

Product Name:	Calm Lotion		
Product Batch:	CL00203PH		
Certificate ID Number:	EVIO: 2007ELP0091.2505		
Date Tested:	07/24/2020		

Cannabinoid Profile & Potency Liquid Tincture:						
D9-THC:	9.91mg/unit					
CBD:	308mg/unit					
CBDA	12.9mg/unit					
CBDV:	LOQ					
CBG:	LOQ					
CBC:	10.9mg/unit					
CBN:	LOQ					
Total Count:	mg/unit					
Total THC:	9.91mg/unit					
Total CBD:	308mg/unit					
Manufactured by: Palme	Manufactured by: Palmetto Synergistic Research					

#### Manufacturer Date: 07/24/2020

Elemental Analysis:	Pass
Microbiological Contaminants:	Pass
Pathogenic Bacterial Contaminants:	Pass
Mycotoxin Testing:	Pass
Pesticide Analysis:	Pass
Terpene Profile:	Please see the full lab for multiple terpene profiles.
Analysis of Volatile Organic Compounds:	Pass

This product has been reviewed by ProVerde. The product contains less than 0.3% THC per the Farm Bill of 2018. This product is not intended to diagnose, treat, cure or prevent any disease. The FDA has not evaluated this product.



Quality Approval									
Prepared By/Date		Approved By/Date							
Mark Van	Date Signed: 8/18/2020	Direct of Operations David Newsom Quality Assurance Peter Girolamo	DocuSigned by: Date Signed: David MWSOM8/18/2020 489756D981174A2 DocuSigned by: Date Signed: Peter Givolame8/18/2020 17117FDA4E4B4C3						

This product has been approved by our Quality Assurance Team, Peter Girolamo. Our Director of Operations has reviewed the product and approves the product. This product passes our requirements for distribution to consumers.

This product has been reviewed by ProVerde. The product contains less than 0.3% THC per the Farm Bill of 2018. This product is not intended to diagnose, treat, cure or prevent any disease. The FDA has not evaluated this product.



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Calm Lotion Palmetto Synergistic Research Info Only- Edibles/Infused Project

Confident Cannabis ID: 2007ELP0091.2505 Sample ID: P200634-03 Matrix: Cannabinoid Product (solid) METRC Batch #: Sampling Method/SOP: Client Date Sampled: NA Date Accepted: 08/06/20 Harvest/Process Lot ID:



Batch ID: CL00203PH Batch Size (g): Unit for Sale: 100mL bottle Harvest/Production Date:

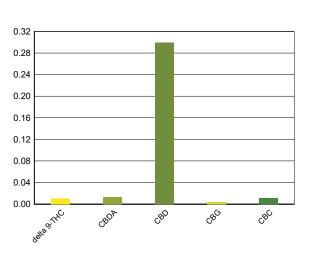
## **Cannabinoid Analysis**

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 Date/Time Extracted:
 07/24/20
 13:31
 Analysis Method/SOP: SOPT.40.023

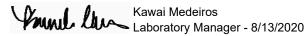
 Date/Time Analyzed:
 07/24/20
 15:34
 Sample mass: 99.14g/ 100mL bottle

Date/Time Analyz	200. 07724720 7	5.54	
Cannabinoids	LOQ(%)	mg/g	mg/unit
Total THC ((THCA*0.8	77)+∆9THC)	0.10	9.91
Total CBD ((CBDA*0	.877)+CBD)	3.10	308
THCA	0.005	< LOQ	< LOQ
delta 9-THC	0.005	0.10	9.91
delta 8-THC	0.005	< LOQ	< LOQ
THCV	0.005	< LOQ	< LOQ
CBGA	0.005	< LOQ	< LOQ
CBDA	0.005	0.13	12.9
CBD	0.005	2.99	296
CBDV	0.005	< LOQ	< LOQ
CBN	0.005	< LOQ	< LOQ
CBG	0.005	< LOQ	< LOQ
CBC	0.005	0.11	10.9
THCV-A	0.005	< LOQ	< LOQ
CBDV-A	0.005	< LOQ	< LOQ
Sum of tested Cannabinoids	0.005	3.34	331



**Cannabinoid Profile** 

"Total THC" and "Total CBD" are calculated values and are an Oregon reporting requirement (OAR 333-064-0100). For Cannabinoid analysis, only delta 9-THC, THCA, CBD, CBDA are ORELAP accredited analytes. Cannabinoid values reported for plant matter are dry weight corrected; Oregon Water Activity action level is 0.65Aw and Oregon Moisture Content action level is 15%, Samples above limit will be highlighted RED; FD = Field Duplicate; LOQ = Limit of Quantitation.



