

**Manifest:** 2212090001  
**Sample Id:** 1A-GHEMP-2212090001-0001  
**Sample Name:** Urb: Strawberry Sweet Lozenges  
**Sample Type:** Infused (edible)  
**Client Id:** CID-50374  
**Client:** Lifted Made  
**Address:** 5511 95th Ave, , Kenosha, WI 53144

**Test Performed:** Hemp Lab  
**Intended Use:** Oral Consumption or Audited Product  
**Report No:** MT-2212090001-V1  
**Receive Date:** 2022-12-09  
**Test Date:** 2022-12-12  
**Report Date:** 2022-12-13  
**Sample Condition:** Good  
**Method Reference:** GH-OP-17

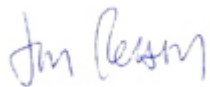
### Scope

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

| Metals  | LOD (ppm) | LOQ (ppm) | Parts Per Million (ppm) |
|---------|-----------|-----------|-------------------------|
| Arsenic | 0.007     | 0.025     | ND                      |
| Cadmium | 0.003     | 0.010     | ND                      |
| Lead    | 0.003     | 0.010     | ND                      |
| Mercury | 0.0009    | 0.003     | ND                      |

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

Laboratory Comments:



Jon Person Client Relations Manager

2022-12-13

Date

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• Wheat Ridge CO 80033 •  
• ISO/IEC 17025:2017 Accredited •  
• (303) 955-4934 •



# Gobi Hemp

## Analytical Report - Certificate of Analysis



**Manifest:** 2212090001  
**Sample Id:** 1A-GHEMP-2212090001-0002  
**Sample Name:** Urb: Watermelon Sweet Lozenges  
**Sample Type:** Infused (edible)  
**Client Id:** CID-50374  
**Client:** Lifted Made  
**Address:** 5511 95th Ave, , Kenosha, WI 53144

**Test Performed:** Hemp Lab  
**Intended Use:** Oral Consumption or Audited Product  
**Report No:** MT-2212090001-V1  
**Receive Date:** 2022-12-09  
**Test Date:** 2022-12-12  
**Report Date:** 2022-12-13  
**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.

| Metals  | LOD (ppm) | LOQ (ppm) | Parts Per Million (ppm) |
|---------|-----------|-----------|-------------------------|
| Arsenic | 0.007     | 0.025     | ND                      |
| Cadmium | 0.003     | 0.010     | ND                      |
| Lead    | 0.003     | 0.010     | ND                      |
| Mercury | 0.0009    | 0.003     | ND                      |

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

Laboratory Comments:

2022-12-13

Jon Person Client Relations Manager

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**Manifest:** 2212090001  
**Sample Id:** 1A-GHEMP-2212090001-0003  
**Sample Name:** Urb: Kiwi Sweet Lozenges  
**Sample Type:** Infused (edible)  
**Client Id:** CID-50374  
**Client:** Lifted Made  
**Address:** 5511 95th Ave, , Kenosha, WI 53144

**Test Performed:** Hemp Lab  
**Intended Use:** Oral Consumption or Audited Product  
**Report No:** MT-2212090001-V1  
**Receive Date:** 2022-12-09  
**Test Date:** 2022-12-12  
**Report Date:** 2022-12-13  
**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.

| Metals  | LOD (ppm) | LOQ (ppm) | Parts Per Million (ppm) |
|---------|-----------|-----------|-------------------------|
| Arsenic | 0.007     | 0.025     | ND                      |
| Cadmium | 0.003     | 0.010     | ND                      |
| Lead    | 0.003     | 0.010     | ND                      |
| Mercury | 0.0009    | 0.003     | ND                      |

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

Laboratory Comments:



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2022-12-13

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# Gobi Hemp

## Analytical Report - Certificate of Analysis



**Manifest:** 2212090001  
**Sample Id:** 1A-GHEMP-2212090001-0004  
**Sample Name:** Urb: Grape Sweet Lozenges  
**Sample Type:** Infused (edible)  
**Client Id:** CID-50374  
**Client:** Lifted Made  
**Address:** 5511 95th Ave, , Kenosha, WI 53144

