

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: SA28312)

Testing Location:	Customer ID: 710	Order ID: OR8814	Sample Type: Primary
OKC	Sunday Extracts	Lot Number:	Matrix: Concentrate
3680 E. I-240 Service Rd.	1006 North Oak St	Not Entered	Mass: g
Oklahoma City, OK 73135	Guthrie, OK 73044	Batch Number:	Date Collected: 06/08/2020
License: LAAA-4Y4X-Z72Z	License: PAAA-NYIN-5SXK	RDS602CDMK1065	Date Received: 06/08/2020
Cultivar (Strain) or Sample I	Date Completed: 06/17/2020		

*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		82.2	822
Δ9-THC		2.48	25
CBGa		0.871	9
TOTAL CB	D	0.637	6
TOTAL TH	С	74.6	746
TOTAL CANNAB	INOIDS	86.6	866
Terpenes (To	p 5)	<u>(%)</u>	<u>µg/g</u>
β-Caryophyll	ene	0.813	8135
Guaiol		0.172	1724
Linalool		0.149	1490
β-Myrcene		0.138	1376
d-Limonen	е	0.135	1355
TOTAL TERPH	ENES	1.95	19456
Contaminants	PASS/FAIL	Sample Picture Up	on Receipt
Heavy Metals.	PASS		

Contaminants	PASS/FAIL
Heavy Metals:	PASS
Microbiology:	PASS
Mycotoxins:	PASS
Pesticides:	PASS
Residual Solvents:	PASS
Visual Inspection:	PASS





Scan the QR code to verify results.

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Cultivar (Strain) or Sample	Date Completed: 06/17/2020		

Cultivar (Strain) or Sample Description: Chem Dawg x Master Kush

CANNABINOID (POTENCY) PROFILE

Method: HPLC/DAD

Analysis Date/Time: 06/09/2020 1323 Analyst: WW

CBD

CBDa

CBDv

THCa

THCv

CBC

CBG

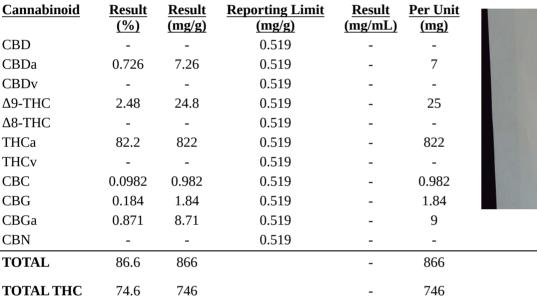
CBGa

CBN

Instrument: Agilent 1100

6

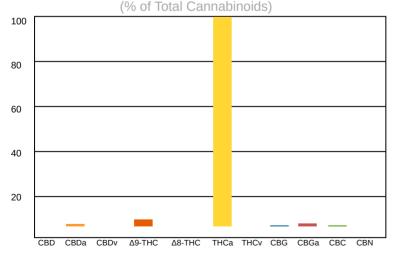
Moisture Content (%): -Water Activity (aw): -



TOTAL CBD 0.637

Cannabinoid Distribution

6



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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



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Cultivar (Strain) or Sample 1	Date Completed: 06/17/2020		

Cultivar (Strain) or	Sample Description	n: Chem Da	awg x Master Kush	Date Completed: 06/17/2020	
TERPENOID PROFILE					
Analysis Date/Time: 06/09/2020 1323Method: HS/GC/FID (International Analyst: DJAnalyst: DJInstrument: Agilent 6890		nod: HS/GC/FID (Internal Method rument: Agilent 6890	1-002) Deviations from SOP: None		
<u>Terpene</u>	<u>Result</u> (µg/g)	<u>Result</u> (%)			
α-Bisabolol	-	-		and the second se	
Camphene	274	0.0274		6	
δ-3-Carene	-	-		ching and but matter	
β-Caryophyllene	8135	0.813		A rate -	
Caryophyllene oxide	1195	0.12	1 - C		
p-Cymene	469	0.0469			
Eucalyptol	-	-			
Geraniol	-	-	•		
Guaiol	1724	0.172		Abbreviations: HS - Headspace, GC -	
α-Humulene	-	-		Gas Chromatography, MS - Mass Spectrometry, RL - Reporting Limit	
Isopulegol	730	0.073		This information is provided as a	
d-Limonene	1355	0.135	s	service and makes no claims of efficacy	
Linalool	1490	0.149		and/or safety of this product. Results are applicable only for the	
β-Myrcene	1376	0.138	s	sample(s) analyzed and for the specific	
cis-Nerolidol	-	-		analysis conducted. This report is for informational	
trans-Nerolidol	-	-	I	purposes only and should not be used to	
α-Ocimene	413	0.0413		diagnose, treat, or prevent any nedical-related symptoms.	
β-Ocimene	512	0.0512		The statements and results herein have	
α-Pinene	600	0.06	Г	not been approved and/or endorsed by	
β-Pinene	471	0.0471	t	he FDA.	
α-Terpinene	257	0.0257			
γ-Terpinene	-	-			
Terpinolene	456	0.0456	"	'-'' Not detected above RL.	
TOTAL	19456	1.95		Reporting Limit (µg/g): 20.7	







