

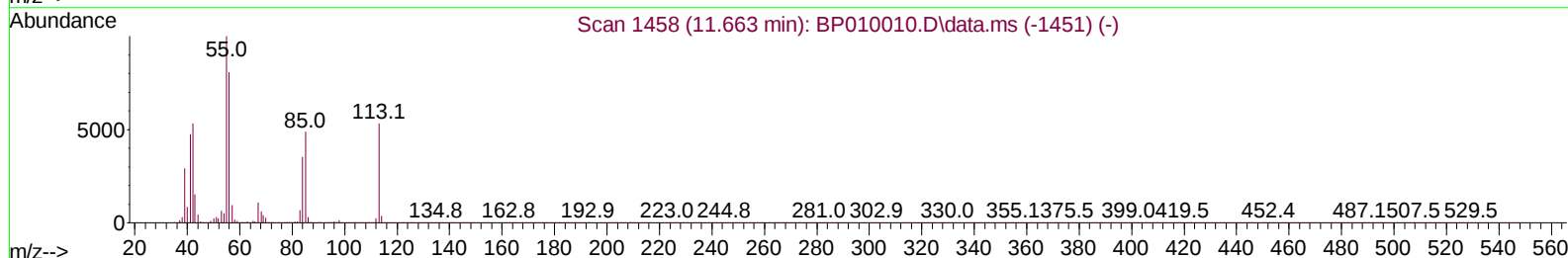
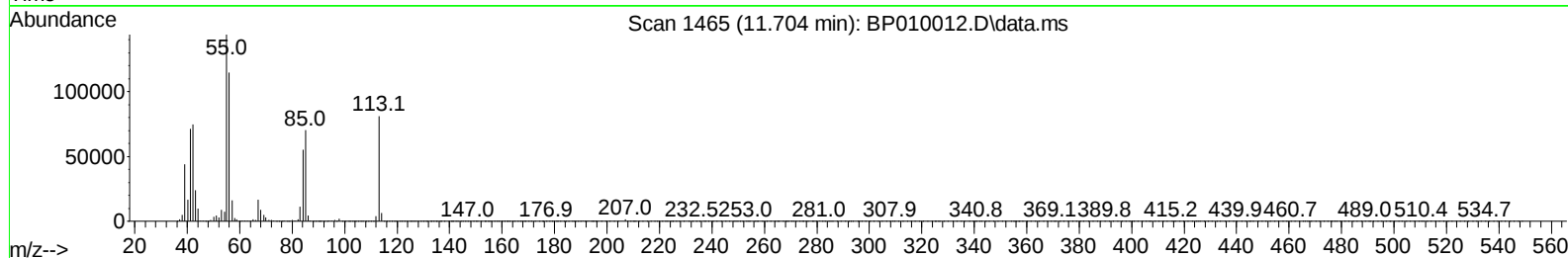
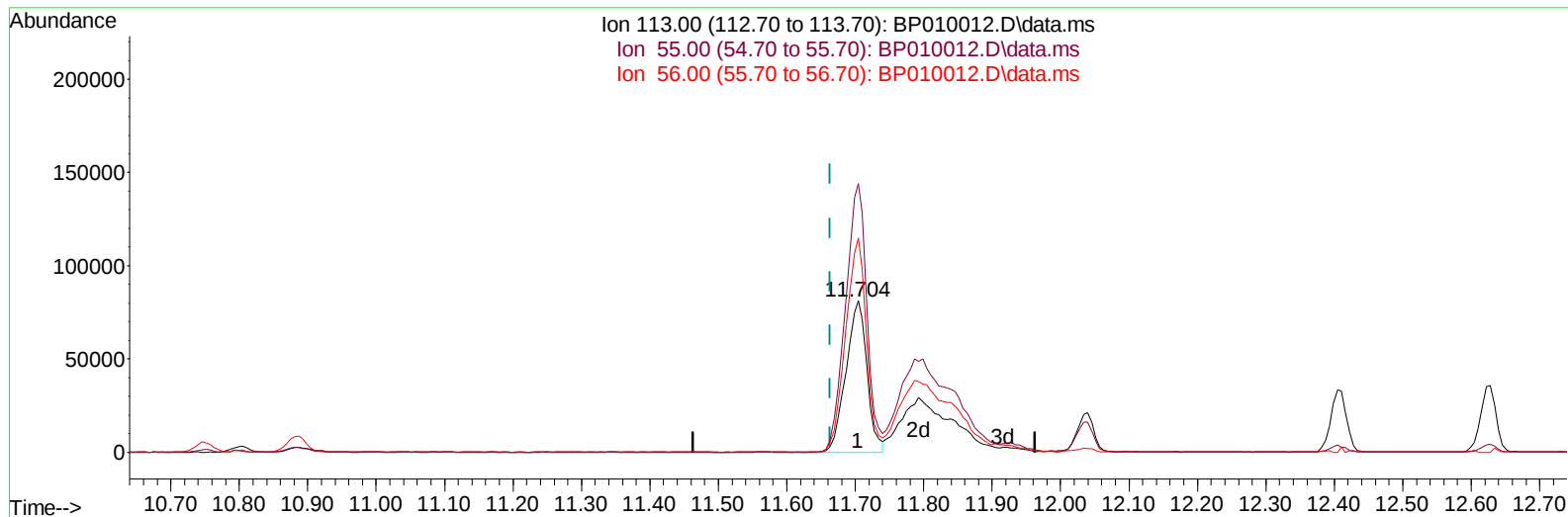
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042122\  
 Data File : BP010012.D  
 Acq On : 21 Apr 2022 13:19  
 Operator : CG/JU  
 Sample : SST08027  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SST080627