**Test Performed:** Hemp Lab  
**Intended Use:** Oral Consumption or Audited Product  
**Report No:** MT-2212090001-V1  
**Receive Date:** 2022-12-09  
**Test Date:** 2022-12-12  
**Report Date:** 2022-12-13  
**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.

| Metals  | LOD (ppm) | LOQ (ppm) | Parts Per Million (ppm) |
|---------|-----------|-----------|-------------------------|
| Arsenic | 0.007     | 0.025     | ND                      |
| Cadmium | 0.003     | 0.010     | ND                      |
| Lead    | 0.003     | 0.010     | ND                      |
| Mercury | 0.0009    | 0.003     | ND                      |

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

Laboratory Comments:

Jon Person Client Relations Manager

2022-12-13

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# Gobi Hemp

## Analytical Report - Certificate of Analysis



**Manifest:** 2212090001  
**Sample Id:** 1A-GHEMP-2212090001-0001  
**Sample Name:** Urb: Strawberry Sweet Lozenges  
**Sample Type:** Infused (edible)  
**Client Id:** CID-50374  
**Client:** Lifted Made  
**Address:** 5511 95th Ave, , Kenosha, WI 53144

**Test Performed:** Hemp Lab  
**Report No:** R-2212090001-V1  
**Receive Date:** 2022-12-09  
**Test Date:** 2022-12-09  
**Report Date:** 2022-12-12  
**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; T - trace; ULOQ - upper limit of quantitation; LOD - limit of detection; LOQ - limit of quantitation

Laboratory Comments:

2022-12-12

Jon Person Client Relations Manager

Date

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# Gobi Hemp

## Analytical Report - Certificate of Analysis



**Manifest:** 2212090001  
**Sample Id:** 1A-GHEMP-2212090001-0002  
**Sample Name:** Urb: Watermelon Sweet Lozenges  
**Sample Type:** Infused (edible)  
**Client Id:** CID-50374  
**Client:** Lifted Made  
**Address:** 5511 95th Ave, , Kenosha, WI 53144

**Test Performed:** Hemp Lab  
**Report No:** R-2212090001-V1  
**Receive Date:** 2022-12-09  
**Test Date:** 2022-12-09  
**Report Date:** 2022-12-12  
**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; T - trace; ULOQ - upper limit of quantitation; LOD - limit of detection; LOQ - limit of quantitation

Laboratory Comments:

2022-12-12

Jon Person Client Relations Manager

Date

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**Manifest:** 2212090001  
**Sample Id:** 1A-GHEMP-2212090001-0003  
**Sample Name:** Urb: Kiwi Sweet Lozenges  
**Sample Type:** Infused (edible)  
**Client Id:** CID-50374  
**Client:** Lifted Made  
**Address:** 5511 95th Ave, , Kenosha, WI 53144

**Test Performed:** Hemp Lab  
**Report No:** R-2212090001-V1  
**Receive Date:** 2022-12-09  
**Test Date:** 2022-12-09  
**Report Date:** 2022-12-12  
**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; T - trace; ULOQ - upper limit of quantitation; LOD - limit of detection; LOQ - limit of quantitation

Laboratory Comments:

2022-12-12

Jon Person Client Relations Manager

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• Wheat Ridge CO 80033 •  
• (303) 955-4934 •





**Manifest:** 2212090001  
**Sample Id:** 1A-GHEMP-2212090001-0004  
**Sample Name:** Urb: Grape Sweet Lozenges  
**Sample Type:** Infused (edible)  
**Client Id:** CID-50374  
**Client:** Lifted Made  
**Address:** 5511 95th Ave, , Kenosha, WI 53144

**Test Performed:** Hemp Lab  
**Report No:** R-2212090001-V1  
**Receive Date:** 2022-12-09  
**Test Date:** 2022-12-09  
**Report Date:** 2022-12-12  
**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; T - trace; ULOQ - upper limit of quantitation; LOD - limit of detection; LOQ - limit of quantitation

Laboratory Comments:

2022-12-12

Jon Person Client Relations Manager

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• Wheat Ridge CO 80033 •  
• (303) 955-4934 •





