

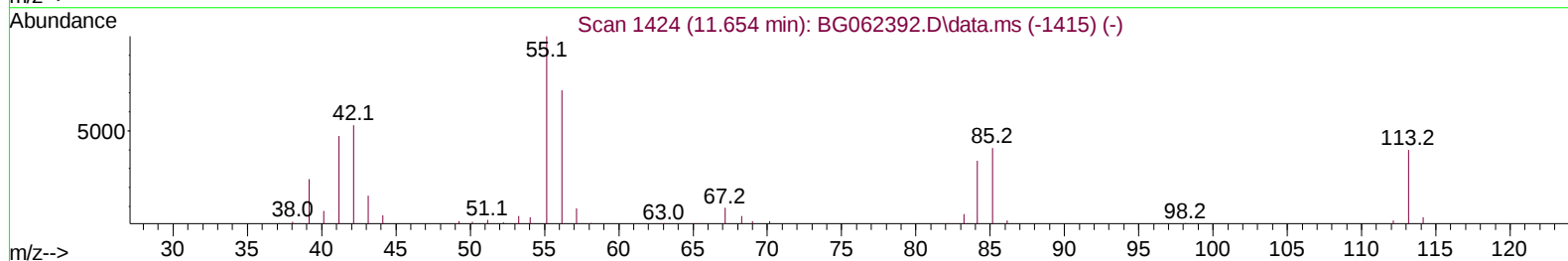
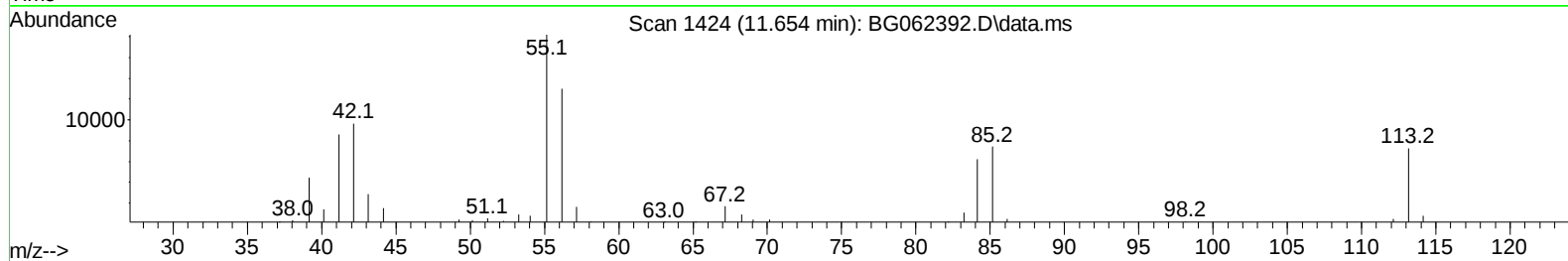
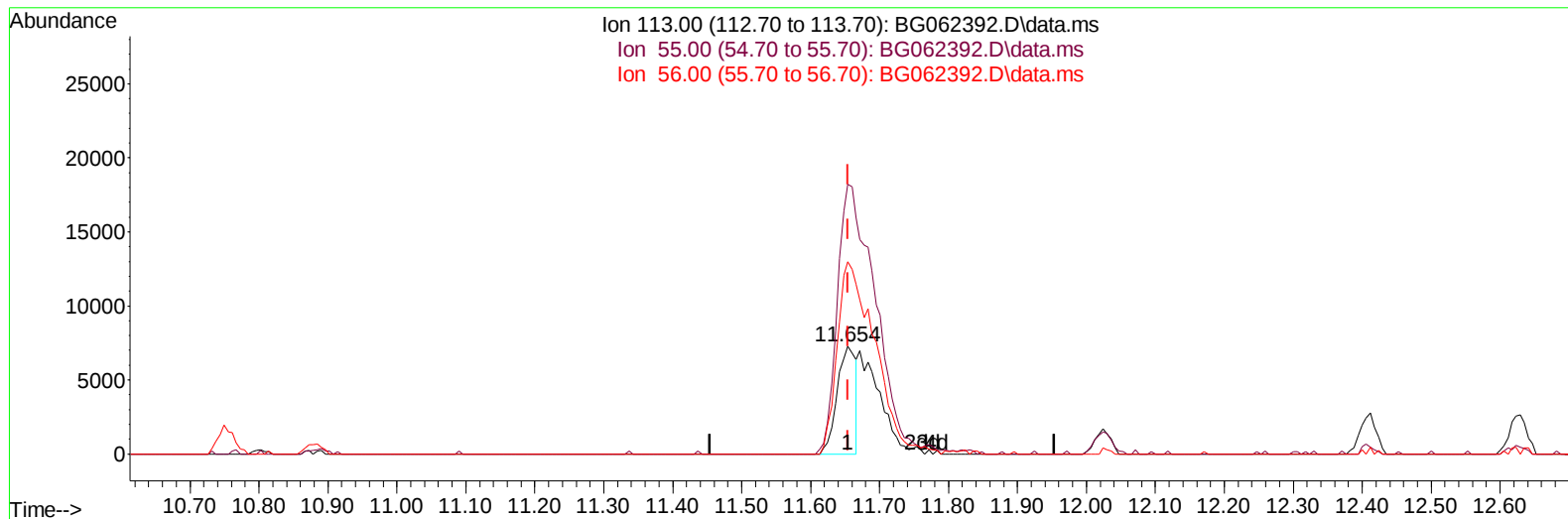
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG081424\  
Data File : BG062392.D  
Acq On : 14 Aug 2024 12:51  
Operator : MA/JU  
Sample : SSTD02048  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SSTD020433