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# Calm Lotion

Palmetto Synergistic Research

Date/Time Extracted:

Info Only- Edibles/Infused Project Sample ID: P200634-03 METRC Batch #: Matrix: Cannabinoid Product

08/12/20 15:47

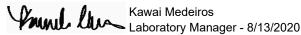
#### **Date Sampled: NA**

Date Accepted: 08/06/20 Batch ID: CL00203PH Batch Size: Sampling Method/SOP: Client

Terpene Analysis Analysis Method/SOP: SOP.T.40.092

Date/Time Analyzed:	08/13/20 0	9:48						
Analyte	LOQ (mg/g	LOQ (mg/g)Mass (mg/g) Mass (%)		Analyte	LOQ (mg/	ˈɡ͡ʃlass (mg/g)	Mass (%)	
alpha-Pinene	0.020	0.033	0.0033	beta-Pinene	0.020	< LOQ	< LOQ	
Camphene	0.020	< LOQ	< LOQ	Sabinene	0.020	< LOQ	< LOQ	
Sabinene hydrate	0.020	< LOQ	< LOQ	beta-Myrcene	0.020	0.080	0.008	
p-Mentha-1,5-diene	0.020	< LOQ	< LOQ	(+)-3-Carene	0.020	< LOQ	< LOQ	
alpha-Terpinene	0.020	< LOQ	< LOQ	gamma-Terpinene	0.020	< LOQ	< LOQ	
Limonene	0.020	2.47	0.247	Eucalyptol	0.020	0.159	0.0159	
Guaiol	0.020	< LOQ	< LOQ	Terpinolene	0.020	< LOQ	< LOQ	
Linalool	0.020	2.35	0.235	Camphor	0.020	0.093	0.0093	
(+)-Camphor	0.020	0.103	0.0103	(-)-Camphor	0.020	0.087	0.0087	
Isopulegol	0.020	< LOQ	< LOQ	Isoborneol	0.020	< LOQ	< LOQ	
Borneol	0.020	0.106	0.0106	Hexahydrothymol	0.020	< LOQ	< LOQ	
Geraniol	0.020	< LOQ	< LOQ	(+)-Pulegone	0.020	< LOQ	< LOQ	
Nerol	0.020	0.028	0.0028	cis-Nerolidol	0.020	< LOQ	< LOQ	
trans-Nerolidol	0.020	< LOQ	< LOQ	Geranyl acetate	0.020	0.047	0.0047	
alpha-Cedrene	0.020	< LOQ	< LOQ	trans-Caryophyllene	0.020	0.114	0.0114	
Caryophyllene Oxide	0.020	0.035	0.0035	alpha-Humulene	0.020	< LOQ	< LOQ	
Valencene	0.020	< LOQ	< LOQ	alpha-Farnesene	0.020	< LOQ	< LOQ	
beta-Farnesene	0.020	< LOQ	< LOQ	Cedrol	0.020	< LOQ	< LOQ	
alpha-Bisabolol	0.020	0.045	0.0045	Fenchone	0.020	< LOQ	< LOQ	
Fenchyl Alcohol	0.020	< LOQ	< LOQ	trans, beta- Ocimene	0.020	0.170	0.017	
beta, cis- Ocimene	0.020	0.071	0.0071	Terpineol	0.020	0.121	0.0121	
				Total (Sum):		6.11	0.61	

Analysis performed on GCMS with confirmation ion identification. Terpene analysis is not ORELAP accredited. Results reported as wet weight, or as is. LOQ = Limit of Quantitation.



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# Calm Lotion

Palmetto Synergistic ResearchInfo Only- Edibles/Infused ProjectSample ID: P200634-03METRC Batch #:

Matrix: Cannabinoid Product

#### Date Sampled: NA

Date Accepted: 08/06/20 Batch ID: CL00203PH Batch Size: Sampling Method/SOP: Client

#### Pesticides

Date/Time Analyzed: 8/6/2020 11:08:24PM

Date/Time Extracted: 08/06/20 13:41 Analysis Method/SOP: SOP.T.40.050 / SOP.T.40.051

Analyte	LOQ	Action Level	Result	Units	Туре
Abamectin	0.250	0.5	< LOQ	ppm	
Acephate	0.200	0.4	< LOQ	ppm	Organophosphate insecticide
Acequinocyl	1.00	2	< LOQ	ppm	
Acetamiprid	0.100	0.2	< LOQ	ppm	Neonicotinoid instecticide
Aldicarb	0.200	0.4	< LOQ	ppm	Carbamate insecticide
Azoxystrobin	0.100	0.2	< LOQ	ppm	
Bifenazate	0.100	0.2	< LOQ	ppm	Unclassified insecticide
Bifenthrin	0.100	0.2	< LOQ	ppm	
Boscalid	0.200	0.4	< LOQ	ppm	Anilide fungicide
Carbaryl	0.100	0.2	< LOQ	ppm	Carbamate insecticide
Carbofuran	0.100	0.2	< LOQ	ppm	Carbamate insecticide
Chlorantraniliprole	0.100	0.2	< LOQ	ppm	Anthranilic diamide insecticide
Chlorfenapyr	0.500	1	< LOQ	ppm	Pyrazole insecticide
Chlorpyrifos	0.100	0.2	< LOQ	ppm	Organophosphate insecticide
Clofentezine	0.100	0.2	< LOQ	ppm	
Cyfluthrin	0.500	1	< LOQ	ppm	
Cypermethrin	0.500	1	< LOQ	ppm	
Daminozide	0.500	1	< LOQ	ppm	
DVP (Dichlorvos)	0.500	1	< LOQ	ppm	
Diazinon	0.100	0.2	< LOQ	ppm	Organophosphate insecticide
Dimethoate	0.100	0.2	< LOQ	ppm	
Ethoprophos	0.100	0.2	< LOQ	ppm	
Etofenprox	0.200	0.4	< LOQ	ppm	
Etoxazole	0.100	0.2	< LOQ	ppm	Unclassified miticide
enoxycarb	0.100	0.2	< LOQ	ppm	
enpyroximate	0.200	0.4	< LOQ	ppm	
Fipronil	0.200	0.4	< LOQ	ppm	Pyrazole insecticide
Flonicamid	0.500	1	< LOQ	ppm	Pyridinecarboxamide insecticide
Fludioxonil	0.200	0.4	< LOQ	ppm	non-systemic fungicide
Hexythiazox	0.500	1	< LOQ	ppm	
mazalil	0.100	0.2	< LOQ	ppm	Azole fungicide
midacloprid	0.200	0.4	< LOQ	ppm	Neonicotinoid insectide
Kresoxim-methyl	0.200	0.4	< LOQ	ppm	
Malathion	0.100	0.2	< LOQ	ppm	
Vletalaxyl	0.100	0.2	< LOQ	ppm	
Methiocarb	0.100	0.2	< LOQ	ppm	Carbamate insecticide

