LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Title : SVOA CALIBRATION

Signal : TIC: BP010063.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 4.340 | 208 | 213 | 226 | rVB | 2295582 | 3708059 | 0.93% | 0.173% |
| 2 | 5.534 | 409 | 416 | 429 | rVB | 3861220 | 6073388 | 1.52% | 0.284% |
| 3 | 6.622 | 586 | 601 | 610 | rBV | 2342248 | 4292322 | 1.07% | 0.201% |
| 4 | 7.110 | 677 | 684 | 696 | rBV | 773810 | 1520016 | 0.38% | 0.071% |
| 5 | 8.552 | 917 | 929 | 941 | rBV2 | 1061489 | 2553829 | 0.64% | 0.119% |
| 6 | 8.787 | 960 | 969 | 971 | rBV3 | 724659 | 1696519 | 0.42% | 0.079% |
| 7 | 8.822 | 972 | 975 | 987 | rVB | 1206295 | 2134268 | 0.53% | 0.100% |
| 8 | 9.369 | 1030 | 1068 | 1071 | rBV2 | 24823895 | 193427050 | 48.40% | 9.037% |
| 9 | 9.469 | 1077 | 1085 | 1092 | rVB | 10872829 | 18889165 | 4.73% | 0.882% |
| 10 | 9.534 | 1092 | 1096 | 1112 | rVB2 | 867304 | 1688241 | 0.42% | 0.079% |
| 11 | 9.775 | 1128 | 1137 | 1144 | rBV2 | 19898107 | 36759028 | 9.20% | 1.717% |
| 12 | 9.881 | 1152 | 1155 | 1158 | rVV | 1144160 | 1668748 | 0.42% | 0.078% |
| 13 | 10.034 | 1170 | 1181 | 1191 | rBV | 5062591 | 14892103 | 3.73% | 0.696% |
| 14 | 10.181 | 1200 | 1206 | 1214 | rVB2 | 2135202 | 4352168 | 1.09% | 0.203% |
| 15 | 10.322 | 1214 | 1230 | 1236 | rBV | 12305621 | 37456251 | 9.37% | 1.750% |
| 16 | 10.381 | 1237 | 1240 | 1257 | rVB | 1441757 | 3283599 | 0.82% | 0.153% |
| 17 | 10.575 | 1264 | 1273 | 1280 | rBV | 7618210 | 17049677 | 4.27% | 0.797% |
| 18 | 10.757 | 1297 | 1304 | 1309 | rBV2 | 1371929 | 2338023 | 0.59% | 0.109% |
| 19 | 11.128 | 1356 | 1367 | 1372 | rBV | 19610969 | 40448234 | 10.12% | 1.890% |
| 20 | 11.216 | 1375 | 1382 | 1387 | rBV | 2337451 | 3831638 | 0.96% | 0.179% |
| 21 | 11.281 | 1387 | 1393 | 1405 | rVV | 2501765 | 6414370 | 1.61% | 0.300% |
| 22 | 11.398 | 1405 | 1413 | 1426 | rVV | 7727711 | 14540515 | 3.64% | 0.679% |
| 23 | 11.522 | 1426 | 1434 | 1443 | rVB | 3424157 | 5745431 | 1.44% | 0.268% |
| 24 | 12.151 | 1533 | 1541 | 1544 | rVV | 3561630 | 6348852 | 1.59% | 0.297% |
| 25 | 12.193 | 1544 | 1548 | 1557 | rVB2 | 3875642 | 6228352 | 1.56% | 0.291% |
| 26 | 12.410 | 1569 | 1585 | 1589 | rBV | 19582155 | 39232921 | 9.82% | 1.833% |
| 27 | 12.469 | 1589 | 1595 | 1601 | rVV | 15225626 | 26301999 | 6.58% | 1.229% |
| 28 | 12.557 | 1601 | 1610 | 1615 | rVV | 1130498 | 2188319 | 0.55% | 0.102% |
| 29 | 12.810 | 1645 | 1653 | 1660 | rBV3 | 1350654 | 3054138 | 0.76% | 0.143% |
| 30 | 12.951 | 1669 | 1677 | 1684 | rBV | 1655511 | 3069316 | 0.77% | 0.143% |
| 31 | 13.051 | 1684 | 1694 | 1699 | rVV | 5196187 | 8324472 | 2.08% | 0.389% |
| 32 | 13.322 | 1729 | 1740 | 1751 | rVB | 13570738 | 33754755 | 8.45% | 1.577% |
| 33 | 13.598 | 1765 | 1787 | 1791 | rBV | 40406224 | 203422212 | 50.91% | 9.504% |
| 34 | 13.710 | 1801 | 1806 | 1810 | rBV2 | 1201812 | 1662695 | 0.42% | 0.078% |
| 35 | 13.757 | 1810 | 1814 | 1819 | rVV2 | 1096553 | 2324517 | 0.58% | 0.109% |
| 36 | 13.804 | 1819 | 1822 | 1829 | rVB | 1386156 | 1853914 | 0.46% | 0.087% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 13.992 | 1848 | 1854 | 1860 | rBV | 4615364 | 6589443 | 1.65% | 0.308% |
| 38 | 14.151 | 1873 | 1881 | 1889 | rVB | 2930431 | 4536743 | 1.14% | 0.212% |
| 39 | 14.281 | 1898 | 1903 | 1907 | rVV | 1869125 | 3070419 | 0.77% | 0.143% |
| 40 | 14.334 | 1907 | 1912 | 1920 | rVB | 3568207 | 6027004 | 1.51% | 0.282% |
| 41 | 14.463 | 1926 | 1934 | 1939 | rVV2 | 1770538 | 3219549 | 0.81% | 0.150% |
| 42 | 14.534 | 1939 | 1946 | 1950 | rVV | 11065458 | 17144438 | 4.29% | 0.801% |
| 43 | 14.581 | 1950 | 1954 | 1959 | rVV | 1479067 | 2350845 | 0.59% | 0.110% |
| 44 | 14.634 | 1959 | 1963 | 1968 | rVB | 1917508 | 2628984 | 0.66% | 0.123% |
| 45 | 14.786 | 1984 | 1989 | 1995 | rVV4 | 963972 | 1954082 | 0.49% | 0.091% |
| 46 | 15.569 | 2112 | 2122 | 2128 | rBV | 2562498 | 4328993 | 1.08% | 0.202% |
| 47 | 15.892 | 2166 | 2177 | 2188 | rVB | 3637578 | 6307723 | 1.58% | 0.295% |
| 48 | 15.992 | 2188 | 2194 | 2197 | rBV | 3823214 | 5659883 | 1.42% | 0.264% |
| 49 | 16.033 | 2197 | 2201 | 2214 | rVB4 | 4182176 | 10546020 | 2.64% | 0.493% |
| 50 | 16.145 | 2215 | 2220 | 2221 | rBV | 1652499 | 2192448 | 0.55% | 0.102% |
| 51 | 16.245 | 2234 | 2237 | 2241 | rVV2 | 2201271 | 3102752 | 0.78% | 0.145% |
| 52 | 16.286 | 2241 | 2244 | 2249 | rVB | 3793037 | 5385827 | 1.35% | 0.252% |
| 53 | 16.333 | 2249 | 2252 | 2257 | rVB | 1991918 | 2564637 | 0.64% | 0.120% |
| 54 | 16.545 | 2281 | 2288 | 2294 | rBV4 | 771060 | 1795667 | 0.45% | 0.084% |
| 55 | 16.739 | 2312 | 2321 | 2322 | rBV2 | 929919 | 2248122 | 0.56% | 0.105% |
| 56 | 16.810 | 2327 | 2333 | 2344 | rVB | 4930490 | 9389354 | 2.35% | 0.439% |
| 57 | 17.133 | 2345 | 2388 | 2389 | rBV3 | 39692096 | 399606566 | 100.00% | 18.669% |
| 58 | 17.269 | 2407 | 2411 | 2415 | rVB | 1423572 | 1898333 | 0.48% | 0.089% |
| 59 | 17.363 | 2421 | 2427 | 2431 | rBV | 4208567 | 6260193 | 1.57% | 0.292% |
| 60 | 17.457 | 2438 | 2443 | 2447 | rBV | 2481807 | 3960457 | 0.99% | 0.185% |
| 61 | 17.657 | 2458 | 2477 | 2478 | rBV2 | 43530488 | 201099963 | 50.32% | 9.395% |
| 62 | 17.710 | 2483 | 2486 | 2491 | rVB | 3914190 | 5224438 | 1.31% | 0.244% |
| 63 | 18.310 | 2583 | 2588 | 2595 | rBV | 2863786 | 4392850 | 1.10% | 0.205% |
| 64 | 18.480 | 2604 | 2617 | 2621 | rVB | 12888750 | 36200323 | 9.06% | 1.691% |
| 65 | 18.933 | 2691 | 2694 | 2699 | rVB | 1400170 | 1899411 | 0.48% | 0.089% |
| 66 | 19.074 | 2713 | 2718 | 2728 | rBV | 6572018 | 14454063 | 3.62% | 0.675% |
| 67 | 19.216 | 2736 | 2742 | 2746 | rVB3 | 872614 | 1625099 | 0.41% | 0.076% |
| 68 | 19.522 | 2790 | 2794 | 2797 | rBV | 1249650 | 1718783 | 0.43% | 0.080% |
| 69 | 19.669 | 2811 | 2819 | 2823 | rBV | 2732224 | 4265698 | 1.07% | 0.199% |
| 70 | 19.722 | 2823 | 2828 | 2832 | rBV2 | 1971420 | 3583857 | 0.90% | 0.167% |
| 71 | 19.910 | 2852 | 2860 | 2868 | rVB | 5633668 | 11648514 | 2.91% | 0.544% |
| 72 | 20.063 | 2878 | 2886 | 2893 | rBV | 2705921 | 5178376 | 1.30% | 0.242% |
| 73 | 20.204 | 2905 | 2910 | 2916 | rBV | 3039682 | 4189494 | 1.05% | 0.196% |
| 74 | 20.333 | 2927 | 2932 | 2939 | rBV4 | 2973404 | 5248118 | 1.31% | 0.245% |
| 75 | 20.504 | 2956 | 2961 | 2965 | rBV | 4982824 | 6386994 | 1.60% | 0.298% |
| 76 | 20.639 | 2976 | 2984 | 2987 | rBV | 2031808 | 3878640 | 0.97% | 0.181% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 77 | 20.721 | 2989 | 2998 | 3004 | rVV2 | 32827739 | 73380004 | 18.36% | 3.428% |
| 78 | 20.774 | 3004 | 3007 | 3017 | rVB | 1629898 | 2475280 | 0.62% | 0.116% |
| 79 | 21.063 | 3051 | 3056 | 3065 | rVV2 | 2622916 | 5749323 | 1.44% | 0.269% |
| 80 | 21.145 | 3067 | 3070 | 3074 | rVV | 1774703 | 2561728 | 0.64% | 0.120% |
| 81 | 21.210 | 3075 | 3081 | 3086 | rVV | 5774660 | 10191967 | 2.55% | 0.476% |
| 82 | 21.280 | 3086 | 3093 | 3096 | rVV2 | 17946855 | 33181311 | 8.30% | 1.550% |
| 83 | 21.321 | 3096 | 3100 | 3102 | rVV | 19190411 | 26251875 | 6.57% | 1.226% |
| 84 | 21.345 | 3102 | 3104 | 3108 | rVV | 11244867 | 13338145 | 3.34% | 0.623% |
| 85 | 21.398 | 3108 | 3113 | 3124 | rVB2 | 16461408 | 23879779 | 5.98% | 1.116% |
| 86 | 21.580 | 3140 | 3144 | 3151 | rBV2 | 1464086 | 3276291 | 0.82% | 0.153% |
| 87 | 21.710 | 3161 | 3166 | 3173 | rBV | 6477118 | 10404107 | 2.60% | 0.486% |
| 88 | 21.927 | 3197 | 3203 | 3206 | rBV2 | 2837040 | 5632260 | 1.41% | 0.263% |
| 89 | 21.968 | 3207 | 3210 | 3221 | rVB2 | 3228178 | 9201033 | 2.30% | 0.430% |
| 90 | 22.157 | 3231 | 3242 | 3251 | rVB4 | 8707146 | 28609355 | 7.16% | 1.337% |
| 91 | 22.263 | 3252 | 3260 | 3266 | rVB | 4144203 | 11773188 | 2.95% | 0.550% |
| 92 | 22.392 | 3267 | 3282 | 3283 | rBV | 33673998 | 106483441 | 26.65% | 4.975% |
| 93 | 23.192 | 3413 | 3418 | 3425 | rVB | 2187838 | 3733340 | 0.93% | 0.174% |
| 94 | 23.615 | 3483 | 3490 | 3497 | rVB | 5822183 | 11587318 | 2.90% | 0.541% |
| 95 | 24.198 | 3577 | 3589 | 3595 | rBV | 13750661 | 37838701 | 9.47% | 1.768% |
| 96 | 24.392 | 3614 | 3622 | 3632 | rVB | 6813394 | 15864469 | 3.97% | 0.741% |
| 97 | 25.533 | 3810 | 3816 | 3830 | rVB | 2869307 | 7549488 | 1.89% | 0.353% |
| 98 | 26.927 | 4045 | 4053 | 4060 | rBV4 | 1235203 | 4616551 | 1.16% | 0.216% |
| 99 | 27.533 | 4138 | 4156 | 4170 | rVB | 7634010 | 36908933 | 9.24% | 1.724% |
| 100 | 27.827 | 4181 | 4206 | 4214 | rBV2 | 12189170 | 69710103 | 17.44% | 3.257% |

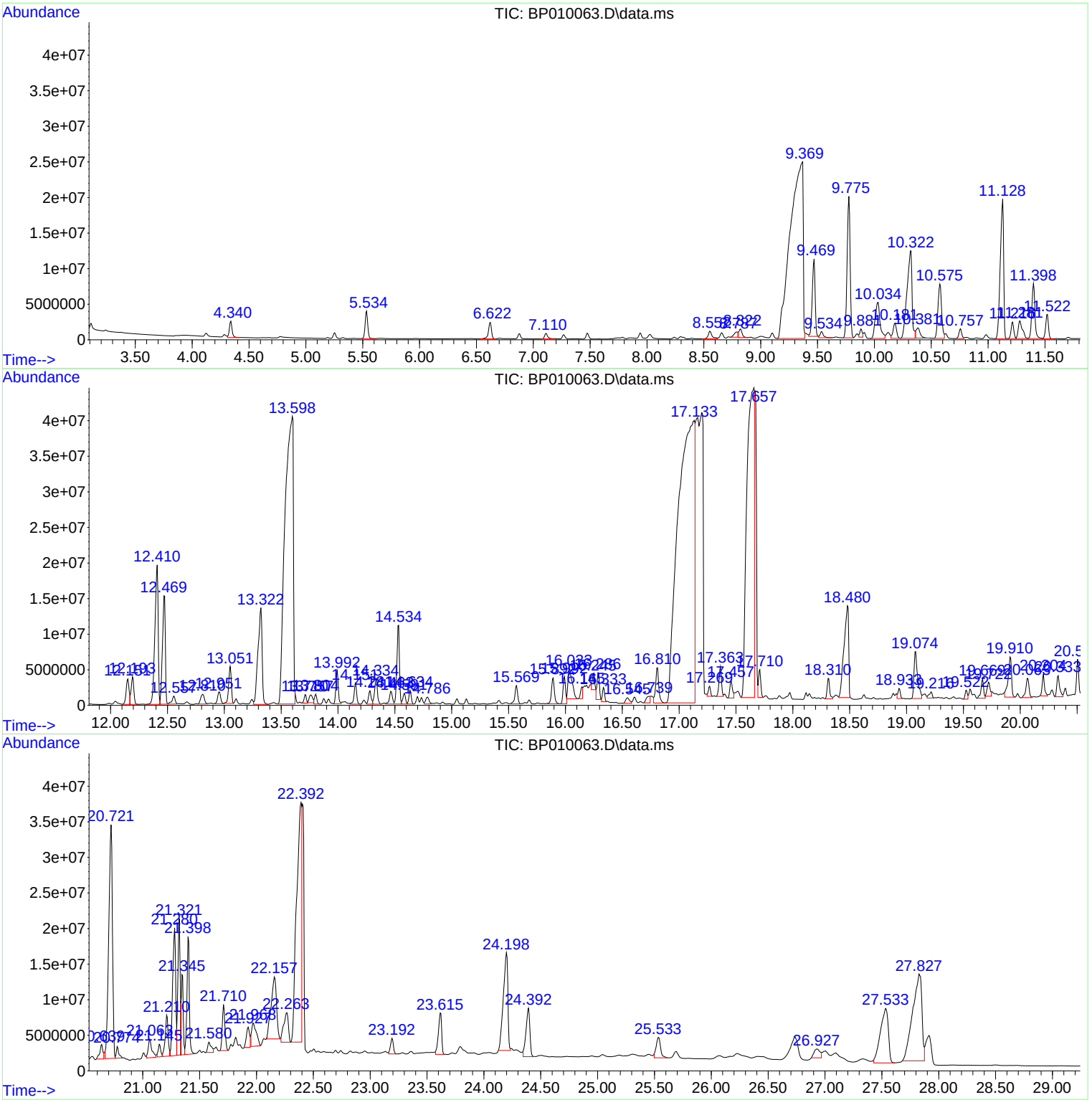
Sum of corrected areas: 2140490197

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

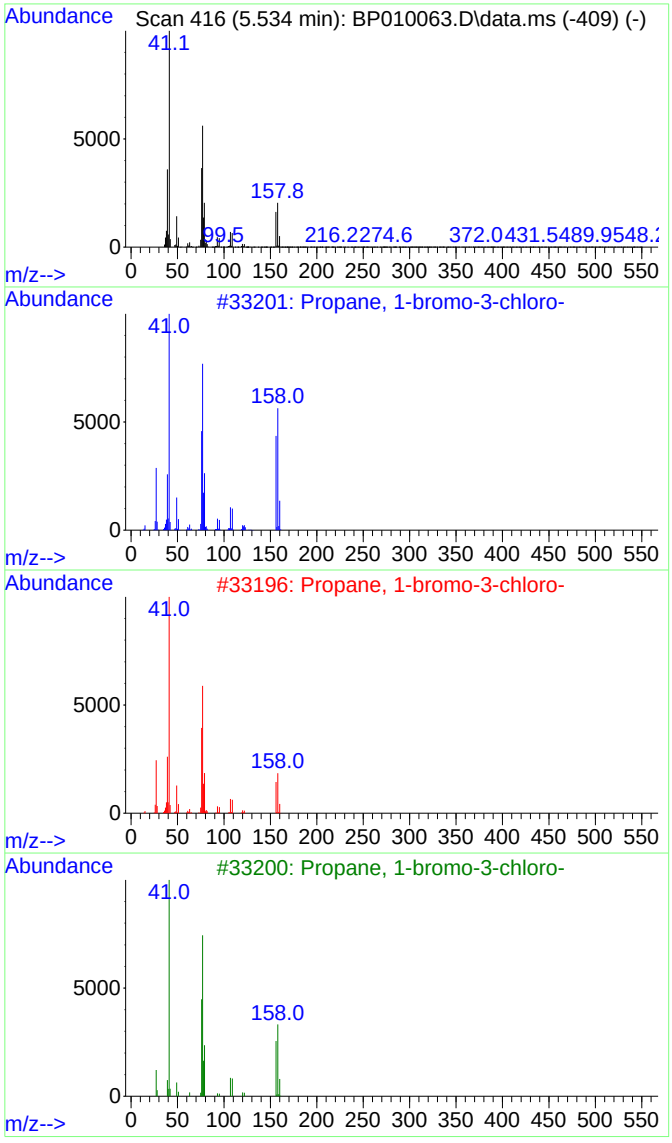
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

