

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM051122\  
 Data File : BM035012.D  
 Acq On : 12 May 2022 07:56  
 Operator : CG/JU  
 Sample : N2707-06MSD  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 EXL87MSD

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 05/12/2022  
 Supervised By :Yogesh Patel 05/17/2022

Quant Time: May 12 08:33:45 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM051022.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 10 16:40:09 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.916  | 152  | 138773   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.722 | 136  | 652530   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.551 | 164  | 424936   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.292 | 188  | 838723   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.468 | 240  | 636667   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 23.850 | 264  | 595122   | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.363  | 96   | 19150    | 5.773  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.775  | 84   | 114415   | 12.514 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.081  | 99   | 106654   | 8.884  | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.245  | 67   | 344150   | 45.671 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.451  | 132  | 346612   | 36.741 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.628  | 113  | 226844   | 23.176 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.075  | 128  | 248246   | 49.637 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.798  | 143  | 264237   | 48.985 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.339 | 165  | 434417   | 43.032 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.851 | 131  | 609999   | 42.438 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.963 | 166  | 1757683  | 54.846 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.245 | 160  | 1991112  | 48.965 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.739 | 143  | 65460    | 11.955 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.545 | 176  | 1460040  | 54.002 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.662 | 200  | 292459   | 55.335 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.392 | 188  | 2209329  | 54.142 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.680 | 212  | 2205603  | 55.479 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.703 | 264  | 1788817  | 57.512 | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.399  | 88   | 20719    | 6.273  | ng/uL  | 94       |
| 5) Pyridine                   | 3.793  | 79   | 137212   | 14.220 | ng/u1# | 88       |
| 6) Benzaldehyde               | 7.057  | 77   | 398355m  | 59.455 | ng/u1  |          |
| 8) Phenol                     | 7.104  | 94   | 126438   | 10.580 | ng/u1  | 94       |
| 10) Bis(2-Chloroethyl)ether   | 7.339  | 93   | 455492   | 48.155 | ng/u1  | 96       |
| 12) 2-Chlorophenol            | 7.481  | 128  | 369816   | 38.438 | ng/u1  | 99       |
| 13) 2-Methylphenol            | 8.363  | 108  | 264194   | 29.167 | ng/u1  | 96       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.457  | 45   | 697691   | 45.116 | ng/u1  | 99       |
| 16) Acetophenone              | 8.739  | 105  | 779254   | 51.290 | ng/u1  | 98       |
| 17) N-Nitroso-di-n-propyla... | 8.728  | 70   | 420939   | 51.946 | ng/u1  | 95       |
| 18) 4-Methylphenol            | 8.692  | 108  | 249922   | 25.255 | ng/u1  | 100      |
| 19) Hexachloroethane          | 9.004  | 117  | 190680   | 46.719 | ng/u1  | 96       |
| 22) Nitrobenzene              | 9.116  | 77   | 603698   | 48.644 | ng/u1  | 97       |
| 23) Isophorone                | 9.651  | 82   | 1175774  | 51.562 | ng/u1  | 98       |
| 25) 2-Nitrophenol             | 9.833  | 139  | 291431   | 50.280 | ng/u1  | 89       |
| 26) 2,4-Dimethylphenol        | 9.892  | 107  | 401198   | 33.097 | ng/u1  | 99       |
| 27) Bis(2-Chloroethoxy)met... | 10.133 | 93   | 657384   | 46.508 | ng/u1  | 99       |
| 29) 2,4-Dichlorophenol        | 10.369 | 162  | 448844   | 45.701 | ng/u1  | 98       |
| 30) Naphthalene               | 10.769 | 128  | 1635839  | 47.717 | ng/u1  | 99       |
| 32) 4-Chloroaniline           | 10.875 | 127  | 639753   | 44.473 | ng/u1  | 99       |
| 33) Hexachlorobutadiene       | 11.069 | 225  | 294727   | 45.785 | ng/u1  | 98       |
| 34) Caprolactam               | 11.657 | 113  | 22881    | 7.791  | ng/u1  | 95       |
| 35) 4-Chloro-3-methylphenol   | 11.998 | 107  | 490558   | 46.464 | ng/u1  | 98       |
| 36) 2-Methylnaphthalene       | 12.380 | 142  | 1183690  | 49.939 | ng/u1  | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM051122\  
 Data File : BM035012.D  
 Acq On : 12 May 2022 07:56  
 Operator : CG/JU  
 Sample : N2707-06MSD  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 EXL87MSD

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 05/12/2022  
 Supervised By :Yogesh Patel 05/17/2022