Kawai Medeiros Laboratory Manager - 8/13/2020

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## **Calm Lotion**

Palmetto Synergistic Research Info Only- Edibles/Infused Project

Sample ID: P200634-03 METRC Batch #: Matrix: Cannabinoid Product

Date/Time Extracted: 08/06/20 13:41

#### Date Sampled: NA

Date Accepted: 08/06/20 Batch ID: CL00203PH Batch Size: Sampling Method/SOP: Client

#### Pesticides

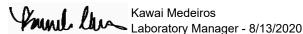
Date/Time Analyzed: 8/6/2020 11:08:24PM

Analysis Method/SOP:	SOP.T.40.050 / SOP.T.40.051

Analyte	LOQ	Action Level	Result	Units	Туре
Methomyl	0.200	0.4	< LOQ	ppm	Carbamate insecticide
Methyl parathion	0.100	0.2	< LOQ	ppm	
MGK-264	0.100	0.2	< LOQ	ppm	
Myclobutanil	0.100	0.2	< LOQ	ppm	Azole fungicide
Naled	0.250	0.5	< LOQ	ppm	
Oxamyl	0.500	1	< LOQ	ppm	Carbamate insecticide
Paclobutrazol	0.200	0.4	< LOQ	ppm	Azole plant growth regulator
Permethrins	0.100	0.2	< LOQ	ppm	
Phosmet	0.100	0.2	< LOQ	ppm	Organophosphate insecticide
Piperonyl butoxide	1.00	2	< LOQ	ppm	
Prallethrin	0.100	0.2	< LOQ	ppm	
Propiconazole	0.200	0.4	< LOQ	ppm	
Propoxur	0.100	0.2	< LOQ	ppm	Carbamate insecticide
Pyrethrins	0.500	1	< LOQ	ppm	
Pyridaben	0.100	0.2	< LOQ	ppm	Unclassified insecticide
Spinosad	0.100	0.2	< LOQ	ppm	Spinosyn insecticide
Spiromesifen	0.100	0.2	< LOQ	ppm	Keto-enol insecticide
Spirotetramat	0.100	0.2	< LOQ	ppm	Keto-enol insecticide
Spiroxamine	0.200	0.4	< LOQ	ppm	Unclassified fungicide
Tebuconazole	0.200	0.4	< LOQ	ppm	
Thiacloprid	0.100	0.2	< LOQ	ppm	
Thiamethoxam	0.100	0.2	< LOQ	ppm	Neonicotinoid insectide
Trifloxystrobin	0.100	0.2	< LOQ	ppm	Strobin fungicide

Results above the action level fail Oregon state testing requirements and will be highlighted RED.

LOQ= Limit of Quantitation; PPM= Parts per million; ND= Not detected; NT= Not tested; AC= Above calibration range. PASS/FAIL status based on OAR 333-007. Pesticide testing performed in conjunction with EVIO Labs Medford, an ORELAP accredited laboratory.



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# Calm Lotion

Palmetto Synergistic Research

Info Only- Edibles/Infused Project

Sample ID: P200634-03 METRC Batch #:

Matrix: Cannabinoid Product

#### **Date Sampled: NA**

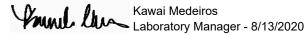
Date Accepted: 08/06/20 Batch ID: CL00203PH

Batch Size:

