

Lot#22c0165011

issue date 2/17/22 12:01 PM

This Product Has Been Tested and Complies with 7USC1639o(1)

Stillwater Laboratories

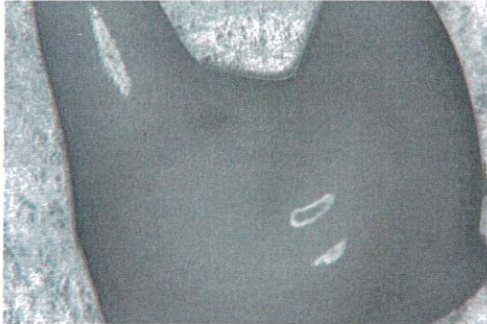
13142
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total cannabinoids per mL
10.2mg

total THC‡ ND
total CBD‡ 9.4mg

Incoming Inspection MSP-7.5.1.2

DESCRIPTION: Topical sample received 2/15/2022 2:44:22 PM in a client-labeled bottle, by commercial courier per Method 7.3.1.1. and as described in the Montana METRC Lab User Guide. Labeled Lot.



- caryophyllene 
- humulene 
- terpinolene 
- ocimene 
- beta pinene 
- alpha pinene 
- limonene 
- myrcene 
- linalool 

SECURITY FEATURE: WATERMARK MUST MATCH CERTIFICATE ID AND ISSUE DATE

Potency	MSP-7.5.1.4	per mL	LOD	LOQ	error
total cannabinoids		10.2mg	0.02	0.06	±0.24mg
total THC‡		ND	0.02	0.06	±0.06mg
total THC (THC+THCa)		ND	0.02	0.06	±0.06mg
total CBD‡		9.4mg	0.02	0.06	±0.23mg
total CBD (CBD+CBDA)		9.4mg	0.02	0.06	±0.23mg
tetrahydrocannabinolic acid (THCa)		ND	0.02	0.06	±0.06mg
Δ9-tetrahydrocannabinol (Δ9 THC)		ND	0.02	0.06	±0.06mg
Δ8-tetrahydrocannabinol (Δ8 THC)*		ND	0.03	0.08	±0.08mg
tetrahydrocannabivarin (THCv)		ND	0.02	0.06	±0.06mg
cannabidiolic acid (CBDA)		ND	0.02	0.05	±0.05mg
cannabidiol (CBD)		9.4mg	0.02	0.06	±0.23mg
cannabidivarin (CBDv)		<LOQ	0.02	0.06	±0.06mg
cannabigerolic acid (CBGA)		ND	0.02	0.05	±0.05mg
cannabigerol (CBG)		0.3mg	0.01	0.02	±0.02mg
cannabinol (CBN)		0.3mg	0.01	0.03	±0.04mg
cannabichromene (CBC)		0.2mg	0.02	0.06	±0.06mg

Terpenes	MSP-7.5.1.6	LOD	LOQ	error
total terpenes	ND	0.0008	0.0025	±0.0025%
linalool	ND	<0.0010	0.0012	±0.0012%
β-myrcene	ND	0.0009	0.0027	±0.0027%
D-limonene	ND	0.0005	0.0016	±0.0016%
α-pinene	ND	<0.0010	0.0010	±0.0010%
β-pinene	ND	0.0007	0.0022	±0.0022%
ocimene	ND	0.0015	0.0045	±0.0045%
terpinolene	ND	0.0011	0.0032	±0.0032%
α-humulene	ND	0.0007	0.0022	±0.0022%
β-caryophyllene	ND	0.0020	0.0061	±0.0061%
α-bisabolol	ND	0.0016	0.0049	±0.0049%
camphene	ND	0.0009	0.0028	±0.0028%
Δ3-carene	ND	0.0040	0.0121	±0.0121%
caryophyllene oxide	ND	0.0026	0.0078	±0.0078%
para-cymene	ND	0.0152	0.0457	±0.0457%
eucalyptol	ND	0.0021	0.0064	±0.0064%
geraniol	ND	0.0064	0.0191	±0.0191%
guaial	ND	0.0020	0.0061	±0.0061%
isopulegol	ND	0.0020	0.0060	±0.0060%
cis-nerolidol	ND	0.0034	0.0102	±0.0102%
trans-nerolidol	ND	0.0017	0.0051	±0.0051%
α-terpinene	ND	0.0011	0.0032	±0.0032%

Pass / Fail Criteria

Microbial (Plating) MSP-7.5.1.10

FAIL: no failures
PASS: E.coli, Salmonella sp., molds

Mycotoxins MSP-7.5.1.8

FAIL: no failures
PASS: Ochratoxin A, Aflatoxin B1B2G1G2, Aflatoxin B1, Aflatoxin B2, Aflatoxin G1.

