

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG062724\  
 Data File : BG061771.D  
 Acq On : 28 Jun 2024 14:10  
 Operator : MA/JU  
 Sample : P2934-17  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 A4CQ3

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 06/29/2024  
 Supervised By :mohammad ahmed 06/29/2024

Quant Time: Jun 29 00:23:45 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG062024.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jun 27 12:23:53 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.059  | 152  | 73121    | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.868 | 136  | 334096   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.687 | 164  | 212979   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.431 | 188  | 456553m  | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.690 | 240  | 340412   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.910 | 264  | 398377   | 20.000 | ng/u1  | 0.00     |
|                               |        |      |          |        |        |          |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.459  | 96   | 10729    | 5.300  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.876  | 84   | 122746   | 20.057 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.207  | 99   | 204808   | 25.801 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.384  | 67   | 127869   | 25.841 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.583  | 132  | 147047   | 26.995 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.753  | 113  | 157558   | 25.675 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.223  | 128  | 74071    | 32.652 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.945  | 143  | 83581    | 36.378 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.480 | 165  | 157035   | 28.865 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 11.003 | 131  | 188536   | 22.890 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.088 | 166  | 472503   | 28.830 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.381 | 160  | 553399   | 30.810 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.869 | 143  | 65758    | 23.362 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.680 | 176  | 412045   | 28.718 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.786 | 200  | 28968m   | 14.838 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.531 | 188  | 618995   | 29.702 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.810 | 212  | 627812   | 33.017 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.693 | 264  | 587996   | 30.903 | ng/u1  | 0.00     |
|                               |        |      |          |        |        |          |
| Target Compounds              |        |      |          |        | Qvalue |          |
| 72) Phenanthrene              | 17.472 | 178  | 155360m  | 6.468  | ng/u1  |          |
| 74) Anthracene                | 17.566 | 178  | 29693    | 1.209  | ng/u1  | 97       |
| 80) Fluoranthene              | 19.481 | 202  | 219463   | 9.947  | ng/u1  | 97       |
| 82) Pyrene                    | 19.840 | 202  | 212585   | 9.062  | ng/u1  | 95       |
| 85) Benzo(a)anthracene        | 21.673 | 228  | 99158    | 4.150  | ng/u1  | 93       |
| 87) Chrysene                  | 21.737 | 228  | 97256    | 4.301  | ng/u1  | 95       |
| 90) Benzo(b)fluoranthene      | 23.894 | 252  | 128614   | 5.371  | ng/u1# | 97       |
| 91) Benzo(k)fluoranthene      | 23.952 | 252  | 46266m   | 1.879  | ng/u1  |          |
| 93) Benzo(a)pyrene            | 24.757 | 252  | 97045    | 4.209  | ng/u1# | 93       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.571 | 276  | 72674m   | 2.530  | ng/u1  |          |
| 96) Benzo(g,h,i)perylene      | 29.716 | 276  | 66656m   | 2.878  | ng/u1  |          |
| -----                         |        |      |          |        |        |          |

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

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