Sampling Method/SOP: Client

Analysis Method/SOP: SOP.T.40.031           sor-Butane         250         5000         < LOQ			R	esidual S	olvents	
Dutaties         250         250         1 COU         ppm         Analysis Method/SOP: SOP.T.40.031           iso-Butane         250         6000         < LOQ         ppm         Analysis Method/SOP: SOP.T.40.031           iso-Butane         174         290         < LOQ         ppm         and iso-butanes (CAS# 106-97-8). and iso-butane (CAS# 106-97-8).           2-Methylpentane         174         290         < LOQ         ppm         4. Total hexanes are calculated as           2.2.Dimethylbutane         174         290         < LOQ         ppm         4. Total hexanes are calculated as           2.2.Dimethylbutane         174         290         < LOQ         ppm         2methylpentane (CAS# 10-54-3),           2.3.Joinethylbutane         174         290         < LOQ         ppm         3. methylpentane (CAS# 10-85-3),           3.1.2.Dimethylbutane         174         290         < LOQ         ppm         3. methylpentane (CAS# 10-85-3),           1.2.Dimethylbenzene         1300         6000         < LOQ         ppm         3. methylpentane (CAS# 10-82-3),           1.3.Dimethylbenzene         1302         2170         < LOQ         ppm         and neo-pentane (CAS# 10-82-0),           1.3.Dimethylbenzene         1302         2170         < LOQ	Analyte	LOQ	Action Level	Result	Units	Date/Time Extracted: 07/31/20 09:28
Analysis Method/SOP: SOP.7.40.031           sor-Butane         250         5000         < LOQ         ppm         Analysis Method/SOP: SOP.7.40.031           sor-Butane         174         220         < LOQ         ppm         3 - Total butanes are calculated as sum of n-butanes (CAS# 106-97-8) and iso-butanes (CAS# 107-83-6).           Adethylpentane         174         220         < LOQ         ppm         4 - Total hexanes are calculated as sum of n-butanes (CAS# 110-54-3).           2.2-Dimethylbutane         174         290         < LOQ         ppm         sum of n-butanes (CAS# 110-54-3).           2.2-Dimethylbutane         174         290         < LOQ         ppm         sum of n-bexane (CAS# 110-54-3).           2.2-Dimethylbutane         174         290         < LOQ         ppm         sum of n-bexane (CAS# 110-54-3).           2.2-Dimethylbutane         174         290         < LOQ         ppm         sum of n-bexane (CAS# 110-54-3).           2.2-Dimethylbutane         174         290         < LOQ         ppm         2.3-dimethylbutane (CAS# 178-78-2).           2.2-Dimethylbutane         174         290         < LOQ         ppm         3-methylpentane (CAS# 108-40.).           1.2-Dimethylbenzene         1902         2170         < LOQ         ppm         3-methylpentane (C	Butanes	250	5000 <sup>3</sup>	< LOQ	ppm	Date/Time Analyzed: 07/31/20 12:23
Nexanes         174         290         4         LOQ         ppm         3- Total butanes are calculated as sum of n-butanes (CAS# 106-87-8) and iso-butanes (CAS# 106-87-8)         and iso-butanes (CAS# 106-87-8)         and iso-butanes (CAS# 106-87-8)         and iso-butanes (CAS# 106-87-8)         and iso-butanes (CAS# 110-84-3)         2-           2.2-Dimethylbutane         174         290         < LOQ	n-Butane	250	5000			Analysis Method/SOP: SOP.T.40.031
Hexanes         1/4         290         *         < LOQ         ppm         sum of n-butanes (CAS# 106-97-8) and iso-butane (CAS# 75-28-5)           2-Methylpentane         174         290         < LOQ	iso-Butane	250	5000	< LOQ	ppm	• Total but an an a shadata daa
n-Hexane       174       290       < LOQ	Hexanes	174	290 4	< LOQ	ppm	
2-Methylpentane         1/4         200         < LOQ	n-Hexane	174	290	< LOQ	ppm	( )
2.2 Dimethylbutane         174         290         < LOQ	2-Methylpentane	174	290	< LOQ	ppm	
2.2-Dimethylbutane         174         290         < LOQ	3-Methylpentane	174	290	< LOQ	ppm	4 - Total hexanes are calculated as
2.3-Dimethylbutane       174       290       < LOQ	2,2-Dimethylbutane	174	290	< LOQ	ppm	
Pentanes         1400         5000         *         < LOQ         ppm         3-methylpentane (CAS# 96-14-0), 2,2-dimethylbutane (CAS# 75-83-2), 3,2-dimethylbutane (CAS# 75-83-2), 2,2-dimethylbutane (CAS# 75-83-2), 3,3-dimethylbutane (CAS# 75-83-2), 3,3-dimethylbutane (CAS# 78-78-3)           Neopentane         250         5000         < LOQ	2,3-Dimethylbutane	174	290	< LOQ	ppm	
so-Pentane         1400         5000         < LOQ         ppm         2,3-dimethylbutane (CAS# 79-29-8)           Neopentane         250         5000         < LOQ	Pentanes	1400	5000 5	< LOQ	ppm	
Neopentane         250         5000         < LOQ         ppm           Xylenes         1302         2170         < LOQ	n-Pentane	1400	5000	< LOQ	ppm	2,2-dimethylbutane (CAS# 75-83-2),
Xylenes         1302         2170         < LOQ         ppm         5 - Total pentanes are calculated as           1.2-Dimethylbenzene         1302         2170         < LOQ	iso-Pentane	1400	5000	< LOQ	ppm	2,3-dimethylbutane (CAS# 79-29-8)
JDimethylbenzene         1302         2170         < LOQ         ppm         sum of n-pentane (CAS# 109-66-0), iso-pentane (CAS# 78-78-4), and neo-pentane (CAS# 78-78-4), and neo-pentane (CAS# 78-78-4), and neo-pentane (CAS# 78-78-4), and neo-pentane (CAS# 463-82-1)           Xylenes MP         1302         2170         < LOQ	Neopentane	250	5000	< LOQ	ppm	
1.3-Dimethylbenzene       1302       2170       < LOQ	Xylenes	1302	2170	< LOQ	ppm	5 - Total pentanes are calculated as
1.4-Dimethylbenzene       1302       2170       < LOQ	1,2-Dimethylbenzene	1302	2170	< LOQ	ppm	
1,4-Chined Hysich Zehro       1210       1200       ppm         2,4-Propanol (IPA)       1400       5000       < LOQ	1,3-Dimethylbenzene	1302	2170	< LOQ	ppm	,
Both Market         CLOQ         ppm         6 - Total xylenes are calculated as           22-Propanol (IPA)         1400         5000         < LOQ	1,4-Dimethylbenzene	1302	2170	< LOQ	ppm	and neo-pentane (CAS# 463-82-1)
Entry benzene       1302       1VA       < LOQ	Xylenes MP	1302	2170	< LOQ	ppm	• Tetal and an an and a lood at a days
2-Proparior (IPA)         1400         5000         < LOQ         ppm         1,3-dimethylbenzene (CAS# 106-42-3), and 1-4-dimethylbenzene (CAS# 106-42-3), and 1-4-dimethylbenzene (CAS# 106-42-3),           Acetonitrile         246         410         < LOQ	Ethyl benzene	1302	NA	< LOQ	ppm	5
Acetonic14005000< LOQppmand 1-4-dimethylbenzene (CAS# 106-42-3)Acetonitrile246410< LOQ	2-Propanol (IPA)	1400	5000	< LOQ	ppm	
Acctonitrile         246         410         < LOQ         ppm           Benzene         1.2         2         < LOQ	Acetone	1400	5000	< LOQ	ppm	
Methanol       1000       3000       < LOQ	Acetonitrile	246	410	< LOQ	ppm	
Methanol         1000         3000         < LOQ         ppm         OAR-333-007-0410.           Propane         250         5000         < LOQ	Benzene	1.2	2	< LOQ	ppm	7 - Ethanol is not regulated under
Propane         250         5000         < LOQ         ppm           Toluene         534         890         < LOQ	Methanol	1000	3000	< LOQ	ppm	
Dichloromethane       360       600       < LOQ	Propane	250	5000	< LOQ	ppm	
1,4-Dioxane       228       380       < LOQ	Toluene	534	890	< LOQ	ppm	
2-Butanol       1400       5000       < LOQ	Dichloromethane	360	600	< LOQ	ppm	
2-Ethoxyethanol       96       160       < LOQ	1,4-Dioxane	228	380	< LOQ	ppm	
Cumene       42       70       < LOQ	2-Butanol	1400	5000	< LOQ	ppm	
Cyclohexane       2278       3880       < LOQ	2-Ethoxyethanol	96	160	< LOQ	ppm	
Ethyl acetate14005000< LOQppmEthyl acetate14005000< LOQ	Cumene	42	70	< LOQ	ppm	
Ethyl ether14005000< LOQppmEthylene glycol558620< LOQ	Cyclohexane	2278	3880	< LOQ	ppm	
Ethylene glycol558620< LOQppmEthylene oxide3050< LOQ	Ethyl acetate	1400	5000	< LOQ	ppm	
Ethylene oxide3050< LOQppmHeptane14005000< LOQ	Ethyl ether		5000	< LOQ	ppm	
Heptane         1400         5000         < LOQ         ppm           Isopropyl acetate         1400         5000         < LOQ	Ethylene glycol		620	< LOQ	ppm	
Isopropyl acetate 1400 5000 < LOQ ppm Tetrahydrofuran 432 720 < LOQ ppm	Ethylene oxide	30	50	< LOQ	ppm	
Tetrahydrofuran 432 720 < LOQ ppm	Heptane	1400	5000	< LOQ	ppm	
	Isopropyl acetate	1400	5000		ppm	
Ethanol 1400 NA 7 < LOQ ppm	Tetrahydrofuran	432		< LOQ	ppm	
	Ethanol	1400	NA 7	< LOQ	ppm	