# Gobi Hemp

## Pesticide Residues Report - Certificate of Analysis



**Manifest:** 2212090001  
**Sample Id:** 1A-GHEMP-2212090001-0001  
**Sample Name:** Urb: Strawberry Sweet Lozenges  
**Sample Type:** Infused (edible)  
**Client Id:** CID-50374  
**Client:** Lifted Made  
**Address:** 5511 95th Ave, , Kenosha, WI 53144

**Test Performed:** Hemp Lab  
**Report No:** PE-2212090001-V1  
**Receive Date:** 2022-12-09  
**Test Date:** 2022-12-12  
**Report Date:** 2022-12-13  
**Sample Condition:** Good  
**Method Reference:** GH-OP-11

### Scope

The content of 60 pesticides were quantified using liquid chromatography coupled to multiple mass spectrometry (LC-MS2) equipped with electrospray ionization (ESI) in positive mode after sample extraction using methodology based on AOAC 2007 and EN 15662 standard procedures. Identification was based on the retention time of each compound and the product mass generated using single reaction monitoring (SRM), and quantitation was determined using external standard calibration.

| Analyte                 | Reporting Level µg/g | µg/g |
|-------------------------|----------------------|------|
| Avermectin B1a          | 0.1                  | ND   |
| Acephate                | 0.1                  | ND   |
| Acetamiprid             | 0.1                  | ND   |
| Aldicarb                | 0.1                  | ND   |
| Azoxystrobin            | 0.1                  | ND   |
| Bifenazate              | 0.1                  | ND   |
| Bifenthrin              | 0.1                  | ND   |
| Boscalid                | 0.1                  | ND   |
| Captan                  | 0.1                  | ND   |
| Carbaryl                | 0.1                  | ND   |
| Carbofuran              | 0.1                  | ND   |
| Chlorantraniliprole     | 0.1                  | ND   |
| Chlordane               | 0.1                  | ND   |
| Chlorpyrifos            | 0.1                  | ND   |
| Clofentazine            | 0.1                  | ND   |
| Coumaphos               | 0.1                  | ND   |
| Baythroid (Cyfluthrin)* | 0.1                  | NT   |
| Cypermethrin*           | 0.1                  | NT   |
| Dichlorvos              | 0.1                  | ND   |
| Diazinon                | 0.1                  | ND   |
| Dimethoate              | 0.1                  | ND   |
| Dimethomorph*           | 0.1                  | ND   |
| Prophos                 | 0.1                  | ND   |
| Etofenprox              | 0.1                  | ND   |
| Etoxazole               | 0.1                  | ND   |
| Fenhexamid              | 0.1                  | ND   |
| Fenoxycarb              | 0.1                  | ND   |
| Fenpyroximate           | 0.1                  | ND   |
| Fipronil                | 0.1                  | ND   |
| Flonicamid              | 0.1                  | ND   |
| Fludioxonil             | 0.1                  | ND   |

| Analyte                 | Reporting Level µg/g | µg/g |
|-------------------------|----------------------|------|
| Hexythiazox             | 0.1                  | ND   |
| Imazilil                | 0.1                  | ND   |
| Imidacloprid            | 0.1                  | ND   |
| Kresoxim Methyl         | 0.1                  | ND   |
| Malathion               | 0.1                  | ND   |
| Metalaxyl               | 0.1                  | ND   |
| Methiocarb              | 0.1                  | ND   |
| Methomyl                | 0.1                  | ND   |
| Mevinphos*              | 0.1                  | ND   |
| MGK-264                 | 0.1                  | NT   |
| Myclobutanil            | 0.1                  | ND   |
| Oxamyl                  | 0.1                  | ND   |
| Paclobutrazol           | 0.1                  | ND   |
| Pentachloronitrobenzene | 0.1                  | ND   |
| Permethrin*             | 0.1                  | ND   |
| Imidan(Phosmet)         | 0.1                  | ND   |
| Piperonyl Butoxide      | 0.1                  | ND   |
| Propiconazole           | 0.1                  | ND   |
| Propuxor                | 0.1                  | ND   |
| Pyrethrin*              | 0.1                  | ND   |
| Pyridaben               | 0.1                  | ND   |
| Spinetoram              | 0.1                  | ND   |
| Spinosad*               | 0.1                  | ND   |
| Spiromefesin            | 0.1                  | ND   |
| Spirotetramat           | 0.1                  | ND   |
| Spiroxamine             | 0.1                  | ND   |
| Tebuconazole            | 0.1                  | ND   |
| Thiacloprid             | 0.1                  | ND   |
| Thiamethoxam            | 0.1                  | ND   |
| Trifloxystrobin         | 0.1                  | ND   |

