

484 - Deathstar RCO 8.13.20

Farmer's Friend Extracts
 6451 NE Colwood Wy
 Portland, OR 97218
 503-442-8653

Sample Type: Extracts
Sample Date: 8/17/2020
Analysis Date: 8/18/2020
Report Date: 8/19/2020

Metric Batch ID:
 1A4010300016315000029962
Metric Sample ID:
 1A4010300016315000029985

Harvest/Process Date: 8/13/2020
Report ID: LS-200819-17
Sample Plan ID: SP-200817-5-A
Sample Procedure: 160721_LAB-SOP_SampleCollection-v008

Potency

Potency Analysis Date: 8/18/2020
 Potency Batch ID: CAN_081820C
 Potency Method: JAOAC 2015.1



68.2% Total THC
0.0854% Total CBD



Samples: CPW-GZN-XND, BXS-DDX-RNP

| Analyte | Description | LOQ | RPD | Min. | Max. | Avg. | Unit: % |
|------------------|-------------------------------|------|-------|-------|-------|----------------|---------|
| Δ9THC | Delta-9 Tetrahydrocannabinol | 0.16 | 2.89 | 67.2 | 69.2 | 68.2 | |
| THCA | Tetrahydrocannabinolic acid | 0.16 | 0.00 | ND | <LOQ | <LOQ | |
| CBD | Cannabidiol | 0.16 | 6.47 | ND | 0.171 | 0.0854 | |
| CBDA | Cannabidiolic acid | 0.16 | 0.00 | ND | <LOQ | <LOQ | |
| Δ8THC | Delta-8 Tetrahydrocannabinol* | 0.16 | 0.00 | ND | ND | ND | |
| THCV | Tetrahydrocannabivarin* | 0.16 | 3.22 | 3.81 | 3.93 | 3.87 | |
| CBG | Cannabigerol* | 0.16 | 7.44 | 0.954 | 1.03 | 0.991 | |
| CBGA | Cannabigerolic acid* | 0.16 | 0.00 | ND | <LOQ | <LOQ | |
| CBC | Cannabichromene* | 0.16 | 0.125 | 2.15 | 2.16 | 2.16 | |
| CBCA | Cannabichromenic acid* | 0.16 | 0.00 | ND | ND | ND | |
| CBN | Cannabinol | 0.16 | 42.6 | 0.356 | 0.549 | 0.452 | |
| Total THC | Δ9THC + (THCA × 0.877) | | 2.89 | 67.2 | 69.2 | 68.2 | |
| Total CBD | CBD + (CBDA × 0.877) | | 6.47 | ND | 0.171 | 0.0854 | |
| Total | | | 2.39 | 74.5 | 77.1 | 75.8 | |

Compliance

| | | | |
|-------------------|---------------|--------------------------|--|
| Pesticides | Within limits | Analysis Date: 8/18/2020 | Pass  |
| Solvents | Within limits | Analysis Date: 8/19/2020 | Pass  |
| Potency | Within limits | Analysis Date: 8/18/2020 | Pass  |


 Bryce Kidd, Ph.D.
 Lab Director


 Aaron Troyer
 Chief Science Officer



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Terpene Analysis Date: 8/18/2020
Terpene Batch ID: TRP_081820A

Method: JAOAC 2015.1
Unit: %

| Analyte | Avg. | Notes |
|---------------------|----------|-------|
| β-Caryophyllene | 3.88% | |
| Humulene | 1.54% | |
| α-Bisabolol | 0.343% | |
| β-Myrcene | 0.328% | |
| Limonene | 0.317% | |
| α-Terpineol | 0.180% | |
| Linalool | 0.173% | |
| Selinadiene | 0.159% | |
| Fenchol | 0.153% | |
| β-Pinene | 0.0847% | |
| Sabinene | 0.0755% | |
| β-Ocimene | 0.0582% | |
| α-Pinene | 0.0525% | |
| Caryophyllene Oxide | 0.0488% | |
| α-Cedrene | 0.0346% | |
| Eucalyptol | 0.0331% | |
| Terpinolene | 0.0211% | |
| Borneol | 0.0171% | |
| α-Terpinene | 0.0100% | |
| γ-Terpinene | 0.00912% | |
| α-Ocimene | 0.00841% | |
| Camphene | 0.00737% | |
| Fenchone | 0.00651% | |
| Δ3-Carene | 0.00621% | |
| Azulene | ND | |
| Camphore | ND | |
| Cedrol | ND | |
| Cymene | ND | |
| Geraniol | ND | |