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Cultivar (Strain) or Sample Description: Chem Dawg x Master Kush

Analysis Date/Time: 06/09/2020 1323 Analyst: CC

Instrument: Agilent 6890

RESIDUAL SOLVENT PROFILE

Deviations from SOP:

None



Solvent	

<u>Solvent</u>	<u>Result</u> (µg/g)	<u>Action Level</u> <u>(μg/g)</u>
Acetone (67-64-1)	-	1000
Benzene (71-43-2)	-	2
n-Butane (106-97-2)	590.418	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
2-Methylpentane (107-83-5)	-	60
3-Methylpentane (96-10-0)	-	60
n-Pentane (109-66-0)	-	1000
n-Propane (74-98-6)	-	1000
Toluene (108-88-3)	-	180
o-Xylene (95-47-6)	-	430
m,p-Xylene (108-38-3 or 106-42-3)	-	430
Xylenes* (1330-20-7)	-	430

Met	hod: U	JSI) <	467	'>
. .					

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

<u>Solvent</u>	<u>Synonym(s)</u>	Solvent	Synonym(s)	Solvent	<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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Cultivar (Strain) or Sample Description: Chem Dawg x Master Kush

Analysis Date/Time: 06/09/2020 1323 Analyst: HL

PESTICIDES PROFILE

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

Deviations from SOP:

None

Pesticide	<u>Result</u> (µg/g)	<u>Action Level</u> (μg/g)	<u>Color Key</u>
Abamectin (71751-41-2)	-	0.5	$\frac{\mathbf{RESULT} < 1/2 \text{ AL}}{1/2 \text{ AL} < \mathbf{RESULT} < \mathbf{AL}}$
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	$\frac{1}{1/2}$ of AL
Imidacloprid (138261-41-3)	-	0.5	
Malathion (121-75-5)	-	0.5	"-" not detected above reporting limit
Myclobutanil (88671-89-0)	-	0.5	reporting mint
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	<u>Synonym(s)</u>
Cyfluthrin	Baythroid	Myclobutanil	Systhane	Propiconazole	Tilt
DDVP	Dichlorvos	Naled	Dibrom	Propoxur	Baygon
Ethoprophos	Prophos	Phosmet	Imidan		

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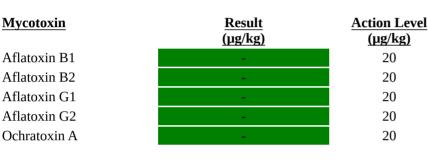
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MYCOTOXIN PROFILE

Analysis Date/Time: 06/09/2020 1323 Analyst: HL

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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HEAVY METAL PROFILE

Analysis Date/Time: 06/09/2020 1323 Analyst: CC

Method: ICP/MS Instrument: PerkinElmer Elan 9000 **Deviations from SOP:** None

Heavy Metal Arsenic (As)

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

<u>Result</u>	Action Level
<u>(µg/kg)</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: 06/09/2020 1323 Analyst: DJ

Method: Hardy Diagnostics CompactDry Instrument: Thermo Incubator

Deviations from SOP: None

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

Action Level (CFU/g) 1 10000 1 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

RESULT < 1/2 Al 1/2 AL < RESULT < AL**RESULT > AL**

Reporting Limit (CFU/g) 1

"-" not detected above reporting limit

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