LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** EXF01

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-BP080822.M
Title : SVOA CALIBRATION

Signal : TIC: BP011343.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 3.564 | 91 | 98 | 119 | rBV2 | 7294167 | 13826837 | 6.64% | 0.641% |
| 2 | 3.828 | 130 | 143 | 153 | rBV2 | 17908964 | 35965840 | 17.27% | 1.667% |
| 3 | 4.052 | 173 | 181 | 189 | rVV | 22754318 | 44207549 | 21.22% | 2.049% |
| 4 | 4.487 | 244 | 255 | 276 | rBV | 4084158 | 9092107 | 4.37% | 0.422% |
| 5 | 4.881 | 315 | 322 | 325 | rBV | 5625477 | 8734826 | 4.19% | 0.405% |
| 6 | 4.934 | 325 | 331 | 337 | rVB | 25345376 | 44195020 | 21.22% | 2.049% |
| 7 | 4.999 | 337 | 342 | 349 | rVB2 | 3675846 | 5491612 | 2.64% | 0.255% |
| 8 | 5.093 | 349 | 358 | 366 | rBV3 | 1437973 | 3057932 | 1.47% | 0.142% |
| 9 | 5.222 | 367 | 380 | 387 | rBV | 12027999 | 20155346 | 9.68% | 0.934% |
| 10 | 6.281 | 548 | 560 | 568 | rVB4 | 16007725 | 33827590 | 16.24% | 1.568% |
| 11 | 6.599 | 609 | 614 | 620 | rVB | 4865792 | 7227679 | 3.47% | 0.335% |
| 12 | 6.846 | 633 | 656 | 665 | rVV4 | 1861787 | 7321314 | 3.51% | 0.339% |
| 13 | 7.675 | 790 | 797 | 807 | rVV | 6726813 | 11238224 | 5.40% | 0.521% |
| 14 | 7.910 | 829 | 837 | 846 | rBV | 7070785 | 12987247 | 6.24% | 0.602% |
| 15 | 8.022 | 846 | 856 | 862 | rVV | 9400003 | 19943171 | 9.57% | 0.925% |
| 16 | 8.216 | 879 | 889 | 895 | rBV3 | 7628795 | 14708498 | 7.06% | 0.682% |
| 17 | 8.322 | 896 | 907 | 913 | rBV | 3385387 | 8695495 | 4.17% | 0.403% |
| 18 | 8.381 | 913 | 917 | 923 | rVB | 2333002 | 3945752 | 1.89% | 0.183% |
| 19 | 8.487 | 924 | 935 | 943 | rVB | 6077107 | 11233851 | 5.39% | 0.521% |
| 20 | 8.581 | 946 | 951 | 957 | rVB | 2163291 | 3438111 | 1.65% | 0.159% |
| 21 | 8.705 | 957 | 972 | 980 | rBV | 3924629 | 15441132 | 7.41% | 0.716% |
| 22 | 8.852 | 991 | 997 | 1001 | rBV | 2184331 | 3531579 | 1.70% | 0.164% |
| 23 | 8.969 | 1001 | 1017 | 1032 | rBV3 | 7235703 | 41593983 | 19.97% | 1.928% |
| 24 | 9.099 | 1032 | 1039 | 1044 | rVB | 13219168 | 21900796 | 10.51% | 1.015% |
| 25 | 9.163 | 1044 | 1050 | 1053 | rBV | 1980252 | 3667430 | 1.76% | 0.170% |
| 26 | 9.263 | 1054 | 1067 | 1085 | rVB | 9619125 | 37951568 | 18.22% | 1.759% |
| 27 | 9.416 | 1085 | 1093 | 1094 | rBV | 20302257 | 33259077 | 15.97% | 1.542% |
| 28 | 9.540 | 1105 | 1114 | 1117 | rBV2 | 3066075 | 6296584 | 3.02% | 0.292% |
| 29 | 9.828 | 1130 | 1163 | 1181 | rBV3 | 23009806 | 208295363 | 100.00% | 9.657% |
| 30 | 9.999 | 1181 | 1192 | 1193 | rBV | 9995384 | 24722190 | 11.87% | 1.146% |
| 31 | 10.469 | 1260 | 1272 | 1277 | rBV | 13005588 | 23305598 | 11.19% | 1.080% |
| 32 | 10.699 | 1300 | 1311 | 1320 | rVB | 13510391 | 29335482 | 14.08% | 1.360% |
| 33 | 10.822 | 1321 | 1332 | 1334 | rBV | 18465320 | 43008831 | 20.65% | 1.994% |
| 34 | 10.904 | 1341 | 1346 | 1353 | rVB | 1759276 | 3649525 | 1.75% | 0.169% |
| 35 | 11.046 | 1353 | 1370 | 1373 | rBV2 | 33297276 | 112551583 | 54.03% | 5.218% |
| 36 | 11.140 | 1373 | 1386 | 1390 | rVV2 | 30707540 | 115701733 | 55.55% | 5.364% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

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-BP080822.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 11.187 | 1390 | 1394 | 1396 | rVV2 | 2875670 | 5116505 | 2.46% | 0.237% |
| 38 | 11.216 | 1396 | 1399 | 1412 | rVB | 4174446 | 6740974 | 3.24% | 0.313% |
| 39 | 11.522 | 1441 | 1451 | 1457 | rBV | 2018176 | 5278981 | 2.53% | 0.245% |
| 40 | 11.740 | 1481 | 1488 | 1495 | rVV | 7588349 | 15038679 | 7.22% | 0.697% |
| 41 | 11.834 | 1495 | 1504 | 1508 | rVV4 | 8249210 | 18345575 | 8.81% | 0.850% |
| 42 | 11.881 | 1508 | 1512 | 1519 | rVV | 7434987 | 11842670 | 5.69% | 0.549% |
| 43 | 12.093 | 1530 | 1548 | 1553 | rVV | 15993094 | 30584440 | 14.68% | 1.418% |
| 44 | 12.216 | 1553 | 1569 | 1573 | rVV2 | 33270422 | 121273575 | 58.22% | 5.622% |
| 45 | 12.287 | 1573 | 1581 | 1586 | rVV3 | 2308142 | 6758599 | 3.24% | 0.313% |
| 46 | 12.357 | 1586 | 1593 | 1606 | rVV3 | 2922443 | 8166912 | 3.92% | 0.379% |
| 47 | 12.475 | 1606 | 1613 | 1615 | rVV | 2529119 | 5685849 | 2.73% | 0.264% |
| 48 | 12.516 | 1615 | 1620 | 1627 | rVV2 | 6418107 | 14216032 | 6.82% | 0.659% |
| 49 | 12.657 | 1635 | 1644 | 1655 | rVV2 | 9123344 | 21716426 | 10.43% | 1.007% |
| 50 | 12.757 | 1655 | 1661 | 1667 | rVV4 | 3307946 | 8591829 | 4.12% | 0.398% |
| 51 | 12.816 | 1667 | 1671 | 1681 | rVB2 | 2991053 | 5146651 | 2.47% | 0.239% |
| 52 | 12.998 | 1694 | 1702 | 1708 | rVB2 | 5133491 | 12318512 | 5.91% | 0.571% |
| 53 | 13.281 | 1733 | 1750 | 1751 | rBV2 | 34776455 | 133725489 | 64.20% | 6.199% |
| 54 | 13.451 | 1774 | 1779 | 1788 | rVB | 9471361 | 12880174 | 6.18% | 0.597% |
| 55 | 13.616 | 1802 | 1807 | 1819 | rVV | 2431765 | 4061100 | 1.95% | 0.188% |
| 56 | 13.763 | 1821 | 1832 | 1842 | rVV2 | 20885788 | 47708942 | 22.90% | 2.212% |
| 57 | 13.892 | 1848 | 1854 | 1862 | rVB2 | 9440979 | 15301484 | 7.35% | 0.709% |
| 58 | 13.998 | 1868 | 1872 | 1876 | rVV | 1774148 | 2798903 | 1.34% | 0.130% |
| 59 | 14.081 | 1876 | 1886 | 1892 | rVB | 13337325 | 23048344 | 11.07% | 1.069% |
| 60 | 14.192 | 1899 | 1905 | 1907 | rBV | 3186071 | 4830449 | 2.32% | 0.224% |
| 61 | 14.263 | 1911 | 1917 | 1920 | rVV | 14029963 | 21940546 | 10.53% | 1.017% |
| 62 | 14.357 | 1927 | 1933 | 1939 | rVV2 | 5525241 | 8657053 | 4.16% | 0.401% |
| 63 | 14.434 | 1939 | 1946 | 1953 | rVV3 | 3784025 | 6171495 | 2.96% | 0.286% |
| 64 | 14.616 | 1970 | 1977 | 1981 | rBV | 14395988 | 20958083 | 10.06% | 0.972% |
| 65 | 14.669 | 1981 | 1986 | 1990 | rVV3 | 2369373 | 4978995 | 2.39% | 0.231% |
| 66 | 14.757 | 1995 | 2001 | 2010 | rVB2 | 1759440 | 3857525 | 1.85% | 0.179% |
| 67 | 14.945 | 2030 | 2033 | 2041 | rVB2 | 1873505 | 3163921 | 1.52% | 0.147% |
| 68 | 15.216 | 2070 | 2079 | 2083 | rBV3 | 1538918 | 5180721 | 2.49% | 0.240% |
| 69 | 15.281 | 2086 | 2090 | 2093 | rVB2 | 2442601 | 3672254 | 1.76% | 0.170% |
| 70 | 15.716 | 2150 | 2164 | 2167 | rBV | 2395358 | 6268159 | 3.01% | 0.291% |
| 71 | 15.839 | 2180 | 2185 | 2188 | rBV | 2222788 | 3756928 | 1.80% | 0.174% |
| 72 | 16.104 | 2226 | 2230 | 2238 | rVB | 6060283 | 10359961 | 4.97% | 0.480% |
| 73 | 16.334 | 2264 | 2269 | 2275 | rVB2 | 3103309 | 5165939 | 2.48% | 0.239% |
| 74 | 16.728 | 2312 | 2336 | 2340 | rBV6 | 29779242 | 125226477 | 60.12% | 5.805% |
| 75 | 17.010 | 2378 | 2384 | 2388 | rVB2 | 2752884 | 5882505 | 2.82% | 0.273% |
| 76 | 17.157 | 2405 | 2409 | 2414 | rBV | 2089307 | 3044170 | 1.46% | 0.141% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

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-BP080822.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 77 | 17.310 | 2426 | 2435 | 2440 | rBV | 25912589 | 44385347 | 21.31% | 2.058% |
| 78 | 17.363 | 2440 | 2444 | 2448 | rVB | 3086847 | 3928441 | 1.89% | 0.182% |
| 79 | 18.869 | 2693 | 2700 | 2703 | rVV2 | 2034751 | 5043214 | 2.42% | 0.234% |
| 80 | 19.398 | 2785 | 2790 | 2795 | rBV | 2577260 | 3947132 | 1.89% | 0.183% |
| 81 | 19.486 | 2795 | 2805 | 2813 | rVV | 20020380 | 32936386 | 15.81% | 1.527% |
| 82 | 20.304 | 2933 | 2944 | 2946 | rBV | 27608064 | 48819126 | 23.44% | 2.263% |
| 83 | 20.327 | 2946 | 2948 | 2951 | rVV | 19840410 | 22301362 | 10.71% | 1.034% |
| 84 | 20.357 | 2951 | 2953 | 2957 | rVV3 | 4488400 | 6746895 | 3.24% | 0.313% |
| 85 | 20.469 | 2965 | 2972 | 2980 | rVB2 | 16031727 | 24823904 | 11.92% | 1.151% |
| 86 | 21.027 | 3062 | 3067 | 3070 | rBV3 | 1969886 | 2829785 | 1.36% | 0.131% |
| 87 | 21.069 | 3070 | 3074 | 3077 | rBV2 | 2724664 | 3705740 | 1.78% | 0.172% |
| 88 | 21.151 | 3083 | 3088 | 3100 | rBV2 | 3087091 | 5162891 | 2.48% | 0.239% |
| 89 | 21.869 | 3206 | 3210 | 3218 | rVB | 2388234 | 3770675 | 1.81% | 0.175% |
| 90 | 22.063 | 3236 | 3243 | 3250 | rVB | 23767552 | 40395877 | 19.39% | 1.873% |
| 91 | 23.145 | 3420 | 3427 | 3431 | rBV | 3480152 | 7473876 | 3.59% | 0.346% |
| 92 | 23.180 | 3431 | 3433 | 3436 | rVV | 2474100 | 3749806 | 1.80% | 0.174% |
| 93 | 23.221 | 3436 | 3440 | 3447 | rVB | 5277437 | 8513452 | 4.09% | 0.395% |
| 94 | 23.327 | 3452 | 3458 | 3463 | rBV3 | 1753036 | 3392031 | 1.63% | 0.157% |
| 95 | 23.410 | 3467 | 3472 | 3478 | rVB | 2193913 | 4038493 | 1.94% | 0.187% |
| 96 | 25.027 | 3740 | 3747 | 3759 | rVB | 1664496 | 3859543 | 1.85% | 0.179% |
| 97 | 26.121 | 3928 | 3933 | 3949 | rVB | 1937996 | 5910706 | 2.84% | 0.274% |
| 98 | 26.298 | 3955 | 3963 | 3969 | rBV | 1148150 | 3359067 | 1.61% | 0.156% |
| 99 | 26.827 | 4039 | 4053 | 4070 | rVB | 4145005 | 14906386 | 7.16% | 0.691% |
| 100 | 27.080 | 4079 | 4096 | 4103 | rBV | 7916362 | 30071802 | 14.44% | 1.394% |

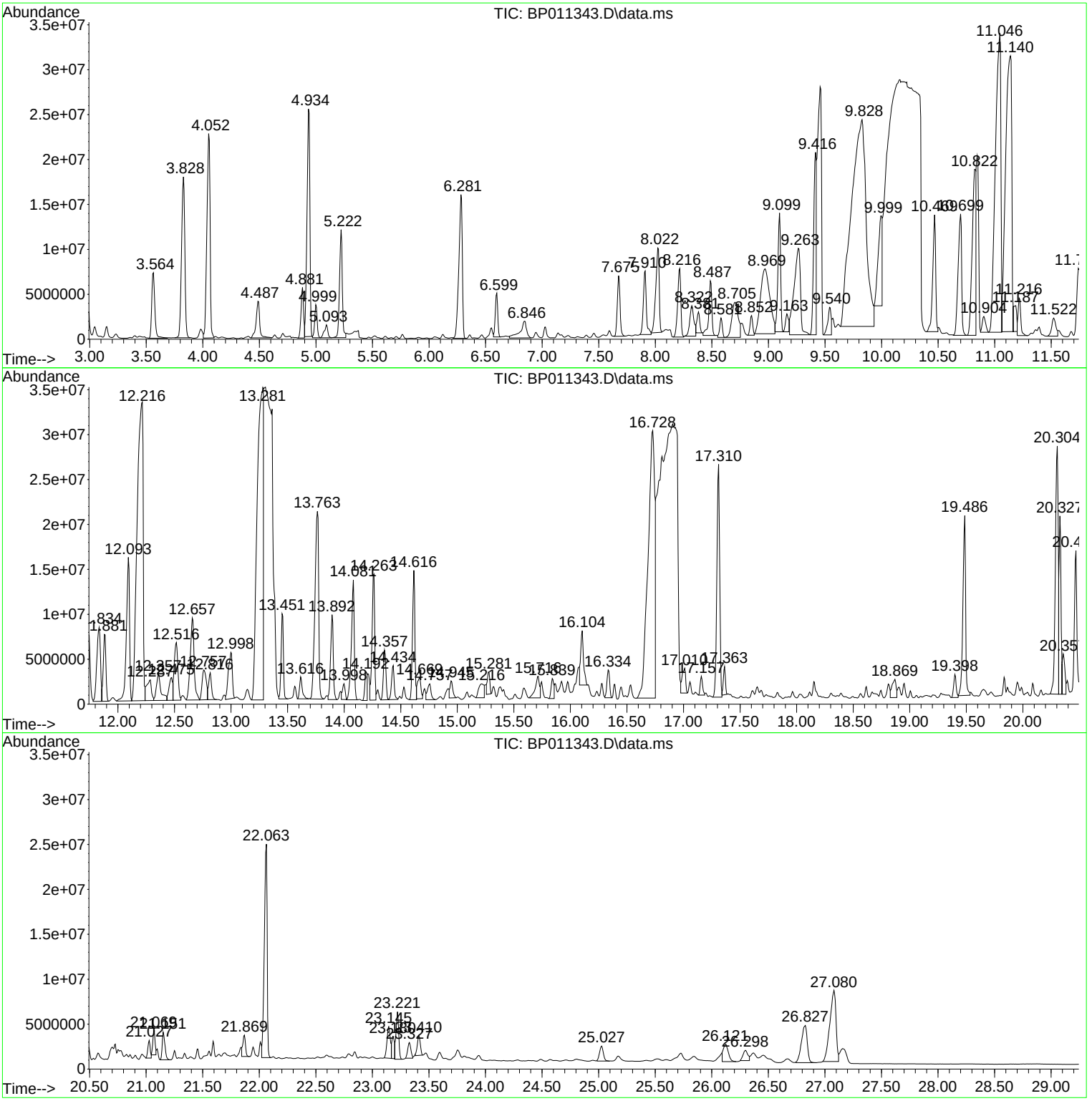
Sum of corrected areas: 2157040218

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

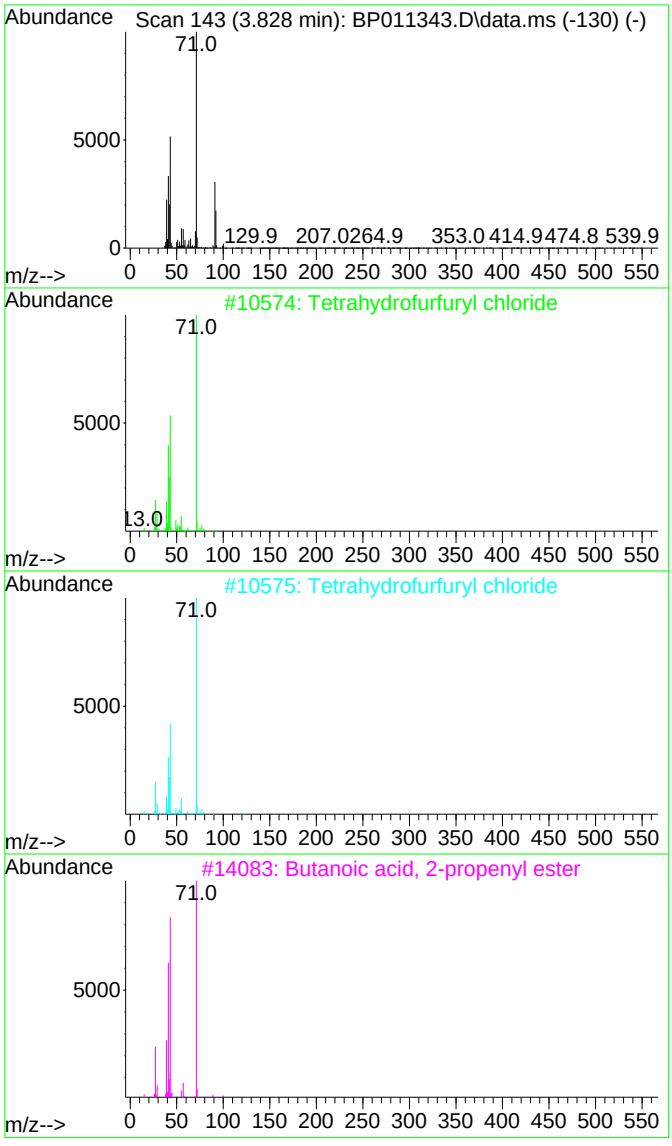
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

