

**SROAZ LLC**

Tempe, AZ 85281

Lic. #00000109ESVM44878444

Harvest Dates: 10/25/2023

**Space Dust - Space Berry**

Concentrates & Extracts, Diamonds, Extraction Method: CO2

Dispensary License #: ; Manufacturing License #: 00000109ESVM44878444; Cultivation License #: 00000077DCPS00216601

**Sample: 2404TLL0133.0657**

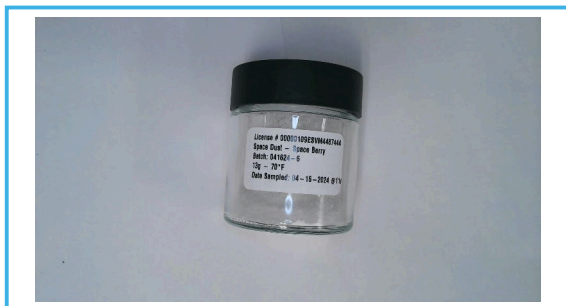
Strain: Space Berry

Parent Batch #: ; Batch#: 041624-6; Batch Size: g

Sample Received: 04/17/2024; Report Created: 04/23/2024; Expires: 04/23/2025

Manufacturing Date: 04/16/2024

Sampling: ; Environment:



## Safety

|                           |                           |                                     |
|---------------------------|---------------------------|-------------------------------------|
| <b>Pass</b><br>Pesticides | <b>Pass</b><br>Microbials | <b>Pass</b><br>Mycotoxins           |
| <b>Pass</b><br>Solvents   | <b>Pass</b><br>Metals     | <b>Not Tested</b><br>Foreign Matter |

## Cannabinoids

TPL\_Potency\_01

|                            |                        |   |
|----------------------------|------------------------|---|
| <b>85.31%</b><br>Total THC | <b>ND</b><br>Total CBD | <b>97.28%</b><br>Total Cannabinoids<br>Q3 |
|----------------------------|------------------------|---|

| Analyte      | LOQ  | Mass         | Mass         | Qualifier |
|--------------|------|--------------|--------------|-----------|
|              | %    | %            | mg/g         |           |
| THCa         | 0.20 | 97.28        | 972.8        | D1        |
| Δ9-THC       | 0.20 | ND           | ND           | D1        |
| Δ8-THC       | 0.20 | ND           | ND           | D1        |
| THCV         | 0.20 | ND           | ND           | D1        |
| CBDa         | 0.20 | ND           | ND           | D1        |
| CBD          | 0.20 | ND           | ND           | D1        |
| CBDV         | 0.20 | ND           | ND           | D1        |
| CBN          | 0.20 | ND           | ND           | D1        |
| CBGa         | 0.20 | ND           | ND           | D1        |
| CBG          | 0.20 | ND           | ND           | D1        |
| CBC          | 0.20 | ND           | ND           | D1        |
| <b>Total</b> |      | <b>97.28</b> | <b>972.8</b> |           |

Total THC = THCa \* 0.877 + Δ9-THC  
Total CBD = CBDa \* 0.877 + CBD  
Instrument: HPLC-DAD; ; Method: TPL\_Potency\_01

## Terpenes

TPL\_Terpenes\_01

|  |  |  |  |
|--|--|--|--|
|  |  |  |  |
|--|--|--|--|

| Analyte | LOQ | Mass | Mass | Qualifier |
|---------|-----|------|------|-----------|
|---------|-----|------|------|-----------|

Notes:

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**Pesticides TPL\_Pesticides\_01**

