

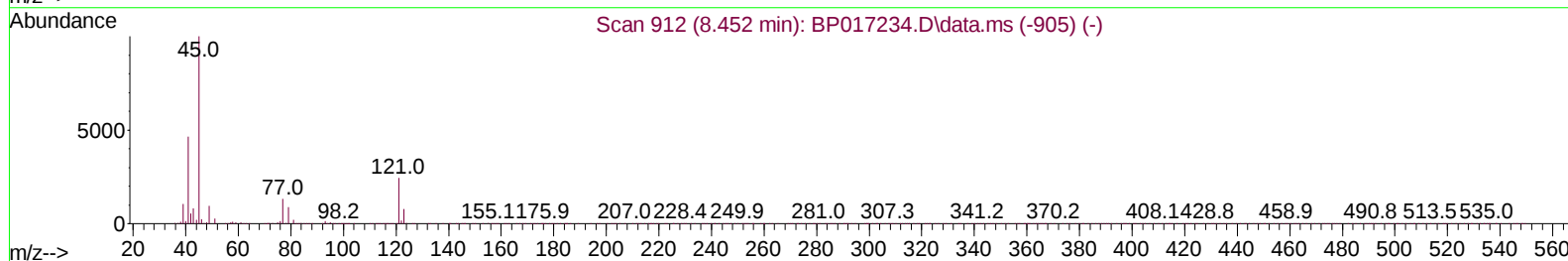
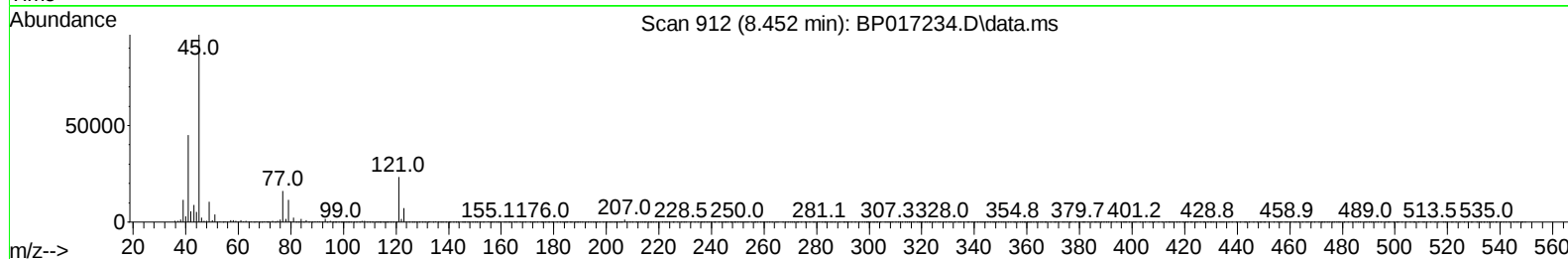
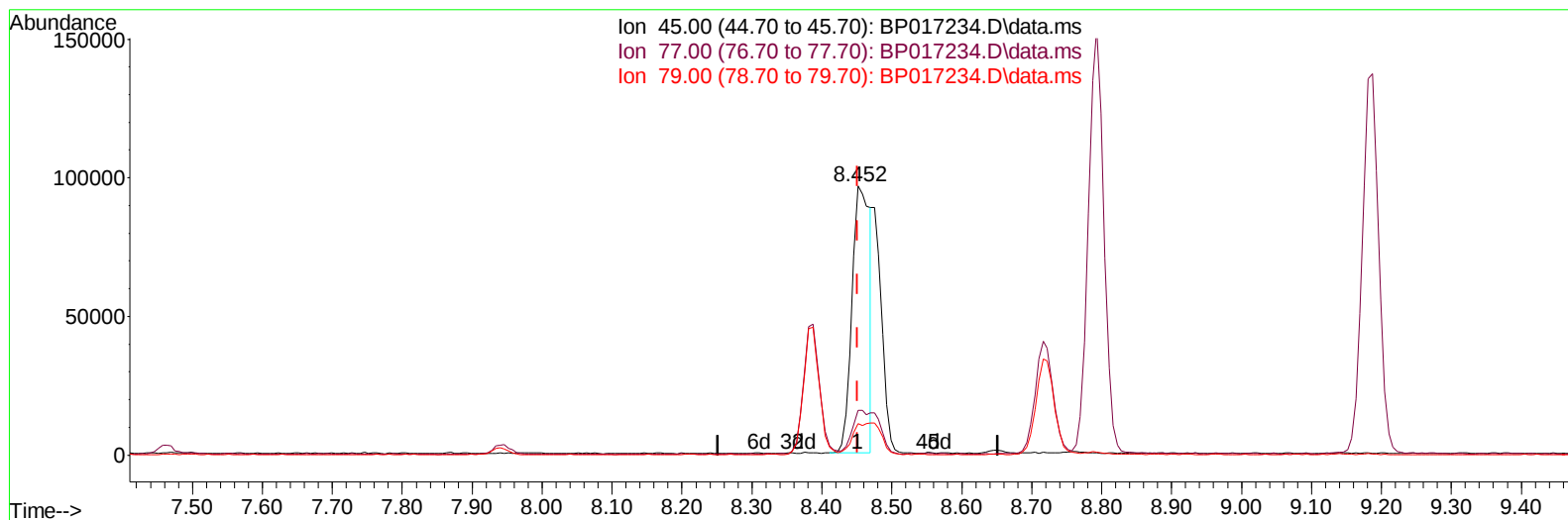
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092023\  
 Data File : BP017234.D  
 Acq On : 20 Sep 2023 12:17  
 Operator : MA/JU  
 Sample : SSTD02057  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD020619

