Detected
**>ULOL** Above upper limit of linearity
**CFU/g** Colony Forming Units per 1 gram
**TNTC** Too Numerous to Count

PJLA Testing Acc. #85368

DEA license: RP0611043
ISO/IEC 17025:2017 Acc. 85368

Scan the QR code to verify authenticity.

Authorized Signature

Brandon Starr

Brandon Starr, Quality Assurance Manager
Thu, 28 May 2026 10:14:37 -0700

Full Panel on Bulk Ingredients

# Gobi Hemp

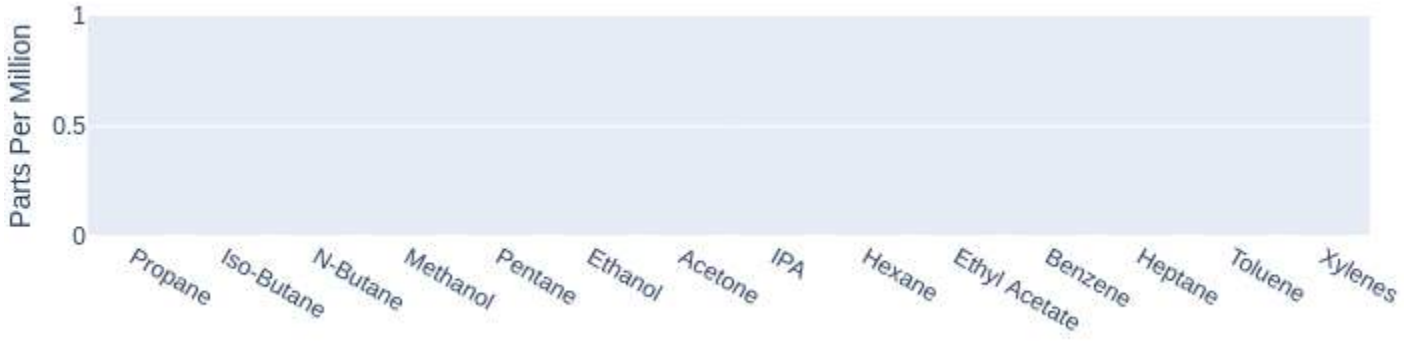
## Amended Report For: Analytical Report 2605120002-V1 - CDPHE Certified Certificate of Analysis

| | | | |
|---|---|---|---|
| **Manifest:** | 2605120002 | **Test Performed:** | RSA |
| **Sample ID:** | 1A-GHEMP-2605120002-0001 | **Report No:** | A-R-2605120002-V1 |
| **Sample Name:** | 344026031632C - FM-942-811 | **Receive Date:** | 2026-05-12 |
| **Sample Type:** | D8 - Distillate | **Test Date:** | 2026-05-14 |
| **Client ID:** | CID-00207 | **Report Date:** | 2026-05-28 |
| **Client:** | FM Labs | **Sample Condition:** | Good |
| **Address:** | 1285 Factory Circle, Fort Lupton, CO 80621 | **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:**

2026-05-28

Walter Marsh Lead Research Lab Analyst — Date



● Gobi Hemp ●
● 3940 Youngfield St. ● Wheat Ridge CO 80033 ● ISO/IEC 17025:2017 Accredited ● (303)456-2040 ●

PJLA Testing Accreditation #103051

# Gobi Hemp

## Amended Report For: Pesticide Residues

## Report 2605120002-V1 - Certificate of Analysis

| | | | |
|---|---|---|---|
| **Manifest:** | 2605120002 | **Test Performed:** | Pesticide |
| **Sample ID:** | 1A-GHEMP-2605120002-0001 | **Report No:** | A-PE-2605120002-V1 |
| **Sample Name:** | 344026031632C - FM-942-811 | **Receive Date:** | 2026-05-12 |
| **Sample Type:** | D8 - Distillate | **Test Date:** | 2026-05-14 |
| **Client ID:** | CID-00207 | **Report Date:** | 2026-05-28 |
| **Client:** | FM Labs | **Sample Condition:** | Good |
| **Facility Address:** | 1285 Factory Circle, Fort Lupton, CO 80621 | **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