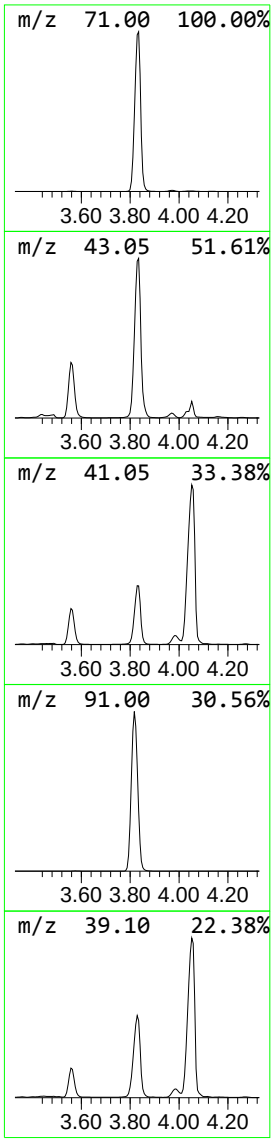
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 2 Tetrahydrofurfuryl chloride Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 3.828 | 64.01 ng/ul | 35965800 | 1,4-Dichlorobenzene-d4 | 7.593 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Tetrahydrofurfuryl chloride | 120 | C5H9ClO | 003003-84-7 | 59 |
| 2 | Tetrahydrofurfuryl chloride | 120 | C5H9ClO | 003003-84-7 | 59 |
| 3 | Butanoic acid, 2-propenyl ester | 128 | C7H12O2 | 002051-78-7 | 50 |
| 4 | 2-Ethyltetrahydrofuran | 100 | C6H12O | 1010431-64-0 | 50 |
| 5 | Butanoic acid, 2-propenyl ester | 128 | C7H12O2 | 002051-78-7 | 50 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

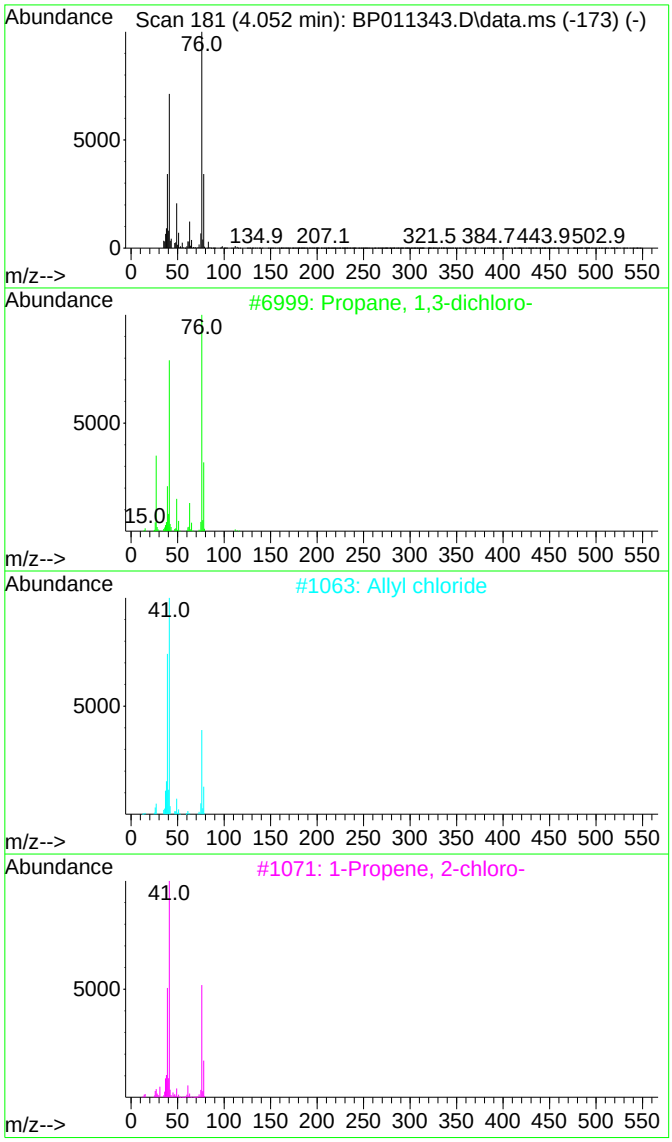
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

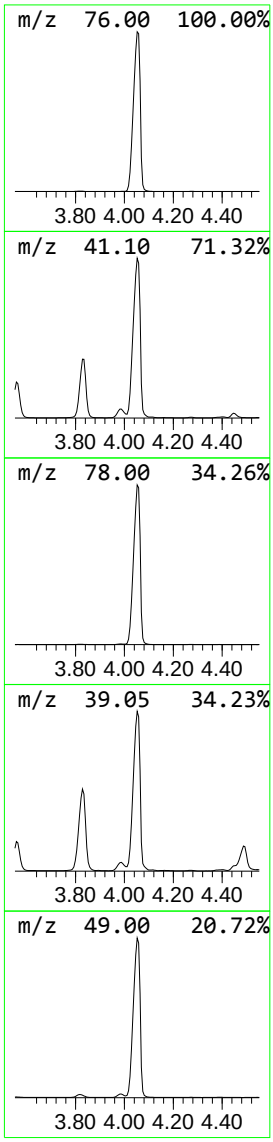
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 3 Propane, 1,3-dichloro- Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.052 | 78.67 ng/ul | 44207500 | 1,4-Dichlorobenzene-d4 | 7.593 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 1,3-dichloro- | 112 | C3H6Cl2 | 000142-28-9 | 91 |
| 2 | Allyl chloride | 76 | C3H5Cl | 000107-05-1 | 72 |
| 3 | 1-Propene, 2-chloro- | 76 | C3H5Cl | 000557-98-2 | 64 |
| 4 | Propane, 1,3-dichloro- | 112 | C3H6Cl2 | 000142-28-9 | 64 |
| 5 | 1-Propene, 2-chloro- | 76 | C3H5Cl | 000557-98-2 | 59 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

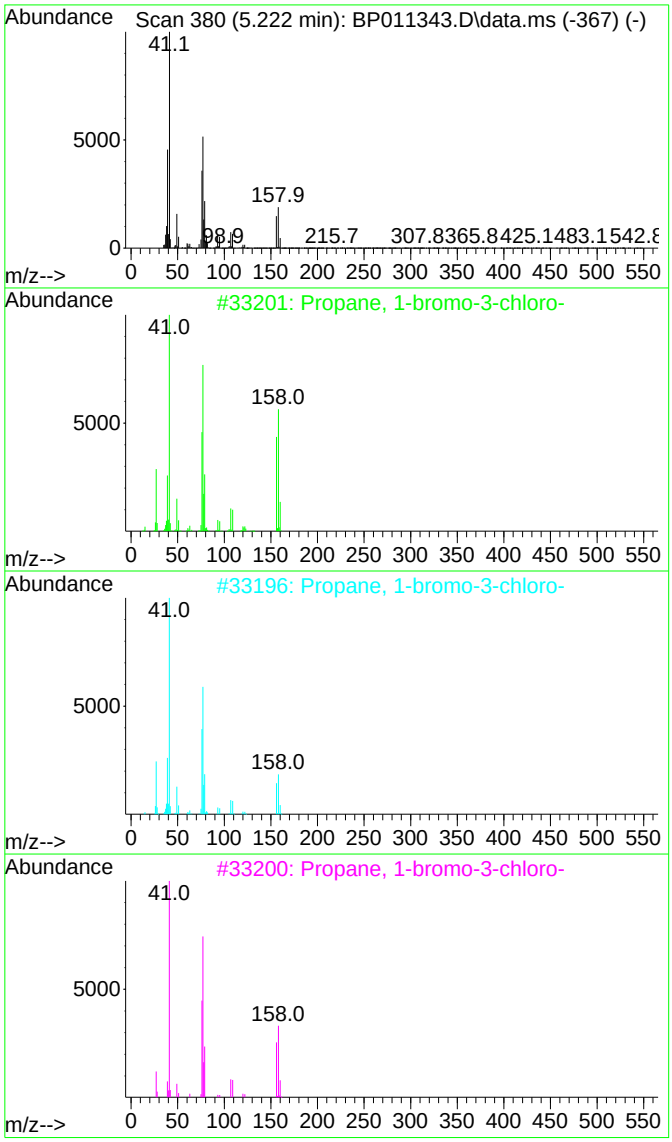
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

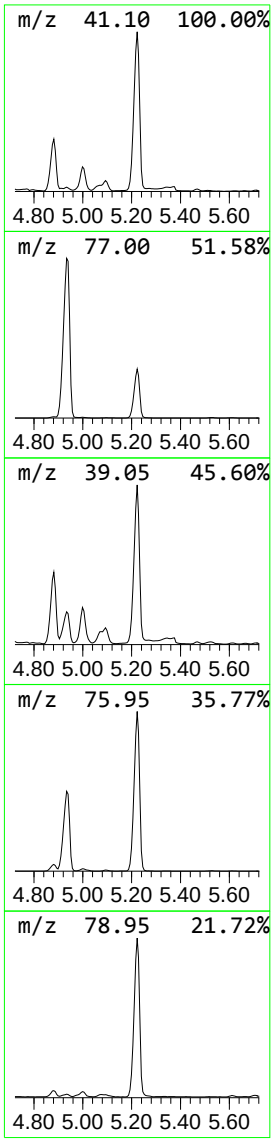
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 9 Propane, 1-bromo-3-chloro- Concentration Rank 23

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.222 | 35.87 ng/ul | 20155300 | 1,4-Dichlorobenzene-d4 | 7.593 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 1-bromo-3-chloro- | 156 | C3H6BrCl | 000109-70-6 | 95 |
| 2 | Propane, 1-bromo-3-chloro- | 156 | C3H6BrCl | 000109-70-6 | 95 |
| 3 | Propane, 1-bromo-3-chloro- | 156 | C3H6BrCl | 000109-70-6 | 60 |
| 4 | Propane, 1-bromo-3-chloro- | 156 | C3H6BrCl | 000109-70-6 | 55 |
| 5 | Butane, 2-chloro-2-methyl- | 106 | C5H11Cl | 000594-36-5 | 45 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

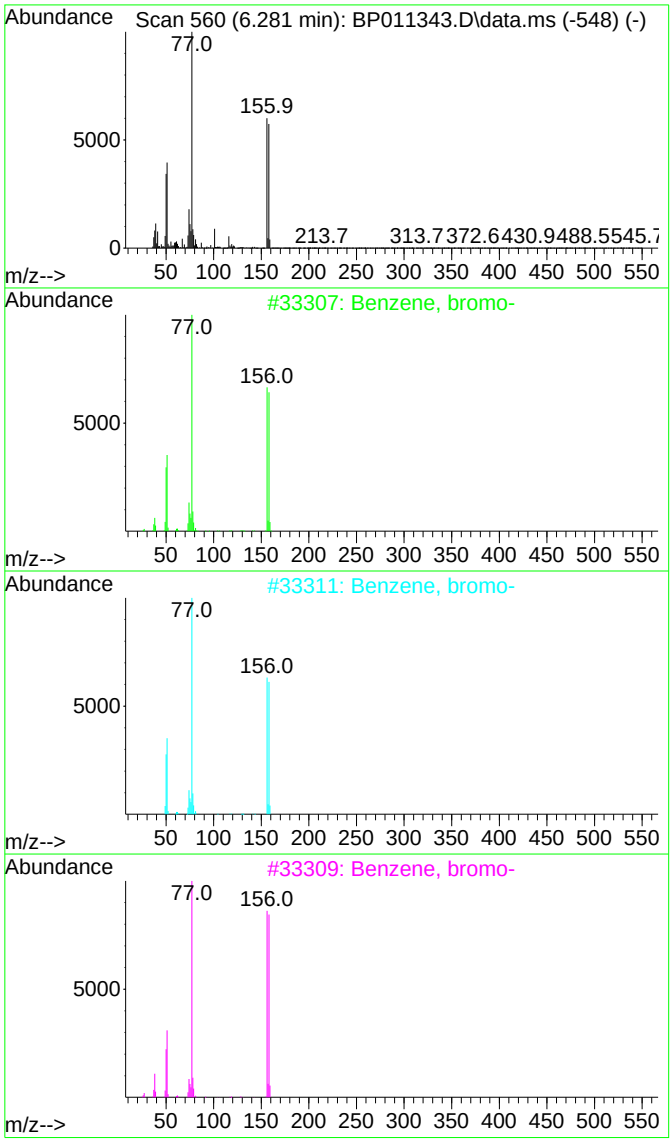
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

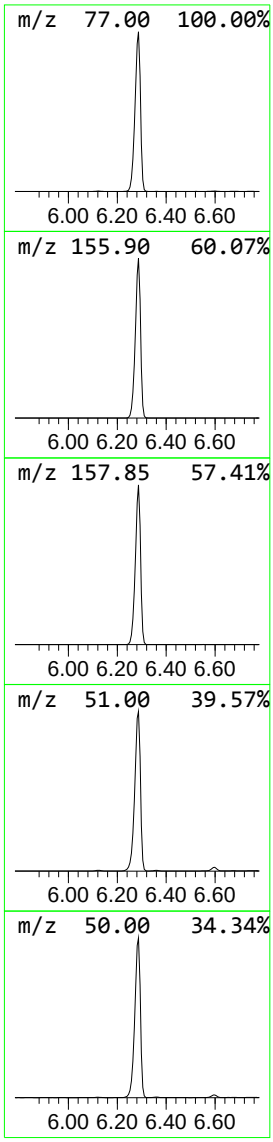
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 10 Benzene, bromo- Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.281 | 60.20 ng/ul | 33827600 | 1,4-Dichlorobenzene-d4 | 7.593 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, bromo- | 156 | C6H5Br | 000108-86-1 | 96 |
| 2 | Benzene, bromo- | 156 | C6H5Br | 000108-86-1 | 96 |
| 3 | Benzene, bromo- | 156 | C6H5Br | 000108-86-1 | 94 |
| 4 | Benzene, bromo- | 156 | C6H5Br | 000108-86-1 | 91 |
| 5 | Benzene, bromo- | 156 | C6H5Br | 000108-86-1 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

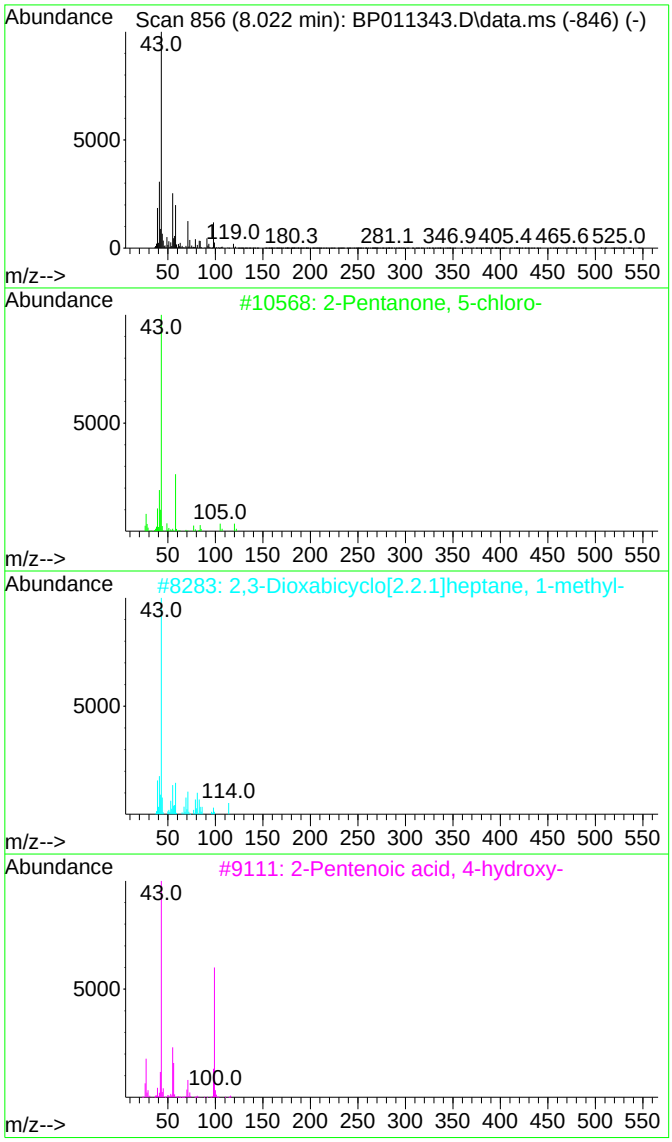
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

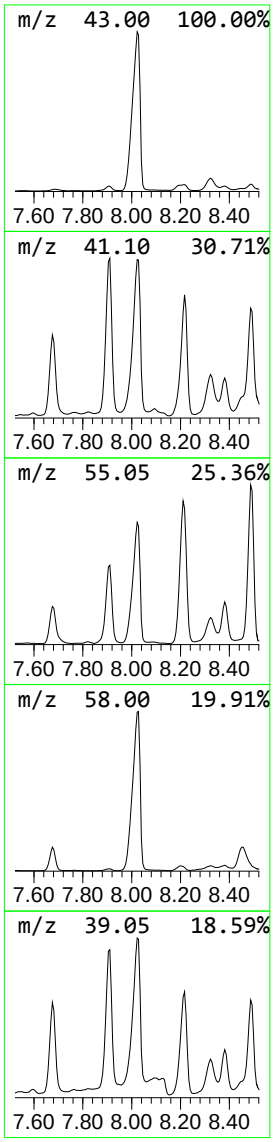
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 14 unknown-01 Concentration Rank 24

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.022 | 35.49 ng/ul | 19943200 | 1,4-Dichlorobenzene-d4 | 7.593 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Pentanone, 5-chloro- | 120 | C5H9ClO | 005891-21-4 | 37 |
| 2 | 2,3-Dioxabicyclo[2.2.1]heptane, ... | 114 | C6H10O2 | 1010142-34-6 | 33 |
| 3 | 2-Pentenoic acid, 4-hydroxy- | 116 | C5H8O3 | 028525-83-9 | 25 |
| 4 | Butane | 58 | C4H10 | 000106-97-8 | 23 |
| 5 | 3-Hexene-2,5-diol | 116 | C6H12O2 | 007319-23-5 | 23 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

