

19173A1

Anandapure, LLC
 350 Greenwood Avenue
 Reedsport, OR 97467
 302-981-1558

Sample Type: Other
 Sample Date: 6/25/2019
 Analysis Date: 6/26/2019
 Report Date: 6/28/2019

Metric Batch ID:
 Metric Sample ID:

Harvest/Process Date:
 Report ID:
LS-190628-34

Potency



Potency Analysis Date: 6/26/2019
 Potency Batch ID: CAN_062619B
 Potency Method: JAOAC 2015.1

<LOQ Total THC
19.8 mg/g Total CBD

Samples: NBM-ZSF-ZZN

| Analyte | Description | LOQ | RPD (%) | Min. | Max. | Conc. | Unit: mg/g |
|------------------|-------------------------------|------|---------|------|------|-------|------------|
| Δ9THC | Delta-9 Tetrahydrocannabinol | 0.80 | - | - | - | ND | |
| THCA | Tetrahydrocannabinolic acid | 0.80 | - | - | - | <LOQ | |
| CBD | Cannabidiol | 0.80 | - | - | - | 4.72 | |
| CBDA | Cannabidiolic acid | 0.80 | - | - | - | 17.2 | |
| Δ8THC | Delta-8 Tetrahydrocannabinol* | 0.80 | - | - | - | ND | |
| THCV | Tetrahydrocannabivarin* | 0.80 | - | - | - | ND | |
| CBG | Cannabigerol* | 0.80 | - | - | - | ND | |
| CBGA | Cannabigerolic acid* | 0.80 | - | - | - | ND | |
| CBC | Cannabichromene* | 0.80 | - | - | - | <LOQ | |
| CBCA | Cannabichromenic acid* | 0.80 | - | - | - | <LOQ | |
| CBN | Cannabinol | 0.80 | - | - | - | ND | |
| Total THC | Δ9THC + (THCA × 0.877) | | - | - | - | <LOQ | |
| Total CBD | CBD + (CBDA × 0.877) | | - | - | - | 19.8 | |
| Total | | | - | - | - | 21.9 | |

Compliance

| | | | |
|------------|---------------|--------------------------|--|
| Pesticides | Within limits | Analysis Date: 6/27/2019 | Pass  |
| Solvents | Within limits | Analysis Date: 6/26/2019 | Pass  |


 Ian Eustis
 Lab Director


 Aaron Troyer
 Chief Science Officer

This data cannot be used for OLCC or OHA compliance for usable marijuana or marijuana products and is provided for Research and Development purposes only.



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Pesticides Sample Data

Pesticides Analysis Date: 6/27/2019
 Pesticides Batch ID: PST_062719A

Method: EN 15662
 Unit: µg/g (ppm)

