

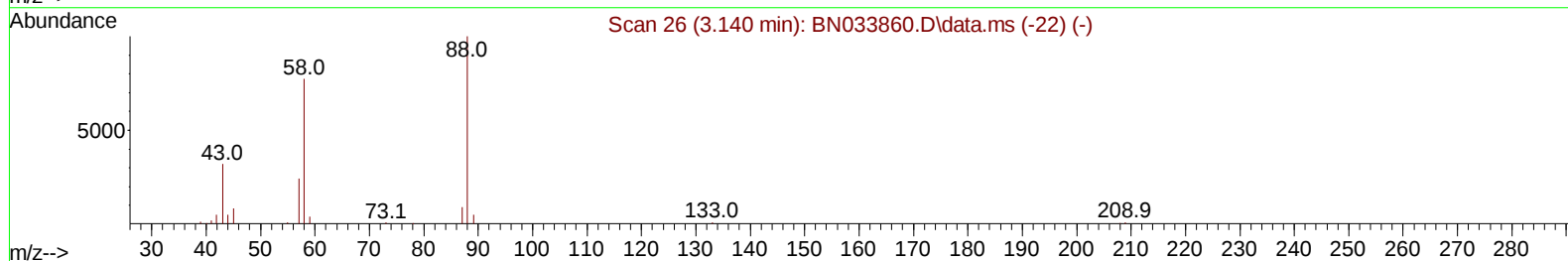
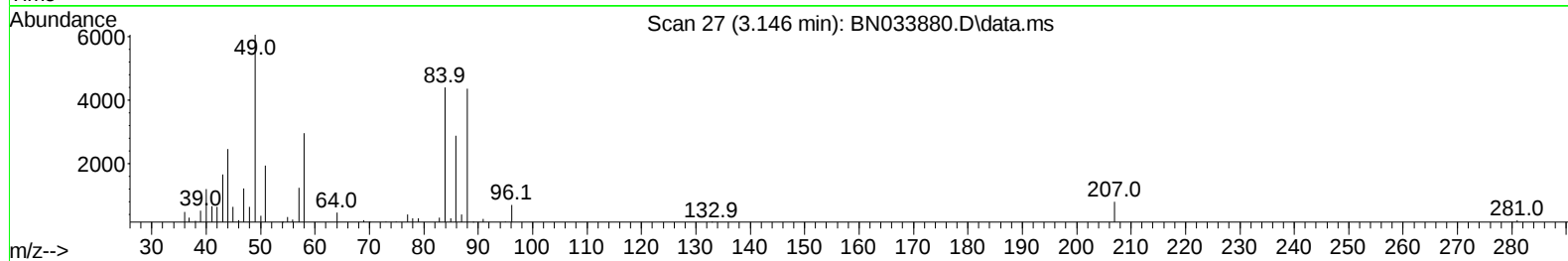
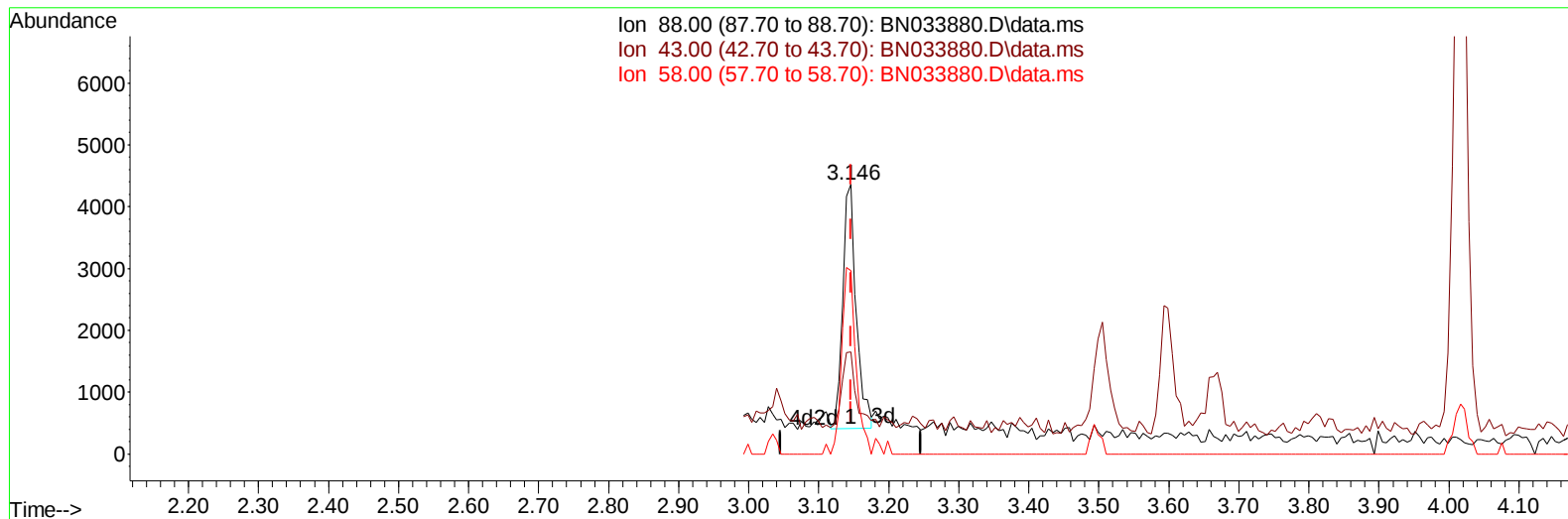
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091224\  
Data File : BN033880.D  
Acq On : 12 Sep 2024 13:17  
Operator : JU/RC  
Sample : P3391-02  
Misc : SFAM-MDL-Q3-WATER-02  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
MDL-WATER-QT3-2024-06

