

Oklahoma City Location 3680 E I-240 Service Rd Oklahoma City, OK 73135 (405) 595-0344 **Tulsa Location** 14720 E Admiral Pl, Ste H Tulsa, OK 74108 (918) 340-7139

CERTIFICATE OF ANALYSIS

Order Type: Medical Cannabis

Customer ID: 194

Harvest/Extract Lot: None Harvest/Extract Batch:

Order ID: OR2019-4655

Customer Name: Higher Health

HH-19-1-276-288

Cultivar (Strain): Oil HH191276288

Lab ID: SA2019-14844 **Date Received:** 10/17/2019

Sample Matrix: Concentrate

Sample Date: 10/17/2019

Jale Received: 10/1//

Date Completed: 10/24/2019

Remarks:

CANNABINOID (POTENCY) PROFILE

Analysis Date/Time: 10/21/2019 1234

Method: 2019-10-18T20:36:50Z

30

Moisture Content (%): -Water Activity (aw): -

Analyst: OL

TOTAL CBD

Instrument: Agilent 1100

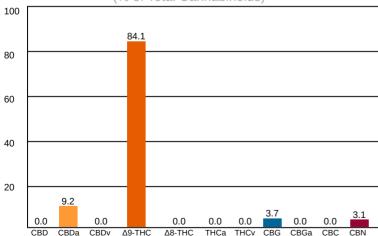
Cannabinoid	Result (%)	Result (mg/g)	Reporting Limit (mg/g)	Result (mg/mL)	Per Unit (mg)	300
CDD	(70)	(IIIS/S)		(IIIg/IIIL)	(IIIg)	
CBD	-	-	0.517	-	-	
CBDa	3.43	34.3	0.517	-	34	
CBDv	-	-	0.517	-	-	H - 14 - 1
Δ9-ΤΗС	31.4	314	0.517	-	314	76 - 94
$\Delta 8$ -THC	-	-	0.517	-	-	10- 400
THCa	-	-	0.517	-	-	
THCv	-	-	0.517	-	-	
CBC	-	-	0.517	-	-	
CBG	1.37	13.7	0.517	-	14	
CBGa	-	-	0.517	-	-	
CBN	1.17	11.7	0.517	-	12	
TOTAL	37.4	374		-	374	UNIT MASS (g): 1
TOTAL THC	31.4	314		-	314	"-" Not detected above RL.

Cannabinoid Distribution

30

3.01

(% of Total Cannabinoids)



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Deviations from standard operating procedure: None

Recoveries for all analyte standards: 90-110% Replicate Uncertainties: <5% RSD, <20% RPD Sample/Reagent Blanks: <RL for all analytes

Values for plant matter are adjusted for moisture content.

Total THC = (THCa x 0.877) + $\Delta 9$ -THC Total CBD = (CBDa x 0.877) + CBD

Percentage results are reported by mass. mg/g results are reported as mass component per mass material.

Abbreviations: UV - Ultraviolet, HPLC - High Pressure Liquid Chromatography, RL - Reporting Limit, RPD -Relative Percent Difference, RSD - Relative Standard Deviation





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Deviations from SOP:

Remarks:

TERPENOID PROFILE

Analysis Date/Time: 10/21/2019 1234

Analyst: OL

Guaiol

Method: HS/GC/FID (Internal Method-002)

Instrument: Agilent 6890

None



Terpene Result Result $(\mu g/g)$ (%) α-Bisabolol 9920 0.992 Camphene

δ-3-Carene **β-**Caryophyllene 7620 0.762

Caryophyllene 8260 oxide

p-Cymene Eucalyptol Geraniol

> 1920 0.192 1480

α-Humulene Isopulegol d-Limonene

Linalool **β-Myrcene** 5120

cis-Nerolidol trans-Nerolidol 5320

α-Ocimene **β-Ocimene** α-Pinene

β-Pinene α-Terpinene y-Terpinene

Terpinolene **TOTAL** 39640

0.826

0.148

0.512

0.532

3.96



Abbreviations: HS - Headspace, GC -Gas Chromatography, MS - Mass Spectrometry, RL - Reporting Limit

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"-" Not detected above RL.