Results above the action level fail Oregon state testing requirements and will be highlighted RED. LOQ=Limit of Quantitation; PPM=Parts per million; ND=Not detected; NT=Not tested; AC=Above calibration range. PASS/FAIL status based on OAR 333-007.



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# Calm Lotion Date Sampled: NA Palmetto Synergistic Research Date Accepted: 08/06/20 Info Only- Edibles/Infused Project Batch ID: CL00203PH Sample ID: P200634-03 METRC Batch #: Batch Size: Matrix: Cannabinoid Product (solid) Sampling Method/SOP: Client

Yeast and Mold Enumeration

Date/Time Extracted: 07/30/20 10:20 Date/Time Analyzed: 08/05/20 13:54 Analysis Method/SOP: \*\*\* DEFAULT

Total Colonies: 0.00 CFU/g

#### About Your Yeast and Mold Results

Botanical materials often have total yeast and mold counts between 1,500 - 7,500 CFU/g. Products that have undergone exposure to solvents, such as alcohol tinctures or concentrated materials extracted with butane, propane, hexane, carbon dioxide, or other organic solvents will typically feature total yeast and mold counts at 0 CFU/g.

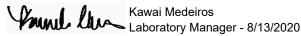
The American Herbal Pharmacoepia recommends herbal products contain no greater than 10,000 CFU/g of total yeasts and molds. Results above 10,000 CFU/g will be highlighted **Red**. Counts greater than 25,000 CFU/g are designated as "**TNTC**" or "Too numerous to count."

