

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM071116\  
 Data File : BM006403.D  
 Acq On : 11 Jul 2016 20:00  
 Operator : UM/SJ  
 Sample : H3914-12MSD  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 A4TS1MSD

Quant Time: Jul 12 01:00:20 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM070216.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jul 12 00:59:35 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.63  | 152  | 130550   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.41 | 136  | 539076   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.28 | 164  | 292801   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.03 | 188  | 641872   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.23 | 240  | 745494   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.43 | 264  | 800274   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.16  | 96  | 1942    | 0.64  | ng/uL | 0.01 |
| 5) Phenol-d5                   | 6.81  | 99  | 89441   | 7.32  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.97  | 67  | 225952  | 31.33 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.16  | 132 | 237872  | 25.42 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.34  | 113 | 156458  | 15.82 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.79  | 128 | 143965  | 33.51 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.50  | 143 | 158441  | 32.54 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.04 | 165 | 257365  | 29.17 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.56 | 131 | 4342    | 0.47  | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.69 | 166 | 842907  | 34.50 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 13.97 | 160 | 1039630 | 34.36 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.50 | 143 | 29159   | 6.30  | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.28 | 176 | 718946  | 34.13 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.41 | 200 | 113879  | 30.28 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.13 | 188 | 1104604 | 36.07 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.43 | 212 | 1220711 | 34.34 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.30 | 264 | 1366042 | 36.74 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |         |       |        | Qvalue |
|--------------------------------|-------|-----|---------|-------|--------|--------|
| 6) Phenol                      | 6.83  | 94  | 78325   | 6.17  | ng/ul  | 96     |
| 10) 2-Chlorophenol             | 7.20  | 128 | 198428  | 20.49 | ng/ul  | 98     |
| 15) N-Nitroso-di-n-propylamine | 8.43  | 70  | 196870  | 22.86 | ng/ul  | 96     |
| 33) 4-Chloro-3-methylphenol    | 11.71 | 107 | 214747  | 20.15 | ng/ul  | 96     |
| 49) Acenaphthene               | 14.34 | 153 | 556888  | 27.60 | ng/ul  | 98     |
| 52) 4-Nitrophenol              | 14.52 | 109 | 26509   | 5.69  | ng/ul  | 93     |
| 54) 2,4-Dinitrotoluene         | 14.66 | 165 | 196557  | 26.14 | ng/ul  | 98     |
| 68) Pentachlorophenol          | 16.69 | 266 | 112370  | 26.65 | ng/ul  | 97     |
| 77) Pyrene                     | 19.46 | 202 | 1280103 | 27.38 | ng/ul# | 94     |

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

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