| Analyte | Avg. | Notes |
|------------------|--------------|-------|
| Geranyl Acetate | ND | |
| Guaiol | ND | |
| Isoborneol | ND | |
| Isopulegol | ND | |
| Nerol | ND | |
| Pulegone | ND | |
| Sabinene Hydrate | ND | |
| Valencene | ND | |
| cis-Nerolidol | ND | |
| trans-Nerolidol | ND | |
| α-Phellandrene | ND | |
| β-Farnesene 1 | ND | |
| β-Farnesene 2 | ND | |
| γ-Terpineol | ND | |
| Total | 7.55% | |

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Pesticides Analysis Date: 8/18/2020
 Pesticides Batch ID: PST_081820B

Method: EN 15662
 Unit: µg/g (ppm)

Pass 

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

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Sample Plan ID: SP-200817-5-A
Sample Procedure: 160721_LAB-SOP_SampleCollection-v008

Pesticides Quality Control Data

Pesticides QC Analysis Date: 8/18/2020
 Pesticides QC Batch ID: PST_081820B

Method: EN 15662
 Unit: µg/g (ppm)

| Analyte | Blank | LOQ | LCS | LCS Spike | LCS Rec (%) | Limits (%) | Notes | Analyte | Blank | LOQ | LCS | LCS Spike | LCS Rec (%) | Limits (%) | Notes |
|---------------------|-------|-----|-------|-----------|-------------|------------|-------|--------------------|-------|-----|-------|-----------|-------------|------------|-------|
| Abamectin | ND | 0.1 | 0.777 | 1.00 | 77.7 | 50 - 150 | | Metalaxyl | ND | 0.1 | 0.945 | 1.00 | 94.5 | 50 - 150 | |
| Acephate | ND | 0.1 | 0.866 | 1.00 | 86.6 | 50 - 150 | | Methiocarb | ND | 0.1 | 0.954 | 1.00 | 95.4 | 50 - 150 | |
| Acequinocyl | ND | 1.5 | 1.02 | 1.00 | 102 | 50 - 150 | | Methomyl | ND | 0.1 | 0.935 | 1.00 | 93.5 | 50 - 150 | |
| Acetamiprid | ND | 0.1 | 0.883 | 1.00 | 88.3 | 50 - 150 | | Methyl Parathion | ND | 0.2 | 3.79 | 1.00 | 379 | 30 - 150 | HB |
| Aldicarb | ND | 0.1 | 0.940 | 1.00 | 94.0 | 50 - 150 | | MGK-264 | ND | 0.2 | 0.179 | 0.600 | 29.9 | 50 - 150 | LR |
| Azoxystrobin | ND | 0.1 | 0.920 | 1.00 | 92.0 | 50 - 150 | | Myclobutanil | ND | 0.1 | 0.905 | 1.00 | 90.5 | 50 - 150 | |
| Bifenazate | ND | 0.1 | 1.11 | 1.00 | 111 | 50 - 150 | | Naled | ND | 0.2 | 0.969 | 1.00 | 96.9 | 50 - 150 | |
| Bifenthrin | ND | 0.1 | 1.07 | 1.00 | 107 | 50 - 150 | | Oxamyl | ND | 0.1 | 0.975 | 1.00 | 97.5 | 50 - 150 | |
| Boscalid | ND | 0.1 | 0.861 | 1.00 | 86.1 | 50 - 150 | | Paclobutrazol | ND | 0.1 | 0.881 | 1.00 | 88.1 | 50 - 150 | |
| Carbaryl | ND | 0.1 | 0.954 | 1.00 | 95.4 | 50 - 150 | | Permethrins | ND | 0.1 | 1.36 | 1.00 | 136 | 50 - 150 | |
