1 of 8

1:1:1 CBG

Sample ID: SA-230727-25056

Batch: 080123G111

Type: Finished Product - Ingestible

Matrix: Edible - Gummy Unit Mass (g): 5.38143

Received: 08/14/2023 Completed: 08/25/2023 Client

SunFlora Inc. 600 8th Ave W, STE 400 Palmetto, FL 34221

USA



Summary

Test Cannabinoids Heavy Metals Microbials Mycotoxins Pesticides **Residual Solvents** Terpenes

Date Tested 08/16/2023 08/23/2023 08/18/2023 08/18/2023 08/18/2023 08/25/2023 08/21/2023

Status Tested Tested Tested Tested Tested Tested Tested

0.178 % Total Δ9-THC 0.235 % CBD

0.700 % **Total Cannabinoids**

Not Tested Moisture Content

Not Tested Foreign Matter

Internal Standard Normalization

Yes

Cannabinoids by HPLC-PDA, LC-MS/MS, and/or GC-MS/MS

| Analyte | LOD (%) | LOQ (%) | Result (%) | Result (mg/unit) |
|--------------|------------|------------|---|---------------------|
| CBC | 0.00095 | 0.00284 | 0.0590 | 3.18 |
| CBCA | 0.00181 | 0.00543 | ND | ND |
| CBCV | 0.0006 | 0.0018 | ND | ND |
| CBD | 0.00081 | 0.00242 | 0.235 | 12.6 |
| CBDA | 0.00043 | 0.0013 | ND | ND |
| CBDV | 0.00061 | 0.00182 | 0.00435 | 0.234 |
| CBDVA | 0.00021 | 0.00063 | ND | ND |
| CBG | 0.00057 | 0.00172 | 0.201 | 10.8 |
| CBGA | 0.00049 | 0.00147 | ND | ND |
| CBL | 0.00112 | 0.00335 | 0.00365 | 0.196 |
| CBLA | 0.00124 | 0.00371 | ND | ND |
| CBN | 0.00056 | 0.00169 | <loq< td=""><td><loq< td=""></loq<></td></loq<> | <loq< td=""></loq<> |
| CBNA | 0.0006 | 0.00181 | ND | ND |
| СВТ | 0.0018 | 0.0054 | 0.0198 | 1.06 |
| Δ8-ΤΗС | 0.00104 | 0.00312 | <loq< td=""><td><loq< td=""></loq<></td></loq<> | <loq< td=""></loq<> |
| Δ9-ΤΗС | 0.00076 | 0.00227 | 0.178 | 9.58 |
| Δ9-ΤΗCΑ | 0.00084 | 0.00251 | ND | ND |
| Δ9-THCV | 0.00069 | 0.00206 | ND | ND |
| Δ9-THCVA | 0.00062 | 0.00186 | ND | ND |
| Total Δ9-THC | | | 0.178 | 9.58 |
| Total | | | 0.700 | 37.7 |

ND = Not Detected; NT = Not Tested; LOD = Limit of Detection; LOQ = Limit of Quantitation; RL = Reporting Limit; Δ = Delta; Total Δ9-THC = Δ9-THCA * 0.877 + Δ9-THC; Total CBD = CBDA * 0.877 + CBD;

Generated By: Ryan Bellone CCO

Date: 08/25/2023

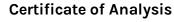
Tested By: Nicholas Howard Scientist Date: 08/16/2023



Accreditation #108651









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2 of 8

1:1:1 CBG

Sample ID: SA-230727-25056

Batch: 080123G111

Type: Finished Product - Ingestible

Matrix: Edible - Gummy Unit Mass (g): 5.38143 Received: 08/14/2023 Completed: 08/25/2023 Client