**Pass**

| Analyte             | LOQ  | Limit | Mass | Status | Qualifier | Analyte         | LOQ  | Limit | Mass | Status | Qualifier |
|---------------------|------|-------|------|--------|-----------|-----------------|------|-------|------|--------|-----------|
|                     | PPM  | PPM   | PPM  |        |           |                 | PPM  | PPM   | PPM  |        |           |
| Abamectin           | 0.24 | 0.50  | ND   | Pass   | V1        | Hexythiazox     | 0.48 | 1.00  | ND   | Pass   |           |
| Acephate            | 0.19 | 0.40  | ND   | Pass   | V1        | Imazalil        | 0.10 | 0.20  | ND   | Pass   |           |
| Acetamiprid         | 0.10 | 0.20  | ND   | Pass   |           | Imidacloprid    | 0.19 | 0.40  | ND   | Pass   |           |
| Aldicarb            | 0.19 | 0.40  | ND   | Pass   | L1        | Kresoxim        | 0.19 | 0.40  | ND   | Pass   |           |
| Azoxystrobin        | 0.10 | 0.20  | ND   | Pass   |           | Methyl          |      |       | ND   | Pass   |           |
| Bifenazate          | 0.10 | 0.20  | ND   | Pass   | L1 V1     | Malathion       | 0.10 | 0.20  | ND   | Pass   |           |
| Bifenthrin          | 0.10 | 0.20  | ND   | Pass   |           | Metalaxyl       | 0.10 | 0.20  | ND   | Pass   |           |
| Boscalid            | 0.19 | 0.40  | ND   | Pass   | L1        | Methiocarb      | 0.10 | 0.20  | ND   | Pass   | L1        |
| Carbaryl            | 0.10 | 0.20  | ND   | Pass   | L1 V1     | Methomyl        | 0.19 | 0.40  | ND   | Pass   | L1 V1     |
| Carbofuran          | 0.10 | 0.20  | ND   | Pass   | L1 V1     | Myclobutanil    | 0.10 | 0.20  | ND   | Pass   |           |
| Chlorantraniliprole | 0.10 | 0.20  | ND   | Pass   | L1        | Naled           | 0.24 | 0.50  | ND   | Pass   | L1 V1     |
| Chlorfenapyr        | 0.48 | 1.00  | ND   | Pass   |           | Oxamyl          | 0.48 | 1.00  | ND   | Pass   | L1 V1     |
| Chlorpyrifos        | 0.10 | 0.20  | ND   | Pass   | L1 V1     | Pacllobutrazol  | 0.19 | 0.40  | ND   | Pass   |           |
| Clofentezine        | 0.10 | 0.20  | ND   | Pass   |           | Permethrin      | 0.10 | 0.20  | ND   | Pass   | M2        |
| Cyfluthrin          | 0.48 | 1.00  | ND   | Pass   |           | Phosmet         | 0.10 | 0.20  | ND   | Pass   | L1        |
| Cypermethrin        | 0.48 | 1.00  | ND   | Pass   |           | Piperonyl       |      |       |      |        |           |
| Daminozide          | 0.48 | 1.00  | ND   | Pass   |           | Butoxide        | 0.96 | 2.00  | <LOQ | Pass   |           |
| Diazinon            | 0.10 | 0.20  | ND   | Pass   | L1 V1     | Prallethrin     | 0.10 | 0.20  | ND   | Pass   |           |
| Dichlorvos          | 0.05 | 0.10  | ND   | Pass   | V1        | Propiconazole   | 0.19 | 0.40  | ND   | Pass   | M2        |
| Dimethoate          | 0.10 | 0.20  | ND   | Pass   | L1 V1     | Propoxur        | 0.10 | 0.20  | ND   | Pass   | L1 V1     |
| Ethoprophos         | 0.10 | 0.20  | ND   | Pass   | L1        | Pyrethrins      | 0.48 | 1.00  | ND   | Pass   |           |
| Etofenprox          | 0.19 | 0.40  | ND   | Pass   | M2        | Pyridaben       | 0.10 | 0.20  | ND   | Pass   |           |
| Etoxazole           | 0.10 | 0.20  | ND   | Pass   |           | Spinosad        | 0.10 | 0.20  | ND   | Pass   |           |
| Fenoxycarb          | 0.10 | 0.20  | ND   | Pass   | V1        | Spiromesifen    | 0.10 | 0.20  | ND   | Pass   |           |
| Fenpyroximate       | 0.19 | 0.40  | ND   | Pass   |           | Spirotetramat   | 0.10 | 0.20  | ND   | Pass   |           |
| Fipronil            | 0.19 | 0.40  | ND   | Pass   |           | Spiroxamine     | 0.19 | 0.40  | ND   | Pass   |           |
| Fonicamid           | 0.48 | 1.00  | ND   | Pass   |           | Tebuconazole    | 0.19 | 0.40  | ND   | Pass   |           |
| Fludioxonil         | 0.19 | 0.40  | ND   | Pass   |           | Thiacloprid     | 0.10 | 0.20  | ND   | Pass   |           |
|                     |      |       |      |        |           | Thiamethoxam    | 0.10 | 0.20  | ND   | Pass   |           |
|                     |      |       |      |        |           | Trifloxystrobin | 0.10 | 0.20  | ND   | Pass   |           |