Reporting Limit (µg/g): 20





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RESIDUAL SOLVENT PROFILE

Analysis Date/Time: 10/21/2019 1234

Analyst: OL

2-Methylpentane (107-83-5)

3-Methylpentane (96-10-0)

m,p-Xylene (108-38-3 or 106-42-3)

n-Pentane (109-66-0)

n-Propane (74-98-6)

Toluene (108-88-3)

o-Xylene (95-47-6)

Xylenes* (1330-20-7)

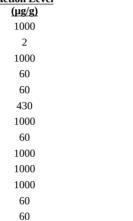
Method: USP <467> **Instrument:** Agilent 6890

Deviations from SOP:

None

<i>y</i>		
Solvent	<u>Result</u> (μg/g)	Action Level (μg/g)
Acetone (67-64-1)	-	1000
Benzene (71-43-2)	-	2
n-Butane (106-97-2)	-	1000
2,2-Dimethylbutane (75-83-2)	-	60
2,3-Dimethylbutane (79-29-8)	-	60
Ethyl benzene (100-41-4)	-	430
n-Heptane (142-82-5)	-	1000
n-Hexane (110-54-3)	-	60
Isobutane (75-28-5)	-	1000
Isopropanol (67-63-0)	-	1000
2-Methylbutane (78-78-4)	-	1000

DMSO



1000

1000

180

430

430

430

Color Key

RESULT < 1/2 AL
1/2 AL < RESULT < AL
RESULT > AL

Reporting Limit (μg/g) 1/2 of AL

"-" not detected above reporting limit

"*" - o,m,p-Xylene and Ethylbenzene



Solvent	Synonym(s)	Solvent	Synonym(s)	Solvent	Synonym(s)
Acetonitrile	Methyl Cyanide, ACN	2-Ethoxyethanol	Cellosolve, Ethyl glycol	Methanol	Methyl alcohol
1-Butanol	n-Butanol, Butyl	Ethyl ether	Diethyl ether, Ether	2-Methylbutane	Isopentane
1 Battarior	Alcohol	Ethyl acetate	EtOAc	Methylene chloride	Dichloromethane
2-Butanol	sec-Butyl alcohol	Ethyl benzene	Phenylethane	2-Methylpentane	Isohexane
2-Butanone	Methyl ethyl ketone, MEK	Ethylene glycol	1,2-Ethanediol	1-Pentanol	n-Amyl alcohol
1.2 Di		Ethylene oxide	Oxirane	1-Propanol	Propyl alcohol
1,2-Dimethoxyethane	Monoglyme	Isobutane	2-Methylpropane	Tetrahydrofuran	THF
2,3-Dimethylbutane	Neohexane	Icopropanol	2-Propanol, IPA	Tetramethylene sulfone	Sulfolane
2,3-Dimethylbutane	Diisopropyl	Isopropanol	1 ,	<i>y</i>	
N,N-Dimethylformamide	DMF	Isopropyl Acetate	Acetic acid isopropyl ester	Xylene	Dimethylbenzene

Abbreviations: HS-Headspace, GC-Gas Chromatography, MS-Mass Spectrometry, RL-Reporting Limit, AL-Action Level CAS-Chemical Abstract Services

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Dimethysufoxide



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Analyst: OL

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Analysis Date/Time: 10/21/2019 1234

Lab ID: SA2019-14844 **Date Received:** 10/17/2019 Sample Matrix: Concentrate **Date Completed:** 10/24/2019

Remarks:

PESTICIDES PROFILE

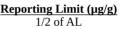
Method: LC/MS/MS and GC/MS **Deviations from SOP:**