SunFlora Inc. 600 8th Ave W, STE 400 Palmetto, FL 34221

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Terpenes by GC-MS

| Analyte | LOD (%) | LOQ (%) | Result (%) | Analyte | LOD (%) | LOQ (%) | Result (%) |
|---------------------|------------|------------|---|------------------------|------------|------------|---------------------|
| α -Bisabolol | 0.0002 | 0.001 | 0.00816 | Limonene | 0.0002 | 0.001 | 0.0149 |
| (+)-Borneol | 0.0002 | 0.001 | <loq< th=""><th>Linalool</th><th>0.0002</th><th>0.001</th><th>0.00165</th></loq<> | Linalool | 0.0002 | 0.001 | 0.00165 |
| Camphene | 0.0002 | 0.001 | <loq< th=""><th>β-myrcene</th><th>0.0002</th><th>0.001</th><th>0.00169</th></loq<> | β-myrcene | 0.0002 | 0.001 | 0.00169 |
| Camphor | 0.0004 | 0.002 | ND | Nerol | 0.0002 | 0.001 | ND |
| 3-Carene | 0.0002 | 0.001 | <loq< th=""><th>cis-Nerolidol</th><th>0.0002</th><th>0.001</th><th>ND</th></loq<> | cis-Nerolidol | 0.0002 | 0.001 | ND |
| β-Caryophyllene | 0.0002 | 0.001 | 0.00395 | trans-Nerolidol | 0.0002 | 0.001 | 0.00157 |
| Caryophyllene Oxide | 0.0002 | 0.001 | 0.00307 | Ocimene | 0.0002 | 0.001 | ND |
| α -Cedrene | 0.0002 | 0.001 | ND | α -Phellandrene | 0.0002 | 0.001 | <loq< th=""></loq<> |
| Cedrol | 0.0002 | 0.001 | ND | α-Pinene | 0.0002 | 0.001 | <loq< th=""></loq<> |
| Eucalyptol | 0.0002 | 0.001 | <loq< th=""><th>β-Pinene</th><th>0.0002</th><th>0.001</th><th><loq< th=""></loq<></th></loq<> | β-Pinene | 0.0002 | 0.001 | <loq< th=""></loq<> |
| Fenchone | 0.0004 | 0.002 | <loq< th=""><th>Pulegone</th><th>0.0002</th><th>0.001</th><th>ND</th></loq<> | Pulegone | 0.0002 | 0.001 | ND |
| Fenchyl Alcohol | 0.0002 | 0.001 | <loq< th=""><th>Sabinene</th><th>0.0002</th><th>0.001</th><th>ND</th></loq<> | Sabinene | 0.0002 | 0.001 | ND |
| Geraniol | 0.0002 | 0.001 | ND | Sabinene Hydrate | 0.0002 | 0.001 | ND |
| Geranyl Acetate | 0.0002 | 0.001 | 0.00125 | α-Terpinene | 0.0002 | 0.001 | ND |
| Guaiol | 0.0002 | 0.001 | 0.00215 | γ-Terpinene | 0.0002 | 0.001 | <loq< th=""></loq<> |
| Hexahydrothymol | 0.0002 | 0.001 | ND | α-Terpineol | 0.0001 | 0.0005 | 0.00055 |
| α -Humulene | 0.0002 | 0.001 | 0.0022 | γ-Terpineol | 0.0001 | 0.0005 | ND |
| Isoborneol | 0.0002 | 0.001 | ND | Terpinolene | 0.0002 | 0.001 | 0.001 |
| Isopulegol | 0.0002 | 0.001 | ND | Valencene | 0.0002 | 0.001 | ND |
| | | | | Total Terpenes (%) | | | 0.0486 |

ND = Not Detected; NT = Not Tested; LOD = Limit of Detection; LOQ = Limit of Quantitation; P = Pass; F = Fail; RL = Reporting Limit











Generated By: Ryan Bellone CCO

Date: 08/25/2023

Tested By: Jasper van Heemst Principal Scientist Date: 08/21/2023





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Certificate of Analysis

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1:1:1 CBG

Sample ID: SA-230727-25056 Batch: 080123G111

Type: Finished Product - Ingestible

Matrix: Edible - Gummy Unit Mass (g): 5.38143 Received: 08/14/2023 Completed: 08/25/2023 Client