Manual IntegrationsAPPROVED

Quant Time: Sep 20 13:29:28 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 20 13:27:02 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/21/2023  
 Supervised By :mohammad ahmed 09/21/2023



TIC: BP017234.D\data.ms

(14) 2,2'-oxybis(1-Chloropropane)

8.452min ( 0.000) 12.96 ng/u1

response 173641

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 45.00 | 100.00 | 100.00 |
| 77.00 | 16.60  | 16.62  |
| 79.00 | 11.80  | 11.79  |
| 0.00  | 0.00   | 0.00   |

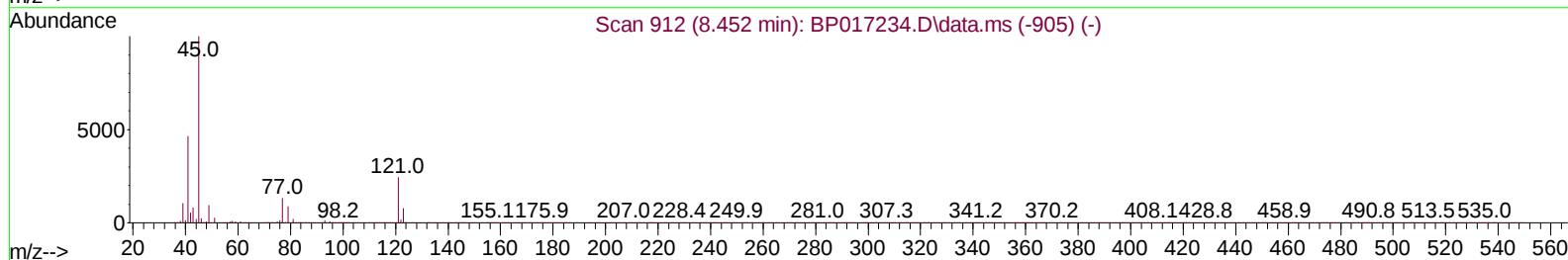
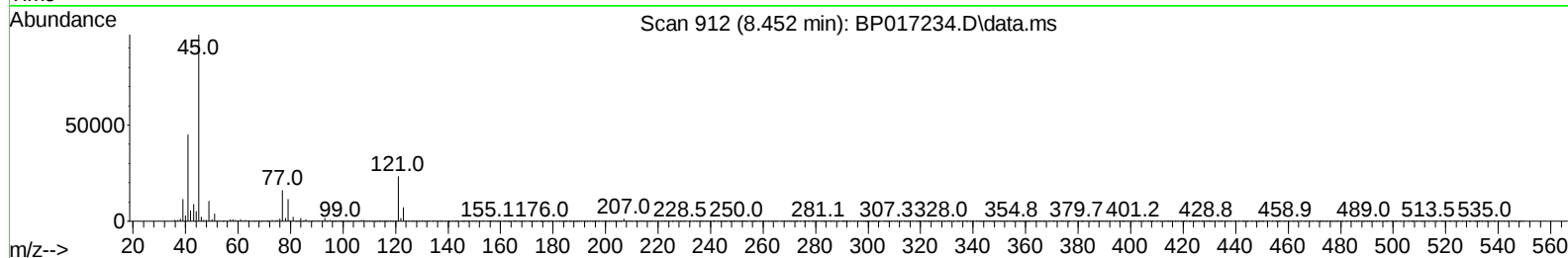
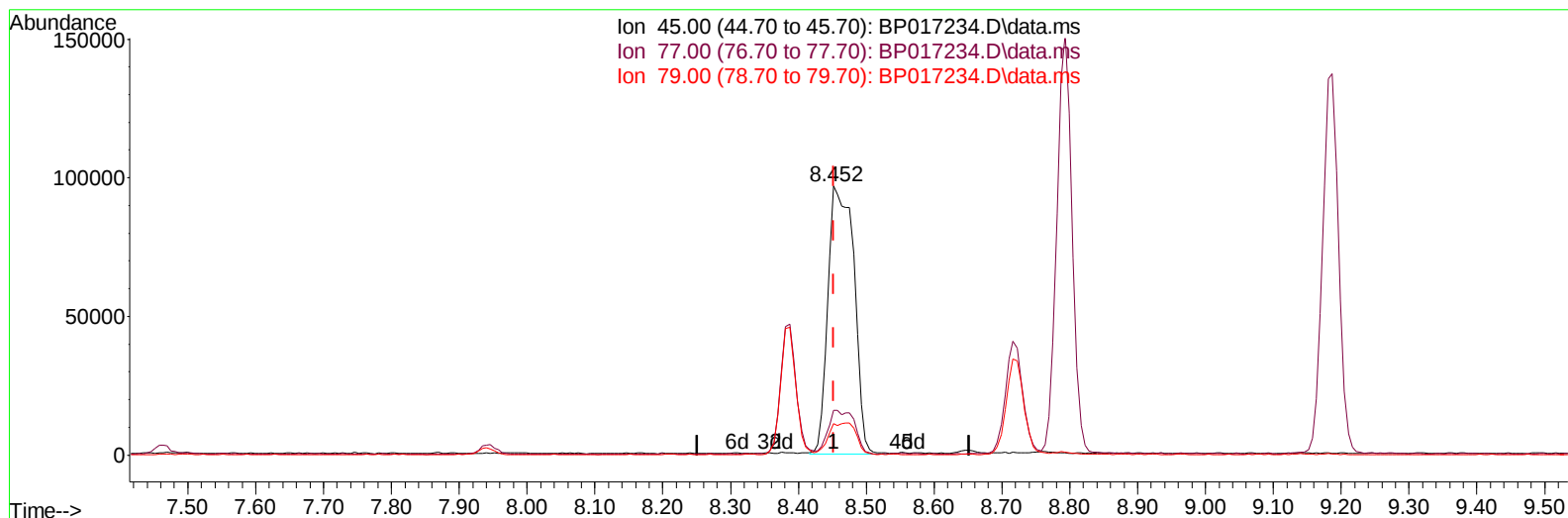
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092023\  
 Data File : BP017234.D  
 Acq On : 20 Sep 2023 12:17  
 Operator : MA/JU  
 Sample : SSTD02057  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD020619

