

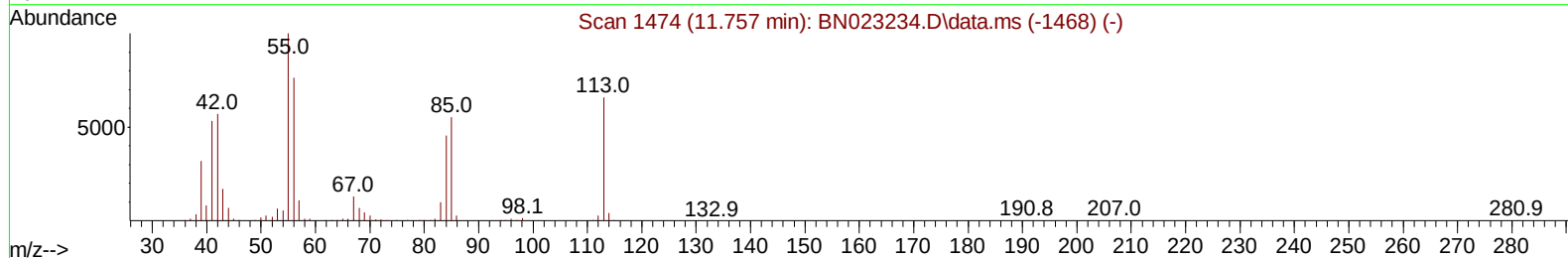
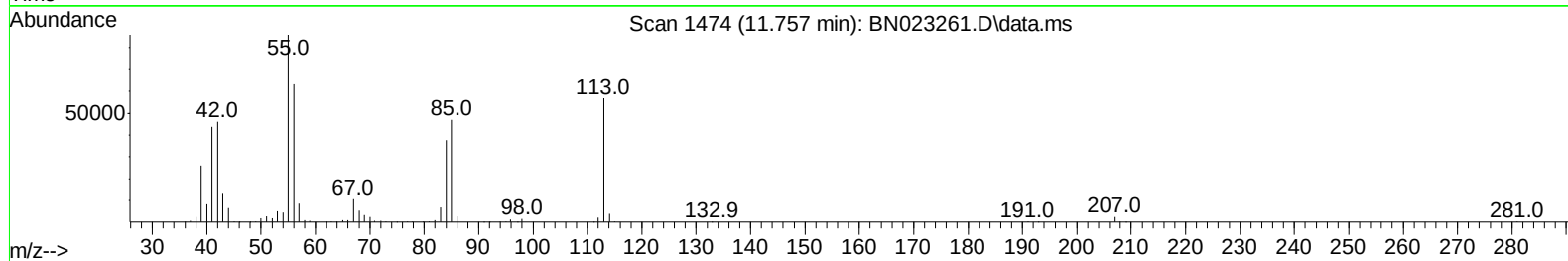
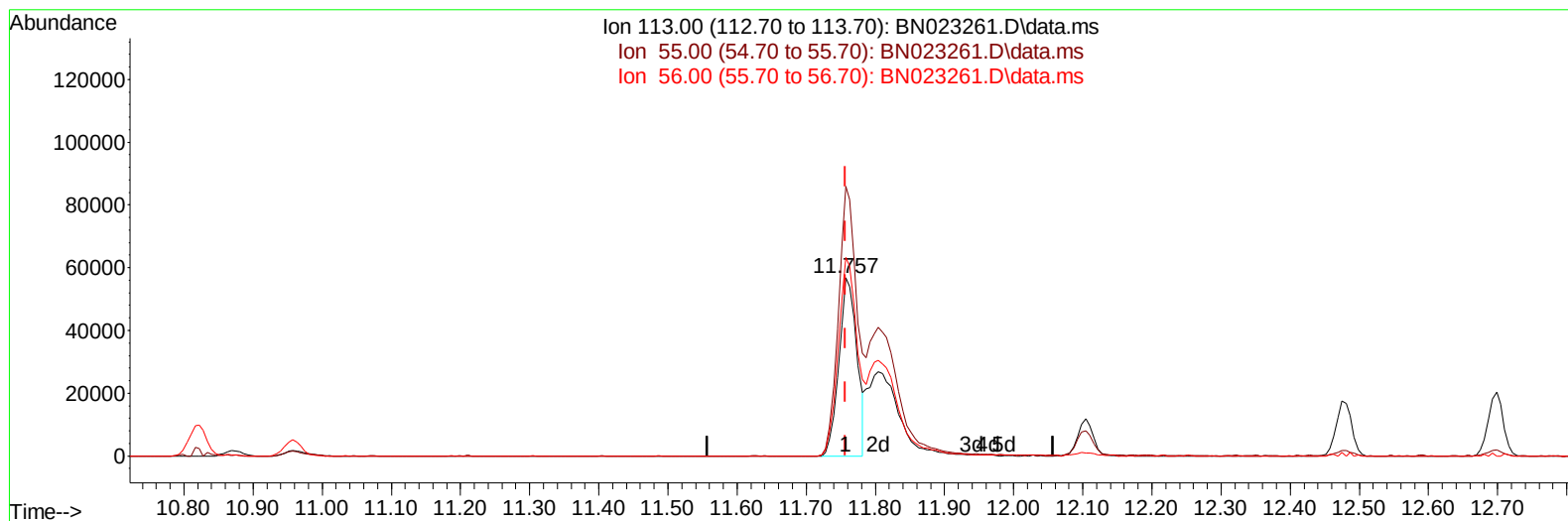
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN121622\  
 Data File : BN023261.D  
 Acq On : 17 Dec 2022 04:48  
 Operator : CG/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 29 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Dec 17 05:27:18 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN121522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Dec 15 23:22:36 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/19/2022  
 Supervised By :mohammad ahmed 12/19/2022