Moisture MSP-7.5.1.3
not required / not requested

Metals MSP-7.5.1.7

FAIL: no failures
PASS: Arsenic, Cadmium, Lead, Mercury

Pesticides MSP-7.5.1.8

FAIL: no failures
PASS: Abamectin, Acephate, Acequinocyl, Acetamiprid, Aldicarb, Azoxystrobin, Bifenazate, Bifenthrin, Boscalid, Carbaryl, Carbofuran, Chloanthraniliprole, Chlorpyrifos, Clofentezine, Coumaphos, Cyfluthrin, Cypermethrin, Daminozide, Dichlorvos, Diazinon, Dimethoate, Etoxazole, Fenoxycarb, Fenpyroximate, Fipronil, Flonicamid, Fludioxonil, Hexythiazox, Imazalil, Imidacloprid, Malathion, Metalaxyl, Methiocarb, Methomyl, Mevinphos, Myclobutanil, Naled, Oxamyl, Pacllobutrazol, Permethrin, Phosmet, Piperonylbutoxide, Prallethrin, Propiconazole, Propoxur, Pyrethrin, Pyridaben, Spinetoram, Spinosad, Spiromesifen, Spirotetramat, Spiroxamine, Tebuconazole, Thiocloprid, Thiamethoxam, Trifloxystrobin

Solvents MSP-7.5.1.7

FAIL: no failures
PASS: Acetone, Acetonitrile, Benzene, Butane, Chloroform, Cyclohexane, Ethanol, Heptane, Hexane, Isopropyl alcohol, Methanol, Pentane, Propane, Toluene, Xylenes

Analysis Location: L-00001

Certified by:



https://customer.a2la.org/index.cfm?event=directory_detail&tabPID=423635B2-5128-4C6F-871A-419DCF43B0D7

Stillwater Laboratories Inc.

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These results are only valid for the samples tested. • Potency (cannabinoid concentration) is calculated as: [cannabinoid] = [cannabinoid]_{HPLC} x volume_{dilution}/m_{dry}. •• Decarboxyated cannabinoid concentration is calculated XXX_{total} = 0.877 x XXX_a + XXX ••• Standards are used to calibrate the resulting data and estimate error using a standard estimate of error method; LOD is the limit of detection (3.3s), LOQ is the limit of quantification (3xLOD), and experimental error is calculated from weighing, dilution, and interpolation error using the formula s_g² = Σ (∂f/∂i)²s_i² where i is the contributor to error. The 95% confidence range is calculated from: (concentration) ± t_{CL90} x s_g. Sampling error is not considered in error calculations. ND = not detected (< LOD), NT = not tested, NL = no limit, NA = not applicable. ‡ = decarbed, * = analyte is off-scope.

Corganics CBD Pain Cream

Lot#22c0165011

Report Version: 1
Analysis Location: L-00001

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rec'd date 2/15/2022 2:44:22 PM
issue date 2/17/2022 12:01:16 PM

Methods and Instruments

MSP-7.3.1.1	BAL-05	2/15/2022	L-00001	MSP-7.5.1.7	QP2020/HS20	2/16/2022	L-00001
MSP-7.5.1.2	YSC HD801m12	2/16/2022		MSP-7.5.1.8	LCMS8060	2/16/2022	
MSP-7.5.1.3	6MOC63u	2/16/2022		MSP-7.5.1.8	LCMS8060	2/16/2022	
MSP-7.5.1.4	LC-2030C	2/16/2022	L-00001	MSP-7.5.1.10A	Hardy Diag	2/16/2022	
MSP-7.5.1.6	QP2020/HS20	2/16/2022	L-00001	MSP-7.5.1.11	ICPMS2030	2/16/2022	

Mycotoxins

MSP-7.5.1.8	limit	LOD	LOQ	error	result	
Ochratoxin A	ND	20 ppb	0.3	0.8	±0.8 ppb	PASS
Aflatoxin B1B2G1G2	ND	20 ppb	0.3	0.8	±0.8 ppb	PASS

