

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082422\  
 Data File : BP011559.D  
 Acq On : 25 Aug 2022 11:18  
 Operator : CG/JU  
 Sample : PB147280BS  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS280

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 08/26/2022  
 Supervised By :mohammad ahmed 08/27/2022

Quant Time: Aug 26 01:02:43 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 25 02:14:54 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.546  | 152  | 68032    | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.334 | 136  | 305905   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.222 | 164  | 191978   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 16.992 | 188  | 389498   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.128 | 240  | 367829   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 23.416 | 264  | 369694   | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.064  | 96   | 14111    | 6.437  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.470  | 84   | 178184   | 31.745 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 6.734  | 99   | 217936   | 34.859 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.899  | 67   | 140773   | 34.658 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.081  | 132  | 180244   | 39.199 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.275  | 113  | 182827   | 36.536 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 8.716  | 128  | 88824    | 40.332 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.428  | 143  | 94595    | 42.205 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.963  | 165  | 165780   | 40.070 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.487 | 131  | 237623   | 34.503 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.651 | 166  | 524642   | 38.638 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 13.910 | 160  | 628723   | 41.661 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.457 | 143  | 95402    | 36.467 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.228 | 176  | 436853   | 38.242 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.369 | 200  | 84917    | 40.347 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.092 | 188  | 656939   | 38.083 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.357 | 212  | 737658   | 39.637 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.269 | 264  | 704012   | 39.148 | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.093  | 88   | 28210    | 12.341 | ng/uL  | 98       |
| 5) Pyridine                   | 3.487  | 79   | 185473   | 33.104 | ng/u1  | 95       |
| 6) Benzaldehyde               | 6.705  | 77   | 150414   | 52.426 | ng/u1  | 99       |
| 8) Phenol                     | 6.764  | 94   | 223106   | 34.727 | ng/u1  | 96       |
| 10) Bis(2-Chloroethyl)ether   | 6.993  | 93   | 170619   | 34.045 | ng/u1  | 97       |
| 12) 2-Chlorophenol            | 7.111  | 128  | 182464   | 36.877 | ng/u1  | 98       |
| 13) 2-Methylphenol            | 8.005  | 108  | 165727   | 35.015 | ng/u1  | 98       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.087  | 45   | 258969   | 34.881 | ng/u1  | 99       |
| 16) Acetophenone              | 8.381  | 105  | 289399   | 34.766 | ng/u1  | 100      |
| 17) N-Nitroso-di-n-propyla... | 8.375  | 70   | 151695   | 36.636 | ng/u1  | 97       |
| 18) 4-Methylphenol            | 8.334  | 108  | 184162   | 35.146 | ng/u1  | 94       |
| 19) Hexachloroethane          | 8.611  | 117  | 85900    | 38.273 | ng/u1  | 97       |
| 22) Nitrobenzene              | 8.758  | 77   | 235280   | 37.902 | ng/u1  | 97       |
| 23) Isophorone                | 9.287  | 82   | 408899   | 37.525 | ng/u1  | 99       |
| 25) 2-Nitrophenol             | 9.463  | 139  | 102613   | 40.126 | ng/u1  | 97       |
| 26) 2,4-Dimethylphenol        | 9.528  | 107  | 224028   | 36.896 | ng/u1  | 98       |
| 27) Bis(2-Chloroethoxy)met... | 9.769  | 93   | 240022   | 36.441 | ng/u1  | 98       |
| 29) 2,4-Dichlorophenol        | 9.993  | 162  | 164708   | 38.900 | ng/u1  | 97       |
| 30) Naphthalene               | 10.387 | 128  | 582791   | 36.082 | ng/u1  | 99       |
| 32) 4-Chloroaniline           | 10.510 | 127  | 236451   | 33.138 | ng/u1  | 100      |
| 33) Hexachlorobutadiene       | 10.663 | 225  | 105641   | 39.192 | ng/u1  | 98       |
| 34) Caprolactam               | 11.316 | 113  | 55571m   | 36.197 | ng/u1  |          |
| 35) 4-Chloro-3-methylphenol   | 11.652 | 107  | 207651   | 38.790 | ng/u1  | 99       |
| 36) 2-Methylnaphthalene       | 12.016 | 142  | 391148   | 37.222 | ng/u1  | 96       |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082422\  
 Data File : BP011559.D  
 Acq On : 25 Aug 2022 11:18  
 Operator : CG/JU  
 Sample : PB147280BS  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS280