PJLA Testing Accreditation #103051

# Gobi Hemp

## Amended Report For: Pesticide Residues

## Report 2605120002-V1

| Pesticide | Limits (ppm) | | Result (ppm) | |
|---|---|---|---|---|
| | Regulatory | Reporting* | | |
| Abamectin | | 0.10000 | ND | LCMS |
| Acephate | | 0.10000 | ND | LCMS |
| Acequinocyl | | 0.10000 | ND | LCMS |
| Acetamiprid | | 0.10000 | ND | LCMS |
| Aldicarb | | 0.10000 | ND | LCMS |
| Allethrin | | 0.10000 | ND | LCMS |
| Atrazine | | 0.10000 | ND | LCMS |
| Azadirachtin | | 0.50000 | ND | LCMS |
| Azoxystrobin | | 0.10000 | ND | LCMS |
| Benzovindiflupyr | | 0.10000 | ND | LCMS |
| Bifenazate | | 0.10000 | ND | LCMS |
| Bifenthrin | | 1.00000 | ND | LCMS |
| Boscalid | | 0.10000 | ND | LCMS |
| Buprofezin | | 0.10000 | ND | LCMS |
| Carbaryl | | 0.10000 | ND | LCMS |
| Carbofuran | | 0.10000 | ND | LCMS |
| Chlorantraniliprole | | 0.10000 | ND | LCMS |
| Chlorphenapyr | | 0.10000 | ND | GCMS |
| Chlorpyrifos | | 0.10000 | ND | LCMS |
| Clofentezine | | 0.10000 | ND | LCMS |
| Clothianidin | | 0.10000 | ND | LCMS |
| Coumaphos | | 0.10000 | ND | LCMS |
| Cyantraniliprole | | 0.10000 | ND | LCMS |
| Cyfluthrin | | 0.20000 | ND | GCMS |
| Cypermethrin | | 0.25000 | ND | GCMS |
| Cyprodinil | | 0.10000 | ND | LCMS |
| Daminozide | | 0.10000 | ND | LCMS |
| Deltamethrin | | 0.50000 | ND | LCMS |
| Diazinon | | 0.10000 | ND | LCMS |
| Dichlorvos | | 0.10000 | ND | GCMS |
| Dimethoate | | 0.10000 | ND | LCMS |
| Dimethomorph | | 0.10000 | ND | LCMS |
| Dinotefuran | | 0.10000 | ND | LCMS |
| Diuron | | 0.10000 | ND | LCMS |

| Pesticide | Limits (ppm) | | Result (ppm) | |
|---|---|---|---|---|
| | Regulatory | Reporting* | | |
| Dodemorph | | 0.10000 | ND | LCMS |
| Endosulfan sulfate | | 0.10000 | ND | GCMS |
| Endosulfan-alpha | | 0.20000 | ND | GCMS |
| Endosulfan-beta | | 0.10000 | ND | GCMS |
| Ethoprophos | | 0.10000 | ND | LCMS |
| Etofenprox | | 0.10000 | ND | LCMS |
| Etoxazole | | 0.10000 | ND | LCMS |
| Etridiazole | | 0.10000 | ND | GCMS |
| Fenhexamid | | 0.12500 | ND | LCMS |
| Fenoxycarb | | 0.10000 | ND | LCMS |
| Fenpyroximate | | 0.10000 | ND | LCMS |
| Fensulfothion | | 0.10000 | ND | LCMS |
| Fenthion | | 0.10000 | ND | GCMS |
| Fenvalerate | | 0.10000 | ND | GCMS |
| Fipronil | | 0.10000 | ND | LCMS |
| Flonicamid | | 0.10000 | ND | LCMS |
| Fludioxonil | | 0.10000 | ND | LCMS |
| Fluopyram | | 0.10000 | ND | LCMS |
| Hexythiazox | | 0.10000 | ND | LCMS |
| Imazalil | | 0.10000 | ND | LCMS |
| Imidacloprid | | 0.10000 | ND | LCMS |
| Iprodione | | 0.50000 | ND | LCMS |
| Kinoprene | | 0.10000 | ND | GCMS |
| Kresoxim-methyl | | 0.10000 | ND | LCMS |
| MGK-264 | | 0.10000 | ND | GCMS |
| Malathion | | 0.10000 | ND | LCMS |
| Metalaxyl | | 0.10000 | ND | LCMS |
| Methiocarb | | 0.10000 | ND | LCMS |
| Methomyl | | 0.10000 | ND | LCMS |
| Methoprene | | 2.00000 | ND | LCMS |
| Mevinphos | | 0.10000 | ND | LCMS |
| Myclobutanil | | 0.10000 | ND | LCMS |
| Naled | | 0.10000 | ND | LCMS |
| Novaluron | | 0.10000 | ND | LCMS |