TIC: BN023261.D\data.ms

**(34) Caprolactam**

**11.757min (+ 0.000) 10.73 ng/ul**

**response 102238**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 154.60 | 151.33 |
| 56.00  | 114.60 | 111.64 |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN121622\  
 Data File : BN023261.D  
 Acq On : 17 Dec 2022 04:48  
 Operator : CG/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 29 Sample Multiplier: 1

Instrument :

BNA\_N

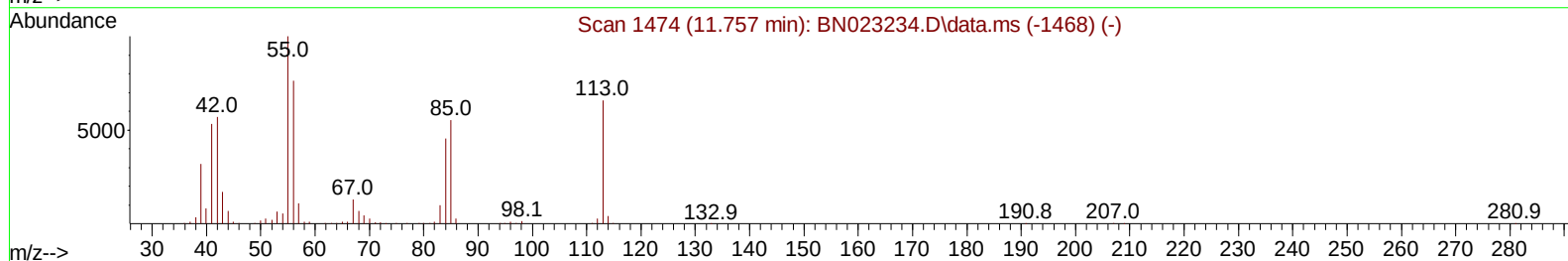
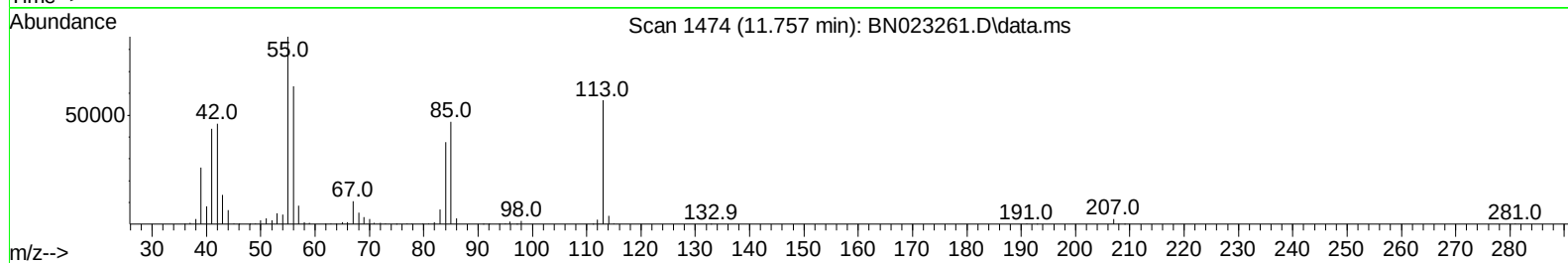
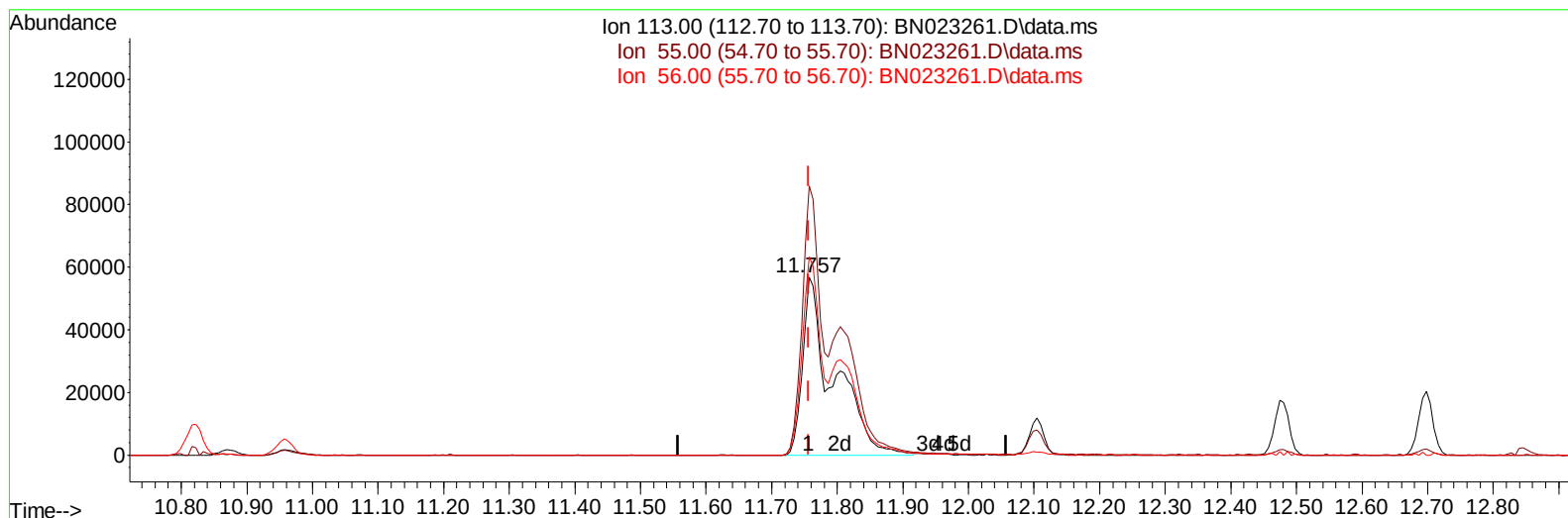
LabSampleId :

SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Dec 17 05:27:18 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN121522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Dec 15 23:22:36 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/19/2022  
 Supervised By :mohammad ahmed 12/19/2022



TIC: BN023261.D\data.ms

**(34) Caprolactam**

11.757min (+ 0.000) 19.60 ng/ul m

response 186702

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 154.60 | 151.33 |
| 56.00  | 114.60 | 111.64 |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN121622\  
 Data File : BN023261.D  
 Acq On : 17 Dec 2022 04:48  
 Operator : CG/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 29 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Dec 17 05:27:18 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN121522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Dec 15 23:22:36 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/19/2022  
 Supervised By :mohammad ahmed 12/19/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.999  | 152  | 334091   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.822 | 136  | 1611709  | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.651 | 164  | 1059782  | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.398 | 188  | 2432520  | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.586 | 240  | 2190150  | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 24.039 | 264  | 2216533  | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.346  | 96   | 58231    | 7.211  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.775  | 84   | 464773   | 19.498 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.158  | 99   | 604268   | 19.447 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.322  | 67   | 356962   | 19.741 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.528  | 132  | 485734   | 20.746 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.716  | 113  | 508795   | 20.394 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.169  | 128  | 259164   | 20.812 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.899  | 143  | 277639   | 21.169 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.434 | 165  | 508642   | 20.543 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.957 | 131  | 803853   | 20.177 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.063 | 166  | 1628330  | 20.503 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.345 | 160  | 1859723  | 20.880 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.845 | 143  | 340032   | 19.170 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.639 | 176  | 1371459  | 20.272 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.757 | 200  | 281378   | 19.650 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.498 | 188  | 2212915  | 20.299 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.792 | 212  | 2611867  | 21.923 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.880 | 264  | 2290594  | 21.084 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.381  | 88   | 61357    | 7.332  | ng/uL  | 96       |
| 5) Pyridine                   | 3.793  | 79   | 457908   | 18.800 | ng/ul  | 98       |
| 6) Benzaldehyde               | 7.128  | 77   | 238676   | 21.444 | ng/ul  | 99       |
| 8) Phenol                     | 7.181  | 94   | 611883   | 19.495 | ng/ul  | 98       |
| 10) Bis(2-Chloroethyl)ether   | 7.416  | 93   | 486764   | 19.967 | ng/ul  | 100      |
| 12) 2-Chlorophenol            | 7.558  | 128  | 487605   | 20.277 | ng/ul  | 97       |
| 13) 2-Methylphenol            | 8.446  | 108  | 476249   | 19.943 | ng/ul  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.540  | 45   | 618475   | 20.031 | ng/ul  | 99       |
| 16) Acetophenone              | 8.834  | 105  | 785798   | 20.818 | ng/ul  | 98       |
| 17) N-Nitroso-di-n-propyla... | 8.816  | 70   | 402435   | 21.291 | ng/ul  | 100      |
| 18) 4-Methylphenol            | 8.775  | 108  | 532715   | 20.101 | ng/ul  | 95       |
| 19) Hexachloroethane          | 9.087  | 117  | 191864   | 20.716 | ng/ul  | 97       |
| 22) Nitrobenzene              | 9.216  | 77   | 597155   | 19.999 | ng/ul  | 100      |
| 23) Isophorone                | 9.746  | 82   | 1155694  | 20.520 | ng/ul  | 99       |
| 25) 2-Nitrophenol             | 9.928  | 139  | 301949   | 21.108 | ng/ul  | 99       |
| 26) 2,4-Dimethylphenol        | 9.987  | 107  | 586083   | 20.241 | ng/ul  | 98       |
| 27) Bis(2-Chloroethoxy)met... | 10.228 | 93   | 709009   | 19.680 | ng/ul  | 100      |
| 29) 2,4-Dichlorophenol        | 10.463 | 162  | 502572   | 20.513 | ng/ul  | 99       |
| 30) Naphthalene               | 10.869 | 128  | 1700566  | 19.750 | ng/ul  | 99       |
| 32) 4-Chloroaniline           | 10.981 | 127  | 804841   | 20.249 | ng/ul  | 98       |
| 33) Hexachlorobutadiene       | 11.157 | 225  | 287877   | 20.610 | ng/ul  | 99       |
| 34) Caprolactam               | 11.757 | 113  | 186702m  | 19.599 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 12.104 | 107  | 582483   | 20.990 | ng/ul  | 96       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN121622\  
 Data File : BN023261.D  
 Acq On : 17 Dec 2022 04:48  
 Operator : CG/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 29 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 LabSampleId :  
 SSTDCCC020