Pass 

| Analyte | NBM-ZSF-ZZN | Limits | LOQ | Notes | Status | Analyte | NBM-ZSF-ZZN | Limits | LOQ | Notes | Status |
|---------------------|-------------|--------|-----|-------|--------|--------------------|-------------|--------|-----|-------|--------|
| Abamectin | <LOQ | 0.5 | 0.1 | - | Pass | Metalaxyl | <LOQ | 0.2 | 0.1 | - | Pass |
| Acephate | <LOQ | 0.4 | 0.1 | - | Pass | Methiocarb | <LOQ | 0.2 | 0.1 | - | Pass |
| Acequinocyl | <LOQ | 2.0 | 1.5 | - | Pass | Methomyl | <LOQ | 0.4 | 0.1 | - | Pass |
| Acetamiprid | <LOQ | 0.2 | 0.1 | - | Pass | Methyl Parathion | <LOQ | 0.2 | 0.2 | - | Pass |
| Aldicarb | <LOQ | 0.4 | 0.1 | - | Pass | MGK-264 | <LOQ | 0.2 | 0.2 | - | Pass |
| Azoxystrobin | <LOQ | 0.2 | 0.1 | - | Pass | Myclobutanil | <LOQ | 0.2 | 0.1 | - | Pass |
| Bifenazate | <LOQ | 0.2 | 0.1 | - | Pass | Naled | <LOQ | 0.5 | 0.2 | - | Pass |
| Bifenthrin | <LOQ | 0.2 | 0.1 | - | Pass | Oxamyl | <LOQ | 1.0 | 0.1 | - | Pass |
| Boscalid | <LOQ | 0.4 | 0.1 | - | Pass | Paclobutrazol | <LOQ | 0.4 | 0.1 | - | Pass |
| Carbaryl | <LOQ | 0.2 | 0.1 | - | Pass | Permethrins | <LOQ | 0.2 | 0.1 | - | Pass |
| Carbofuran | <LOQ | 0.2 | 0.1 | - | Pass | Phosmet | <LOQ | 0.2 | 0.1 | - | Pass |
| Chlorantraniliprole | <LOQ | 0.2 | 0.1 | - | Pass | Piperonyl Butoxide | <LOQ | 2.0 | 0.1 | - | Pass |
| Chlorfenapyr | <LOQ | 1.0 | 0.1 | - | Pass | Prallethrin | <LOQ | 0.2 | 0.1 | - | Pass |
| Chlorpyrifos | <LOQ | 0.2 | 0.1 | - | Pass | Propiconazole | <LOQ | 0.4 | 0.1 | - | Pass |
| Clofentezine | <LOQ | 0.2 | 0.1 | - | Pass | Propoxur | <LOQ | 0.2 | 0.1 | - | Pass |
| Cyfluthrin | <LOQ | 1.0 | 0.5 | - | Pass | Pyrethrins | <LOQ | 1.0 | 0.5 | - | Pass |
| Cypermethrin | <LOQ | 1.0 | 0.1 | - | Pass | Pyridaben | <LOQ | 0.2 | 0.1 | - | Pass |
| Daminozide | <LOQ | 1.0 | 0.5 | - | Pass | Spinosad | <LOQ | 0.2 | 0.1 | - | Pass |
| Diazinon | <LOQ | 0.2 | 0.1 | - | Pass | Spiromesifen | <LOQ | 0.2 | 0.1 | - | Pass |
| Dichlorvos (DDVP) | <LOQ | 1.0 | 0.5 | - | Pass | Spirotetramat | <LOQ | 0.2 | 0.1 | - | Pass |
| Dimethoate | <LOQ | 0.2 | 0.1 | - | Pass | Spiroxamine | <LOQ | 0.4 | 0.1 | - | Pass |
| Ethoprophos | <LOQ | 0.2 | 0.1 | - | Pass | Tebuconazole | <LOQ | 0.4 | 0.1 | - | Pass |
| Etofenprox | <LOQ | 0.4 | 0.1 | - | Pass | Thiacloprid | <LOQ | 0.2 | 0.1 | - | Pass |
| Etoxazole | <LOQ | 0.2 | 0.1 | - | Pass | Thiamethoxam | <LOQ | 0.2 | 0.1 | - | Pass |
| Fenoxycarb | <LOQ | 0.2 | 0.1 | - | Pass | Trifloxystrobin | <LOQ | 0.2 | 0.1 | - | Pass |
| Fenpyroximate | <LOQ | 0.4 | 0.1 | - | Pass | | | | | | |
| Fipronil | <LOQ | 0.4 | 0.1 | - | Pass | | | | | | |
| Flonicamid | <LOQ | 1.0 | 0.1 | - | Pass | | | | | | |
| Fludioxonil | <LOQ | 0.4 | 0.1 | - | Pass | | | | | | |
| Hexythiazox | <LOQ | 1.0 | 0.1 | - | Pass | | | | | | |
| Imazalil | <LOQ | 0.2 | 0.1 | - | Pass | | | | | | |
| Imidacloprid | <LOQ | 0.4 | 0.1 | - | Pass | | | | | | |
| Kresoxim-methyl | <LOQ | 0.4 | 0.1 | - | Pass | | | | | | |
| Malathion | <LOQ | 0.2 | 0.1 | - | Pass | | | | | | |