Quant Time: Aug 26 01:02:43 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 25 02:14:54 2022  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 08/26/2022  
 Supervised By :mohammad ahmed 08/27/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 37) 1-Methylnaphthalene       | 12.234 | 142  | 397157   | 37.263 | ng/ul  | 97       |
| 39) 1,2,4,5-Tetrachloroben... | 12.387 | 216  | 180211   | 36.984 | ng/ul  | 99       |
| 40) Hexachlorocyclopentadiene | 12.357 | 237  | 96325    | 31.864 | ng/ul  | 99       |
| 41) 2,4,6-Trichlorophenol     | 12.640 | 196  | 121036   | 38.594 | ng/ul  | 97       |
| 42) 2,4,5-Trichlorophenol     | 12.710 | 196  | 133964   | 38.374 | ng/ul  | 100      |
| 43) 1,1'-Biphenyl             | 13.046 | 154  | 517033   | 36.144 | ng/ul  | 99       |
| 44) 2-Chloronaphthalene       | 13.087 | 162  | 392725   | 36.433 | ng/ul  | 99       |
| 45) 2-Nitroaniline            | 13.310 | 65   | 149898   | 40.744 | ng/ul  | 97       |
| 47) Dimethylphthalate         | 13.699 | 163  | 511214   | 36.427 | ng/ul  | 98       |
| 48) 2,6-Dinitrotoluene        | 13.822 | 165  | 100358   | 39.776 | ng/ul  | 100      |
| 50) Acenaphthylene            | 13.940 | 152  | 659537   | 36.787 | ng/ul  | 99       |
| 51) 3-Nitroaniline            | 14.151 | 138  | 108866   | 39.613 | ng/ul  | 98       |
| 52) Acenaphthene              | 14.287 | 153  | 436886   | 36.382 | ng/ul  | 99       |
| 53) 2,4-Dinitrophenol         | 14.363 | 184  | 57459    | 37.207 | ng/ul  | 99       |
| 55) 4-Nitrophenol             | 14.475 | 109  | 119163   | 38.229 | ng/ul  | 88       |
| 56) Dibenzofuran              | 14.628 | 168  | 602853   | 36.405 | ng/ul  | 99       |
| 57) 2,4-Dinitrotoluene        | 14.616 | 165  | 151579   | 40.131 | ng/ul  | 91       |
| 58) 2,3,4,6-Tetrachlorophenol | 14.857 | 232  | 107884   | 39.361 | ng/ul  | 97       |
| 59) Diethylphthalate          | 15.075 | 149  | 578729   | 37.651 | ng/ul  | 99       |
| 61) Fluorene                  | 15.281 | 166  | 506068   | 37.356 | ng/ul  | 98       |
| 62) 4-Chlorophenyl-phenyle... | 15.287 | 204  | 231675   | 37.855 | ng/ul  | 98       |
| 63) 4-Nitroaniline            | 15.328 | 138  | 122392m  | 46.105 | ng/ul  |          |
| 66) 4,6-Dinitro-2-methylph... | 15.381 | 198  | 81472    | 37.422 | ng/ul# | 89       |
| 67) N-Nitrosodiphenylamine    | 15.504 | 169  | 426930   | 37.361 | ng/ul  | 99       |
| 68) 4-Bromophenyl-phenylether | 16.187 | 248  | 132012   | 37.968 | ng/ul  | 95       |
| 69) Hexachlorobenzene         | 16.287 | 284  | 152807   | 38.134 | ng/ul  | 99       |
| 70) Atrazine                  | 16.481 | 200  | 153350   | 35.959 | ng/ul  | 96       |
| 71) Pentachlorophenol         | 16.645 | 266  | 92977    | 37.421 | ng/ul  | 99       |
| 72) Phenanthrene              | 17.040 | 178  | 776923   | 36.852 | ng/ul  | 99       |
| 74) Anthracene                | 17.128 | 178  | 789440   | 37.116 | ng/ul  | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 13.004 | 216  | 180903   | 36.131 | ng/uL  | 99       |
| 76) Pentachlorobenzene        | 14.540 | 250  | 179871   | 37.860 | ng/uL  | 100      |
| 77) Carbazole                 | 17.410 | 167  | 726248   | 36.302 | ng/ul  | 99       |
| 78) Di-n-butylphthalate       | 17.987 | 149  | 990369   | 40.200 | ng/ul  | 99       |
| 80) Fluoranthene              | 19.034 | 202  | 876472   | 38.297 | ng/ul  | 96       |
| 82) Pyrene                    | 19.386 | 202  | 912631   | 37.951 | ng/ul  | 98       |
| 83) Butylbenzylphthalate      | 20.286 | 149  | 446560   | 41.396 | ng/ul  | 95       |
| 84) 3,3'-Dichlorobenzidine    | 21.051 | 252  | 284458   | 40.851 | ng/ul  | 99       |
| 85) Benzo(a)anthracene        | 21.110 | 228  | 853458   | 38.084 | ng/ul  | 98       |
| 86) Bis(2-ethylhexyl)phtha... | 21.051 | 149  | 672763   | 42.143 | ng/ul  | 97       |
| 87) Chrysene                  | 21.163 | 228  | 839009   | 38.112 | ng/ul  | 99       |
| 89) Di-n-octyl phthalate      | 21.945 | 149  | 1138430  | 38.308 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene      | 22.722 | 252  | 867582   | 37.727 | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene      | 22.763 | 252  | 838050   | 38.188 | ng/ul  | 98       |
| 93) Benzo(a)pyrene            | 23.316 | 252  | 846614   | 37.708 | ng/ul  | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 25.774 | 276  | 928698   | 37.085 | ng/ul  | 99       |
| 95) Dibenzo(a,h)anthracene    | 25.792 | 278  | 808471   | 37.494 | ng/ul  | 97       |
| 96) Benzo(g,h,i)perylene      | 26.498 | 276  | 695320   | 32.641 | ng/ul  | 99       |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082422\  
 Data File : BP011559.D  
 Acq On : 25 Aug 2022 11:18  
 Operator : CG/JU  
 Sample : PB147280BS  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :

BNA\_P

ClientSampleId :

SLCS280

Manual Integrations

APPROVED

Reviewed By :Jagrut Upadhyay 08/26/2022  
 Supervised By :mohammad ahmed 08/27/2022

Quant Time: Aug 26 01:02:43 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 25 02:14:54 2022  
 Response via : Initial Calibration

