

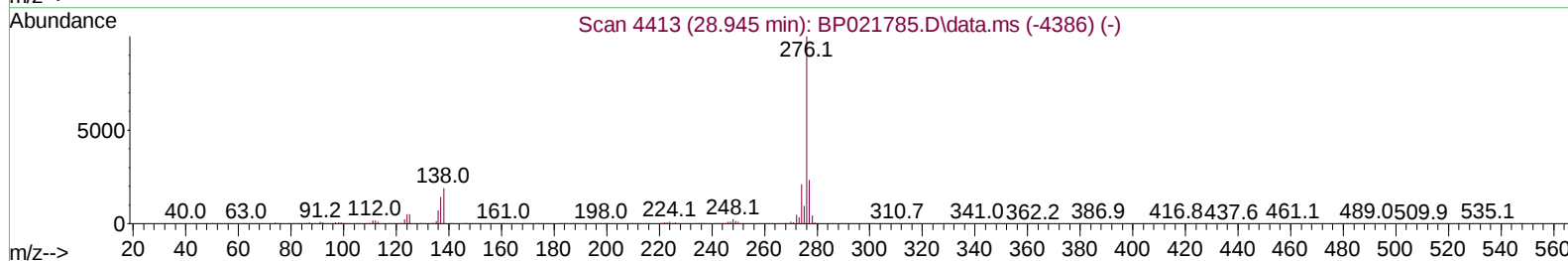
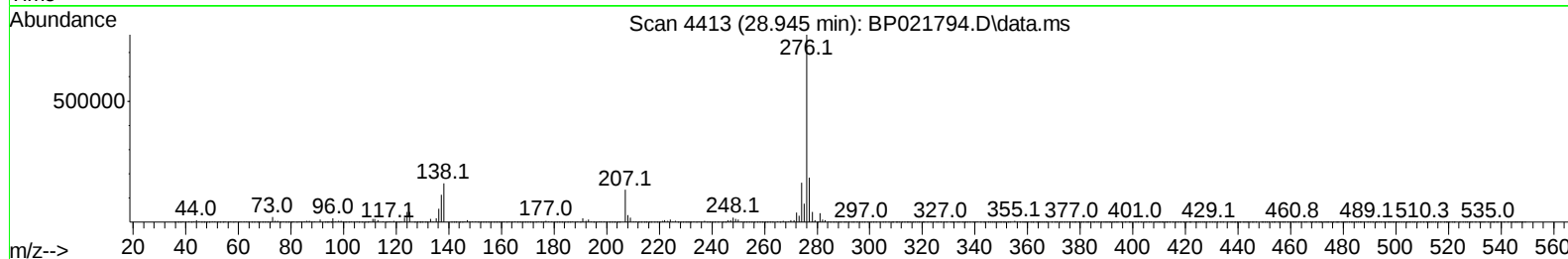
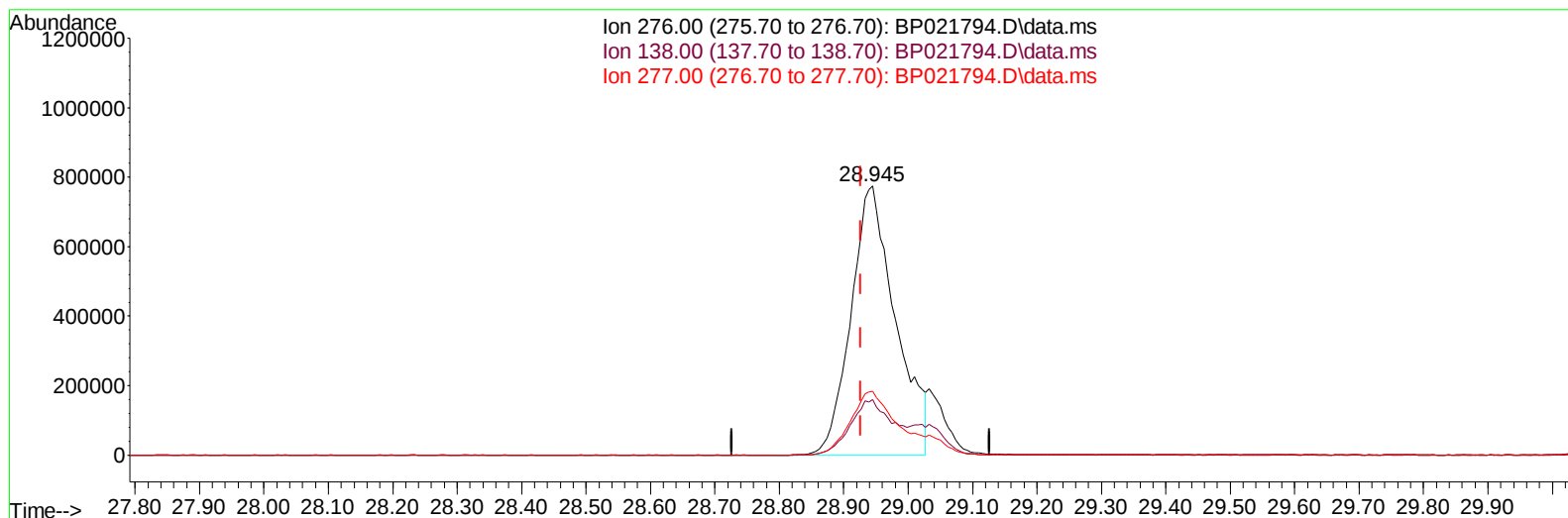
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP090324\  
Data File : BP021794.D  
Acq On : 03 Sep 2024 19:13  
Operator : RC/JU  
Sample : P3722-02MS  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCYD5MS