Manual IntegrationsAPPROVED

Quant Time: Dec 17 05:27:18 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN121522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Dec 15 23:22:36 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/19/2022  
 Supervised By :mohammad ahmed 12/19/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 36) 2-Methylnaphthalene       | 12.481 | 142  | 1178193  | 20.303 | ng/ul | 97       |
| 37) 1-Methylnaphthalene       | 12.698 | 142  | 1212174  | 20.270 | ng/ul | 96       |
| 39) 1,2,4,5-Tetrachloroben... | 12.845 | 216  | 609164   | 20.322 | ng/ul | 97       |
| 40) Hexachlorocyclopentadiene | 12.828 | 237  | 353862   | 19.626 | ng/ul | 97       |
| 41) 2,4,6-Trichlorophenol     | 13.087 | 196  | 430924   | 20.900 | ng/ul | 96       |
| 42) 2,4,5-Trichlorophenol     | 13.151 | 196  | 475346   | 20.831 | ng/ul | 98       |
| 43) 1,1'-Biphenyl             | 13.487 | 154  | 1602277  | 20.008 | ng/ul | 99       |
| 44) 2-Chloronaphthalene       | 13.528 | 162  | 1254575  | 20.118 | ng/ul | 99       |
| 45) 2-Nitroaniline            | 13.734 | 65   | 380048   | 21.030 | ng/ul | 98       |
| 47) Dimethylphthalate         | 14.110 | 163  | 1623651  | 20.261 | ng/ul | 99       |
| 48) 2,6-Dinitrotoluene        | 14.228 | 165  | 339866   | 21.905 | ng/ul | 99       |
| 50) Acenaphthylene            | 14.375 | 152  | 2007773  | 20.499 | ng/ul | 98       |
| 51) 3-Nitroaniline            | 14.557 | 138  | 315675   | 19.981 | ng/ul | 95       |
| 52) Acenaphthene              | 14.716 | 153  | 1365854  | 20.059 | ng/ul | 99       |
| 53) 2,4-Dinitrophenol         | 14.757 | 184  | 181892   | 18.398 | ng/ul | 95       |
| 55) 4-Nitrophenol             | 14.857 | 109  | 255660   | 19.488 | ng/ul | 94       |
| 56) Dibenzofuran              | 15.051 | 168  | 1902530  | 20.068 | ng/ul | 100      |
| 57) 2,4-Dinitrotoluene        | 15.010 | 165  | 507563   | 22.161 | ng/ul | 95       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.275 | 232  | 414441   | 20.978 | ng/ul | 98       |
| 59) Diethylphthalate          | 15.469 | 149  | 1640551  | 20.579 | ng/ul | 99       |
| 61) Fluorene                  | 15.698 | 166  | 1526060  | 19.878 | ng/ul | 100      |
| 62) 4-Chlorophenyl-phenyle... | 15.692 | 204  | 728292   | 19.793 | ng/ul | 99       |
| 63) 4-Nitroaniline            | 15.722 | 138  | 302677   | 19.673 | ng/ul | 98       |
| 66) 4,6-Dinitro-2-methylph... | 15.775 | 198  | 279270   | 19.977 | ng/ul | 99       |
| 67) N-Nitrosodiphenylamine    | 15.904 | 169  | 1336034  | 19.845 | ng/ul | 99       |