| Pesticide | Limits (ppm) | | Result (ppm) | |
|---|---|---|---|---|
| | Regulatory | Reporting* | | |
| Oxamyl | | 1.50000 | ND | LCMS |
| Paclobutrazol | | 0.10000 | ND | LCMS |
| Parathion-methyl | | 0.10000 | ND | GCMS |
| Permethrins | | 0.50000 | ND | LCMS |
| Phenothrin | | 0.10000 | ND | LCMS |
| Phosmet | | 0.10000 | ND | LCMS |
| Piperonyl butoxide | | 1.00000 | ND | LCMS |
| Pirimicarb | | 0.10000 | ND | LCMS |
| Prallethrin | | 0.10000 | ND | LCMS |
| Propiconazole | | 0.10000 | ND | LCMS |
| Propoxur | | 0.10000 | ND | LCMS |
| Pyraclostrobin | | 0.10000 | ND | LCMS |
| Pyrethrins | | 0.10000 | ND | LCMS |
| Pyridaben | | 0.10000 | ND | LCMS |
| Pyriproxyfen | | 0.10000 | ND | LCMS |
| Quintozene | | 0.10000 | ND | GCMS |
| Resmethrin | | 0.10000 | ND | LCMS |
| Spinetoram | | 0.10000 | ND | LCMS |
| Spinosad | | 0.10000 | ND | LCMS |
| Spirodiclofen | | 0.25000 | ND | LCMS |
| Spiromesifen | | 3.00000 | ND | LCMS |
| Spirotetramat | | 0.10000 | ND | LCMS |
| Spiroxamine | | 0.10000 | ND | LCMS |
| Tebuconazole | | 0.10000 | ND | LCMS |
| Tebufenozide | | 0.10000 | ND | LCMS |
| Teflubenzuron | | 0.10000 | ND | LCMS |
| Tetrachlorvinphos | | 0.10000 | ND | LCMS |
| Tetramethrin | | 0.10000 | ND | LCMS |
| Thiabendazole | | 0.10000 | ND | LCMS |
| Thiacloprid | | 0.10000 | ND | LCMS |
| Thiamethoxam | | 0.10000 | ND | LCMS |
| Thiophanate-methyl | | 0.10000 | ND | LCMS |
| Trifloxystrobin | | 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.
**

2026-05-28

Walter Marsh Lead Research Lab Analyst — Date



● Gobi Hemp ●
● 3940 Youngfield St. ● Wheat Ridge CO 80033 ● ISO/IEC 17025:2017 Accredited ● (303)456-2040 ●

PJLA Testing Accreditation #103051

# Gobi Hemp

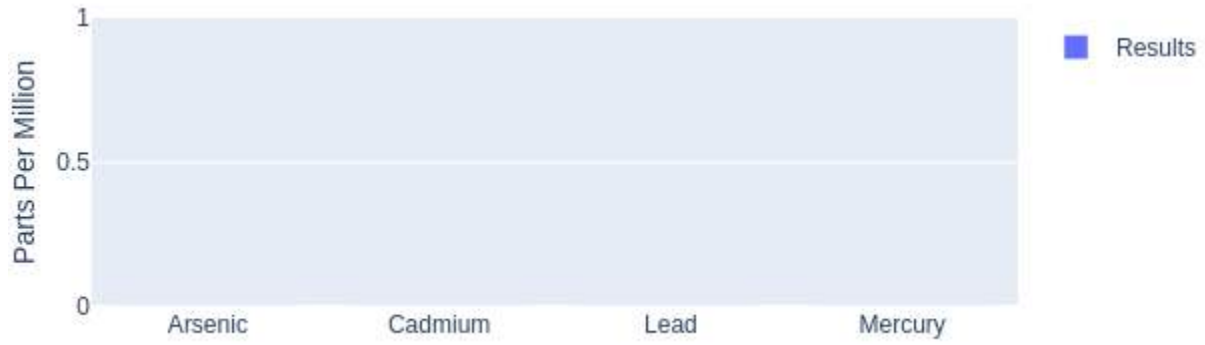
## Amended Report For: Analytical Report 2605120002-V1 - CDPHE Certified Certificate of Analysis