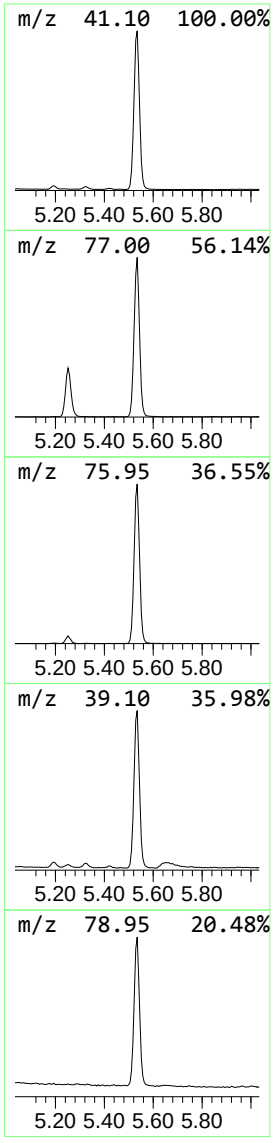
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 2 Propane, 1-bromo-3-chloro- Concentration Rank 28

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.534 | 47.56 ng/ul | 6073390 | 1,4-Dichlorobenzene-d4 | 7.940 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 1-bromo-3-chloro- | 156 | C3H6BrCl | 000109-70-6 | 97 |
| 2 | Propane, 1-bromo-3-chloro- | 156 | C3H6BrCl | 000109-70-6 | 97 |
| 3 | Propane, 1-bromo-3-chloro- | 156 | C3H6BrCl | 000109-70-6 | 94 |
| 4 | Propane, 1-bromo-3-chloro- | 156 | C3H6BrCl | 000109-70-6 | 76 |
| 5 | Acetic acid, chloro-, anhydride | 170 | C4H4Cl2O3 | 000541-88-8 | 43 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

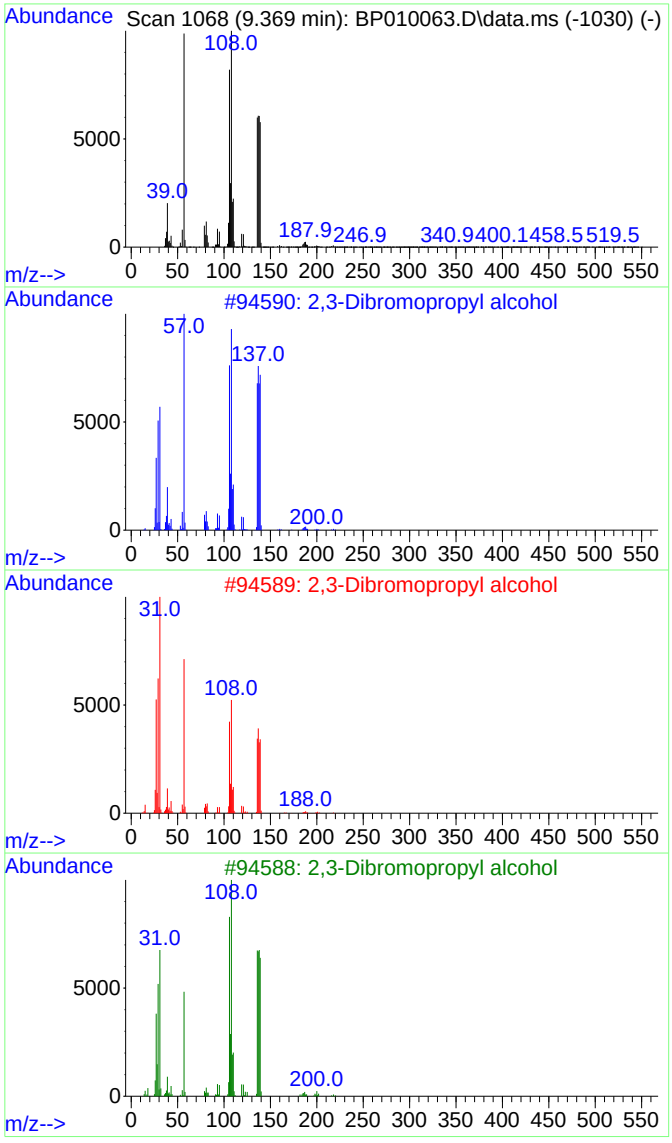
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

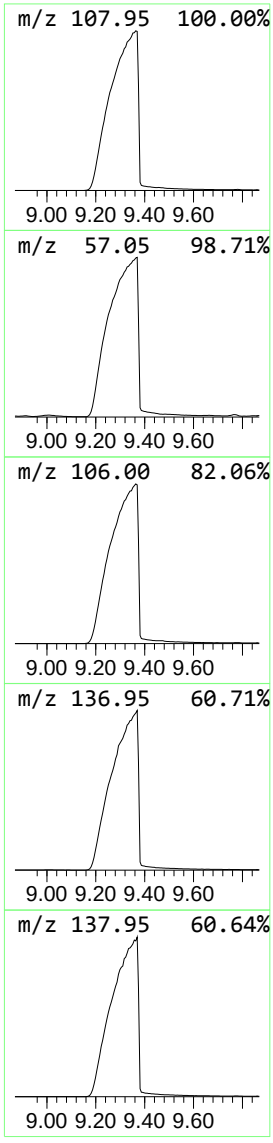
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 7 2,3-Dibromopropyl alcohol Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.369 | 1654.62 ng/ul | 193427000 | Naphthalene-d8 | 10.757 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2,3-Dibromopropyl alcohol | 216 | C3H6Br2O | 000096-13-9 | 99 |
| 2 | 2,3-Dibromopropyl alcohol | 216 | C3H6Br2O | 000096-13-9 | 97 |
| 3 | 2,3-Dibromopropyl alcohol | 216 | C3H6Br2O | 000096-13-9 | 95 |
| 4 | Vinyl bromide | 106 | C2H3Br | 000593-60-2 | 35 |
| 5 | 4,5,6,7-Tetrahydro-3-methylbenzo... | 136 | C9H12O | 1000426-71-8 | 25 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

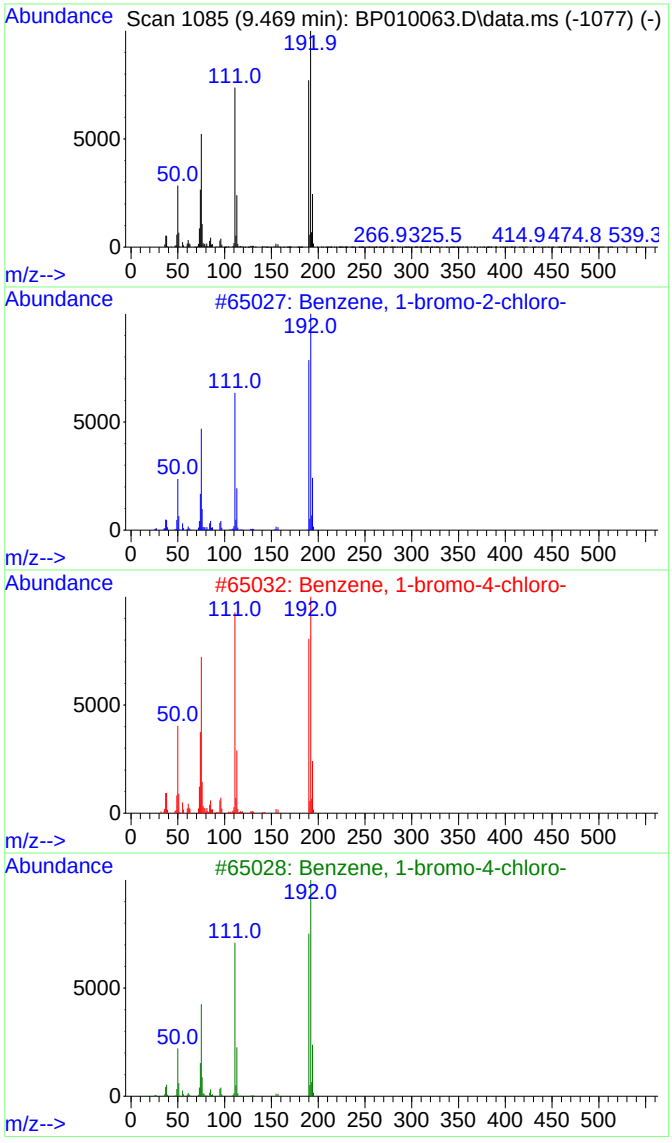
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

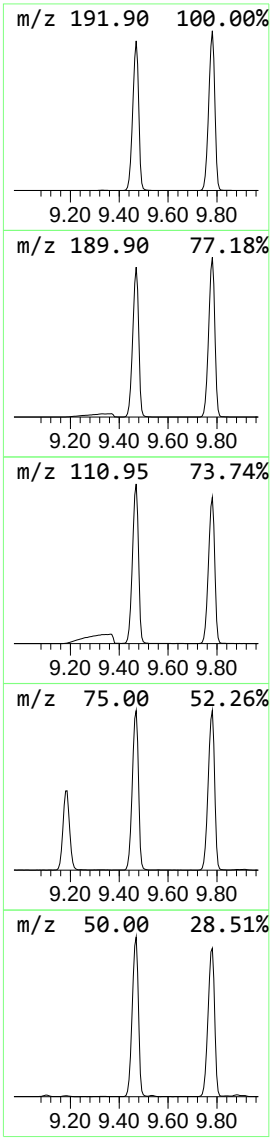
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 8 Benzene, 1-bromo-2-chloro- Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.469 | 161.58 ng/ul | 18889200 | Naphthalene-d8 | 10.757 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1-bromo-2-chloro- | 190 | C6H4BrCl | 000694-80-4 | 98 |
| 2 | Benzene, 1-bromo-4-chloro- | 190 | C6H4BrCl | 000106-39-8 | 98 |
| 3 | Benzene, 1-bromo-4-chloro- | 190 | C6H4BrCl | 000106-39-8 | 98 |
| 4 | Benzene, 1-bromo-4-chloro- | 190 | C6H4BrCl | 000106-39-8 | 96 |
| 5 | 4-Chlorophenylselenol | 192 | C6H5ClSe | 016645-10-6 | 96 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

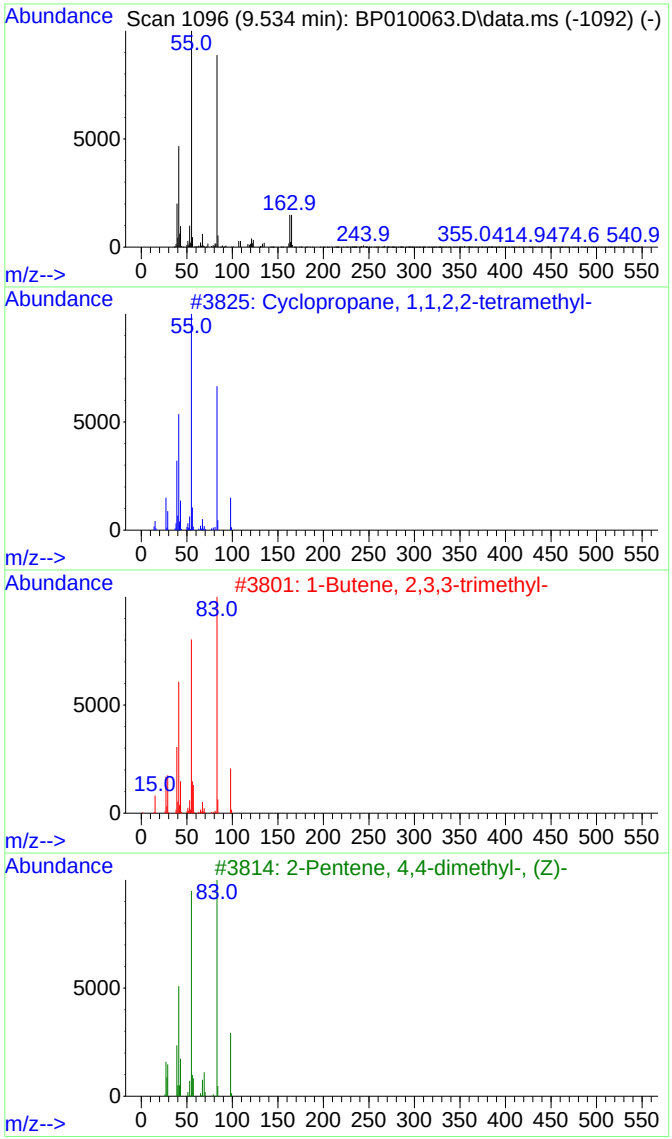
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

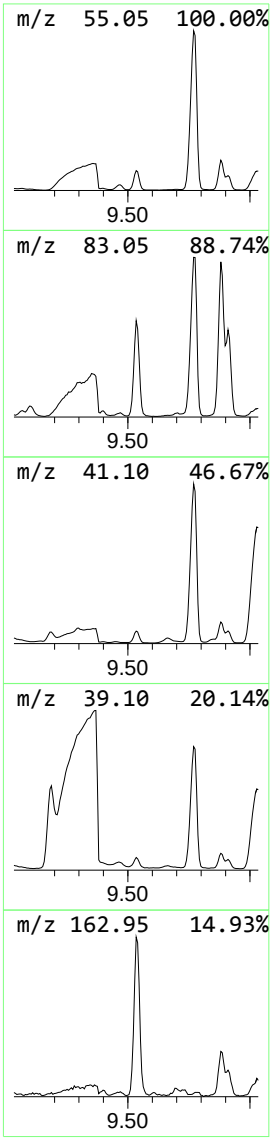
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 9 (DEL) Alkane: Cyclic9.534 Concentration Rank 54

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.534 | 14.44 ng/ul | 1688240 | Naphthalene-d8 | 10.757 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclopropane, 1,1,2,2-tetramethyl- | 98 | C7H14 | 004127-47-3 | 59 |
| 2 | 1-Butene, 2,3,3-trimethyl- | 98 | C7H14 | 000594-56-9 | 59 |
| 3 | 2-Pentene, 4,4-dimethyl-, (Z)- | 98 | C7H14 | 000762-63-0 | 59 |
| 4 | Cyclopropane, 2-bromo-1,1,3-trim... | 162 | C6H11Br | 036617-00-2 | 59 |
| 5 | 2-Methyl-2-bromohept-6-en-3-one | 204 | C8H13BrO | 1000424-28-0 | 59 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

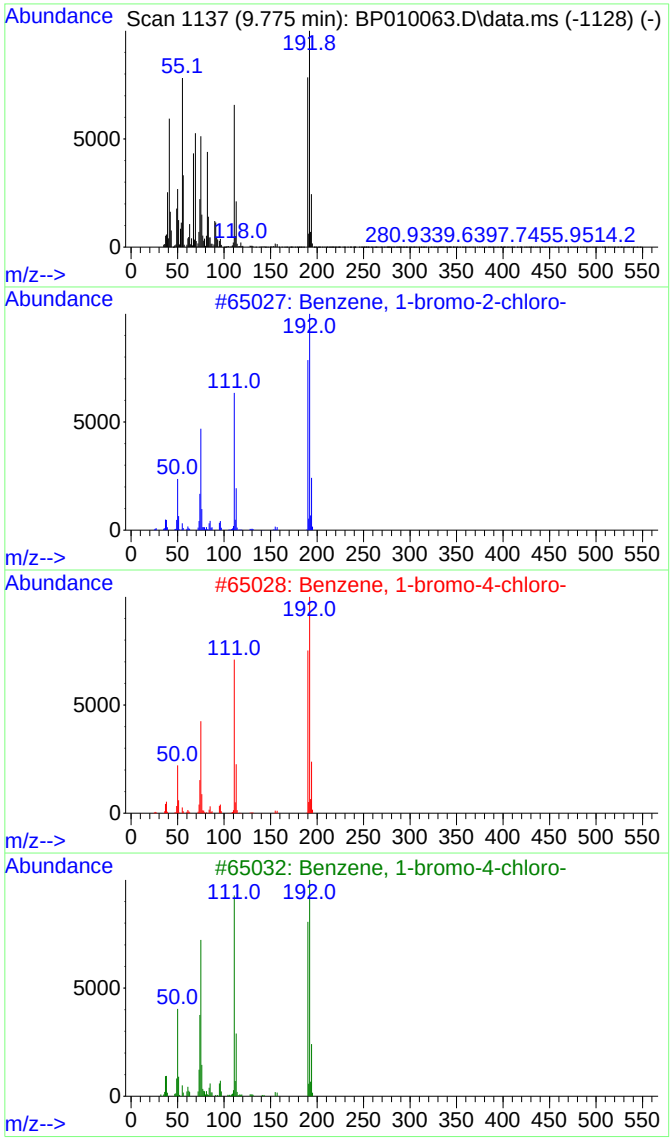
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

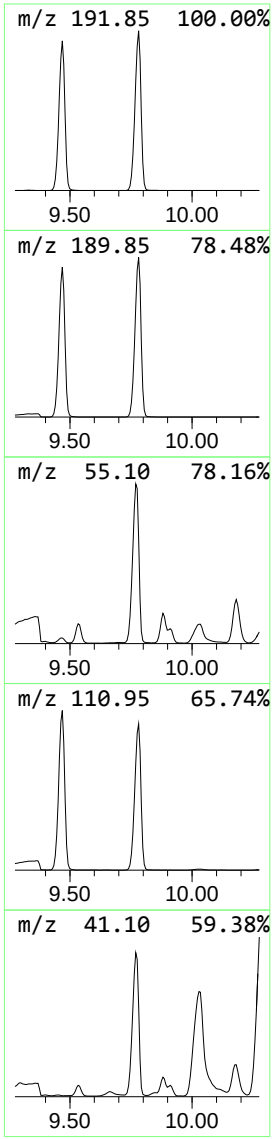
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 10 Benzene, 1-bromo-4-chloro- Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.775 | 314.44 ng/ul | 36759000 | Naphthalene-d8 | 10.757 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1-bromo-2-chloro- | 190 | C6H4BrCl | 000694-80-4 | 98 |
| 2 | Benzene, 1-bromo-4-chloro- | 190 | C6H4BrCl | 000106-39-8 | 98 |
| 3 | Benzene, 1-bromo-4-chloro- | 190 | C6H4BrCl | 000106-39-8 | 97 |
| 4 | 4-Chlorophenylselenol | 192 | C6H5ClSe | 016645-10-6 | 94 |
| 5 | Benzene, 1-bromo-3-chloro- | 190 | C6H4BrCl | 000108-37-2 | 93 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

