

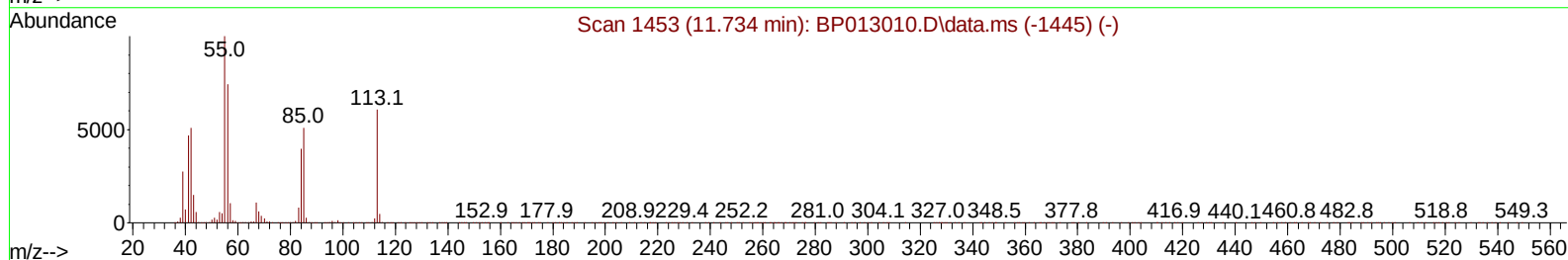
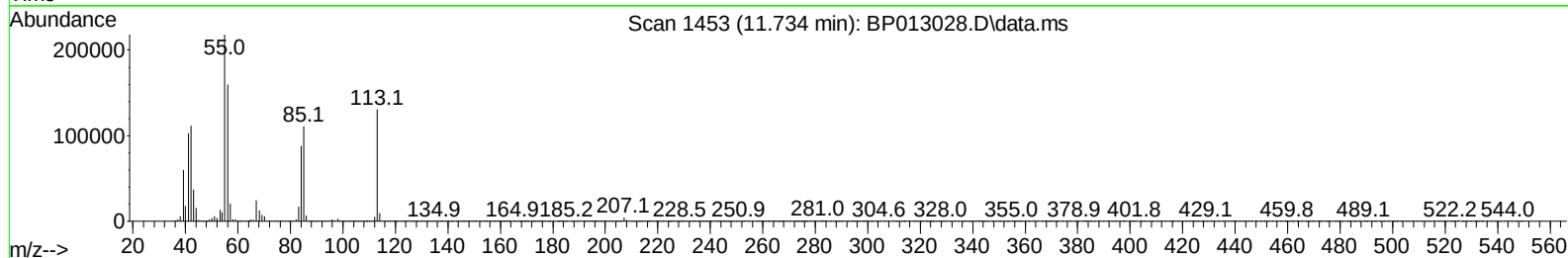
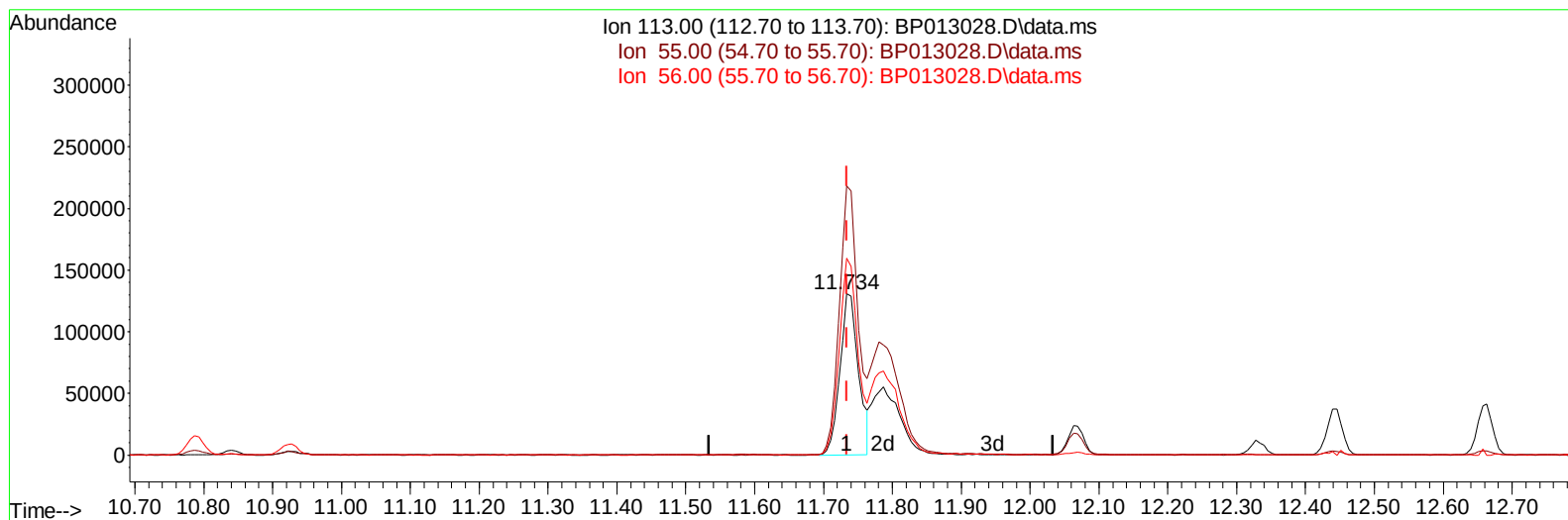
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP120822\  
 Data File : BP013028.D  
 Acq On : 09 Dec 2022 20:49  
 Operator : CG/JU  
 Sample : PB149458BS  
 Misc :  
 ALS Vial : 55 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS458

