

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP053123\  
 Data File : BP015370.D  
 Acq On : 31 May 2023 21:38  
 Operator : CG/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD020734

Quant Time: May 31 23:40:26 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP051123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed May 31 14:24:45 2023  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Yogesh Patel 06/01/2023  
 Supervised By :mohammad ahmed 06/01/2023

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.105  | 152  | 213002   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.934 | 136  | 953034   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.745 | 164  | 622189   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.504 | 188  | 1417222  | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.580 | 240  | 1338029  | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.133 | 264  | 1437021  | 20.000 | ng/u1  | -0.01    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.411  | 96   | 40048    | 7.305  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.852  | 84   | 288243   | 18.549 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.258  | 99   | 364549   | 18.433 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.422  | 67   | 220319   | 19.289 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.628  | 132  | 278429   | 19.148 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.822  | 113  | 300259   | 19.053 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.281  | 128  | 149133   | 19.914 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.010 | 143  | 174622   | 20.956 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.552 | 165  | 307618   | 20.458 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.069 | 131  | 428019   | 19.239 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.151 | 166  | 949768   | 20.201 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.440 | 160  | 1066789  | 19.839 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.945 | 143  | 171514   | 18.787 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.740 | 176  | 801878   | 20.212 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.863 | 200  | 178368   | 19.971 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.604 | 188  | 1260307  | 19.421 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.822 | 212  | 1474164  | 18.996 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.969 | 264  | 1381716  | 19.369 | ng/u1  | -0.01    |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.452  | 88   | 44149    | 7.495  | ng/uL# | 95       |
| 5) Pyridine                   | 3.875  | 79   | 298173   | 18.450 | ng/u1  | 94       |
| 6) Benzaldehyde               | 7.234  | 77   | 136450   | 25.102 | ng/u1  | 98       |
| 8) Phenol                     | 7.287  | 94   | 378979   | 18.688 | ng/u1# | 89       |
| 10) Bis(2-Chloroethyl)ether   | 7.522  | 93   | 303707   | 18.719 | ng/u1  | 97       |
| 12) 2-Chlorophenol            | 7.664  | 128  | 291103   | 18.927 | ng/u1  | 97       |
| 13) 2-Methylphenol            | 8.552  | 108  | 295484   | 18.852 | ng/u1  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.640  | 45   | 396909   | 19.526 | ng/u1  | 99       |
| 16) Acetophenone              | 8.940  | 105  | 486660   | 19.787 | ng/u1# | 92       |
| 17) N-Nitroso-di-n-propyla... | 8.922  | 70   | 257010   | 19.924 | ng/u1  | 91       |
| 18) 4-Methylphenol            | 8.881  | 108  | 321501   | 18.636 | ng/u1  | 96       |
| 19) Hexachloroethane          | 9.193  | 117  | 124521   | 19.828 | ng/u1  | 86       |
| 22) Nitrobenzene              | 9.322  | 77   | 385889   | 20.863 | ng/u1  | 95       |
| 23) Isophorone                | 9.852  | 82   | 746132   | 19.626 | ng/u1# | 98       |
| 25) 2-Nitrophenol             | 10.040 | 139  | 185875   | 20.342 | ng/u1# | 84       |
| 26) 2,4-Dimethylphenol        | 10.099 | 107  | 374018   | 20.512 | ng/u1  | 94       |
| 27) Bis(2-Chloroethoxy)met... | 10.334 | 93   | 436656   | 18.877 | ng/u1  | 98       |
| 29) 2,4-Dichlorophenol        | 10.575 | 162  | 305391   | 20.107 | ng/u1  | 99       |
| 30) Naphthalene               | 10.987 | 128  | 988242   | 19.073 | ng/u1  | 99       |
| 32) 4-Chloroaniline           | 11.099 | 127  | 441668   | 19.316 | ng/u1  | 100      |
| 33) Hexachlorobutadiene       | 11.263 | 225  | 203703   | 22.253 | ng/u1  | 98       |
| 34) Caprolactam               | 11.881 | 113  | 109649m  | 19.149 | ng/u1  |          |
| 35) 4-Chloro-3-methylphenol   | 12.210 | 107  | 357546   | 20.849 | ng/u1  | 95       |
| 36) 2-Methylnaphthalene       | 12.581 | 142  | 682713   | 19.556 | ng/u1  | 98       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP053123\  
 Data File : BP05370.D  
 Acq On : 31 May 2023 21:38  
 Operator : CG/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SSTD020734

Quant Time: May 31 23:40:26 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP051123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed May 31 14:24:45 2023  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Yogesh Patel 06/01/2023  
 Supervised By :mohammad ahmed 06/01/2023

