

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM092922\  
 Data File : BM036800.D  
 Acq On : 29 Sep 2022 14:15  
 Operator : CG/JU  
 Sample : SSTD08082  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD080030

Quant Time: Sep 30 01:13:38 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM092922.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Sep 30 01:10:23 2022  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 09/30/2022  
 Supervised By :mohammad ahmed 10/03/2022

| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| -----                         |        |      |          |         |        |          |
| Internal Standards            |        |      |          |         |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.987  | 152  | 329832   | 20.000  | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.798 | 136  | 1499022  | 20.000  | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.616 | 164  | 930171   | 20.000  | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.351 | 188  | 1878720  | 20.000  | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.527 | 240  | 1959300  | 20.000  | ng/u1  | 0.01     |
| 88) Perylene-d12              | 23.944 | 264  | 1908589  | 20.000  | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |         |        |          |
| 3) 1,4-Dioxane-d8             | 3.369  | 96   | 283681   | 32.242  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.793  | 84   | 2310832  | 90.056  | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.151  | 99   | 2833827  | 96.142  | ng/u1  | 0.01     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.316  | 67   | 1688173  | 94.310  | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.516  | 132  | 2003663  | 94.420  | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.704  | 113  | 2113867  | 97.058  | ng/u1  | 0.01     |
| 21) Nitrobenzene-d5           | 9.157  | 128  | 1032722  | 101.785 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.881  | 143  | 1071290  | 117.878 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.416 | 165  | 1924716  | 91.385  | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.939 | 131  | 3019475  | 86.228  | ng/u1  | 0.01     |
| 46) Dimethylphthalate-d6      | 14.033 | 166  | 5423035  | 86.444  | ng/u1  | 0.01     |
| 49) Acenaphthylene-d8         | 14.310 | 160  | 6488203  | 87.190  | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.822 | 143  | 1184505  | 97.346  | ng/u1  | 0.02     |
| 60) Fluorene-d10              | 15.604 | 176  | 4765617  | 84.767  | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.727 | 200  | 931128   | 125.269 | ng/u1  | 0.01     |
| 73) Anthracene-d10            | 17.457 | 188  | 7348998  | 85.553  | ng/u1  | 0.01     |
| 81) Pyrene-d10                | 19.739 | 212  | 8350451  | 86.632  | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.797 | 264  | 8275640  | 88.751  | ng/u1  | 0.02     |
| Target Compounds              |        |      |          |         |        |          |
|                               |        |      |          |         | Qvalue |          |
| 2) 1,4-Dioxane                | 3.405  | 88   | 310002   | 32.708  | ng/uL  | 95       |
| 5) Pyridine                   | 3.816  | 79   | 2338220  | 90.699  | ng/u1  | 100      |
| 6) Benzaldehyde               | 7.122  | 77   | 1100801  | 77.123  | ng/u1  | 98       |
| 8) Phenol                     | 7.175  | 94   | 2994853  | 94.669  | ng/u1  | 99       |
| 10) Bis(2-Chloroethyl)ether   | 7.410  | 93   | 2254548  | 91.598  | ng/u1  | 98       |
| 12) 2-Chlorophenol            | 7.551  | 128  | 2125932  | 92.611  | ng/u1  | 99       |
| 13) 2-Methylphenol            | 8.434  | 108  | 2184954  | 93.966  | ng/u1  | 98       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.522  | 45   | 2930842  | 100.465 | ng/u1  | 99       |
| 16) Acetophenone              | 8.822  | 105  | 3442745  | 91.133  | ng/u1  | 99       |
| 17) N-Nitroso-di-n-propyla... | 8.816  | 70   | 1825810  | 100.164 | ng/u1  | 97       |
| 18) 4-Methylphenol            | 8.775  | 108  | 2384994  | 94.852  | ng/u1  | 97       |
| 19) Hexachloroethane          | 9.069  | 117  | 832576   | 89.544  | ng/u1  | 98       |
| 22) Nitrobenzene              | 9.198  | 77   | 2610354  | 96.900  | ng/u1  | 98       |
| 23) Isophorone                | 9.739  | 82   | 5164325  | 94.022  | ng/u1  | 99       |
| 25) 2-Nitrophenol             | 9.910  | 139  | 1203162  | 107.186 | ng/u1  | 96       |
| 26) 2,4-Dimethylphenol        | 9.969  | 107  | 2509145  | 90.630  | ng/u1  | 98       |
| 27) Bis(2-Chloroethoxy)met... | 10.210 | 93   | 2988997  | 91.299  | ng/u1  | 100      |
| 29) 2,4-Dichlorophenol        | 10.445 | 162  | 1984913  | 89.645  | ng/u1  | 100      |
| 30) Naphthalene               | 10.851 | 128  | 6817561  | 83.486  | ng/u1  | 99       |
| 32) 4-Chloroaniline           | 10.963 | 127  | 3109166  | 86.000  | ng/u1  | 100      |
| 33) Hexachlorobutadiene       | 11.128 | 225  | 1041674  | 76.921  | ng/u1  | 99       |
| 34) Caprolactam               | 11.775 | 113  | 683541m  | 95.169  | ng/u1  |          |
| 35) 4-Chloro-3-methylphenol   | 12.081 | 107  | 2325871  | 95.160  | ng/u1  | 99       |
| 36) 2-Methylnaphthalene       | 12.451 | 142  | 4691294  | 86.224  | ng/u1  | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM092922\  
 Data File : BM036800.D  
 Acq On : 29 Sep 2022 14:15  
 Operator : CG/JU  
 Sample : SSTD08082  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD080030