Manual IntegrationsAPPROVED

Quant Time: Sep 12 14:01:06 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Sep 11 01:58:59 2024  
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/13/2024  
Supervised By :mohammad ahmed 09/14/2024



TIC: BN033880.D\data.ms

(2) 1,4-Dioxane

3.146min (-0.000) 1.23 ng/uL

response 5334

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 88.00 | 100.00 | 100.00 |
| 43.00 | 30.90  | 37.92# |
| 58.00 | 80.70  | 68.02  |
| 0.00  | 0.00   | 0.00   |

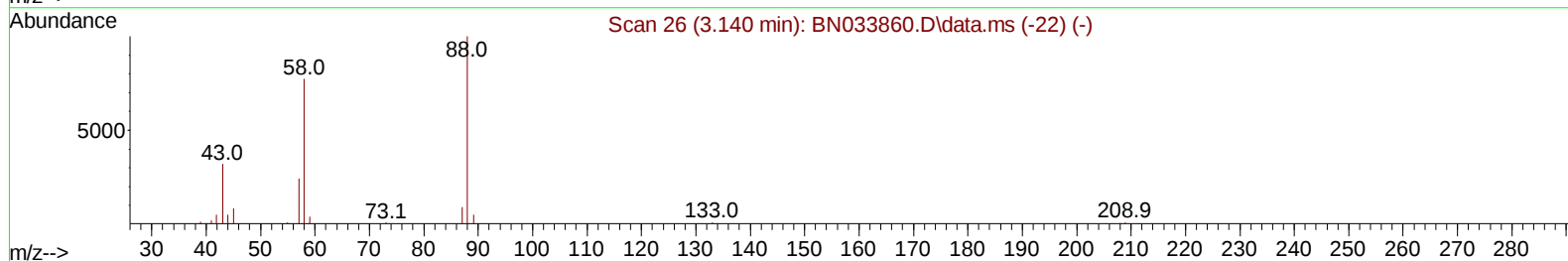
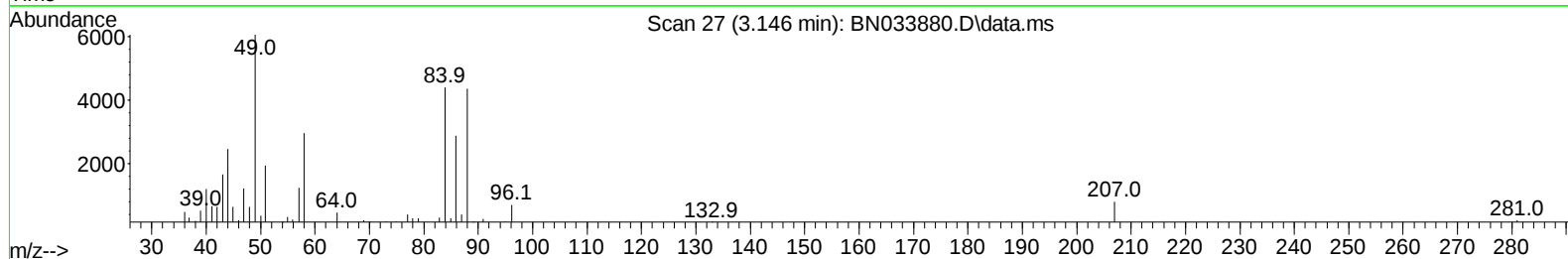
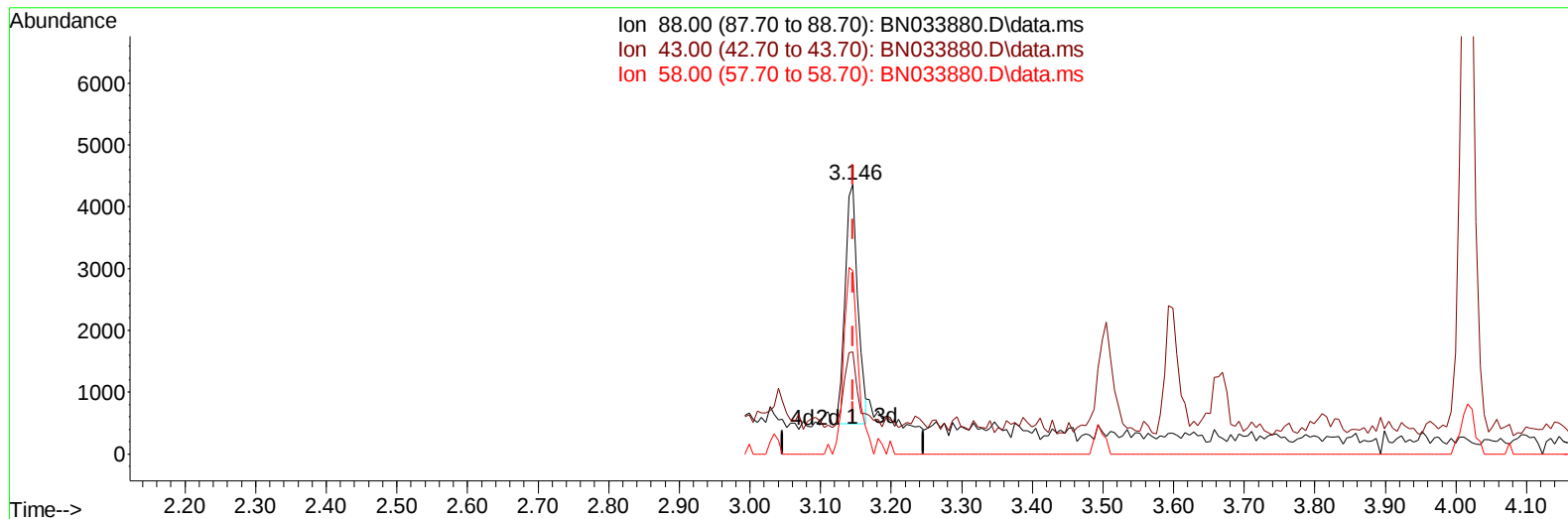
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091224\  
Data File : BN033880.D  
Acq On : 12 Sep 2024 13:17  
Operator : JU/RC  
Sample : P3391-02  
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Instrument :  
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ClientSampleId :  
MDL-WATER-QT3-2024-06

## Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 09/13/2024  
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Quant Time: Sep 12 14:01:06 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Sep 11 01:58:59 2024  
Response via : Initial Calibration