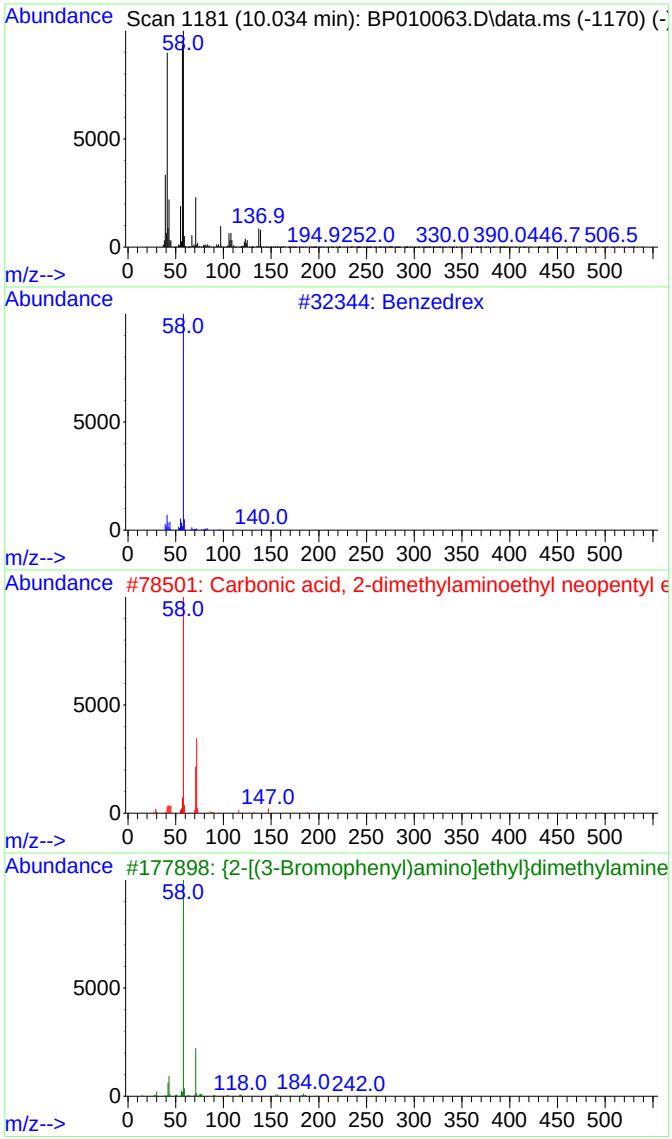
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

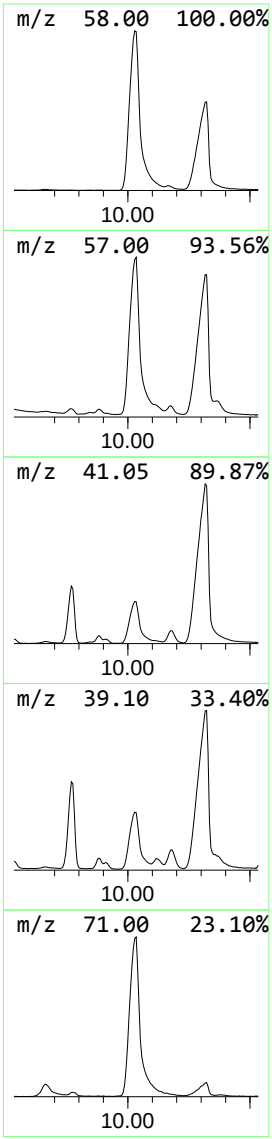
**********

Peak Number 11 unknown-01 Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.034 | 127.39 ng/ul | 14892100 | Naphthalene-d8 | 10.757 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzedrex | 155 | C10H21N | 000101-40-6 | 43 |
| 2 | Carbonic acid, 2-dimethylaminoet... | 203 | C10H21NO3 | 1000331-43-1 | 40 |
| 3 | {2-[(3-Bromophenyl)amino]ethyl}d... | 284 | C12H17BrN2O | 1010506-64-8 | 38 |
| 4 | Phenol, 4-(2-aminopropoxy)-3,5-d... | 195 | C11H17NO2 | 053566-99-7 | 38 |
| 5 | 1,2-Bis(2-dimethylaminoethoxy)et... | 204 | C10H24N2O2 | 003065-46-1 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

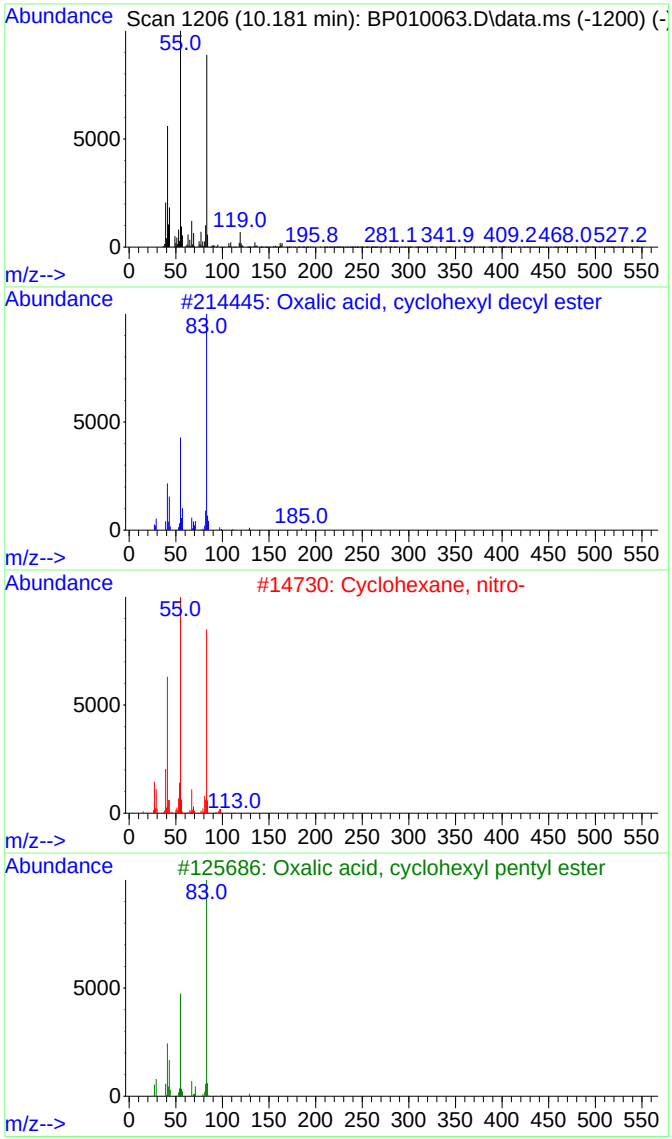
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

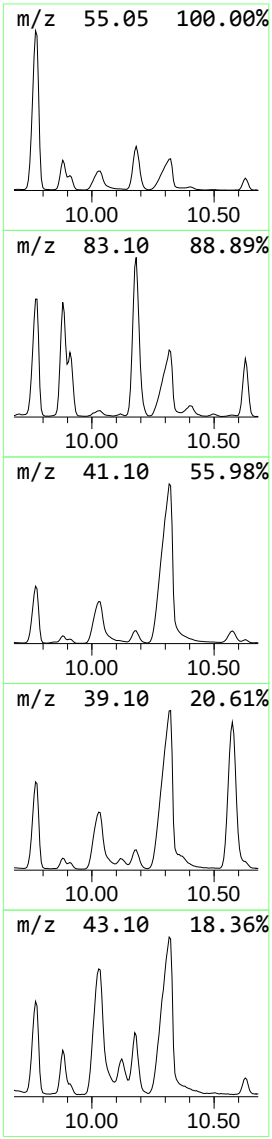
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 12 Oxalic acid, cyclohexyl dec... Concentration Rank 31

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.181 | 37.23 ng/ul | 4352170 | Naphthalene-d8 | 10.757 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Oxalic acid, cyclohexyl decyl ester | 312 | C18H32O4 | 1000309-31-2 | 78 |
| 2 | Cyclohexane, nitro- | 129 | C6H11NO2 | 001122-60-7 | 78 |
| 3 | Oxalic acid, cyclohexyl pentyl e... | 242 | C13H22O4 | 1000309-30-6 | 78 |
| 4 | Oxalic acid, cyclohexyl octyl ester | 284 | C16H28O4 | 1000309-31-0 | 78 |
| 5 | Acetaldehyde 2E-hexenyl 3Z-hexen... | 226 | C14H26O2 | 1000431-45-5 | 78 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

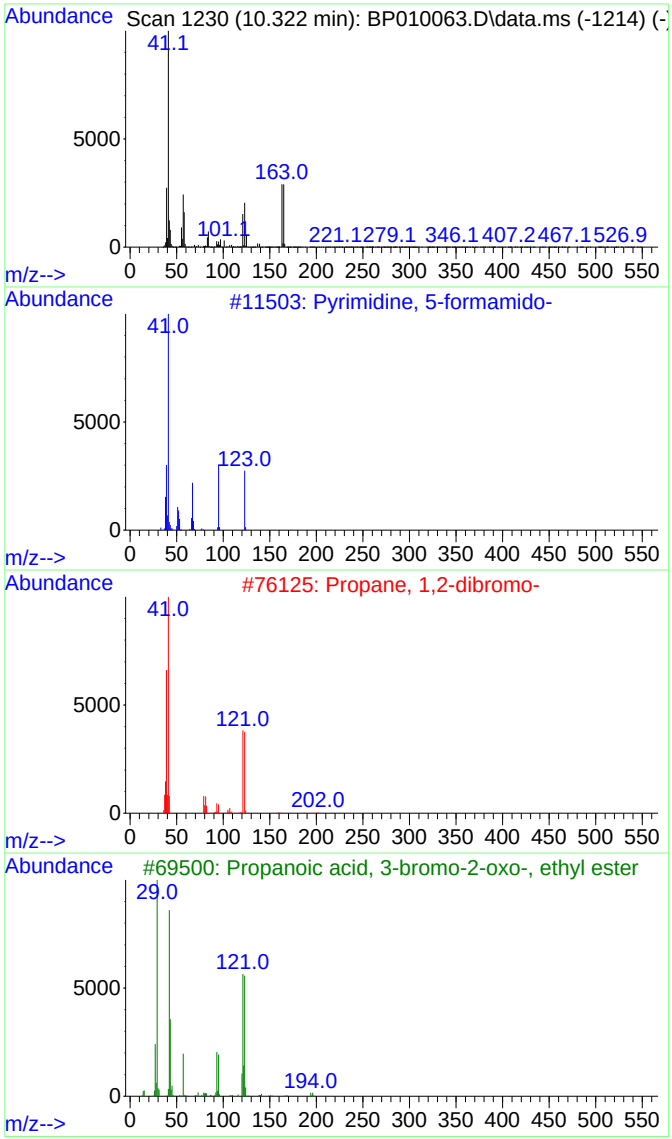
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

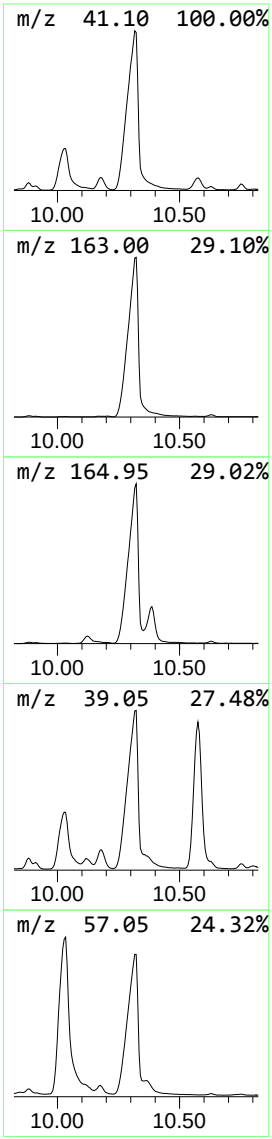
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 13 unknown-02 Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.322 | 320.41 ng/ul | 37456300 | Naphthalene-d8 | 10.757 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pyrimidine, 5-formamido- | 123 | C5H5N3O | 056621-84-2 | 9 |
| 2 | Propane, 1,2-dibromo- | 200 | C3H6Br2 | 000078-75-1 | 9 |
| 3 | Propanoic acid, 3-bromo-2-oxo-, ... | 194 | C5H7BrO3 | 000070-23-5 | 9 |
| 4 | 1-Propene, 3-(ethenyloxy)- | 84 | C5H8O | 003917-15-5 | 7 |
| 5 | o-Allylhydroxylamine | 73 | C3H7NO | 006542-54-7 | 5 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

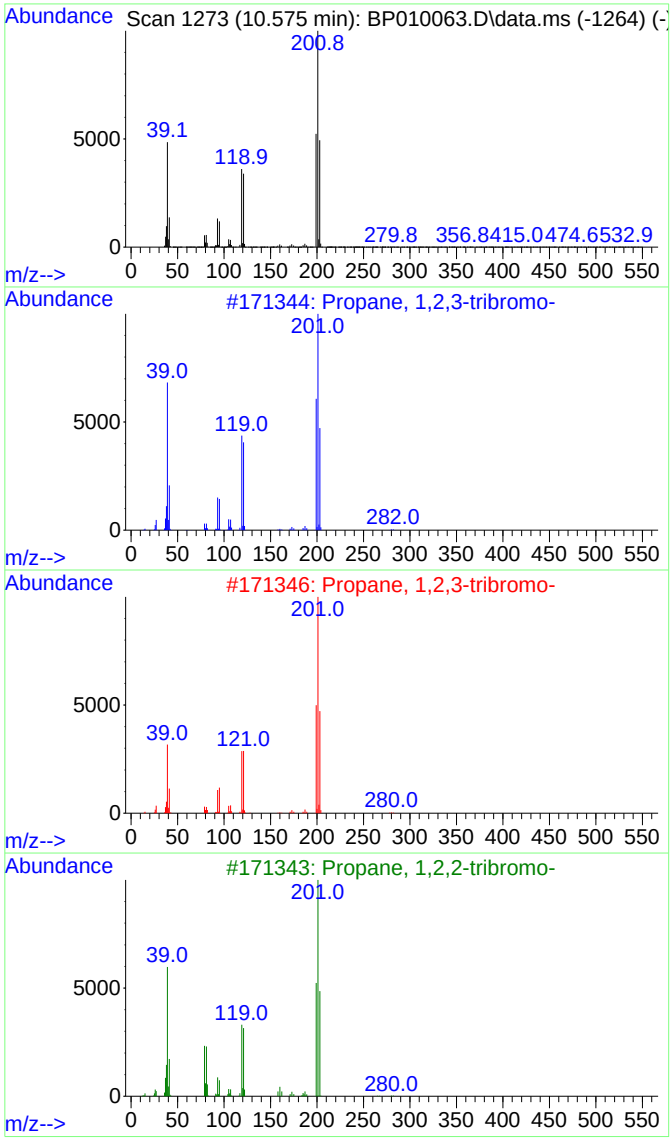
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

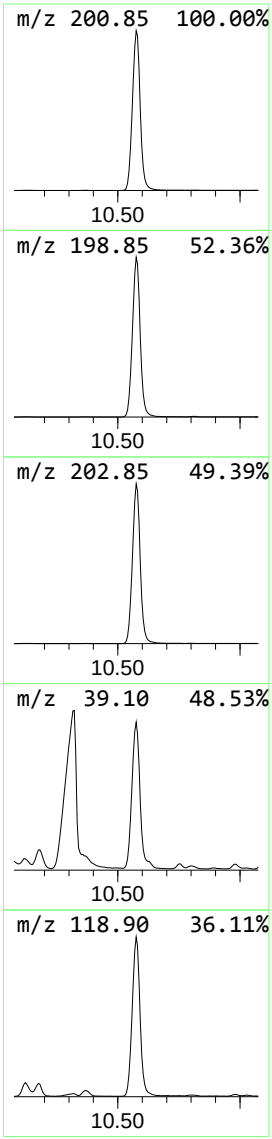
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 14 Propane, 1,2,3-tribromo- Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.575 | 145.85 ng/ul | 17049700 | Naphthalene-d8 | 10.757 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 96 |
| 2 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 96 |
| 3 | Propane, 1,2,2-tribromo- | 278 | C3H5Br3 | 014476-30-3 | 38 |
| 4 | 4-Bromobenzaldoxime | 199 | C7H6BrNO | 034158-73-1 | 9 |
| 5 | 4-(Difluoromethoxy)-3-methoxyben... | 232 | C9H10F2N2O3 | 446267-81-8 | 7 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

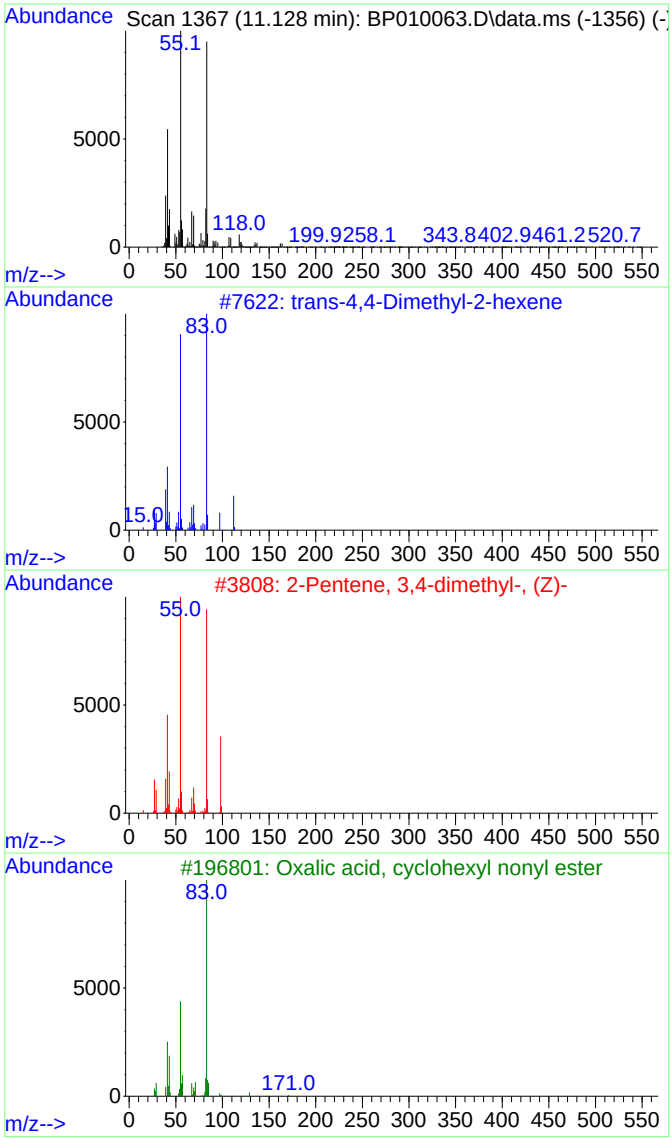
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

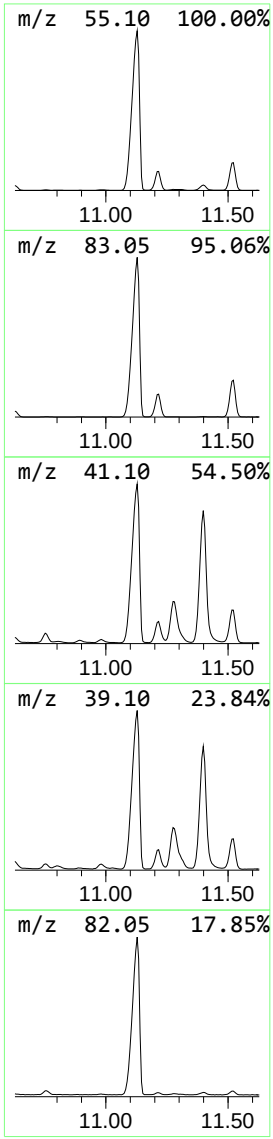
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.128 | 346.00 ng/ul | 40448200 | Naphthalene-d8 | 10.757 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | trans-4,4-Dimethyl-2-hexene | 112 | C8H16 | 019550-83-5 | 64 |
| 2 | 2-Pentene, 3,4-dimethyl-, (Z)- | 98 | C7H14 | 004914-91-4 | 64 |
| 3 | Oxalic acid, cyclohexyl nonyl ester | 298 | C17H30O4 | 1000309-31-1 | 64 |
| 4 | 2-Pentene, 4,4-dimethyl-, (E)- | 98 | C7H14 | 000690-08-4 | 64 |
| 5 | 3,3-Dimethylhex-5-en-2-one | 126 | C8H14O | 1000424-30-3 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