Manual IntegrationsAPPROVED

Quant Time: Aug 14 15:14:52 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG081424.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Aug 14 15:13:32 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/15/2024  
Supervised By :mohammad ahmed 08/16/2024



TIC: BG062392.D\data.ms

**(34) Caprolactam**

11.654min ( 0.000) 10.56 ng/ul

response 13696

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 249.80 | 249.83 |
| 56.00  | 178.20 | 178.23 |
| 0.00   | 0.00   | 0.00   |

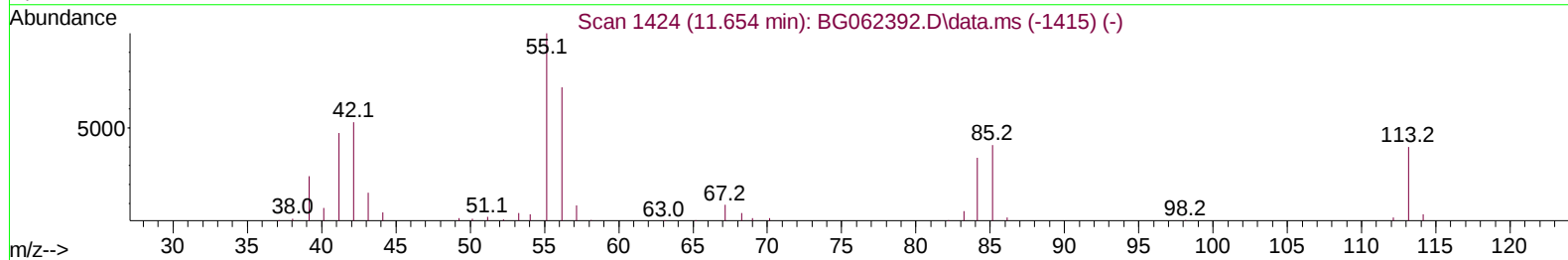
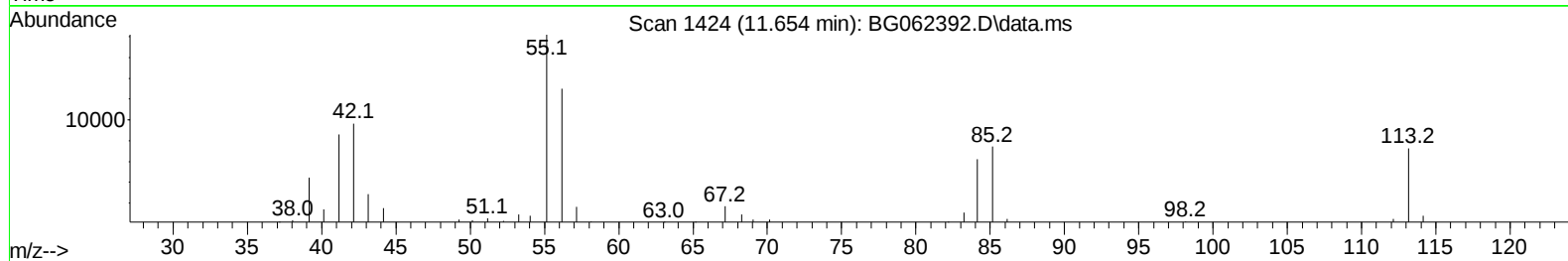
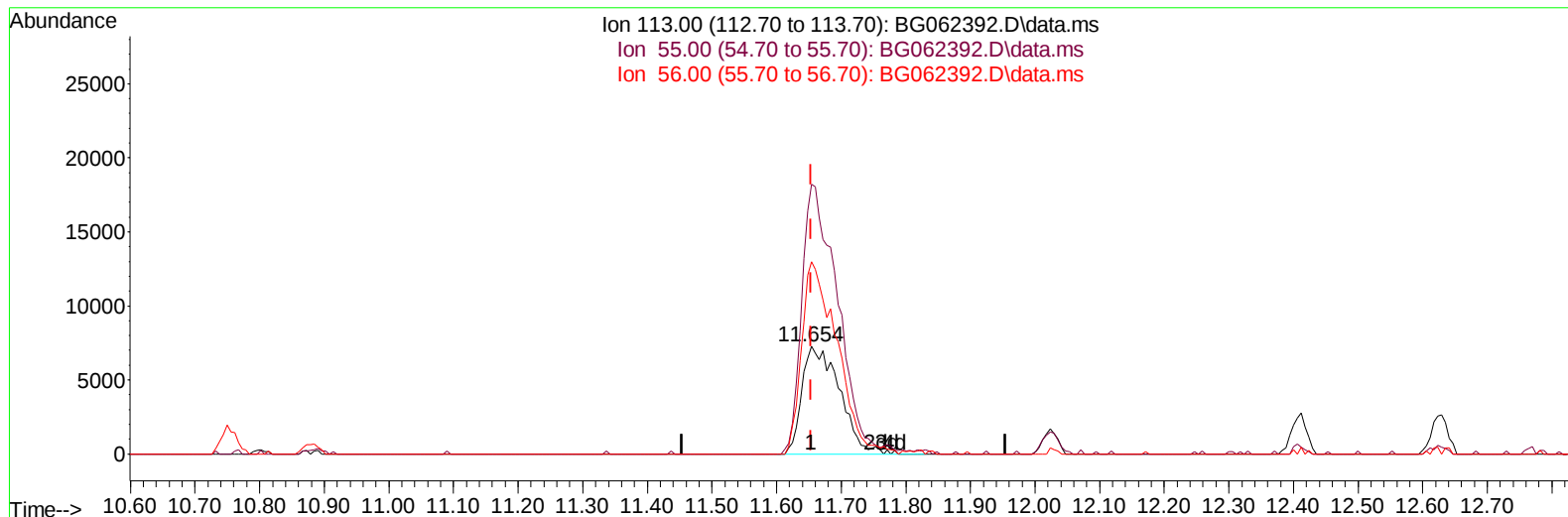
Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG081424\  
Data File : BG062392.D  
Acq On : 14 Aug 2024 12:51  
Operator : MA/JU  
Sample : SSTD02048  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SSTD020433