#### Yeasts vs Molds

Yeasts and molds are both broad types of fungi. Yeasts are unicellular and reproduce by budding, creating a small smooth apperance, whereas molds are multicellular and grow through fungal strands called hyphae, creating a fuzzy appearance often associated with mold.

Yeasts and molds are commonly found on natural products, and not all are harmful. Nevertheless, yeasts and molds, as well as their spores, can cause lung irritation, facilitate allergic reactions, or even present life-threatening conditions for immuno-compromised consumers. For instance, the dark mold, *Aspergillus*, can produce toxic chemical byproducts which can be harmful to human health. *Aspergillus* spores can lodge in small crevaces in the lungs and grow, leading to a potentially life-threatening condition called Aspergillosis.

A simple total yeast and mold count can be a great way to monitor for potential health hazards in botanical products and help ensure the safety of consumers.



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## **Quality Control**

#### Batch: M20H027 - SOP.T.30.060 Pesticide Prep

Blank(M20H027-BLK	(1)	E	Extracted: 08/06/20 13:41			Analyzed: 08/06/20 15:06		
Analyte	Result	LOQ	Recovery Limits	Analyte	Result	LOQ	Recovery Limits	
Methyl parathion	< LOQ	0.100 (ppm)	< LOQ	MGK-264	< LOQ	0.100 (ppm)	< LOQ	
Chlorfenapyr	< LOQ	0.500 (ppm)	< LOQ	Cyfluthrin	< LOQ	0.500 (ppm)	< LOQ	
Cypermethrin	< LOQ	0.500 (ppm)	< LOQ	Abamectin	< LOQ	0.250 (ppm)	< LOQ	
Acephate	< LOQ	0.200 (ppm)	< LOQ	Acequinocyl	< LOQ	1.00 (ppm)	< LOQ	
cetamiprid	< LOQ	0.100 (ppm)	< LOQ	Aldicarb	< LOQ	0.200 (ppm)	< LOQ	
zoxystrobin	< LOQ	0.100 (ppm)	< LOQ	Bifenazate	< LOQ	0.100 (ppm)	< LOQ	
ifenthrin	< LOQ	0.100 (ppm)	< LOQ	Boscalid	< LOQ	0.200 (ppm)	< LOQ	
arbaryl	< LOQ	0.100 (ppm)	< LOQ	Carbofuran	< LOQ	0.100 (ppm)	< LOQ	
hlorantraniliprole	< LOQ	0.100 (ppm)	< LOQ	Chlorpyrifos	< LOQ	0.100 (ppm)	< LOQ	
lofentezine	< LOQ	0.100 (ppm)	< LOQ	Daminozide	< LOQ	0.500 (ppm)	< LOQ	
DVP (Dichlorvos)	< LOQ	0.500 (ppm)	< LOQ	Diazinon	< LOQ	0.100 (ppm)	< LOQ	
imethoate	< LOQ	0.100 (ppm)	< LOQ	Ethoprophos	< LOQ	0.100 (ppm)	< LOQ	
tofenprox	< LOQ	0.200 (ppm)	< LOQ	Etoxazole	< LOQ	0.100 (ppm)	< LOQ	
enoxycarb	< LOQ	0.100 (ppm)	< LOQ	Fenpyroximate	< LOQ	0.200 (ppm)	< LOQ	
pronil	< LOQ	0.200 (ppm)	< LOQ	Flonicamid	< LOQ	0.500 (ppm)	< LOQ	
udioxonil	< LOQ	0.200 (ppm)	< LOQ	Hexythiazox	< LOQ	0.500 (ppm)	< LOQ	
nazalil	< LOQ	0.100 (ppm)	< LOQ	Imidacloprid	< LOQ	0.200 (ppm)	< LOQ	
resoxim-methyl	< LOQ	0.200 (ppm)	< LOQ	Malathion	< LOQ	0.100 (ppm)	< LOQ	
letalaxyl	< LOQ	0.100 (ppm)	< LOQ	Methiocarb	< LOQ	0.100 (ppm)	< LOQ	
lethomyl	< LOQ	0.200 (ppm)	< LOQ	Myclobutanil	< LOQ	0.100 (ppm)	< LOQ	
aled	< LOQ	0.250 (ppm)	< LOQ	Oxamyl	< LOQ	0.500 (ppm)	< LOQ	
aclobutrazol	< LOQ	0.200 (ppm)	< LOQ	Permethrins	< LOQ	0.100 (ppm)	< LOQ	
hosmet	< LOQ	0.100 (ppm)	< LOQ	Piperonyl butoxide	< LOQ	1.00 (ppm)	< LOQ	
rallethrin	< LOQ	0.100 (ppm)	< LOQ	Propiconazole	< LOQ	0.200 (ppm)	< LOQ	
ropoxur	< LOQ	0.100 (ppm)	< LOQ	Pyridaben	< LOQ	0.100 (ppm)	< LOQ	
yrethrins	< LOQ	0.500 (ppm)	< LOQ	Spinosad	< LOQ	0.100 (ppm)	< LOQ	
piromesifen	< LOQ	0.100 (ppm)	< LOQ	Spirotetramat	< LOQ	0.100 (ppm)	< LOQ	
piroxamine	< LOQ	0.200 (ppm)	< LOQ	Tebuconazole	< LOQ	0.200 (ppm)	< LOQ	
niacloprid	< LOQ	0.100 (ppm)	< LOQ	Thiamethoxam	< LOQ	0.100 (ppm)	< LOQ	
rifloxystrobin	< LOQ	0.100 (ppm)	< LOQ					
LCS(M20H027-BS1)		Ex	ctracted: 08/0	6/20 13:41	Analyzed: 08/06	6/20 15:33		

