

Data Path : Z:\HPCHEM1\BNA G\DATA\BG051116\  
 Data File : BG021853.D  
 Acq On : 11 May 2016 21:00  
 Operator : UM/SJ  
 Sample : H2937-20  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 D9X98

Manual Integrations  
 APPROVED

sohil  
 5/12/2016 7:28:02 PM

Quant Time: May 12 08:19:42 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG050916.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu May 12 07:33:29 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.19  | 152  | 85260    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.01 | 136  | 430565   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.81 | 164  | 250773   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.55 | 188  | 500311   | 20.00 | ng/ul | 0.00     |
| 76) Chrysene-d12          | 21.84 | 240  | 482671   | 20.00 | ng/ul | 0.00     |
| 84) Perylene-d12          | 25.19 | 264  | 481133   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.57  | 96  | 5447   | 3.08  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.33  | 99  | 109919 | 17.37 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.50  | 67  | 54454  | 17.36 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.72  | 132 | 101419 | 20.65 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.88  | 113 | 90307  | 16.99 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.35  | 128 | 51860  | 18.32 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.08 | 143 | 54437  | 30.45 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.62 | 165 | 101473 | 20.07 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.14 | 131 | 123802 | 22.73 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 14.20 | 166 | 321699 | 18.26 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.50 | 160 | 438943 | 17.55 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 15.00 | 143 | 44285m | 14.67 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.79 | 176 | 297035 | 16.02 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.91 | 200 | 33654  | 24.28 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.65 | 188 | 398771 | 16.37 | ng/ul | 0.00 |
| 77) Pyrene-d10                 | 19.92 | 212 | 419186 | 18.55 | ng/ul | 0.00 |
| 88) Benzo(a)pyrene-d12         | 24.95 | 264 | 395865 | 17.32 | ng/ul | 0.02 |

## Target Compounds

|                            |       |     |        |      |        | Ovalue |
|----------------------------|-------|-----|--------|------|--------|--------|
| 45) Dimethylphthalate      | 14.25 | 163 | 79431  | 4.43 | ng/ul  | 98     |
| 75) Fluoranthene           | 19.59 | 202 | 86059  | 2.77 | ng/ul# | 96     |
| 78) Pyrene                 | 19.95 | 202 | 73593  | 2.64 | ng/ul  | 98     |
| 81) Benzo(a)anthracene     | 21.82 | 228 | 58611  | 2.18 | ng/ul  | 98     |
| 83) Chrysene               | 21.89 | 228 | 47836  | 1.80 | ng/ul  | 95     |
| 86) Benzo(b)fluoranthene   | 24.12 | 252 | 86384  | 3.14 | ng/ul# | 88     |
| 87) Benzo(k)fluoranthene   | 24.16 | 252 | 36719m | 1.25 | ng/ul  |        |
| 89) Benzo(a)pyrene         | 25.02 | 252 | 56279  | 2.05 | ng/ul# | 87     |
| 90) Indeno(1,2,3-cd)pyrene | 29.01 | 276 | 41360m | 1.27 | ng/ul  |        |
| 92) Benzo(g,h,i)perylene   | 30.21 | 276 | 32139m | 1.20 | ng/ul  |        |

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

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