NT - not tested; ND - not detected above Reporting Level; T - trace; \* Total of Isomers

### Lab Comments:

Jon Person Client Relations Manager

2022-12-13

Date

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- (720)560-9299 •



# Gobi Hemp

## Pesticide Residues Report - Certificate of Analysis



**Manifest:** 2212090001  
**Sample Id:** 1A-GHEMP-2212090001-0002  
**Sample Name:** Urb: Watermelon Sweet Lozenges  
**Sample Type:** Infused (edible)  
**Client Id:** CID-50374  
**Client:** Lifted Made  
**Address:** 5511 95th Ave, , Kenosha, WI 53144

**Test Performed:** Hemp Lab  
**Report No:** PE-2212090001-V1  
**Receive Date:** 2022-12-09  
**Test Date:** 2022-12-12  
**Report Date:** 2022-12-13  
**Sample Condition:** Good  
**Method Reference:** GH-OP-11

### Scope

The content of 60 pesticides were quantified using liquid chromatography coupled to multiple mass spectrometry (LC-MS2) equipped with electrospray ionization (ESI) in positive mode after sample extraction using methodology based on AOAC 2007 and EN 15662 standard procedures. Identification was based on the retention time of each compound and the product mass generated using single reaction monitoring (SRM), and quantitation was determined using external standard calibration.

| Analyte                 | Reporting Level µg/g | µg/g |
|-------------------------|----------------------|------|
| Avermectin B1a          | 0.1                  | ND   |
| Acephate                | 0.1                  | ND   |
| Acetamiprid             | 0.1                  | ND   |
| Aldicarb                | 0.1                  | ND   |
| Azoxystrobin            | 0.1                  | ND   |
| Bifenazate              | 0.1                  | ND   |
| Bifenthrin              | 0.1                  | ND   |
| Boscalid                | 0.1                  | ND   |
| Captan                  | 0.1                  | ND   |
| Carbaryl                | 0.1                  | ND   |
| Carbofuran              | 0.1                  | ND   |
| Chlorantraniliprole     | 0.1                  | ND   |
| Chlordane               | 0.1                  | ND   |
| Chlorpyrifos            | 0.1                  | ND   |
| Clofentazine            | 0.1                  | ND   |
| Coumaphos               | 0.1                  | ND   |
| Baythroid (Cyfluthrin)* | 0.1                  | NT   |
| Cypermethrin*           | 0.1                  | NT   |
| Dichlorvos              | 0.1                  | ND   |
| Diazinon                | 0.1                  | ND   |
| Dimethoate              | 0.1                  | ND   |
| Dimethomorph*           | 0.1                  | ND   |
| Prophos                 | 0.1                  | ND   |
| Etofenprox              | 0.1                  | ND   |
| Etoxazole               | 0.1                  | ND   |
| Fenhexamid              | 0.1                  | ND   |
| Fenoxycarb              | 0.1                  | ND   |
| Fenpyroximate           | 0.1                  | ND   |
| Fipronil                | 0.1                  | ND   |
| Flonicamid              | 0.1                  | ND   |
| Fludioxonil             | 0.1                  | ND   |