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 Report Date: 6/28/2019

Metric Batch ID:
 Metric Sample ID:

Harvest/Process Date:
 Report ID:
LS-190628-34



Pesticides Quality Control Data

Pesticides QC Analysis Date: 6/27/2019
 Pesticides QC Batch ID: PST_062719A

Method: EN 15662
 Unit: µg/g (ppm)

Laboratory Pesticides Quality Control Results

| Method: EN 15662 | | | | Units: ppm (µg/g) | | | | Batch ID: PST_062719A | | | | | | | | | |
|---------------------|--------------|-----|-------|-------------------|-----------|----------|----------|-----------------------|--------------------|--------------|-----|-------|------------|-----------|----------|----------|-------|
| Pesticide | Blank Result | LOQ | Notes | LCS Result | LCS Spike | LCS% Rec | Limits | Notes | Pesticide | Blank Result | LOQ | Notes | LCS Result | LCS Spike | LCS% Rec | Limits | Notes |
| Abamectin | nd | 0.1 | | 0.9 | 1.0 | 87 | 50 - 150 | | Imazalil | nd | 0.1 | | 0.9 | 1.0 | 89 | 50 - 150 | |
| Acephate | nd | 0.1 | | 1.1 | 1.0 | 105 | 50 - 150 | | Imidacloprid | nd | 0.1 | | 1.1 | 1.0 | 110 | 50 - 150 | |
| Acequinocyl | nd | 1.0 | | 0.5 | 1.0 | 54 | 50 - 150 | | Kresoxim-methyl | nd | 0.1 | | 1.2 | 1.0 | 115 | 50 - 150 | |
| Acetamiprid | nd | 0.1 | | 1.2 | 1.0 | 115 | 50 - 150 | | Malathion | nd | 0.1 | | 1.2 | 1.0 | 124 | 50 - 150 | |
| Aldicarb | nd | 0.1 | | 1.0 | 1.0 | 101 | 50 - 150 | | Metaxyl | nd | 0.1 | | 1.1 | 1.0 | 110 | 50 - 150 | |
| Azoxystrobin | nd | 0.1 | | 1.1 | 1.0 | 114 | 50 - 150 | | Methiocarb | nd | 0.1 | | 1.1 | 1.0 | 108 | 50 - 150 | |
| Bifenthrin | nd | 0.1 | | 1.0 | 1.0 | 95 | 50 - 150 | | Methomyl | nd | 0.1 | | 1.1 | 1.0 | 105 | 50 - 150 | |
| Bifenazate | nd | 0.1 | | 1.3 | 1.0 | 129 | 50 - 150 | | Methyl Parathion | nd | 0.1 | | 0.5 | 1.0 | 53 | 30 - 150 | |
| Boscalid | nd | 0.1 | | 1.2 | 1.0 | 117 | 50 - 150 | | MGK-264 | nd | 0.2 | | 0.6 | 0.6 | 107 | 50 - 150 | |
| Carbaryl | nd | 0.1 | | 1.1 | 1.0 | 112 | 50 - 150 | | Myclobutanil | nd | 0.1 | | 1.1 | 1.0 | 113 | 50 - 150 | |
| Carbofuran | nd | 0.1 | | 1.0 | 1.0 | 104 | 50 - 150 | | Naled | nd | 0.1 | | 0.9 | 1.0 | 95 | 50 - 150 | |