| | | | |
|---|---|---|---|
| **Manifest:** | 2605120002 | **Test Performed:** | Elemental impurities |
| **Sample ID:** | 1A-GHEMP-2605120002-0001 | **Intended Use:** | Inhaled or Audited Product |
| **Sample Name:** | 344026031632C - FM-942-811 | **Report No:** | A-MT-2605120002-V1 |
| **Sample Type:** | D8 - Distillate | **Receive Date:** | 2026-05-12 |
| **Client ID:** | CID-00207 | **Test Date:** | 2026-05-15 |
| **Client:** | FM Labs | **Report Date:** | 2026-05-28 |
| **Address:** | 1285 Factory Circle, Fort Lupton, CO 80621 | **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:**

2026-05-28

Walter Marsh Lead Research Lab Analyst — Date





PJLA Testing Accreditation #103051

# Gobi Hemp

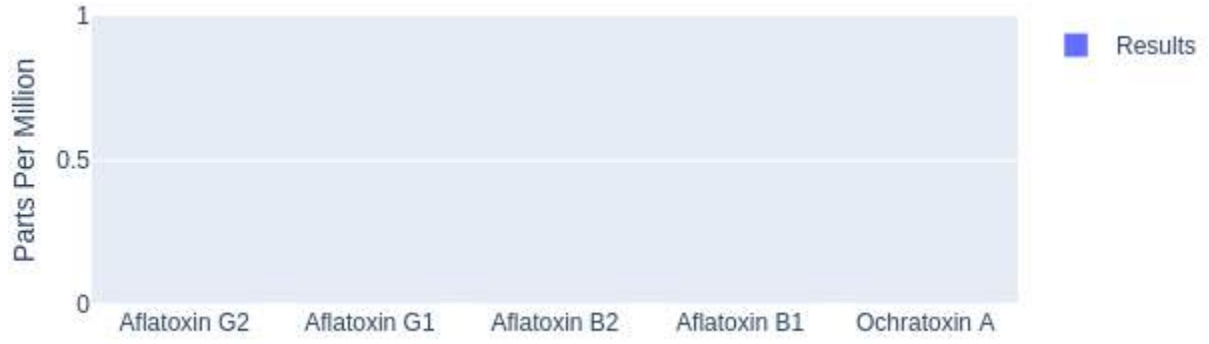
## Amended Report For: Analytical Report 2605120002-V1 - CDPHE Certified Certificate of Analysis

| | | | |
|---|---|---|---|
| **Manifest:** | 2605120002 | **Test Performed:** | Mycotoxin |
| **Sample ID:** | 1A-GHEMP-2605120002-0001 | **Report No:** | A-R-2605120002-V1 |
| **Sample Name:** | 344026031632C - FM-942-811 | **Receive Date:** | 2026-05-12 |
| **Sample Type:** | D8 - Distillate | **Test Date:** | 2026-05-15 |
| **Client ID:** | CID-00207 | **Report Date:** | 2026-05-28 |
| **Client:** | FM Labs | **Sample Condition:** | Good |
| **Address:** | 1285 Factory Circle, Fort Lupton, CO 80621 | **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 Lead Research Lab Analyst

