



Juniper Analytics LLC

1334 NE 2nd St
Bend, OR
(541) 382-3796
<http://www.juniperanalyticsllc.com>
ORELAP: License # 10035537931

CBD WELLNESS

Tucson, AZ 85743
520-429-9618

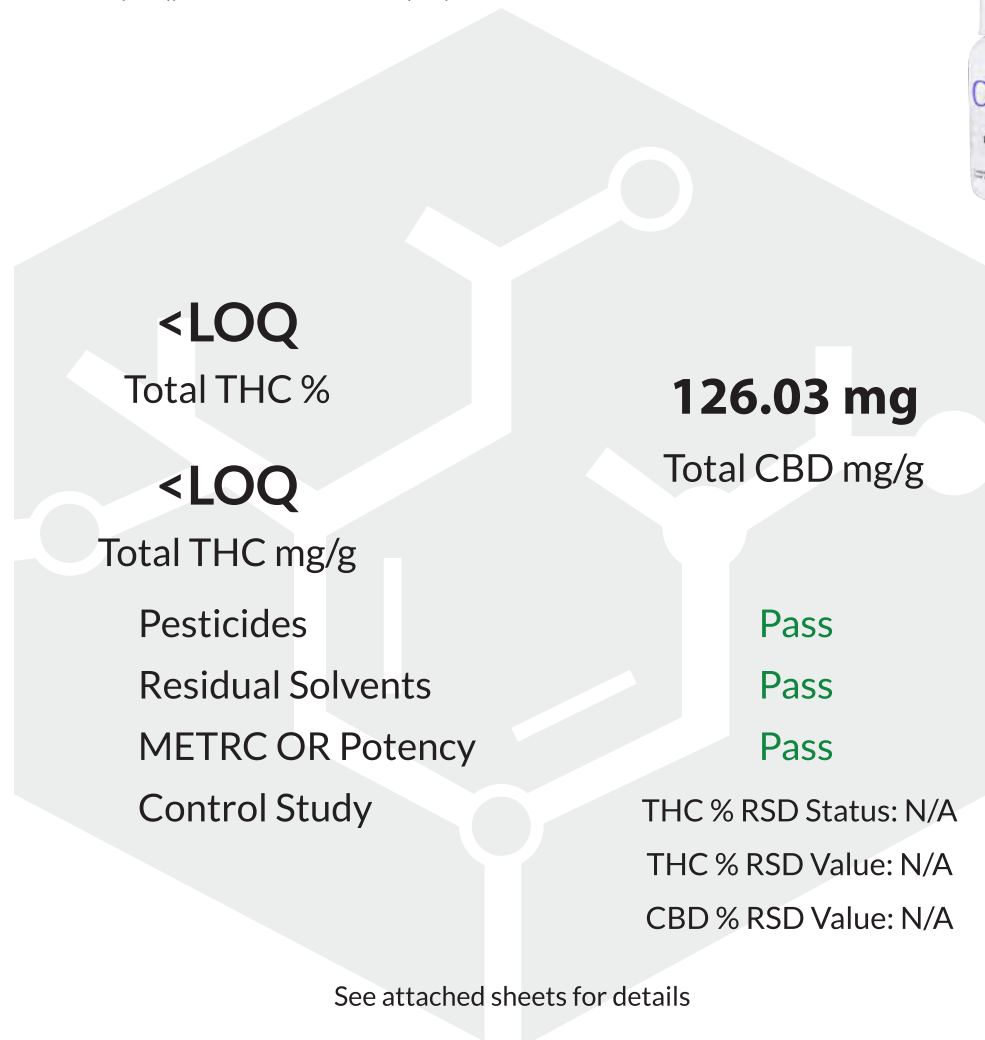
Strain: 125mg Original Liniment (28.35g)

Product Batch: 125OrigLini74

Total CBD: 126.03mg / 4.445mg/ml

License #: 00000077DCPS00216601 (MED), 00000067ESBS89254298 (REC)

Report Date: 07/06/2021
Intake Date: 06/21/2021
Harvest/Prod. Date: 06/14/2021



This product has been tested by Juniper Analytics LLC, using valid testing methodologies and a quality system as required by state law. Values reported relate only to the product tested and batched under the batch number identified above. Juniper Analytics LLC makes no claims as to the efficacy, safety, or other risks associated with any detected or non-detected level of any compounds reported herein. This Certificate must not be altered, and shall not be reproduced except in full, without the written approval of Juniper Analytics LLC.