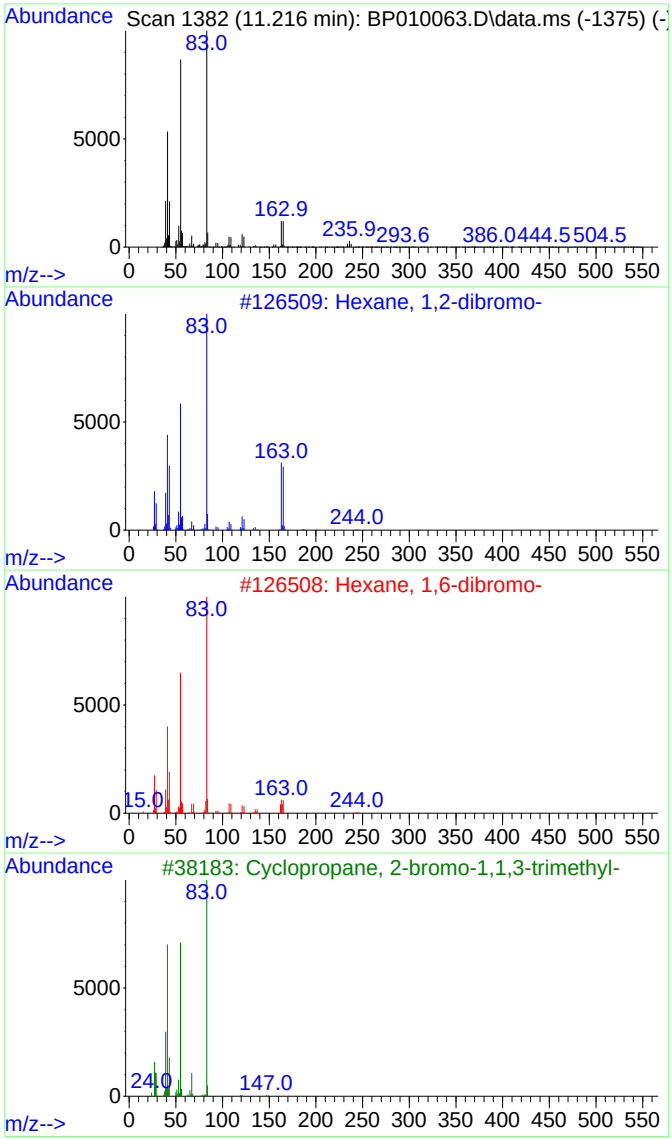
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

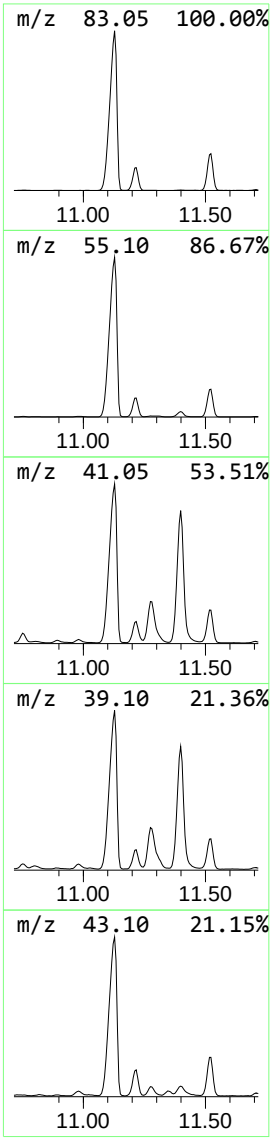
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 16 Hexane, 1,2-dibromo- Concentration Rank 34

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.216 | 32.78 ng/ul | 3831640 | Naphthalene-d8 | 10.757 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexane, 1,2-dibromo- | 242 | C6H12Br2 | 000624-20-4 | 90 |
| 2 | Hexane, 1,6-dibromo- | 242 | C6H12Br2 | 000629-03-8 | 72 |
| 3 | Cyclopropane, 2-bromo-1,1,3-trim... | 162 | C6H11Br | 036617-00-2 | 53 |
| 4 | Oxalic acid, cyclohexyl propyl e... | 214 | C11H18O4 | 1000309-30-3 | 53 |
| 5 | 1-Pentyn-3-ol | 84 | C5H8O | 004187-86-4 | 53 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

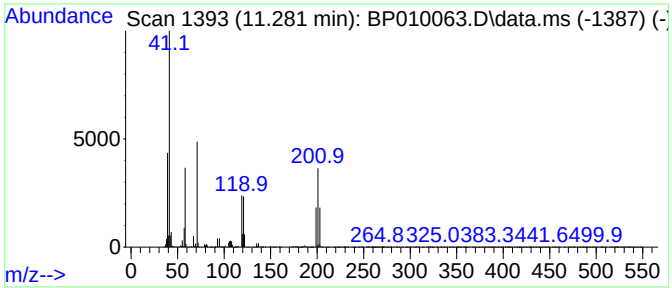
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

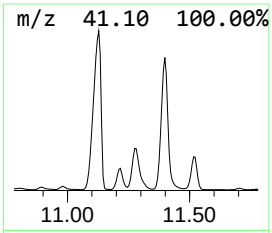
**********

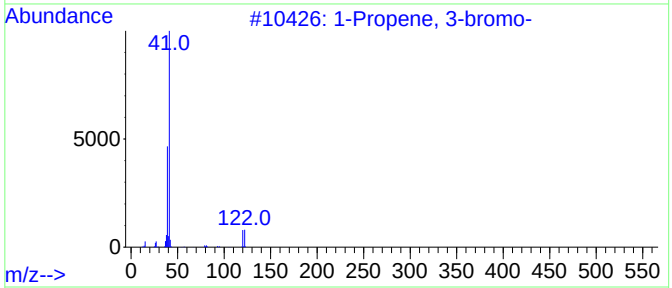
Peak Number 17 unknown-03 Concentration Rank 22

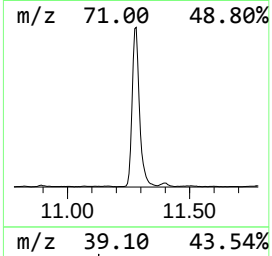
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.281 | 54.87 ng/ul | 6414370 | Naphthalene-d8 | 10.757 |

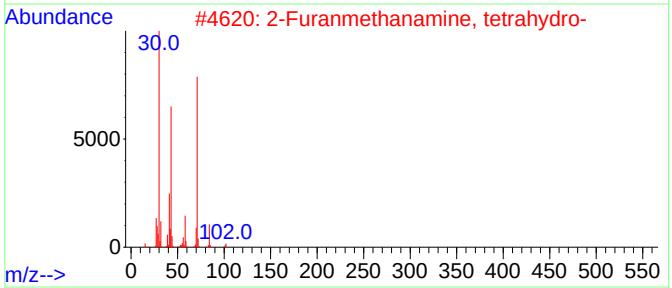
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Propene, 3-bromo- | 120 | C3H5Br | 000106-95-6 | 9 |
| 2 | 2-Furanmethanamine, tetrahydro- | 101 | C5H11NO | 004795-29-3 | 7 |
| 3 | Oxirane, propyl- | 86 | C5H10O | 001003-14-1 | 7 |
| 4 | Benzeneethanamine | 121 | C8H11N | 000064-04-0 | 4 |
| 5 | Butane, 1-bromo-2-methyl- | 150 | C5H11Br | 005973-11-5 | 4 |

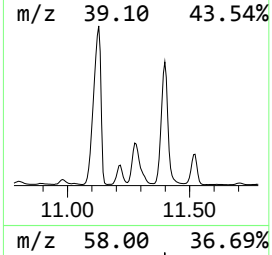


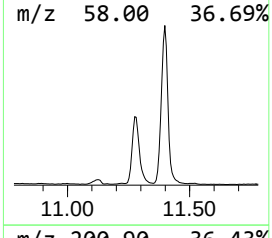



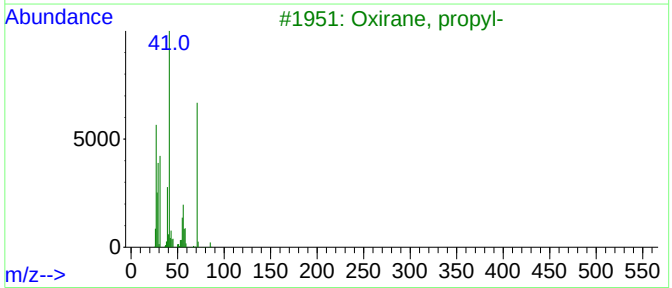



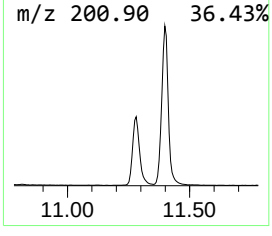



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

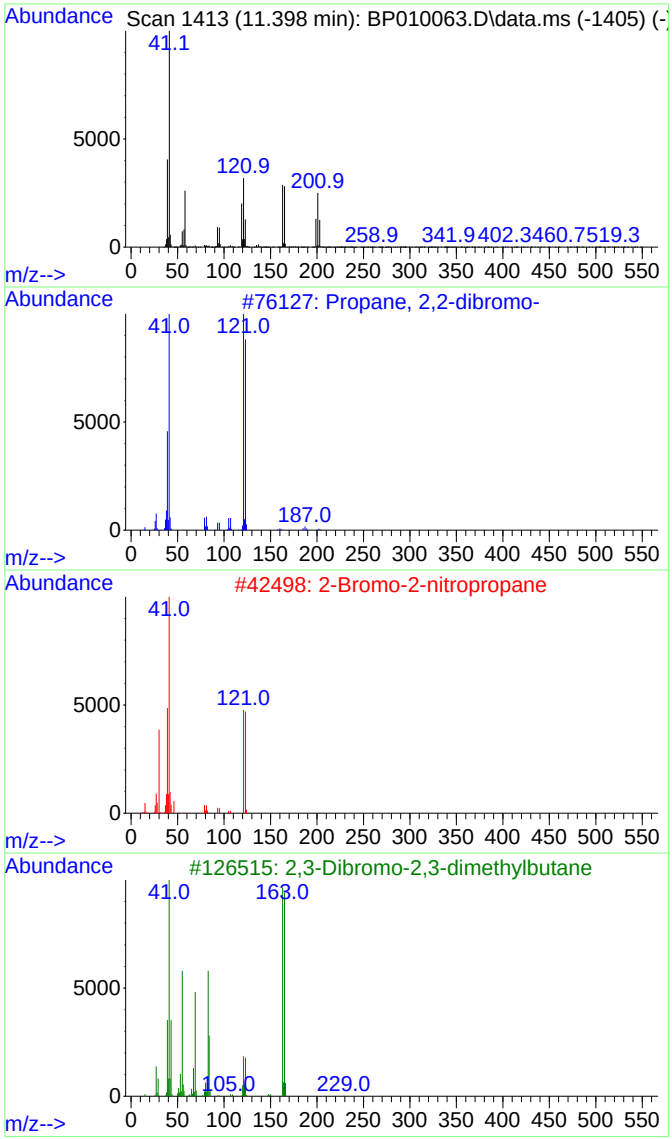
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

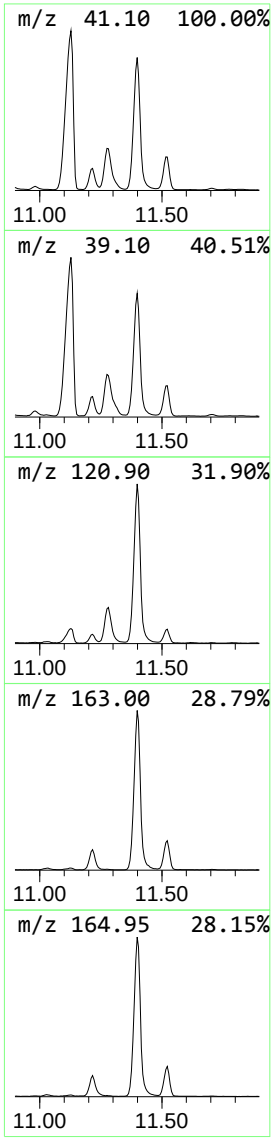
**********

Peak Number 18 unknown-04 Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.398 | 124.38 ng/ul | 14540500 | Naphthalene-d8 | 10.757 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 2,2-dibromo- | 200 | C3H6Br2 | 000594-16-1 | 7 |
| 2 | 2-Bromo-2-nitropropane | 167 | C3H6BrNO2 | 005447-97-2 | 4 |
| 3 | 2,3-Dibromo-2,3-dimethylbutane | 242 | C6H12Br2 | 000594-81-0 | 4 |
| 4 | Benzisoxazole-2-acetic acid, hyd... | 191 | C9H9N3O2 | 055244-10-5 | 4 |
| 5 | Benzeneethanamine | 121 | C8H11N | 000064-04-0 | 3 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

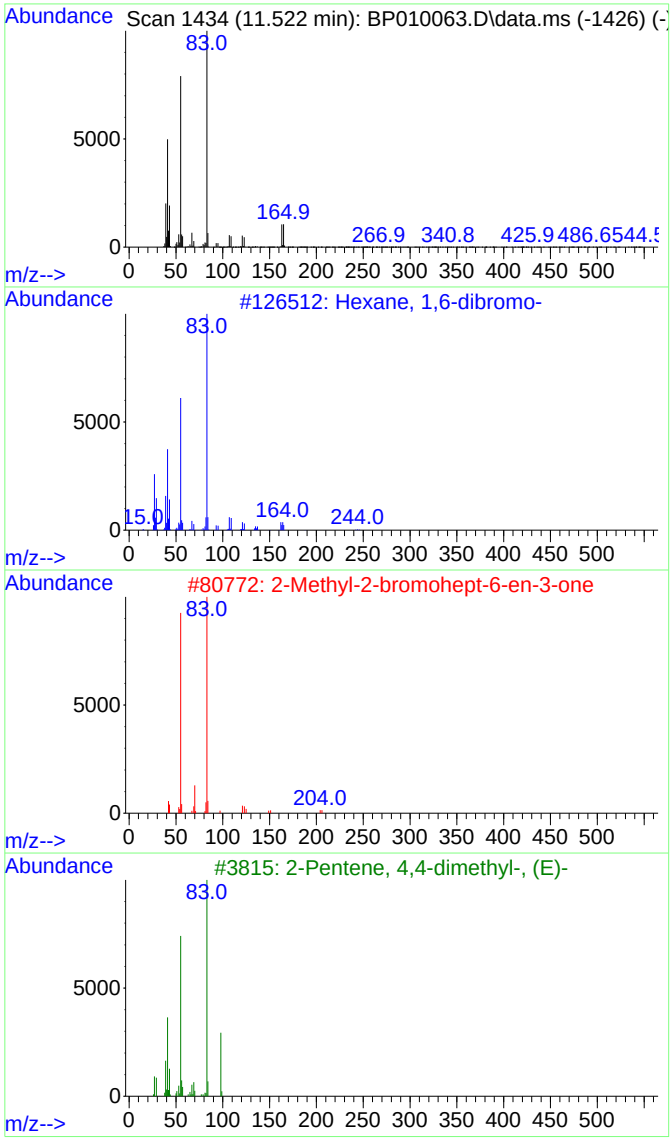
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

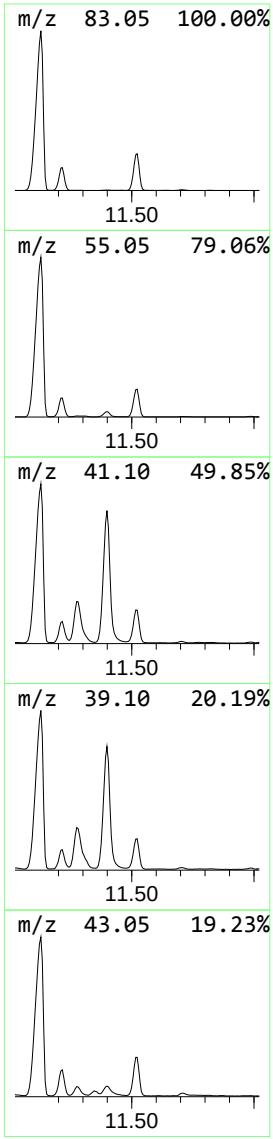
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 19 Hexane, 1,6-dibromo- Concentration Rank 27

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.522 | 49.15 ng/ul | 5745430 | Naphthalene-d8 | 10.757 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexane, 1,6-dibromo- | 242 | C6H12Br2 | 000629-03-8 | 78 |
| 2 | 2-Methyl-2-bromohept-6-en-3-one | 204 | C8H13BrO | 1000424-28-0 | 59 |
| 3 | 2-Pentene, 4,4-dimethyl-, (E)- | 98 | C7H14 | 000690-08-4 | 53 |
| 4 | 1-Butene, 2,3,3-trimethyl- | 98 | C7H14 | 000594-56-9 | 53 |
| 5 | Cyclopropane, 2-bromo-1,1,3-trim... | 162 | C6H11Br | 036617-00-2 | 53 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

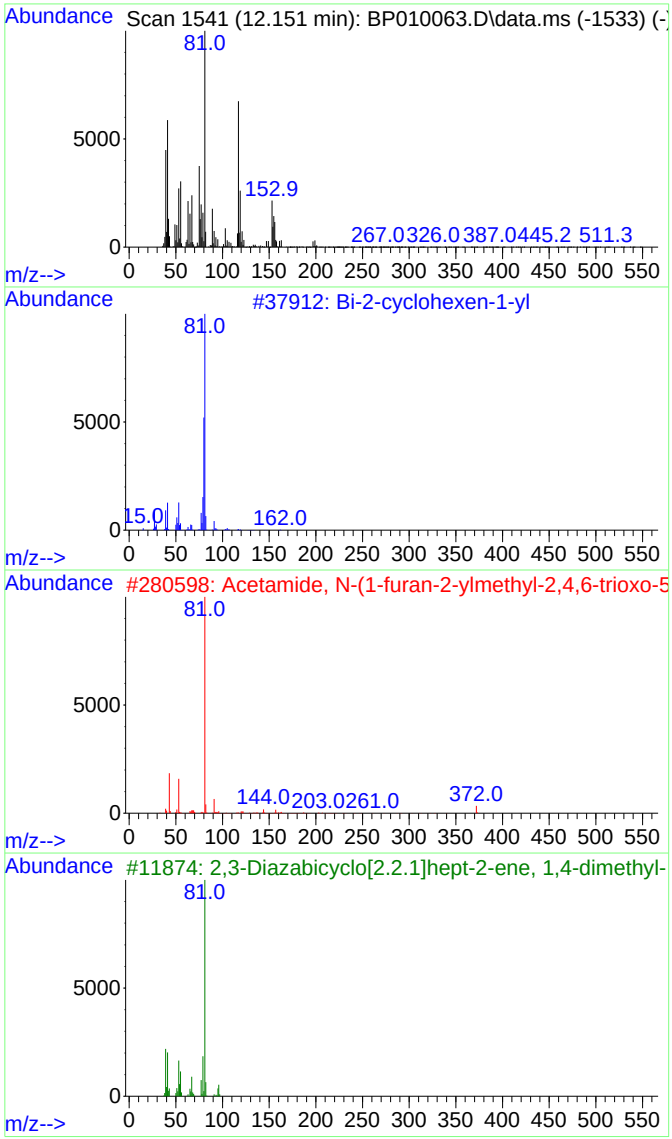
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