Manual IntegrationsAPPROVED

Quant Time: Dec 09 21:59:56 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Dec 08 00:35:07 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/10/2022  
 Supervised By :mohammad ahmed 12/10/2022



TIC: BP013028.D\data.ms

**(34) Caprolactam**

**11.734min (-0.000) 18.28 ng/ul**

**response 244230**

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 177.40 | 166.93 |
| 56.00  | 124.50 | 122.05 |
| 0.00   | 0.00   | 0.00   |

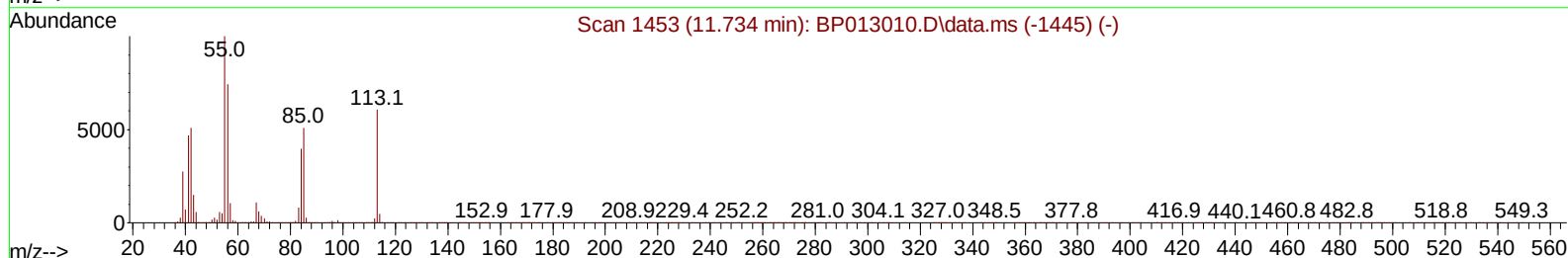
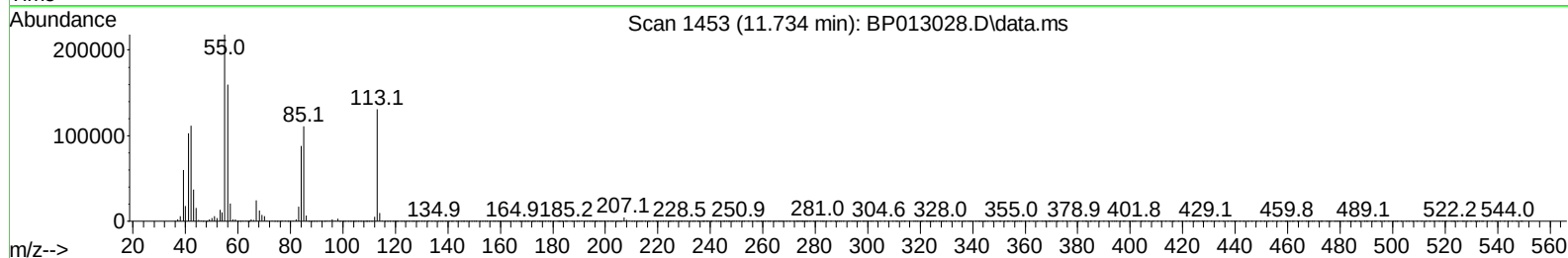
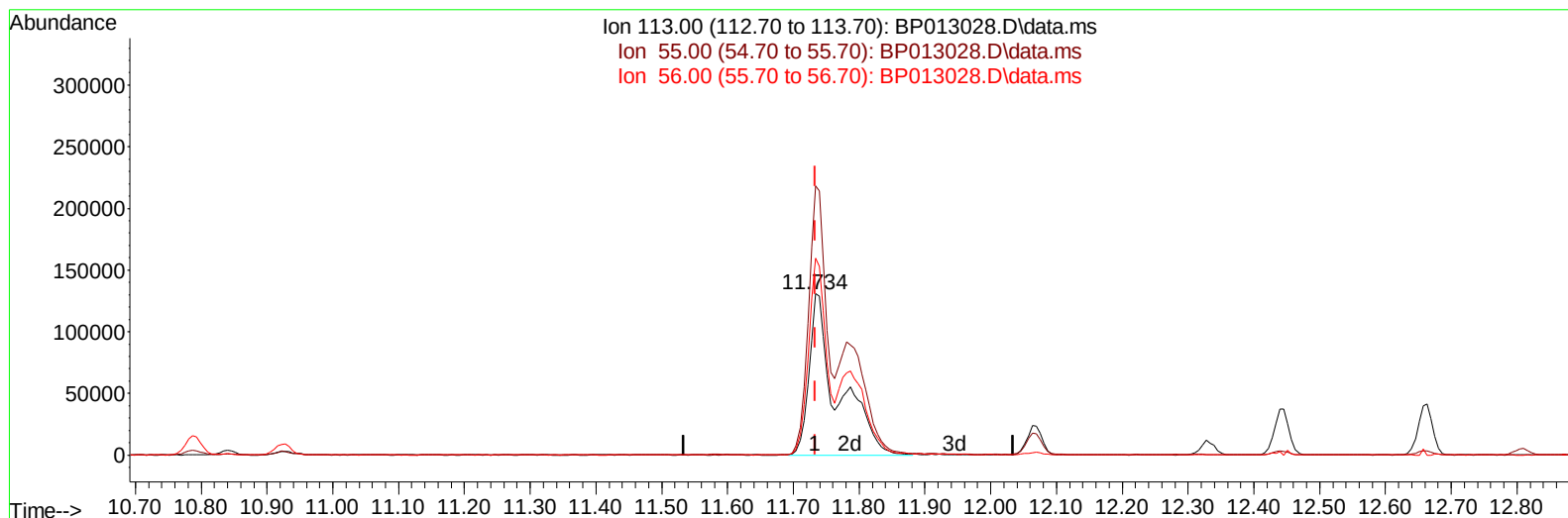
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP120822\  
 Data File : BP013028.D  
 Acq On : 09 Dec 2022 20:49  
 Operator : CG/JU  
 Sample : PB149458BS  
 Misc :  
 ALS Vial : 55 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS458

