

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112723\  
 Data File : BP018445.D  
 Acq On : 28 Nov 2023 17:33  
 Operator : MA/JU  
 Sample : PB157489BS  
 Misc :  
 ALS Vial : 53 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SLCS489

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 12/01/2023  
 Supervised By :mohammad ahmed 12/01/2023

Quant Time: Nov 29 01:30:00 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112223.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 29 01:22:40 2023  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.858  | 152  | 490518   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.658 | 136  | 1850053  | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.493 | 164  | 984085   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.245 | 188  | 1717155  | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.428 | 240  | 1742187  | 20.000 | ng/u1  | -0.01    |
| 88) Perylene-d12              | 24.086 | 264  | 2061357  | 20.000 | ng/u1  | -0.05    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.264  | 96   | 70620    | 5.392  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.681  | 84   | 912745   | 25.361 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.022  | 99   | 1156124  | 25.790 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.193  | 67   | 665436   | 24.824 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.387  | 132  | 997728   | 26.508 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.569  | 113  | 881260   | 24.360 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.016  | 128  | 444389   | 29.573 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.740  | 143  | 433586   | 31.243 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.275 | 165  | 922286   | 27.546 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.793 | 131  | 1104174  | 24.452 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.910 | 166  | 2178683  | 24.881 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.187 | 160  | 2673264  | 26.793 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.687 | 143  | 309598   | 25.611 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.487 | 176  | 1848744  | 25.242 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.604 | 200  | 218163   | 29.845 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.345 | 188  | 2436572  | 26.548 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.598 | 212  | 2865281  | 20.665 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.904 | 264  | 3318510  | 27.768 | ng/u1  | -0.05    |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 2) 1,4-Dioxane                | 3.299  | 88   | 149115   | 10.618 | ng/uL# | 89       |
| 5) Pyridine                   | 3.705  | 79   | 959933   | 26.471 | ng/u1  | 91       |
| 6) Benzaldehyde               | 6.993  | 77   | 653087   | 27.530 | ng/u1  | 95       |
| 8) Phenol                     | 7.052  | 94   | 1192385  | 27.116 | ng/u1  | 93       |
| 10) Bis(2-Chloroethyl)ether   | 7.287  | 93   | 1012693  | 27.647 | ng/u1  | 95       |
| 12) 2-Chlorophenol            | 7.417  | 128  | 1032712  | 27.349 | ng/u1  | 95       |
| 13) 2-Methylphenol            | 8.305  | 108  | 903357   | 26.375 | ng/u1  | 99       |
| 14) 2,2'-oxybis(1-Chloropr... | 8.393  | 45   | 1189365  | 22.843 | ng/u1# | 95       |
| 16) Acetophenone              | 8.681  | 105  | 1420899  | 25.668 | ng/u1  | 94       |
| 17) N-Nitroso-di-n-propyla... | 8.675  | 70   | 639950   | 20.694 | ng/u1  | 92       |
| 18) 4-Methylphenol            | 8.634  | 108  | 950320   | 25.982 | ng/u1  | 98       |
| 19) Hexachloroethane          | 8.934  | 117  | 431615   | 27.371 | ng/u1  | 96       |
| 22) Nitrobenzene              | 9.058  | 77   | 983763   | 27.944 | ng/u1  | 95       |
| 23) Isophorone                | 9.593  | 82   | 1863879  | 23.755 | ng/u1  | 97       |
| 25) 2-Nitrophenol             | 9.769  | 139  | 502417   | 32.177 | ng/u1  | 91       |
| 26) 2,4-Dimethylphenol        | 9.834  | 107  | 737001   | 21.023 | ng/u1  | 98       |
| 27) Bis(2-Chloroethoxy)met... | 10.075 | 93   | 1246294  | 26.300 | ng/u1  | 99       |
| 29) 2,4-Dichlorophenol        | 10.305 | 162  | 915826   | 28.556 | ng/u1  | 98       |
| 30) Naphthalene               | 10.705 | 128  | 3154267  | 28.190 | ng/u1  | 100      |
| 32) 4-Chloroaniline           | 10.816 | 127  | 1039135  | 24.386 | ng/u1  | 99       |
| 33) Hexachlorobutadiene       | 10.993 | 225  | 696069   | 31.822 | ng/u1  | 99       |
| 34) Caprolactam               | 11.599 | 113  | 224174   | 24.256 | ng/u1  | 87       |
| 35) 4-Chloro-3-methylphenol   | 11.946 | 107  | 784075   | 22.866 | ng/u1  | 94       |
| 36) 2-Methylnaphthalene       | 12.316 | 142  | 2050545  | 25.894 | ng/u1  | 99       |