| 68) 4-Bromophenyl-phenylether | 16.586 | 248  | 474602   | 20.220 | ng/ul | 96       |
| 69) Hexachlorobenzene         | 16.698 | 284  | 560692   | 20.432 | ng/ul | 98       |
| 70) Atrazine                  | 16.857 | 200  | 506295   | 20.580 | ng/ul | 99       |
| 71) Pentachlorophenol         | 17.045 | 266  | 362610   | 20.011 | ng/ul | 98       |
| 72) Phenanthrene              | 17.439 | 178  | 2584249  | 20.077 | ng/ul | 100      |
| 74) Anthracene                | 17.533 | 178  | 2550657  | 19.860 | ng/ul | 98       |
| 75) 1,2,3,4-Tetrachloroben... | 13.451 | 216  | 609854   | 20.160 | ng/uL | 98       |
| 76) Pentachlorobenzene        | 14.969 | 250  | 637641   | 20.801 | ng/uL | 96       |
| 77) Carbazole                 | 17.798 | 167  | 2355572  | 19.918 | ng/ul | 100      |
| 78) Di-n-butylphthalate       | 18.363 | 149  | 2896038  | 20.624 | ng/ul | 99       |
| 80) Fluoranthene              | 19.457 | 202  | 3115247  | 22.483 | ng/ul | 97       |
| 82) Pyrene                    | 19.822 | 202  | 3178772  | 21.839 | ng/ul | 96       |
| 83) Butylbenzylphthalate      | 20.710 | 149  | 1359529  | 22.571 | ng/ul | 100      |
| 84) 3,3'-Dichlorobenzidine    | 21.498 | 252  | 993477   | 20.642 | ng/ul | 98       |
| 85) Benzo(a)anthracene        | 21.569 | 228  | 2948453  | 20.094 | ng/ul | 95       |
| 86) Bis(2-ethylhexyl)phtha... | 21.486 | 149  | 1961934  | 23.036 | ng/ul | 100      |
| 87) Chrysene                  | 21.621 | 228  | 2815299  | 20.330 | ng/ul | 97       |
| 89) Di-n-octyl phthalate      | 22.433 | 149  | 3318335  | 24.721 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 23.286 | 252  | 3010152  | 21.531 | ng/ul | 97       |
| 91) Benzo(k)fluoranthene      | 23.339 | 252  | 2870466  | 21.575 | ng/ul | 100      |
| 93) Benzo(a)pyrene            | 23.933 | 252  | 2503450  | 20.837 | ng/ul | 98       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.604 | 276  | 2863108  | 18.811 | ng/ul | 98       |
| 95) Dibenzo(a,h)anthracene    | 26.627 | 278  | 2349472  | 18.932 | ng/ul | 97       |
| 96) Benzo(g,h,i)perylene      | 27.398 | 276  | 2211961  | 18.368 | ng/ul | 98       |

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

**Instrument :**  
BNA\_N  
**LabSampleId :**  
SSTDCCC020

**Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 12/19/2022  
Supervised By :mohammad ahmed 12/19/2022

Manual IntegrationsAPPROVED

Quant Time: Dec 17 05:27:18 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN121522.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Dec 15 23:22:36 2022  
Response via : Initial Calibration