| Analyte                 | Reporting Level µg/g | µg/g |
|-------------------------|----------------------|------|
| Hexythiazox             | 0.1                  | ND   |
| Imazilil                | 0.1                  | ND   |
| Imidacloprid            | 0.1                  | ND   |
| Kresoxim Methyl         | 0.1                  | ND   |
| Malathion               | 0.1                  | ND   |
| Metalaxyl               | 0.1                  | ND   |
| Methiocarb              | 0.1                  | ND   |
| Methomyl                | 0.1                  | ND   |
| Mevinphos*              | 0.1                  | ND   |
| MGK-264                 | 0.1                  | NT   |
| Myclobutanil            | 0.1                  | ND   |
| Oxamyl                  | 0.1                  | ND   |
| Paclobutrazol           | 0.1                  | ND   |
| Pentachloronitrobenzene | 0.1                  | ND   |
| Permethrin*             | 0.1                  | ND   |
| Imidan(Phosmet)         | 0.1                  | ND   |
| Piperonyl Butoxide      | 0.1                  | ND   |
| Propiconazole           | 0.1                  | ND   |
| Propuxor                | 0.1                  | ND   |
| Pyrethrin*              | 0.1                  | ND   |
| Pyridaben               | 0.1                  | ND   |
| Spinetoram              | 0.1                  | ND   |
| Spinosad*               | 0.1                  | ND   |
| Spiromefesin            | 0.1                  | ND   |
| Spirotetramat           | 0.1                  | ND   |
| Spiroxamine             | 0.1                  | ND   |
| Tebuconazole            | 0.1                  | ND   |
| Thiacloprid             | 0.1                  | ND   |
| Thiamethoxam            | 0.1                  | ND   |
| Trifloxystrobin         | 0.1                  | ND   |

NT - not tested; ND - not detected above Reporting Level; T - trace; \* Total of Isomers

### Lab Comments:

Jon Person Client Relations Manager

2022-12-13

Date

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- Gobi Hemp •
- 3940 Youngfield St. Wheat Ridge CO 80033 •
- (720)560-9299 •



# Gobi Hemp

## Pesticide Residues Report - Certificate of Analysis



**Manifest:** 2212090001  
**Sample Id:** 1A-GHEMP-2212090001-0003  
**Sample Name:** Urb: Kiwi Sweet Lozenges  
**Sample Type:** Infused (edible)  
**Client Id:** CID-50374  
**Client:** Lifted Made  
**Address:** 5511 95th Ave, , Kenosha, WI 53144

**Test Performed:** Hemp Lab  
**Report No:** PE-2212090001-V1  
**Receive Date:** 2022-12-09  
**Test Date:** 2022-12-12  
**Report Date:** 2022-12-13  
**Sample Condition:** Good  
**Method Reference:** GH-OP-11

### Scope

The content of 60 pesticides were quantified using liquid chromatography coupled to multiple mass spectrometry (LC-MS2) equipped with electrospray ionization (ESI) in positive mode after sample extraction using methodology based on AOAC 2007 and EN 15662 standard procedures. Identification was based on the retention time of each compound and the product mass generated using single reaction monitoring (SRM), and quantitation was determined using external standard calibration.

| Analyte                 | Reporting Level µg/g | µg/g |
|-------------------------|----------------------|------|
| Avermectin B1a          | 0.1                  | ND   |
| Acephate                | 0.1                  | ND   |
| Acetamiprid             | 0.1                  | ND   |
| Aldicarb                | 0.1                  | ND   |
| Azoxystrobin            | 0.1                  | ND   |
| Bifenazate              | 0.1                  | ND   |
| Bifenthrin              | 0.1                  | ND   |
| Boscalid                | 0.1                  | ND   |
| Captan                  | 0.1                  | ND   |
| Carbaryl                | 0.1                  | ND   |
| Carbofuran              | 0.1                  | ND   |
| Chlorantraniliprole     | 0.1                  | ND   |
| Chlordane               | 0.1                  | ND   |
| Chlorpyrifos            | 0.1                  | ND   |
| Clofentazine            | 0.1                  | ND   |
| Coumaphos               | 0.1                  | ND   |
| Baythroid (Cyfluthrin)* | 0.1                  | NT   |
| Cypermethrin*           | 0.1                  | NT   |
| Dichlorvos              | 0.1                  | ND   |
| Diazinon                | 0.1                  | ND   |
| Dimethoate              | 0.1                  | ND   |
| Dimethomorph*           | 0.1                  | ND   |
| Prophos                 | 0.1                  | ND   |
| Etofenprox              | 0.1                  | ND   |
| Etoxazole               | 0.1                  | ND   |
| Fenhexamid              | 0.1                  | ND   |
| Fenoxycarb              | 0.1                  | ND   |
| Fenpyroximate           | 0.1                  | ND   |
| Fipronil                | 0.1                  | ND   |
| Flonicamid              | 0.1                  | ND   |
| Fludioxonil             | 0.1                  | ND   |