Microbial (Plating)

MSP-7.5.1.10	limit	LOD	LOQ	error	result	
E.coli	ND	0CFU	0.0	0.1	±0.1CFU	PASS
Salmonella sp.	ND	0CFU	0.0	0.1	±0.1CFU	PASS
molds	ND	10000CFU	1.6	4.8	±4.8CFU	PASS

Solvents

MSP-7.5.1.7	limit	LOD	LOQ	error	result	
Acetone	ND	5000 ppm	0.7	2.1	±2.1 ppm	PASS
Acetonitrile	ND	410 ppm	0.6	1.9	±1.9 ppm	PASS
Benzene	ND	0 ppm	0.0	0.1	±0.1 ppm	PASS
Butane	ND	5000 ppm	1.4	4.2	±4.2 ppm	PASS
Chloroform	ND	0 ppm	0.1	0.2	±0.2 ppm	PASS
Cyclohexane	ND	0 ppm	0.5	1.6	±1.6 ppm	PASS
Ethanol	ND	10000 ppm	0.7	2.2	±2.2 ppm	PASS
Heptane	ND	5000 ppm	0.4	1.3	±1.3 ppm	PASS
Hexane	ND	290 ppm	0.5	1.6	±1.6 ppm	PASS
Isopropyl alcohol	ND	5000 ppm	0.6	1.9	±1.9 ppm	PASS
Methanol	ND	3000 ppm	0.5	1.6	±1.6 ppm	PASS
Pentane	ND	5000 ppm	0.2	0.6	±0.6 ppm	PASS
Propane	ND	5000 ppm	0.5	1.6	±1.6 ppm	PASS
Toluene	ND	890 ppm	0.3	0.9	±0.9 ppm	PASS
Xylenes	ND	2170 ppm	0.3	1.0	±1.0 ppm	PASS

Metals

MSP-7.5.1.7	limit	LOD	LOQ	error	result	
Arsenic	ND	1500 ppb	7.8	23.4	±23.4 ppt	PASS
Cadmium	ND	500 ppb	8.4	25.2	±25.2 ppt	PASS
Lead	ND	500 ppb	13.1	39.2	±39.2 ppt	PASS
Mercury	ND	300 ppb	6.6	19.7	±19.7 ppt	PASS

These results are only valid for the samples tested. • Standards are used to calibrate the resulting data and estimate error using a standard estimate of error method; LOD is the limit of detection (3.3s_g), LOQ is the limit of quantification (3xLOD), and experimental error is calculated from weighing, dilution, and interpolation error using the formula s_g² = Σ(∂f/∂i)²s_i² where i is the contributor to error. The 95% confidence range is calculated from: (concentration) ± t_{CL90} × s_g. Sampling error is not considered in error calculations. ND = not detected (< LOD), NT = not tested, NL = no limit, NA = not applicable.