Ben Armstrong
QA Review

Confident Cannabis
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Cannabinoid Profile

<LOQ

Total THC

126.03 mg

Total CBD mg/g

<LOQ

Total THC mg/g

| Analyte | LOQ mg/g | Mass % | Mass mg/g |
|----------------|-------------|---------------|---------------|
| THCa | 1.000 | <LOQ | <LOQ |
| Δ 9-THC | 1.000 | <LOQ | <LOQ |
| Δ 8-THC | 1.000 | <LOQ | <LOQ |
| THCV | 1.000 | <LOQ | <LOQ |
| CBDa | 1.000 | <LOQ | <LOQ |
| CBD | 1.000 | 0.4445 | 126.03 |
| CBN | 1.000 | <LOQ | <LOQ |
| CBGa | 1.000 | <LOQ | <LOQ |
| CBG | 1.000 | <LOQ | <LOQ |
| CBC | 1.000 | <LOQ | <LOQ |
| Total | | 0.4445 | 126.03 |

QC Batch ID: 125OrigLini74
Total THC = THCa * 0.877 + Δ 9-THC
Total CBD = CBDa * 0.877 + CBD

Instrument: HPLC/DAD; Method: JA-Potency-Proprietary; LOQ = Limit of Quantitation; The reported result is based on a sample weight with the applicable moisture content for that sample; Unless otherwise stated all quality control samples performed within specifications established by the Laboratory. Potency Analysis (Oregon Compliance Standard OAR 333-007-0430).

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Pass

| Analyte | LOQ | Limit | Results | Status | Analyte | LOQ | Limit | Results | Status |
|---------------------|------|-------|---------|--------|--------------------|------|-------|---------|--------|
| | PPM | PPM | PPM | | | PPM | PPM | PPM | |
| Abamectin | 0.25 | 0.50 | <LOQ | Pass | Imazalil | 0.10 | 0.20 | <LOQ | Pass |
| Acephate | 0.20 | 0.40 | <LOQ | Pass | Imidacloprid | 0.20 | 0.40 | <LOQ | Pass |
| Acequinocyl | 1.00 | 2.00 | <LOQ | Pass | Kresoxim Methyl | 0.20 | 0.40 | <LOQ | Pass |
| Acetamiprid | 0.10 | 0.20 | <LOQ | Pass | Malathion | 0.10 | 0.20 | <LOQ | Pass |
| Aldicarb | 0.20 | 0.40 | <LOQ | Pass | Metalaxyl | 0.10 | 0.20 | <LOQ | Pass |
| Azoxystrobin | 0.10 | 0.20 | <LOQ | Pass | Methiocarb | 0.10 | 0.20 | <LOQ | Pass |
| Bifenazate | 0.10 | 0.20 | <LOQ | Pass | Methomyl | 0.20 | 0.40 | <LOQ | Pass |
| Bifenthrin | 0.10 | 0.20 | <LOQ | Pass | Methyl Parathion | 0.10 | 0.20 | <LOQ | Pass |
| Boscalid | 0.20 | 0.40 | <LOQ | Pass | MGK-264 | 0.10 | 0.20 | <LOQ | Pass |
| Carbaryl | 0.10 | 0.20 | <LOQ | Pass | Myclobutanil | 0.10 | 0.20 | <LOQ | Pass |
| Carbofuran | 0.10 | 0.20 | <LOQ | Pass | Naled | 0.25 | 0.50 | <LOQ | Pass |
| Chlorantraniliprole | 0.10 | 0.20 | <LOQ | Pass | Oxamyl | 0.50 | 1.00 | <LOQ | Pass |
| Chlorfenapyr | 0.50 | 1.00 | <LOQ | Pass | Paclobotrazol | 0.20 | 0.40 | <LOQ | Pass |
| Chlorpyrifos | 0.10 | 0.20 | <LOQ | Pass | Permethrins | 0.10 | 0.20 | <LOQ | Pass |
| Clofentezine | 0.10 | 0.20 | <LOQ | Pass | Phosmet | 0.10 | 0.20 | <LOQ | Pass |
| Cyfluthrin | 0.50 | 1.00 | <LOQ | Pass | Piperonyl Butoxide | 1.00 | 2.00 | <LOQ | Pass |
| Cypermethrin | 0.50 | 1.00 | <LOQ | Pass | Prallethrin | 0.10 | 0.20 | <LOQ | Pass |
| Daminozide | 0.50 | 1.00 | <LOQ | Pass | Propiconazole | 0.20 | 0.40 | <LOQ | Pass |
| DDVP | 0.50 | 1.00 | <LOQ | Pass | Propoxur | 0.10 | 0.20 | <LOQ | Pass |
| Diazinon | 0.10 | 0.20 | <LOQ | Pass | Pyrethrins | 0.50 | 1.00 | <LOQ | Pass |
| Dimethoate | 0.10 | 0.20 | <LOQ | Pass | Pyridaben | 0.10 | 0.20 | <LOQ | Pass |
| Ethoprophos | 0.10 | 0.20 | <LOQ | Pass | Spinosad | 0.10 | 0.20 | <LOQ | Pass |
| Etofenprox | 0.20 | 0.40 | <LOQ | Pass | Spiromesifen | 0.10 | 0.20 | <LOQ | Pass |
| Etoxazole | 0.10 | 0.20 | <LOQ | Pass | Spirotetramat | 0.10 | 0.20 | <LOQ | Pass |
| Fenoxycarb | 0.10 | 0.20 | <LOQ | Pass | Spiroxamine | 0.20 | 0.40 | <LOQ | Pass |
| Fenpyroximate | 0.20 | 0.40 | <LOQ | Pass | Tebuconazole | 0.20 | 0.40 | <LOQ | Pass |
| Fipronil | 0.20 | 0.40 | <LOQ | Pass | Thiacloprid | 0.10 | 0.20 | <LOQ | Pass |
| Flonicamid | 0.50 | 1.00 | <LOQ | Pass | Thiamethoxam | 0.10 | 0.20 | <LOQ | Pass |
| Fludioxonil | 0.20 | 0.40 | <LOQ | Pass | Trifloxystrobin | 0.10 | 0.20 | <LOQ | Pass |
| Hexythiazox | 0.50 | 1.00 | <LOQ | Pass | | | | | |