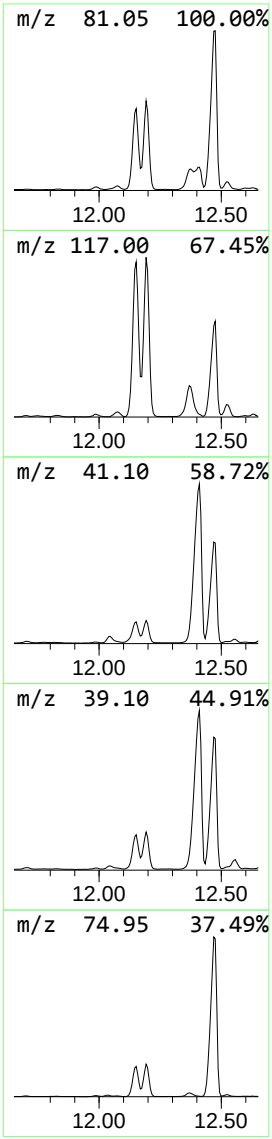
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 20 unknown-05 Concentration Rank 23

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.151 | 54.31 ng/ul | 6348850 | Naphthalene-d8 | 10.757 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Bi-2-cyclohexen-1-yl | 162 | C12H18 | 001541-20-4 | 25 |
| 2 | Acetamide, N-(1-furan-2-ylmethyl... | 372 | C14H11F3N4O5 | 1000303-51-7 | 22 |
| 3 | 2,3-Diazabicyclo[2.2.1]hept-2-en... | 124 | C7H12N2 | 071312-54-4 | 22 |
| 4 | 2-Hexen-4-yn-1-ol, (E)- | 96 | C6H8O | 053497-80-6 | 16 |
| 5 | Ethanethioic acid, S-(2-furanylm... | 156 | C7H8O2S | 013678-68-7 | 16 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

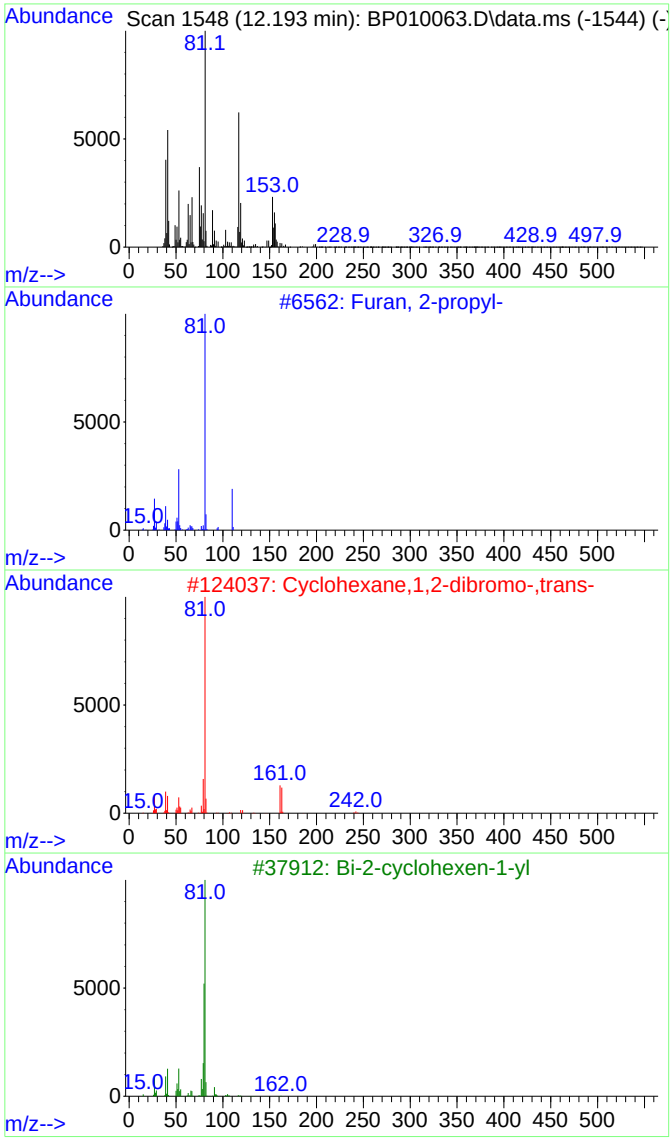
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

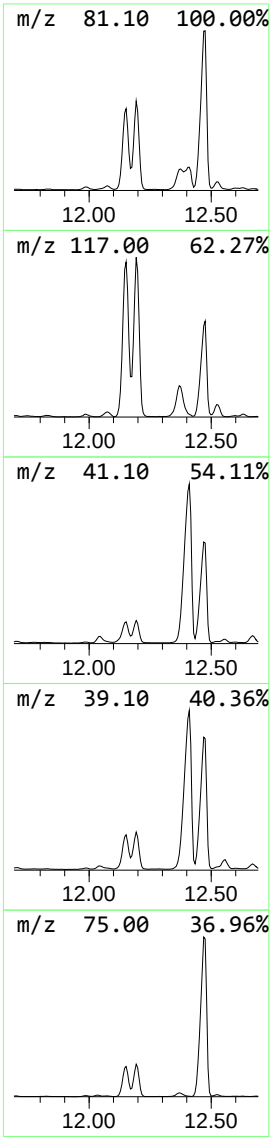
**********

Peak Number 21 unknown-06 Concentration Rank 25

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.193 | 53.28 ng/ul | 6228350 | Naphthalene-d8 | 10.757 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Furan, 2-propyl- | 110 | C7H10O | 004229-91-8 | 22 |
| 2 | Cyclohexane,1,2-dibromo-,trans- | 240 | C6H10Br2 | 007429-37-0 | 22 |
| 3 | Bi-2-cyclohexen-1-yl | 162 | C12H18 | 001541-20-4 | 22 |
| 4 | Bis(2-furfuryl)disulfide | 226 | C10H10O2S2 | 004437-20-1 | 12 |
| 5 | Benzyl nitrile | 117 | C8H7N | 000140-29-4 | 11 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

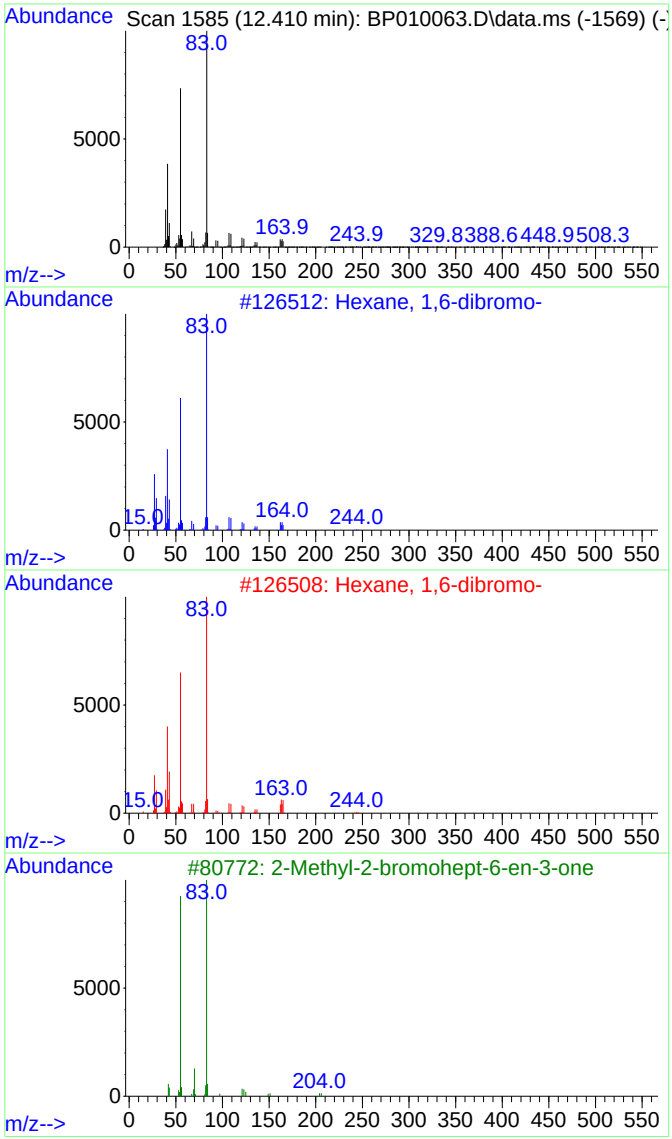
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

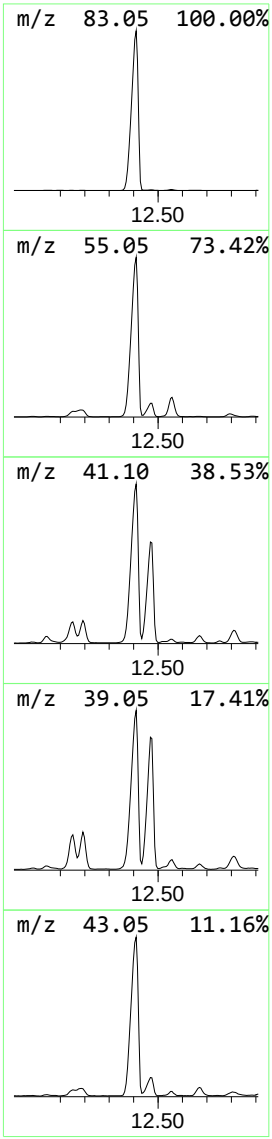
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 22 2-Methyl-2-bromohept-6-en-3... Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.410 | 335.61 ng/ul | 39232900 | Naphthalene-d8 | 10.757 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexane, 1,6-dibromo- | 242 | C6H12Br2 | 000629-03-8 | 97 |
| 2 | Hexane, 1,6-dibromo- | 242 | C6H12Br2 | 000629-03-8 | 90 |
| 3 | 2-Methyl-2-bromohept-6-en-3-one | 204 | C8H13BrO | 1000424-28-0 | 64 |
| 4 | Cumenyl angelate, o- | 218 | C14H18O2 | 1000383-67-2 | 64 |
| 5 | Cumenyl tiglate, m- | 218 | C14H18O2 | 1000383-67-1 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

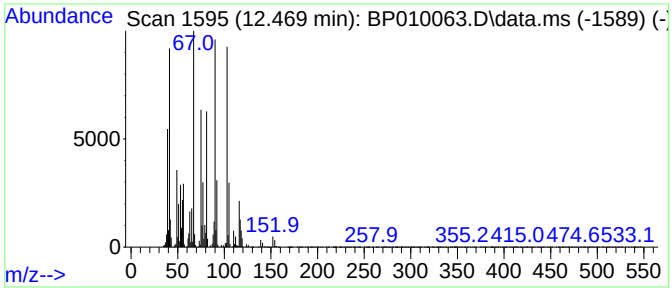
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

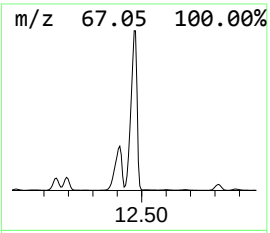
**********

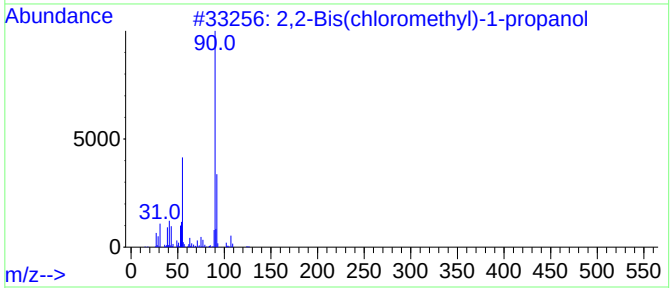
Peak Number 23 unknown-07 Concentration Rank 10

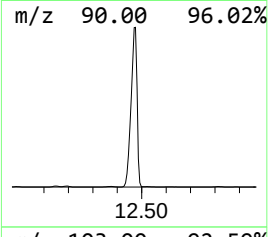
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.469 | 224.99 ng/ul | 26302000 | Naphthalene-d8 | 10.757 |

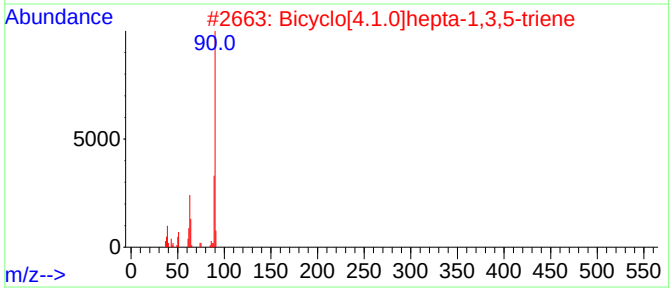
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2,2-Bis(chloromethyl)-1-propanol | 156 | C5H10Cl2O | 005355-54-4 | 38 |
| 2 | Bicyclo[4.1.0]hepta-1,3,5-triene | 90 | C7H6 | 004646-69-9 | 38 |
| 3 | Thiourea, methyl- | 90 | C2H6N2S | 000598-52-7 | 35 |
| 4 | Thiourea, methyl- | 90 | C2H6N2S | 000598-52-7 | 35 |
| 5 | 1,3,5-Trichloro-4-methyl-2-penta... | 202 | C6H9Cl3O | 1000140-03-8 | 28 |

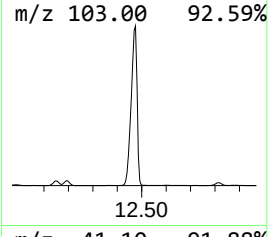


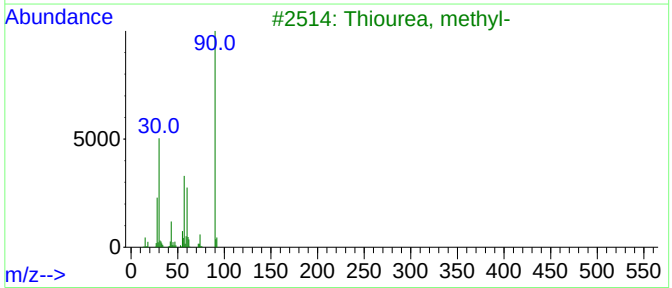



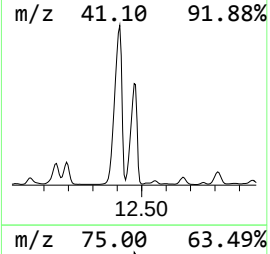



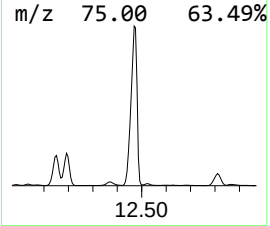



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

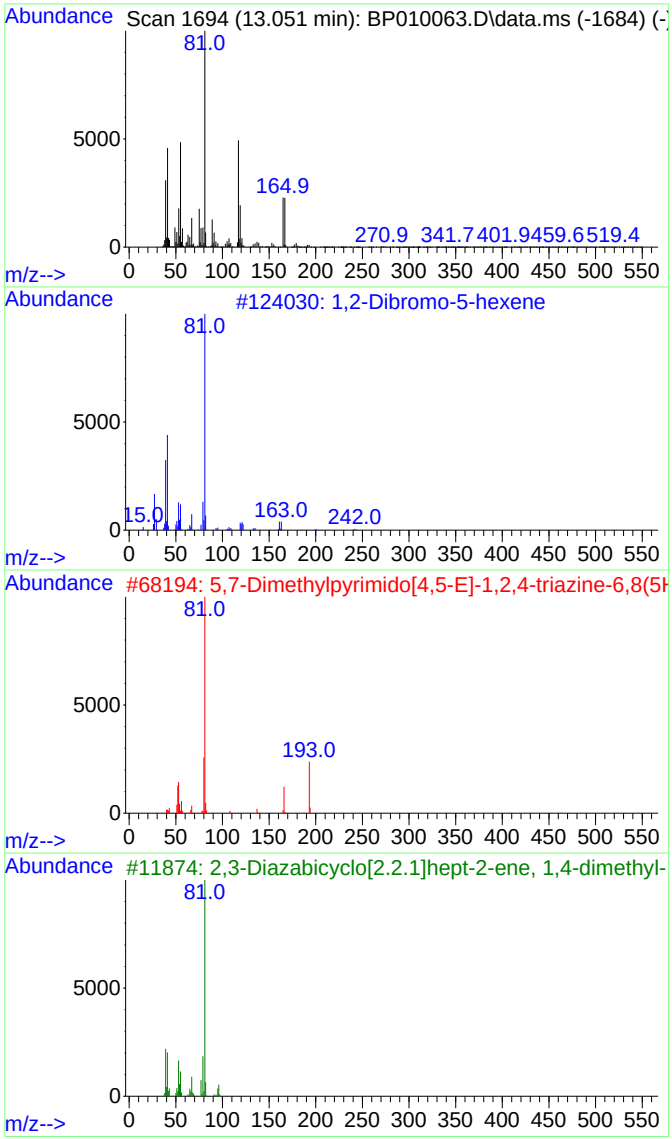
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

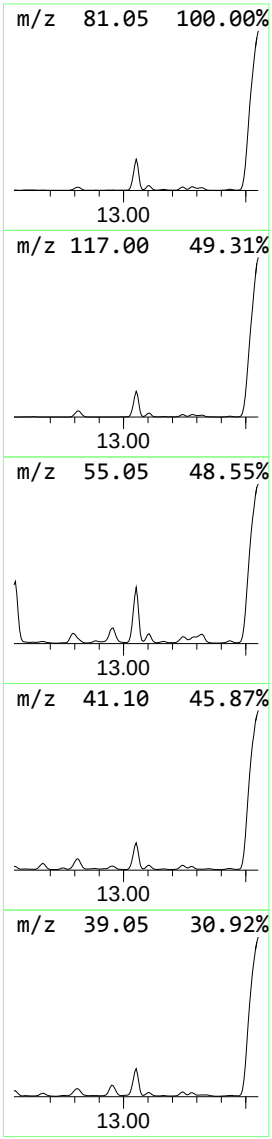
**********

Peak Number 27 unknown-08 Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.051 | 70.82 ng/ul | 8324470 | Acenaphthene-d10 | 14.581 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,2-Dibromo-5-hexene | 240 | C6H10Br2 | 004285-48-7 | 27 |
| 2 | 5,7-Dimethylpyrimido[4,5-E]-1,2,... | 193 | C7H7N5O2 | 016044-79-4 | 22 |
| 3 | 2,3-Diazabicyclo[2.2.1]hept-2-en... | 124 | C7H12N2 | 071312-54-4 | 22 |
| 4 | 2,4-Dimethyl 1,4-pentadiene | 96 | C7H12 | 004161-65-3 | 22 |
| 5 | 2,4-Heptadienal, (E,E)- | 110 | C7H10O | 004313-03-5 | 22 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