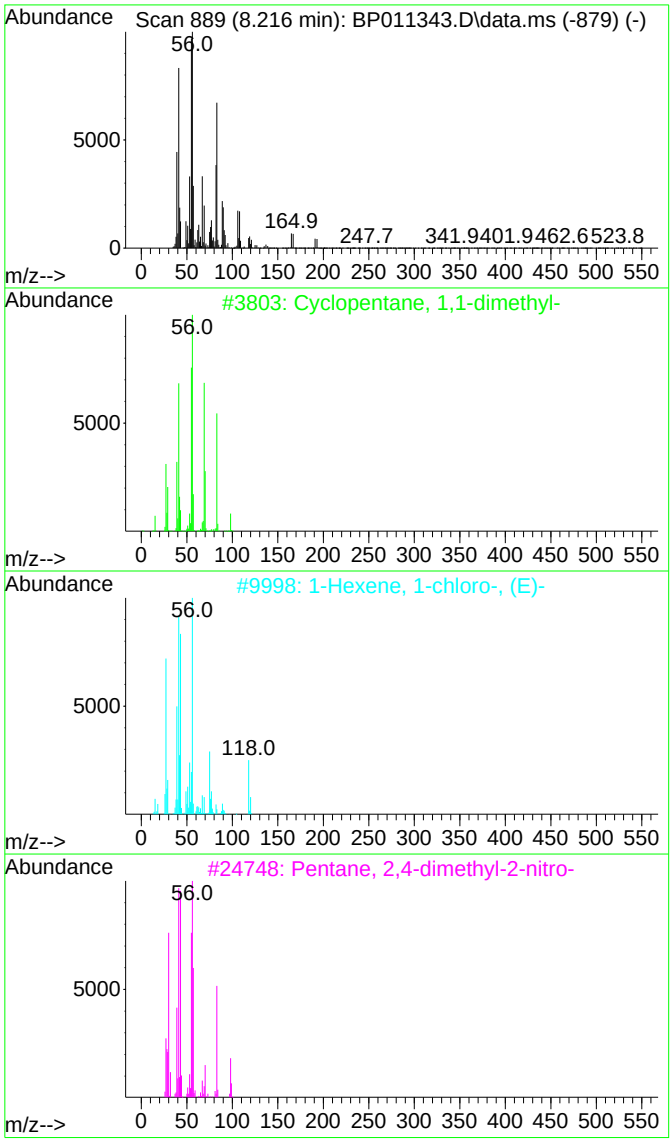
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

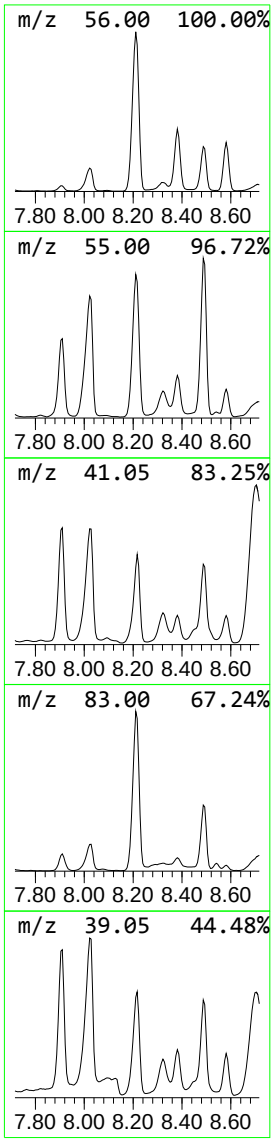
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 15 (DEL) Alkane: Cyclic8.216 Concentration Rank 31

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.216 | 26.18 ng/ul | 14708500 | 1,4-Dichlorobenzene-d4 | 7.593 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclopentane, 1,1-dimethyl- | 98 | C7H14 | 001638-26-2 | 38 |
| 2 | 1-Hexene, 1-chloro-, (E)- | 118 | C6H11Cl | 050586-19-1 | 27 |
| 3 | Pentane, 2,4-dimethyl-2-nitro- | 145 | C7H15NO2 | 000597-45-5 | 25 |
| 4 | Hexane, 1,6-dichloro- | 154 | C6H12Cl2 | 002163-00-0 | 25 |
| 5 | Cyclopropane, 2-chloro-1-ethyl-1... | 118 | C6H11Cl | 343926-09-0 | 16 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

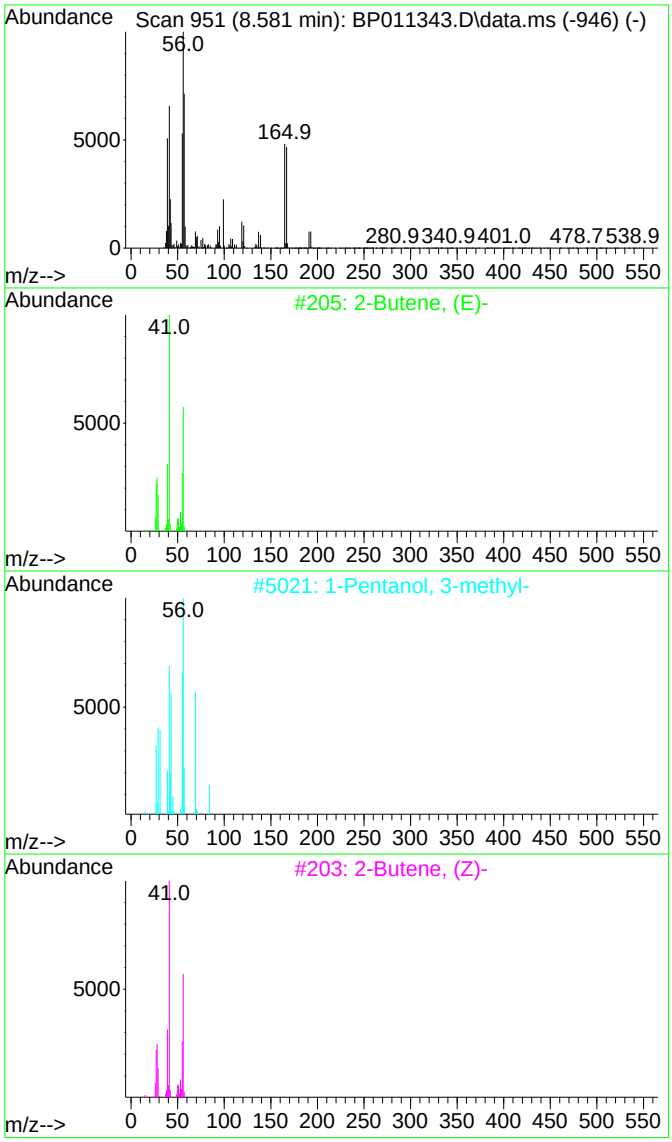
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

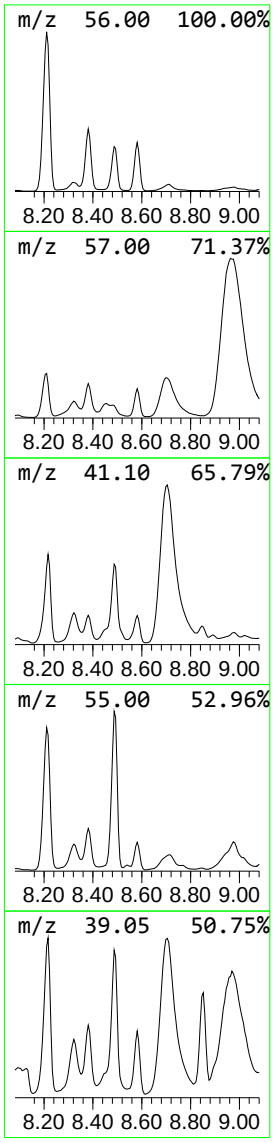
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 70

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.581 | 6.12 ng/ul | 3438110 | 1,4-Dichlorobenzene-d4 | 7.593 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Butene, (E)- | 56 | C4H8 | 000624-64-6 | 35 |
| 2 | 1-Pentanol, 3-methyl- | 102 | C6H14O | 000589-35-5 | 22 |
| 3 | 2-Butene, (Z)- | 56 | C4H8 | 000590-18-1 | 17 |
| 4 | Acetic acid, bromo-, butyl ester | 194 | C6H11BrO2 | 018991-98-5 | 12 |
| 5 | 1,3,2-Dioxaborinane, 2-ethyl-5,5... | 142 | C7H15BO2 | 057186-59-1 | 12 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

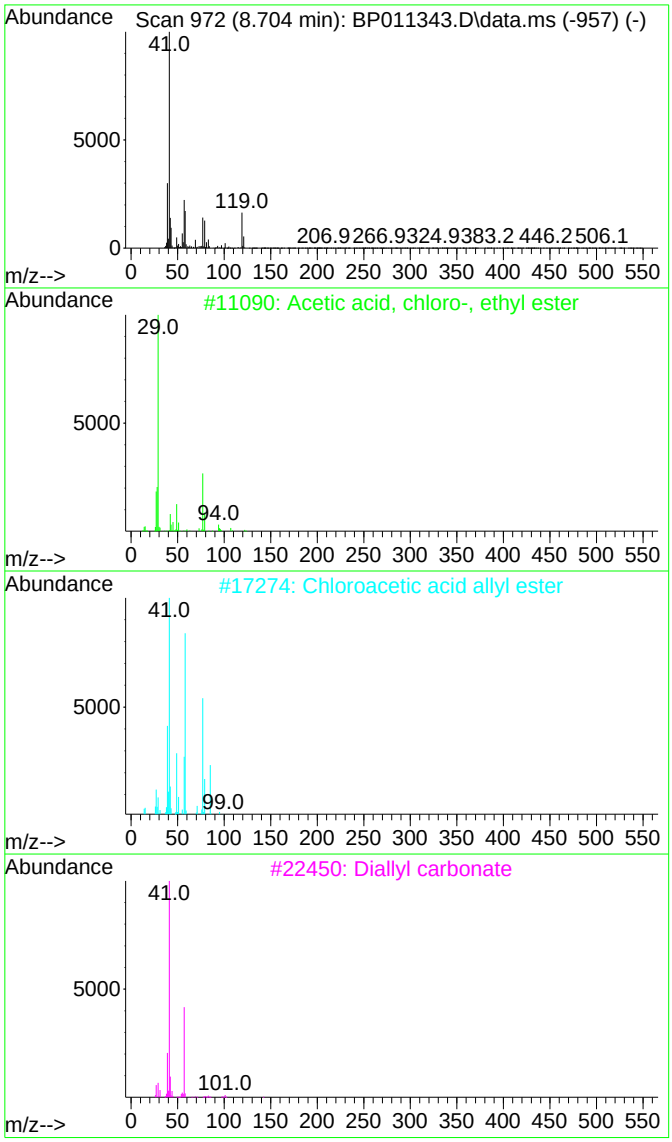
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

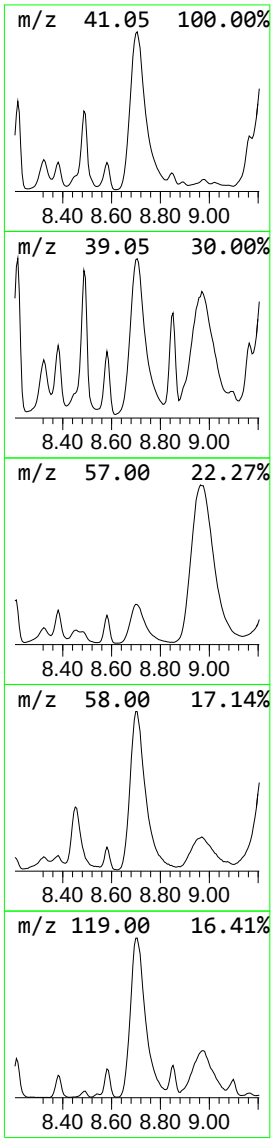
**********

Peak Number 18 unknown-02 Concentration Rank 29

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.704 | 27.48 ng/ul | 15441100 | 1,4-Dichlorobenzene-d4 | 7.593 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Acetic acid, chloro-, ethyl ester | 122 | C4H7ClO2 | 000105-39-5 | 9 |
| 2 | Chloroacetic acid allyl ester | 134 | C5H7ClO2 | 002916-14-5 | 9 |
| 3 | Diallyl carbonate | 142 | C7H10O3 | 015022-08-9 | 9 |
| 4 | o-Allylhydroxylamine | 73 | C3H7NO | 006542-54-7 | 5 |
| 5 | 2-Carbomethoxyaziridine | 101 | C4H7NO2 | 005950-34-5 | 4 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

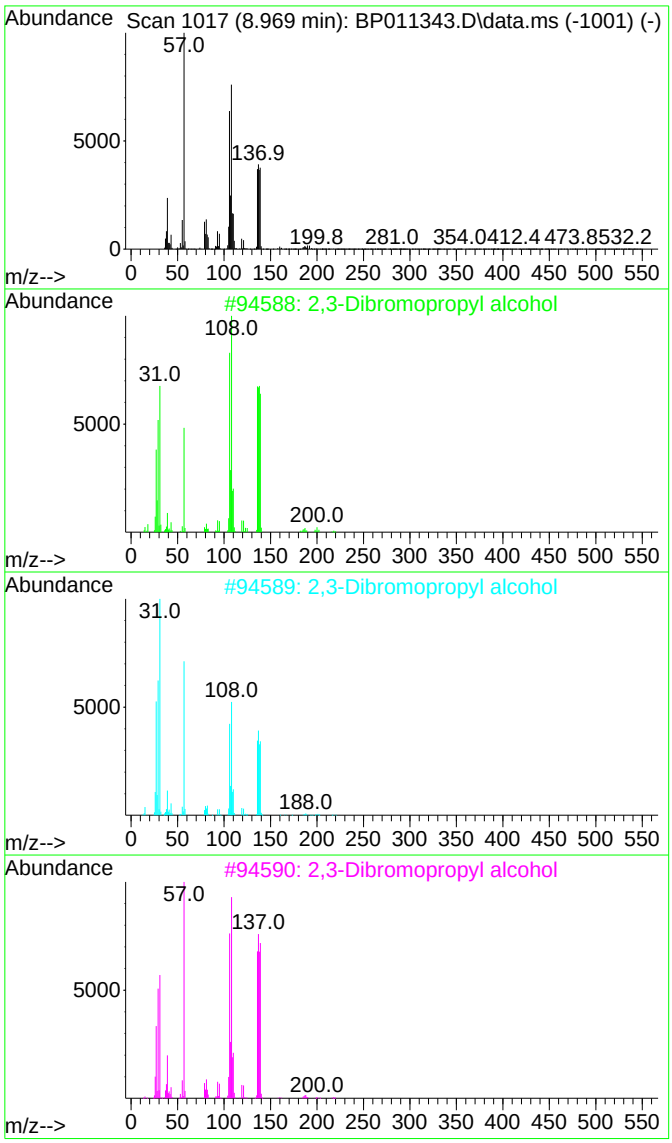
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

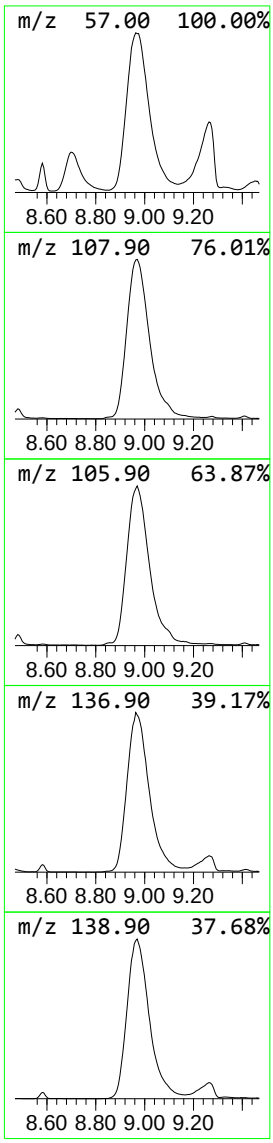
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 20 2,3-Dibromopropyl alcohol Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.969 | 74.02 ng/ul | 41594000 | 1,4-Dichlorobenzene-d4 | 7.593 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2,3-Dibromopropyl alcohol | 216 | C3H6Br2O | 000096-13-9 | 86 |
| 2 | 2,3-Dibromopropyl alcohol | 216 | C3H6Br2O | 000096-13-9 | 70 |
| 3 | 2,3-Dibromopropyl alcohol | 216 | C3H6Br2O | 000096-13-9 | 70 |
| 4 | Vinyl bromide | 106 | C2H3Br | 000593-60-2 | 38 |
| 5 | Vinyl bromide | 106 | C2H3Br | 000593-60-2 | 25 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

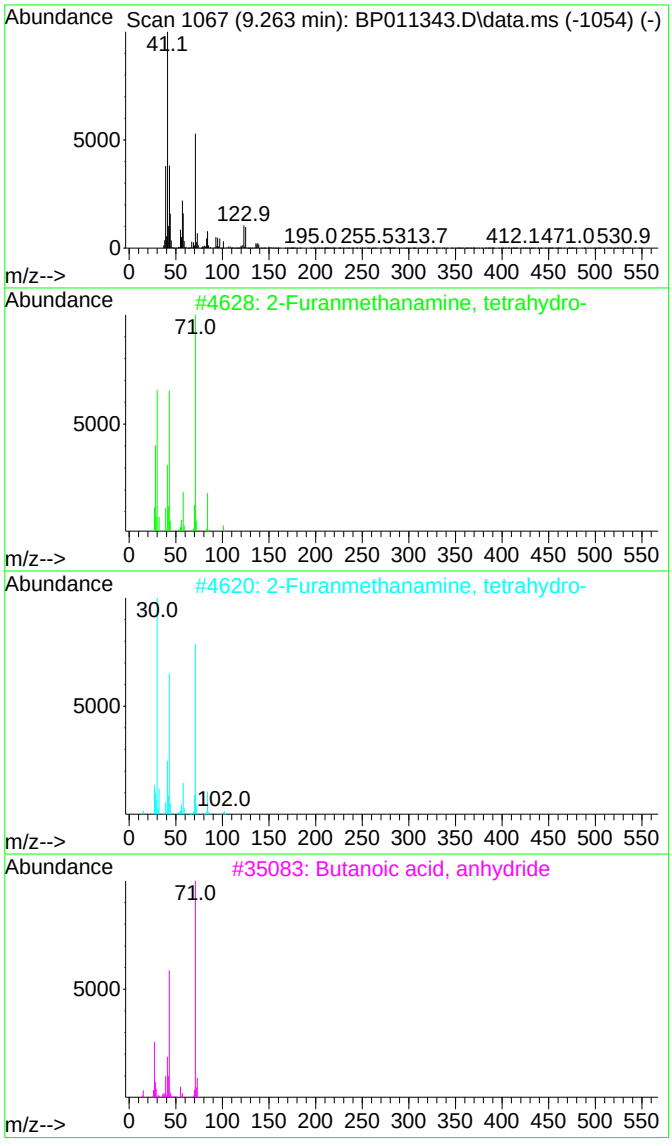
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