LCS(M20H027-BS1)		E	Extracted: 08/06/20 13:41			Analyzed: 08/06/20 15:33		
Analyte	% Recovery	LOQ	Recovery Limits	Analyte	% Recovery	LOQ	Recovery Limits	
Methyl parathion	137	0.100 (ppm)	50-150	MGK-264	157	0.100 (ppm)	50-150	
Chlorfenapyr	171	0.500 (ppm)	50-150	Cyfluthrin	79.7	0.500 (ppm)	50-150	
Cypermethrin	64.2	0.500 (ppm)	50-150	Abamectin	56.4	0.250 (ppm)	50-150	
Acephate	71.6	0.200 (ppm)	50-150	Acequinocyl		1.00 (ppm)	50-150	

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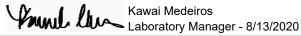
## **Quality Control**

#### Batch: M20H027 - SOP.T.30.060 Pesticide Prep (Continued)

LCS(M20H027-BS1)		Extracted: 08/06/20 13:41			Analyzed: 08/06/20 20:02		
Analyte	% Recovery	Recovery LOQ Limits Analyte		Analyte	% Recovery LOQ		Recovery Limits
Acetamiprid	83.9	0.100 (ppm)	50-150	Aldicarb	76.6	0.200 (ppm)	50-150
Azoxystrobin	82.0	0.100 (ppm)	50-150	Bifenazate	93.3	0.100 (ppm)	50-150
Bifenthrin	84.4	0.100 (ppm)	50-150	Boscalid	120	0.200 (ppm)	50-150
Carbaryl	53.5	0.100 (ppm)	50-150	Carbofuran	58.7	0.100 (ppm)	50-150
Chlorantraniliprole	154	0.100 (ppm)	50-150	Chlorpyrifos	127	0.100 (ppm)	50-150
Clofentezine	104	0.100 (ppm)	50-150	Daminozide	72.6	0.500 (ppm)	50-150
DVP (Dichlorvos)	77.6	0.500 (ppm)	50-150	Diazinon	151	0.100 (ppm)	50-150
Dimethoate	80.2	0.100 (ppm)	50-150	Ethoprophos	61.1	0.100 (ppm)	50-150
Etofenprox	118	0.200 (ppm)	50-150	Etoxazole	154	0.100 (ppm)	50-150
enoxycarb	107	0.100 (ppm)	50-150	Fenpyroximate	103	0.200 (ppm)	50-150
ipronil	101	0.200 (ppm)	50-150	Flonicamid	46.7	0.500 (ppm)	50-150
Iudioxonil	93.9	0.200 (ppm)	50-150	Hexythiazox	85.5	0.500 (ppm)	50-150
mazalil	143	0.100 (ppm)	50-150	Imidacloprid	54.7	0.200 (ppm)	50-150
Kresoxim-methyl	115	0.200 (ppm)	50-150	Malathion	123	0.100 (ppm)	50-150
letalaxyl	88.1	0.100 (ppm)	50-150	Methiocarb	78.5	0.100 (ppm)	50-150
/lethomyl	108	0.200 (ppm)	50-150	Myclobutanil	120	0.100 (ppm)	50-150
laled	76.1	0.250 (ppm)	50-150	Oxamyl	85.1	0.500 (ppm)	50-150
Paclobutrazol	118	0.200 (ppm)	50-150	Permethrins		0.100 (ppm)	50-150
Phosmet	127	0.100 (ppm)	50-150	Piperonyl butoxide	102	1.00 (ppm)	50-150
Prallethrin	200	0.100 (ppm)	50-150	Propiconazole	145	0.200 (ppm)	50-150
Propoxur	60.8	0.100 (ppm)	50-150	Pyridaben	78.9	0.100 (ppm)	50-150
Pyrethrins	91.7	0.500 (ppm)	50-150	Spinosad	126	0.100 (ppm)	50-150
piromesifen	154	0.100 (ppm)	50-150	Spirotetramat	130	0.100 (ppm)	50-150
Spiroxamine	171	0.200 (ppm)	50-150	Tebuconazole	147	0.200 (ppm)	50-150
Thiacloprid	86.0	0.100 (ppm)	50-150	Thiamethoxam	77.9	0.100 (ppm)	50-150
Frifloxystrobin	118	0.100 (ppm)	50-150				

