

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

SUMMARY OF ANALYSIS (SAMPLE ID: SA27232)

Testing Location:	Customer ID: 710	Order ID: OR8487	Sample Type: Primary
OKC	Sunday Extracts	Lot Number:	Matrix: Concentrate
3680 E. I-240 Service Rd.	1006 North Oak St	Not Entered	Mass: 2g
Oklahoma City, OK 73135	Guthrie, OK 73044	Batch Number:	Date Collected: 05/18/2020
License: LAAA-4Y4X-Z72Z	License: PAAA-NYIN-5SXK	RDS507MDMC900	Date Received: 05/18/2020
Cultivar (Strain) or Sample I	Date Completed: 05/22/2020		

*This page is simply a summary of the analysis performed. For analytical details, please consult the individual Certificate(s) of Analysis for each of the specific test(s) performed.

*For Oklahoma, with the new OMMA rules to be effective April 1, 2020, limits on moisture are proposed at 15% and water activity at 0.65aw.

Moisture Content (%)	PASS/FAIL	Water Activity (aw)	PASS/FAIL	
Not Tested	N/A	Not Tested	N/A	
Cannabinoids (Гор 3)	<u>(%)</u>	mg/g	
THCa		78.3	783	
Δ9-THC		0.718	7	
CBDa		0.245	2	
TOTAL CB	D	0.215	2	
TOTAL TH	C	69.4	694	
TOTAL CANNAB	INOIDS	79.2	792	
<u>Terpenes (To</u>	<u>o 5)</u>	<u>(%)</u>	<u>µg/g</u>	
β-Caryophylle	ene	2.494	24943	
α-Humulen	e	1.308	13079	
β-Myrcene		0.496	4964	
d-Limonen	2	0.428	4283	
β-Pinene		0.269	2687	
TOTAL TERPE	INES	5.77	57683	
Contaminants	PASS/FAIL	Sample Picture Up	on Receipt	
Heavy Metals:	PASS			
Microbiology:	PASS		The second second	
			and the second	

Mycotoxins: Pesticides: Residual Solvents: Visual Inspection:

Sample Picture Upon Receipt	
No.24 Martin Martin BIDA Martin Martin	



Scan the QR code to verify results.

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PASS

PASS

PASS PASS

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Cultivar (Strain) or Sample 1	Date Completed: 05/22/2020		

Cultivar (Strain) or Sample Description: Mandarin Dream/ Mandarin Cookie

Reporting Limit

(mg/g)

0.505

0.505

0.505

0.505

0.505

0.505

0.505

0.505

0.505

0.505

0.505

Δ8-THC THCa THCv CBG CBGa CBC

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CANNABINOID (POTENCY) PROFILE

Result

(mg/mL)

Analysis Date/Time: 05/22/2020 1400 Analyst: WW

Result

(%)

0.245

0.718

78.3

79.2

69.4

0.215

Δ9-THC

Result

(mg/g)

2.45

_

7.18

783

792

694

2

Cannabinoid Distribution

(% of Total Cannabinoids)

Cannabinoid

CBD

CBDa

CBDv

 $\Delta 9$ -THC

 $\Delta 8$ -THC

THCa

THCv

CBC

CBG

CBGa

CBN

TOTAL

TOTAL THC

TOTAL CBD

100

80

60

40

20

Method: HPLC/DAD Instrument: Agilent 1100

Per Unit

(mg)

2

7

783

_ 792

694

2

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



UNIT MASS (g): 1

"-" Not detected above RL.

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) + Δ 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



medical-related symptoms.

CBD

CBDa CBDv

REPORT OF LABORATORY ANALYSIS

CBN

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Page 2 of 8 05/22/2020



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

Deviations from SOP:

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Cultivar (Strain) or Sample 1	Date Completed: 05/22/2020		

Analysis Date/Time: 05/22/2020 1400

TERPENOID PROFILE

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

Analyst: DJ		Instr	rument: Agilent 6890	None
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)		
α-Bisabolol	384	0.0384		
Camphene	1566	0.157		and the second s
δ-3-Carene	-	-		nulad Dan rayuda sabar
β-Caryophyllene	24943	2.49		151/174 4.046/10
Caryophyllene oxide	-	-		
p-Cymene	1242	0.124		
Eucalyptol	-	-		
Geraniol	-	-		
Guaiol	729	0.0729		Abbreviations: HS - Headspace, GC -
α-Humulene	13079	1.31		Gas Chromatography, MS - Mass Spectrometry, RL - Reporting Limit
Isopulegol	-	-		This information is provided as a
d-Limonene	4283	0.428		service and makes no claims of efficacy
Linalool	-	-		and/or safety of this product. Results are applicable only for the
β-Myrcene	4964	0.496		sample(s) analyzed and for the specific
cis-Nerolidol	-	-		analysis conducted.
trans-Nerolidol	-	-		This report is for informational purposes only and should not be used to
α-Ocimene	-	-		diagnose, treat, or prevent any
β-Ocimene	1560	0.156		medical-related symptoms. The statements and results herein have
α-Pinene	2246	0.225		not been approved and/or endorsed by
β-Pinene	2687	0.269		the FDA.
α-Terpinene	-	-		
γ-Terpinene	-	-		
Terpinolene	-	-		"-" Not detected above RL.
TOTAL	57683	5.77		Reporting Limit (µg/g): 20.2







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Cultivar (Strain) or Sample Description: Mandarin Dream/ Mandarin Cookie

Action Level

<u>(µg/g)</u>

1000

2

1000 60

60

430

1000

60

1000

Analysis Date/Time: 05/22/2020 1400 .