Manual IntegrationsAPPROVED

Quant Time: Apr 21 14:25:12 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 21 13:57:49 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/22/2022  
 Supervised By :mohammad ahmed 04/26/2022



TIC: BP010012.D\data.ms

**(34) Caprolactam**

**11.704min (+ 0.041) 47.38 ng/ul**

**response 172287**

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 113.00 | 100.00 | 100.00  |
| 55.00  | 72.70  | 177.33# |
| 56.00  | 85.80  | 141.44# |
| 0.00   | 0.00   | 0.00    |

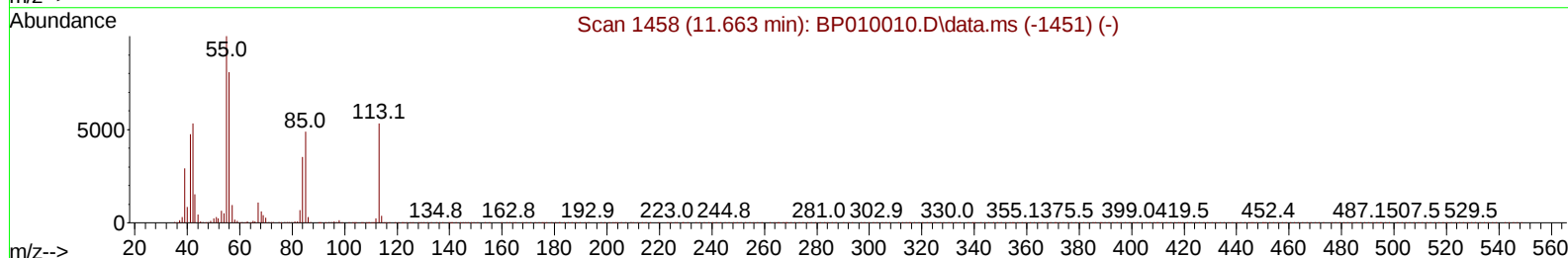