Manual IntegrationsAPPROVED

Quant Time: Sep 04 04:01:34 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP082324.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Aug 29 23:56:59 2024  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/04/2024  
Supervised By :mohammad ahmed 09/07/2024



TIC: BP021794.D\data.ms

(94) Indeno(1,2,3-cd)pyrene

28.945min (+ 0.018) 35.16 ng/u1

response 3710216

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 18.60  | 20.69  |
| 277.00 | 23.80  | 23.82  |
| 0.00   | 0.00   | 0.00   |

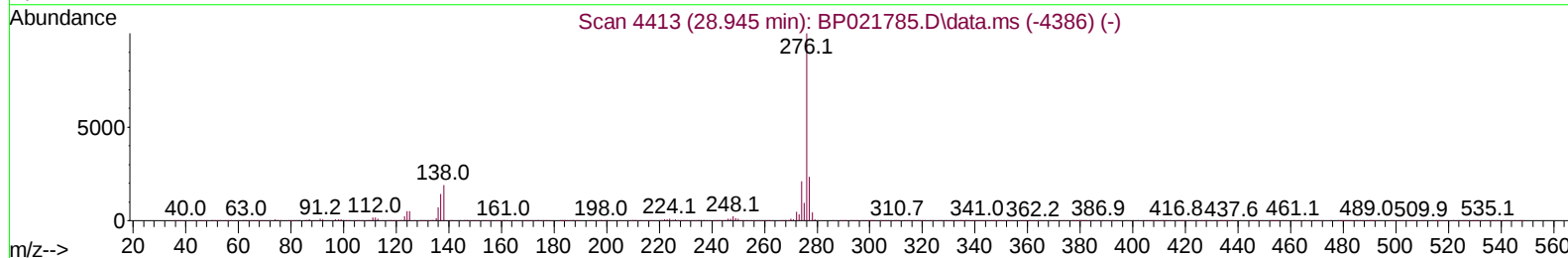
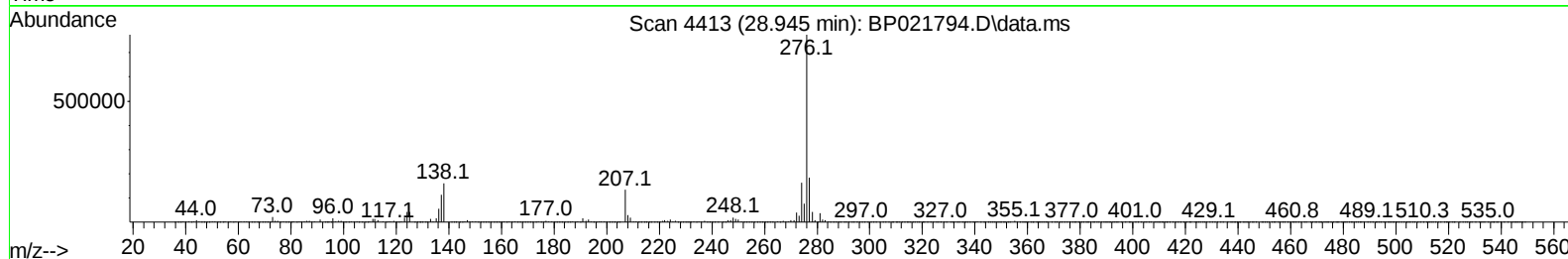
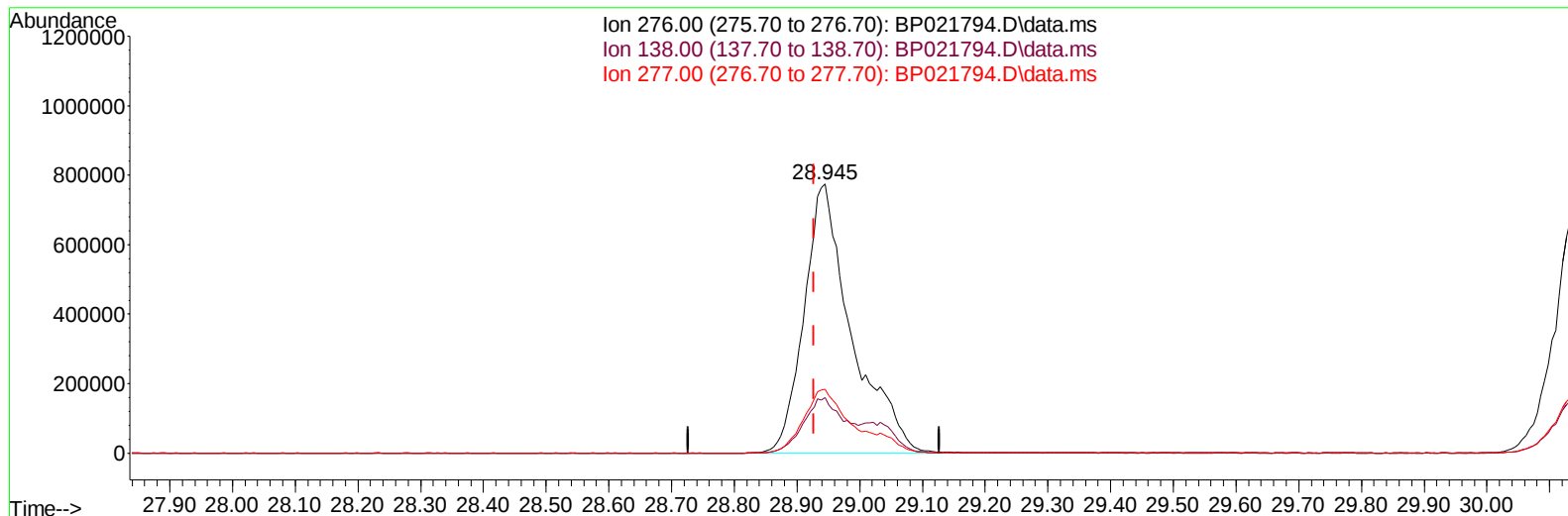
Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP090324\  
 Data File : BP021794.D  
 Acq On : 03 Sep 2024 19:13  
 Operator : RC/JU  
 Sample : P3722-02MS  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DCYD5MS