| Chlorantraniliprole | nd | 0.1 | | 1.3 | 1.0 | 128 | 50 - 150 | | Oxamyl | nd | 0.1 | | 1.1 | 1.0 | 107 | 50 - 150 | |
| Chlorfenapyr | nd | 0.1 | | 1.0 | 1.0 | 95 | 50 - 150 | | Paclobutrazol | nd | 0.1 | | 0.9 | 1.0 | 94 | 50 - 150 | |
| Chlorpyrifos | nd | 0.1 | | 1.0 | 1.0 | 99 | 50 - 150 | | Permethrin | nd | 0.1 | | 1.0 | 1.0 | 99 | 50 - 150 | |
| Clofentezine | nd | 0.1 | | 0.9 | 1.0 | 91 | 50 - 150 | | Phosmet | nd | 0.1 | | 1.1 | 1.0 | 112 | 50 - 150 | |
| Cyfluthrin | nd | 0.5 | | 1.3 | 1.0 | 126 | 50 - 150 | | Piperonyl Butoxide | nd | 0.1 | | 0.9 | 1.0 | 90 | 50 - 150 | |
| Cypermethrin | nd | 0.1 | | 0.9 | 1.0 | 91 | 50 - 150 | | Prallethrin | nd | 0.1 | | 1.1 | 1.0 | 109 | 50 - 150 | |
| Daminozide | nd | 0.5 | | nd | 1.0 | | 10 - 150 | LR | Propiconazole | nd | 0.1 | | 1.2 | 1.0 | 125 | 50 - 150 | |
| Diazinon | nd | 0.1 | | 1.0 | 1.0 | 104 | 50 - 150 | | Propoxur | nd | 0.1 | | 1.0 | 1.0 | 105 | 50 - 150 | |
| Dichlorvos | nd | 0.5 | | 1.1 | 1.0 | 113 | 50 - 150 | | Pyrethrins | nd | 0.2 | | 0.9 | 1.0 | 95 | 50 - 150 | |
| Dimethoate | nd | 0.1 | | 1.1 | 1.0 | 105 | 50 - 150 | | Pyridaben | nd | 0.1 | | 0.9 | 1.0 | 91 | 50 - 150 | |
| Ethoprophos | nd | 0.1 | | 1.1 | 1.0 | 106 | 50 - 150 | | Spinosad A kps | nd | 0.1 | | 0.7 | 1.0 | 67 | 50 - 150 | |
| Etofenprox | nd | 0.1 | | 0.9 | 1.0 | 93 | 50 - 150 | | Spinosad D kps | nd | 0.1 | | 0.1 | 0.1 | 70 | 50 - 150 | |
| Etoxazole | nd | 0.1 | | 0.8 | 1.0 | 85 | 50 - 150 | | Spiromesifen | nd | 0.1 | | 1.0 | 1.0 | 97 | 50 - 150 | |
| Fenoxycarb | nd | 0.1 | | 1.1 | 1.0 | 114 | 50 - 150 | | Spirotetramat | nd | 0.1 | | 1.1 | 1.0 | 110 | 50 - 150 | |
| Fenpyroximate | nd | 0.1 | | 0.9 | 1.0 | 89 | 50 - 150 | | Spiroxamine | nd | 0.1 | | 0.7 | 1.0 | 71 | 50 - 150 | |
| Fipronil | nd | 0.1 | | 1.3 | 1.0 | 128 | 50 - 150 | | Tebuconazole | nd | 0.1 | | 1.3 | 1.0 | 128 | 50 - 150 | |
| Flonicamid | nd | 0.1 | | 1.0 | 1.0 | 98 | 50 - 150 | | Thiacloprid | nd | 0.1 | | 1.2 | 1.0 | 119 | 50 - 150 | |
| Fludioxonil | nd | 0.1 | | 1.0 | 1.0 | 103 | 50 - 150 | | Thiamethoxam | nd | 0.1 | | 1.3 | 1.0 | 132 | 50 - 150 | |
| Hexythiazox | nd | 0.1 | | 0.8 | 1.0 | 80 | 50 - 150 | | Trifloxystrobin | nd | 0.1 | | 1.0 | 1.0 | 101 | 50 - 150 | |