| Analyte                 | Reporting Level µg/g | µg/g |
|-------------------------|----------------------|------|
| Hexythiazox             | 0.1                  | ND   |
| Imazilil                | 0.1                  | ND   |
| Imidacloprid            | 0.1                  | ND   |
| Kresoxim Methyl         | 0.1                  | ND   |
| Malathion               | 0.1                  | ND   |
| Metalaxyl               | 0.1                  | ND   |
| Methiocarb              | 0.1                  | ND   |
| Methomyl                | 0.1                  | ND   |
| Mevinphos*              | 0.1                  | ND   |
| MGK-264                 | 0.1                  | NT   |
| Myclobutanil            | 0.1                  | ND   |
| Oxamyl                  | 0.1                  | ND   |
| Paclobutrazol           | 0.1                  | ND   |
| Pentachloronitrobenzene | 0.1                  | ND   |
| Permethrin*             | 0.1                  | ND   |
| Imidan(Phosmet)         | 0.1                  | ND   |
| Piperonyl Butoxide      | 0.1                  | ND   |
| Propiconazole           | 0.1                  | ND   |
| Propuxor                | 0.1                  | ND   |
| Pyrethrin*              | 0.1                  | ND   |
| Pyridaben               | 0.1                  | ND   |
| Spinetoram              | 0.1                  | ND   |
| Spinosad*               | 0.1                  | ND   |
| Spiromefesin            | 0.1                  | ND   |
| Spirotetramat           | 0.1                  | ND   |
| Spiroxamine             | 0.1                  | ND   |
| Tebuconazole            | 0.1                  | ND   |
| Thiacloprid             | 0.1                  | ND   |
| Thiamethoxam            | 0.1                  | ND   |
| Trifloxystrobin         | 0.1                  | ND   |

NT - not tested; ND - not detected above Reporting Level; T - trace; \* Total of Isomers

### Lab Comments:

Jon Person Client Relations Manager

2022-12-13

Date

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# Gobi Hemp

## Pesticide Residues Report - Certificate of Analysis



**Manifest:** 2212090001  
**Sample Id:** 1A-GHEMP-2212090001-0004  
**Sample Name:** Urb: Grape Sweet Lozenges  
**Sample Type:** Infused (edible)  
**Client Id:** CID-50374  
**Client:** Lifted Made  
**Address:** 5511 95th Ave, , Kenosha, WI 53144

**Test Performed:** Hemp Lab  
**Report No:** PE-2212090001-V1  
**Receive Date:** 2022-12-09  
**Test Date:** 2022-12-12  
**Report Date:** 2022-12-13  
**Sample Condition:** Good  
**Method Reference:** GH-OP-11

### Scope

The content of 60 pesticides were quantified using liquid chromatography coupled to multiple mass spectrometry (LC-MS2) equipped with electrospray ionization (ESI) in positive mode after sample extraction using methodology based on AOAC 2007 and EN 15662 standard procedures. Identification was based on the retention time of each compound and the product mass generated using single reaction monitoring (SRM), and quantitation was determined using external standard calibration.

| Analyte                 | Reporting Level µg/g | µg/g |
|-------------------------|----------------------|------|
| Avermectin B1a          | 0.1                  | ND   |
| Acephate                | 0.1                  | ND   |
| Acetamiprid             | 0.1                  | ND   |
| Aldicarb                | 0.1                  | ND   |
| Azoxystrobin            | 0.1                  | ND   |
| Bifenazate              | 0.1                  | ND   |
| Bifenthrin              | 0.1                  | ND   |
| Boscalid                | 0.1                  | ND   |
| Captan                  | 0.1                  | ND   |
| Carbaryl                | 0.1                  | ND   |
| Carbofuran              | 0.1                  | ND   |
| Chlorantraniliprole     | 0.1                  | ND   |
| Chlordane               | 0.1                  | ND   |
| Chlorpyrifos            | 0.1                  | ND   |
| Clofentazine            | 0.1                  | ND   |
| Coumaphos               | 0.1                  | ND   |
| Baythroid (Cyfluthrin)* | 0.1                  | NT   |
| Cypermethrin*           | 0.1                  | NT   |
| Dichlorvos              | 0.1                  | ND   |
| Diazinon                | 0.1                  | ND   |
| Dimethoate              | 0.1                  | ND   |
| Dimethomorph*           | 0.1                  | ND   |
| Prophos                 | 0.1                  | ND   |
| Etofenprox              | 0.1                  | ND   |
| Etoxazole               | 0.1                  | ND   |
| Fenhexamid              | 0.1                  | ND   |
| Fenoxycarb              | 0.1                  | ND   |
| Fenpyroximate           | 0.1                  | ND   |
| Fipronil                | 0.1                  | ND   |
| Flonicamid              | 0.1                  | ND   |
| Fludioxonil             | 0.1                  | ND   |