Manual IntegrationsAPPROVED

Quant Time: Dec 09 21:59:56 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Dec 08 00:35:07 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/10/2022  
 Supervised By :mohammad ahmed 12/10/2022



TIC: BP013028.D\data.ms

**(34) Caprolactam**

11.734min (-0.000) 29.87 ng/ul m

response 399062

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100.00 | 100.00 |
| 55.00  | 177.40 | 166.93 |
| 56.00  | 124.50 | 122.05 |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP120822\  
 Data File : BP013028.D  
 Acq On : 09 Dec 2022 20:49  
 Operator : CG/JU  
 Sample : PB149458BS  
 Misc :  
 ALS Vial : 55 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS458

Manual IntegrationsAPPROVED

Quant Time: Dec 09 21:59:56 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Dec 08 00:35:07 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/10/2022  
 Supervised By :mohammad ahmed 12/10/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.975  | 152  | 580299   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.787 | 136  | 2489730  | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.616 | 164  | 1633274  | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.369 | 188  | 3703574  | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.457 | 240  | 3282012  | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 23.945 | 264  | 3035681  | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.352  | 96   | 73662    | 5.427  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.781  | 84   | 995871   | 25.830 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.128  | 99   | 1435773  | 30.376 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.299  | 67   | 836138   | 29.325 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.505  | 132  | 1145481  | 30.785 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.687  | 113  | 1161549  | 29.992 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.146  | 128  | 574945   | 31.430 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.869  | 143  | 675036   | 32.778 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.404 | 165  | 1256688  | 31.416 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.928 | 131  | 1520841  | 26.931 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.022 | 166  | 3703028  | 29.500 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.310 | 160  | 4195102  | 30.415 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.810 | 143  | 712624   | 31.561 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.610 | 176  | 3188088  | 30.021 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.728 | 200  | 706367   | 29.638 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.469 | 188  | 5065505  | 30.317 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.698 | 212  | 5908680  | 31.365 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.786 | 264  | 4705169  | 31.096 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.387  | 88   | 150787   | 10.643 | ng/uL  | 97       |
| 5) Pyridine                   | 3.799  | 79   | 1009988  | 26.358 | ng/ul  | 99       |
| 6) Benzaldehyde               | 7.110  | 77   | 704940   | 38.088 | ng/ul  | 99       |
| 8) Phenol                     | 7.157  | 94   | 1478253  | 30.932 | ng/ul  | 100      |
| 10) Bis(2-Chloroethyl)ether   | 7.393  | 93   | 1144338  | 29.174 | ng/ul  | 100      |
| 12) 2-Chlorophenol            | 7.534  | 128  | 1172517  | 30.729 | ng/ul  | 99       |
| 13) 2-Methylphenol            | 8.416  | 108  | 1129701  | 30.321 | ng/ul  | 98       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.504  | 45   | 1629779  | 30.279 | ng/ul  | 99       |
| 16) Acetophenone              | 8.804  | 105  | 1746657  | 29.413 | ng/ul  | 99       |
| 17) N-Nitroso-di-n-propyla... | 8.793  | 70   | 883681   | 29.592 | ng/ul  | 99       |
| 18) 4-Methylphenol            | 8.752  | 108  | 1234888  | 29.923 | ng/ul  | 100      |
| 19) Hexachloroethane          | 9.057  | 117  | 449650   | 29.489 | ng/ul  | 93       |
| 22) Nitrobenzene              | 9.187  | 77   | 1284188  | 29.917 | ng/ul  | 98       |
| 23) Isophorone                | 9.716  | 82   | 2579798  | 29.722 | ng/ul  | 100      |
| 25) 2-Nitrophenol             | 9.899  | 139  | 695217   | 31.156 | ng/ul  | 98       |
| 26) 2,4-Dimethylphenol        | 9.957  | 107  | 1281710  | 29.439 | ng/ul  | 100      |
| 27) Bis(2-Chloroethoxy)met... | 10.193 | 93   | 1569255  | 28.276 | ng/ul  | 99       |
| 29) 2,4-Dichlorophenol        | 10.434 | 162  | 1217770  | 31.078 | ng/ul  | 99       |
| 30) Naphthalene               | 10.840 | 128  | 3799295  | 29.178 | ng/ul  | 100      |
| 32) 4-Chloroaniline           | 10.951 | 127  | 1287896  | 23.047 | ng/ul  | 99       |
| 33) Hexachlorobutadiene       | 11.122 | 225  | 805743   | 29.949 | ng/ul  | 99       |
| 34) Caprolactam               | 11.734 | 113  | 399062m  | 29.869 | ng/ul  |          |
| 35) 4-Chloro-3-methylphenol   | 12.069 | 107  | 1255447  | 31.395 | ng/ul  | 97       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP120822\  
 Data File : BP013028.D  
 Acq On : 09 Dec 2022 20:49  
 Operator : CG/JU  
 Sample : PB149458BS  
 Misc :  
 ALS Vial : 55 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS458