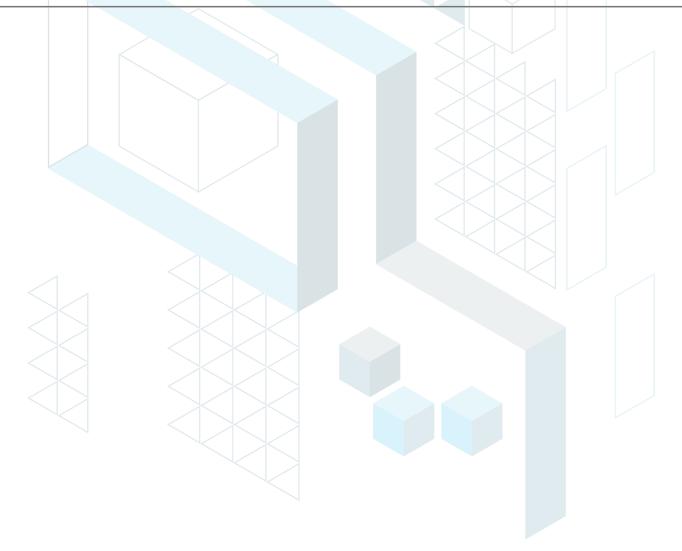
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USA

Heavy Metals by ICP-MS

| Analyte | LOD (ppb) | LOQ (ppb) | Result (ppb) |
|---------|-----------|-----------|---------------------|
| Arsenic | 2 | 20 | <loq< td=""></loq<> |
| Cadmium | 1 | 20 | ND |
| Lead | 2 | 20 | <rl< td=""></rl<> |
| Mercury | 12 | 50 | ND |

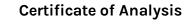
ND = Not Detected; NT = Not Tested; LOD = Limit of Detection; LOQ = Limit of Quantitation; P = Pass; F = Fail; RL = Reporting Limit



Generated By: Ryan Bellone CCO Date: 08/25/2023

Tested By: Chris Farman Scientist Date: 08/23/2023







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1:1:1 CBG

Sample ID: SA-230727-25056

Batch: 080123G111

Type: Finished Product - Ingestible

Matrix: Edible - Gummy Unit Mass (g): 5.38143 Received: 08/14/2023 Completed: 08/25/2023 Client