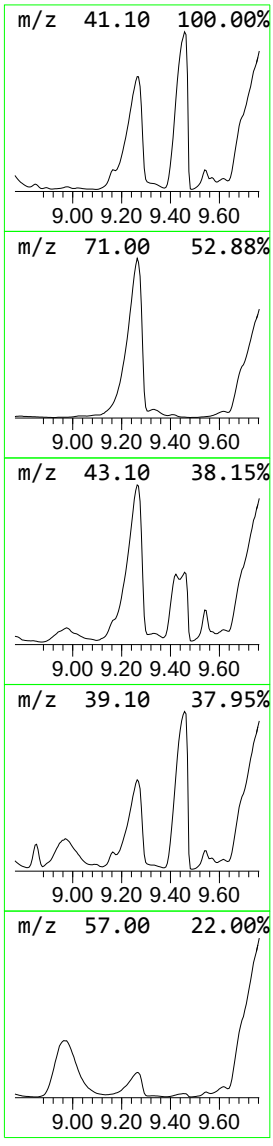
**********

Peak Number 22 unknown-03 Concentration Rank 26

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.263 | 32.57 ng/ul | 37951600 | Naphthalene-d8 | 10.428 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Furanmethanamine, tetrahydro- | 101 | C5H11NO | 004795-29-3 | 49 |
| 2 | 2-Furanmethanamine, tetrahydro- | 101 | C5H11NO | 004795-29-3 | 38 |
| 3 | Butanoic acid, anhydride | 158 | C8H14O3 | 000106-31-0 | 38 |
| 4 | dl-Mevalonic acid lactone | 130 | C6H10O3 | 000674-26-0 | 38 |
| 5 | Furan-2-carbonyl chloride, tetra... | 134 | C5H7ClO2 | 052449-98-6 | 35 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

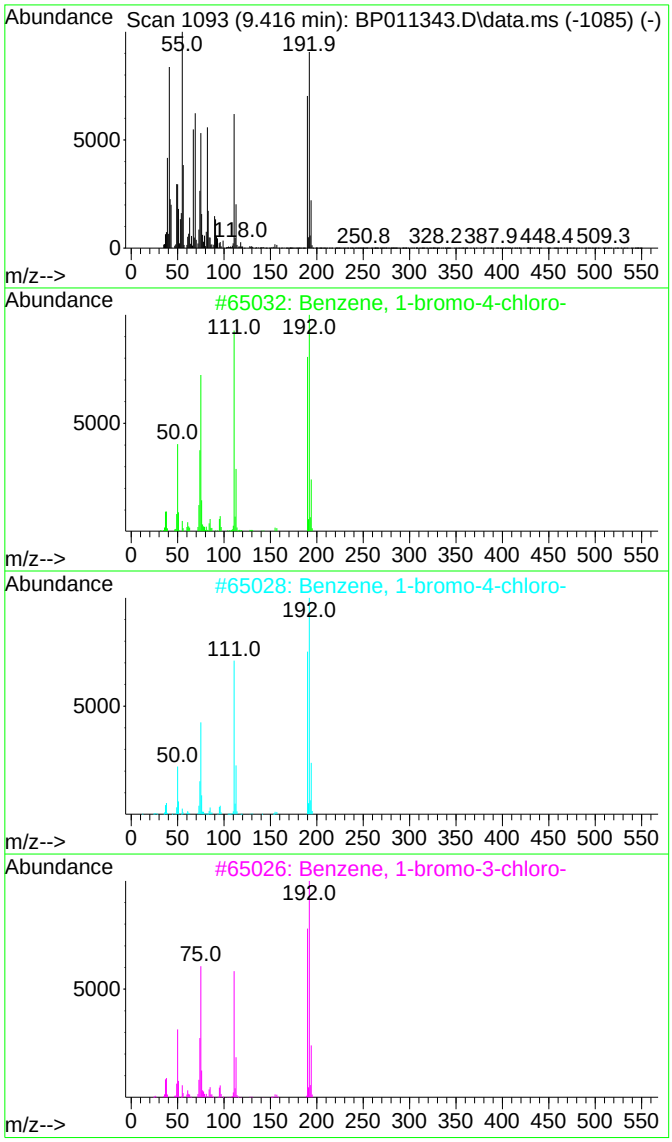
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

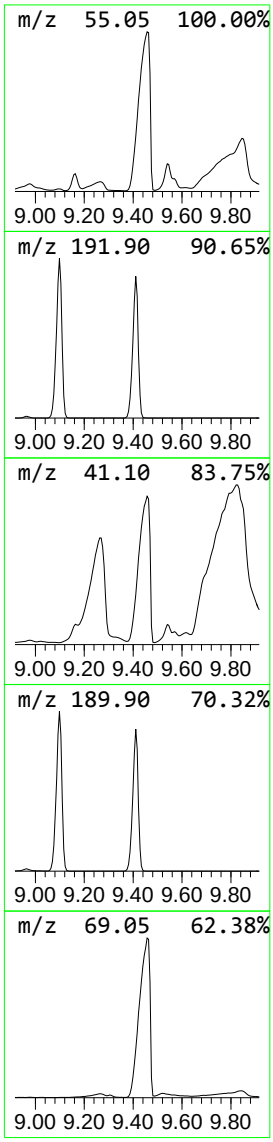
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 23 Benzene, 1-bromo-4-chloro- Concentration Rank 28

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.416 | 28.54 ng/ul | 33259100 | Naphthalene-d8 | 10.428 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1-bromo-4-chloro- | 190 | C6H4BrCl | 000106-39-8 | 97 |
| 2 | Benzene, 1-bromo-4-chloro- | 190 | C6H4BrCl | 000106-39-8 | 91 |
| 3 | Benzene, 1-bromo-3-chloro- | 190 | C6H4BrCl | 000108-37-2 | 89 |
| 4 | Benzene, 1-bromo-2-chloro- | 190 | C6H4BrCl | 000694-80-4 | 89 |
| 5 | Benzene, 1-bromo-4-chloro- | 190 | C6H4BrCl | 000106-39-8 | 87 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

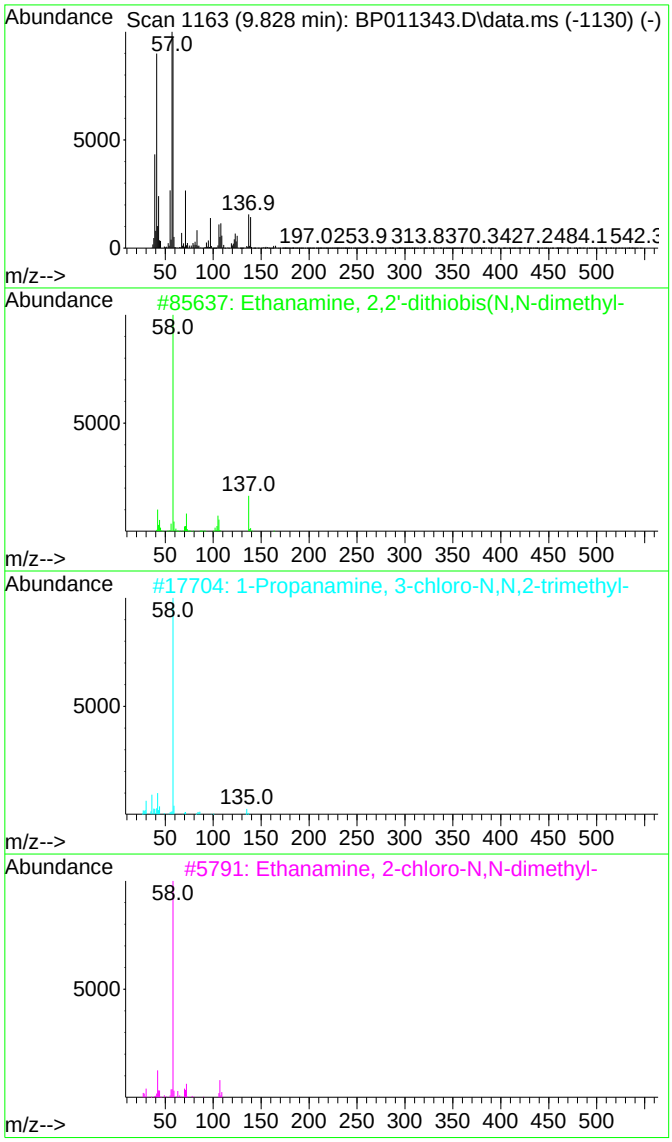
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

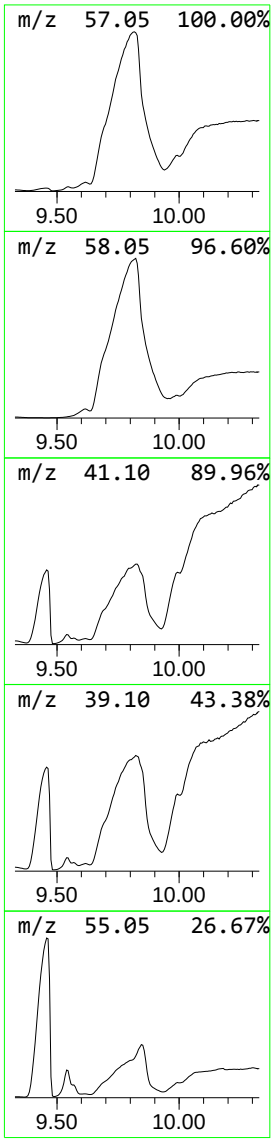
**********

Peak Number 24 unknown-04 Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.828 | 178.75 ng/ul | 208295000 | Naphthalene-d8 | 10.428 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Ethanamine, 2,2'-dithiobis(N,N-d... | 208 | C8H20N2S2 | 001072-11-3 | 37 |
| 2 | 1-Propanamine, 3-chloro-N,N,2-tr... | 135 | C6H14ClN | 023349-86-2 | 30 |
| 3 | Ethanamine, 2-chloro-N,N-dimethyl- | 107 | C4H10ClN | 000107-99-3 | 27 |
| 4 | Bis(2-(Dimethylamino)ethyl) ether | 160 | C8H20N2O | 003033-62-3 | 25 |
| 5 | Carbonic acid, 2-dimethylaminoet... | 203 | C10H21NO3 | 1000331-43-1 | 25 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

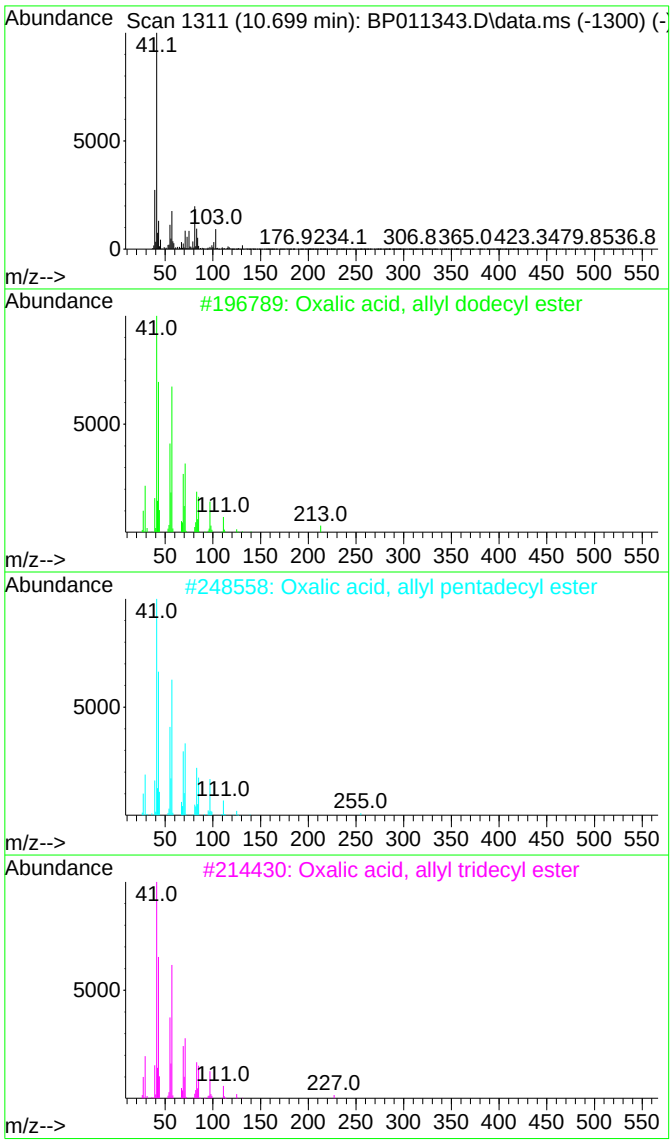
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

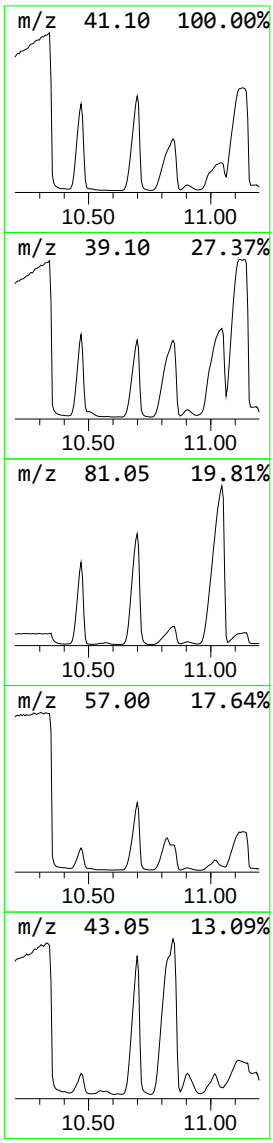
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 26 unknown-05 Concentration Rank 33

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.698 | 25.17 ng/ul | 29335500 | Naphthalene-d8 | 10.428 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Oxalic acid, allyl dodecyl ester | 298 | C17H30O4 | 1000309-24-0 | 12 |
| 2 | Oxalic acid, allyl pentadecyl ester | 340 | C20H36O4 | 1000309-24-3 | 12 |
| 3 | Oxalic acid, allyl tridecyl ester | 312 | C18H32O4 | 1000309-24-1 | 12 |
| 4 | Oxalic acid, allyl hexadecyl ester | 354 | C21H38O4 | 1000309-24-4 | 12 |
| 5 | 1-Propene, 3-(ethenyloxy)- | 84 | C5H8O | 003917-15-5 | 9 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

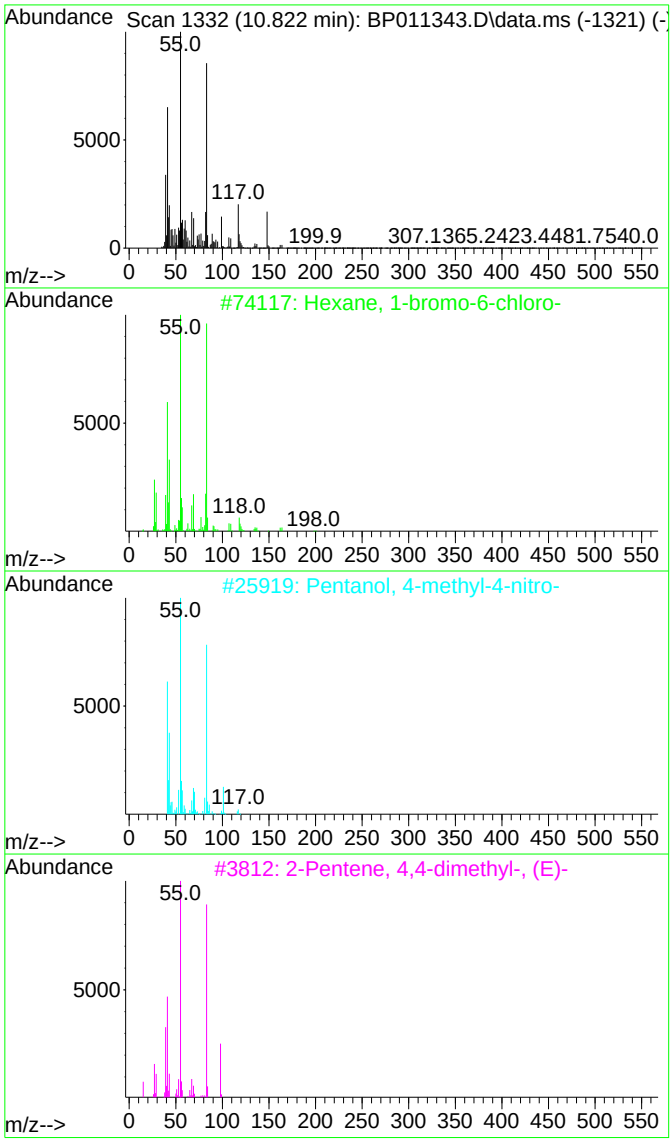
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

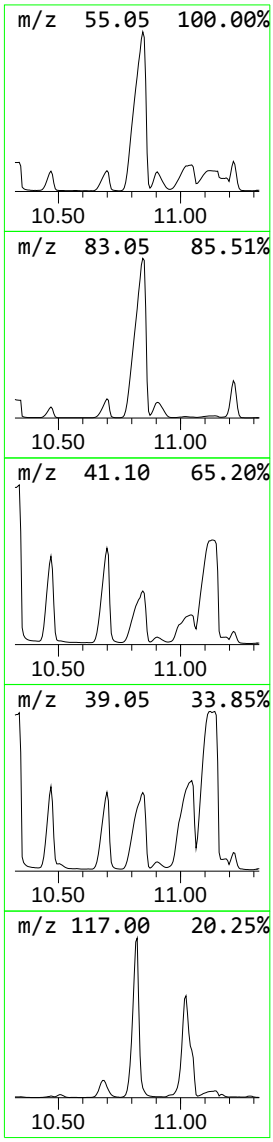
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 27 Hexane, 1-bromo-6-chloro- Concentration Rank 22

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.822 | 36.91 ng/ul | 43008800 | Naphthalene-d8 | 10.428 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexane, 1-bromo-6-chloro- | 198 | C6H12BrCl | 006294-17-3 | 92 |
| 2 | Pentanol, 4-methyl-4-nitro- | 147 | C6H13NO3 | 005215-92-9 | 50 |
| 3 | 2-Pentene, 4,4-dimethyl-, (E)- | 98 | C7H14 | 000690-08-4 | 47 |
| 4 | Benzene, 2-chloro-1,4-bis(cycloh... | 364 | C20H25ClO4 | 294874-04-7 | 47 |
| 5 | Cyclohexane, bromo- | 162 | C6H11Br | 000108-85-0 | 47 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