Manual IntegrationsAPPROVED

Quant Time: Aug 14 15:14:52 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG081424.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Aug 14 15:13:32 2024  
Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/15/2024  
Supervised By :mohammad ahmed 08/16/2024



TIC: BG062392.D\data.ms

(34) Caprolactam

11.654min ( 0.000) 22.60 ng/ul m

response 29316

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 249.80 | 249.83 |
| 56.00  | 178.20 | 178.23 |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG081424\  
 Data File : BG062392.D  
 Acq On : 14 Aug 2024 12:51  
 Operator : MA/JU  
 Sample : SST002048  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SST0020433

Manual IntegrationsAPPROVED

Quant Time: Aug 14 15:18:10 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG081424.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 14 15:13:32 2024  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/15/2024  
 Supervised By :mohammad ahmed 08/16/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.959  | 152  | 43047    | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.755 | 136  | 211186   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.579 | 164  | 168321   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.317 | 188  | 412397   | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.564 | 240  | 414377   | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 24.654 | 264  | 478376   | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.418  | 96   | 9772     | 9.358  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.823  | 84   | 68735    | 21.424 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.119  | 99   | 83790    | 20.257 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.283  | 67   | 55234    | 21.551 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.489  | 132  | 59481    | 20.230 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.658  | 113  | 72155    | 20.848 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.110  | 128  | 27742    | 21.026 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.833  | 143  | 26763    | 21.354 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.367 | 165  | 70620    | 19.935 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.879 | 131  | 108637   | 20.864 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.986 | 166  | 270378   | 20.746 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.274 | 160  | 297159   | 20.245 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.762 | 143  | 41432    | 21.995 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.572 | 176  | 242156   | 20.696 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.684 | 200  | 32004    | 23.766 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.417 | 188  | 398415   | 20.128 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.702 | 212  | 503979   | 20.407 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.443 | 264  | 506385   | 19.876 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.453  | 88   | 9125     | 8.136  | ng/uL  | 100      |
| 5) Pyridine                   | 3.847  | 79   | 73879    | 22.075 | ng/ul  | 100      |
| 6) Benzaldehyde               | 7.095  | 77   | 55142    | 24.685 | ng/ul  | 100      |
| 8) Phenol                     | 7.142  | 94   | 88223    | 20.199 | ng/ul  | 100      |
| 10) Bis(2-Chloroethyl)ether   | 7.377  | 93   | 67362    | 20.874 | ng/ul  | 100      |
| 12) 2-Chlorophenol            | 7.524  | 128  | 61975    | 19.949 | ng/ul  | 100      |
| 13) 2-Methylphenol            | 8.394  | 108  | 68055    | 20.710 | ng/ul  | 100      |
| 14) 2,2'-oxybis(1-Chloropr... | 8.476  | 45   | 139478   | 21.077 | ng/ul  | 100      |
| 16) Acetophenone              | 8.776  | 105  | 120832   | 21.820 | ng/ul  | 100      |
| 17) N-Nitroso-di-n-propyla... | 8.758  | 70   | 65494    | 22.241 | ng/ul  | 100      |
| 18) 4-Methylphenol            | 8.717  | 108  | 77524    | 20.916 | ng/ul  | 100      |
| 19) Hexachloroethane          | 9.040  | 117  | 24714    | 20.999 | ng/ul  | 100      |
| 22) Nitrobenzene              | 9.151  | 77   | 77065    | 20.713 | ng/ul  | 100      |
| 23) Isophorone                | 9.674  | 82   | 186579   | 20.811 | ng/ul  | 100      |
| 25) 2-Nitrophenol             | 9.862  | 139  | 30660    | 20.728 | ng/ul  | 100      |
| 26) 2,4-Dimethylphenol        | 9.921  | 107  | 83853    | 19.780 | ng/ul  | 100      |
| 27) Bis(2-Chloroethoxy)met... | 10.162 | 93   | 102173   | 20.277 | ng/ul  | 100      |
| 29) 2,4-Dichlorophenol        | 10.397 | 162  | 72761    | 20.175 | ng/ul  | 100      |
| 30) Naphthalene               | 10.802 | 128  | 247185   | 20.557 | ng/ul  | 100      |
| 32) 4-Chloroaniline           | 10.902 | 127  | 106627   | 20.958 | ng/ul  | 100      |
| 33) Hexachlorobutadiene       | 11.096 | 225  | 53934    | 19.851 | ng/ul  | 100      |
| 34) Caprolactam               | 11.654 | 113  | 29316m   | 22.596 | ng/ul  | 100      |
| 35) 4-Chloro-3-methylphenol   | 12.024 | 107  | 87977    | 21.550 | ng/ul  | 100      |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG081424\  
 Data File : BG062392.D  
 Acq On : 14 Aug 2024 12:51  
 Operator : MA/JU  
 Sample : SST002048  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SST0020433

## Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 08/15/2024  
 Supervised By :mohammad ahmed 08/16/2024