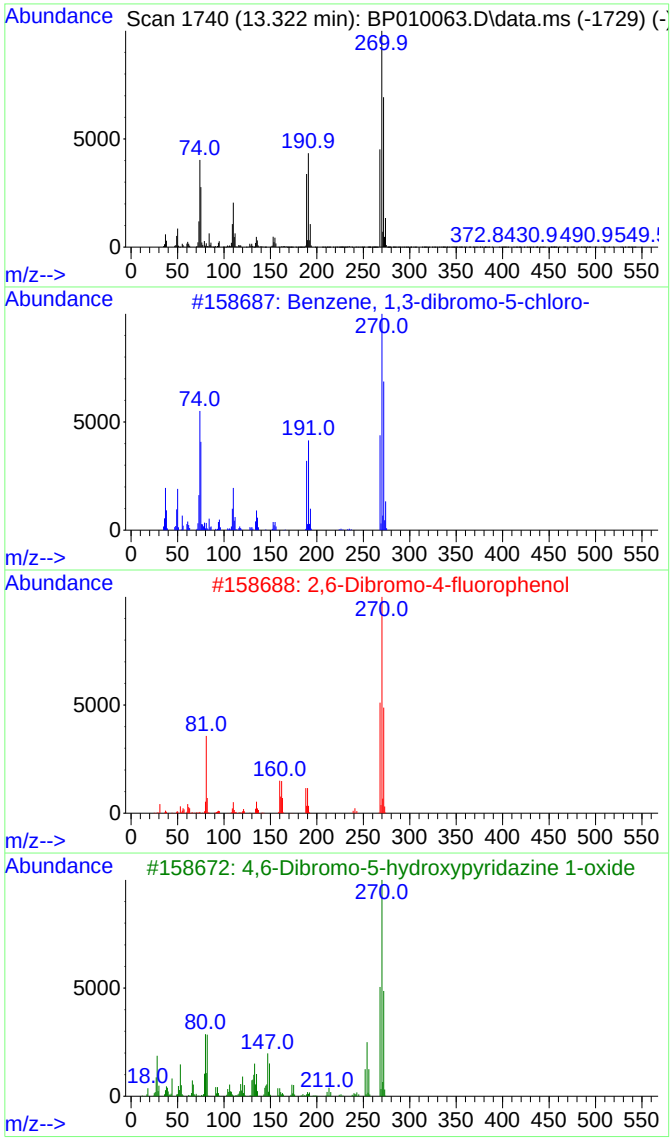
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

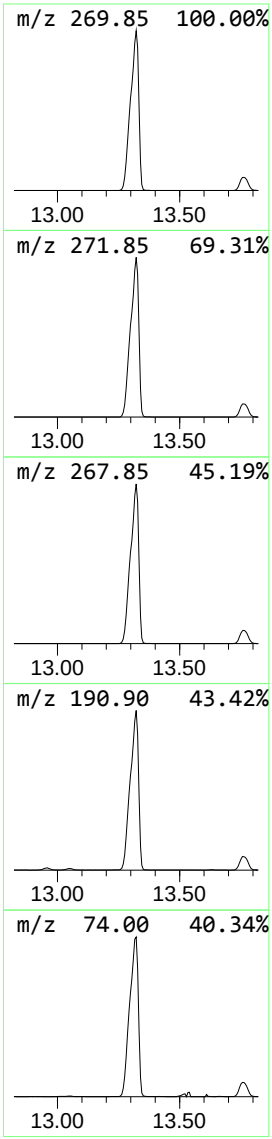
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 28 Benzene, 1,3-dibromo-5-chloro- Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.322 | 287.17 ng/ul | 33754800 | Acenaphthene-d10 | 14.581 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1,3-dibromo-5-chloro- | 268 | C6H3Br2Cl | 014862-52-3 | 94 |
| 2 | 2,6-Dibromo-4-fluorophenol | 268 | C6H3Br2FO | 000344-20-7 | 43 |
| 3 | 4,6-Dibromo-5-hydroxypyridazine ... | 268 | C4H2Br2N2O2 | 019971-61-0 | 43 |
| 4 | 4,5-Dibromo-2-furoic acid | 268 | C5H2Br2O3 | 002434-03-9 | 25 |
| 5 | 6-Bromo-7-methylbenzo(b)thiophen... | 270 | C10H7BrO2S | 001678-69-9 | 25 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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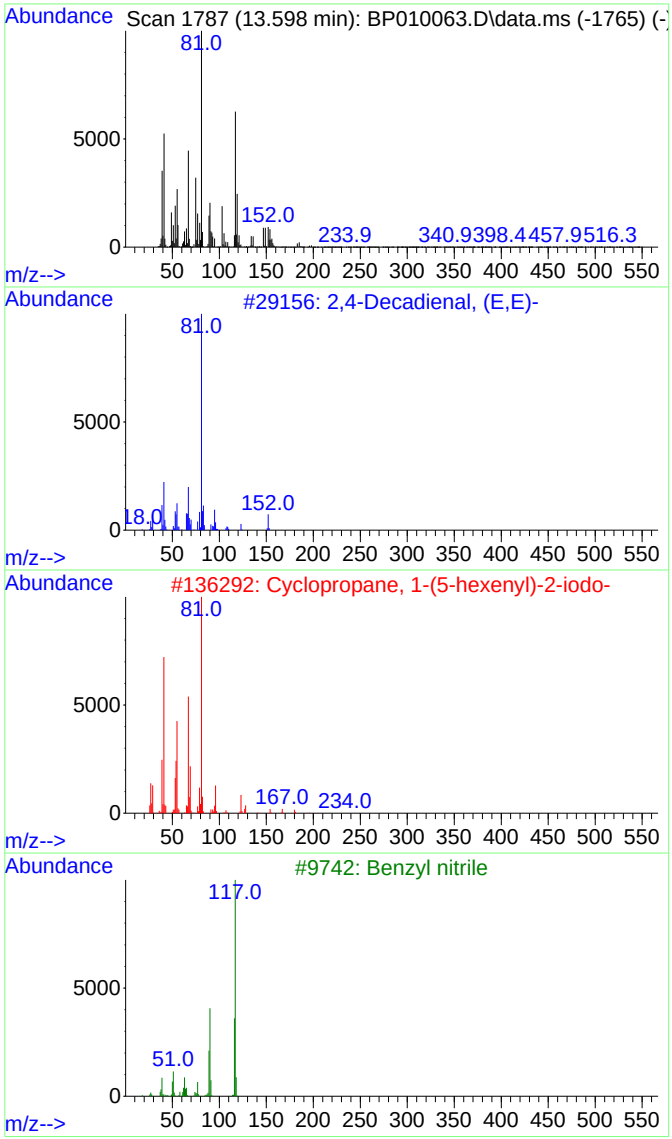
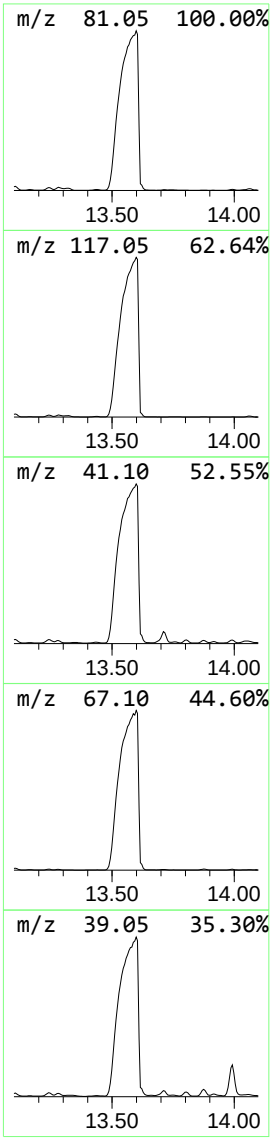
Peak Number 29 unknown-09 Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.598 | 1730.63 ng/ul | 203422000 | Acenaphthene-d10 | 14.581 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2,4-Decadienal, (E,E)- | 152 | C10H16O | 025152-84-5 | 38 |
| 2 | Cyclopropane, 1-(5-hexenyl)-2-iodo- | 250 | C9H15I | 074685-40-8 | 35 |
| 3 | Benzyl nitrile | 117 | C8H7N | 000140-29-4 | 25 |
| 4 | 2-Furanacetaldehyde, .alpha.-pro... | 152 | C9H12O2 | 031681-26-2 | 22 |
| 5 | 1,2-Dibromo-5-hexene | 240 | C6H10Br2 | 004285-48-7 | 22 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

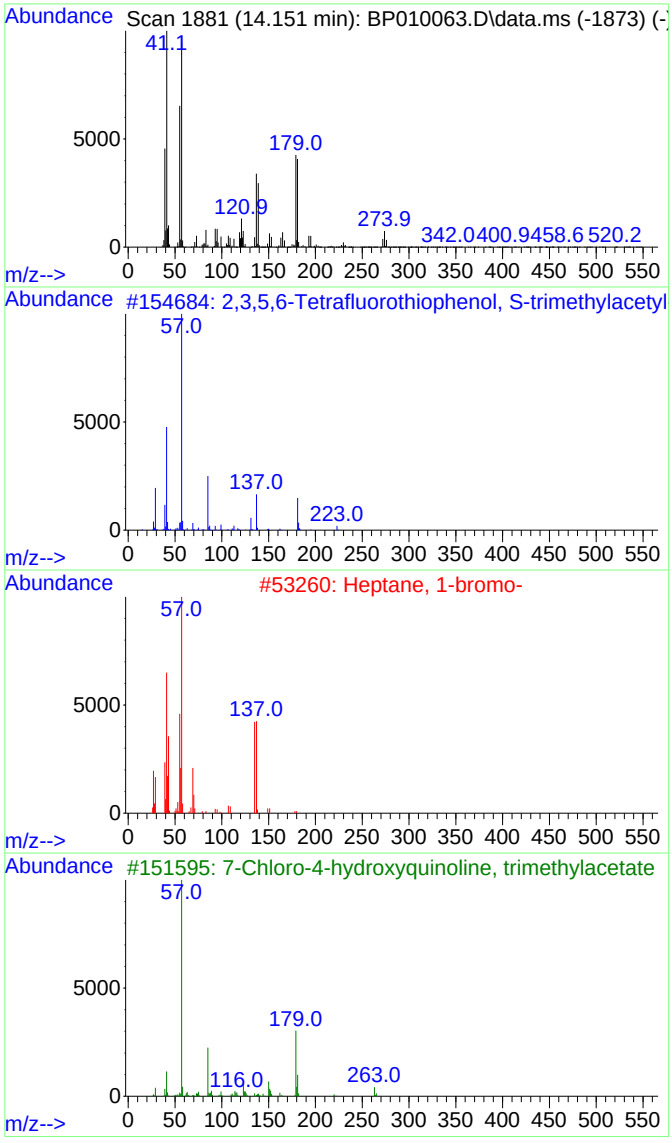
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

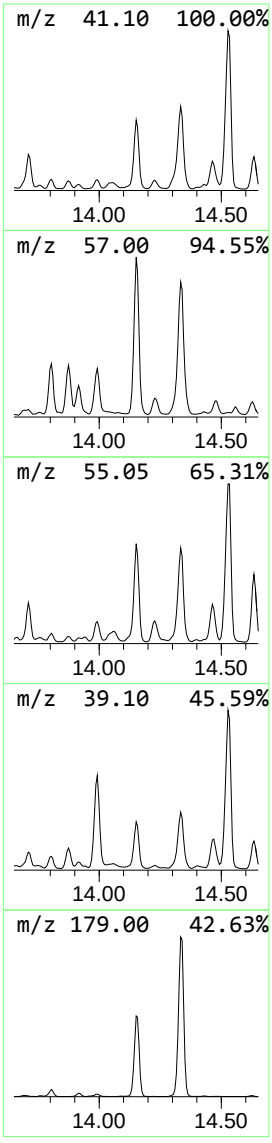
**********

Peak Number 33 unknown-10 Concentration Rank 30

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.151 | 38.60 ng/ul | 4536740 | Acenaphthene-d10 | 14.581 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2,3,5,6-Tetrafluorothiophenol, S... | 266 | C11H10F4OS | 1000470-48-7 | 10 |
| 2 | Heptane, 1-bromo- | 178 | C7H15Br | 000629-04-9 | 10 |
| 3 | 7-Chloro-4-hydroxyquinoline, tri... | 263 | C14H14ClNO2 | 1000463-55-2 | 10 |
| 4 | 1,2-Dipropylcyclopropene-3-carbo... | 168 | C10H16O2 | 001605-38-5 | 10 |
| 5 | Butane, 2,3-dibromo- | 214 | C4H8Br2 | 005408-86-6 | 10 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

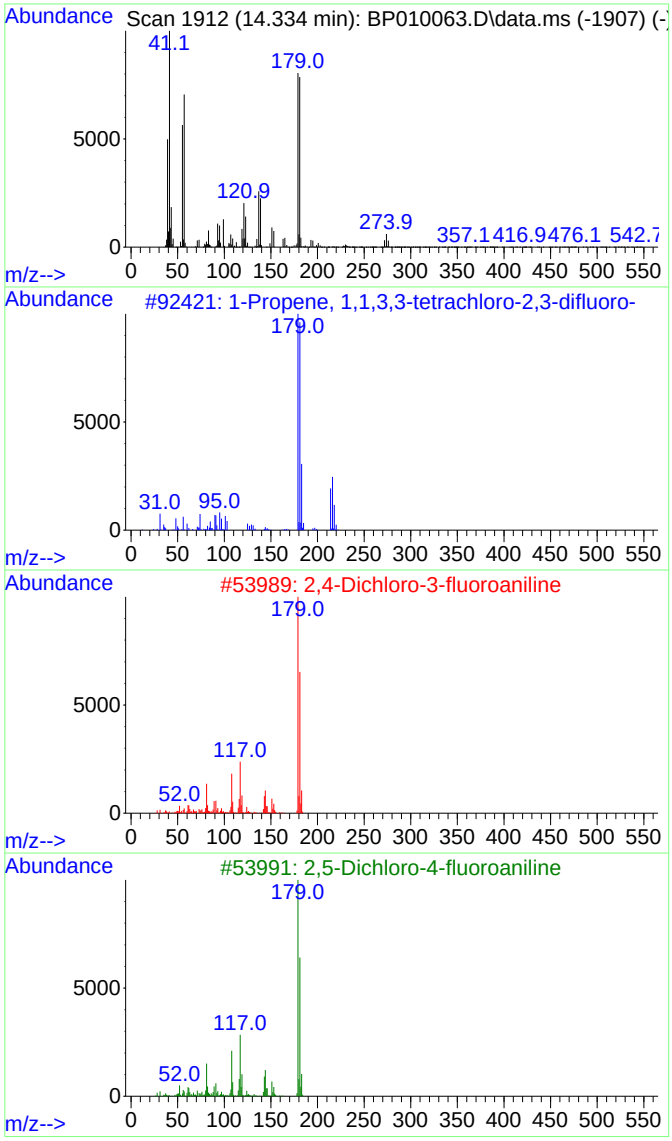
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

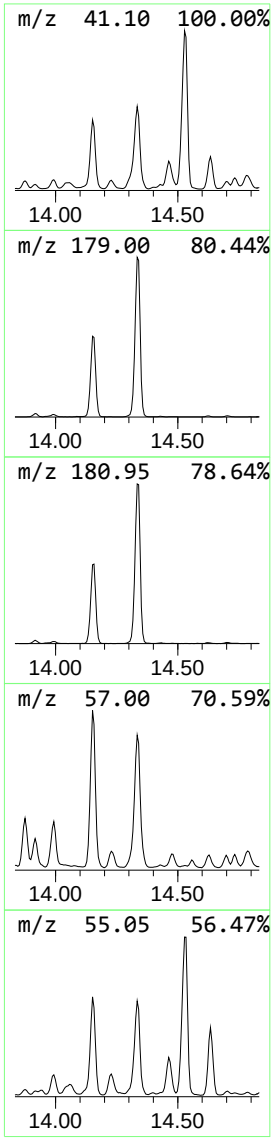
**********

Peak Number 34 unknown-11 Concentration Rank 26

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.334 | 51.28 ng/ul | 6027000 | Acenaphthene-d10 | 14.581 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Propene, 1,1,3,3-tetrachloro-2... | 214 | C3Cl4F2 | 000815-16-7 | 47 |
| 2 | 2,4-Dichloro-3-fluoroaniline | 179 | C6H4Cl2FN | 000443-93-6 | 35 |
| 3 | 2,5-Dichloro-4-fluoroaniline | 179 | C6H4Cl2FN | 002729-37-5 | 35 |
| 4 | N-(2,4-Dichloro-3-fluorophenyl)a... | 221 | C8H6Cl2FNO | 1000497-83-4 | 35 |
| 5 | 4,5-Dichloro-2-fluoroaniline | 179 | C6H4Cl2FN | 002729-36-4 | 35 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

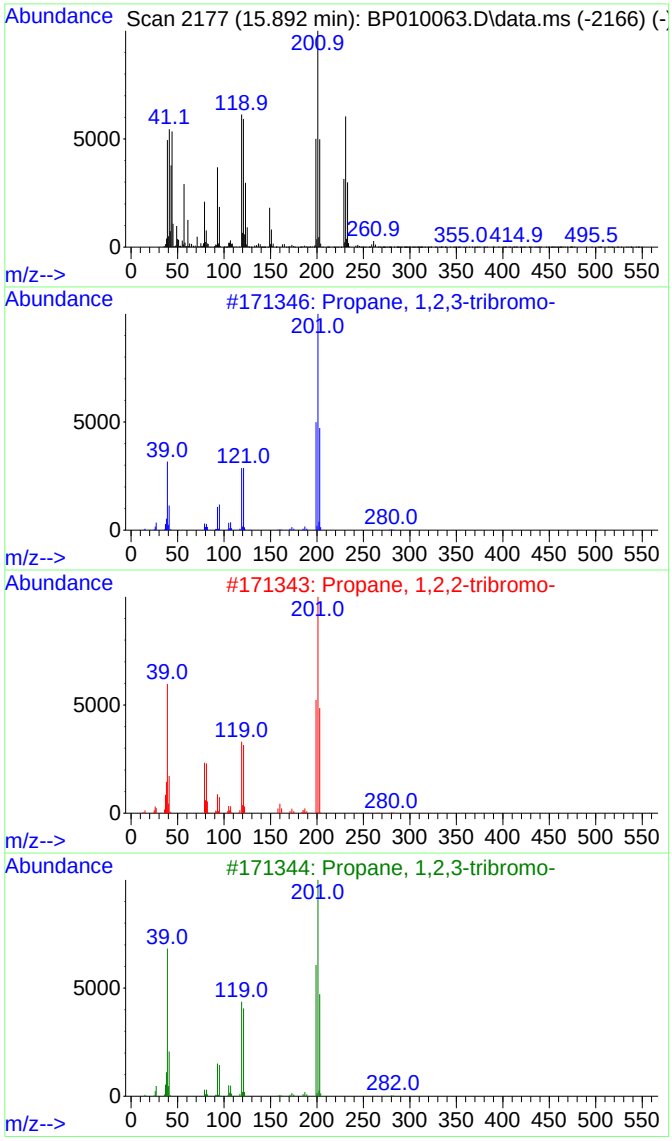
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