Instrument: LC-QQQ ; Method: TPL\_Pesticides\_01



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## Heavy Metals Pass

| Analyte | LOQ   | Limit  | Mass | Status | Qualifier |
|---------|-------|--------|------|--------|-----------|
|         | PPB   | PPB    | PPB  |        |           |
| Arsenic | 200.0 | 400.0  | ND   | Pass   |           |
| Cadmium | 200.0 | 400.0  | <LOQ | Pass   |           |
| Lead    | 500.0 | 1000.0 | <LOQ | Pass   |           |
| Mercury | 100.0 | 200.0  | <LOQ | Pass   | L1        |

LOQ=Limit of Quantitation. The reported result is based on a simple weight with the applicable moisture content for that sample. Unless otherwise stated, all quality control samples performed within specifications established by the Laboratory. Instrument: ICPMS; Method: AOAC 2021.03

## Residual Solvents Pass

| Analyte           | LOQ    | Limit  | Mass | Status | Qualifier |
|-------------------|--------|--------|------|--------|-----------|
|                   | PPM    | PPM    | PPM  |        |           |
| Acetone           | 460.0  | 1000.0 | ND   | Pass   |           |
| Acetonitrile      | 188.6  | 410.0  | ND   | Pass   |           |
| Benzene           | 0.9    | 2.0    | ND   | Pass   |           |
| Butanes           | 575.0  | 5000.0 | ND   | Pass   |           |
| Chloroform        | 27.6   | 60.0   | ND   | Pass   |           |
| Dichloromethane   | 276.0  | 600.0  | ND   | Pass   |           |
| Ethanol           | 2299.9 | 5000.0 | ND   | Pass   |           |
| Ethyl-Acetate     | 2299.9 | 5000.0 | ND   | Pass   |           |
| Ethyl-Ether       | 2299.9 | 5000.0 | ND   | Pass   |           |
| Heptane           | 2299.9 | 5000.0 | ND   | Pass   |           |
| Hexanes           | 133.4  | 290.0  | ND   | Pass   |           |
| Isopropyl-Acetate | 2299.9 | 5000.0 | ND   | Pass   |           |
| Methanol          | 1379.9 | 3000.0 | ND   | Pass   |           |
| Pentanes          | 133.4  | 5000.0 | ND   | Pass   |           |
| 2-Propanol        | 2299.9 | 5000.0 | ND   | Pass   |           |
| Toluene           | 409.4  | 890.0  | ND   | Pass   |           |
| Xylenes           | 92.0   | 2170.0 | ND   | Pass   |           |

Performed by GCMS-HS SOP-004. Methods used per AZDHS R9-17-404.03 and the solvent limits set by AZDHS R9-17 Table 3.1. AZDHS approved method for residual solvents by GCMS-HS for all listed analytes. Subcontracted through DVT Registration Certificate Identification Number : 0000031LRCHX78341676