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Harvest/Process Date:
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Residual Solvents Sample Data

Solvents Analysis Date: 6/26/2019
 Solvents Batch ID: RES_062619A

Method: EPA 5021A
 Unit: µg/g (ppm)

Pass 

| Analyte | NBM-ZSF-ZZN | RPD (%) | Limits | LOQ | Notes | Status |
|--------------------------|-------------|---------|--------|-------|-------|--------|
| 1,4-Dioxane | <LOQ | 0.00 | 380.0 | 50.0 | - | Pass |
| 2-Butanol | <LOQ | 0.00 | 5000.0 | 250.0 | - | Pass |
| 2-Ethoxyethanol | <LOQ | 0.00 | 160.0 | 50.0 | - | Pass |
| Acetone | <LOQ | 0.00 | 5000.0 | 250.0 | - | Pass |
| Acetonitrile | <LOQ | 0.00 | 410.0 | 50.0 | - | Pass |
| Benzene | <LOQ | 0.00 | 2.0 | 2.0 | - | Pass |
| Butanes | <LOQ | 0.00 | 5000.0 | 250.0 | - | Pass |
| Cumene | <LOQ | 0.00 | 70.0 | 50.0 | - | Pass |
| Cyclohexane | <LOQ | 0.00 | 3880.0 | 50.0 | - | Pass |
| Ethyl Acetate | <LOQ | 0.00 | 5000.0 | 250.0 | - | Pass |
| Ethyl Ether | <LOQ | 0.00 | 5000.0 | 250.0 | - | Pass |
| Ethylene Glycol | <LOQ | 0.00 | 620.0 | 250.0 | - | Pass |
| Ethylene Oxide | <LOQ | 0.00 | 50.0 | 50.0 | - | Pass |
| Heptane | <LOQ | 0.00 | 5000.0 | 250.0 | - | Pass |
| Hexanes | <LOQ | 0.00 | 290.0 | 50.0 | - | Pass |
| Isopropanol (2-Propanol) | <LOQ | 0.00 | 5000.0 | 50.0 | - | Pass |
| Isopropyl Acetate | <LOQ | 0.00 | 5000.0 | 250.0 | - | Pass |
| Methanol | <LOQ | 0.00 | 3000.0 | 250.0 | - | Pass |
| Dichloromethane | <LOQ | 0.00 | 600.0 | 50.0 | - | Pass |
| Pentanes | <LOQ | 0.00 | 5000.0 | 250.0 | - | Pass |
| Propane | <LOQ | 0.00 | 5000.0 | 250.0 | - | Pass |
| Tetrahydrofuran | <LOQ | 0.00 | 720.0 | 50.0 | - | Pass |
| Toluene | <LOQ | 0.00 | 890.0 | 50.0 | - | Pass |
| Xylenes | <LOQ | 0.00 | 2170.0 | 50.0 | - | Pass |

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Residual Solvents Quality Control Data

Solvents QC Analysis Date: 6/26/2019
 Solvents QC Batch ID: RES_062619A

Method: EPA 5021A
 Unit: µg/g (ppm)

