

Farmer's Friend Extracts 6451 NE Colwood Wy Portland, OR 97218 503-442-8653 Sample Type: Extracts
Sample Date: 11/3/2020
Analysis Date: 11/4/2020
Report Date: 11/5/2020

Metrc Batch ID: 1A4010300016315000027948 Metrc Sample ID: 1A4010300016315000028008 Harvest/Process Date: 11/2/2020 Report ID: LS-201105-18 Sample Plan ID:SP-201104-3-A Sample Procedure: 160721_LAB-SOP_SampleCollection-v008

Potency

Potency Analysis Date: 11/4/2020 Potency Batch ID: CAN_110420C Potency Method: JAOAC 2015.1

79.2%

Total THC

0.101%

Total CBD

Samples: TBH-XST-SHH, PSR-ZBM-DRG



| Analyte | Description | LOQ | RPD | Min. | Max. | Avg. | |
|--------------|-------------------------------|------|------|-------|--|------------------------------|---|
| Д9ТНС | Delta-9 Tetrahydrocannabinol | 0.16 | 1.01 | 78.8 | 79.6 | 79.2 | |
| THCA | Tetrahydrocannabinolic acid | 0.16 | 0.00 | ND | ND | ND | _ |
| CBD | Cannabidiol | 0.16 | 23.3 | ND | 0.202 | 0.101 | • |
| CBDA | Cannabidiolic acid | 0.16 | 0.00 | ND | <l0q< td=""><td><l0q< td=""><td></td></l0q<></td></l0q<> | <l0q< td=""><td></td></l0q<> | |
| Δ8ΤΗC | Delta-8 Tetrahydrocannabinol* | 0.16 | 0.00 | ND | ND | ND | |
| THCV | Tetrahydrocannabivarin* | 0.16 | 7.85 | 2.44 | 2.64 | 2.54 | - |
| CBG | Cannabigerol* | 0.16 | 21.2 | 1.53 | 1.89 | 1.71 | • |
| CBGA | Cannabigerolic acid* | 0.16 | 33.7 | 0.418 | 0.587 | 0.502 | • |
| CBC | Cannabichromene* | 0.16 | 6.38 | 0.874 | 0.932 | 0.903 | • |
| CBCA | Cannabichromenic acid* | 0.16 | 0.00 | ND | ND | ND | _ |
| CBN | Cannabinol | 0.16 | 53.0 | 0.170 | 0.293 | 0.232 | • |
| Total THC | Δ9THC + (THCA × 0.877) | | 1.01 | 78.8 | 79.6 | 79.2 | _ |
| Total CBD | CBD + (CBDA × 0.877) | | 23.3 | ND | 0.202 | 0.101 | • |
| Total | | | 1.96 | 84.3 | 86.2 | 85.2 | |

Compliance

| Pesticides | Within limits | Analysis Date: 11/4/2020 | Pass 🕢 |
|------------|---------------|--------------------------|--------|
| Solvents | Within limits | Analysis Date: 11/5/2020 | Pass 🕢 |
| Potency | Within limits | Analysis Date: 11/4/2020 | Pass 🕢 |

Bryce Kidd, Ph.D.
Lab Director

Aaron Troyer //
Chief Science Officer





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

Sample Type: Extracts Sample Date: 11/3/2020 Analysis Date: 11/4/2020 Report Date: 11/5/2020

1A4010300016315000027948 Metrc Sample ID: 1A4010300016315000028008

Terpene Analysis Date: 11/4/2020

Terpene Batch ID: TRP 110420A

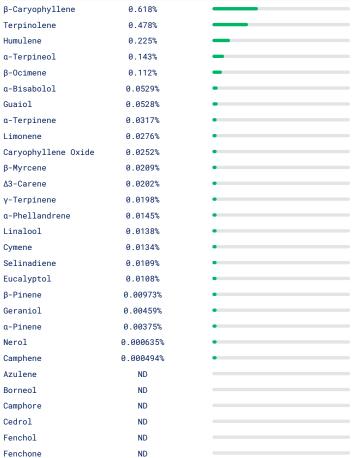
Harvest/Process Date: 11/2/2020 Report ID: LS-201105-18 Sample Plan ID:SP-201104-3-A Sample Procedure: 160721 LAB-SOP_SampleCollection-v008