QC Batch ID: 125OrigLini74

Instrument: LC-MS/MS; Method: AOAC 2007.1-Mod; LOQ = Limit of Quantitation; The reported result is based on wet sample weight for that sample; Unless otherwise stated all quality control samples performed within specifications established by the Laboratory.Pesticide Analysis (Oregon Compliance Standard OAR 333-008-1190).

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Pass

| Analyte | LOQ | Limit | Results | RPD | Analyte | LOQ | Limit | Results | RPD |
|-------------------|-----|---------|---------|-----|---------------------|------------|----------------|----------------|-----|
| | PPM | PPM | PPM | | | PPM | PPM | PPM | |
| 1,4-Dioxane | 100 | 380.00 | <LOQ | - | Toluene | 100 | 890.00 | <LOQ | - |
| 2-Butanol | 500 | 5000.00 | <LOQ | - | Pentanes | 500 | 5000.00 | <LOQ | - |
| 2-Ethoxy-Ethanol | 100 | 160.00 | <LOQ | - | n-Pentane | | | <LOQ | - |
| 2-Propanol IPA | 500 | 5000.00 | <LOQ | - | Isopentane | | | <LOQ | - |
| Acetone | 500 | 5000.00 | <LOQ | - | Neopentane | | | <LOQ | - |
| Acetonitrile | 100 | 410.00 | <LOQ | - | Butanes | 500 | 5000.00 | <LOQ | - |
| Benzene | 1 | 2.00 | <LOQ | - | n-Butane | | | <LOQ | - |
| Cumene | 50 | 70.00 | <LOQ | - | Isobutane | | | <LOQ | - |
| Cyclohexane | 500 | 3880.00 | <LOQ | - | Hexanes | 50 | 290.00 | <LOQ | - |
| Dichloromethane | 100 | 600.00 | <LOQ | - | n-Hexane | | | <LOQ | - |
| Ethyl-Acetate | 500 | 5000.00 | <LOQ | - | 2-Methyl-Pentane | | | <LOQ | - |
| Ethyl-Ether | 500 | 5000.00 | <LOQ | - | 3-Methyl-Pentane | | | <LOQ | - |
| Ethylene Glycol | 300 | 620.00 | <LOQ | - | 2,2-Dimethyl-Butane | | | <LOQ | - |
| Ethylene Oxide | 20 | 50.00 | <LOQ | - | 2,3-Dimethyl-Butane | | | <LOQ | - |
| Heptane | 500 | 5000.00 | <LOQ | - | Xylenes | 300 | 2170.00 | <LOQ | - |
| Isopropyl-Acetate | 500 | 5000.00 | <LOQ | - | m-Xylene | | | <LOQ | - |
| Methanol | 500 | 3000.00 | <LOQ | - | o-Xylene | | | <LOQ | - |
| Propane | 500 | 5000.00 | <LOQ | - | p-Xylene | | | <LOQ | - |
| Tetrahydrofuran | 100 | 720.00 | <LOQ | - | Ethyl-Benzene | | | <LOQ | - |

METRC RPD Status: N/A

Tentatively Identified Compounds: Peak 1: Hits 1-3: 2-methylhexane; Peak 2: Hit 1: 3-methylhexane; Peak 3: Hit 1: 2-propanone, 1-(acetyloxy)-

QC Batch ID: 125OrigLini74

Instrument: GC/MS; Method: USP 467 - Modified; LOQ = Limit of Quantitation; The reported result is based on wet sample weight for that sample; Unless otherwise stated all quality control samples performed within specifications established by the Laboratory. Residual Solvent Analysis (Oregon Compliance Standard OAR 333-007-0410).

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