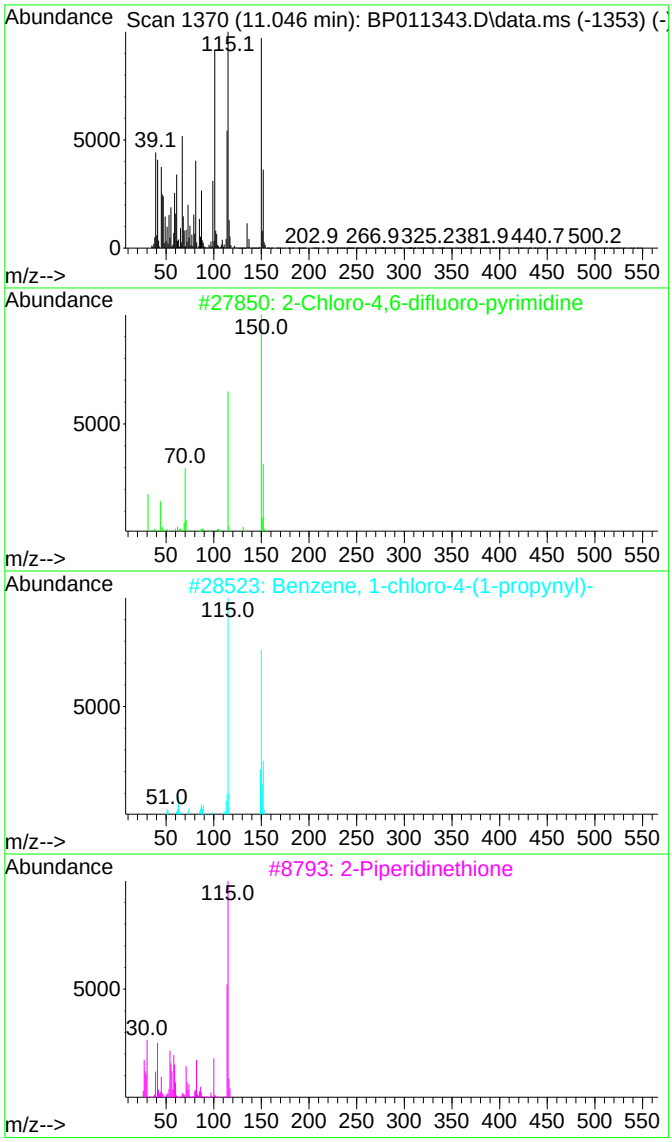
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

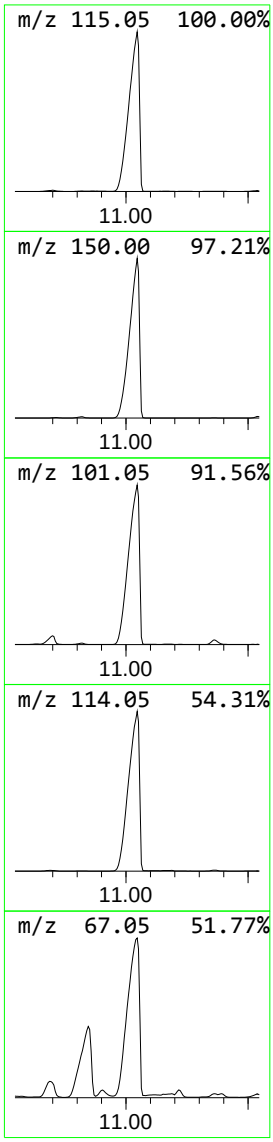
**********

Peak Number 28 unknown-06 Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.046 | 96.59 ng/ul | 112552000 | Naphthalene-d8 | 10.428 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Chloro-4,6-difluoro-pyrimidine | 150 | C4HClF2N2 | 038953-30-9 | 35 |
| 2 | Benzene, 1-chloro-4-(1-propynyl)- | 150 | C9H7Cl | 002809-65-6 | 14 |
| 3 | 2-Piperidinethione | 115 | C5H9NS | 013070-01-4 | 14 |
| 4 | 3(2H)-Isothiazolone, 2-methyl- | 115 | C4H5NOS | 002682-20-4 | 14 |
| 5 | 1-Methyl-1H-pyrazolo[3,4-d]pyrim... | 150 | C6H6N4O | 005334-56-5 | 10 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

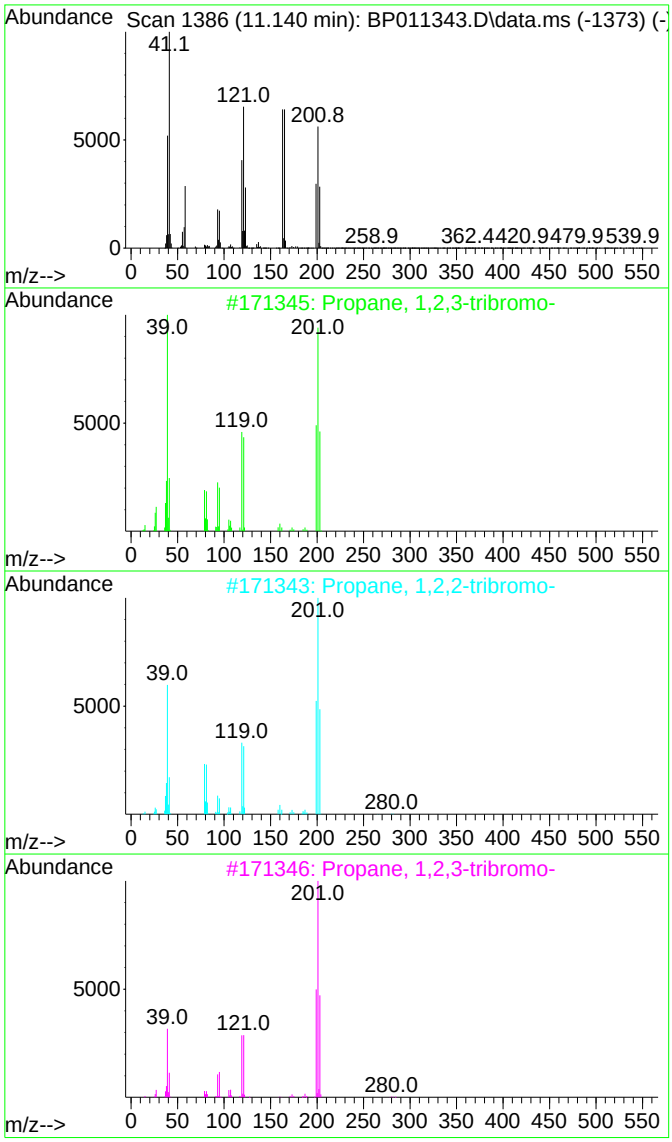
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

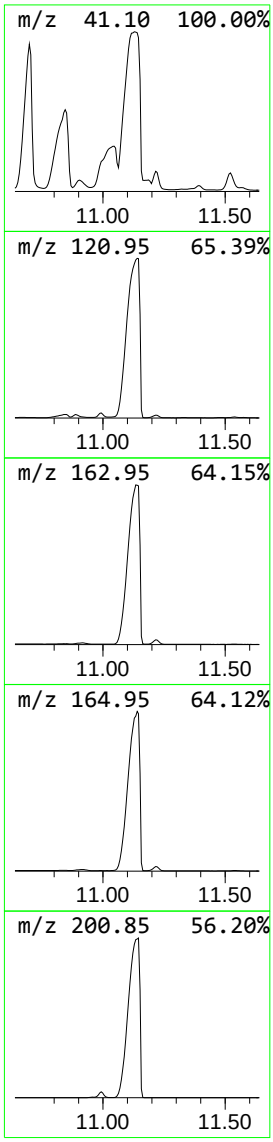
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 29 unknown-07 Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.140 | 99.29 ng/ul | 115702000 | Naphthalene-d8 | 10.428 |

| 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 | 23 |
| 3 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 16 |
| 4 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 12 |
| 5 | 2(3H)-Benzothiazolethione, 5-chl... | 201 | C7H4ClNS2 | 005331-91-9 | 9 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

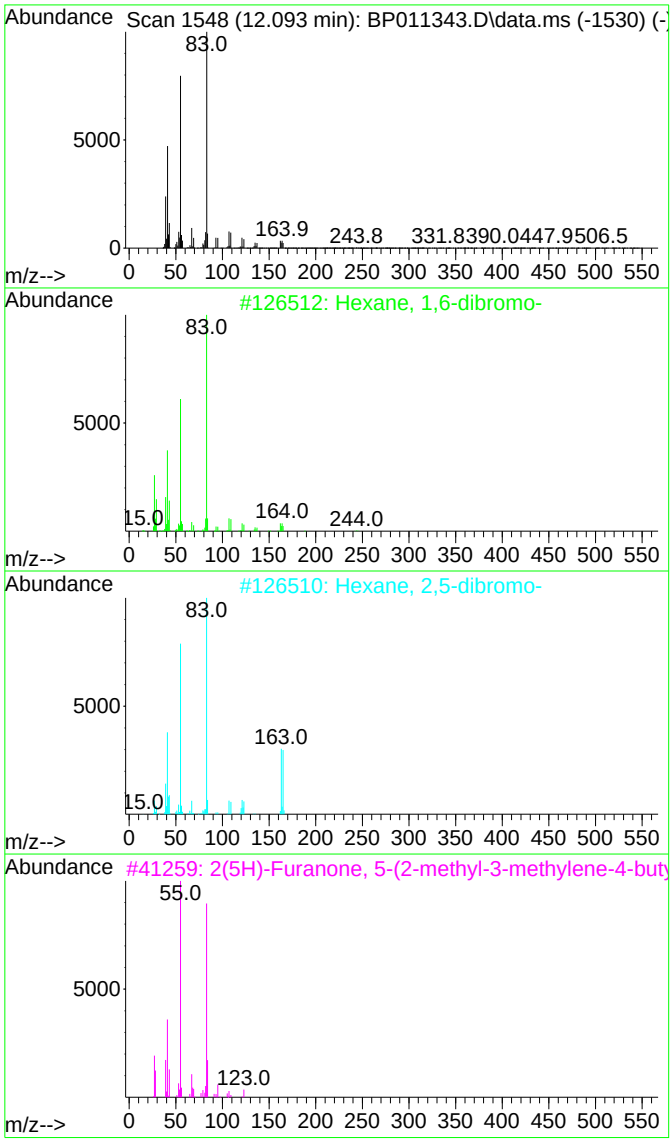
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

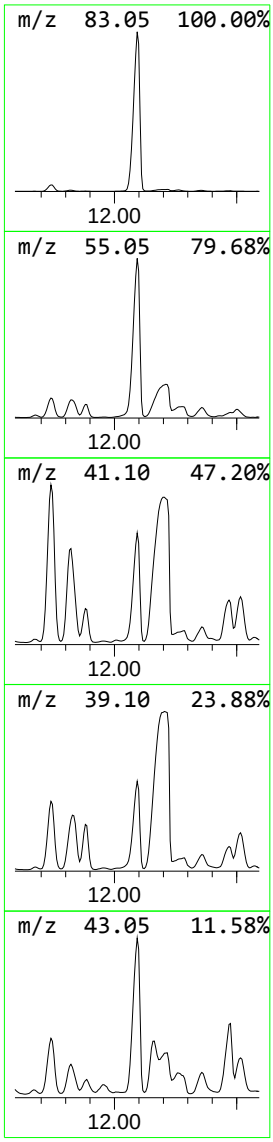
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 35 Hexane, 1,6-dibromo- Concentration Rank 30

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.093 | 26.25 ng/ul | 30584400 | Naphthalene-d8 | 10.428 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexane, 1,6-dibromo- | 242 | C6H12Br2 | 000629-03-8 | 94 |
| 2 | Hexane, 2,5-dibromo- | 242 | C6H12Br2 | 024774-58-1 | 83 |
| 3 | 2(5H)-Furanone, 5-(2-methyl-3-me... | 166 | C10H14O2 | 1000155-85-8 | 72 |
| 4 | Cyclohexane, bromo- | 162 | C6H11Br | 000108-85-0 | 68 |
| 5 | Oxalic acid, cyclohexyl hexyl ester | 256 | C14H24O4 | 1000309-30-8 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

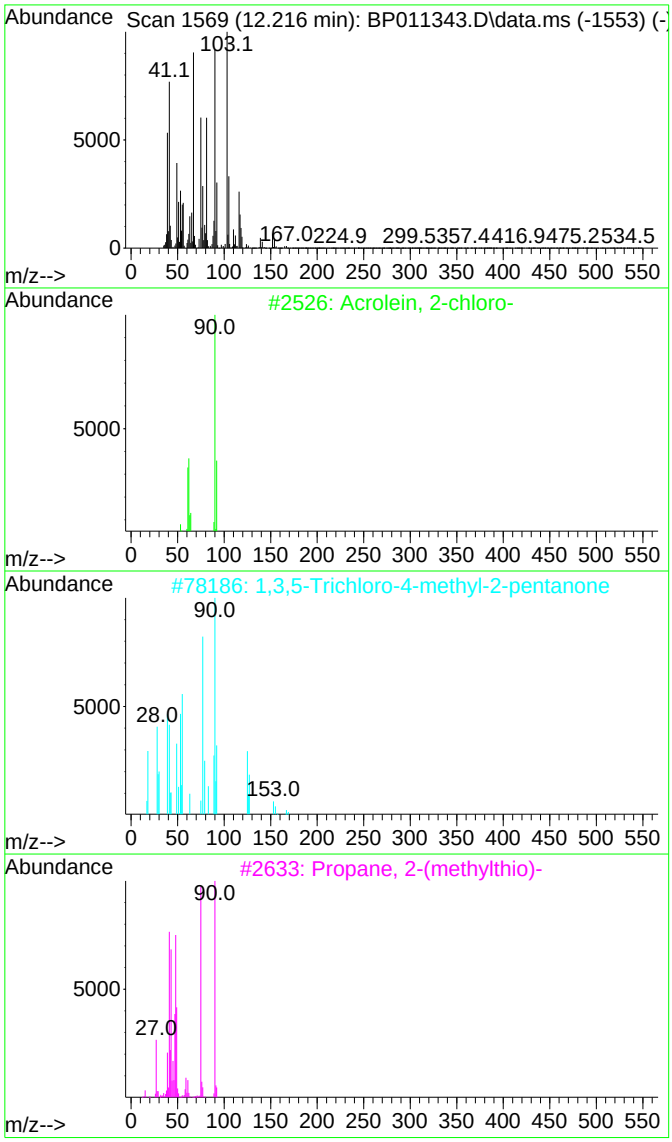
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

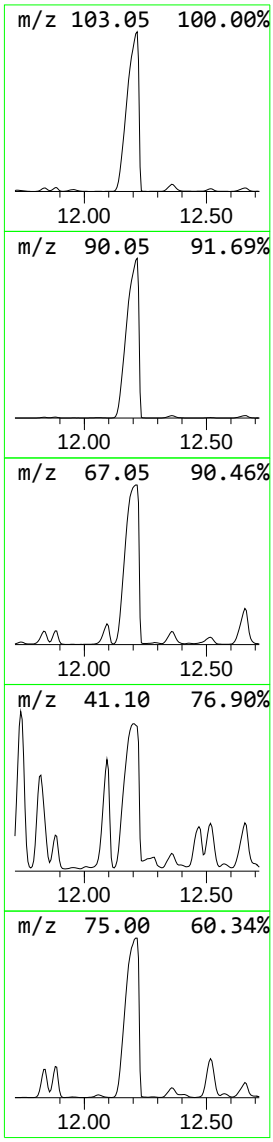
**********

Peak Number 36 unknown-08 Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.216 | 104.07 ng/ul | 121274000 | Naphthalene-d8 | 10.428 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Acrolein, 2-chloro- | 90 | C3H3ClO | 000683-51-2 | 37 |
| 2 | 1,3,5-Trichloro-4-methyl-2-penta... | 202 | C6H9Cl3O | 1000140-03-8 | 33 |
| 3 | Propane, 2-(methylthio)- | 90 | C4H10S | 001551-21-9 | 32 |
| 4 | Bicyclo[4.1.0]hepta-1,3,5-triene | 90 | C7H6 | 004646-69-9 | 30 |
| 5 | Thiourea, methyl- | 90 | C2H6N2S | 000598-52-7 | 27 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

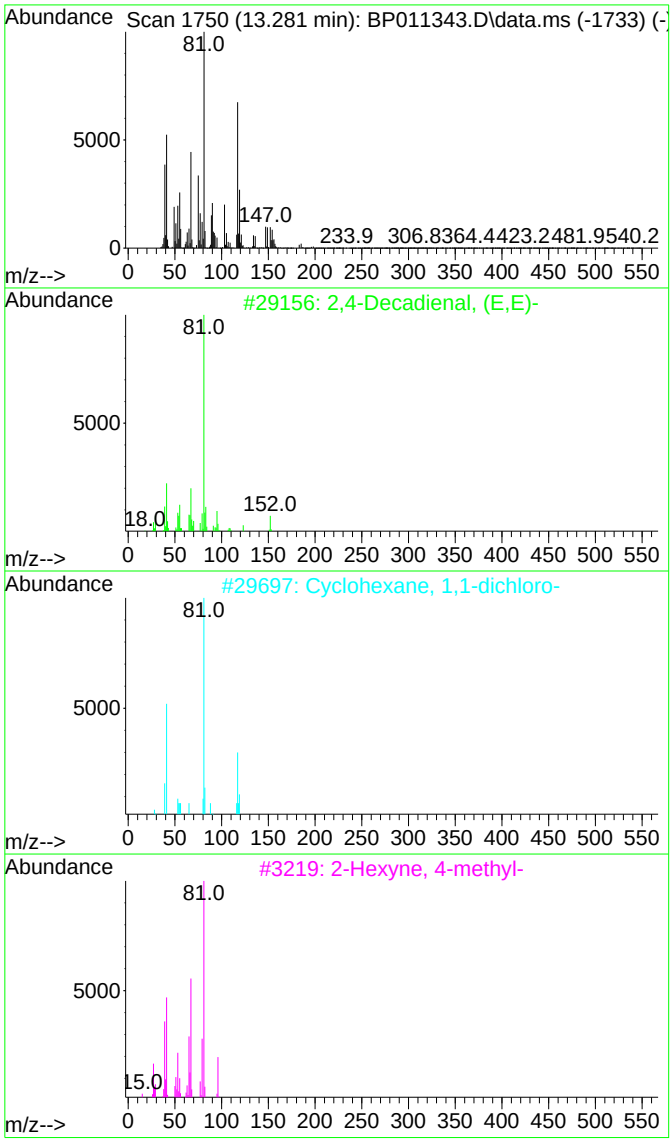
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

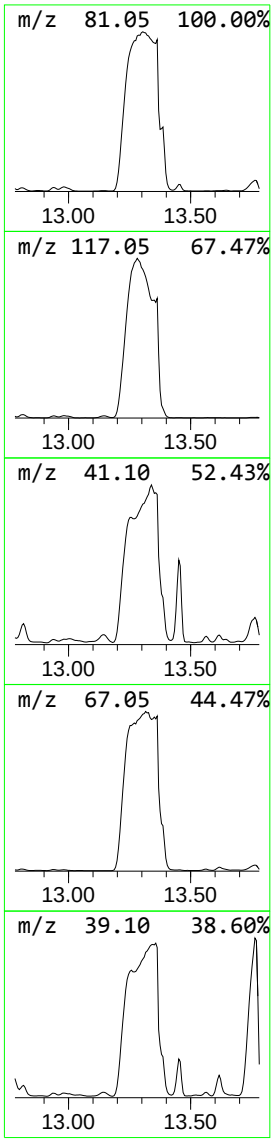
**********

Peak Number 45 unknown-09 Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.281 | 121.90 ng/ul | 133725000 | Acenaphthene-d10 | 14.298 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2,4-Decadienal, (E,E)- | 152 | C10H16O | 025152-84-5 | 27 |
| 2 | Cyclohexane, 1,1-dichloro- | 152 | C6H10Cl2 | 002108-92-1 | 16 |
| 3 | 2-Hexyne, 4-methyl- | 96 | C7H12 | 020198-49-6 | 12 |
| 4 | 1-Octyne | 110 | C8H14 | 000629-05-0 | 12 |
| 5 | 1-Ethyl-5-methylcyclopentene | 110 | C8H14 | 097797-57-4 | 12 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