Method: JAOAC 2015.1

Unit: %



| Sample Data | | | |
|-----------------|--------|-------|--|
| Analyte | Avg. | Notes | |
| β-Caryophyllene | 0.618% | | |
| Terpinolene | 0.478% | | |



| Analyte | Avg. | Notes |
|------------------|-------|-------|
| Geranyl Acetate | ND | |
| Isoborneol | ND | |
| Isopulegol | ND | |
| Pulegone | ND | |
| Sabinene | ND | |
| Sabinene Hydrate | ND | |
| Valencene | ND | |
| cis-Nerolidol | ND | |
| trans-Nerolidol | ND | |
| α-Cedrene | ND | |
| α-Ocimene | ND | |
| β-Farnesene 1 | ND | |
| β-Farnesene 2 | ND | |
| γ-Terpineol | ND | |
| Total | 1.91% | |



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1A4010300016315000027948 Metrc Sample ID: 1A4010300016315000028008 Harvest/Process Date: 11/2/2020 Report ID: LS-201105-18 Sample Plan ID:SP-201104-3-A Sample Procedure: 160721 LAB-SOP_SampleCollection-v008



Pesticides Analysis Date: 11/4/2020 Pesticides Batch ID: PST 110420C

Method: EN 15662 Unit: µg/g (ppm)

Pass 🕢

| Analyte | TBH-XST-SHH | PSR-ZBM-DRG | Limits | LOQ | Notes | Status |
|---------------------|-------------|-------------|--------|-----|-------|--------|
| Abamectin | ND | ND | 0.5 | 0.1 | | Pass |
| Acephate | ND | ND | 0.4 | 0.1 | | Pass |
| Acequinocyl | ND | ND | 2.0 | 1.5 | | Pass |
| Acetamiprid | ND | ND | 0.2 | 0.1 | | Pass |
| Aldicarb | ND | ND | 0.4 | 0.1 | | Pass |
| Azoxystrobin | ND | ND | 0.2 | 0.1 | | Pass |
| Bifenazate | ND | ND | 0.2 | 0.1 | | Pass |
| Bifenthrin | ND | ND | 0.2 | 0.1 | | Pass |
| Boscalid | ND | ND | 0.4 | 0.1 | | Pass |
| Carbaryl | ND | ND | 0.2 | 0.1 | | Pass |
| Carbofuran | ND | ND | 0.2 | 0.1 | | Pass |
| Chlorantraniliprole | ND | ND | 0.2 | 0.1 | | Pass |
| Chlorfenapyr | ND | ND | 1.0 | 0.1 | | Pass |
| Chlorpyrifos | ND | ND | 0.2 | 0.1 | | Pass |
| Clofentezine | ND | ND | 0.2 | 0.1 | | Pass |
| Cyfluthrin | ND | ND | 1.0 | 0.5 | | Pass |
| Cypermethrin | ND | ND | 1.0 | 0.1 | | Pass |
| Daminozide | ND | ND | 1.0 | 0.5 | | Pass |
| Diazinon | ND | ND | 0.2 | 0.1 | | Pass |
| Dichlorvos (DDVP) | ND | ND | 1.0 | 0.5 | | Pass |
| Dimethoate | ND | ND | 0.2 | 0.1 | | Pass |
| Ethoprophos | ND | ND | 0.2 | 0.1 | | Pass |
| Etofenprox | ND | ND | 0.4 | 0.1 | | Pass |
| Etoxazole | ND | ND | 0.2 | 0.1 | | Pass |
| Fenoxycarb | 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 |