#### Batch: P20H040 - SOP.T.40.092 PDX Terpenoid Analysis via GC-MS

Blank(P20H040-BLK1)		Extracted: 08/12/20 15:47			Analyzed: 08/13/20 09:48		
Analyte	Result	LOQ	Recovery Limits	Analyte	Result	LOQ	Recovery Limits
alpha-Pinene	< LOQ	0.200 (mg/g)	< LOQ	beta-Pinene	< LOQ	0.200 (mg/g)	< LOQ
Camphene	< LOQ	0.200 (mg/g)	< LOQ	Sabinene	< LOQ	0.200 (mg/g)	< LOQ
Sabinene hydrate	< LOQ	0.200 (mg/g)	< LOQ	beta-Myrcene	< LOQ	0.200 (mg/g)	< LOQ
p-Mentha-1,5-diene	< LOQ	0.200 (mg/g)	< LOQ	(+)-3-Carene	< LOQ	0.200 (mg/g)	< LOQ
alpha-Terpinene	< LOQ	0.200 (mg/g)	< LOQ	gamma-Terpinene	< LOQ	0.200 (mg/g)	< LOQ
Limonene	< LOQ	0.200 (mg/g)	< LOQ	Eucalyptol	< LOQ	0.200 (mg/g)	< LOQ
Guaiol	< LOQ	0.200 (mg/g)	< LOQ	Terpinolene	< LOQ	0.200 (mg/g)	< LOQ



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## **Quality Control**

#### Batch: P20H040 - SOP.T.40.092 PDX Terpenoid Analysis via GC-MS (Continued)

Blank(P20H040-BLK1)		Extracted: 08/12/20 15:47			Analyzed: 08/13/20 09:48		
Analyte	Result	LOQ	Recovery Limits	Analyte	Result	LOQ	Recovery Limits
Linalool	< LOQ	0.200 (mg/g)	< LOQ	Camphor	< LOQ	0.200 (mg/g)	< LOQ
(+)-Camphor	< LOQ	0.200 (mg/g)	< LOQ	(-)-Camphor	< LOQ	0.200 (mg/g)	< LOQ
lsopulegol	< LOQ	0.200 (mg/g)	< LOQ	Isoborneol	< LOQ	0.200 (mg/g)	< LOQ
Borneol	< LOQ	0.200 (mg/g)	< LOQ	Hexahydrothymol	< LOQ	0.200 (mg/g)	< LOQ
Geraniol	< LOQ	0.200 (mg/g)	< LOQ	(+)-Pulegone	< LOQ	0.200 (mg/g)	< LOQ
Verol	< LOQ	0.200 (mg/g)	< LOQ	cis-Nerolidol	< LOQ	0.200 (mg/g)	< LOQ
rans-Nerolidol	< LOQ	0.200 (mg/g)	< LOQ	Geranyl acetate	< LOQ	0.200 (mg/g)	< LOQ
Ipha-Cedrene	< LOQ	0.200 (mg/g)	< LOQ	trans-Caryophyllene	< LOQ	0.200 (mg/g)	< LOQ
Caryophyllene Oxide	< LOQ	0.200 (mg/g)	< LOQ	alpha-Humulene	< LOQ	0.200 (mg/g)	< LOQ
/alencene	< LOQ	0.200 (mg/g)	< LOQ	alpha-Farnesene	< LOQ	0.200 (mg/g)	< LOQ
oeta-Farnesene	< LOQ	0.200 (mg/g)	< LOQ	Cedrol	< LOQ	0.200 (mg/g)	< LOQ
alpha-Bisabolol	< LOQ	0.200 (mg/g)	< LOQ	Fenchone	< LOQ	0.200 (mg/g)	< LOQ
enchyl Alcohol	< LOQ	0.200 (mg/g)	< LOQ	trans, beta- Ocimene	< LOQ	0.200 (mg/g)	< LOQ
eta, cis- Ocimene	< LOQ	0.200 (mg/g)	< LOQ	Terpineol	< LOQ	0.200 (mg/g)	< LOQ

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## Certificate of Analysis For R+D Use Only



<b>EVIO</b> LAI	BS	Mycotoxin Analy	/sis Report	R&D Use only. Not fo Compliance		
Palmetto S	Synergistic Resea	rch EVIO Se	h EVIO Sample ID: Product Name:		34-03	
	Info Only	Produ			otion	
Batch ID Batch Size				Ordered: Sampled: Completed:	7/24/2020 NA 8/4/2020	
Mycotoxin Ana	alysis					
Analyte	LOQ (ug/mL)	Results (ug/mL)				
Aflatoxin B1	0.025	<loq< td=""><td></td><td></td><td></td></loq<>				
Aflatoxin B2	0.025	<loq< td=""><td></td><td></td><td></td></loq<>				
Aflatoxin G1	0.025	<loq< td=""><td></td><td></td><td></td></loq<>				
Aflatoxin G2	0.025	<loq< td=""><td></td><td></td><td></td></loq<>				
Ochratoxin A	0.100	<loq< td=""><td></td><td></td><td></td></loq<>				
Mycotoxin Analytica	l Batch ID : N	Л20Н006				
<i>Votes:</i> LCS recoveries for a Applicable.		e recoveries <20% RSD; Sample and	d solvent blanks <loq< td=""><td>(or ND); LOQ = Limit of (</td><td>Quantitation; NA = Not</td></loq<>	(or ND); LOQ = Limit of (	Quantitation; NA = Not	
<b>EVIO</b> LABS	540 E. Vilas Rd., Suite F Central Point, OR 97502	St	Stoph			
	www.eviolabs.com		Stephanie Moon Lab Director			
	541.668.7444	proval from EVIO Labs, Inc., and Kenevir Resear	Lab Dire	ctor		