Manual IntegrationsAPPROVED

Quant Time: Dec 09 21:59:56 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Dec 08 00:35:07 2022  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/10/2022  
 Supervised By :mohammad ahmed 12/10/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 36) 2-Methylnaphthalene       | 12.445 | 142  | 2657912  | 29.754 | ng/ul | 99       |
| 37) 1-Methylnaphthalene       | 12.663 | 142  | 2656728  | 28.920 | ng/ul | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 12.810 | 216  | 1589699  | 30.227 | ng/ul | 100      |
| 40) Hexachlorocyclopentadiene | 12.787 | 237  | 713161   | 20.838 | ng/ul | 100      |
| 41) 2,4,6-Trichlorophenol     | 13.045 | 196  | 1042942  | 30.857 | ng/ul | 99       |
| 42) 2,4,5-Trichlorophenol     | 13.116 | 196  | 1150416  | 31.687 | ng/ul | 98       |
| 43) 1,1'-Biphenyl             | 13.445 | 154  | 3527601  | 28.758 | ng/ul | 99       |
| 44) 2-Chloronaphthalene       | 13.492 | 162  | 2843592  | 29.454 | ng/ul | 99       |
| 45) 2-Nitroaniline            | 13.698 | 65   | 801043   | 34.222 | ng/ul | 97       |
| 47) Dimethylphthalate         | 14.069 | 163  | 3655475  | 29.382 | ng/ul | 99       |
| 48) 2,6-Dinitrotoluene        | 14.187 | 165  | 760831   | 32.907 | ng/ul | 99       |
| 50) Acenaphthylene            | 14.334 | 152  | 4366557  | 29.447 | ng/ul | 100      |
| 51) 3-Nitroaniline            | 14.522 | 138  | 789250   | 36.151 | ng/ul | 97       |
| 52) Acenaphthene              | 14.681 | 153  | 3026735  | 29.455 | ng/ul | 100      |
| 53) 2,4-Dinitrophenol         | 14.722 | 184  | 462014   | 29.233 | ng/ul | 98       |
| 55) 4-Nitrophenol             | 14.822 | 109  | 487868   | 31.666 | ng/ul | 99       |
| 56) Dibenzofuran              | 15.010 | 168  | 4292546  | 29.381 | ng/ul | 100      |
| 57) 2,4-Dinitrotoluene        | 14.975 | 165  | 1105009  | 32.902 | ng/ul | 98       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.239 | 232  | 1008286  | 30.433 | ng/ul | 96       |
| 59) Diethylphthalate          | 15.434 | 149  | 3646165  | 30.148 | ng/ul | 99       |
| 61) Fluorene                  | 15.663 | 166  | 3486371  | 29.754 | ng/ul | 99       |
| 62) 4-Chlorophenyl-phenyle... | 15.651 | 204  | 1845889  | 29.790 | ng/ul | 98       |
| 63) 4-Nitroaniline            | 15.686 | 138  | 843711   | 43.918 | ng/ul | 98       |
| 66) 4,6-Dinitro-2-methylph... | 15.739 | 198  | 689586   | 29.608 | ng/ul | 98       |