| Analyte | TBH-XST-SHH | PSR-ZBM-DRG | Limits | LOQ | Notes | Status |
|--------------------|-------------|-------------|--------|-----|-------|--------|
| Metalaxyl | ND | ND | 0.2 | 0.1 | | Pass |
| Methiocarb | ND | ND | 0.2 | 0.1 | | Pass |
| Methomyl | ND | ND | 0.4 | 0.1 | | Pass |
| Methyl Parathion | ND | ND | 0.2 | 0.2 | | Pass |
| MGK-264 | ND | ND | 0.2 | 0.2 | | Pass |
| Myclobutanil | ND | ND | 0.2 | 0.1 | | Pass |
| Naled | ND | ND | 0.5 | 0.2 | | Pass |
| Oxamyl | ND | ND | 1.0 | 0.1 | | Pass |
| Paclobutrazol | ND | ND | 0.4 | 0.1 | | Pass |
| Permethrins | ND | ND | 0.2 | 0.1 | | Pass |
| Phosmet | ND | ND | 0.2 | 0.1 | | Pass |
| Piperonyl Butoxide | ND | ND | 2.0 | 1.0 | | Pass |
| Prallethrin | ND | ND | 0.2 | 0.1 | | Pass |
| Propiconazole | ND | ND | 0.4 | 0.1 | | Pass |
| Propoxur | ND | ND | 0.2 | 0.1 | | Pass |
| Pyrethrins | ND | ND | 1.0 | 0.5 | | Pass |
| Pyridaben | ND | ND | 0.2 | 0.1 | | Pass |
| Spinosad | ND | ND | 0.2 | 0.1 | | Pass |
| Spiromesifen | ND | ND | 0.2 | 0.1 | | Pass |
| Spirotetramat | ND | ND | 0.2 | 0.1 | | Pass |
| Spiroxamine | ND | ND | 0.4 | 0.1 | | Pass |
| Tebuconazole | ND | ND | 0.4 | 0.1 | | Pass |
| Thiacloprid | ND | ND | 0.2 | 0.1 | | Pass |
| Thiamethoxam | ND | ND | 0.2 | 0.1 | | Pass |
| Trifloxystrobin | ND | ND | 0.2 | 0.1 | | Pass |
| | | | | | | |



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Metrc Batch ID: 1A4010300016315000027948 Metrc Sample ID: 1A4010300016315000028008 Harvest/Process Date: 11/2/2020 Report ID: LS-201105-18 Sample Plan ID:SP-201104-3-A Sample Procedure: 160721_LAB-SOP_SampleCollection-v008



Pesticides Quality Control Data

| Analyte | Blank | LOQ | LCS | LCS Spike | LCS Rec (%) | Limits (%) | Notes |
|---------------------|-------|-----|-------|-----------|-------------|------------|-------|
| Abamectin | ND | 0.1 | 1.41 | 1.00 | 141 | 50 - 150 | |
| Acephate | ND | 0.1 | 1.11 | 1.00 | 111 | 50 - 150 | |
| Acequinocyl | ND | 1.5 | 1.87 | 1.00 | 187 | 50 - 150 | HB |
| Acetamiprid | ND | 0.1 | 1.03 | 1.00 | 103 | 50 - 150 | |
| Aldicarb | ND | 0.1 | 0.988 | 1.00 | 98.8 | 50 - 150 | |
| Azoxystrobin | ND | 0.1 | 1.33 | 1.00 | 133 | 50 - 150 | |
| Bifenazate | ND | 0.1 | 1.76 | 1.00 | 176 | 50 - 150 | HB |
| Bifenthrin | ND | 0.1 | 1.31 | 1.00 | 131 | 50 - 150 | |
| Boscalid | ND | 0.1 | 1.18 | 1.00 | 118 | 50 - 150 | |
| Carbaryl | ND | 0.1 | 0.939 | 1.00 | 93.9 | 50 - 150 | |
| Carbofuran | ND | 0.1 | 1.02 | 1.00 | 102 | 50 - 150 | |
| Chlorantraniliprole | ND | 0.1 | 1.30 | 1.00 | 130 | 50 - 150 | |
| Chlorfenapyr | ND | 0.1 | 1.21 | 1.00 | 121 | 50 - 150 | |
| Chlorpyrifos | ND | 0.1 | 1.02 | 1.00 | 102 | 50 - 150 | |
| Clofentezine | ND | 0.1 | 1.18 | 1.00 | 118 | 50 - 150 | |
| Cyfluthrin | ND | 0.5 | 1.46 | 1.00 | 146 | 50 - 150 | |
| Cypermethrin | ND | 0.1 | 1.43 | 1.00 | 143 | 50 - 150 | |
| Daminozide | ND | 0.5 | 0.569 | 1.00 | 56.9 | 10 - 150 | |
| Diazinon | ND | 0.1 | 1.09 | 1.00 | 109 | 50 - 150 | |
| Dichlorvos (DDVP) | ND | 0.5 | 0.917 | 1.00 | 91.7 | 50 - 150 | |
| Dimethoate | ND | 0.1 | 1.08 | 1.00 | 108 | 50 - 150 | |
| Ethoprophos | ND | 0.1 | 1.17 | 1.00 | 117 | 50 - 150 | |
| Etofenprox | ND | 0.1 | 1.29 | 1.00 | 129 | 50 - 150 | |
| Etoxazole | ND | 0.1 | 1.11 | 1.00 | 111 | 50 - 150 | |
| Fenoxycarb | ND | 0.1 | 1.30 | 1.00 | 130 | 50 - 150 | |
| Fenpyroximate | ND | 0.1 | 1.53 | 1.00 | 153 | 50 - 150 | |
| Fipronil | ND | 0.1 | 1.30 | 1.00 | 130 | 50 - 150 | |
| Flonicamid | ND | 0.1 | 1.04 | 1.00 | 104 | 50 - 150 | |
| Fludioxonil | ND | 0.1 | 1.24 | 1.00 | 124 | 50 - 150 | |
| Hexythiazox | ND | 0.1 | 0.631 | 1.00 | 63.1 | 50 - 150 | |
| Imazalil | ND | 0.1 | 1.32 | 1.00 | 132 | 50 - 150 | |
| Imidacloprid | ND | 0.1 | 1.09 | 1.00 | 109 | 50 - 150 | |
| Kresoxim-methyl | ND | 0.1 | 1.20 | 1.00 | 120 | 50 - 150 | |
| Malathion | ND | 0.1 | 1.10 | 1.00 | 110 | 50 - 150 | |
| | | | | | | | |

