

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG060624\  
 Data File : BG061517.D  
 Acq On : 6 Jun 2024 19:46  
 Operator : MA/JU  
 Sample : P2623-11DL 5X  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BHBR6DL

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 06/07/2024  
 Supervised By :mohammad ahmed 06/08/2024

Quant Time: Jun 06 22:57:38 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG052124.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jun 06 12:15:45 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.876  | 152  | 37198    | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.667 | 136  | 130542   | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.509 | 164  | 91427    | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.259 | 188  | 209177   | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.519 | 240  | 229453   | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.609 | 264  | 265285   | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 0.000  | 96   | 0d       | 0.000  | ng/uL  |          |
| 4) Pyridine-d5                | 3.769  | 84   | 3809     | 1.611  | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 7.071  | 99   | 14423    | 4.722  | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.206  | 67   | 9594     | 4.997  | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.424  | 132  | 12327    | 5.453  | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.605  | 113  | 10034    | 3.852  | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.033  | 128  | 6241     | 6.209  | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.756  | 143  | 7063     | 6.004  | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.314 | 165  | 13719    | 5.982  | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.808 | 131  | 7615     | 2.538  | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.916 | 166  | 43538    | 6.140  | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.204 | 160  | 45657    | 6.193  | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.768 | 143  | 3515     | 4.010  | ng/u1  | 0.01     |
| 60) Fluorene-d10              | 15.502 | 176  | 37706    | 6.159  | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.655 | 200  | 3848     | 3.042  | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.359 | 188  | 60622    | 6.574  | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.656 | 212  | 69140    | 5.869  | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.398 | 264  | 61430    | 4.817  | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 30) Naphthalene               | 10.714 | 128  | 10211    | 1.482  | ng/u1  | 96       |
| 52) Acenaphthene              | 14.574 | 153  | 12690    | 2.366  | ng/u1  | 90       |
| 56) Dibenzofuran              | 14.909 | 168  | 17989    | 2.346  | ng/u1  | 96       |
| 61) Fluorene                  | 15.561 | 166  | 16604    | 2.503  | ng/u1  | 91       |
| 72) Phenanthrene              | 17.306 | 178  | 325547   | 32.065 | ng/u1  | 97       |
| 74) Anthracene                | 17.394 | 178  | 71730    | 6.833  | ng/u1  | 98       |
| 77) Carbazole                 | 17.670 | 167  | 30123    | 3.528  | ng/u1  | 98       |
| 80) Fluoranthene              | 19.321 | 202  | 495003   | 35.024 | ng/u1  | 99       |
| 82) Pyrene                    | 19.686 | 202  | 330600   | 22.535 | ng/u1  | 98       |
| 85) Benzo(a)anthracene        | 21.501 | 228  | 244985   | 15.474 | ng/u1  | 97       |
| 86) Bis(2-ethylhexyl)phtha... | 21.407 | 149  | 15303    | 2.051  | ng/u1# | 98       |
| 87) Chrysene                  | 21.566 | 228  | 214464   | 14.711 | ng/u1  | 98       |
| 90) Benzo(b)fluoranthene      | 23.634 | 252  | 268306   | 16.987 | ng/u1  | 98       |
| 91) Benzo(k)fluoranthene      | 23.693 | 252  | 93868m   | 5.854  | ng/u1  |          |
| 93) Benzo(a)pyrene            | 24.462 | 252  | 108618   | 7.242  | ng/u1  | 98       |
| 94) Indeno(1,2,3-cd)pyrene    | 28.117 | 276  | 88448    | 4.877  | ng/u1# | 97       |
| 95) Dibenzo(a,h)anthracene    | 28.158 | 278  | 32529m   | 2.175  | ng/u1  |          |
| -----                         |        |      |          |        |        |          |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG060624\  
 Data File : BG061517.D  
 Acq On : 6 Jun 2024 19:46  
 Operator : MA/JU  
 Sample : P2623-11DL 5X  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BHBR6DL

Quant Time: Jun 06 22:57:38 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG052124.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jun 06 12:15:45 2024  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 06/07/2024  
 Supervised By :mohammad ahmed 06/08/2024

