

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG120621\  
 Data File : BG051392.D  
 Acq On : 7 Dec 2021 21:23  
 Operator : CG/JU  
 Sample : M4868-11MEDL2 25X  
 Misc :  
 ALS Vial : 50 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGKP6MEDL2

Quant Time: Dec 08 02:23:42 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG112321.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Dec 03 15:23:09 2021  
 Response via : Initial Calibration

### Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 12/08/2021  
 Supervised By :mohammad ahmed 12/15/2021

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 8.186  | 152  | 29640    | 20.000 | ng/u1  | -0.02    |
| 20) Naphthalene-d8            | 11.012 | 136  | 129887   | 20.000 | ng/u1  | -0.02    |
| 38) Acenaphthene-d10          | 14.819 | 164  | 84621    | 20.000 | ng/u1  | -0.01    |
| 64) Phenanthrene-d10          | 17.569 | 188  | 173828   | 20.000 | ng/u1  | -0.01    |
| 79) Chrysene-d12              | 21.870 | 240  | 151896   | 20.000 | ng/u1  | -0.01    |
| 88) Perylene-d12              | 25.266 | 264  | 150502   | 20.000 | ng/u1  | -0.02    |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 0.000  | 96   | 0        | 0.000  | ng/uL  |          |
| 4) Pyridine-d5                | 0.000  | 84   | 0d       | 0.000  | ng/u1  |          |
| 7) Phenol-d5                  | 7.357  | 99   | 2017     | 0.689  | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.510  | 67   | 1577     | 0.857  | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.728  | 132  | 1702     | 0.807  | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.920  | 113  | 1231     | 0.521  | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 9.379  | 128  | 692      | 0.631  | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 10.101 | 143  | 753      | 0.609  | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.683 | 165  | 1300m    | 0.619  | ng/u1  | 0.03     |
| 31) 4-Chloroaniline-d4        | 11.176 | 131  | 1507m    | 0.491  | ng/u1  | 0.01     |
| 46) Dimethylphthalate-d6      | 14.214 | 166  | 5914     | 0.908  | ng/u1  | -0.01    |
| 49) Acenaphthylene-d8         | 14.520 | 160  | 8124     | 0.989  | ng/u1  | -0.01    |
| 54) 4-Nitrophenol-d4          | 0.000  | 143  | 0        | 0.000  | ng/u1  |          |
| 60) Fluorene-d10              | 15.818 | 176  | 6163     | 1.051  | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 0.000  | 200  | 0        | 0.000  | ng/u1  |          |
| 73) Anthracene-d10            | 17.675 | 188  | 9892     | 1.190  | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.954 | 212  | 10556    | 1.149  | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 25.031 | 264  | 8632     | 1.074  | ng/u1  | -0.01    |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 30) Naphthalene               | 11.065 | 128  | 18563    | 2.627  | ng/u1  | 99       |
| 36) 2-Methylnaphthalene       | 12.663 | 142  | 17596    | 3.660  | ng/u1  | 99       |
| 37) 1-Methylnaphthalene       | 12.874 | 142  | 7167     | 1.449  | ng/u1# | 89       |
| 83) Butylbenzylphthalate      | 20.836 | 149  | 187471   | 40.837 | ng/u1  | 99       |
| 86) Bis(2-ethylhexyl)phtha... | 21.705 | 149  | 184724   | 27.963 | ng/u1  | 99       |
| 89) Di-n-octyl phthalate      | 22.974 | 149  | 33204    | 3.045  | ng/u1  | 100      |

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

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