| Analyte                 | Reporting Level µg/g | µg/g |
|-------------------------|----------------------|------|
| Hexythiazox             | 0.1                  | ND   |
| Imazilil                | 0.1                  | ND   |
| Imidacloprid            | 0.1                  | ND   |
| Kresoxim Methyl         | 0.1                  | ND   |
| Malathion               | 0.1                  | ND   |
| Metalaxyl               | 0.1                  | ND   |
| Methiocarb              | 0.1                  | ND   |
| Methomyl                | 0.1                  | ND   |
| Mevinphos*              | 0.1                  | ND   |
| MGK-264                 | 0.1                  | NT   |
| Myclobutanil            | 0.1                  | ND   |
| Oxamyl                  | 0.1                  | ND   |
| Paclobutrazol           | 0.1                  | ND   |
| Pentachloronitrobenzene | 0.1                  | ND   |
| Permethrin*             | 0.1                  | ND   |
| Imidan(Phosmet)         | 0.1                  | ND   |
| Piperonyl Butoxide      | 0.1                  | ND   |
| Propiconazole           | 0.1                  | ND   |
| Propuxor                | 0.1                  | ND   |
| Pyrethrin*              | 0.1                  | ND   |
| Pyridaben               | 0.1                  | ND   |
| Spinetoram              | 0.1                  | ND   |
| Spinosad*               | 0.1                  | ND   |
| Spiromefesin            | 0.1                  | ND   |
| Spirotetramat           | 0.1                  | ND   |
| Spiroxamine             | 0.1                  | ND   |
| Tebuconazole            | 0.1                  | ND   |
| Thiacloprid             | 0.1                  | ND   |
| Thiamethoxam            | 0.1                  | ND   |
| Trifloxystrobin         | 0.1                  | ND   |

NT - not tested; ND - not detected above Reporting Level; T - trace; \* Total of Isomers

### Lab Comments:

Jon Person Client Relations Manager

2022-12-13

Date

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# Gobi Hemp

## Analytical Report - Certificate of Analysis



**Manifest:** 2212090001  
**Sample Id:** 1A-GHEMP-2212090001-0001  
**Sample Name:** Urb: Strawberry Sweet Lozenges  
**Sample Type:** Infused (edible)  
**Client Id:** CID-50374  
**Client:** Lifted Made  
**Address:** 5511 95th Ave, , Kenosha, WI 53144

**Test Performed:** Hemp Lab  
**Report No:** R-2212090001-V1  
**Receive Date:** 2022-12-09  
**Test Date:** 2022-12-12  
**Report Date:** 2022-12-13  
**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       | 135       | 372       | ND                      |
| Iso-Butane    | 82        | 490       | ND                      |
| N-Butane      | 107       | 490       | ND                      |
| Methanol      | 38        | 120       | ND                      |
| Pentane       | 73        | 100       | ND                      |
| Ethanol       | 50        | 200       | ND                      |
| Acetone       | 82        | 200       | ND                      |
| IPA           | 40        | 200       | ND                      |
| Hexane        | 25        | 50        | ND                      |
| Ethyl Acetate | 57        | 200       | ND                      |
| Benzene       | 0.65      | 1         | ND                      |
| Heptane       | 137       | 200       | ND                      |
| Toluene       | 75        | 100       | ND                      |
| Xylenes       | 112       | 200       | ND                      |

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

Laboratory Comments:

2022-12-13

Jon Person Client Relations Manager

Date

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# Gobi Hemp

## Analytical Report - Certificate of Analysis



**Manifest:** 2212090001  
**Sample Id:** 1A-GHEMP-2212090001-0002  
**Sample Name:** Urb: Watermelon Sweet Lozenges  
**Sample Type:** Infused (edible)  
**Client Id:** CID-50374  
**Client:** Lifted Made  
**Address:** 5511 95th Ave, , Kenosha, WI 53144