2026-05-28
Date





PJLA Testing Accreditation #103051

# Gobi Hemp

## Amended Report For: Microbial Contaminant Report 2605120002-V1 - Certificate of Analysis

| | | | |
|---|---|---|---|
| **Manifest:** | 2605120002 | **Test Performed:** | Microbial |
| **Sample ID:** | 1A-GHEMP-2605120002-0001 | **Report No:** | A-M-2605120002-V1 |
| **Sample Name:** | 344026031632C - FM-942-811 | **Receive Date:** | 2026-05-12 |
| **Sample Type:** | D8 - Distillate | **Test Date:** | 2026-05-12 |
| **Client ID:** | CID-00207 | **Report Date:** | 2026-05-28 |
| **Client:** | FM Labs | **Sample Condition:** | Good |
| **Address:** | 1285 Factory Circle, Fort Lupton, CO 80621 | **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 |
|---|---|
| *Salmonella spp.* | 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





PJLA Testing Accreditation #103051

**KCA Laboratories**
232 North Plaza Drive
Nicholasville, KY 40356

+1-833-KCA-LABS
https://kcalabs.com
KDA Lic.# P_0058

**Certificate of Analysis**



# D9 Distillate - Naturally Derived From Hemp

Sample ID: SA-250627-64260
Batch: 09DST240_032525
Type: In-Process Material
Matrix: Concentrate - Distillate
Unit Mass (g):

Collected: 03/26/2025
Received: 03/27/2025
Completed: 04/10/2025

**Client**
TribeTokes: 55 Madison Avenue
Suite 400, Morristown NJ 07960, USA
team@tribetokes.com,
844-77-TRIBE (87423)

## Summary

| Test | Date Tested | Status |
|---|---|---|
| Cannabinoids | 04/08/2025 | Tested |
| Heavy Metals | 04/10/2025 | Tested |
| Pesticides | 04/07/2025 | Tested |
| Residual Solvents | 04/07/2025 | Tested |

| 88.7 % | 88.7 % | 94.4 % | Not Tested | Not Tested | Yes |
|---|---|---|---|---|---|
| Total Δ9-THC | Δ9-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 | 0.213 | 2.13 |
| CBCA | 0.0181 | 0.0543 | ND | ND |
| CBCV | 0.006 | 0.018 | ND | ND |
| CBD | 0.0081 | 0.0242 | 0.303 | 3.04 |
| CBDA | 0.0043 | 0.013 | ND | ND |
| CBDV | 0.0061 | 0.0182 | ND | ND |
| CBDVA | 0.0021 | 0.0063 | ND | ND |
| CBG | 0.0057 | 0.0172 | 2.65 | 26.5 |
| CBGA | 0.0049 | 0.0147 | ND | ND |
| CBL | 0.0112 | 0.0335 | 0.137 | 1.37 |
| CBLA | 0.0124 | 0.0371 | ND | ND |
| CBN | 0.0056 | 0.0169 | 0.724 | 7.24 |
| CBNA | 0.006 | 0.0181 | ND | ND |
| CBT | 0.018 | 0.054 | 1.24 | 12.4 |
| Δ4,8-iso-THC | 0.0067 | 0.02 | ND | ND |
| Δ8-iso-THC | 0.0067 | 0.02 | ND | ND |
| Δ8-THC | 0.0104 | 0.0312 | ND | ND |
| Δ8-THCV | 0.0067 | 0.02 | ND | ND |
| Δ9-THC | 0.0076 | 0.0227 | 88.7 | 887 |
| Δ9-THCA | 0.0084 | 0.0251 | ND | ND |
| Δ9-THCV | 0.0069 | 0.0206 | 0.444 | 4.44 |
| Δ9-THCVA | 0.0062 | 0.0186 | ND | ND |
| exo-THC | 0.0067 | 0.02 | ND | ND |
| **Total Δ9-THC** | | | **88.7** | **887** |
| **Total** | | | **94.4** | **944** |

ND = Not Detected; NT = Not Tested; 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: Ryan Bellone
Commercial Director
Date: 07/02/2025

Tested By: Scott Caudill
Laboratory Manager
Date: 04/08/2025

ISO/IEC 17025:2017 Accredited
Accreditation #108651

**KCA Laboratories**
232 North Plaza Drive
Nicholasville, KY 40356

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https://kcalabs.com
KDA Lic.# P_0058