SunFlora Inc. 600 8th Ave W, STE 400 Palmetto, FL 34221

USA

Pesticides by LC-MS/MS

| Analyte | LOD (ppb) | LOQ (ppb) | Result (ppb) | Analyte | LOD (ppb) | LOQ (ppb) | Result (ppb) |
|----------------------|--------------|--------------|-----------------|--------------------|--------------|--------------|-----------------|
| Acephate | 30 | 100 | ND | Hexythiazox | 30 | 100 | ND |
| Acetamiprid | 30 | 100 | ND | Imazalil | 30 | 100 | ND |
| Aldicarb | 30 | 100 | ND | Imidacloprid | 30 | 100 | ND |
| Azoxystrobin | 30 | 100 | ND | Kresoxim methyl | 30 | 100 | ND |
| Bifenazate | 30 | 100 | ND | Malathion | 30 | 100 | ND |
| Bifenthrin | 30 | 100 | ND | Metalaxyl | 30 | 100 | ND |
| Boscalid | 30 | 100 | ND | Methiocarb | 30 | 100 | ND |
| Carbaryl | 30 | 100 | ND | Methomyl | 30 | 100 | ND |
| Carbofuran | 30 | 100 | ND | Mevinphos | 30 | 100 | ND |
| Chloranthraniliprole | 30 | 100 | ND | Myclobutanil | 30 | 100 | ND |
| Chlorfenapyr | 30 | 100 | ND | Oxamyl | 30 | 100 | ND |
| Chlorpyrifos | 30 | 100 | ND | Paclobutrazol | 30 | 100 | ND |
| Clofentezine | 30 | 100 | ND | Permethrin | 30 | 100 | ND |
| Coumaphos | 30 | 100 | ND | Phosmet | 30 | 100 | ND |
| Daminozide | 30 | 100 | ND | Piperonyl Butoxide | 30 | 100 | ND |
| Diazinon | 30 | 100 | ND | Prallethrin | 30 | 100 | ND |
| Dichlorvos | 30 | 100 | ND | Propiconazole | 30 | 100 | ND |
| Dimethoate | 30 | 100 | ND | Propoxur | 30 | 100 | ND |
| Dimethomorph | 30 | 100 | ND | Pyrethrins | 30 | 100 | ND |
| Ethoprophos | 30 | 100 | ND | Pyridaben | 30 | 100 | ND |
| Etofenprox | 30 | 100 | ND | Spinetoram | 30 | 100 | ND |
| Etoxazole | 30 | 100 | ND | Spinosad | 30 | 100 | ND |
| Fenhexamid | 30 | 100 | ND | Spiromesifen | 30 | 100 | ND |
| Fenoxycarb | 30 | 100 | ND | Spirotetramat | 30 | 100 | ND |
| Fenpyroximate | 30 | 100 | ND | Spiroxamine | 30 | 100 | ND |
| Fipronil | 30 | 100 | ND | Tebuconazole | 30 | 100 | ND |
| Flonicamid | 30 | 100 | ND | Thiacloprid | 30 | 100 | ND |
| Fludioxonil | 30 | 100 | ND | Thiamethoxam | 30 | 100 | ND |
| | | | | Trifloxystrobin | 30 | 100 | ND |

ND = Not Detected; NT = Not Tested; LOD = Limit of Detection; LOQ = Limit of Quantitation; P = Pass; F = Fail; RL = Reporting Limit

Generated By: Ryan Bellone CCO

Date: 08/25/2023

Tested By: Jasper van Heemst Principal Scientist Date: 08/18/2023





KCA Laboratories 232 North Plaza Drive

Nicholasville, KY 40356

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Certificate of Analysis

5 of 8

1:1:1 CBG

Sample ID: SA-230727-25056 Batch: 080123G111

Type: Finished Product - Ingestible Matrix: Edible - Gummy

Unit Mass (g): 5.38143

Received: 08/14/2023 Completed: 08/25/2023 Client

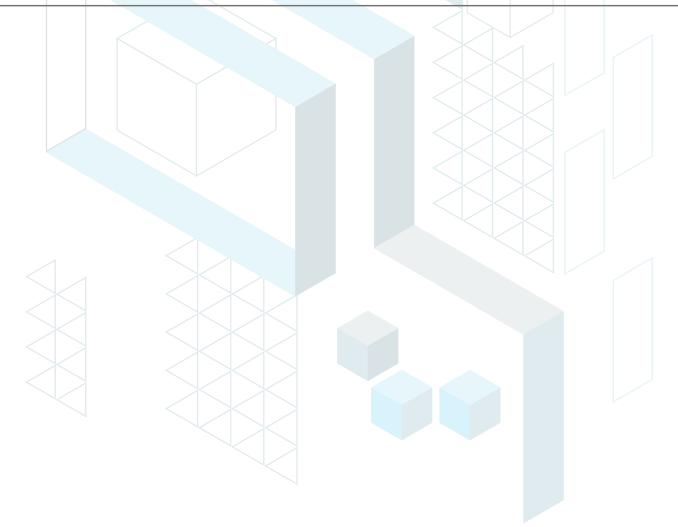
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USA