Quant Time: Sep 30 01:13:38 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM092922.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Sep 30 01:10:23 2022  
 Response via : Initial Calibration

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 09/30/2022  
 Supervised By :mohammad ahmed 10/03/2022

| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| 37) 1-Methylnaphthalene       | 12.669 | 142  | 4662083  | 85.235  | ng/u1  | 100      |
| 39) 1,2,4,5-Tetrachloroben... | 12.816 | 216  | 2164346  | 82.209  | ng/u1  | 99       |
| 40) Hexachlorocyclopentadiene | 12.792 | 237  | 1238960  | 90.300  | ng/u1  | 98       |
| 41) 2,4,6-Trichlorophenol     | 13.057 | 196  | 1488042  | 92.310  | ng/u1  | 99       |
| 42) 2,4,5-Trichlorophenol     | 13.128 | 196  | 1606430  | 92.949  | ng/u1  | 99       |
| 43) 1,1'-Biphenyl             | 13.457 | 154  | 5824885  | 84.950  | ng/u1  | 99       |
| 44) 2-Chloronaphthalene       | 13.498 | 162  | 4527422  | 84.144  | ng/u1  | 99       |
| 45) 2-Nitroaniline            | 13.704 | 65   | 1591337  | 120.160 | ng/u1  | 99       |
| 47) Dimethylphthalate         | 14.080 | 163  | 5703688  | 85.709  | ng/u1  | 100      |
| 48) 2,6-Dinitrotoluene        | 14.198 | 165  | 1196838  | 110.620 | ng/u1  | 99       |
| 50) Acenaphthylene            | 14.339 | 152  | 7197145  | 84.595  | ng/u1  | 99       |
| 51) 3-Nitroaniline            | 14.527 | 138  | 1244562  | 95.912  | ng/u1  | 97       |
| 52) Acenaphthene              | 14.680 | 153  | 5047544  | 84.792  | ng/u1  | 99       |
| 53) 2,4-Dinitrophenol         | 14.727 | 184  | 706095   | 147.081 | ng/u1  | 96       |
| 55) 4-Nitrophenol             | 14.839 | 109  | 963345   | 92.615  | ng/u1  | 94       |
| 56) Dibenzofuran              | 15.016 | 168  | 6577204  | 82.271  | ng/u1  | 98       |
| 57) 2,4-Dinitrotoluene        | 14.980 | 165  | 1705827  | 108.451 | ng/u1  | 95       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.239 | 232  | 1321186  | 91.321  | ng/u1  | 93       |
| 59) Diethylphthalate          | 15.439 | 149  | 5859889  | 87.963  | ng/u1  | 99       |
| 61) Fluorene                  | 15.663 | 166  | 5601557  | 82.269  | ng/u1  | 100      |
| 62) 4-Chlorophenyl-phenyle... | 15.651 | 204  | 2608296  | 80.100  | ng/u1  | 97       |
| 63) 4-Nitroaniline            | 15.692 | 138  | 1116090  | 88.527  | ng/u1  | 99       |
| 66) 4,6-Dinitro-2-methylph... | 15.745 | 198  | 980804   | 116.932 | ng/u1# | 93       |
| 67) N-Nitrosodiphenylamine    | 15.869 | 169  | 4742249  | 85.703  | ng/u1  | 99       |
