

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120820\  
 Data File : BG047669.D  
 Acq On : 9 Dec 2020 11:01  
 Operator : CG/JU  
 Sample : L4921-05DL 10X  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 C0FW4DL

Manual Integrations  
 APPROVED

mohammad  
 12/9/2020 5:40:05 PM

Quant Time: Dec 09 12:20:48 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG120720MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Dec 08 17:39:41 2020  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|--------|--------|----------|
| 1) 1,4-Dichlorobenzene-d4      | 8.22  | 152  | 27041    | 20.00  | ng/ul  | 0.00     |
| 18) Naphthalene-d8             | 11.04 | 136  | 98958    | 20.00  | ng/ul  | 0.00     |
| 36) Acenaphthene-d10           | 14.83 | 164  | 81756    | 20.00  | ng/ul  | 0.00     |
| 62) Phenanthrene-d10           | 17.57 | 188  | 220410   | 20.00  | ng/ul  | 0.00     |
| 78) Chrysene-d12               | 21.85 | 240  | 242535   | 20.00  | ng/ul  | 0.00     |
| 86) Perylene-d12               | 25.21 | 264  | 244222   | 20.00  | ng/ul  | 0.00     |
| System Monitoring Compounds    |       |      |          |        |        |          |
| 3) 1,4-Dioxane-d8              | 0.00  | 96   | 0        | 0.00   | ng/uL  |          |
| 5) Phenol-d5                   | 7.35  | 99   | 1622     | 0.70   | ng/ul  | 0.00     |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.53  | 67   | 4244     | 3.57   | ng/ul  | 0.00     |
| 9) 2-Chlorophenol-d4           | 7.73  | 132  | 4576     | 2.61   | ng/ul  | 0.00     |
| 13) 4-Methylphenol-d8          | 8.90  | 113  | 2367     | 1.29   | ng/ul  | 0.00     |
| 19) Nitrobenzene-d5            | 9.39  | 128  | 2653     | 3.29   | ng/ul  | 0.00     |
| 22) 2-Nitrophenol-d4           | 10.10 | 143  | 3614     | 3.72   | ng/ul  | 0.00     |
| 26) 2,4-Dichlorophenol-d3      | 10.65 | 165  | 6345     | 3.00   | ng/ul  | 0.00     |
| 29) 4-Chloroaniline-d4         | 0.00  | 131  | 0d       | 0.00   | ng/ul  |          |
| 44) Dimethylphthalate-d6       | 14.23 | 166  | 20043    | 2.82   | ng/ul  | 0.00     |
| 47) Acenaphthylene-d8          | 14.53 | 160  | 30769    | 3.79   | ng/ul  | 0.00     |
| 52) 4-Nitrophenol-d4           | 15.00 | 143  | 538      | 0.51   | ng/ul  | 0.00     |
| 58) Fluorene-d10               | 15.82 | 176  | 20622    | 3.15   | ng/ul  | 0.00     |
| 63) 4,6-Dinitro-2-methylphenol | 0.00  | 200  | 0d       | 0.00   | ng/ul  |          |
| 71) Anthracene-d10             | 17.67 | 188  | 42284m   | 3.85   | ng/ul  | 0.00     |
| 79) Pyrene-d10                 | 19.94 | 212  | 51064    | 3.56   | ng/ul  | 0.00     |
| 90) Benzo(a)pyrene-d12         | 24.98 | 264  | 53837    | 3.77   | ng/ul  | 0.00     |
| Target Compounds               |       |      |          |        |        |          |
| 20) Nitrobenzene               | 9.42  | 77   | 23951    | 9.381  | ng/ul# | 89       |
| 37) 1,2,4,5-Tetrachlorobenzene | 13.04 | 216  | 11896    | 3.661  | ng/ul# | 81       |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.65 | 216  | 36011    | 10.427 | ng/uL  | 94       |

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

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