TIC: BN033880.D\data.ms

## (2) 1,4-Dioxane

3.146min (-0.000) 1.12 ng/uL m

response 4888

| Ion   | Exp%   | Act%   |
|-------|--------|--------|
| 88.00 | 100.00 | 100.00 |
| 43.00 | 30.90  | 37.92# |
| 58.00 | 80.70  | 68.02  |
| 0.00  | 0.00   | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN091224\  
 Data File : BN033880.D  
 Acq On : 12 Sep 2024 13:17  
 Operator : JU/RC  
 Sample : P3391-02  
 Misc : SFAM-MDL-Q3-WATER-02  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 MDL-WATER-QT3-2024-06

## Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 09/13/2024  
 Supervised By :mohammad ahmed 09/14/2024

Quant Time: Sep 12 14:04:17 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 11 01:58:59 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.493  | 152  | 168208   | 20.000 | ng/ul  | 0.00     |
| 20) Naphthalene-d8            | 10.251 | 136  | 722024   | 20.000 | ng/ul  | 0.00     |
| 38) Acenaphthene-d10          | 14.139 | 164  | 456605   | 20.000 | ng/ul  | 0.00     |
| 64) Phenanthrene-d10          | 16.892 | 188  | 1004415  | 20.000 | ng/ul  | 0.00     |
| 79) Chrysene-d12              | 21.133 | 240  | 1011696  | 20.000 | ng/ul  | 0.00     |
| 88) Perylene-d12              | 23.910 | 264  | 1070608  | 20.000 | ng/ul  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.111  | 96   | 20737    | 5.031  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.499  | 84   | 283260   | 23.119 | ng/ul  | 0.00     |
| 7) Phenol-d5                  | 6.681  | 99   | 338210   | 22.655 | ng/ul  | -0.01    |
| 9) Bis-(2-Chloroethyl)eth...  | 6.840  | 67   | 208186   | 22.477 | ng/ul  | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.034  | 132  | 270743   | 23.040 | ng/ul  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.204  | 113  | 245558   | 19.991 | ng/ul  | 0.00     |
| 21) Nitrobenzene-d5           | 8.634  | 128  | 135523   | 23.549 | ng/ul  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.346  | 143  | 143097   | 24.180 | ng/ul  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.881  | 165  | 243124   | 21.815 | ng/ul  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.393 | 131  | 338419   | 20.057 | ng/ul  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.563 | 166  | 793587   | 23.256 | ng/ul  | 0.00     |
| 49) Acenaphthylene-d8         | 13.828 | 160  | 883914   | 23.026 | ng/ul  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.357 | 143  | 148233   | 21.167 | ng/ul  | 0.00     |
| 60) Fluorene-d10              | 15.139 | 176  | 678808   | 23.472 | ng/ul  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.263 | 200  | 119855   | 20.014 | ng/ul  | 0.00     |
| 73) Anthracene-d10            | 16.992 | 188  | 914856   | 19.373 | ng/ul  | 0.00     |
| 81) Pyrene-d10                | 19.298 | 212  | 1244248  | 22.073 | ng/ul  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.721 | 264  | 1311092  | 23.840 | ng/ul  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.146  | 88   | 4888m    | 1.124  | ng/uL  |          |
| 5) Pyridine                   | 3.516  | 79   | 57127    | 4.556  | ng/ul# | 74       |
| 6) Benzaldehyde               | 6.646  | 77   | 43087    | 5.468  | ng/ul  | 97       |
| 8) Phenol                     | 6.710  | 94   | 68703    | 4.420  | ng/ul  | 98       |
| 10) Bis(2-Chloroethyl)ether   | 6.934  | 93   | 57308    | 4.744  | ng/ul  | 97       |
| 12) 2-Chlorophenol            | 7.063  | 128  | 56973    | 4.731  | ng/ul  | 97       |
| 13) 2-Methylphenol            | 7.940  | 108  | 46369    | 3.991  | ng/ul  | 100      |
| 14) 2,2'-oxybis(1-Chloropr... | 8.016  | 45   | 84182    | 4.725  | ng/ul  | 97       |
| 16) Acetophenone              | 8.304  | 105  | 88283    | 4.537  | ng/ul  | 96       |
| 17) N-Nitroso-di-n-propyla... | 8.299  | 70   | 44410    | 4.609  | ng/ul  | 98       |
| 18) 4-Methylphenol            | 8.263  | 108  | 52243    | 4.041  | ng/ul  | 99       |
| 19) Hexachloroethane          | 8.552  | 117  | 23926    | 4.711  | ng/ul  | 95       |
| 22) Nitrobenzene              | 8.675  | 77   | 68133    | 4.730  | ng/ul  | 98       |
| 23) Isophorone                | 9.199  | 82   | 117166   | 4.411  | ng/ul  | 98       |
| 25) 2-Nitrophenol             | 9.381  | 139  | 30192    | 4.559  | ng/ul  | 98       |
| 26) 2,4-Dimethylphenol        | 9.451  | 107  | 24919    | 1.829  | ng/ul  | 90       |
| 27) Bis(2-Chloroethoxy)met... | 9.687  | 93   | 74092    | 4.533  | ng/ul  | 99       |
| 29) 2,4-Dichlorophenol        | 9.910  | 162  | 51124    | 4.530  | ng/ul  | 95       |
| 30) Naphthalene               | 10.304 | 128  | 185047   | 4.685  | ng/ul  | 98       |
| 32) 4-Chloroaniline           | 10.416 | 127  | 70395    | 4.234  | ng/ul  | 95       |
| 33) Hexachlorobutadiene       | 10.593 | 225  | 35140    | 4.854  | ng/ul  | 96       |
| 34) Caprolactam               | 11.210 | 113  | 27888    | 6.207  | ng/ul  | 98       |
| 35) 4-Chloro-3-methylphenol   | 11.557 | 107  | 54367    | 4.092  | ng/ul  | 99       |

