

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
 Data File : BG061842.D  
 Acq On : 2 Jul 2024 11:42  
 Operator : MA/JU  
 Sample : P2963-10  
 Misc :  
 ALS Vial : 39 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 A4CR3

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 07/02/2024  
 Supervised By :mohammad ahmed 07/04/2024

Quant Time: Jul 02 12:18:03 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  | 68426    | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.868 | 136  | 314950   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.687 | 164  | 199874   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.431 | 188  | 422934   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.691 | 240  | 313964   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.916 | 264  | 380020   | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.465  | 96   | 7137     | 3.768  | ng/uL  | 0.01     |
| 4) Pyridine-d5                | 3.882  | 84   | 85947    | 15.007 | ng/u1  | 0.01     |
| 7) Phenol-d5                  | 7.213  | 99   | 196788   | 26.492 | ng/u1  | 0.01     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.384  | 67   | 142825   | 30.843 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.589  | 132  | 158788   | 31.150 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.765  | 113  | 166894   | 29.063 | ng/u1  | 0.01     |
| 21) Nitrobenzene-d5           | 9.223  | 128  | 85741    | 40.094 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.951  | 143  | 102920   | 47.519 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.492 | 165  | 180205   | 35.138 | ng/u1  | 0.01     |
| 31) 4-Chloroaniline-d4        | 11.003 | 131  | 88550    | 11.404 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 14.094 | 166  | 592046   | 38.492 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.387 | 160  | 674502   | 40.015 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.887 | 143  | 59749    | 22.619 | ng/u1  | 0.02     |
| 60) Fluorene-d10              | 15.680 | 176  | 521185   | 38.706 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.792 | 200  | 42873    | 23.706 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.531 | 188  | 768261   | 39.795 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.816 | 212  | 779112   | 44.426 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.693 | 264  | 768803   | 42.358 | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 72) Phenanthrene              | 17.472 | 178  | 116372   | 5.230  | ng/u1  | 96       |
| 74) Anthracene                | 17.566 | 178  | 25379    | 1.115  | ng/u1# | 92       |
| 80) Fluoranthene              | 19.476 | 202  | 210884   | 10.364 | ng/u1# | 95       |
| 82) Pyrene                    | 19.840 | 202  | 200726   | 9.277  | ng/u1  | 96       |
| 85) Benzo(a)anthracene        | 21.673 | 228  | 98388    | 4.464  | ng/u1  | 95       |
| 86) Bis(2-ethylhexyl)phtha... | 21.579 | 149  | 74006    | 6.391  | ng/u1  | 94       |
| 87) Chrysene                  | 21.738 | 228  | 108909   | 5.223  | ng/u1  | 97       |
| 90) Benzo(b)fluoranthene      | 23.888 | 252  | 153263m  | 6.709  | ng/u1  |          |
| 91) Benzo(k)fluoranthene      | 23.953 | 252  | 39680    | 1.689  | ng/u1# | 90       |
| 93) Benzo(a)pyrene            | 24.758 | 252  | 105743   | 4.808  | ng/u1  | 98       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.577 | 276  | 78882m   | 2.878  | ng/u1  |          |
| 96) Benzo(g,h,i)perylene      | 29.734 | 276  | 77327m   | 3.500  | ng/u1  |          |
| -----                         |        |      |          |        |        |          |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG070124\  
Data File : BG061842.D  
Acq On : 2 Jul 2024 11:42  
Operator : MA/JU  
Sample : P2963-10  
Misc :  
ALS Vial : 39 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
A4CR3

Manual Integrations  
APPROVED

Reviewed By :Jagrut Upadhyay 07/02/2024  
Supervised By :mohammad ahmed 07/04/2024

Quant Time: Jul 02 12:18:03 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