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 37) 1-Methylnaphthalene       | 12.799 | 142  | 692547   | 19.754 | ng/ul# | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 12.946 | 216  | 389337   | 20.923 | ng/ul  | 99       |
| 40) Hexachlorocyclopentadiene | 12.922 | 237  | 252359   | 20.931 | ng/ul  | 99       |
| 41) 2,4,6-Trichlorophenol     | 13.187 | 196  | 267341   | 20.346 | ng/ul  | 96       |
| 42) 2,4,5-Trichlorophenol     | 13.263 | 196  | 285393   | 20.032 | ng/ul  | 99       |
| 43) 1,1'-Biphenyl             | 13.581 | 154  | 947069   | 19.619 | ng/ul  | 99       |
| 44) 2-Chloronaphthalene       | 13.628 | 162  | 748063   | 19.843 | ng/ul  | 98       |
| 45) 2-Nitroaniline            | 13.834 | 65   | 247494   | 22.944 | ng/ul  | 88       |
| 47) Dimethylphthalate         | 14.198 | 163  | 978975   | 19.994 | ng/ul  | 99       |
| 48) 2,6-Dinitrotoluene        | 14.328 | 165  | 209906   | 21.638 | ng/ul  | 100      |
| 50) Acenaphthylene            | 14.469 | 152  | 1171171  | 19.520 | ng/ul  | 99       |
| 51) 3-Nitroaniline            | 14.657 | 138  | 172993   | 19.805 | ng/ul  | 94       |
| 52) Acenaphthene              | 14.810 | 153  | 809152   | 19.746 | ng/ul  | 98       |
| 53) 2,4-Dinitrophenol         | 14.869 | 184  | 161461   | 25.068 | ng/ul# | 87       |
| 55) 4-Nitrophenol             | 14.963 | 109  | 169534   | 25.248 | ng/ul# | 78       |
| 56) Dibenzofuran              | 15.145 | 168  | 1131665  | 19.855 | ng/ul  | 97       |
| 57) 2,4-Dinitrotoluene        | 15.116 | 165  | 302490   | 21.549 | ng/ul  | 94       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.369 | 232  | 257444   | 21.411 | ng/ul  | 99       |
| 59) Diethylphthalate          | 15.557 | 149  | 1000331  | 20.483 | ng/ul  | 100      |
| 61) Fluorene                  | 15.798 | 166  | 921019   | 20.108 | ng/ul  | 99       |
| 62) 4-Chlorophenyl-phenyle... | 15.781 | 204  | 463439   | 20.707 | ng/ul  | 94       |
| 63) 4-Nitroaniline            | 15.816 | 138  | 150917   | 20.007 | ng/ul# | 82       |
| 66) 4,6-Dinitro-2-methylph... | 15.881 | 198  | 182288   | 19.776 | ng/ul# | 99       |
| 67) N-Nitrosodiphenylamine    | 15.998 | 169  | 796557   | 19.115 | ng/ul  | 99       |
| 68) 4-Bromophenyl-phenylether | 16.681 | 248  | 301225   | 20.635 | ng/ul  | 97       |
| 69) Hexachlorobenzene         | 16.804 | 284  | 355479   | 20.607 | ng/ul  | 98       |
| 70) Atrazine                  | 16.951 | 200  | 321869   | 20.897 | ng/ul  | 98       |
| 71) Pentachlorophenol         | 17.151 | 266  | 221759   | 20.551 | ng/ul  | 98       |
| 72) Phenanthrene              | 17.545 | 178  | 1482919  | 19.215 | ng/ul  | 99       |
| 74) Anthracene                | 17.639 | 178  | 1501316  | 19.238 | ng/ul  | 99       |
| 75) 1,2,3,4-Tetrachloroben... | 13.551 | 216  | 401528   | 19.982 | ng/ul  | 99       |
| 76) Pentachlorobenzene        | 15.063 | 250  | 410369   | 20.339 | ng/ul  | 98       |
| 77) Carbazole                 | 17.904 | 167  | 1365271  | 19.951 | ng/ul  | 100      |
| 78) Di-n-butylphthalate       | 18.434 | 149  | 1739997  | 19.745 | ng/ul  | 99       |
| 80) Fluoranthene              | 19.498 | 202  | 1822292  | 19.266 | ng/ul  | 96       |
| 82) Pyrene                    | 19.851 | 202  | 1858309  | 19.025 | ng/ul  | 96       |
| 83) Butylbenzylphthalate      | 20.704 | 149  | 833337   | 19.245 | ng/ul  | 99       |
| 84) 3,3'-Dichlorobenzidine    | 21.486 | 252  | 626963   | 21.335 | ng/ul  | 98       |
| 85) Benzo(a)anthracene        | 21.563 | 228  | 1820456  | 19.677 | ng/ul  | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.457 | 149  | 1225413  | 19.657 | ng/ul  | 99       |
| 87) Chrysene                  | 21.622 | 228  | 1714224  | 19.383 | ng/ul  | 98       |
| 89) Di-n-octyl phthalate      | 22.427 | 149  | 2117290  | 21.513 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 23.345 | 252  | 1861095  | 20.429 | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene      | 23.398 | 252  | 1722486  | 19.691 | ng/ul  | 98       |
| 93) Benzo(a)pyrene            | 24.021 | 252  | 1530599  | 19.008 | ng/ul  | 98       |
| 94) Indeno(1,2,3-cd)pyrene    | 26.839 | 276  | 1810965  | 17.369 | ng/ul  | 96       |
| 95) Dibenzo(a,h)anthracene    | 26.851 | 278  | 1503522  | 17.567 | ng/ul  | 97       |
| 96) Benzo(g,h,i)perylene      | 27.674 | 276  | 1458176  | 17.151 | ng/ul  | 96       |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP053123\  
Data File : BP015370.D  
Acq On : 31 May 2023 21:38  
Operator : CG/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SSTD020734

Quant Time: May 31 23:40:26 2023  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP051123.MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed May 31 14:24:45 2023  
Response via : Initial Calibration

Manual Integrations  
APPROVED

Reviewed By :Yogesh Patel 06/01/2023  
Supervised By :mohammad ahmed 06/01/2023