Laboratory Residual Solvent Quality Control Results

Method: EPA 5021A

Units: µg/mL

Batch ID: RES_062619A

Matrix Blank / LCS Results

| Analyte | Blank Result | Blank Limit | Notes | LCS Result | LCS Spike | LCS% Rec | Limits | Notes |
|---------------------------|--------------|-------------|-------|------------|-----------|----------|----------|-------|
| 1,4-Dioxane | < LOQ | 50 | | 871 | 1000 | 87 | 70 - 130 | |
| 2-Butanol | < LOQ | 50 | | 839 | 1000 | 84 | 70 - 130 | |
| 2-Ethoxyethanol | < LOQ | 50 | | 882 | 1000 | 88 | 70 - 130 | |
| Acetone | < LOQ | 50 | | 884 | 1000 | 88 | 70 - 130 | |
| Acetonitrile | < LOQ | 50 | | 834 | 1000 | 83 | 70 - 130 | |
| Benzene | < LOQ | 2 | | 15 | 20 | 75 | 70 - 130 | |
| Butanes | | | | | | | | |
| <i>Butane</i> | < LOQ | 50 | | 921 | 1000 | 92 | 70 - 130 | |
| <i>Isobutane</i> | < LOQ | 50 | | 938 | 1000 | 94 | 70 - 130 | |
| Cyclohexane | < LOQ | 50 | | 881 | 1000 | 88 | 70 - 130 | |
| Ethyl acetate | < LOQ | 50 | | 850 | 1000 | 85 | 70 - 130 | |
| Ethyl ether | < LOQ | 50 | | 902 | 1000 | 90 | 70 - 130 | |
| Ethylbenzene | < LOQ | 50 | | 843 | 1000 | 84 | 70 - 130 | |
| Ethylene glycol | < LOQ | 250 | | 877 | 1000 | 88 | 70 - 130 | |
| Ethylene oxide | < LOQ | 50 | | 856 | 1000 | 86 | 70 - 130 | |
| Heptane | < LOQ | 50 | | 832 | 1000 | 83 | 70 - 130 | |
| Hexanes | | | | | | | | |
| <i>n-Hexane</i> | < LOQ | 50 | | 894 | 1000 | 89 | 70 - 130 | |
| <i>2-Methylpentane</i> | < LOQ | 50 | | 851 | 1000 | 85 | 70 - 130 | |
| <i>3-Methylpentane</i> | < LOQ | 50 | | 893 | 1000 | 89 | 70 - 130 | |
| <i>2,2-Dimethylbutane</i> | < LOQ | 50 | | 874 | 1000 | 87 | 70 - 130 | |
| <i>2,3-Dimethylbutane</i> | < LOQ | 50 | | 811 | 1000 | 81 | 70 - 130 | |
| Isopropanol | < LOQ | 50 | | 847 | 1000 | 85 | 70 - 130 | |
| Isopropyl acetate | < LOQ | 50 | | 836 | 1000 | 84 | 70 - 130 | |
| Cumene | < LOQ | 50 | | 828 | 1000 | 83 | 70 - 130 | |
| Methanol | < LOQ | 50 | | 893 | 1000 | 89 | 70 - 130 | |
| Dichloromethane | < LOQ | 50 | | 844 | 1000 | 84 | 70 - 130 | |
| Pentanes | | | | | | | | |
| <i>Pentane</i> | < LOQ | 50 | | 861 | 1000 | 86 | 70 - 130 | |
| <i>Isopentane</i> | < LOQ | 50 | | 883 | 1000 | 88 | 70 - 130 | |
| <i>Neopentane</i> | < LOQ | 50 | | 960 | 1000 | 96 | 70 - 130 | |
| Propane | < LOQ | 50 | | 983 | 1000 | 98 | 70 - 130 | |
| Tetrahydrofuran | < LOQ | 50 | | 830 | 1000 | 83 | 70 - 130 | |
| Toluene | < LOQ | 50 | | 857 | 1000 | 86 | 70 - 130 | |
| Xylenes | | | | | | | | |
| <i>m-Xylene</i> | < LOQ | 50 | | 824 | 1000 | 82 | 70 - 130 | |
| <i>o/p-Xylene</i> | < LOQ | 50 | | 1675 | 2000 | 84 | 70 - 130 | |

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Metric Sample ID:Harvest/Process Date:
Report ID:
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Qualifier Flag Descriptions

| | |
|------------|--|
| J | Reported result is an estimate - the value is less than the minimum calibration level but greater than the estimated detection limit (EDL) |
| U | The analyte was not detected in the sample at the estimated detection limit (EDL) |
| E | Exceeds calibration range |
| D | Dilution data - result was obtained from the analysis of a dilution |
| B | Analyte found in sample and associated blank |
| C | Co-eluting compound |
| R | Relative Percent Difference (RPD) outside control limits |
| NR | Analyte not reported because of problems in sample preparation or analysis |
| ND | Non-Detect |
| X | Results from reinjection/repeat/re-column data |
| EMC | Estimated maximum possible concentration - indicates that a peak is detected but did not meet the method required criteria |
| M | Manual integration |
| PS | Peaks split |
| HB | Control acceptance criteria are exceeded high and the associated sample is below the detection limit |
| LB | Control acceptance criteria are exceeded low and the associated sample exceeds the regulatory limit |
| ME | Marginal Exceedance |
| LR | Low Recovery Analyte |
| LOQ | Limit of Quantitation |