Quant Time: Aug 14 15:18:10 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG081424.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 14 15:13:32 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 36) 2-Methylnaphthalene       | 12.406 | 142  | 174548   | 20.896 | ng/ul | 100      |
| 37) 1-Methylnaphthalene       | 12.629 | 142  | 181095   | 21.021 | ng/ul | 100      |
| 39) 1,2,4,5-Tetrachloroben... | 12.776 | 216  | 111397   | 20.285 | ng/ul | 100      |
| 40) Hexachlorocyclopentadiene | 12.758 | 237  | 53342    | 23.546 | ng/ul | 100      |
| 41) 2,4,6-Trichlorophenol     | 13.011 | 196  | 67880    | 21.270 | ng/ul | 100      |
| 42) 2,4,5-Trichlorophenol     | 13.081 | 196  | 71038    | 20.858 | ng/ul | 100      |
| 43) 1,1'-Biphenyl             | 13.416 | 154  | 257621   | 20.382 | ng/ul | 100      |
| 44) 2-Chloronaphthalene       | 13.457 | 162  | 197912   | 20.184 | ng/ul | 100      |
| 45) 2-Nitroaniline            | 13.651 | 65   | 53794    | 21.300 | ng/ul | 100      |
| 47) Dimethylphthalate         | 14.033 | 163  | 271988   | 20.773 | ng/ul | 100      |
| 48) 2,6-Dinitrotoluene        | 14.145 | 165  | 38291    | 21.381 | ng/ul | 100      |
| 50) Acenaphthylene            | 14.303 | 152  | 313687   | 20.066 | ng/ul | 100      |
| 51) 3-Nitroaniline            | 14.474 | 138  | 44554    | 21.435 | ng/ul | 100      |
| 52) Acenaphthene              | 14.644 | 153  | 236077   | 20.214 | ng/ul | 100      |
| 53) 2,4-Dinitrophenol         | 14.679 | 184  | 20253    | 25.136 | ng/ul | 100      |
| 55) 4-Nitrophenol             | 14.779 | 109  | 34750    | 21.066 | ng/ul | 100      |
| 56) Dibenzofuran              | 14.979 | 168  | 333077   | 20.640 | ng/ul | 100      |
| 57) 2,4-Dinitrotoluene        | 14.932 | 165  | 61369    | 24.225 | ng/ul | 100      |
| 58) 2,3,4,6-Tetrachlorophenol | 15.202 | 232  | 67234    | 21.586 | ng/ul | 100      |
| 59) Diethylphthalate          | 15.402 | 149  | 276685   | 20.898 | ng/ul | 100      |
| 61) Fluorene                  | 15.625 | 166  | 265108   | 20.885 | ng/ul | 100      |
| 62) 4-Chlorophenyl-phenyle... | 15.619 | 204  | 136277   | 20.842 | ng/ul | 100      |
| 63) 4-Nitroaniline            | 15.631 | 138  | 49103    | 24.010 | ng/ul | 100      |
| 66) 4,6-Dinitro-2-methylph... | 15.696 | 198  | 35914    | 23.315 | ng/ul | 100      |
| 67) N-Nitrosodiphenylamine    | 15.831 | 169  | 241678   | 19.722 | ng/ul | 100      |