**Test Performed:** Hemp Lab  
**Report No:** R-2212090001-V1  
**Receive Date:** 2022-12-09  
**Test Date:** 2022-12-12  
**Report Date:** 2022-12-13  
**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       | 135       | 372       | ND                      |
| Iso-Butane    | 82        | 490       | ND                      |
| N-Butane      | 107       | 490       | ND                      |
| Methanol      | 38        | 120       | ND                      |
| Pentane       | 73        | 100       | ND                      |
| Ethanol       | 50        | 200       | ND                      |
| Acetone       | 82        | 200       | ND                      |
| IPA           | 40        | 200       | ND                      |
| Hexane        | 25        | 50        | ND                      |
| Ethyl Acetate | 57        | 200       | ND                      |
| Benzene       | 0.65      | 1         | ND                      |
| Heptane       | 137       | 200       | ND                      |
| Toluene       | 75        | 100       | ND                      |
| Xylenes       | 112       | 200       | ND                      |

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

Laboratory Comments:

Jon Person Client Relations Manager

2022-12-13

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# Gobi Hemp

## Analytical Report - Certificate of Analysis



**Manifest:** 2212090001  
**Sample Id:** 1A-GHEMP-2212090001-0003  
**Sample Name:** Urb: Kiwi Sweet Lozenges  
**Sample Type:** Infused (edible)  
**Client Id:** CID-50374  
**Client:** Lifted Made  
**Address:** 5511 95th Ave, , Kenosha, WI 53144

**Test Performed:** Hemp Lab  
**Report No:** R-2212090001-V1  
**Receive Date:** 2022-12-09  
**Test Date:** 2022-12-12  
**Report Date:** 2022-12-13  
**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       | 135       | 372       | ND                      |
| Iso-Butane    | 82        | 490       | ND                      |
| N-Butane      | 107       | 490       | ND                      |
| Methanol      | 38        | 120       | ND                      |
| Pentane       | 73        | 100       | ND                      |
| Ethanol       | 50        | 200       | ND                      |
| Acetone       | 82        | 200       | ND                      |
| IPA           | 40        | 200       | ND                      |
| Hexane        | 25        | 50        | ND                      |
| Ethyl Acetate | 57        | 200       | ND                      |
| Benzene       | 0.65      | 1         | ND                      |
| Heptane       | 137       | 200       | ND                      |
| Toluene       | 75        | 100       | ND                      |
| Xylenes       | 112       | 200       | ND                      |

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

Laboratory Comments:

2022-12-13

Jon Person Client Relations Manager

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# Gobi Hemp

## Analytical Report - Certificate of Analysis



**Manifest:** 2212090001  
**Sample Id:** 1A-GHEMP-2212090001-0004  
**Sample Name:** Urb: Grape Sweet Lozenges  
**Sample Type:** Infused (edible)  
**Client Id:** CID-50374  
**Client:** Lifted Made  
**Address:** 5511 95th Ave, , Kenosha, WI 53144

**Test Performed:** Hemp Lab  
**Report No:** R-2212090001-V1  
**Receive Date:** 2022-12-09  
**Test Date:** 2022-12-12  
**Report Date:** 2022-12-13  
**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       | 135       | 372       | ND                      |
| Iso-Butane    | 82        | 490       | ND                      |
| N-Butane      | 107       | 490       | ND                      |
| Methanol      | 38        | 120       | ND                      |
| Pentane       | 73        | 100       | ND                      |
| Ethanol       | 50        | 200       | ND                      |
| Acetone       | 82        | 200       | ND                      |
| IPA           | 40        | 200       | ND                      |
| Hexane        | 25        | 50        | ND                      |
| Ethyl Acetate | 57        | 200       | ND                      |
| Benzene       | 0.65      | 1         | ND                      |
| Heptane       | 137       | 200       | ND                      |
| Toluene       | 75        | 100       | ND                      |
| Xylenes       | 112       | 200       | ND                      |

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

Laboratory Comments:

Jon Person Client Relations Manager

2022-12-13

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