Mycotoxins by LC-MS/MS

| Analyte | LOD (p) | bb) LOQ (ppb) | Result (ppb) | |
|--------------|---------|---------------|--------------|--|
| B1 | 1 | 5 | ND | |
| B2 | 1 | 5 | ND | |
| G1 | 1 | 5 | ND | |
| G2 | 1 | 5 | ND | |
| Ochratoxin A | 1 | 5 | ND | |
| | | | | |

ND = Not Detected; NT = Not Tested; LOD = Limit of Detection; LOQ = Limit of Quantitation; P = Pass; F = Fail; RL = Reporting Limit



Generated By: Ryan Bellone CCO

Date: 08/25/2023

Tested By: Jasper van Heemst Principal Scientist Date: 08/18/2023





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Certificate of Analysis

6 of 8

1:1:1 CBG

Sample ID: SA-230727-25056 Batch: 080123G111 Type: Finished Product - Ingestible

Matrix: Edible - Gummy Unit Mass (g): 5.38143 Received: 08/14/2023 Completed: 08/25/2023 Client

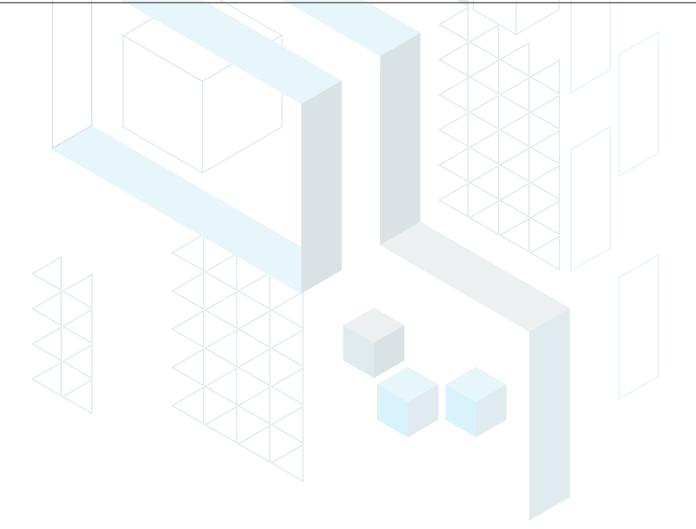
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USA

Microbials by PCR and Plating

| Analyte | LOD (CFU/g) | Result (CFU/g) | Result (Qualitative) |
|--------------------------------------|-------------|----------------|-------------------------|
| Total aerobic count | 1 | ND | |
| Total coliforms | 1 | ND | |
| Generic E. coli | 1 | ND | |
| Salmonella spp. | 1 | | Not Detected per 1 gram |
| Shiga-toxin producing E. coli (STEC) | 1 | | Not Detected per 1 gram |
| | | | |

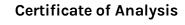
ND = Not Detected; NT = Not Tested; LOD = Limit of Detection; LOQ = Limit of Quantitation; CFU = Colony Forming Units; P = Pass; F = Fail; RL = Reporting Limit



Generated By: Ryan Bellone CCO

Date: 08/25/2023

Tested By: Jade Pinkston Microbiology Technician Date: 08/18/2023





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1:1:1 CBG

Sample ID: SA-230727-25056

Batch: 080123G111

Type: Finished Product - Ingestible

Matrix: Edible - Gummy Unit Mass (g): 5.38143

Received: 08/14/2023 Completed: 08/25/2023 Client

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Residual Solvents by HS-GC-MS