| 68) 4-Bromophenyl-phenylether | 16.512 | 248  | 92780    | 19.926 | ng/ul | 100      |
| 69) Hexachlorobenzene         | 16.630 | 284  | 108886   | 20.007 | ng/ul | 100      |
| 70) Atrazine                  | 16.782 | 200  | 100990   | 21.140 | ng/ul | 100      |
| 71) Pentachlorophenol         | 16.970 | 266  | 57200    | 19.546 | ng/ul | 100      |
| 72) Phenanthrene              | 17.364 | 178  | 468115   | 20.213 | ng/ul | 100      |
| 74) Anthracene                | 17.452 | 178  | 482708   | 20.297 | ng/ul | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 13.381 | 216  | 115086   | 19.069 | ng/ul | 100      |
| 76) Pentachlorobenzene        | 14.897 | 250  | 121428   | 19.087 | ng/ul | 100      |
| 77) Carbazole                 | 17.716 | 167  | 410829   | 20.973 | ng/ul | 100      |
| 78) Di-n-butylphthalate       | 18.286 | 149  | 491789   | 21.092 | ng/ul | 100      |
| 80) Fluoranthene              | 19.367 | 202  | 573651   | 20.480 | ng/ul | 100      |
| 82) Pyrene                    | 19.731 | 202  | 613239   | 20.557 | ng/ul | 100      |
| 83) Butylbenzylphthalate      | 20.624 | 149  | 204693   | 20.038 | ng/ul | 100      |
| 84) 3,3'-Dichlorobenzidine    | 21.464 | 252  | 201527   | 21.432 | ng/ul | 100      |
| 85) Benzo(a)anthracene        | 21.546 | 228  | 611523   | 20.122 | ng/ul | 100      |
| 86) Bis(2-ethylhexyl)phtha... | 21.470 | 149  | 303394   | 19.383 | ng/ul | 100      |
| 87) Chrysene                  | 21.611 | 228  | 578892   | 20.169 | ng/ul | 100      |
| 89) Di-n-octyl phthalate      | 22.645 | 149  | 513066   | 19.064 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 23.679 | 252  | 572419   | 19.779 | ng/ul | 100      |
| 91) Benzo(k)fluoranthene      | 23.743 | 252  | 590945   | 20.004 | ng/ul | 100      |
| 93) Benzo(a)pyrene            | 24.507 | 252  | 546540   | 19.895 | ng/ul | 100      |
| 94) Indeno(1,2,3-cd)pyrene    | 28.149 | 276  | 688216   | 19.448 | ng/ul | 100      |
| 95) Dibenzo(a,h)anthracene    | 28.208 | 278  | 558205   | 19.269 | ng/ul | 100      |
| 96) Benzo(g,h,i)perylene      | 29.236 | 276  | 528236   | 19.602 | ng/ul | 100      |

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

**Instrument :**

BNA\_G

**ClientSampleId :**

SSTD020433

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 08/15/2024

Supervised By :mohammad ahmed 08/16/2024