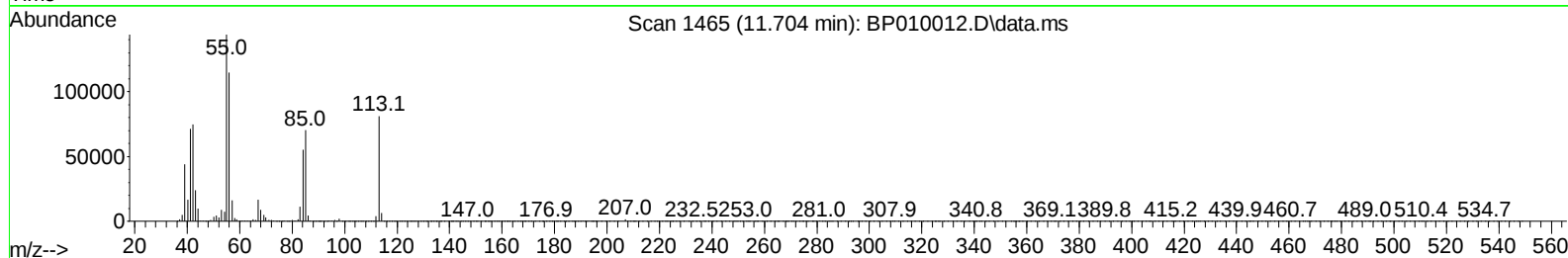
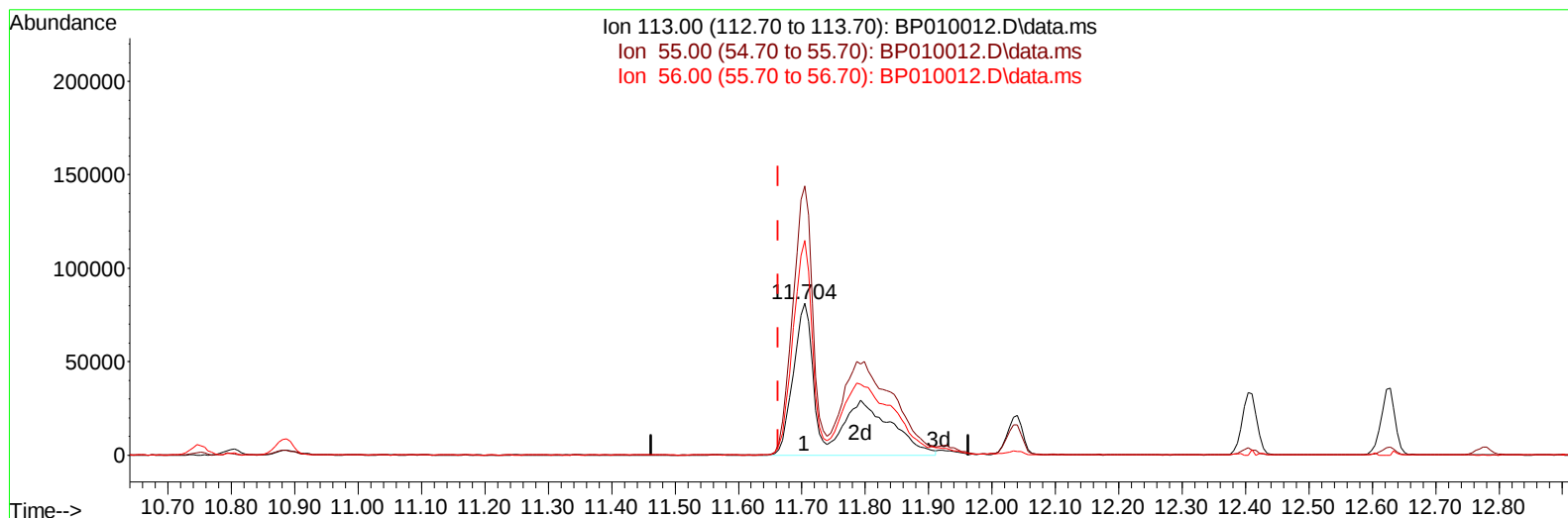
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042122\  
 Data File : BP010012.D  
 Acq On : 21 Apr 2022 13:19  
 Operator : CG/JU  
 Sample : SST08027  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SST080627

Manual IntegrationsAPPROVED

Quant Time: Apr 21 14:25:12 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 21 13:57:49 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/22/2022  
 Supervised By :mohammad ahmed 04/26/2022