Method: USP <467>

Deviations from SOP:

None

Solvent	<u>Result</u> (µg/g)
Acetone (67-64-1)	-
Benzene (71-43-2)	-
n-Butane (106-97-2)	-
2,2-Dimethylbutane (75-83-2)	-
2,3-Dimethylbutane (79-29-8)	-
Ethyl benzene (100-41-4)	-
n-Heptane (142-82-5)	-
n-Hexane (110-54-3)	-
Isobutane (75-28-5)	-
Isopropanol (67-63-0)	-
2-Methylbutane (78-78-4)	-
2-Methylpentane (107-83-5)	-
3-Methylpentane (96-10-0)	-
n-Pentane (109-66-0)	-
n-Propane (74-98-6)	-
Toluene (108-88-3)	-
o-Xylene (95-47-6)	-
m,p-Xylene (108-38-3 or 106-42-3)	-
Xylenes* (1330-20-7)	-

RESIDUAL SOLVENT PROFILE

Instrument: Agilent 6890

Color Key

RESULT < 1/2 AI 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	<u>Synonym(s)</u>	<u>Solvent</u>	<u>Synonym(s)</u>	<u>Solvent</u>	<u>Synonym(s)</u>
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
	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 Dimethermothermo		Ethylene oxide	Oxirane	1-Propanol	Propyl alcohol
1,2-Dimethoxyethane	Monoglyme	Isobutane	2-Methylpropane	Tetrahydrofuran	THF
2,3-Dimethylbutane	Neohexane	Isopropanol	2-Propanol, IPA	Tetramethylene sulfone	Sulfolane
2,3-Dimethylbutane	Diisopropyl	1 1	1 ,	5	
N,N-Dimethylformamide	DMF	Isopropyl Acetate	Acetic acid isopropyl ester	Xylene	Dimethylbenzene
Dimethysufoxide	DMSO				

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

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Kyle W. Felling, Ph.D. ry Dir

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Analysis Date/Time: 05/22/2020 1400 Analyst: DJ

PESTICIDES PROFILE

Method: LC/MS/MS Instrument: Waters Acquity/TQD

Deviations from SOP:

None

<u>Pesticide</u>	<u>Result</u> (µg/g)	Action Level (µg/g)	<u>Color Key</u>
Abamectin (71751-41-2)	-	0.5	$\frac{\text{RESULT} < 1/2 \text{ AL}}{1/2 \text{ AL}}$
Azoxystrobin (131860-33-8)	-	0.5	1/2 AL < RESULT < AL
Bifenazate* (149877-41-8)	-	0.5	RESULT > AL
Etoxazole (153233-91-1)	-	0.5	Reporting Limit (µg/g)
Imazalil (35554-44-0)	-	0.5	1/2 of AL
Imidacloprid (138261-41-3)	-	0.5	
Malathion (121-75-5)	-	0.5	"-" not detected above
Myclobutanil (88671-89-0)	-	0.5	reporting limit
Permethrins* (52645-53-1)	-	0.5	"*" analyzed by GC/MS
Spinosad A (168316-95-8)	-	0.5	(all others analyzed
Spinosad D (168316-95-8)	-	0.5	by LC/MS/MS)
Spiromesifen (283594-90-1)	-	0.5	Permethrins measured as
Spirotetramat (203313-25-1)	-	0.5	the cumulative residue of
Tebuconazole (80443-41-0)	-	0.5	the <i>cis</i> - and <i>trans</i> - permethrin isomers.

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

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

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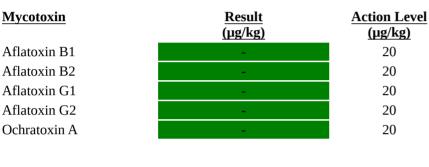
MYCOTOXIN PROFILE

Analysis Date/Time: 05/22/2020 1400 Analyst: DJ

Method: LC/MS/MS Instrument: Waters Acquity/TQD

Deviations from SOP:

None





Abbreviations: EC - Escherichia Coli, CFU - Colony-Forming Unit, RL - Reporting Limit, AL - Action Level

Color Key



Reporting Limit (CFU/g) 1

"-" not detected above reporting limit

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Analysis Date/Time: 05/22/2020 1400 Analyst: CC

HEAVY METAL PROFILE

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



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

Color Key



Reporting Limit (μg/kg) 50

"-" not detected above reporting limit

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MICROBIOLOGICAL PROFILE

Analysis Date/Time: 05/22/2020 1400 Analyst: DJ

Method: Hardy Diagnostics CompactDry **Instrument:** Thermo Incubator

Deviations from SOP: None

Bacteria/Microbe	<u>Result</u> (CFU/g)	<u>Action Level</u> (CFU/g)
Escherichia Coli (E. Coli)	Absent	1
Mold/Yeast	Absent	10000
Salmonella spp.	Absent	1
Staphylococcus Aureus	Absent	1



Abbreviations: EC - Escherichia Coli, CFU - Colony-Forming Unit, RL - Reporting Limit, AL - Action Level, TNTC - Too Numerous To Count, NT - Not Tested Absent - Not Detected Above RL, Present - Detected Above RL

Color Key



Reporting Limit (CFU/g) 1

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