| Carbofuran | ND | 0.1 | 0.890 | 1.00 | 89.0 | 50 - 150 | | Phosmet | ND | 0.1 | 0.881 | 1.00 | 88.1 | 50 - 150 | |
| Chlorantraniliprole | ND | 0.1 | 0.876 | 1.00 | 87.6 | 50 - 150 | | Piperonyl Butoxide | ND | 0.1 | 1.01 | 1.00 | 101 | 50 - 150 | |
| Chlorfenapyr | ND | 0.1 | 0.952 | 1.00 | 95.2 | 50 - 150 | | Prallethrin | ND | 0.1 | 1.08 | 1.00 | 108 | 50 - 150 | |
| Chlorpyrifos | ND | 0.1 | 1.01 | 1.00 | 101 | 50 - 150 | | Propiconazole | ND | 0.1 | 1.19 | 1.00 | 119 | 50 - 150 | |
| Clofentezine | ND | 0.1 | 0.747 | 1.00 | 74.7 | 50 - 150 | | Propoxur | ND | 0.1 | 0.948 | 1.00 | 94.8 | 50 - 150 | |
| Cyfluthrin | ND | 0.5 | 0.902 | 1.00 | 90.2 | 50 - 150 | | Pyrethrins | ND | 0.5 | 0.794 | 1.00 | 79.4 | 50 - 150 | |
| Cypermethrin | ND | 0.1 | 1.18 | 1.00 | 118 | 50 - 150 | | Pyridaben | ND | 0.1 | 1.03 | 1.00 | 103 | 50 - 150 | |
| Daminozide | ND | 0.5 | 0.229 | 1.00 | 22.9 | 10 - 150 | | Spinosad | ND | 0.1 | 1.14 | 1.00 | 114 | 50 - 150 | |
| Diazinon | ND | 0.1 | 0.909 | 1.00 | 90.9 | 50 - 150 | | Spiromesifen | ND | 0.1 | 0.975 | 1.00 | 97.5 | 50 - 150 | |
| Dichlorvos (DDVP) | ND | 0.5 | 1.02 | 1.00 | 102 | 50 - 150 | | Spirotetramat | ND | 0.1 | 0.901 | 1.00 | 90.1 | 50 - 150 | |
| Dimethoate | ND | 0.1 | 0.955 | 1.00 | 95.5 | 50 - 150 | | Spiroxamine | ND | 0.1 | 0.990 | 1.00 | 99.0 | 50 - 150 | |
| Ethoprophos | ND | 0.1 | 0.942 | 1.00 | 94.2 | 50 - 150 | | Tebuconazole | ND | 0.1 | 0.894 | 1.00 | 89.4 | 50 - 150 | |
| Etofenprox | ND | 0.1 | 1.24 | 1.00 | 124 | 50 - 150 | | Thiacloprid | ND | 0.1 | 0.897 | 1.00 | 89.7 | 50 - 150 | |
| Etoxazole | ND | 0.1 | 0.901 | 1.00 | 90.1 | 50 - 150 | | Thiamethoxam | ND | 0.1 | 0.909 | 1.00 | 90.9 | 50 - 150 | |
| Fenoxycarb | ND | 0.1 | 0.915 | 1.00 | 91.5 | 50 - 150 | | Trifloxystrobin | ND | 0.1 | 0.896 | 1.00 | 89.6 | 50 - 150 | |
| Fenpyroximate | ND | 0.1 | 0.805 | 1.00 | 80.5 | 50 - 150 | | | | | | | | | |
| Fipronil | ND | 0.1 | 0.961 | 1.00 | 96.1 | 50 - 150 | | | | | | | | | |
| Flonicamid | ND | 0.1 | 0.946 | 1.00 | 94.6 | 50 - 150 | | | | | | | | | |
| Fludioxonil | ND | 0.1 | 0.996 | 1.00 | 99.6 | 50 - 150 | | | | | | | | | |
| Hexythiazox | ND | 0.1 | 0.588 | 1.00 | 58.8 | 50 - 150 | | | | | | | | | |
| Imazalil | ND | 0.1 | 1.02 | 1.00 | 102 | 50 - 150 | | | | | | | | | |
| Imidacloprid | ND | 0.1 | 0.851 | 1.00 | 85.1 | 50 - 150 | | | | | | | | | |
| Kresoxim-methyl | ND | 0.1 | 0.923 | 1.00 | 92.3 | 50 - 150 | | | | | | | | | |
| Malathion | ND | 0.1 | 0.936 | 1.00 | 93.6 | 50 - 150 | | | | | | | | | |