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 Sample : P3391-02  
 Misc : SFAM-MDL-Q3-WATER-02  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 MDL-WATER-QT3-2024-06

## Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 09/13/2024  
 Supervised By :mohammad ahmed 09/14/2024

Quant Time: Sep 12 14:04:17 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 11 01:58:59 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|-------|--------|----------|
| 36) 2-Methylnaphthalene       | 11.922 | 142  | 124746   | 4.586 | ng/ul  | 95       |
| 37) 1-Methylnaphthalene       | 12.145 | 142  | 126174   | 4.582 | ng/ul  | 98       |
| 39) 1,2,4,5-Tetrachloroben... | 12.304 | 216  | 63250    | 4.798 | ng/ul  | 98       |
| 40) Hexachlorocyclopentadiene | 12.287 | 237  | 37896    | 4.672 | ng/ul  | 99       |
| 41) 2,4,6-Trichlorophenol     | 12.551 | 196  | 38372    | 4.391 | ng/ul  | 98       |
| 42) 2,4,5-Trichlorophenol     | 12.622 | 196  | 40387    | 4.207 | ng/ul  | 98       |
| 43) 1,1'-Biphenyl             | 12.963 | 154  | 161582   | 4.694 | ng/ul  | 100      |
| 44) 2-Chloronaphthalene       | 12.998 | 162  | 127740   | 4.763 | ng/ul  | 98       |
| 45) 2-Nitroaniline            | 13.210 | 65   | 35512    | 4.098 | ng/ul  | 97       |
| 47) Dimethylphthalate         | 13.604 | 163  | 167302   | 4.811 | ng/ul  | 99       |
| 48) 2,6-Dinitrotoluene        | 13.722 | 165  | 29355    | 4.287 | ng/ul  | 94       |
| 50) Acenaphthylene            | 13.851 | 152  | 209446   | 4.986 | ng/ul  | 99       |
| 51) 3-Nitroaniline            | 14.045 | 138  | 29262    | 3.867 | ng/ul# | 94       |
| 52) Acenaphthene              | 14.204 | 153  | 139448   | 4.650 | ng/ul  | 98       |
| 53) 2,4-Dinitrophenol         | 14.257 | 184  | 27102    | 6.728 | ng/ul  | 88       |
| 55) 4-Nitrophenol             | 14.369 | 109  | 27419    | 4.610 | ng/ul# | 88       |
| 56) Dibenzofuran              | 14.539 | 168  | 192792   | 4.701 | ng/ul  | 97       |
| 57) 2,4-Dinitrotoluene        | 14.516 | 165  | 46496    | 4.504 | ng/ul  | 99       |
| 58) 2,3,4,6-Tetrachlorophenol | 14.775 | 232  | 36783    | 4.618 | ng/ul  | 99       |
| 59) Diethylphthalate          | 14.986 | 149  | 177358   | 4.880 | ng/ul  | 99       |
| 61) Fluorene                  | 15.192 | 166  | 163581   | 4.876 | ng/ul  | 95       |
| 62) 4-Chlorophenyl-phenyle... | 15.198 | 204  | 79300    | 4.877 | ng/ul  | 93       |
| 63) 4-Nitroaniline            | 15.216 | 138  | 30748    | 3.948 | ng/ul  | 97       |