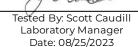
| | J | | | | | | |
|-----------------------|--------------|--------------|--|--------------------------|--------------|--------------|-----------------|
| Analyte | LOD (ppm) | LOQ (ppm) | Result (ppm) | Analyte | LOD (ppm) | LOQ (ppm) | Result (ppm) |
| Acetone | 167 | 500 | ND | Ethylene Glycol | 21 | 62 | ND |
| Acetonitrile | 14 | 41 | ND | Ethylene Oxide | 0.5 | 1 | ND |
| Benzene | 0.5 | 1 | ND | Heptane | 167 | 500 | ND |
| Butane | 167 | 500 | ND | n-Hexane | 10 | 29 | ND |
| 1-Butanol | 167 | 500 | ND | Isobutane | 167 | 500 | ND |
| 2-Butanol | 167 | 500 | ND | Isopropyl Acetate | 167 | 500 | ND |
| 2-Butanone | 167 | 500 | ND | Isopropyl Alcohol | 167 | 500 | ND |
| Chloroform | 2 | 6 | ND | Isopropylbenzene | 167 | 500 | ND |
| Cyclohexane | 129 | 388 | ND | Methanol | 100 | 300 | ND |
| 1,2-Dichloroethane | 0.5 | 1 | ND | 2-Methylbutane | 10 | 29 | ND |
| 1,2-Dimethoxyethane | 4 | 10 | ND | Methylene Chloride | 20 | 60 | ND |
| Dimethyl Sulfoxide | 167 | 500 | ND | 2-Methylpentane | 10 | 29 | ND |
| N,N-Dimethylacetamide | 37 | 109 | ND | 3-Methylpentane | 10 | 29 | ND |
| 2,2-Dimethylbutane | 10 | 29 | ND | n-Pentane | 167 | 500 | ND |
| 2,3-Dimethylbutane | 10 | 29 | ND | 1-Pentanol | 167 | 500 | ND |
| N,N-Dimethylformamide | 30 | 88 | ND | n-Propane | 167 | 500 | ND |
| 2,2-Dimethylpropane | 167 | 500 | ND | 1-Propanol | 167 | 500 | ND |
| 1,4-Dioxane | 13 | 38 | ND | Pyridine | < 7 | 20 | ND |
| Ethanol | 167 | 500 | <rl< td=""><td>Tetrahydrofuran</td><td>24</td><td>72</td><td>ND</td></rl<> | Tetrahydrofuran | 24 | 72 | ND |
| 2-Ethoxyethanol | 6 | 16 | ND | Toluene | 30 | 89 | ND |
| Ethyl Acetate | 167 | 500 | ND | Trichloroethylene | 3 | 8 | ND |
| Ethyl Ether | 167 | 500 | ND | Tetramethylene Sulfone | 6 | 16 | ND |
| Ethylbenzene | 3 | 7 | ND | Xylenes (o-, m-, and p-) | 73 | 217 | ND |
| | | | | | | | |

ND = Not Detected; NT = Not Tested; LOD = Limit of Detection; LOQ = Limit of Quantitation; P = Pass; F = Fail; RL = Reporting Limit

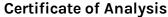


Generated By: Ryan Bellone CCO

Date: 08/25/2023









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8 of 8

1:1:1 CBG

Sample ID: SA-230727-25056

Batch: 080123G111

Type: Finished Product - Ingestible

Matrix: Edible - Gummy Unit Mass (g): 5.38143

Received: 08/14/2023 Completed: 08/25/2023 Client

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Reporting Limit Appendix

Heavy Metals - Colorado CDPHE

| Analyte | Li | mit (ppl |) Analyte | Limit (ppb) |
|---------|----|----------|-----------|-------------|
| Arsenic | | 1500 | Lead | 500 |
| Cadmium | | 500 | Mercury | 1500 |

Microbials -

| Analyte | Limit (CFU/ g) Analyte | Limit (CFU/ g) |
|-----------------|---------------------------|-------------------|
| Total coliforms | 100 Total aerobic count | 100000 |