Manual IntegrationsAPPROVED

Quant Time: Sep 04 04:01:34 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP082324.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 29 23:56:59 2024  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/04/2024  
 Supervised By :mohammad ahmed 09/07/2024



TIC: BP021794.D\data.ms

**(94) Indeno(1,2,3-cd)pyrene**

28.945min (+ 0.018) 38.67 ng/ul m

response 4080579

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 18.60  | 20.69  |
| 277.00 | 23.80  | 23.82  |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP090324\  
 Data File : BP021794.D  
 Acq On : 03 Sep 2024 19:13  
 Operator : RC/JU  
 Sample : P3722-02MS  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DCYD5MS

Manual IntegrationsAPPROVED

Quant Time: Sep 04 04:48:01 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP082324.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 29 23:56:59 2024  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/04/2024  
 Supervised By :mohammad ahmed 09/07/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.775  | 152  | 217530   | 20.000 | ng/ul  | -0.01    |
| 20) Naphthalene-d8            | 10.569 | 136  | 876792   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.422 | 164  | 545423   | 20.000 | ng/ul  | -0.01    |
| 64) Phenanthrene-d10          | 17.228 | 188  | 1198643  | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.680 | 240  | 1227059  | 20.000 | ng/ul  | 0.01     |
| 88) Perylene-d12              | 25.080 | 264  | 1419838  | 20.000 | ng/ul  | 0.02     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.276  | 96   | 32413    | 4.947  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.687  | 84   | 479250   | 25.288 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 6.952  | 99   | 650512   | 28.021 | ng/ul  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.111  | 67   | 354522   | 28.062 | ng/ul  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.317  | 132  | 431695   | 28.803 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.481  | 113  | 511090   | 28.491 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 8.928  | 128  | 212652   | 30.146 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.646  | 143  | 213243   | 29.781 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.187 | 165  | 430308   | 30.496 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.693 | 131  | 571610   | 27.845 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.822 | 166  | 1303701  | 31.238 | ng/ul  | -0.02    |
| 49) Acenaphthylene-d8         | 14.110 | 160  | 1489598  | 31.791 | ng/ul  | -0.01    |
| 54) 4-Nitrophenol-d4          | 14.622 | 143  | 253301   | 31.902 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.434 | 176  | 1122720  | 31.998 | ng/ul  | -0.01    |
| 65) 4,6-Dinitro-2-methylph... | 15.546 | 200  | 186045   | 28.220 | ng/ul  | -0.01    |
| 73) Anthracene-d10            | 17.328 | 188  | 1788278  | 31.436 | ng/ul  | -0.01    |
| 81) Pyrene-d10                | 19.698 | 212  | 2286960  | 32.604 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.851 | 264  | 2582254  | 34.107 | ng/ul  | 0.01     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.305  | 88   | 93658    | 13.450 | ng/uL  | 97       |
| 5) Pyridine                   | 3.705  | 79   | 634856   | 33.523 | ng/ul  | 98       |
| 6) Benzaldehyde               | 6.922  | 77   | 461249   | 38.845 | ng/ul  | 98       |
| 8) Phenol                     | 6.975  | 94   | 841190   | 34.751 | ng/ul  | 97       |
| 10) Bis(2-Chloroethyl)ether   | 7.199  | 93   | 651546   | 34.628 | ng/ul  | 99       |
| 12) 2-Chlorophenol            | 7.346  | 128  | 569967   | 36.232 | ng/ul  | 98       |
| 13) 2-Methylphenol            | 8.216  | 108  | 600410   | 34.383 | ng/ul  | 100      |
| 14) 2,2'-oxybis(1-Chloropr... | 8.293  | 45   | 631833   | 34.791 | ng/ul  | 98       |
| 16) Acetophenone              | 8.593  | 105  | 1001911  | 34.838 | ng/ul  | 99       |
| 17) N-Nitroso-di-n-propyla... | 8.581  | 70   | 456695   | 33.650 | ng/ul  | 97       |
| 18) 4-Methylphenol            | 8.546  | 108  | 668281   | 34.956 | ng/ul  | 99       |
| 19) Hexachloroethane          | 8.858  | 117  | 252530   | 35.902 | ng/ul  | 96       |
| 22) Nitrobenzene              | 8.969  | 77   | 731365   | 37.182 | ng/ul  | 98       |
| 23) Isophorone                | 9.499  | 82   | 1342072  | 34.001 | ng/ul  | 99       |
| 25) 2-Nitrophenol             | 9.675  | 139  | 298019   | 37.565 | ng/ul  | 97       |
| 26) 2,4-Dimethylphenol        | 9.740  | 107  | 670719   | 34.446 | ng/ul  | 98       |
| 27) Bis(2-Chloroethoxy)met... | 9.975  | 93   | 860706   | 34.770 | ng/ul  | 98       |
| 29) 2,4-Dichlorophenol        | 10.210 | 162  | 520042   | 36.677 | ng/ul  | 97       |
| 30) Naphthalene               | 10.616 | 128  | 1781846  | 36.058 | ng/ul  | 100      |
| 32) 4-Chloroaniline           | 10.716 | 127  | 681724   | 32.894 | ng/ul  | 99       |
| 33) Hexachlorobutadiene       | 10.905 | 225  | 337463   | 35.472 | ng/ul  | 98       |
| 34) Caprolactam               | 11.499 | 113  | 220494   | 36.774 | ng/ul  | 97       |
| 35) 4-Chloro-3-methylphenol   | 11.852 | 107  | 694423   | 37.003 | ng/ul  | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP090324\  
 Data File : BP021794.D  
 Acq On : 03 Sep 2024 19:13  
 Operator : RC/JU  
 Sample : P3722-02MS  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DCYD5MS

