

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111920\  
 Data File : BG047365.D  
 Acq On : 19 Nov 2020 23:25  
 Operator : CG/JU  
 Sample : L4710-22  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 C0AY7

Manual Integrations  
 APPROVED

mohammad  
 11/23/2020 1:28:56 PM

Quant Time: Nov 20 07:28:47 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG111920MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Nov 19 14:25:26 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.27  | 152  | 50530    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.10 | 136  | 201523   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.88 | 164  | 141836   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.62 | 188  | 333458   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.90 | 240  | 308156   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 25.30 | 264  | 323408   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.62  | 96  | 3082   | 2.22  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.40  | 99  | 68711  | 13.56 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.59  | 67  | 40839  | 14.13 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.80  | 132 | 48217  | 14.19 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.95  | 113 | 48316  | 12.41 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.44  | 128 | 23167  | 13.92 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.17 | 143 | 26942  | 14.83 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.70 | 165 | 56836  | 13.67 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.22 | 131 | 54618  | 12.99 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 14.28 | 166 | 176641 | 14.51 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.58 | 160 | 222407 | 15.81 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 15.04 | 143 | 23979  | 12.91 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.87 | 176 | 160853 | 14.42 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.96 | 200 | 24503  | 12.91 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.72 | 188 | 259417 | 15.68 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.98 | 212 | 303962 | 16.95 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 25.06 | 264 | 296631 | 15.94 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |       | Ovalue |    |
|--------------------------------|-------|-----|--------|-------|--------|----|
| 45) Dimethylphthalate          | 14.32 | 163 | 80236  | 6.460 | ng/ul  | 96 |
| 70) Phenanthrene               | 17.66 | 178 | 30003  | 1.666 | ng/ul# | 95 |
| 77) Fluoranthene               | 19.65 | 202 | 60230  | 2.648 | ng/ul  | 98 |
| 80) Pyrene                     | 20.01 | 202 | 60595  | 2.776 | ng/ul  | 97 |
| 83) Benzo(a)anthracene         | 21.87 | 228 | 35038  | 1.598 | ng/ul  | 98 |
| 84) Bis(2-ethylhexyl)phthalate | 21.77 | 149 | 12821  | 1.200 | ng/ul  | 98 |
| 85) Chrysene                   | 21.95 | 228 | 40772m | 1.954 | ng/ul  |    |
| 88) Benzo(b)fluoranthene       | 24.21 | 252 | 54799m | 2.400 | ng/ul  |    |
| 91) Benzo(a)pyrene             | 25.14 | 252 | 36015  | 1.769 | ng/ul# | 97 |
| 92) Indeno(1,2,3-cd)pyrene     | 29.18 | 276 | 26598m | 1.074 | ng/ul  |    |
| 94) Benzo(g,h,i)perylene       | 30.41 | 276 | 24510m | 1.200 | ng/ul  |    |

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

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