

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120420\  
 Data File : BG047587.D  
 Acq On : 5 Dec 2020 2:23  
 Operator : CG/JU  
 Sample : L4855-14  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 CB801

Manual Integrations  
 APPROVED

mohammad  
 12/7/2020 2:00:21 PM

Quant Time: Dec 05 03:19:54 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG111920MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Dec 05 02:47:28 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.25  | 152  | 26200    | 20.00 | ng/ul | -0.02    |
| 18) Naphthalene-d8        | 11.08 | 136  | 105958   | 20.00 | ng/ul | -0.02    |
| 36) Acenaphthene-d10      | 14.87 | 164  | 80800    | 20.00 | ng/ul | -0.02    |
| 62) Phenanthrene-d10      | 17.60 | 188  | 184185   | 20.00 | ng/ul | -0.02    |
| 78) Chrysene-d12          | 21.88 | 240  | 169316   | 20.00 | ng/ul | -0.02    |
| 86) Perylene-d12          | 25.27 | 264  | 182309   | 20.00 | ng/ul | -0.03    |

## System Monitoring Compounds

|                                |       |     |        |       |       |       |
|--------------------------------|-------|-----|--------|-------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.59  | 96  | 3543   | 4.92  | ng/uL | -0.03 |
| 5) Phenol-d5                   | 7.38  | 99  | 83048  | 31.61 | ng/ul | -0.02 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.56  | 67  | 43108  | 28.77 | ng/ul | -0.02 |
| 9) 2-Chlorophenol-d4           | 7.78  | 132 | 62822  | 35.65 | ng/ul | -0.02 |
| 13) 4-Methylphenol-d8          | 8.94  | 113 | 64478  | 31.93 | ng/ul | -0.02 |
| 19) Nitrobenzene-d5            | 9.42  | 128 | 29341  | 33.52 | ng/ul | -0.02 |
| 22) 2-Nitrophenol-d4           | 10.14 | 143 | 38877  | 40.71 | ng/ul | -0.03 |
| 26) 2,4-Dichlorophenol-d3      | 10.68 | 165 | 78348  | 35.83 | ng/ul | -0.02 |
| 29) 4-Chloroaniline-d4         | 11.20 | 131 | 66796  | 30.21 | ng/ul | -0.02 |
| 44) Dimethylphthalate-d6       | 14.26 | 166 | 229455 | 33.08 | ng/ul | -0.02 |
| 47) Acenaphthylene-d8          | 14.56 | 160 | 274510 | 34.24 | ng/ul | -0.02 |
| 52) 4-Nitrophenol-d4           | 15.03 | 143 | 33662  | 31.81 | ng/ul | 0.00  |
| 58) Fluorene-d10               | 15.85 | 176 | 208578 | 32.83 | ng/ul | -0.02 |
| 63) 4,6-Dinitro-2-methylphenol | 15.95 | 200 | 46931  | 44.77 | ng/ul | -0.02 |
| 71) Anthracene-d10             | 17.70 | 188 | 318937 | 34.91 | ng/ul | -0.02 |
| 79) Pyrene-d10                 | 19.97 | 212 | 347182 | 35.24 | ng/ul | -0.01 |
| 90) Benzo(a)pyrene-d12         | 25.04 | 264 | 360612 | 34.37 | ng/ul | -0.02 |

## Target Compounds

|                                |       |     |        |       | Ovalue |    |
|--------------------------------|-------|-----|--------|-------|--------|----|
| 4) Benzaldehyde                | 7.38  | 77  | 2270   | 1.463 | ng/ul  | 97 |
| 6) Phenol                      | 7.42  | 94  | 3052   | 1.212 | ng/ul  | 91 |
| 28) Naphthalene                | 11.13 | 128 | 9325   | 1.554 | ng/ul# | 92 |
| 45) Dimethylphthalate          | 14.30 | 163 | 9232   | 1.305 | ng/ul# | 96 |
| 70) Phenanthrene               | 17.64 | 178 | 47038  | 4.729 | ng/ul  | 97 |
| 77) Fluoranthene               | 19.63 | 202 | 97535  | 7.765 | ng/ul# | 96 |
| 80) Pyrene                     | 20.00 | 202 | 83230  | 6.940 | ng/ul# | 89 |
| 83) Benzo(a)anthracene         | 21.86 | 228 | 52215  | 4.334 | ng/ul  | 95 |
| 84) Bis(2-ethylhexyl)phthalate | 21.74 | 149 | 13920  | 2.371 | ng/ul# | 90 |
| 85) Chrysene                   | 21.93 | 228 | 50369m | 4.394 | ng/ul  |    |
| 88) Benzo(b)fluoranthene       | 24.19 | 252 | 69606  | 5.408 | ng/ul  | 99 |
| 89) Benzo(k)fluoranthene       | 24.25 | 252 | 27936m | 2.202 | ng/ul  |    |
| 91) Benzo(a)pyrene             | 25.11 | 252 | 47126  | 4.107 | ng/ul# | 95 |
| 92) Indeno(1,2,3-cd)pyrene     | 29.16 | 276 | 33634m | 2.410 | ng/ul  |    |
| 94) Benzo(g,h,i)perylene       | 30.37 | 276 | 36362m | 3.157 | ng/ul  |    |

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

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