## Microbials Pass

| Analyte | LOQ   | Limit | Result | Status | Qualifier |
|---------|-------|-------|--------|--------|-----------|
|         | CFU/g | CFU/g | CFU/g  |        |           |
| E. Coli | 10    | 100   | <10    | Pass   |           |

| Analyte               | Limit            | Result       | Status | Qualifier |
|-----------------------|------------------|--------------|--------|-----------|
| Salmonella            | Detectable in 1g | Not Detected | Pass   |           |
| Aspergillus           | Detectable in 1g | Not Detected | Pass   |           |
| Aspergillus fumigatus | Detectable in 1g | Not Detected | Pass   |           |
| Aspergillus niger     | Detectable in 1g | Not Detected | Pass   |           |
| Aspergillus flavus    | Detectable in 1g | Not Detected | Pass   |           |
| Aspergillus terreus   | Detectable in 1g | Not Detected | Pass   |           |

Instrument: qPCR/Plating; AOAC Methods 082102, 022202 and 2018.13

## Mycotoxins Pass

| Analyte          | LOQ | Limit | Mass | Status | Qualifier |
|------------------|-----|-------|------|--------|-----------|
|                  | PPB | PPB   | PPB  |        |           |
| B1               | 8.1 | 20.0  | ND   | Pass   | M2        |
| B2               | 8.1 | 20.0  | ND   | Pass   | M2        |
| G1               | 8.1 | 20.0  | ND   | Pass   | M2        |
| G2               | 8.1 | 20.0  | ND   | Pass   | M2        |
| Ochratoxin A     | 8.1 | 20.0  | ND   | Pass   | M2        |
| Total Aflatoxins | 8.1 | 20.0  | ND   | Pass   | M2        |

B1 = Target analyte detected in calibration blank was above LOQ but the concentration of cannabinoid was below LOQ.

B2 = Target analyte detected in calibration blank was above LOQ but was below the maximum allowable concentration.

D1 = The limit of quantitation and the sample results were adjusted to reflect sample dilution,

I1 = The relative intensity of a characteristic ion in a sample analyte exceeded the acceptance criteria with respect to the reference spectra, indicating interference,

L1 = The percent recovery of a laboratory control sample is greater than the acceptance limits in A.A.C 17 R9-17-404.03(K)(2)(C), but the sample's target analytes were not detected above the maximum allowed concentration,

M1 = The recovery from the matrix spike was high, but the recovery from the laboratory control sample was within acceptance criteria,

M2 = The recovery from the matrix spike was low, but the recovery from the laboratory control sample was within acceptance criteria,

M3 = The recovery from the matrix spike was unusable because the analyte concentration was disproportionate to the spike level, but the recovery from the laboratory control sample was within acceptance criteria,

M4 = The analysis of a spiked sample required a dilution such that the spike recovery calculation does not provide useful information, but the recovery from the associated laboratory control sample was within acceptance criteria,

M5 = The analyte concentration was determined by the method of standard addition, in which the standard is added directly to the aliquots of the analyzed sample,

N1 - A description of the variance is described in the final report of testing,

R1 = The relative percent difference for the laboratory control sample and duplicate exceeded the limit in A.A.C 17 R9-17-404.03(K)(3), but the recovery in subsection A.A.C 17 R9-17-404.03 (K)(2) was within accepted criteria,

R2 = The relative percent difference for a sample and duplicated exceeded the limit in subsection A.A.C 17 R9-17-404.03 (O)

Q1 = Sample integrity was not maintained,

Q2 = The sample is heterogeneous and sample homogeneity could not be readily achieved using routine laboratory practices

Q3 = Testing result is for informational purposes only and cannot be used to satisfy dispensary testing requirements in R9-17-317.01(A) or labeling requirements in R9-17-317

V1 = The recovery from continuing calibration verification standards exceeded the acceptance limits denoted in A.A.C 17 R9-17-403.03(J)(1)(b), but the sample's target analytes were not detected above the maximum allowable concentrations for the analytes in the sample.