Manual IntegrationsAPPROVED

Quant Time: Sep 04 04:48:01 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP082324.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 29 23:56:59 2024  
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/04/2024  
 Supervised By :mohammad ahmed 09/07/2024

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| 36) 2-Methylnaphthalene       | 12.228 | 142  | 1189668  | 35.593 | ng/ul | 99       |
| 37) 1-Methylnaphthalene       | 12.446 | 142  | 1190928  | 35.424 | ng/ul | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 12.599 | 216  | 628679   | 36.571 | ng/ul | 99       |
| 40) Hexachlorocyclopentadiene | 12.581 | 237  | 304760   | 30.896 | ng/ul | 99       |
| 41) 2,4,6-Trichlorophenol     | 12.846 | 196  | 418575   | 37.818 | ng/ul | 99       |
| 42) 2,4,5-Trichlorophenol     | 12.910 | 196  | 453167   | 37.918 | ng/ul | 98       |
| 43) 1,1'-Biphenyl             | 13.240 | 154  | 1542238  | 37.374 | ng/ul | 99       |
| 44) 2-Chloronaphthalene       | 13.281 | 162  | 1198579  | 37.045 | ng/ul | 99       |
| 45) 2-Nitroaniline            | 13.487 | 65   | 406057   | 40.061 | ng/ul | 97       |
| 47) Dimethylphthalate         | 13.875 | 163  | 1512008  | 36.182 | ng/ul | 99       |
| 48) 2,6-Dinitrotoluene        | 13.987 | 165  | 298753   | 38.764 | ng/ul | 97       |
| 50) Acenaphthylene            | 14.140 | 152  | 1886266  | 37.431 | ng/ul | 100      |
| 51) 3-Nitroaniline            | 14.316 | 138  | 319446   | 37.930 | ng/ul | 94       |
| 52) Acenaphthene              | 14.487 | 153  | 1338805  | 37.778 | ng/ul | 99       |
| 53) 2,4-Dinitrophenol         | 14.528 | 184  | 154088   | 33.396 | ng/ul | 95       |
| 55) 4-Nitrophenol             | 14.640 | 109  | 303266   | 41.714 | ng/ul | 92       |
| 56) Dibenzofuran              | 14.828 | 168  | 1825559  | 37.235 | ng/ul | 98       |
| 57) 2,4-Dinitrotoluene        | 14.787 | 165  | 451341   | 38.971 | ng/ul | 98       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.057 | 232  | 376131   | 36.529 | ng/ul | 93       |
| 59) Diethylphthalate          | 15.263 | 149  | 1549775  | 36.931 | ng/ul | 99       |
| 61) Fluorene                  | 15.487 | 166  | 1467641  | 37.009 | ng/ul | 98       |
| 62) 4-Chlorophenyl-phenyle... | 15.487 | 204  | 718024   | 36.034 | ng/ul | 99       |
| 63) 4-Nitroaniline            | 15.498 | 138  | 333865   | 41.017 | ng/ul | 93       |
| 66) 4,6-Dinitro-2-methylph... | 15.563 | 198  | 251203   | 34.064 | ng/ul | 98       |