Manual IntegrationsAPPROVED

Quant Time: Sep 20 13:29:28 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 20 13:27:02 2023  
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Reviewed By :Yogesh Patel 09/21/2023  
 Supervised By :mohammad ahmed 09/21/2023



TIC: BP017234.D\data.ms

(14) 2,2'-oxybis(1-Chloropropane)

8.452min ( 0.000) 19.09 ng/ul m

response 255854

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 45.00 | 100.00 | 100.00 |
| 77.00 | 16.60  | 16.62  |
| 79.00 | 11.80  | 11.79  |
| 0.00  | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092023\  
 Data File : BP017234.D  
 Acq On : 20 Sep 2023 12:17  
 Operator : MA/JU  
 Sample : SST002057  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SST0020619

Manual IntegrationsAPPROVED

Quant Time: Sep 20 13:31:39 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 20 13:27:02 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/21/2023  
 Supervised By :mohammad ahmed 09/21/2023

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.940  | 152  | 145105   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.769 | 136  | 586694   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.604 | 164  | 343314   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 17.381 | 188  | 739181   | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.486 | 240  | 716619   | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 24.016 | 264  | 820312   | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.305  | 96   | 31332    | 8.712  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.740  | 84   | 207294   | 20.507 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 7.093  | 99   | 247280   | 20.503 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.281  | 67   | 152296   | 20.337 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.464  | 132  | 197132   | 20.559 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.652  | 113  | 190320   | 19.330 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 9.140  | 128  | 88660    | 21.051 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.858  | 143  | 96861    | 21.459 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.381 | 165  | 183610   | 21.109 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.934 | 131  | 265381   | 20.375 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.022 | 166  | 515384   | 20.612 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 14.304 | 160  | 607181   | 21.280 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.828 | 143  | 87062    | 18.353 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.604 | 176  | 451710   | 20.834 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.734 | 200  | 80780    | 19.215 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 17.481 | 188  | 688286   | 21.268 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.722 | 212  | 808250   | 21.091 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.845 | 264  | 826854   | 20.859 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.340  | 88   | 29111    | 7.774  | ng/uL  | 100      |
| 5) Pyridine                   | 3.758  | 79   | 217583   | 20.776 | ng/ul  | 100      |
| 6) Benzaldehyde               | 7.099  | 77   | 144777   | 22.128 | ng/ul  | 100      |
| 8) Phenol                     | 7.122  | 94   | 247331   | 19.659 | ng/ul  | 100      |
| 10) Bis(2-Chloroethyl)ether   | 7.381  | 93   | 203849   | 20.240 | ng/ul  | 100      |
| 12) 2-Chlorophenol            | 7.493  | 128  | 201236   | 20.661 | ng/ul  | 100      |
| 13) 2-Methylphenol            | 8.387  | 108  | 182516   | 19.480 | ng/ul  | 100      |
| 14) 2,2'-oxybis(1-Chloropr... | 8.452  | 45   | 255854m  | 19.089 | ng/ul  |          |
| 16) Acetophenone              | 8.793  | 105  | 302273   | 19.704 | ng/ul  | 100      |
| 17) N-Nitroso-di-n-propyla... | 8.758  | 70   | 148443   | 19.009 | ng/ul  | 100      |
| 18) 4-Methylphenol            | 8.716  | 108  | 195992   | 18.998 | ng/ul  | 100      |
| 19) Hexachloroethane          | 9.005  | 117  | 82318    | 20.514 | ng/ul  | 100      |
| 22) Nitrobenzene              | 9.187  | 77   | 226554   | 20.742 | ng/ul  | 100      |
| 23) Isophorone                | 9.699  | 82   | 404658   | 19.984 | ng/ul  | 100      |
| 25) 2-Nitrophenol             | 9.887  | 139  | 105556   | 21.137 | ng/ul  | 100      |
| 26) 2,4-Dimethylphenol        | 9.928  | 107  | 208953   | 20.537 | ng/ul  | 100      |
| 27) Bis(2-Chloroethoxy)met... | 10.187 | 93   | 261877   | 20.422 | ng/ul  | 100      |
| 29) 2,4-Dichlorophenol        | 10.410 | 162  | 181779   | 20.658 | ng/ul  | 100      |
| 30) Naphthalene               | 10.816 | 128  | 638775   | 20.932 | ng/ul  | 100      |
| 32) 4-Chloroaniline           | 10.957 | 127  | 253354   | 20.050 | ng/ul  | 100      |
| 33) Hexachlorobutadiene       | 11.046 | 225  | 125950   | 20.984 | ng/ul  | 100      |
| 34) Caprolactam               | 11.775 | 113  | 50423    | 18.871 | ng/ul  | 100      |
| 35) 4-Chloro-3-methylphenol   | 12.057 | 107  | 182970   | 19.610 | ng/ul  | 100      |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092023\  
 Data File : BP017234.D  
 Acq On : 20 Sep 2023 12:17  
 Operator : MA/JU  
 Sample : SST002057  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SST0020619