Pesticides QC Analysis Date: 11/4/2020 Pesticides QC Batch ID: PST_110420C Method: EN 15662 Unit: μg/g (ppm)

| Analyte | Blank | LOQ | LCS | LCS Spike | LCS Rec (%) | Limits (%) | Notes |
|--------------------|-------|-----|-------|-----------|-------------|------------|-------|
| Metalaxyl | ND | 0.1 | 1.23 | 1.00 | 123 | 50 - 150 | |
| Methiocarb | ND | 0.1 | 1.30 | 1.00 | 130 | 50 - 150 | |
| Methomyl | ND | 0.1 | 0.965 | 1.00 | 96.5 | 50 - 150 | |
| Methyl Parathion | ND | 0.2 | ND | 1.00 | 0.00 | 30 - 150 | ND |
| MGK-264 | ND | 0.2 | 0.798 | 0.600 | 133 | 50 - 150 | |
| Myclobutanil | ND | 0.1 | 1.36 | 1.00 | 136 | 50 - 150 | |
| Naled | ND | 0.2 | 1.03 | 1.00 | 103 | 50 - 150 | |
| 0xamy1 | ND | 0.1 | 0.953 | 1.00 | 95.3 | 50 - 150 | |
| Paclobutrazol | ND | 0.1 | 0.835 | 1.00 | 83.5 | 50 - 150 | |
| Permethrins | ND | 0.1 | 1.57 | 1.00 | 157 | 50 - 150 | ME |
| Phosmet | ND | 0.1 | 1.10 | 1.00 | 110 | 50 - 150 | |
| Piperonyl Butoxide | ND | 1.0 | 1.05 | 1.00 | 105 | 50 - 150 | |
| Prallethrin | ND | 0.1 | 1.13 | 1.00 | 113 | 50 - 150 | |
| Propiconazole | ND | 0.1 | 1.37 | 1.00 | 137 | 50 - 150 | |
| Propoxur | ND | 0.1 | 1.06 | 1.00 | 106 | 50 - 150 | |
| Pyrethrins | ND | 0.5 | 0.820 | 1.00 | 82.0 | 50 - 150 | |
| Pyridaben | ND | 0.1 | 1.23 | 1.00 | 123 | 50 - 150 | |
| Spinosad | ND | 0.1 | 1.17 | 1.00 | 117 | 50 - 150 | |
| Spiromesifen | ND | 0.1 | 1.12 | 1.00 | 112 | 50 - 150 | |
| Spirotetramat | ND | 0.1 | 1.33 | 1.00 | 133 | 50 - 150 | |
| Spiroxamine | ND | 0.1 | 1.24 | 1.00 | 124 | 50 - 150 | |
| Tebuconazole | ND | 0.1 | 1.49 | 1.00 | 149 | 50 - 150 | |
| Thiacloprid | ND | 0.1 | 1.01 | 1.00 | 101 | 50 - 150 | |
| Thiamethoxam | ND | 0.1 | 0.996 | 1.00 | 99.6 | 50 - 150 | |
| Trifloxystrobin | ND | 0.1 | 1.31 | 1.00 | 131 | 50 - 150 | |