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Pesticides

MSP-7.5.1.8	limit	LOD	LOQ	error	result	
Abamectin	ND	0.30 ppm	0.005	0.014	±0.014 ppm	PASS
Acephale	ND	5.00 ppm	0.005	0.014	±0.014 ppm	PASS
Acequinocyl	ND	4.00 ppm	0.004	0.012	±0.012 ppm	PASS
Acetamiprid	ND	5.00 ppm	0.003	0.010	±0.010 ppm	PASS
Aldicarb	ND	0.00 ppm	0.001	0.004	±0.004 ppm	PASS
Azoxystrobin	ND	40.00 ppm	0.001	0.004	±0.004 ppm	PASS
Bifenazate	ND	5.00 ppm	0.001	0.003	±0.003 ppm	PASS
Bifenthrin	ND	0.50 ppm	0.001	0.002	±0.002 ppm	PASS
Boscalid	ND	10.00 ppm	0.013	0.040	±0.040 ppm	PASS
Carbaryl	ND	0.50 ppm	0.005	0.016	±0.016 ppm	PASS
Carbofuran	ND	0.00 ppm	0.001	0.003	±0.003 ppm	PASS
Chloanthraniliprole	ND	40.00 ppm	0.013	0.038	±0.038 ppm	PASS
Chlorfenapyr	NT	0.00 ppm				NA
Chlorpyrifos	ND	0.00 ppm	0.026	0.079	±0.079 ppm	PASS
Clofentezine	ND	0.50 ppm	0.005	0.015	±0.015 ppm	PASS
Coumaphos	ND	0.00 ppm	0.003	0.010	±0.010 ppm	PASS
Cyfluthrin	ND	1.00 ppm	0.005	0.014	±0.014 ppm	PASS
Cypermethrin	ND	1.00 ppm	0.003	0.010	±0.010 ppm	PASS
Daminozide	ND	0.00 ppm	0.018	0.054	±0.054 ppm	PASS
Dichlorvos	ND	0.00 ppm	0.009	0.028	±0.028 ppm	PASS
Diazinon	ND	0.20 ppm	0.001	0.002	±0.002 ppm	PASS
Dimethoate	ND	0.00 ppm	0.001	0.004	±0.004 ppm	PASS
Etoazazole	ND	1.50 ppm	0.002	0.007	±0.007 ppm	PASS
Fenoxycarb	ND	0.00 ppm	0.002	0.007	±0.007 ppm	PASS
Fenpyroximate	ND	2.00 ppm	0.001	0.002	±0.002 ppm	PASS
Fipronil	ND	0.00 ppm	0.005	0.015	±0.015 ppm	PASS
Fonicamid	ND	2.00 ppm	0.064	0.192	±0.192 ppm	PASS
Fludioxonil	ND	30.00 ppm	0.004	0.013	±0.013 ppm	PASS
Hexythiazox	ND	2.00 ppm	0.001	0.002	±0.002 ppm	PASS
Imazalil	ND	0.00 ppm	0.004	0.013	±0.013 ppm	PASS
Imidacloprid	ND	3.00 ppm	0.001	0.002	±0.002 ppm	PASS
Malathion	ND	5.00 ppm	0.003	0.010	±0.010 ppm	PASS
Metaxalyl	ND	15.00 ppm	0.005	0.015	±0.015 ppm	PASS
Methiocarb	ND	0.00 ppm	0.002	0.007	±0.007 ppm	PASS
Methomyl	ND	0.10 ppm	<0.001	0.001	±0.001 ppm	PASS
Methyl parathion	NT	0.00 ppm				NA
Mevinphos	ND	0.00 ppm	0.003	0.010	±0.010 ppm	PASS
Myclobutanil	ND	9.00 ppm	0.001	0.002	±0.002 ppm	PASS
Naled	ND	0.50 ppm	0.003	0.010	±0.010 ppm	PASS
Oxamyl	ND	0.20 ppm	0.001	0.004	±0.004 ppm	PASS
Paclobutrazol	ND	0.00 ppm	0.002	0.005	±0.005 ppm	PASS
Permethrin	ND	20.00 ppm	0.006	0.019	±0.019 ppm	PASS
Phosmet	ND	0.20 ppm	0.002	0.006	±0.006 ppm	PASS
Piperonylbutoxide	ND	8.00 ppm	0.007	0.020	±0.020 ppm	PASS
Prallethrin	ND	0.40 ppm	0.002	0.007	±0.007 ppm	PASS
Propiconazole	ND	20.00 ppm	0.002	0.007	±0.007 ppm	PASS
Propoxur	ND	0.00 ppm	0.004	0.011	±0.011 ppm	PASS
Pyrethrin	ND	1.00 ppm	0.002	0.005	±0.005 ppm	PASS
Pyridaben	ND	3.00 ppm	0.001	0.002	±0.002 ppm	PASS
Spinetoram	ND	3.00 ppm	0.002	0.007	±0.007 ppm	PASS
Spinosad	ND	3.00 ppm	0.004	0.013	±0.013 ppm	PASS
Spiromesifen	ND	12.00 ppm	0.002	0.006	±0.006 ppm	PASS
Spirotetramat	ND	13.00 ppm	0.002	0.005	±0.005 ppm	PASS
Spiroxamine	ND	0.00 ppm	0.001	0.002	±0.002 ppm	PASS
Tebuconazole	ND	2.00 ppm	0.003	0.010	±0.010 ppm	PASS
Thiacloprid	ND	0.10 ppm	0.001	0.002	±0.002 ppm	PASS
Thiamethoxam	ND	4.50 ppm	0.002	0.006	±0.006 ppm	PASS
Trifloxystrobin	ND	30.00 ppm	0.001	0.004	±0.004 ppm	PASS

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