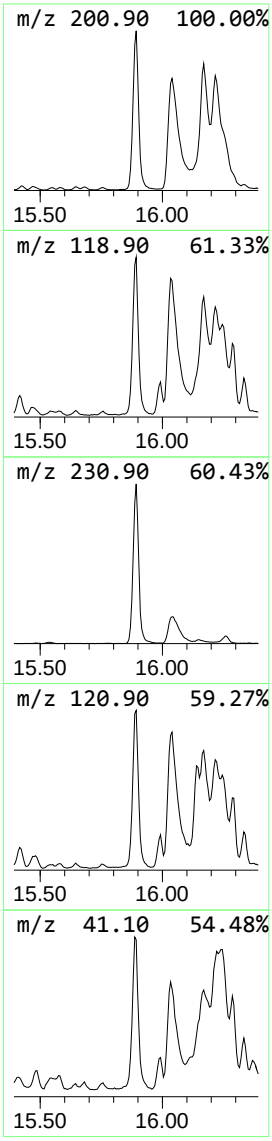
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 37 unknown-12 Concentration Rank 24

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.892 | 53.66 ng/ul | 6307720 | Acenaphthene-d10 | 14.581 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 32 |
| 2 | Propane, 1,2,2-tribromo- | 278 | C3H5Br3 | 014476-30-3 | 32 |
| 3 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 25 |
| 4 | 2(3H)-Benzothiazolethione, 5-chl... | 201 | C7H4ClNS2 | 005331-91-9 | 12 |
| 5 | Pyridine, 3,5-dichloro-2,4,6-tri... | 201 | C5Cl2F3N | 001737-93-5 | 9 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

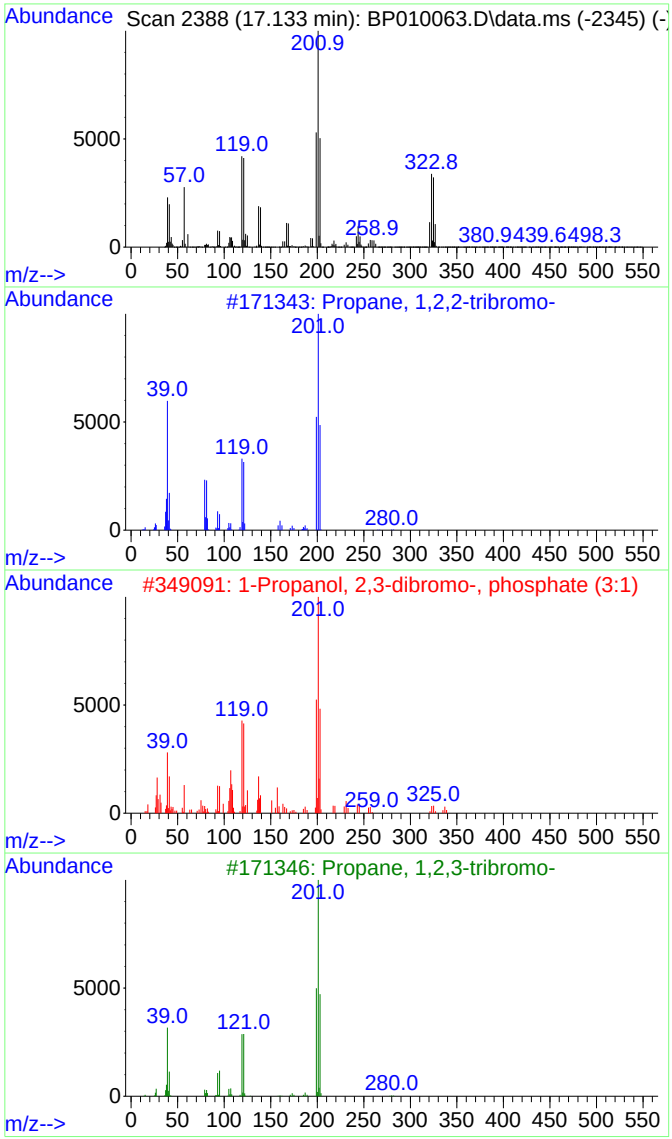
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

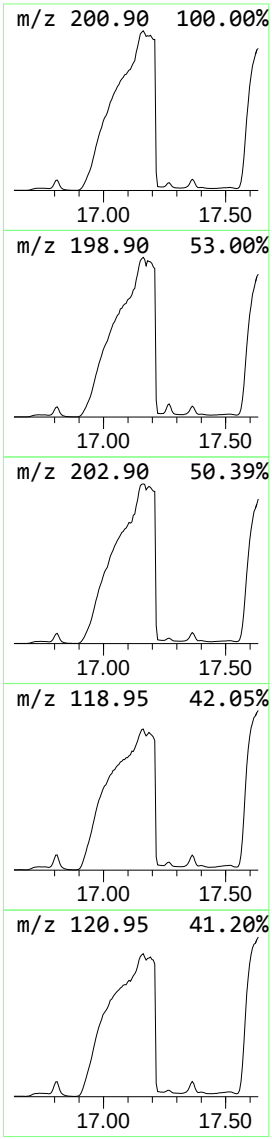
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 47 Propane, 1,2,2-tribromo- Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.133 | 1276.66 ng/ul | 399607000 | Phenanthrene-d10 | 17.357 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 1,2,2-tribromo- | 278 | C3H5Br3 | 014476-30-3 | 53 |
| 2 | 1-Propanol, 2,3-dibromo-, phosph... | 692 | C9H15Br6O4P | 000126-72-7 | 39 |
| 3 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 36 |
| 4 | 2(3H)-Benzothiazolethione, 5-chl... | 201 | C7H4ClNS2 | 005331-91-9 | 16 |
| 5 | 2(3H)-Benzothiazolethione, 5-chl... | 201 | C7H4ClNS2 | 005331-91-9 | 12 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

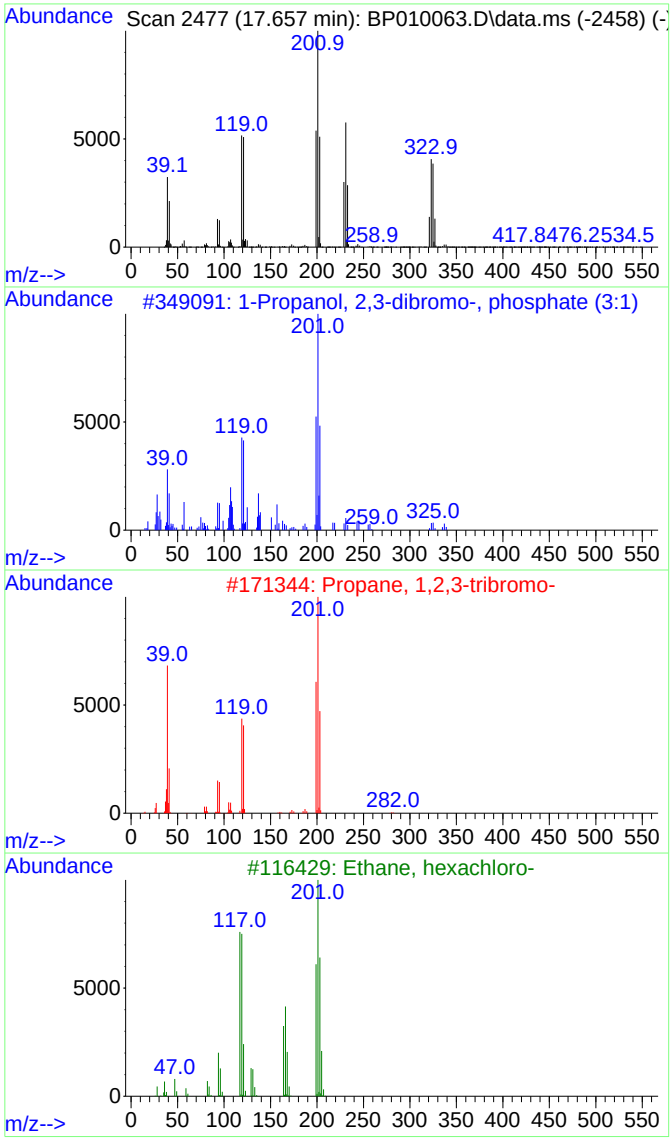
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

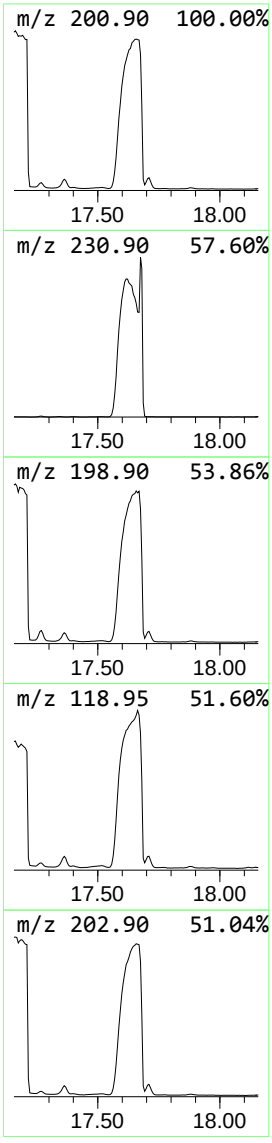
**********

Peak Number 49 unknown-13 Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.657 | 642.47 ng/ul | 201100000 | Phenanthrene-d10 | 17.357 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Propanol, 2,3-dibromo-, phosph... | 692 | C9H15Br6O4P | 000126-72-7 | 38 |
| 2 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 38 |
| 3 | Ethane, hexachloro- | 234 | C2Cl6 | 000067-72-1 | 37 |
| 4 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 11 |
| 5 | Cyclopropane, 1,1-dibromo-2-chloro- | 232 | C3H3Br2Cl | 077628-95-6 | 9 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

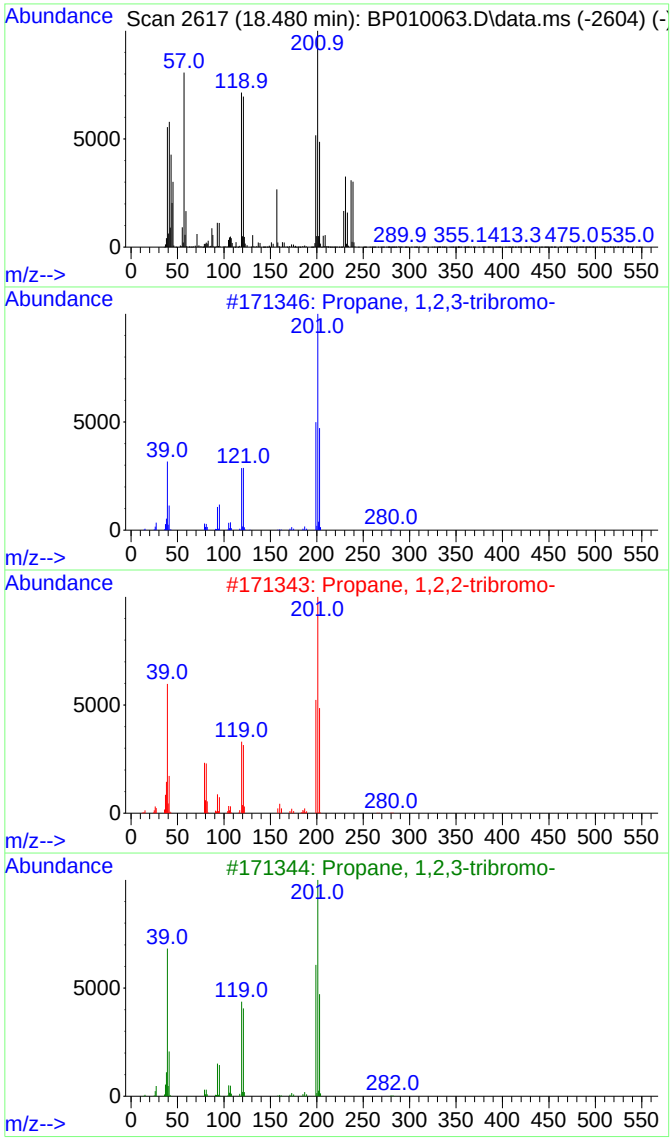
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

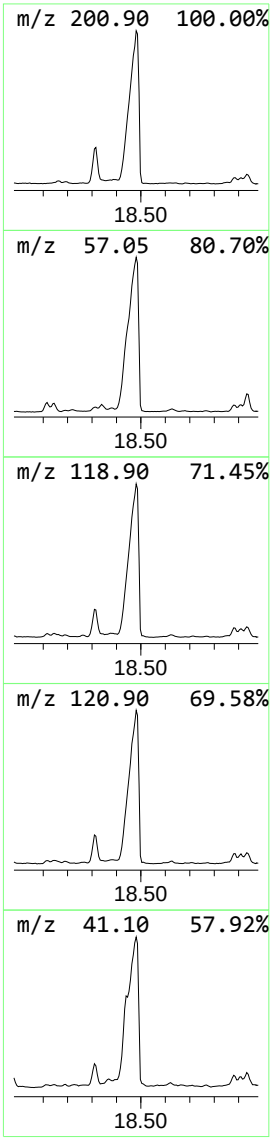
**********

Peak Number 52 unknown-14 Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.480 | 115.65 ng/ul | 36200300 | Phenanthrene-d10 | 17.357 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 47 |
| 2 | Propane, 1,2,2-tribromo- | 278 | C3H5Br3 | 014476-30-3 | 47 |
| 3 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 45 |
| 4 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 13 |
| 5 | 2(3H)-Benzothiazolethione, 5-chl... | 201 | C7H4ClNS2 | 005331-91-9 | 12 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

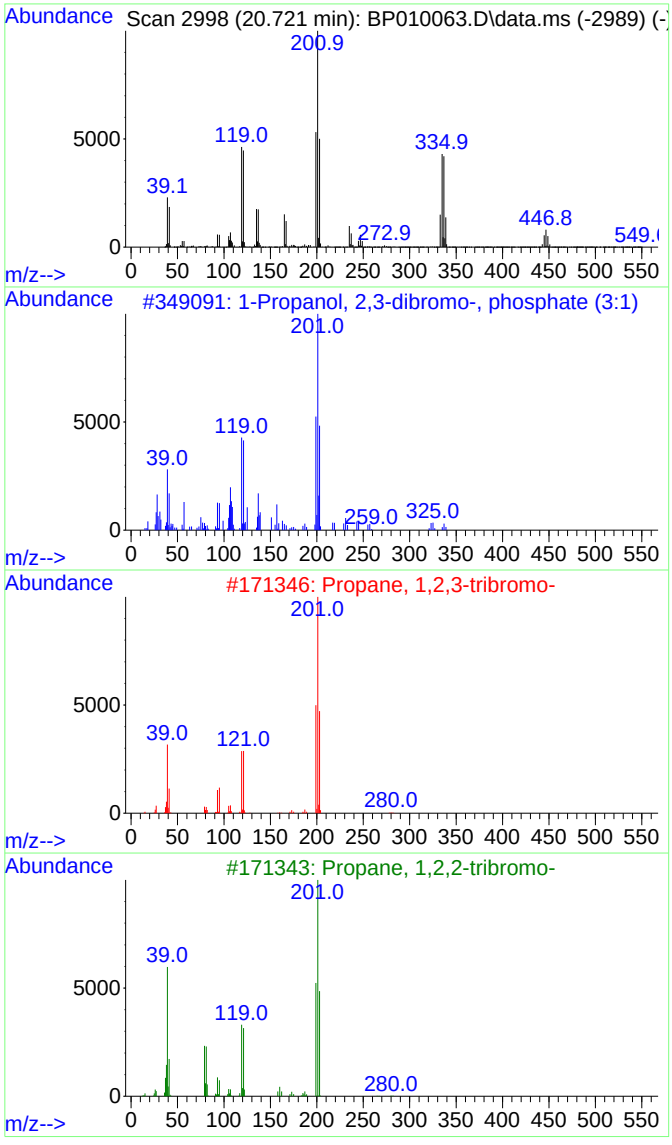
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

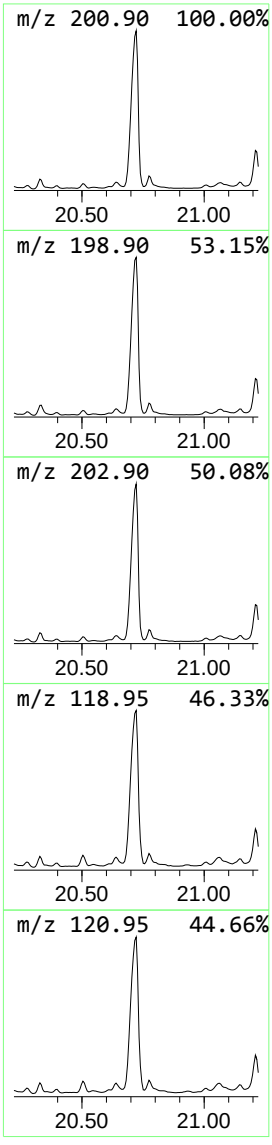
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 61 unknown-15 Concentration Rank 21

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.721 | 61.46 ng/ul | 73380000 | Chrysene-d12 | 21.416 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Propanol, 2,3-dibromo-, phosph... | 692 | C9H15Br6O4P | 000126-72-7 | 40 |
| 2 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 40 |
| 3 | Propane, 1,2,2-tribromo- | 278 | C3H5Br3 | 014476-30-3 | 40 |
| 4 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 32 |
| 5 | 2(3H)-Benzothiazolethione, 5-chl... | 201 | C7H4ClNS2 | 005331-91-9 | 9 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

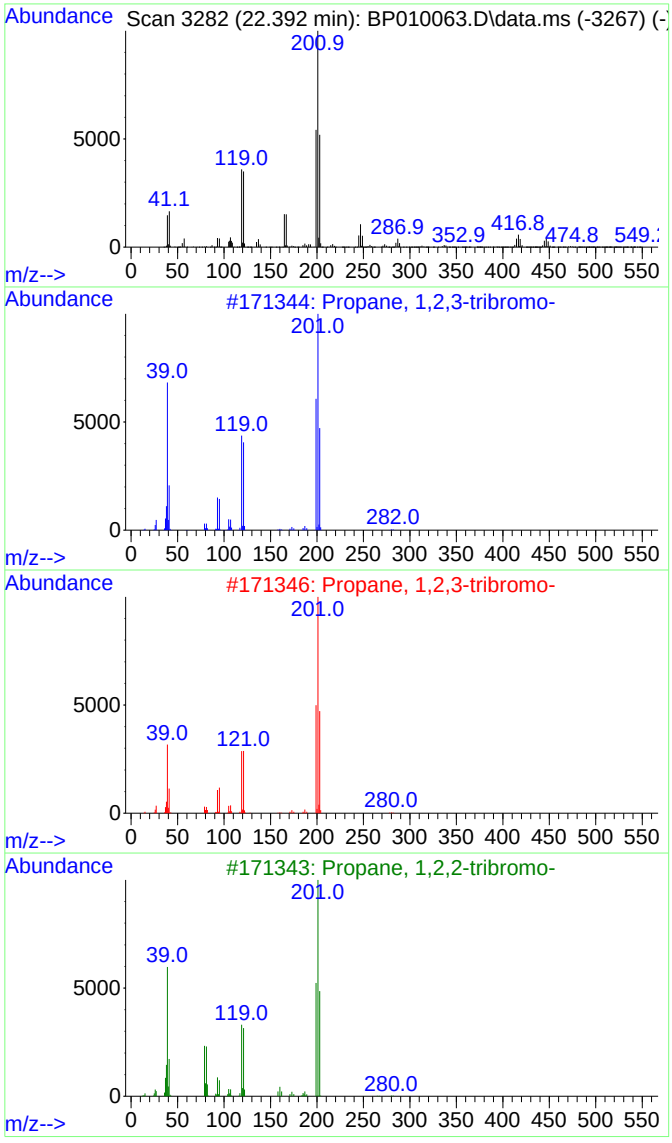
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