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Sample Procedure: 160721_LAB-SOP_SampleCollection-v008

Residual Solvents Sample Data

Solvents Analysis Date: 8/19/2020
 Solvents Batch ID: RES_081820A

Method: EPA 5021A
 Unit: µg/g (ppm)

Pass 

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

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Residual Solvents Quality Control Data

Solvents QC Analysis Date: 8/19/2020
 Solvents QC Batch ID: RES_081820A

Method: EPA 5021A
 Unit: µg/g (ppm)

| Analyte | Blank | LOQ | LCS | LCS Spike | LCS Rec (%) | Limits (%) | Notes |
|--------------------------|-------|-------|------|-----------|-------------|------------|-------|
| 1,4-Dioxane | ND | 50.0 | 1030 | 1000 | 103 | 70 - 130 | |
| 2-Butanol | ND | 250.0 | 923 | 1000 | 92.3 | 70 - 130 | |
| 2-Ethoxyethanol | ND | 50.0 | 1000 | 1000 | 100 | 70 - 130 | |
| Acetone | ND | 250.0 | 1030 | 1000 | 103 | 70 - 130 | |
| Acetonitrile | ND | 50.0 | 897 | 1000 | 89.7 | 70 - 130 | |
| Benzene | ND | 2.0 | 19.0 | 20.0 | 95.1 | 70 - 130 | |
| Butanes | <LOQ | 250.0 | 1670 | 2000 | 83.4 | 70 - 130 | |
| Cumene | ND | 50.0 | 974 | 1000 | 97.4 | 70 - 130 | |
| Cyclohexane | ND | 50.0 | 991 | 1000 | 99.1 | 70 - 130 | |
| Ethyl Acetate | ND | 250.0 | 895 | 1000 | 89.5 | 70 - 130 | |
| Ethyl Ether | ND | 250.0 | 989 | 1000 | 98.9 | 70 - 130 | |
| Ethylene Glycol | ND | 250.0 | 1060 | 1000 | 106 | 70 - 130 | |
| Ethylene Oxide | ND | 50.0 | 875 | 1000 | 87.5 | 70 - 130 | |
| Heptane | ND | 250.0 | 876 | 1000 | 87.6 | 70 - 130 | |
| Hexanes | ND | 50.0 | 4550 | 5000 | 91.0 | 70 - 130 | |
| Isopropanol (2-Propanol) | <LOQ | 50.0 | 937 | 1000 | 93.7 | 70 - 130 | |
| Isopropyl Acetate | ND | 250.0 | 909 | 1000 | 90.9 | 70 - 130 | |
| Methanol | ND | 250.0 | 1080 | 1000 | 108 | 70 - 130 | |
| Dichloromethane | ND | 50.0 | 950 | 1000 | 95.0 | 70 - 130 | |
| Pentanes | ND | 250.0 | 2680 | 3000 | 89.4 | 70 - 130 | |
| Propane | <LOQ | 250.0 | 1380 | 1000 | 138 | 70 - 130 | ME |
| Tetrahydrofuran | ND | 50.0 | 877 | 1000 | 87.7 | 70 - 130 | |
| Toluene | ND | 50.0 | 1010 | 1000 | 101 | 70 - 130 | |
| Xylenes | ND | 50.0 | 5180 | 4000 | 130 | 70 - 130 | |

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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 |