Manual IntegrationsAPPROVED

Quant Time: Sep 20 13:31:39 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP092023.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 20 13:27:02 2023  
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/21/2023  
 Supervised By :mohammad ahmed 09/21/2023

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 36) 2-Methylnaphthalene       | 12.428 | 142  | 417736   | 20.467 | ng/ul | 100      |
| 37) 1-Methylnaphthalene       | 12.646 | 142  | 427679   | 20.665 | ng/ul | 100      |
| 39) 1,2,4,5-Tetrachloroben... | 12.775 | 216  | 228913   | 21.548 | ng/ul | 100      |
| 40) Hexachlorocyclopentadiene | 12.722 | 237  | 113295   | 18.652 | ng/ul | 100      |
| 41) 2,4,6-Trichlorophenol     | 13.034 | 196  | 140771   | 21.060 | ng/ul | 100      |
| 42) 2,4,5-Trichlorophenol     | 13.104 | 196  | 147325   | 20.782 | ng/ul | 100      |
| 43) 1,1'-Biphenyl             | 13.434 | 154  | 545491   | 21.412 | ng/ul | 100      |
| 44) 2-Chloronaphthalene       | 13.487 | 162  | 431856   | 21.311 | ng/ul | 100      |
| 45) 2-Nitroaniline            | 13.722 | 65   | 107022   | 20.060 | ng/ul | 100      |
| 47) Dimethylphthalate         | 14.069 | 163  | 525881   | 20.683 | ng/ul | 100      |
| 48) 2,6-Dinitrotoluene        | 14.210 | 165  | 91554    | 19.995 | ng/ul | 100      |
| 50) Acenaphthylene            | 14.334 | 152  | 705907   | 21.372 | ng/ul | 100      |
| 51) 3-Nitroaniline            | 14.551 | 138  | 97806    | 19.880 | ng/ul | 100      |
| 52) Acenaphthene              | 14.669 | 153  | 460682   | 21.083 | ng/ul | 100      |
| 53) 2,4-Dinitrophenol         | 14.757 | 184  | 48020    | 16.069 | ng/ul | 100      |
| 55) 4-Nitrophenol             | 14.840 | 109  | 68552    | 18.127 | ng/ul | 100      |
| 56) Dibenzofuran              | 15.004 | 168  | 634609   | 21.024 | ng/ul | 100      |
| 57) 2,4-Dinitrotoluene        | 14.998 | 165  | 138959   | 20.380 | ng/ul | 100      |
| 58) 2,3,4,6-Tetrachlorophenol | 15.222 | 232  | 126134   | 20.395 | ng/ul | 100      |
| 59) Diethylphthalate          | 15.422 | 149  | 508395   | 20.366 | ng/ul | 100      |
| 61) Fluorene                  | 15.657 | 166  | 516596   | 20.813 | ng/ul | 100      |
| 62) 4-Chlorophenyl-phenyle... | 15.651 | 204  | 262621   | 20.789 | ng/ul | 100      |
| 63) 4-Nitroaniline            | 15.722 | 138  | 92315    | 20.343 | ng/ul | 100      |
| 66) 4,6-Dinitro-2-methylph... | 15.751 | 198  | 88746    | 19.590 | ng/ul | 100      |
