

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120320\  
 Data File : BG047541.D  
 Acq On : 3 Dec 2020 16:43  
 Operator : CG/JU  
 Sample : L4865-02  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 C0B20

Manual Integrations  
 APPROVED

mohammad  
 12/4/2020 4:30:21 PM

Quant Time: Dec 03 17:53:43 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG111920MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Dec 03 15:14:44 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.25  | 152  | 31479    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.08 | 136  | 119558   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.86 | 164  | 89515    | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.60 | 188  | 240540   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.88 | 240  | 277213   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 25.26 | 264  | 300116   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.58  | 96  | 1829m   | 2.11  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.38  | 99  | 29456   | 9.33  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.56  | 67  | 19008   | 10.56 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.78  | 132 | 25342   | 11.97 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.95  | 113 | 17490   | 7.21  | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.42  | 128 | 12218   | 12.37 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.14 | 143 | 15554   | 14.43 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.68 | 165 | 26977   | 10.93 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.20 | 131 | 26583   | 10.65 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 14.26 | 166 | 98192   | 12.78 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.56 | 160 | 113415  | 12.77 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 15.03 | 143 | 10558   | 9.01  | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.85 | 176 | 92119   | 13.09 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.95 | 200 | 15708   | 11.47 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.69 | 188 | 147890m | 12.39 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.96 | 212 | 201100  | 12.47 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 25.02 | 264 | 229551  | 13.29 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |        | Ovalue |           |
|----------------------------|-------|-----|--------|--------|-----------|
| 6) Phenol                  | 7.41  | 94  | 3603   | 1.191  | ng/ul# 83 |
| 45) Dimethylphthalate      | 14.30 | 163 | 16810  | 2.145  | ng/ul# 97 |
| 70) Phenanthrene           | 17.64 | 178 | 61109  | 4.704  | ng/ul 96  |
| 77) Fluoranthene           | 19.63 | 202 | 187000 | 11.399 | ng/ul 99  |
| 80) Pyrene                 | 19.99 | 202 | 150517 | 7.665  | ng/ul 98  |
| 83) Benzo(a)anthracene     | 21.85 | 228 | 73559  | 3.729  | ng/ul# 90 |
| 85) Chrysene               | 21.92 | 228 | 82268  | 4.383  | ng/ul 96  |
| 88) Benzo(b)fluoranthene   | 24.18 | 252 | 95324  | 4.499  | ng/ul# 95 |
| 89) Benzo(k)fluoranthene   | 24.24 | 252 | 38375m | 1.838  | ng/ul     |
| 91) Benzo(a)pyrene         | 25.10 | 252 | 67252  | 3.560  | ng/ul 98  |
| 92) Indeno(1,2,3-cd)pyrene | 29.13 | 276 | 45299m | 1.971  | ng/ul     |
| 94) Benzo(g,h,i)perylene   | 30.34 | 276 | 38614m | 2.037  | ng/ul     |

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

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