**Certificate of Analysis**



# D9 Distillate - Naturally Derived From Hemp

Sample ID: SA-250627-64260
Batch: 09DST240_032525
Type: In-Process Material
Matrix: Concentrate - Distillate
Unit Mass (g):

Collected: 03/26/2025
Received: 03/27/2025
Completed: 04/10/2025

**Client**
TribeTokes: 55 Madison Avenue
Suite 400, Morristown NJ 07960, USA
team@tribetokes.com,
844-77-TRIBE (87423)

## Heavy Metals by ICP-MS

| Analyte | LOD (ppm) | LOQ (ppm) | Result (ppm) |
|---|---|---|---|
| Arsenic | 0.002 | 0.02 | ND |
| Cadmium | 0.001 | 0.02 | ND |
| Lead | 0.002 | 0.02 | <LOQ |
| Mercury | 0.012 | 0.05 | ND |

ND = Not Detected; NT = Not Tested; LOD = Limit of Detection; LOQ = Limit of Quantitation; P = Pass; F = Fail; RL = Reporting Limit; Values over action limits may be estimates

Generated By: Ryan Bellone
Commercial Director
Date: 07/02/2025

Tested By: Chris Farman
Scientist
Date: 04/10/2025

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team@tribetokes.com,
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## Pesticides by LC-MS/MS and GC-MS/MS

| Analyte | LOD (ppb) | LOQ (ppb) | Result (ppb) | Analyte | LOD (ppb) | LOQ (ppb) | Result (ppb) |
|---|---|---|---|---|---|---|---|
| Abamectin | 30 | 100 | ND | Hexythiazox | 30 | 100 | ND |
| Acephate | 30 | 100 | ND | Imazalil | 30 | 100 | ND |
| Acetamiprid | 30 | 100 | ND | Imidacloprid | 30 | 100 | ND |
| Aldicarb | 30 | 100 | ND | Kresoxim methyl | 30 | 100 | ND |
| Azoxystrobin | 30 | 100 | ND | Malathion | 30 | 100 | ND |
| Bifenazate | 30 | 100 | <LOQ | Metalaxyl | 30 | 100 | ND |
| Bifenthrin | 30 | 100 | ND | Methiocarb | 30 | 100 | ND |
| Boscalid | 30 | 100 | ND | Methomyl | 30 | 100 | ND |
| Carbaryl | 30 | 100 | ND | Mevinphos | 30 | 100 | ND |
| Carbofuran | 30 | 100 | ND | Myclobutanil | 30 | 100 | ND |
| Chloranthraniliprole | 30 | 100 | ND | Naled | 30 | 100 | ND |
| Chlorfenapyr | 30 | 100 | ND | Oxamyl | 30 | 100 | ND |
| Chlorpyrifos | 30 | 100 | ND | Paclobutrazol | 30 | 100 | ND |
| Clofentezine | 30 | 100 | ND | Permethrin | 30 | 100 | ND |
| Coumaphos | 30 | 100 | ND | Phosmet | 30 | 100 | ND |
| Cypermethrin | 30 | 100 | ND | Piperonyl Butoxide | 30 | 100 | ND |
| Daminozide | 30 | 100 | ND | Propiconazole | 30 | 100 | ND |
| Diazinon | 30 | 100 | ND | Propoxur | 30 | 100 | ND |
| Dichlorvos | 30 | 100 | ND | Pyrethrins | 30 | 100 | ND |
| Dimethoate | 30 | 100 | ND | Pyridaben | 30 | 100 | ND |
| Dimethomorph | 30 | 100 | ND | Spinetoram | 30 | 100 | ND |
| Ethoprophos | 30 | 100 | ND | Spinosad | 30 | 100 | ND |
| Etofenprox | 30 | 100 | ND | Spiromesifen | 30 | 100 | ND |
| Etoxazole | 30 | 100 | ND | Spirotetramat | 30 | 100 | ND |
| Fenhexamid | 30 | 100 | ND | Spiroxamine | 30 | 100 | ND |
| Fenoxycarb | 30 | 100 | ND | Tebuconazole | 30 | 100 | ND |
| Fenpyroximate | 30 | 100 | ND | Thiacloprid | 30 | 100 | ND |
| Fipronil | 30 | 100 | ND | Thiamethoxam | 30 | 100 | ND |
| Flonicamid | 30 | 100 | ND | Trifloxystrobin | 30 | 100 | ND |
| Fludioxonil | 30 | 100 | ND | | | | |