| 67) N-Nitrosodiphenylamine    | 15.875 | 169  | 419450   | 21.195 | ng/ul | 100      |
| 68) 4-Bromophenyl-phenylether | 16.557 | 248  | 152690   | 21.000 | ng/ul | 100      |
| 69) Hexachlorobenzene         | 16.634 | 284  | 178116   | 20.727 | ng/ul | 100      |
| 70) Atrazine                  | 16.834 | 200  | 147112   | 20.793 | ng/ul | 100      |
| 71) Pentachlorophenol         | 17.004 | 266  | 104382   | 18.874 | ng/ul | 100      |
| 72) Phenanthrene              | 17.422 | 178  | 805503   | 20.998 | ng/ul | 100      |
| 74) Anthracene                | 17.516 | 178  | 825365   | 21.253 | ng/ul | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 13.393 | 216  | 232234   | 21.829 | ng/uL | 100      |
| 76) Pentachlorobenzene        | 14.898 | 250  | 219143   | 22.498 | ng/uL | 100      |
| 77) Carbazole                 | 17.798 | 167  | 737355   | 21.571 | ng/ul | 100      |
| 78) Di-n-butylphthalate       | 18.310 | 149  | 791859   | 20.684 | ng/ul | 100      |
| 80) Fluoranthene              | 19.392 | 202  | 930744   | 20.834 | ng/ul | 100      |
| 82) Pyrene                    | 19.751 | 202  | 1005211  | 21.183 | ng/ul | 100      |
| 83) Butylbenzylphthalate      | 20.610 | 149  | 347106   | 20.749 | ng/ul | 100      |
| 84) 3,3'-Dichlorobenzidine    | 21.410 | 252  | 318206   | 20.915 | ng/ul | 100      |
| 85) Benzo(a)anthracene        | 21.469 | 228  | 1023049  | 21.175 | ng/ul | 100      |
| 86) Bis(2-ethylhexyl)phtha... | 21.345 | 149  | 498950   | 21.059 | ng/ul | 100      |
| 87) Chrysene                  | 21.527 | 228  | 960793   | 21.330 | ng/ul | 100      |
| 89) Di-n-octyl phthalate      | 22.316 | 149  | 878727   | 19.324 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 23.227 | 252  | 996846   | 20.332 | ng/ul | 100      |
| 91) Benzo(k)fluoranthene      | 23.280 | 252  | 1029210  | 20.917 | ng/ul | 100      |
| 93) Benzo(a)pyrene            | 23.898 | 252  | 974512   | 21.141 | ng/ul | 100      |
| 94) Indeno(1,2,3-cd)pyrene    | 26.692 | 276  | 1219767  | 22.974 | ng/ul | 100      |
| 95) Dibenzo(a,h)anthracene    | 26.715 | 278  | 1015074  | 22.813 | ng/ul | 100      |
| 96) Benzo(g,h,i)perylene      | 27.533 | 276  | 985030   | 23.750 | ng/ul | 100      |

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

**Instrument :**

BNA\_P

**ClientSampleId :**

SSTD020619

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 09/21/2023

Supervised By :mohammad ahmed 09/21/2023

**Instrument :**  
BNA\_P  
**ClientSampleId :**  
SSTD020619

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