| 68) 4-Bromophenyl-phenylether | 16.545 | 248  | 1517051  | 81.925  | ng/u1  | 97       |
| 69) Hexachlorobenzene         | 16.657 | 284  | 1737878  | 79.348  | ng/u1  | 98       |
| 70) Atrazine                  | 16.821 | 200  | 1655468  | 84.783  | ng/u1  | 99       |
| 71) Pentachlorophenol         | 16.998 | 266  | 1137269  | 87.417  | ng/u1  | 99       |
| 72) Phenanthrene              | 17.398 | 178  | 8779118  | 84.837  | ng/u1  | 97       |
| 74) Anthracene                | 17.486 | 178  | 8861059  | 84.244  | ng/u1  | 98       |
| 75) 1,2,3,4-Tetrachloroben... | 13.422 | 216  | 2143018  | 83.400  | ng/uL  | 99       |
| 76) Pentachlorobenzene        | 14.933 | 250  | 2209008  | 79.484  | ng/uL  | 99       |
| 77) Carbazole                 | 17.757 | 167  | 8410254  | 86.362  | ng/u1  | 98       |
| 78) Di-n-butylphthalate       | 18.315 | 149  | 10173561 | 92.170  | ng/u1  | 98       |
| 80) Fluoranthene              | 19.404 | 202  | 10119623 | 85.675  | ng/u1  | 98       |
| 82) Pyrene                    | 19.768 | 202  | 10362117 | 82.877  | ng/u1  | 99       |
| 83) Butylbenzylphthalate      | 20.656 | 149  | 4967536  | 104.852 | ng/u1  | 100      |
| 84) 3,3'-Dichlorobenzidine    | 21.445 | 252  | 3595554  | 93.919  | ng/u1  | 99       |
| 85) Benzo(a)anthracene        | 21.509 | 228  | 10401141 | 84.727  | ng/u1  | 97       |
| 86) Bis(2-ethylhexyl)phtha... | 21.433 | 149  | 7092712  | 102.486 | ng/u1# | 96       |
| 87) Chrysene                  | 21.568 | 228  | 9634729  | 82.685  | ng/u1  | 97       |
| 89) Di-n-octyl phthalate      | 22.368 | 149  | 12009907 | 104.376 | ng/u1  | 100      |
| 90) Benzo(b)fluoranthene      | 23.215 | 252  | 10955402 | 90.518  | ng/u1  | 99       |
| 91) Benzo(k)fluoranthene      | 23.262 | 252  | 10205468 | 84.387  | ng/u1  | 98       |
| 93) Benzo(a)pyrene            | 23.850 | 252  | 9310501  | 87.506  | ng/u1  | 98       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.485 | 276  | 9994601  | 80.439  | ng/u1  | 97       |
| 95) Dibenzo(a,h)anthracene    | 26.503 | 278  | 9201732  | 81.662  | ng/u1  | 99       |
| 96) Benzo(g,h,i)perylene      | 27.262 | 276  | 8339459  | 77.422  | ng/u1  | 99       |

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

Manual Integrations  
APPROVED

Reviewed By :Jagrut Upadhyay 09/30/2022  
Supervised By :mohammad ahmed 10/03/2022

Quant Time: Sep 30 01:13:38 2022  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM092922.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Sep 30 01:10:23 2022  
Response via : Initial Calibration