Residual Solvents - USP 467

| Analyte | Limit (ppm) | Analyte | Limit (ppn |
|-----------------------|-------------|--------------------------|------------|
| Acetone | 5000 | Ethylene Glycol | 620 |
| Acetonitrile | 410 | Ethylene Oxide | 1 |
| Benzene | 2 | Heptane | 5000 |
| Butane | 5000 | n-Hexane | 290 |
| 1-Butanol | 5000 | Isobutane | 5000 |
| 2-Butanol | 5000 | Isopropyl Acetate | 5000 |
| 2-Butanone | 5000 | Isopropyl Alcohol | 5000 |
| Chloroform | 60 | Isopropylbenzene | 5000 |
| Cyclohexane | 3880 | Methanol | 3000 |
| 1,2-Dichloroethane | 5 | 2-Methylbutane | 290 |
| 1,2-Dimethoxyethane | 100 | Methylene Chloride | 600 |
| Dimethyl Sulfoxide | 5000 | 2-Methylpentane | 290 |
| N,N-Dimethylacetamide | 1090 | 3-Methylpentane | 290 |
| 2,2-Dimethylbutane | 290 | n-Pentane | 5000 |
| 2,3-Dimethylbutane | 290 | 1-Pentanol | 5000 |
| N,N-Dimethylformamide | 880 | n-Propane | 5000 |
| 2,2-Dimethylpropane | 5000 | 1-Propanol | 5000 |
| 1,4-Dioxane | 380 | Pyridine | 200 |
| Ethanol | 5000 | Tetrahydrofuran | 720 |
| 2-Ethoxyethanol | 160 | Toluene | 890 |
| Ethyl Acetate | 5000 | Trichloroethylene | 80 |
| Ethyl Ether | 5000 | Tetramethylene Sulfone | 160 |
| Ethylbenzene | 70 | Xylenes (o-, m-, and p-) | 2170 |

Pesticides - CA DCC

| Analyte | Limit (ppb) | Analyte | Limit (ppb) |
|---------------------|-------------|-------------------|-------------|
| Aldicarb | 30 | Imidacloprid | 3000 |
| Azoxystrobin | 40000 | Kresoxim methyl | 1000 |
| Bifenazate | 5000 | Malathion | 5000 |
| Bifenthrin | 500 | Metalaxyl | 15000 |
| Boscalid | 10000 | Methiocarb | 30 |
| Carbaryl | 500 | Methomyl | 100 |
| Carbofuran | 30 | Mevinphos | 30 |
| Chloranthraniliprol | e 40000 | Myclobutanil | 9000 |
| Chlorfenapyr | 30 | Oxamyl | 200 |
| Chlorpyrifos | 30 | Paclobutrazol | 30 |
| Clofentezine | 500 | Permethrin | 20000 |
| Coumaphos | 30 | Phosmet | 200 |
| Daminozide | 30 | Piperonyl Butoxic | le 8000 |
| Diazinon | 200 | Prallethrin | 400 |
| Dichlorvos | 30 | Propiconazole | 20000 |
| Dimethoate | 30 | Propoxur | 30 |
| Dimethomorph | 20000 | Pyrethrins | 1000 |
| Ethoprophos | 30 | Pyridaben | 3000 |
| Etofenprox | 30 | Spinetoram | 3000 |
| Etoxazole | 1500 | Spinosad | 3000 |
| Fenhexamid | 10000 | Spiromesifen | 12000 |
| Fenoxycarb | 30 | Spirotetramat | 13000 |
| Fenpyroximate | 2000 | Spiroxamine | 30 |
| Fipronil | 30 | Tebuconazole | 2000 |
| Flonicamid | 2000 | Thiacloprid | 30 |
| Fludioxonil | 30000 | Thiamethoxam | 4500 |

Mycotoxins - Colorado CDPHE

| Analyte | Limit (ppm) Analyte | Limit (ppm) |
|--------------|---------------------|-------------|
| B1 | 5 B2 | 5 |
| G1 | 5 G2 | 5 |
| Ochratoxin A | 5 | |

Pesticides - CA DCC

| Analyte | Limit (ppb) | Analyte | Limit (ppb) |
|-------------|-------------|-------------|-------------|
| Acephate | 5000 | Hexythiazox | 2000 |
| Acetamiprid | 5000 | Imazalil | 30 |