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Sample Date: 11/3/2020
Analysis Date: 11/4/2020
Report Date: 11/5/2020

Metrc Batch ID: 1A4010300016315000027948 Metrc Sample ID: 1A4010300016315000028008 Harvest/Process Date: 11/2/2020 Report ID: LS-201105-18 Sample Plan ID:SP-201104-3-A Sample Procedure: 160721_LAB-SOP_SampleCollection-v008



Solvents Analysis Date: 11/5/2020 Solvents Batch ID: RES 110420A Method: EPA 5021A Unit: μg/g (ppm) Pass 🕢

| Analyte | TBH-XST-SHH | PSR-ZBM-DRG | 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 | ND | ND | 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 | ND | 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: 11/5/2020 Solvents QC Batch ID: RES_110420A Method: EPA 5021A Unit: μg/g (ppm)

| Analyte | Blank | LOQ | LCS | LCS Spike | LCS Rec (%) | Limits (%) | Notes |
|--------------------------|-------|-------|------|-----------|-------------|------------|-------|
| 1,4-Dioxane | ND | 50.0 | 1050 | 1000 | 105 | 70 - 130 | |
| 2-Butanol | ND | 250.0 | 876 | 1000 | 87.6 | 70 - 130 | |
| 2-Ethoxyethanol | ND | 50.0 | 1070 | 1000 | 107 | 70 - 130 | |
| Acetone | ND | 250.0 | 955 | 1000 | 95.5 | 70 - 130 | |
| Acetonitrile | ND | 50.0 | 932 | 1000 | 93.2 | 70 - 130 | |
| Benzene | ND | 2.0 | 21.7 | 20.0 | 109 | 70 - 130 | |
| Butanes | ND | 250.0 | 1610 | 2000 | 80.5 | 70 - 130 | |
| Cumene | ND | 50.0 | 1090 | 1000 | 109 | 70 - 130 | |
| Cyclohexane | ND | 50.0 | 962 | 1000 | 96.2 | 70 - 130 | |
| Ethyl Acetate | ND | 250.0 | 866 | 1000 | 86.6 | 70 - 130 | |
| Ethyl Ether | ND | 250.0 | 1050 | 1000 | 105 | 70 - 130 | |
| Ethylene Glycol | ND | 250.0 | 1060 | 1000 | 106 | 70 - 130 | |
| Ethylene Oxide | ND | 50.0 | 857 | 1000 | 85.7 | 70 - 130 | |
| Heptane | ND | 250.0 | 850 | 1000 | 85.0 | 70 - 130 | |
| Hexanes | ND | 50.0 | 4520 | 5000 | 90.4 | 70 - 130 | |
| Isopropanol (2-Propanol) | ND | 50.0 | 852 | 1000 | 85.2 | 70 - 130 | |
| Isopropyl Acetate | ND | 250.0 | 861 | 1000 | 86.1 | 70 - 130 | |
| Methanol | ND | 250.0 | 1040 | 1000 | 104 | 70 - 130 | |
| Dichloromethane | ND | 50.0 | 915 | 1000 | 91.5 | 70 - 130 | |
| Pentanes | ND | 250.0 | 2570 | 3000 | 85.7 | 70 - 130 | |
| Propane | ND | 250.0 | 865 | 1000 | 86.5 | 70 - 130 | |
| Tetrahydrofuran | ND | 50.0 | 901 | 1000 | 90.1 | 70 - 130 | |
| Toluene | ND | 50.0 | 972 | 1000 | 97.2 | 70 - 130 | |
| Xylenes | ND | 50.0 | 4030 | 4000 | 101 | 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 |
| В | Analyte found in sample and associated blank |
| С | 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 |
| М | Manual integration |
| PS | Peaks split |
| НВ | 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 |