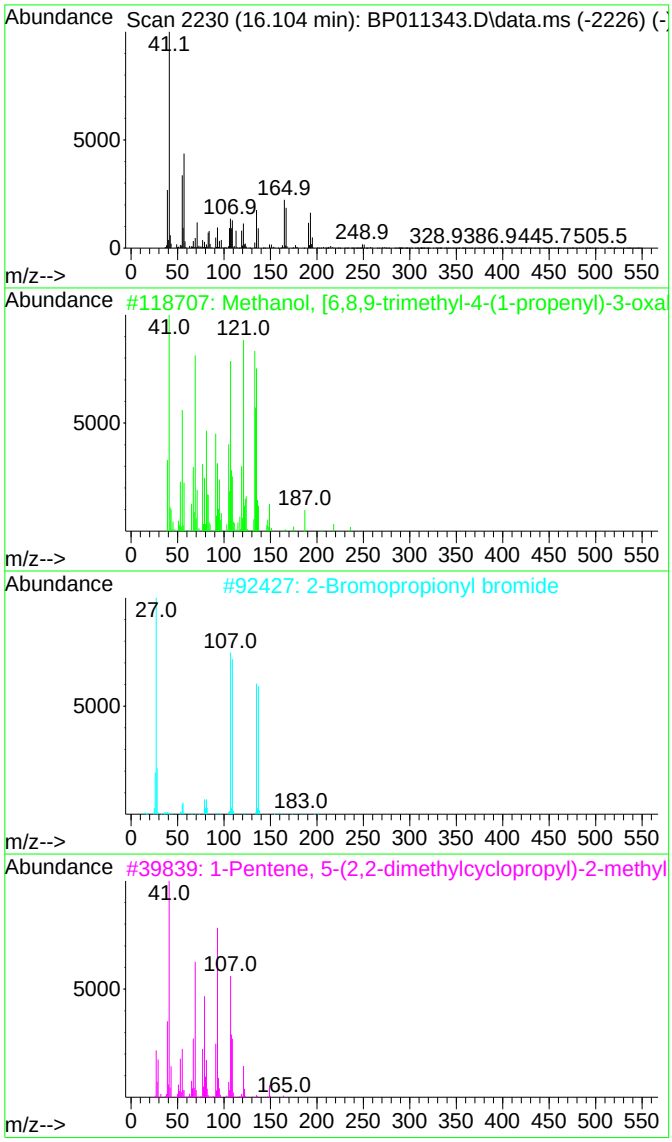
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

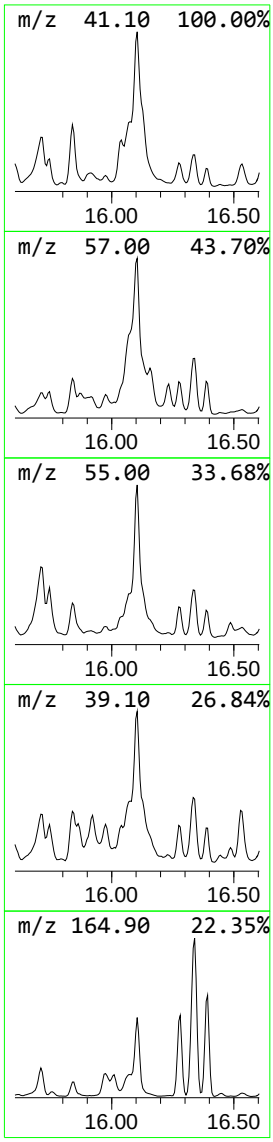
**********

Peak Number 57 unknown-10 Concentration Rank 25

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.104 | 35.22 ng/ul | 10360000 | Phenanthrene-d10 | 17.057 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Methanol, [6,8,9-trimethyl-4-(1-... | 236 | C15H24O2 | 1000277-60-9 | 10 |
| 2 | 2-Bromopropionyl bromide | 214 | C3H4Br2O | 000563-76-8 | 10 |
| 3 | 1-Pentene, 5-(2,2-dimethylcyclop... | 164 | C12H20 | 1000150-39-5 | 9 |
| 4 | [3-(2-Fluorophenyl)-1,2,4-oxadia... | 193 | C9H8FN3O | 1000444-23-9 | 9 |
| 5 | Phenol, 3-(2-aminoethyl)- | 137 | C8H11NO | 000588-05-6 | 9 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

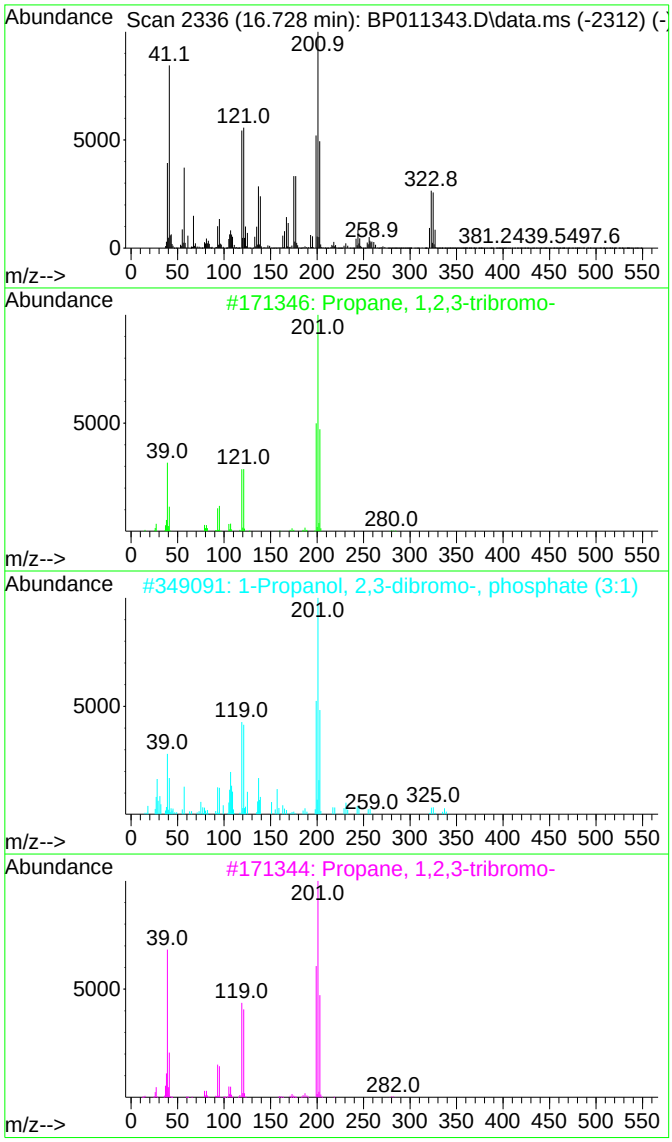
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

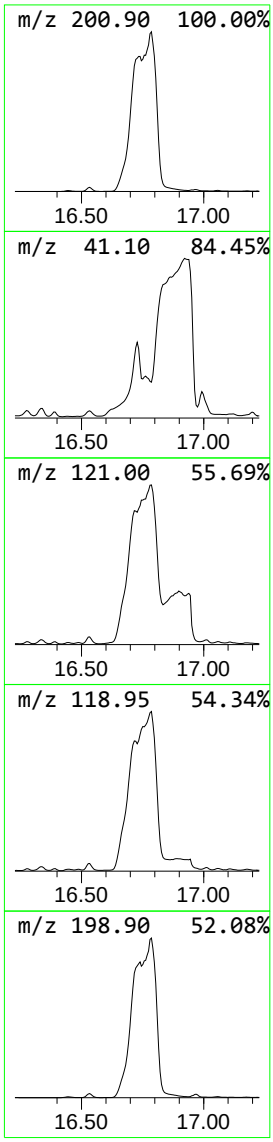
**********

Peak Number 59 unknown-11 Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.728 | 425.76 ng/ul | 125226000 | Phenanthrene-d10 | 17.057 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 43 |
| 2 | 1-Propanol, 2,3-dibromo-, phosph... | 692 | C9H15Br6O4P | 000126-72-7 | 40 |
| 3 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 32 |
| 4 | 2(3H)-Benzothiazolethione, 5-chl... | 201 | C7H4ClNS2 | 005331-91-9 | 30 |
| 5 | 2(3H)-Benzothiazolethione, 5-chl... | 201 | C7H4ClNS2 | 005331-91-9 | 16 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

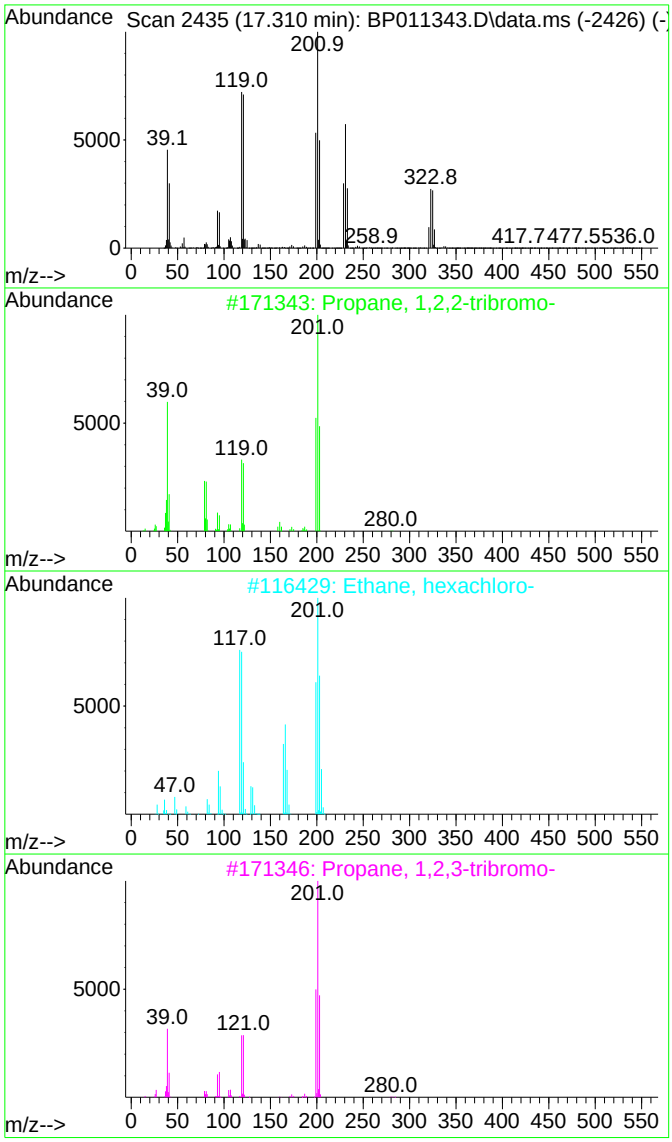
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

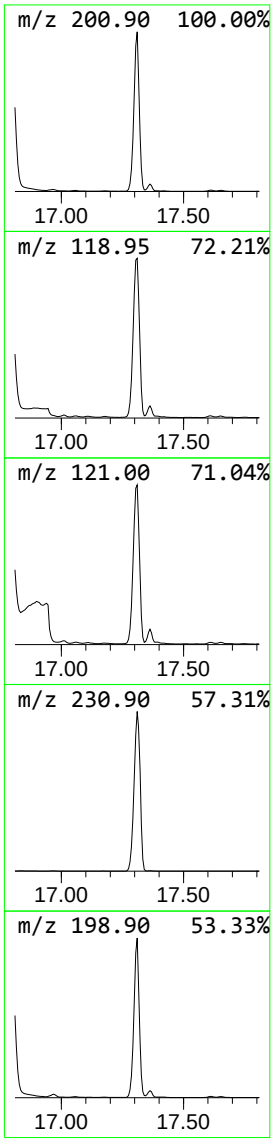
**********

Peak Number 60 unknown-12 Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.310 | 150.91 ng/ul | 44385300 | Phenanthrene-d10 | 17.057 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 1,2,2-tribromo- | 278 | C3H5Br3 | 014476-30-3 | 46 |
| 2 | Ethane, hexachloro- | 234 | C2Cl6 | 000067-72-1 | 38 |
| 3 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 25 |
| 4 | 2(3H)-Benzothiazolethione, 5-chl... | 201 | C7H4ClNS2 | 005331-91-9 | 10 |
| 5 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 10 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

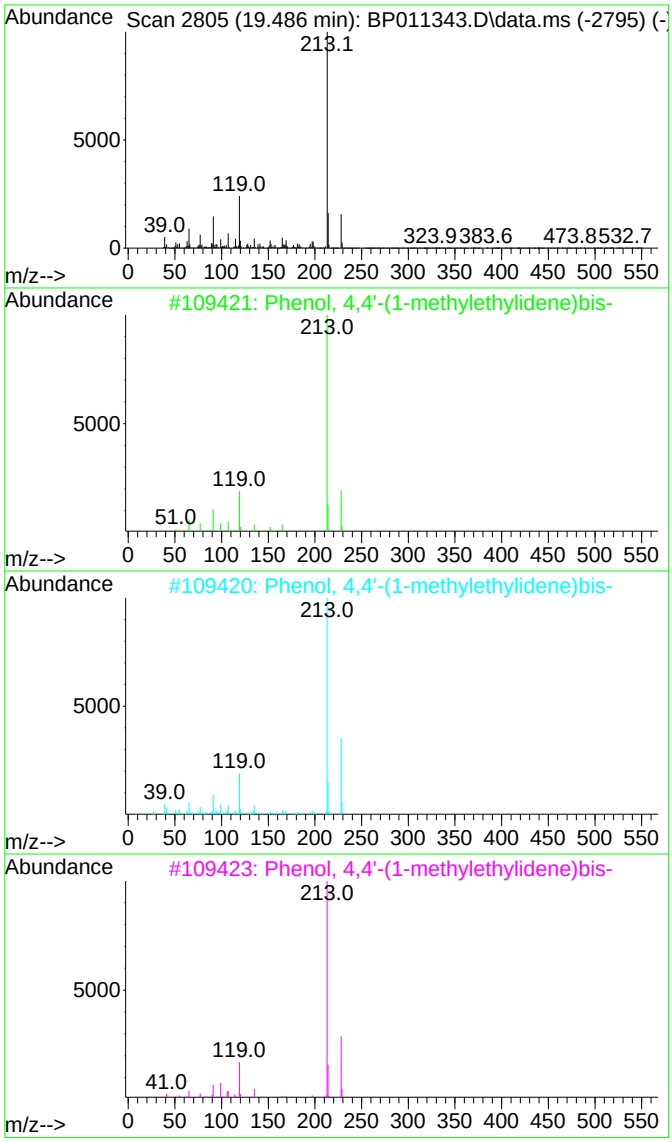
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

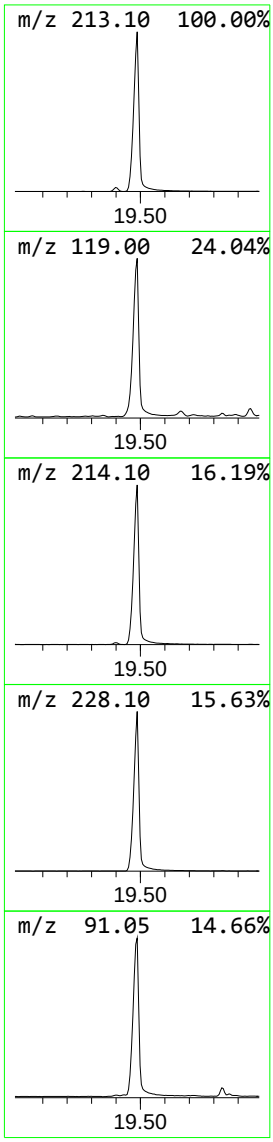
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 63 Phenol, 4,4'-(1-methylethyl... Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.486 | 127.59 ng/ul | 32936400 | Chrysene-d12 | 21.169 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 4,4'-(1-methylethylidene... | 228 | C15H16O2 | 000080-05-7 | 95 |
| 2 | Phenol, 4,4'-(1-methylethylidene... | 228 | C15H16O2 | 000080-05-7 | 94 |
| 3 | Phenol, 4,4'-(1-methylethylidene... | 228 | C15H16O2 | 000080-05-7 | 91 |
| 4 | Phenol, 4,4'-(1-methylethylidene... | 228 | C15H16O2 | 000080-05-7 | 90 |
| 5 | 3,4'-Isopropylidenediphenol | 228 | C15H16O2 | 046765-25-7 | 86 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

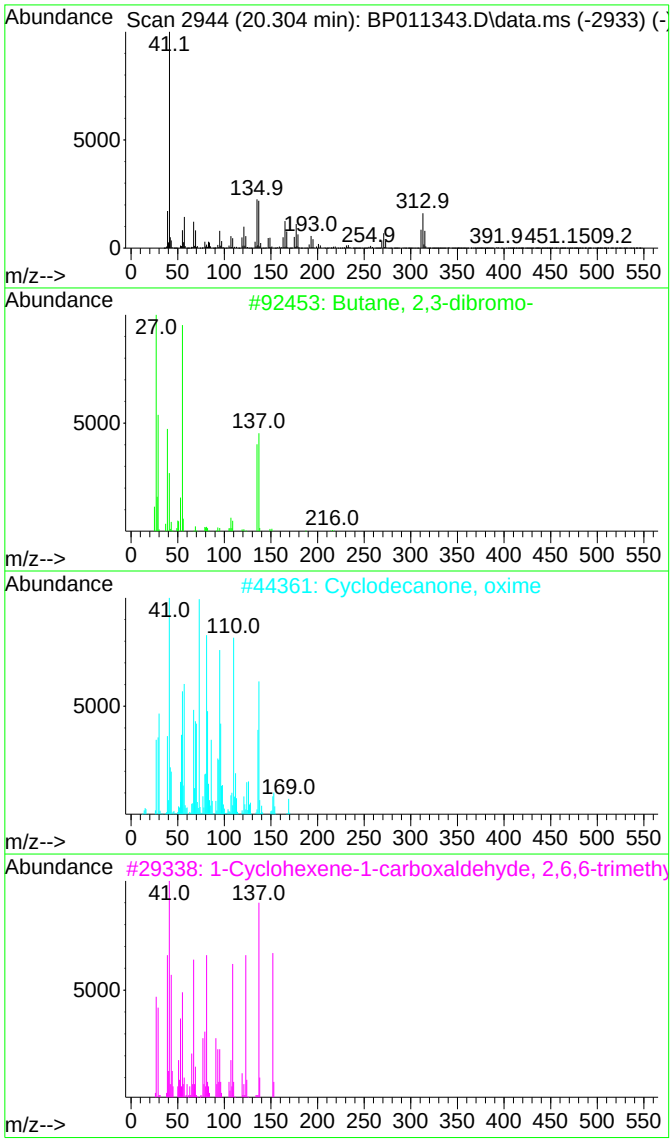
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