| 67) N-Nitrosodiphenylamine    | 15.869 | 169  | 3076552  | 29.963 | ng/ul | 100      |
| 68) 4-Bromophenyl-phenylether | 16.551 | 248  | 1194342  | 29.695 | ng/ul | 99       |
| 69) Hexachlorobenzene         | 16.669 | 284  | 1422882  | 29.794 | ng/ul | 98       |
| 70) Atrazine                  | 16.822 | 200  | 720965   | 18.018 | ng/ul | 100      |
| 71) Pentachlorophenol         | 17.016 | 266  | 846032   | 29.253 | ng/ul | 99       |
| 72) Phenanthrene              | 17.410 | 178  | 5789473  | 30.012 | ng/ul | 100      |
| 74) Anthracene                | 17.504 | 178  | 5841284  | 30.326 | ng/ul | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 13.416 | 216  | 1590600  | 30.102 | ng/uL | 100      |
| 76) Pentachlorobenzene        | 14.934 | 250  | 1571010  | 29.161 | ng/uL | 99       |
| 77) Carbazole                 | 17.775 | 167  | 5327704  | 32.844 | ng/ul | 100      |
| 78) Di-n-butylphthalate       | 18.316 | 149  | 6463738  | 32.727 | ng/ul | 99       |
| 80) Fluoranthene              | 19.374 | 202  | 6916799  | 32.349 | ng/ul | 99       |
| 82) Pyrene                    | 19.727 | 202  | 7091000  | 32.150 | ng/ul | 100      |
| 83) Butylbenzylphthalate      | 20.592 | 149  | 2926077  | 34.340 | ng/ul | 98       |
| 84) 3,3'-Dichlorobenzidine    | 21.369 | 252  | 2270671  | 30.625 | ng/ul | 100      |
| 85) Benzo(a)anthracene        | 21.439 | 228  | 6765458  | 31.052 | ng/ul | 100      |
| 86) Bis(2-ethylhexyl)phtha... | 21.345 | 149  | 4273349  | 35.553 | ng/ul | 99       |
| 87) Chrysene                  | 21.498 | 228  | 6319335  | 30.670 | ng/ul | 100      |
| 89) Di-n-octyl phthalate      | 22.304 | 149  | 7261507  | 42.698 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 23.180 | 252  | 6254332  | 32.309 | ng/ul | 99       |
| 91) Benzo(k)fluoranthene      | 23.233 | 252  | 5880988  | 30.638 | ng/ul | 99       |
| 93) Benzo(a)pyrene            | 23.839 | 252  | 5162316  | 30.627 | ng/ul | 100      |
| 94) Indeno(1,2,3-cd)pyrene    | 26.556 | 276  | 5611876  | 26.729 | ng/ul | 99       |
| 95) Dibenzo(a,h)anthracene    | 26.568 | 278  | 4999867  | 28.763 | ng/ul | 99       |
| 96) Benzo(g,h,i)perylene      | 27.356 | 276  | 4514294  | 26.783 | ng/ul | 99       |

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

**Instrument :**

BNA\_P

**ClientSampleId :**

SLCS458

**Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 12/10/2022

Supervised By :mohammad ahmed 12/10/2022