Instrument: Waters Acquity/TQD

None

Pesticide Result **Action Level** (µg/g) $(\mu g/g)$ Abamectin (71751-41-2) 0.5 Azoxystrobin (131860-33-8) 0.5 Bifenazate* (149877-41-8) 0.5 Etoxazole (153233-91-1) 0.5 Imazalil (35554-44-0) 0.5 Imidacloprid (138261-41-3) 0.5 Malathion (121-75-5) 0.5 Myclobutanil (88671-89-0) 0.5 Permethrins* (52645-53-1) 0.5 Spinosad A (168316-95-8) 0.5 Spinosad D (168316-95-8) 0.5 Spiromesifen (283594-90-1) 0.5 Spirotetramat (203313-25-1) 0.5 Tebuconazole (80443-41-0) 0.5





"-" not detected above reporting limit

"*" analyzed by GC/MS (all others analyzed by LC/MS/MS)

Permethrins measured as the cumulative residue of the cis- and transpermethrin isomers.



Abbreviations: LC - Liquid Chromatography, GC - Gas Chromatography, MS - Mass Spectrometry, RL - Reporting Limit, AL - Action Level, CAS - Chemical Abstract Services

<u>Pesticide</u>	Synonym(s)	<u>Pesticide</u>	Synonym(s)	<u>Pesticide</u>	Synonym(s)
Cyfluthrin	Baythroid	Myclobutanil	Systhane	Propiconazole	Tilt
DDVP	Dichlorvos	Naled	Dibrom	Propoxur	Baygon
Ethoprophos	Prophos	Phosmet	Imidan		

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MYCOTOXIN PROFILE

Analysis Date/Time: 10/21/2019 1234

Analyst: OL

Method: LC/MS/MS

Instrument: Waters Acquity/TQD

Deviations from SOP:

None

Mycotoxin	<u>Result</u> (µg/kg)	<u>Action Level</u> (μg/kg)
Aflatoxin B1	-	20
Aflatoxin B2	-	20
Aflatoxin G1	-	20
Aflatoxin G2	-	20
Ochratoxin A		20
Aflatoxin B2 Aflatoxin G1 Aflatoxin G2	-	20 20 20 20



Abbreviations: EC - Escherichia Coli, CFU - Colony-Forming Unit,

RL - Reporting Limit, AL - Action Level

Color Key

RESULT < 1/2 AI 1/2 AL < RESULT < AL RESULT > AL

Reporting Limit (CFU/g)

"-" not detected above reporting limit

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Remarks:

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HEAVY METAL PROFILE

Analysis Date/Time: 10/21/2019 1234

Analyst: OL

Method: ICP/MS
Instrument: PerkinElmer Elan 9000

Deviations from SOP:

None

Heavy Metal

Arsenic (As)
Cadmium (Cd)
Lead (Pb)
Mercury (Hg)

<u>Result</u> (µg/kg)	<u>Action Level</u> (μg/kg)
-	400
-	440
-	1000
-	200



Abbreviations: ICP - Inductively-Coupled Plasma, OES - Optical Emission Spectroscopy, MS - Mass Spectroscopy, RL - Reporting Limit, AL - Action Level

Color Key

RESULT < 1/2 AL

1/2 AL < RESULT < AL

RESULT > AL

Reporting Limit (μg/kg)
50

"-" not detected above reporting limit

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Remarks:

MICROBIOLOGICAL PROFILE

Analysis Date/Time: 10/21/2019 1234

Method: Hardy Diagnostics CompactDry

Deviations from SOP:

Analyst: OL

Salmonella spp.

Instrument: Thermo Incubator

None

Bacteria/Microbe

Aerobic Plate Count, Total
Escherichia Coli (E. Coli)
Mold
Yeast

Result
(CFU/g)
50
Absent
Absent
Absent
Absent

Action Level (CFU/g) -1 10000 10000 1



Abbreviations: EC - Escherichia Coli, CFU - Colony-Forming Unit, RL - Reporting Limit, AL - Action Level, NT - Not Tested

Color Key

RESULT < 1/2 AL

1/2 AL < RESULT < AL

RESULT > AL

Reporting Limit (CFU/g)

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