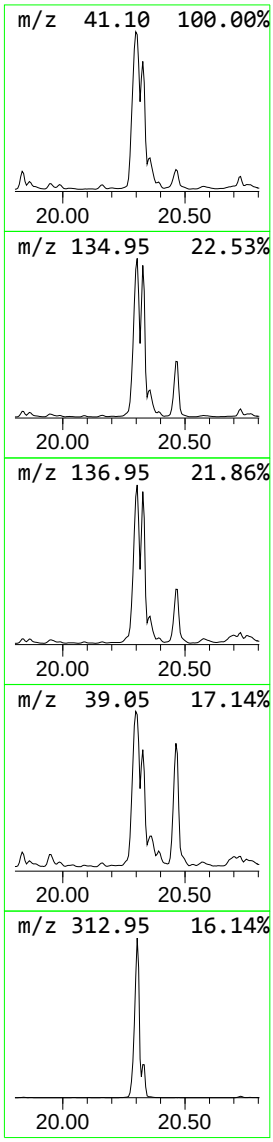
**********

Peak Number 64 unknown-13 Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.304 | 189.12 ng/ul | 48819100 | Chrysene-d12 | 21.169 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Butane, 2,3-dibromo- | 214 | C4H8Br2 | 005408-86-6 | 12 |
| 2 | Cyclodecanone, oxime | 169 | C10H19NO | 002972-01-2 | 10 |
| 3 | 1-Cyclohexene-1-carboxaldehyde, ... | 152 | C10H16O | 000432-25-7 | 9 |
| 4 | Triallyl phosphate | 218 | C9H15O4P | 001623-19-4 | 9 |
| 5 | Stannane, bromodibutyl(1-methyle... | 356 | C11H25BrSn | 055670-18-3 | 9 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

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

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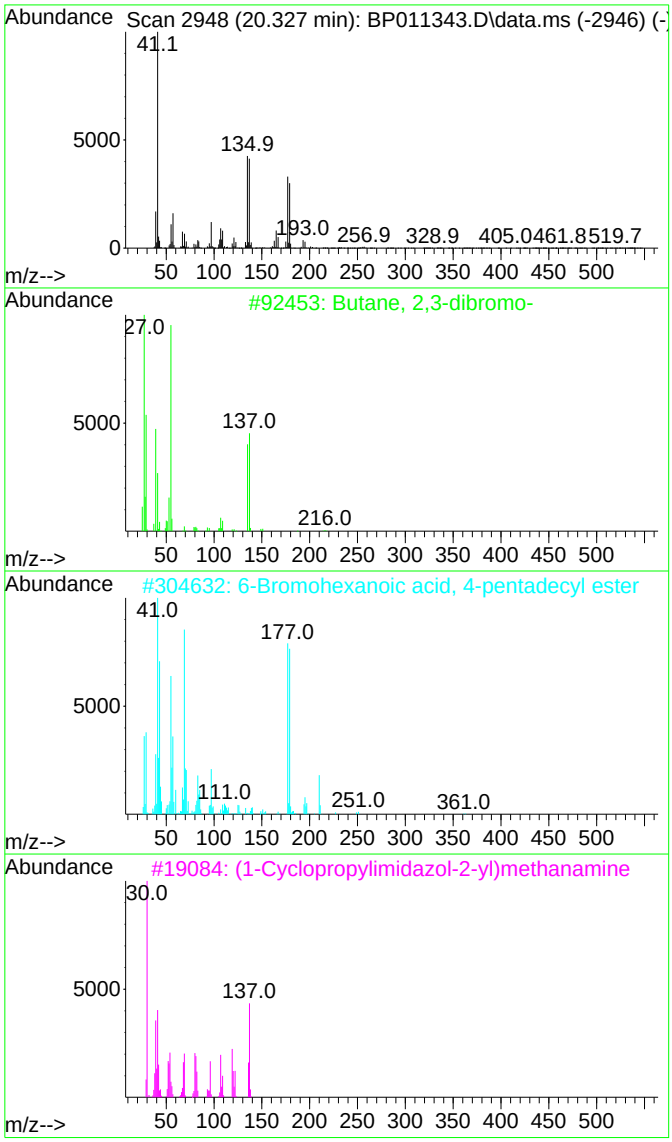
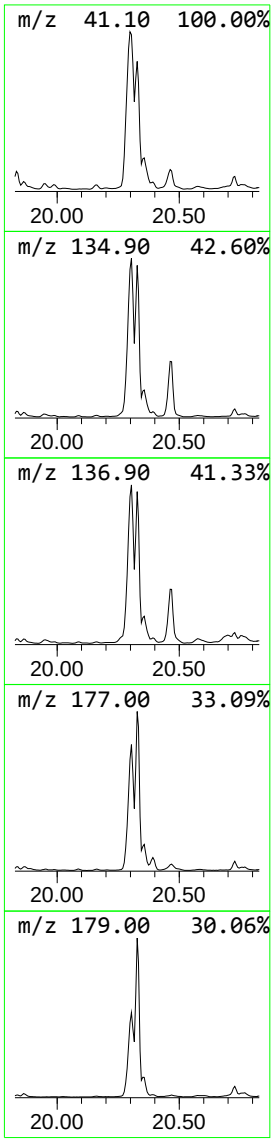
Peak Number 65 unknown-14 Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.327 | 86.39 ng/ul | 22301400 | Chrysene-d12 | 21.169 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Butane, 2,3-dibromo- | 214 | C4H8Br2 | 005408-86-6 | 25 |
| 2 | 6-Bromohexanoic acid, 4-pentadec... | 404 | C21H41BrO2 | 1000282-06-2 | 10 |
| 3 | (1-Cyclopropylimidazol-2-yl)meth... | 137 | C7H11N3 | 1000444-19-0 | 9 |
| 4 | Decane, 1-bromo- | 220 | C10H21Br | 000112-29-8 | 9 |
| 5 | 2-Bromopropionyl bromide | 214 | C3H4Br2O | 000563-76-8 | 9 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

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

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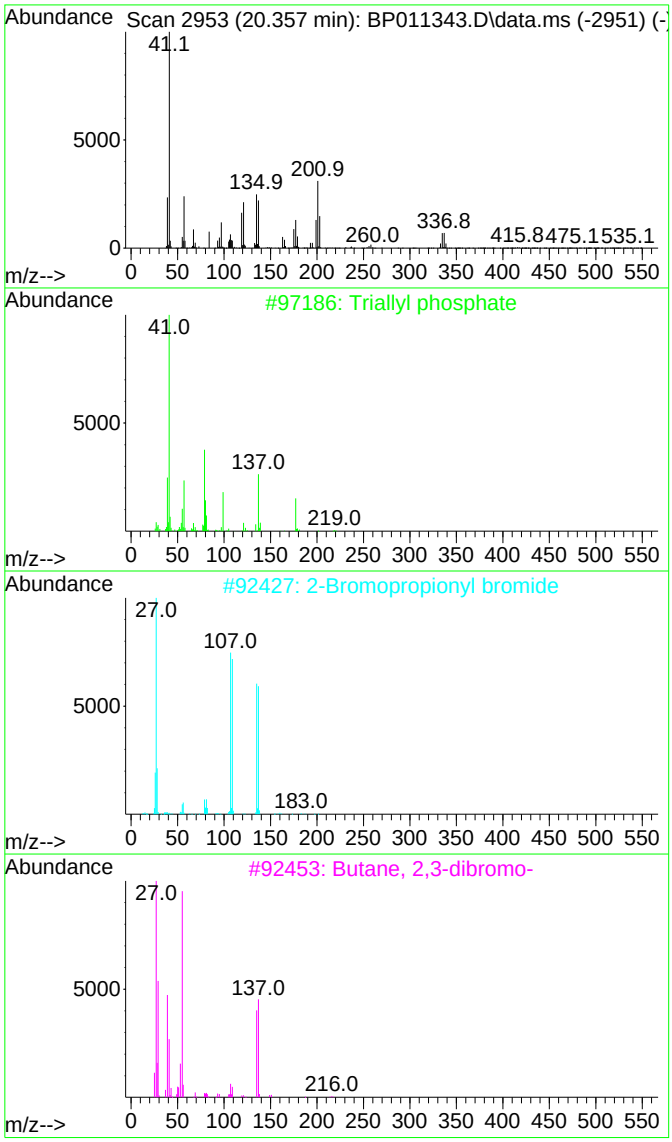
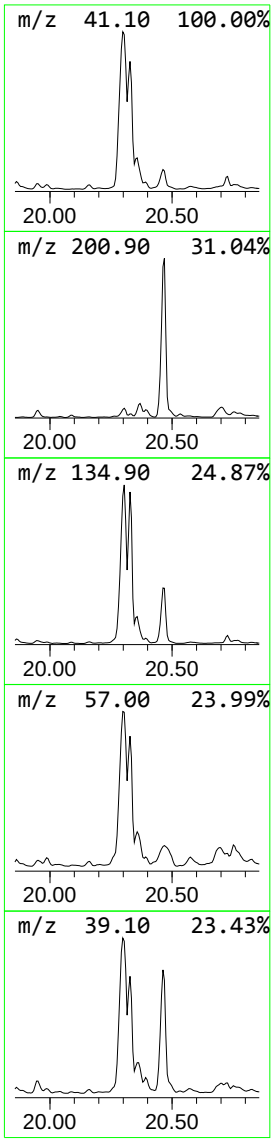
Peak Number 66 unknown-15 Concentration Rank 32

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.357 | 26.14 ng/ul | 6746900 | Chrysene-d12 | 21.169 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Triallyl phosphate | 218 | C9H15O4P | 001623-19-4 | 9 |
| 2 | 2-Bromopropionyl bromide | 214 | C3H4Br2O | 000563-76-8 | 9 |
| 3 | Butane, 2,3-dibromo- | 214 | C4H8Br2 | 005408-86-6 | 8 |
| 4 | Bronopol | 199 | C3H6BrNO4 | 000052-51-7 | 4 |
| 5 | Benzeneethanamine, .beta.-methyl- | 135 | C9H13N | 000582-22-9 | 3 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

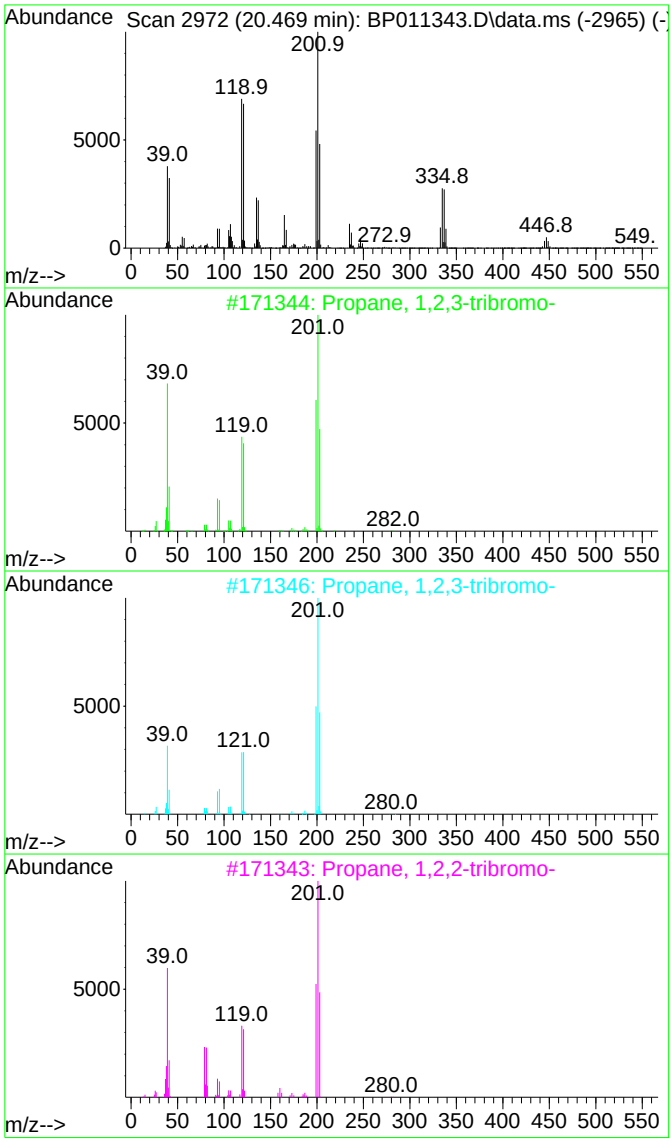
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

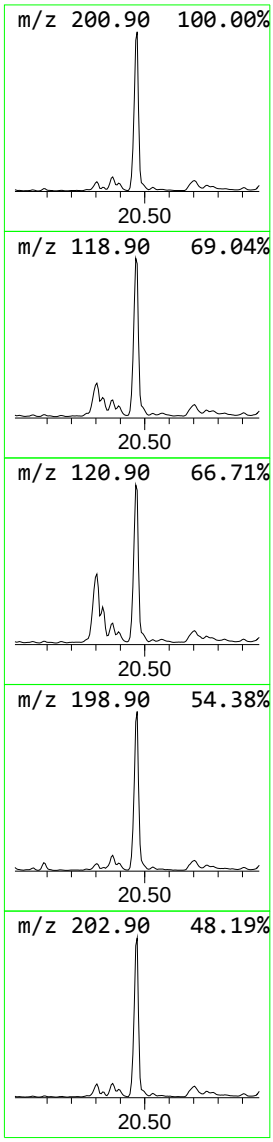
**********

Peak Number 67 unknown-16 Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.469 | 96.16 ng/ul | 24823900 | Chrysene-d12 | 21.169 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 47 |
| 2 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 47 |
| 3 | Propane, 1,2,2-tribromo- | 278 | C3H5Br3 | 014476-30-3 | 38 |
| 4 | 1-Propanol, 2,3-dibromo-, phosph... | 692 | C9H15Br6O4P | 000126-72-7 | 38 |
| 5 | Pyridine, 2,4-dichloro-3,5,6-tri... | 201 | C5Cl2F3N | 054889-43-9 | 16 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

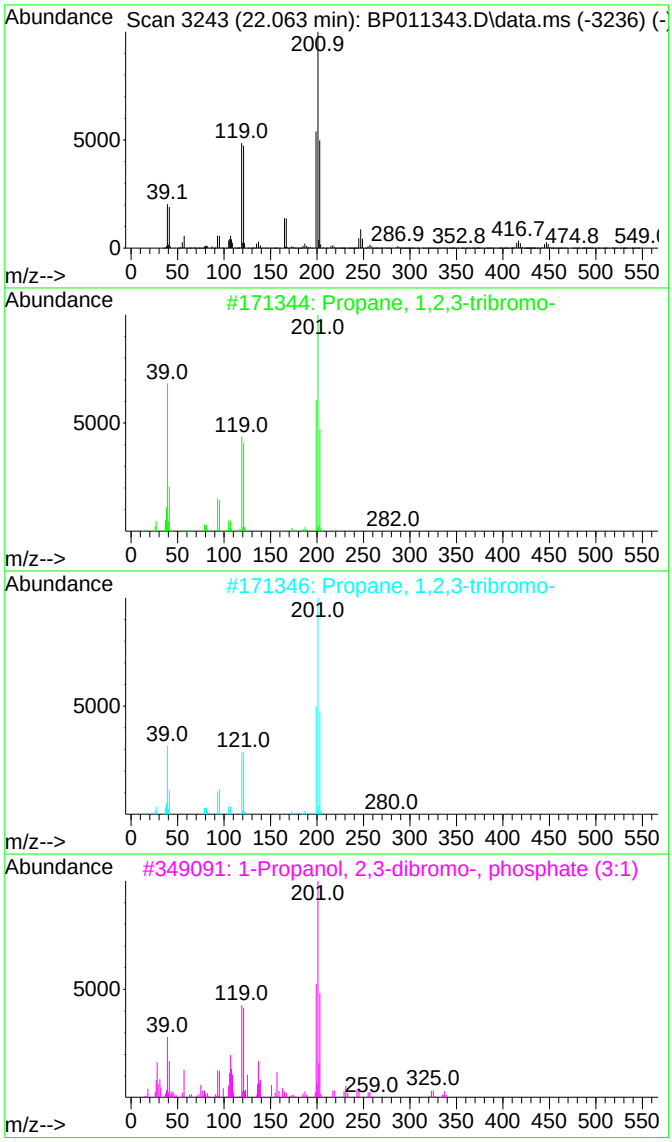
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

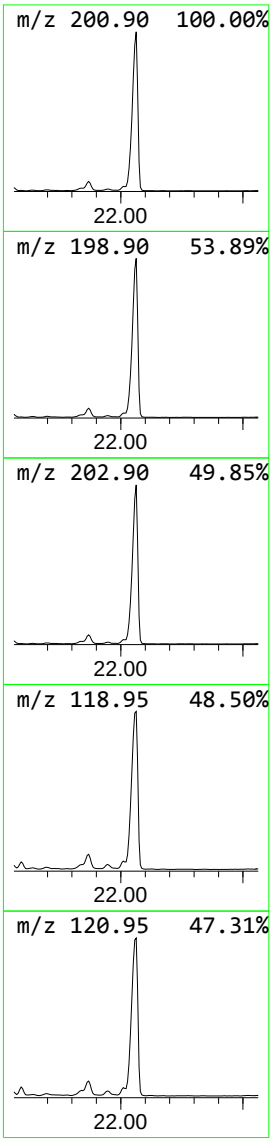
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 71 Propane, 1,2,3-tribromo- Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 22.063 | 156.49 ng/ul | 40395900 | Chrysene-d12 | 21.169 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 90 |
| 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 | Propane, 1,2,2-tribromo- | 278 | C3H5Br3 | 014476-30-3 | 50 |
| 5 | 1,4-Benzenediol, 2-chloro-, 4-(t... | 258 | C12H19ClO2Si | 1600530-39-9 | 25 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

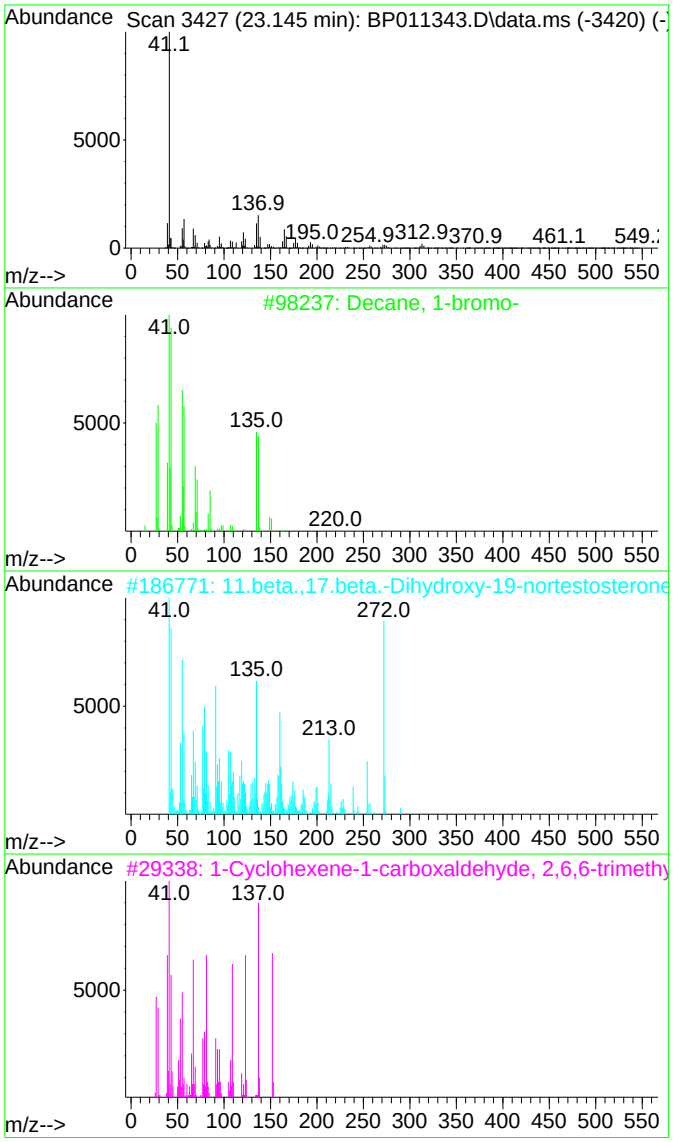
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