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 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| 37) 1-Methylnaphthalene       | 12.534 | 142  | 1999103  | 25.079 | ng/u1  | 100      |
| 39) 1,2,4,5-Tetrachloroben... | 12.681 | 216  | 1181821  | 34.003 | ng/u1  | 97       |
| 40) Hexachlorocyclopentadiene | 12.663 | 237  | 513492   | 26.086 | ng/u1  | 99       |
| 41) 2,4,6-Trichlorophenol     | 12.922 | 196  | 640465   | 30.684 | ng/u1  | 99       |
| 42) 2,4,5-Trichlorophenol     | 12.993 | 196  | 677269   | 30.007 | ng/u1  | 100      |
| 43) 1,1'-Biphenyl             | 13.328 | 154  | 2574679  | 30.028 | ng/u1  | 99       |
| 44) 2-Chloronaphthalene       | 13.369 | 162  | 2048427  | 30.185 | ng/u1  | 98       |
| 45) 2-Nitroaniline            | 13.575 | 65   | 379346   | 25.979 | ng/u1  | 85       |
| 47) Dimethylphthalate         | 13.957 | 163  | 2270914  | 26.505 | ng/u1  | 99       |
| 48) 2,6-Dinitrotoluene        | 14.075 | 165  | 415805   | 28.786 | ng/u1  | 88       |
| 50) Acenaphthylene            | 14.216 | 152  | 3109683  | 27.961 | ng/u1  | 100      |
| 51) 3-Nitroaniline            | 14.398 | 138  | 391589   | 27.285 | ng/u1# | 92       |
| 52) Acenaphthene              | 14.557 | 153  | 2030622  | 27.950 | ng/u1  | 99       |
| 53) 2,4-Dinitrophenol         | 14.598 | 184  | 138933   | 27.569 | ng/u1  | 98       |
| 55) 4-Nitrophenol             | 14.704 | 109  | 223255   | 24.929 | ng/u1  | 91       |
| 56) Dibenzofuran              | 14.893 | 168  | 2716963  | 27.354 | ng/u1  | 98       |
| 57) 2,4-Dinitrotoluene        | 14.857 | 165  | 540083   | 27.668 | ng/u1# | 87       |
| 58) 2,3,4,6-Tetrachlorophenol | 15.116 | 232  | 479822   | 27.070 | ng/u1  | 95       |
| 59) Diethylphthalate          | 15.322 | 149  | 2025639  | 23.536 | ng/u1  | 99       |
| 61) Fluorene                  | 15.540 | 166  | 2087568  | 25.773 | ng/u1  | 100      |
| 62) 4-Chlorophenyl-phenyle... | 15.540 | 204  | 1112338  | 27.044 | ng/u1  | 96       |
| 63) 4-Nitroaniline            | 15.563 | 138  | 379136   | 27.992 | ng/u1  | 92       |
| 66) 4,6-Dinitro-2-methylph... | 15.616 | 198  | 249787   | 30.811 | ng/u1  | 98       |
| 67) N-Nitrosodiphenylamine    | 15.751 | 169  | 1701926  | 29.227 | ng/u1  | 100      |
| 68) 4-Bromophenyl-phenylether | 16.434 | 248  | 643886   | 29.699 | ng/u1  | 97       |
| 69) Hexachlorobenzene         | 16.545 | 284  | 750226   | 30.953 | ng/u1# | 92       |
| 70) Atrazine                  | 16.710 | 200  | 393266   | 22.122 | ng/u1  | 99       |
| 71) Pentachlorophenol         | 16.892 | 266  | 369249   | 29.650 | ng/u1  | 97       |
| 72) Phenanthrene              | 17.287 | 178  | 3023886  | 28.951 | ng/u1  | 100      |
| 74) Anthracene                | 17.381 | 178  | 2949808  | 27.979 | ng/u1  | 100      |
| 75) 1,2,3,4-Tetrachloroben... | 13.287 | 216  | 1138438  | 38.570 | ng/uL  | 100      |
| 76) Pentachlorobenzene        | 14.810 | 250  | 1012628  | 35.319 | ng/uL  | 99       |
| 77) Carbazole                 | 17.651 | 167  | 2625579  | 30.740 | ng/u1  | 99       |
| 78) Di-n-butylphthalate       | 18.222 | 149  | 2908753  | 24.483 | ng/u1  | 100      |
| 80) Fluoranthene              | 19.275 | 202  | 3438499  | 20.758 | ng/u1  | 97       |
| 82) Pyrene                    | 19.628 | 202  | 3623787  | 21.725 | ng/u1  | 98       |
| 83) Butylbenzylphthalate      | 20.569 | 149  | 1254671  | 20.326 | ng/u1  | 95       |
| 84) 3,3'-Dichlorobenzidine    | 21.369 | 252  | 1197396  | 28.417 | ng/u1  | 100      |
| 85) Benzo(a)anthracene        | 21.410 | 228  | 4032015  | 28.398 | ng/u1  | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.457 | 149  | 1881208  | 20.313 | ng/u1  | 99       |
| 87) Chrysene                  | 21.463 | 228  | 3705349  | 28.678 | ng/u1  | 100      |
| 89) Di-n-octyl phthalate      | 22.622 | 149  | 3210336  | 21.154 | ng/u1  | 100      |
| 90) Benzo(b)fluoranthene      | 23.257 | 252  | 4511389  | 30.131 | ng/u1  | 99       |
| 91) Benzo(k)fluoranthene      | 23.316 | 252  | 3955054  | 26.535 | ng/u1  | 99       |
| 93) Benzo(a)pyrene            | 23.963 | 252  | 4044139  | 29.370 | ng/u1  | 99       |
| 94) Indeno(1,2,3-cd)pyrene    | 27.157 | 276  | 4745494m | 29.806 | ng/u1  |          |
| 95) Dibenzo(a,h)anthracene    | 27.327 | 278  | 3973125m | 30.482 | ng/u1  |          |
| 96) Benzo(g,h,i)perylene      | 28.110 | 276  | 3925461m | 30.443 | ng/u1  |          |

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

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APPROVED

Reviewed By :Yogesh Patel 12/01/2023

Supervised By :mohammad ahmed 12/01/2023

Quant Time: Nov 29 01:30:00 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112223.MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 29 01:22:40 2023

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