Quant Time: May 12 08:33:45 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM051022.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 10 16:40:09 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 37) 1-Methylnaphthalene       | 12.598 | 142  | 1231307  | 51.373 | ng/ul  | 98       |
| 39) 1,2,4,5-Tetrachloroben... | 12.751 | 216  | 604408   | 48.361 | ng/ul  | 99       |
| 40) Hexachlorocyclopentadiene | 12.733 | 237  | 286719   | 33.078 | ng/ul  | 98       |
| 41) 2,4,6-Trichlorophenol     | 12.986 | 196  | 434840   | 54.652 | ng/ul  | 98       |
| 42) 2,4,5-Trichlorophenol     | 13.057 | 196  | 474261   | 55.640 | ng/ul  | 98       |
| 43) 1,1'-Biphenyl             | 13.392 | 154  | 1668115  | 49.411 | ng/ul  | 99       |
| 44) 2-Chloronaphthalene       | 13.427 | 162  | 1277684  | 49.696 | ng/ul  | 97       |
| 45) 2-Nitroaniline            | 13.627 | 65   | 458205   | 62.779 | ng/ul  | 92       |
| 47) Dimethylphthalate         | 14.010 | 163  | 1803378  | 56.956 | ng/ul  | 99       |
| 48) 2,6-Dinitrotoluene        | 14.127 | 165  | 388859   | 64.661 | ng/ul  | 93       |
| 50) Acenaphthylene            | 14.274 | 152  | 2182168  | 53.415 | ng/ul  | 98       |
| 51) 3-Nitroaniline            | 14.451 | 138  | 370494   | 67.280 | ng/ul  | 92       |
| 52) Acenaphthene              | 14.615 | 153  | 1451567  | 53.160 | ng/ul  | 99       |
| 53) 2,4-Dinitrophenol         | 14.657 | 184  | 209190   | 59.629 | ng/ul  | 94       |
| 55) 4-Nitrophenol             | 14.751 | 109  | 64200    | 13.333 | ng/ul  | 94       |
| 56) Dibenzofuran              | 14.951 | 168  | 2030439  | 53.992 | ng/ul  | 94       |
| 57) 2,4-Dinitrotoluene        | 14.910 | 165  | 549605   | 64.790 | ng/ul  | 90       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.174 | 232  | 407344   | 60.951 | ng/ul  | 98       |
| 59) Diethylphthalate          | 15.374 | 149  | 1862035  | 57.415 | ng/ul  | 99       |
| 61) Fluorene                  | 15.598 | 166  | 1727476  | 56.152 | ng/ul  | 99       |
| 62) 4-Chlorophenyl-phenyle... | 15.592 | 204  | 834844   | 55.869 | ng/ul  | 99       |
| 63) 4-Nitroaniline            | 15.615 | 138  | 388059   | 79.174 | ng/ul# | 83       |
| 66) 4,6-Dinitro-2-methylph... | 15.674 | 198  | 297043   | 56.763 | ng/ul# | 95       |
| 67) N-Nitrosodiphenylamine    | 15.804 | 169  | 1489733  | 55.170 | ng/ul  | 99       |
| 68) 4-Bromophenyl-phenylether | 16.486 | 248  | 493155   | 56.544 | ng/ul  | 97       |
| 69) Hexachlorobenzene         | 16.604 | 284  | 561323   | 56.879 | ng/ul  | 96       |
| 70) Atrazine                  | 16.757 | 200  | 552225   | 57.534 | ng/ul  | 100      |
| 71) Pentachlorophenol         | 16.945 | 266  | 344021   | 56.160 | ng/ul  | 97       |
| 72) Phenanthrene              | 17.333 | 178  | 2666838  | 56.529 | ng/ul  | 99       |
| 74) Anthracene                | 17.427 | 178  | 2694915  | 56.480 | ng/ul  | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.357 | 216  | 632260   | 47.108 | ng/ul  | 99       |
| 76) Pentachlorobenzene        | 14.874 | 250  | 641824   | 52.206 | ng/ul  | 99       |
| 77) Carbazole                 | 17.692 | 167  | 2408374  | 57.413 | ng/ul  | 99       |
| 78) Di-n-butylphthalate       | 18.262 | 149  | 2996694  | 57.607 | ng/ul  | 99       |
| 80) Fluoranthene              | 19.345 | 202  | 2813804  | 58.877 | ng/ul  | 97       |
| 82) Pyrene                    | 19.709 | 202  | 2822744  | 57.838 | ng/ul  | 99       |
| 83) Butylbenzylphthalate      | 20.609 | 149  | 1178295  | 60.351 | ng/ul  | 99       |
| 84) 3,3'-Dichlorobenzidine    | 21.386 | 252  | 759418   | 63.211 | ng/ul  | 99       |
| 85) Benzo(a)anthracene        | 21.456 | 228  | 2409493  | 58.249 | ng/ul  | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.392 | 149  | 1581750  | 54.675 | ng/ul  | 100      |
| 87) Chrysene                  | 21.509 | 228  | 2319500  | 57.582 | ng/ul  | 98       |
| 89) Di-n-octyl phthalate      | 22.309 | 149  | 2430179  | 49.052 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.127 | 252  | 2404758  | 60.156 | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene      | 23.180 | 252  | 2131190  | 55.988 | ng/ul  | 99       |
| 93) Benzo(a)pyrene            | 23.750 | 252  | 2262527  | 59.369 | ng/ul# | 96       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.332 | 276  | 2640408  | 60.567 | ng/ul# | 94       |
| 95) Dibenzo(a,h)anthracene    | 26.344 | 278  | 2246859  | 60.117 | ng/ul# | 97       |
| 96) Benzo(g,h,i)perylene      | 27.085 | 276  | 2215219  | 61.174 | ng/ul# | 95       |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM051122\  
 Data File : BM035012.D  
 Acq On : 12 May 2022 07:56  
 Operator : CG/JU  
 Sample : N2707-06MSD  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 EXL87MSD

Quant Time: May 12 08:33:45 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM051022.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 10 16:40:09 2022  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 05/12/2022  
 Supervised By :Yogesh Patel 05/17/2022