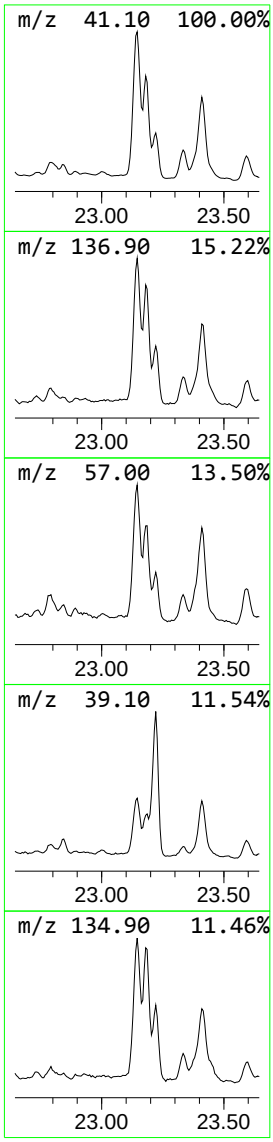
**********

Peak Number 72 unknown-17 Concentration Rank 21

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 23.145 | 37.01 ng/ul | 7473880 | Perylene-d12 | 23.474 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Decane, 1-bromo- | 220 | C10H21Br | 000112-29-8 | 12 |
| 2 | 11.beta.,17.beta.-Dihydroxy-19-n... | 290 | C18H26O3 | 1000213-78-5 | 9 |
| 3 | 1-Cyclohexene-1-carboxaldehyde, ... | 152 | C10H16O | 000432-25-7 | 9 |
| 4 | Ppropiolic acid, 3-(1-hydroxy-2-... | 224 | C13H20O3 | 1000158-61-9 | 9 |
| 5 | 2,3-Dibromopropyl alcohol | 216 | C3H6Br2O | 000096-13-9 | 9 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

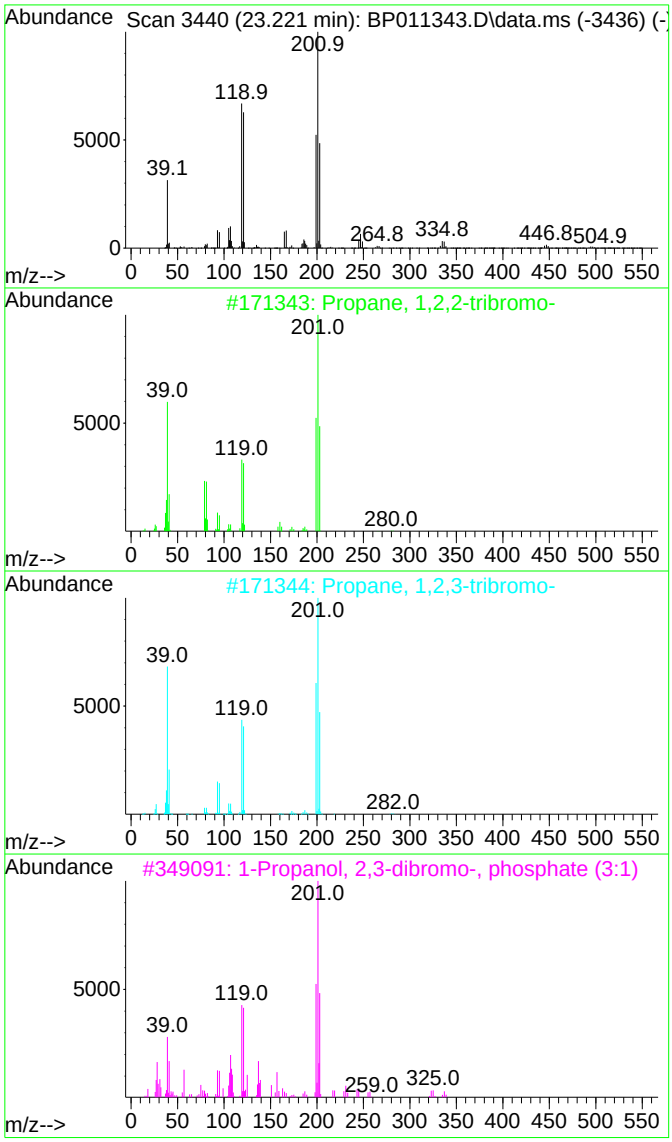
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

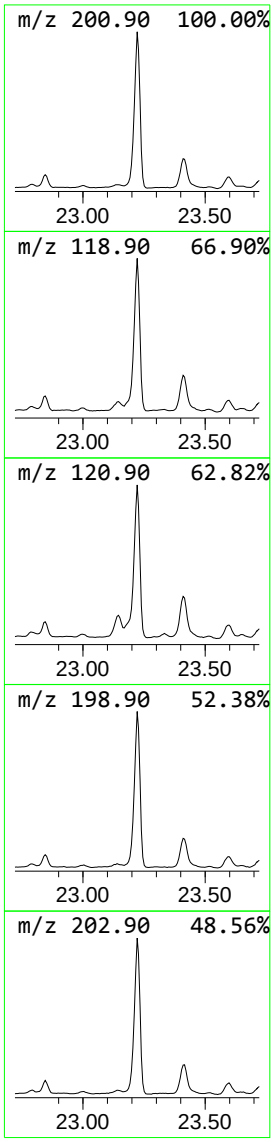
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 74 Propane, 1,2,2-tribromo- Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 23.221 | 42.16 ng/ul | 8513450 | Perylene-d12 | 23.474 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 1,2,2-tribromo- | 278 | C3H5Br3 | 014476-30-3 | 50 |
| 2 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 50 |
| 3 | 1-Propanol, 2,3-dibromo-, phosph... | 692 | C9H15Br6O4P | 000126-72-7 | 43 |
| 4 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 43 |
| 5 | 2(3H)-Benzothiazolethione, 5-chl... | 201 | C7H4ClNS2 | 005331-91-9 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

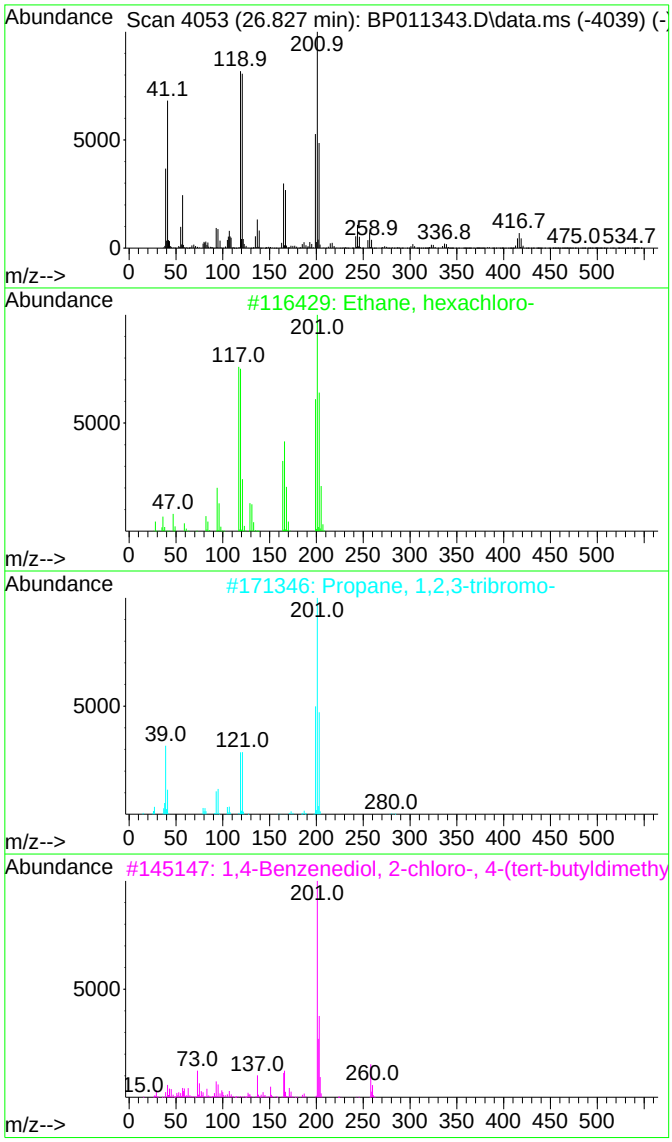
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

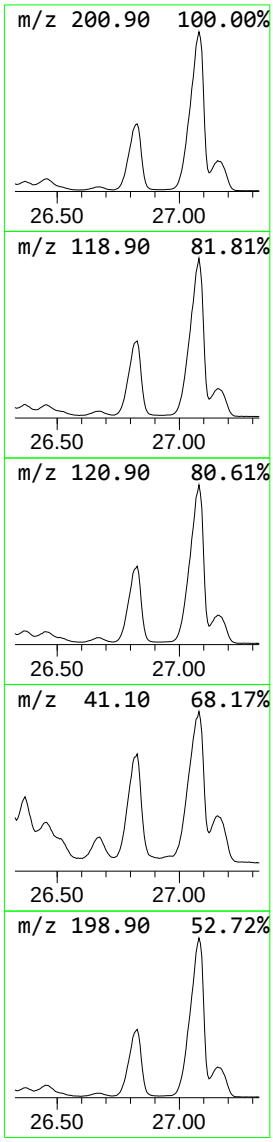
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 79 unknown-18 Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 26.827 | 73.82 ng/ul | 14906400 | Perylene-d12 | 23.474 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Ethane, hexachloro- | 234 | C2Cl6 | 000067-72-1 | 40 |
| 2 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 38 |
| 3 | 1,4-Benzenediol, 2-chloro-, 4-(t... | 258 | C12H19ClO2Si | 1600530-39-9 | 14 |
| 4 | Pyridine, 2,4-dichloro-3,5,6-tri... | 201 | C5Cl2F3N | 054889-43-9 | 12 |
| 5 | 2'-Ethyl-3-[(3-phenylpropionyl)h... | 351 | C21H25N3O2 | 1000225-96-3 | 10 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

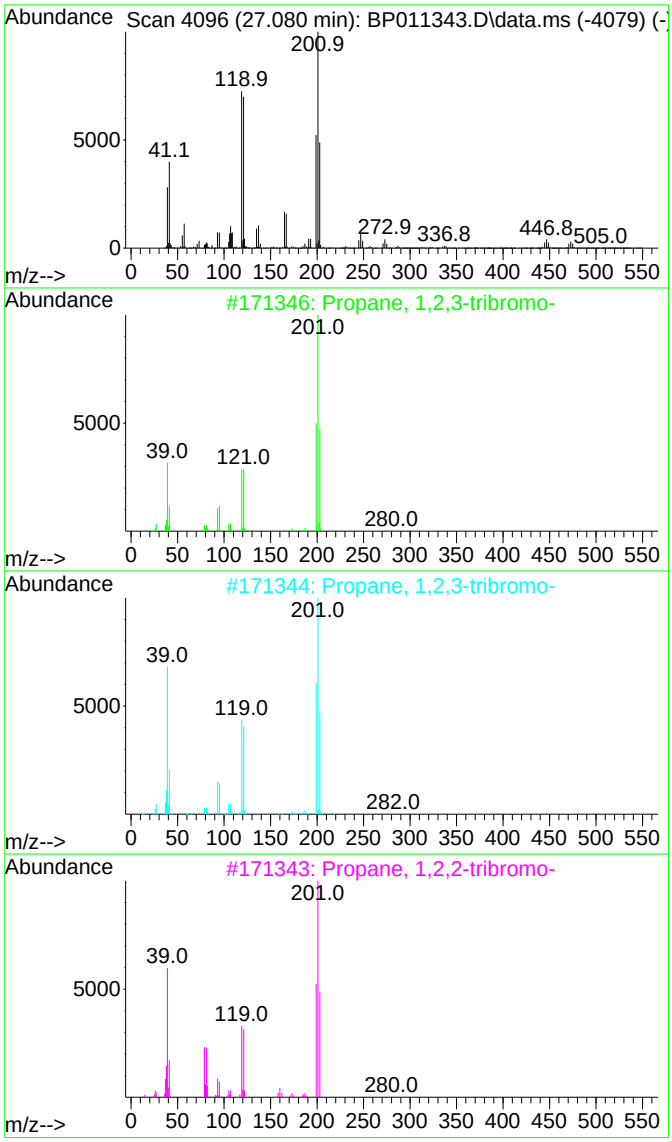
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION

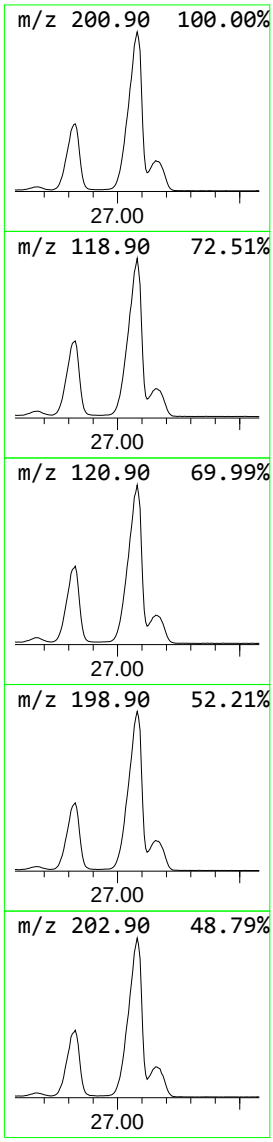
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 80 unknown-19 Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 27.080 | 148.93 ng/ul | 30071800 | Perylene-d12 | 23.474 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 78 |
| 2 | Propane, 1,2,3-tribromo- | 278 | C3H5Br3 | 000096-11-7 | 43 |
| 3 | Propane, 1,2,2-tribromo- | 278 | C3H5Br3 | 014476-30-3 | 39 |
| 4 | 2(3H)-Benzothiazolethione, 5-chl... | 201 | C7H4ClNS2 | 005331-91-9 | 37 |
| 5 | 2(3H)-Benzothiazolethione, 5-chl... | 201 | C7H4ClNS2 | 005331-91-9 | 25 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011343.D
Acq On : 09 Aug 2022 23:04
Operator : CG/JU
Sample : N3956-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EXF01

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.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 |
| Tetrahydrofurfu... | 3.828 | 64.0 | ng/ul | 35965800 | 1 | 7.593 | 11238200 | 20.0 |
| Propane, 1,3-di... | 4.052 | 78.7 | ng/ul | 44207500 | 1 | 7.593 | 11238200 | 20.0 |
| Propane, 1-brom... | 5.222 | 35.9 | ng/ul | 20155300 | 1 | 7.593 | 11238200 | 20.0 |
| Benzene, bromo- | 6.281 | 60.2 | ng/ul | 33827600 | 1 | 7.593 | 11238200 | 20.0 |
| unknown-01 | 8.022 | 35.5 | ng/ul | 19943200 | 1 | 7.593 | 11238200 | 20.0 |
| (DEL) Alkane: C... | 8.216 | 26.2 | ng/ul | 14708500 | 1 | 7.593 | 11238200 | 20.0 |
| (DEL) Alkane: S... | 8.581 | 6.1 | ng/ul | 3438110 | 1 | 7.593 | 11238200 | 20.0 |
| unknown-02 | 8.704 | 27.5 | ng/ul | 15441100 | 1 | 7.593 | 11238200 | 20.0 |
| 2,3-Dibromoprop... | 8.969 | 74.0 | ng/ul | 41594000 | 1 | 7.593 | 11238200 | 20.0 |
| unknown-03 | 9.263 | 32.6 | ng/ul | 37951600 | 2 | 10.428 | 23305600 | 20.0 |
| Benzene, 1-brom... | 9.416 | 28.5 | ng/ul | 33259100 | 2 | 10.428 | 23305600 | 20.0 |
| unknown-04 | 9.828 | 178.8 | ng/ul | 208295000 | 2 | 10.428 | 23305600 | 20.0 |
| unknown-05 | 10.698 | 25.2 | ng/ul | 29335500 | 2 | 10.428 | 23305600 | 20.0 |
| Hexane, 1-bromo... | 10.822 | 36.9 | ng/ul | 43008800 | 2 | 10.428 | 23305600 | 20.0 |
| unknown-06 | 11.046 | 96.6 | ng/ul | 112552000 | 2 | 10.428 | 23305600 | 20.0 |
| unknown-07 | 11.140 | 99.3 | ng/ul | 115702000 | 2 | 10.428 | 23305600 | 20.0 |
| Hexane, 1,6-dib... | 12.093 | 26.3 | ng/ul | 30584400 | 2 | 10.428 | 23305600 | 20.0 |
| unknown-08 | 12.216 | 104.1 | ng/ul | 121274000 | 2 | 10.428 | 23305600 | 20.0 |
| unknown-09 | 13.281 | 121.9 | ng/ul | 133725000 | 3 | 14.298 | 21940500 | 20.0 |
| unknown-10 | 16.104 | 35.2 | ng/ul | 10360000 | 4 | 17.057 | 5882510 | 20.0 |
| unknown-11 | 16.728 | 425.8 | ng/ul | 125226000 | 4 | 17.057 | 5882510 | 20.0 |
| unknown-12 | 17.310 | 150.9 | ng/ul | 44385300 | 4 | 17.057 | 5882510 | 20.0 |
| Phenol, 4,4'-(1... | 19.486 | 127.6 | ng/ul | 32936400 | 5 | 21.169 | 5162890 | 20.0 |
| unknown-13 | 20.304 | 189.1 | ng/ul | 48819100 | 5 | 21.169 | 5162890 | 20.0 |
| unknown-14 | 20.327 | 86.4 | ng/ul | 22301400 | 5 | 21.169 | 5162890 | 20.0 |
| unknown-15 | 20.357 | 26.1 | ng/ul | 6746900 | 5 | 21.169 | 5162890 | 20.0 |
| unknown-16 | 20.469 | 96.2 | ng/ul | 24823900 | 5 | 21.169 | 5162890 | 20.0 |
| Propane, 1,2,3-... | 22.063 | 156.5 | ng/ul | 40395900 | 5 | 21.169 | 5162890 | 20.0 |
| unknown-17 | 23.145 | 37.0 | ng/ul | 7473880 | 6 | 23.474 | 4038490 | 20.0 |
| Propane, 1,2,2-... | 23.221 | 42.2 | ng/ul | 8513450 | 6 | 23.474 | 4038490 | 20.0 |
| unknown-18 | 26.827 | 73.8 | ng/ul | 14906400 | 6 | 23.474 | 4038490 | 20.0 |
| unknown-19 | 27.080 | 148.9 | ng/ul | 30071800 | 6 | 23.474 | 4038490 | 20.0 |