| 67) N-Nitrosodiphenylamine    | 15.704 | 169  | 1278226  | 36.250 | ng/ul | 99       |
| 68) 4-Bromophenyl-phenylether | 16.398 | 248  | 466966   | 35.290 | ng/ul | 96       |
| 69) Hexachlorobenzene         | 16.522 | 284  | 544048   | 34.633 | ng/ul | 98       |
| 70) Atrazine                  | 16.681 | 200  | 481577   | 36.004 | ng/ul | 99       |
| 71) Pentachlorophenol         | 16.875 | 266  | 344542   | 34.270 | ng/ul | 98       |
| 72) Phenanthrene              | 17.269 | 178  | 2435370  | 37.008 | ng/ul | 99       |
| 74) Anthracene                | 17.369 | 178  | 2413384  | 36.200 | ng/ul | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.210 | 216  | 624226   | 35.028 | ng/uL | 97       |
| 76) Pentachlorobenzene        | 14.746 | 250  | 623466   | 33.754 | ng/uL | 98       |
| 77) Carbazole                 | 17.645 | 167  | 2364139  | 39.428 | ng/ul | 100      |
| 78) Di-n-butylphthalate       | 18.245 | 149  | 2749812  | 37.676 | ng/ul | 100      |
| 80) Fluoranthene              | 19.357 | 202  | 3033155  | 37.056 | ng/ul | 99       |
| 82) Pyrene                    | 19.733 | 202  | 3251515  | 37.507 | ng/ul | 100      |
| 83) Butylbenzylphthalate      | 20.681 | 149  | 1331121  | 37.502 | ng/ul | 98       |
| 84) 3,3'-Dichlorobenzidine    | 21.580 | 252  | 928099   | 30.013 | ng/ul | 99       |
| 85) Benzo(a)anthracene        | 21.663 | 228  | 3340731  | 38.273 | ng/ul | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.604 | 149  | 1890553  | 36.087 | ng/ul | 99       |
| 87) Chrysene                  | 21.733 | 228  | 3092115  | 37.641 | ng/ul | 99       |
| 89) Di-n-octyl phthalate      | 22.898 | 149  | 3326298  | 37.011 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene      | 24.004 | 252  | 3467998  | 39.036 | ng/ul | 98       |
| 91) Benzo(k)fluoranthene      | 24.080 | 252  | 3378316  | 38.843 | ng/ul | 98       |
| 93) Benzo(a)pyrene            | 24.921 | 252  | 3190805  | 38.984 | ng/ul | 100      |
| 94) Indeno(1,2,3-cd)pyrene    | 28.945 | 276  | 4080579m | 38.666 | ng/ul |          |
| 95) Dibenzo(a,h)anthracene    | 29.033 | 278  | 3303626  | 38.956 | ng/ul | 100      |
| 96) Benzo(g,h,i)perylene      | 30.133 | 276  | 3229837  | 39.610 | ng/ul | 99       |

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

**Instrument :**

BNA\_P

**ClientSampleId :**

DCYD5MS

**Manual IntegrationsAPPROVED**

Reviewed By :Jagrut Upadhyay 09/04/2024  
Supervised By :mohammad ahmed 09/07/2024