TIC: BP010012.D\data.ms

**(34) Caprolactam**

11.704min (+ 0.041) 88.60 ng/ul m

response 322178

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 113.00 | 100.00 | 100.00  |
| 55.00  | 72.70  | 177.33# |
| 56.00  | 85.80  | 141.44# |
| 0.00   | 0.00   | 0.00    |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042122\  
 Data File : BP010012.D  
 Acq On : 21 Apr 2022 13:19  
 Operator : CG/JU  
 Sample : SST008027  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SST0080627

Manual IntegrationsAPPROVED

Quant Time: Apr 21 14:25:12 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 21 13:57:49 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/22/2022  
 Supervised By :mohammad ahmed 04/26/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.940  | 152  | 154486   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.752 | 136  | 685572   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.575 | 164  | 475900   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.328 | 188  | 1041784  | 20.000 | ng/ul  | # 0.00   |
| 79) Chrysene-d12              | 21.410 | 240  | 1054480  | 20.000 | ng/ul  | # 0.00   |
| 88) Perylene-d12              | 23.868 | 264  | 1004342  | 20.000 | ng/ul  | 0.01     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.370  | 96   | 110180   | 31.830 | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.781  | 84   | 815666   | 85.465 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.111  | 99   | 1172799  | 86.471 | ng/ul  | 0.01     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.269  | 67   | 675105   | 83.610 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.475  | 132  | 885563   | 86.559 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.658  | 113  | 955194   | 84.365 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.105  | 128  | 464332   | 93.921 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.828  | 143  | 535149   | 98.732 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.369 | 165  | 971376   | 85.643 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.887 | 131  | 1333637  | 82.556 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.987 | 166  | 2948900  | 78.884 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.269 | 160  | 3638956  | 82.166 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.781 | 143  | 572702   | 84.475 | ng/ul  | 0.02     |
| 60) Fluorene-d10              | 15.575 | 176  | 2533727  | 80.570 | ng/ul  | 0.01     |
| 65) 4,6-Dinitro-2-methylph... | 15.692 | 200  | 578050   | 93.664 | ng/ul  | 0.01     |
| 73) Anthracene-d10            | 17.434 | 188  | 4090637  | 82.010 | ng/ul  | 0.01     |
| 81) Pyrene-d10                | 19.663 | 212  | 4879640  | 84.730 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.716 | 264  | 4364490  | 83.447 | ng/ul  | 0.02     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.405  | 88   | 113406   | 32.138 | ng/ul# | 22       |
| 5) Pyridine                   | 3.805  | 79   | 843081   | 84.666 | ng/ul# | 46       |
| 6) Benzaldehyde               | 7.075  | 77   | 420177   | 50.515 | ng/ul  | 98       |
| 8) Phenol                     | 7.134  | 94   | 1177662  | 86.746 | ng/ul# | 93       |
| 10) Bis(2-Chloroethyl)ether   | 7.363  | 93   | 878095   | 83.034 | ng/ul# | 77       |
| 12) 2-Chlorophenol            | 7.511  | 128  | 892660   | 86.546 | ng/ul  | 92       |
| 13) 2-Methylphenol            | 8.387  | 108  | 890215   | 84.090 | ng/ul  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.475  | 45   | 1271827  | 81.362 | ng/ul# | 85       |
| 16) Acetophenone              | 8.769  | 105  | 1452656  | 81.434 | ng/ul# | 81       |
| 17) N-Nitroso-di-n-propyla... | 8.769  | 70   | 763656   | 82.572 | ng/ul# | 75       |
| 18) 4-Methylphenol            | 8.728  | 108  | 971245   | 84.719 | ng/ul  | 94       |
| 19) Hexachloroethane          | 9.028  | 117  | 358066   | 82.224 | ng/ul# | 81       |
| 22) Nitrobenzene              | 9.152  | 77   | 1103124  | 86.930 | ng/ul# | 86       |
| 23) Isophorone                | 9.687  | 82   | 2174506  | 83.158 | ng/ul# | 97       |
| 25) 2-Nitrophenol             | 9.863  | 139  | 544442   | 94.179 | ng/ul# | 84       |
| 26) 2,4-Dimethylphenol        | 9.922  | 107  | 1126946  | 82.147 | ng/ul# | 78       |
| 27) Bis(2-Chloroethoxy)met... | 10.157 | 93   | 1230808  | 80.171 | ng/ul# | 98       |
| 29) 2,4-Dichlorophenol        | 10.399 | 162  | 923752   | 84.009 | ng/ul# | 83       |
| 30) Naphthalene               | 10.799 | 128  | 2879507  | 81.668 | ng/ul  | 97       |
| 32) 4-Chloroaniline           | 10.910 | 127  | 1314410  | 82.221 | ng/ul  | 97       |
| 33) Hexachlorobutadiene       | 11.093 | 225  | 631589   | 79.752 | ng/ul  | 95       |
| 34) Caprolactam               | 11.704 | 113  | 322178m  | 88.597 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 12.040 | 107  | 1098279  | 82.430 | ng/ul# | 69       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042122\  
 Data File : BP010012.D  
 Acq On : 21 Apr 2022 13:19  
 Operator : CG/JU  
 Sample : SST008027  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SST0080627

## Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/22/2022  
 Supervised By :mohammad ahmed 04/26/2022

Quant Time: Apr 21 14:25:12 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 21 13:57:49 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 36) 2-Methylnaphthalene       | 12.404 | 142  | 2061863  | 80.044 | ng/ul  | 92       |
| 37) 1-Methylnaphthalene       | 12.628 | 142  | 2086880  | 80.237 | ng/ul# | 95       |
| 39) 1,2,4,5-Tetrachloroben... | 12.775 | 216  | 1205410  | 82.788 | ng/ul  | 94       |
| 40) Hexachlorocyclopentadiene | 12.757 | 237  | 718162   | 71.077 | ng/ul  | 95       |
| 41) 2,4,6-Trichlorophenol     | 13.010 | 196  | 805163   | 86.591 | ng/ul# | 83       |
| 42) 2,4,5-Trichlorophenol     | 13.087 | 196  | 881621   | 87.735 | ng/ul# | 88       |
| 43) 1,1'-Biphenyl             | 13.416 | 154  | 2805695  | 80.801 | ng/ul  | 93       |
| 44) 2-Chloronaphthalene       | 13.457 | 162  | 2187187  | 81.944 | ng/ul  | 93       |
| 45) 2-Nitroaniline            | 13.657 | 65   | 772246   | 98.726 | ng/ul# | 82       |
| 47) Dimethylphthalate         | 14.034 | 163  | 2835498  | 78.809 | ng/ul# | 93       |
| 48) 2,6-Dinitrotoluene        | 14.151 | 165  | 644666   | 97.410 | ng/ul  | 93       |
| 50) Acenaphthylene            | 14.298 | 152  | 3556388  | 81.370 | ng/ul  | 96       |
| 51) 3-Nitroaniline            | 14.481 | 138  | 585152   | 85.050 | ng/ul# | 80       |
| 52) Acenaphthene              | 14.640 | 153  | 2292837  | 80.123 | ng/ul  | 96       |
| 53) 2,4-Dinitrophenol         | 14.687 | 184  | 380108   | 95.019 | ng/ul# | 80       |
| 55) 4-Nitrophenol             | 14.798 | 109  | 487442   | 78.106 | ng/ul# | 47       |
| 56) Dibenzofuran              | 14.975 | 168  | 3337363  | 79.241 | ng/ul  | 97       |
| 57) 2,4-Dinitrotoluene        | 14.940 | 165  | 912976   | 93.948 | ng/ul# | 84       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.204 | 232  | 772864   | 84.355 | ng/ul# | 87       |
| 59) Diethylphthalate          | 15.404 | 149  | 2935565  | 78.837 | ng/ul  | 94       |
| 61) Fluorene                  | 15.628 | 166  | 2730127  | 78.700 | ng/ul  | 90       |
| 62) 4-Chlorophenyl-phenyle... | 15.622 | 204  | 1430335  | 77.544 | ng/ul  | 99       |
| 63) 4-Nitroaniline            | 15.651 | 138  | 560642   | 81.357 | ng/ul# | 79       |
| 66) 4,6-Dinitro-2-methylph... | 15.710 | 198  | 546741   | 89.125 | ng/ul# | 93       |
| 67) N-Nitrosodiphenylamine    | 15.834 | 169  | 2432688  | 81.574 | ng/ul  | 95       |
| 68) 4-Bromophenyl-phenylether | 16.516 | 248  | 921643   | 81.728 | ng/ul  | 92       |
| 69) Hexachlorobenzene         | 16.639 | 284  | 1056026  | 81.439 | ng/ul# | 92       |