| 66) 4,6-Dinitro-2-methylph... | 15.281 | 198  | 29401    | 4.421 | ng/ul# | 92       |
| 67) N-Nitrosodiphenylamine    | 15.410 | 169  | 133854   | 4.447 | ng/ul  | 98       |
| 68) 4-Bromophenyl-phenylether | 16.092 | 248  | 46986    | 4.607 | ng/ul  | 98       |
| 69) Hexachlorobenzene         | 16.198 | 284  | 54419    | 4.597 | ng/ul  | 96       |
| 70) Atrazine                  | 16.381 | 200  | 53389    | 4.896 | ng/ul  | 96       |
| 71) Pentachlorophenol         | 16.545 | 266  | 56469    | 7.162 | ng/ul  | 94       |
| 72) Phenanthrene              | 16.933 | 178  | 264659   | 4.779 | ng/ul  | 98       |
| 74) Anthracene                | 17.022 | 178  | 236269   | 4.178 | ng/ul  | 98       |
| 75) 1,2,3,4-Tetrachloroben... | 12.922 | 216  | 61234    | 4.380 | ng/uL  | 99       |
| 76) Pentachlorobenzene        | 14.457 | 250  | 65825    | 4.625 | ng/uL  | 97       |
| 77) Carbazole                 | 17.298 | 167  | 233483   | 4.486 | ng/ul  | 99       |
| 78) Di-n-butylphthalate       | 17.886 | 149  | 319442   | 4.828 | ng/ul  | 99       |
| 80) Fluoranthene              | 18.963 | 202  | 308081   | 4.675 | ng/ul  | 99       |
| 82) Pyrene                    | 19.327 | 202  | 328463   | 4.636 | ng/ul  | 96       |
| 83) Butylbenzylphthalate      | 20.263 | 149  | 142732   | 4.493 | ng/ul  | 99       |
| 84) 3,3'-Dichlorobenzidine    | 21.051 | 252  | 82767    | 3.642 | ng/ul  | 97       |
| 85) Benzo(a)anthracene        | 21.116 | 228  | 339326   | 4.704 | ng/ul  | 98       |
| 86) Bis(2-ethylhexyl)phtha... | 21.074 | 149  | 216011   | 4.645 | ng/ul  | 97       |
| 87) Chrysene                  | 21.168 | 228  | 316315   | 4.596 | ng/ul  | 98       |
| 89) Di-n-octyl phthalate      | 22.133 | 149  | 349326   | 4.904 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.033 | 252  | 321317   | 4.923 | ng/ul  | 97       |
| 91) Benzo(k)fluoranthene      | 23.092 | 252  | 321939   | 5.000 | ng/ul  | 95       |
| 93) Benzo(a)pyrene            | 23.774 | 252  | 307362   | 4.985 | ng/ul  | 97       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.986 | 276  | 374007   | 4.895 | ng/ul  | 98       |
| 95) Dibenzo(a,h)anthracene    | 27.045 | 278  | 309547   | 5.012 | ng/ul  | 97       |
| 96) Benzo(g,h,i)perylene      | 27.945 | 276  | 300418   | 5.065 | ng/ul  | 97       |

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

**Instrument :**  
BNA\_N  
**ClientSampleId :**  
MDL-WATER-QT3-2024-06

**Manual IntegrationsAPPROVED**

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