ND = Not Detected; NT = Not Tested; LOD = Limit of Detection; LOQ = Limit of Quantitation; P = Pass; F = Fail; RL = Reporting Limit; Values over action limits may be estimates

Generated By: Ryan Bellone
Commercial Director
Date: 07/02/2025

Tested By: Anthony Mattingly
Scientist
Date: 04/07/2025

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**Certificate of Analysis**



# D9 Distillate - Naturally Derived From Hemp

Sample ID: SA-250627-64260
Batch: 09DST240_032525
Type: In-Process Material
Matrix: Concentrate - Distillate
Unit Mass (g):

Collected: 03/26/2025
Received: 03/27/2025
Completed: 04/10/2025

**Client**
TribeTokes: 55 Madison Avenue
Suite 400, Morristown NJ 07960, USA
team@tribetokes.com,
844-77-TRIBE (87423)

## Residual Solvents by HS-GC-MS

| Analyte | LOD (ppm) | LOQ (ppm) | Result (ppm) | Analyte | LOD (ppm) | LOQ (ppm) | Result (ppm) |
|---|---|---|---|---|---|---|---|
| Acetone | 167 | 500 | ND | Ethylene Oxide | 0.5 | 1 | ND |
| Acetonitrile | 14 | 41 | ND | Heptane | 167 | 500 | ND |
| Benzene | 0.5 | 1 | ND | n-Hexane | 10 | 29 | ND |
| Butane | 167 | 500 | ND | Isobutane | 167 | 500 | ND |
| 1-Butanol | 167 | 500 | ND | Isopropyl Acetate | 167 | 500 | ND |
| 2-Butanol | 167 | 500 | ND | Isopropyl Alcohol | 167 | 500 | ND |
| 2-Butanone | 167 | 500 | ND | Isopropylbenzene | 167 | 500 | ND |
| Chloroform | 2 | 6 | ND | Methanol | 100 | 300 | ND |
| Cyclohexane | 129 | 388 | ND | 2-Methylbutane | 10 | 29 | ND |
| 1,2-Dichloroethane | 0.5 | 1 | ND | Methylene Chloride | 20 | 60 | ND |
| 1,2-Dimethoxyethane | 4 | 10 | ND | 2-Methylpentane | 10 | 29 | ND |
| Dimethyl Sulfoxide | 167 | 500 | ND | 3-Methylpentane | 10 | 29 | ND |
| N,N-Dimethylacetamide | 37 | 109 | ND | n-Pentane | 167 | 500 | ND |
| 2,2-Dimethylbutane | 10 | 29 | ND | 1-Pentanol | 167 | 500 | ND |
| 2,3-Dimethylbutane | 10 | 29 | ND | n-Propane | 167 | 500 | ND |
| N,N-Dimethylformamide | 30 | 88 | ND | 1-Propanol | 167 | 500 | ND |
| 2,2-Dimethylpropane | 167 | 500 | ND | Pyridine | 7 | 20 | ND |
| 1,4-Dioxane | 13 | 38 | ND | Tetrahydrofuran | 24 | 72 | ND |
| Ethanol | 167 | 500 | ND | Toluene | 30 | 89 | ND |
| 2-Ethoxyethanol | 6 | 16 | ND | Trichloroethylene | 3 | 8 | ND |
| Ethyl Acetate | 167 | 500 | ND | Xylenes (o-, m-, and p-) | 73 | 217 | ND |
| Ethyl Ether | 167 | 500 | ND | | | | |
| Ethylbenzene | 3 | 7 | ND | | | | |

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Generated By: Ryan Bellone
Commercial Director
Date: 07/02/2025

Tested By: Kelsey Rogers
Scientist
Date: 04/07/2025