| 70) Atrazine                  | 16.792 | 200  | 984836   | 80.007 | ng/ul# | 88       |
| 71) Pentachlorophenol         | 16.981 | 266  | 706600   | 80.505 | ng/ul# | 84       |
| 72) Phenanthrene              | 17.375 | 178  | 4508520  | 81.076 | ng/ul  | 99       |
| 74) Anthracene                | 17.469 | 178  | 4550284  | 80.708 | ng/ul  | 98       |
| 75) 1,2,3,4-Tetrachloroben... | 13.381 | 216  | 1262634  | 83.830 | ng/uL# | 85       |
| 76) Pentachlorobenzene        | 14.898 | 250  | 1208522  | 82.318 | ng/uL  | 92       |
| 77) Carbazole                 | 17.733 | 167  | 4061695  | 79.718 | ng/ul# | 95       |
| 78) Di-n-butylphthalate       | 18.281 | 149  | 5194813  | 82.913 | ng/ul# | 93       |
| 80) Fluoranthene              | 19.333 | 202  | 5581221  | 84.196 | ng/ul  | 97       |
| 82) Pyrene                    | 19.692 | 202  | 5591827  | 83.201 | ng/ul# | 93       |
| 83) Butylbenzylphthalate      | 20.557 | 149  | 2481860  | 90.017 | ng/ul# | 79       |
| 84) 3,3'-Dichlorobenzidine    | 21.322 | 252  | 1516289  | 67.479 | ng/ul# | 96       |
| 85) Benzo(a)anthracene        | 21.398 | 228  | 5480162  | 81.979 | ng/ul  | 95       |
| 86) Bis(2-ethylhexyl)phtha... | 21.316 | 149  | 3608457  | 87.010 | ng/ul# | 82       |
| 87) Chrysene                  | 21.457 | 228  | 5294592  | 82.090 | ng/ul  | 98       |
| 89) Di-n-octyl phthalate      | 22.263 | 149  | 6219148  | 88.381 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.121 | 252  | 5494440  | 82.341 | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene      | 23.174 | 252  | 5184336  | 85.143 | ng/ul# | 98       |
| 93) Benzo(a)pyrene            | 23.774 | 252  | 5092610  | 82.333 | ng/ul  | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.457 | 276  | 5275850  | 81.568 | ng/ul# | 92       |
| 95) Dibenzo(a,h)anthracene    | 26.468 | 278  | 4569166  | 81.468 | ng/ul# | 95       |
| 96) Benzo(g,h,i)perylene      | 27.239 | 276  | 4367941  | 81.901 | ng/ul# | 91       |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

**Instrument :**

BNA\_P

**ClientSampleId :**

SSTD080627

**Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 04/22/2022  
Supervised By :mohammad ahmed 04/26/2022

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042122\  
 Data File : BP010012.D  
 Acq On : 21 Apr 2022 13:19  
 Operator : CG/JU  
 Sample : SST080627  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SST080627

Manual IntegrationsAPPROVED

Quant Time: Apr 21 14:25:12 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP042122.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Apr 21 13:57:49 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/22/2022  
 Supervised By :mohammad ahmed 04/26/2022