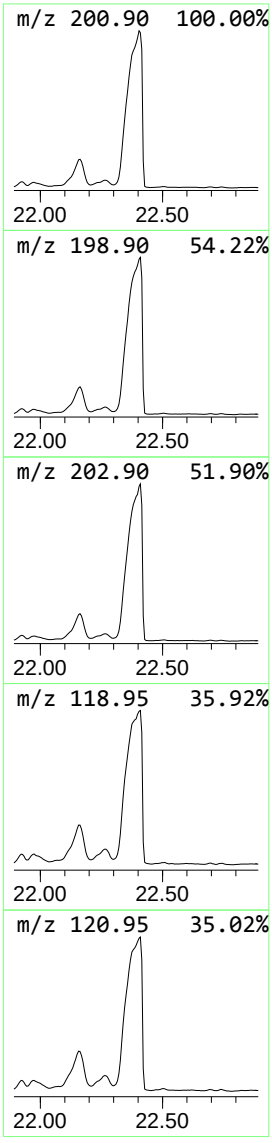
**********

Peak Number 72 unknown-16 Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 22.392 | 89.18 ng/ul | 106483000 | Chrysene-d12 | 21.416 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 80 |
| 2 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 72 |
| 3 | Propane, 1,2,2-tribromo- | 278 | C3H5Br3 | 014476-30-3 | 56 |
| 4 | 2(3H)-Benzothiazolethione, 5-chl... | 201 | C7H4ClNS2 | 005331-91-9 | 25 |
| 5 | Succinic acid, butyl 2-(4-nitrop... | 339 | C16H21NO7 | 1000381-13-5 | 10 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

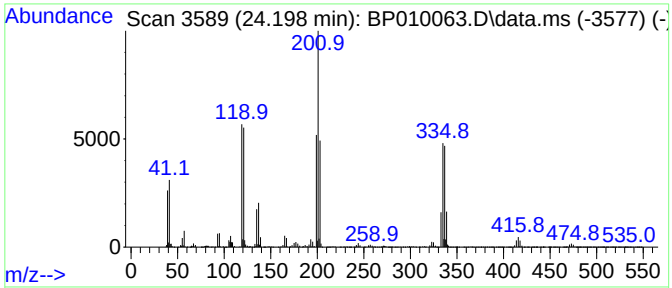
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

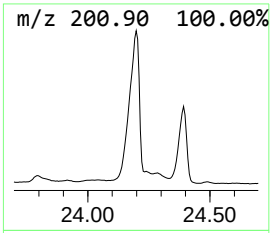
**********

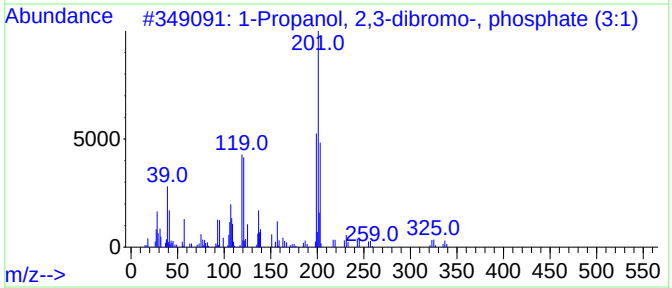
Peak Number 75 unknown-17 Concentration Rank 19

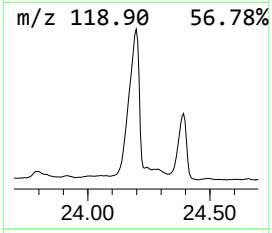
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 24.198 | 65.31 ng/ul | 37838700 | Perylene-d12 | 23.886 |

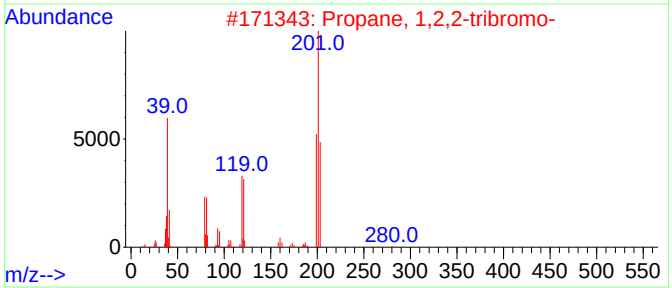
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Propanol, 2,3-dibromo-, phosph... | 692 | C9H15Br6O4P | 000126-72-7 | 45 |
| 2 | Propane, 1,2,2-tribromo- | 278 | C3H5Br3 | 014476-30-3 | 37 |
| 3 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 32 |
| 4 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 25 |
| 5 | 2(3H)-Benzothiazolethione, 5-chl... | 201 | C7H4ClNS2 | 005331-91-9 | 22 |

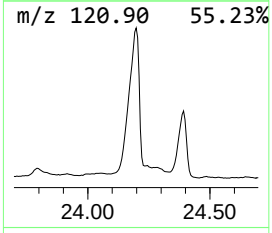


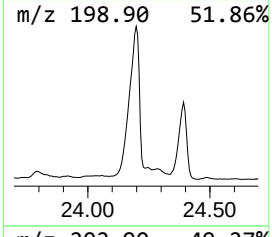



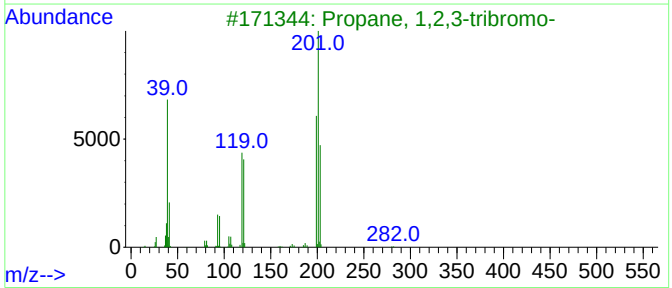



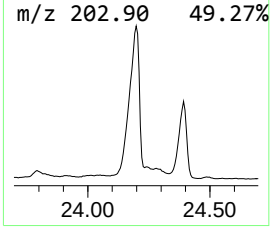



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

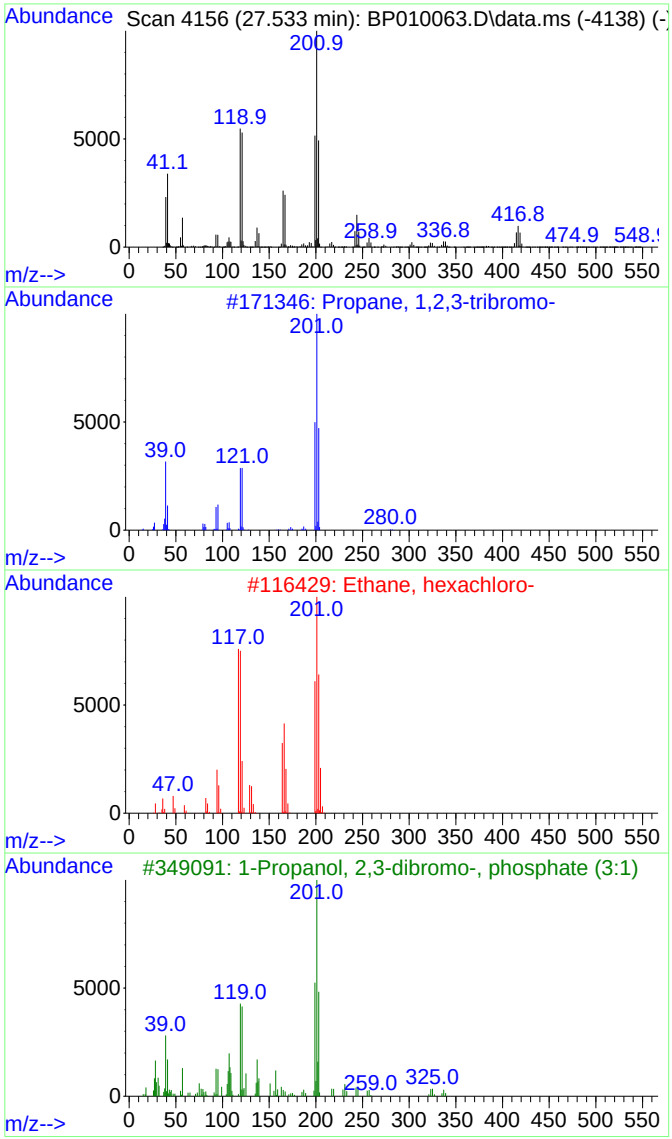
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

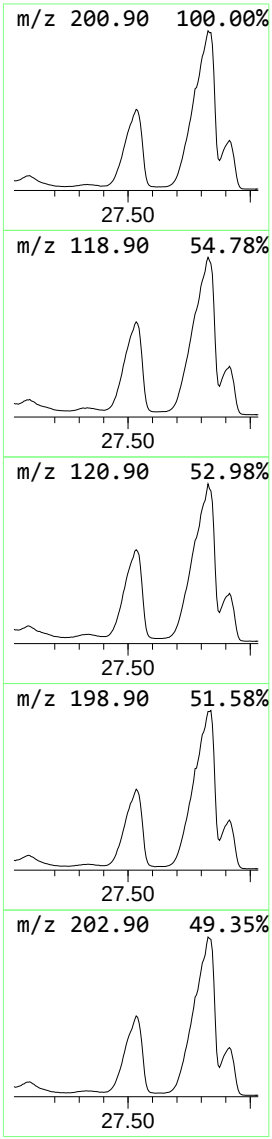
**********

Peak Number 79 unknown-18 Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 27.533 | 63.71 ng/ul | 36908900 | Perylene-d12 | 23.886 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 45 |
| 2 | Ethane, hexachloro- | 234 | C2Cl6 | 000067-72-1 | 45 |
| 3 | 1-Propanol, 2,3-dibromo-, phosph... | 692 | C9H15Br6O4P | 000126-72-7 | 45 |
| 4 | Propane, 1,2,2-tribromo- | 278 | C3H5Br3 | 014476-30-3 | 40 |
| 5 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

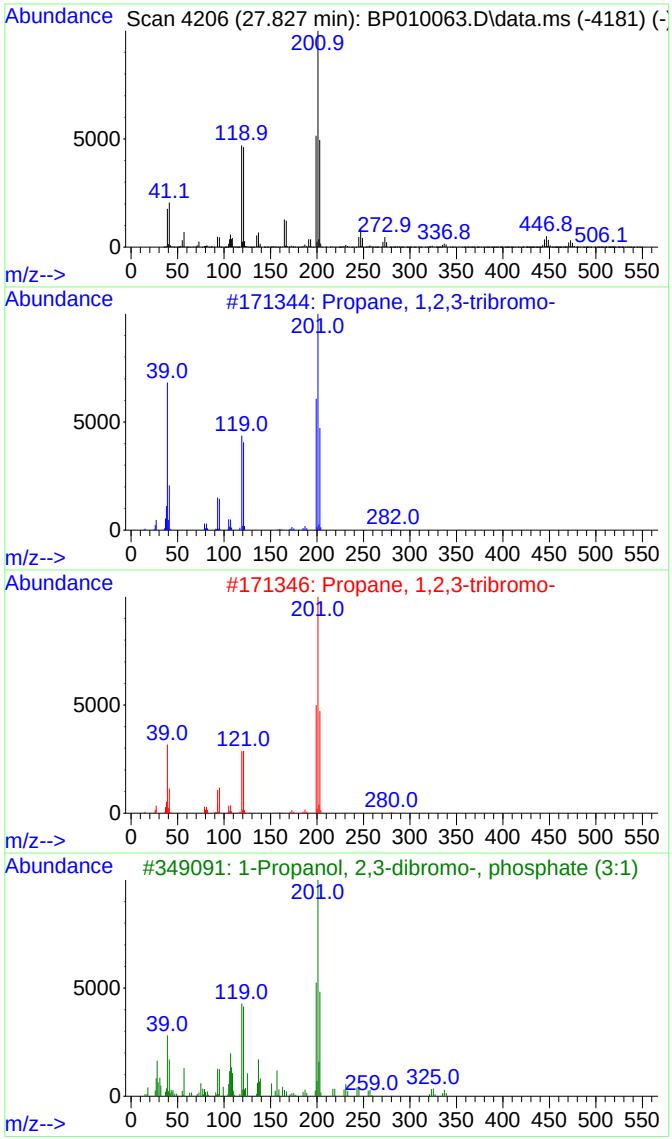
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

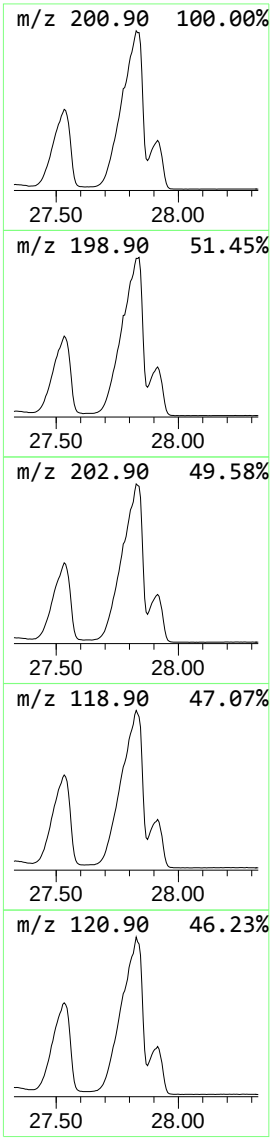
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 80 1-Propanol, 2,3-dibromo-, p... Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 27.827 | 120.32 ng/ul | 69710100 | Perylene-d12 | 23.886 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 86 |
| 2 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 78 |
| 3 | 1-Propanol, 2,3-dibromo-, phosph... | 692 | C9H15Br6O4P | 000126-72-7 | 64 |
| 4 | 2(3H)-Benzothiazolethione, 5-chl... | 201 | C7H4ClNS2 | 005331-91-9 | 43 |
| 5 | Propane, 1,2,2-tribromo- | 278 | C3H5Br3 | 014476-30-3 | 39 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010063.D
Acq On : 22 Apr 2022 20:00
Operator : CG/JU
Sample : N2316-14RE
Misc :
ALS Vial : 59 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
ETNR4RE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

| | | | | | --Internal Standard--- | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| Propane, 1-brom... | 5.534 | 47.6 | ng/ul | 6073390 | 1 | 7.940 | 2553830 | 20.0 |
| 2,3-Dibromoprop... | 9.369 | 1654.6 | ng/ul | 193427000 | 2 | 10.757 | 2338020 | 20.0 |
| Benzene, 1-brom... | 9.469 | 161.6 | ng/ul | 18889200 | 2 | 10.757 | 2338020 | 20.0 |
| (DEL) Alkane: C... | 9.534 | 14.4 | ng/ul | 1688240 | 2 | 10.757 | 2338020 | 20.0 |
| Benzene, 1-brom... | 9.775 | 314.4 | ng/ul | 36759000 | 2 | 10.757 | 2338020 | 20.0 |
| unknown-01 | 10.034 | 127.4 | ng/ul | 14892100 | 2 | 10.757 | 2338020 | 20.0 |
| Oxalic acid, cy... | 10.181 | 37.2 | ng/ul | 4352170 | 2 | 10.757 | 2338020 | 20.0 |
| unknown-02 | 10.322 | 320.4 | ng/ul | 37456300 | 2 | 10.757 | 2338020 | 20.0 |
| Propane, 1,2,3-... | 10.575 | 145.8 | ng/ul | 17049700 | 2 | 10.757 | 2338020 | 20.0 |
| (DEL) Alkane: S... | 11.128 | 346.0 | ng/ul | 40448200 | 2 | 10.757 | 2338020 | 20.0 |
| Hexane, 1,2-dib... | 11.216 | 32.8 | ng/ul | 3831640 | 2 | 10.757 | 2338020 | 20.0 |
| unknown-03 | 11.281 | 54.9 | ng/ul | 6414370 | 2 | 10.757 | 2338020 | 20.0 |
| unknown-04 | 11.398 | 124.4 | ng/ul | 14540500 | 2 | 10.757 | 2338020 | 20.0 |
| Hexane, 1,6-dib... | 11.522 | 49.1 | ng/ul | 5745430 | 2 | 10.757 | 2338020 | 20.0 |
| unknown-05 | 12.151 | 54.3 | ng/ul | 6348850 | 2 | 10.757 | 2338020 | 20.0 |
| unknown-06 | 12.193 | 53.3 | ng/ul | 6228350 | 2 | 10.757 | 2338020 | 20.0 |
| 2-Methyl-2-brom... | 12.410 | 335.6 | ng/ul | 39232900 | 2 | 10.757 | 2338020 | 20.0 |
| unknown-07 | 12.469 | 225.0 | ng/ul | 26302000 | 2 | 10.757 | 2338020 | 20.0 |
| unknown-08 | 13.051 | 70.8 | ng/ul | 8324470 | 3 | 14.581 | 2350850 | 20.0 |
| Benzene, 1,3-di... | 13.322 | 287.2 | ng/ul | 33754800 | 3 | 14.581 | 2350850 | 20.0 |
| unknown-09 | 13.598 | 1730.6 | ng/ul | 203422000 | 3 | 14.581 | 2350850 | 20.0 |
| unknown-10 | 14.151 | 38.6 | ng/ul | 4536740 | 3 | 14.581 | 2350850 | 20.0 |
| unknown-11 | 14.334 | 51.3 | ng/ul | 6027000 | 3 | 14.581 | 2350850 | 20.0 |
| unknown-12 | 15.892 | 53.7 | ng/ul | 6307720 | 3 | 14.581 | 2350850 | 20.0 |
| Propane, 1,2,2-... | 17.133 | 1276.7 | ng/ul | 399607000 | 4 | 17.357 | 6260190 | 20.0 |
| unknown-13 | 17.657 | 642.5 | ng/ul | 201100000 | 4 | 17.357 | 6260190 | 20.0 |
| unknown-14 | 18.480 | 115.7 | ng/ul | 36200300 | 4 | 17.357 | 6260190 | 20.0 |
| unknown-15 | 20.721 | 61.5 | ng/ul | 73380000 | 5 | 21.416 | 23879800 | 20.0 |
| unknown-16 | 22.392 | 89.2 | ng/ul | 106483000 | 5 | 21.416 | 23879800 | 20.0 |
| unknown-17 | 24.198 | 65.3 | ng/ul | 37838700 | 6 | 23.886 | 11587300 | 20.0 |
| unknown-18 | 27.533 | 63.7 | ng/ul | 36908900 | 6 | 23.886 | 11587300 | 20.0 |
| 1-Propanol, 2,3... | 27.827 | 120.3 | ng/ul | 69710100 | 6 | 23.886 | 11587300 | 20.0 |