

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM073016\  
 Data File : BM006771.D  
 Acq On : 30 Jul 2016 13:42  
 Operator : UM/SJ  
 Sample : H4192-01  
 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DA0D8

Quant Time: Aug 01 05:48:18 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM072016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Jul 30 01:37:26 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.60  | 152  | 180560   | 20.00 | ng/ul | 0.00     |
| 7) Naphthalene-d8         | 10.38 | 136  | 806621   | 20.00 | ng/ul | 0.00     |
| 15) Acenaphthene-d10      | 14.26 | 164  | 453480   | 20.00 | ng/ul | 0.00     |
| 23) Phenanthrene-d10      | 17.01 | 188  | 1020138  | 20.00 | ng/ul | 0.00     |
| 29) Chrysene-d12          | 21.21 | 240  | 1158177  | 20.00 | ng/ul | 0.00     |
| 34) Perylene-d12          | 23.41 | 264  | 1187860  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |       |
|--------------------------------|-------|-----|--------|-------|-------|-------|
| 2) 1,4-Dioxane-d8              | 3.13  | 96  | 4351   | 0.97  | ng/uL | 0.00  |
| 3) Phenol-d5                   | 6.79  | 99  | 172681 | 11.24 | ng/ul | 0.00  |
| 4) Bis-(2-Chloroethyl)ether-d  | 6.95  | 67  | 117271 | 12.32 | ng/ul | 0.00  |
| 5) 2-Chlorophenol-d4           | 7.14  | 132 | 140639 | 11.46 | ng/ul | 0.00  |
| 6) 4-Methylphenol-d8           | 8.32  | 113 | 123594 | 10.37 | ng/ul | 0.00  |
| 8) Nitrobenzene-d5             | 8.76  | 128 | 68536  | 11.22 | ng/ul | 0.00  |
| 9) 2-Nitrophenol-d4            | 9.47  | 143 | 72624  | 11.18 | ng/ul | 0.00  |
| 10) 2,4-Dichlorophenol-d3      | 10.02 | 165 | 135893 | 11.08 | ng/ul | 0.00  |
| 12) 4-Chloroaniline-d4         | 10.53 | 131 | 159806 | 11.83 | ng/ul | 0.00  |
| 16) Dimethylphthalate-d6       | 13.67 | 166 | 443440 | 12.47 | ng/ul | 0.00  |
| 17) Acenaphthylene-d8          | 13.95 | 160 | 566825 | 12.49 | ng/ul | 0.00  |
| 20) 4-Nitrophenol-d4           | 14.49 | 143 | 44806  | 7.26  | ng/ul | 0.00  |
| 21) Fluorene-d10               | 15.25 | 176 | 403084 | 12.68 | ng/ul | 0.00  |
| 24) 4,6-Dinitro-2-methylphenol | 15.40 | 200 | 37856  | 7.15  | ng/ul | 0.00  |
| 26) Anthracene-d10             | 17.10 | 188 | 621700 | 12.83 | ng/ul | 0.00  |
| 30) Pyrene-d10                 | 19.41 | 212 | 688436 | 13.05 | ng/ul | 0.00  |
| 37) Benzo(a)pyrene-d12         | 23.27 | 264 | 685657 | 12.55 | ng/ul | -0.01 |

## Target Compounds

|                          |       |     |        |      |        | Qvalue |
|--------------------------|-------|-----|--------|------|--------|--------|
| 25) Phenanthrene         | 17.05 | 178 | 83692  | 1.46 | ng/ul  | 98     |
| 28) Fluoranthene         | 19.07 | 202 | 131761 | 1.95 | ng/ul  | 98     |
| 31) Pyrene               | 19.44 | 202 | 111336 | 1.59 | ng/ul  | 97     |
| 32) Benzo(a)anthracene   | 21.20 | 228 | 72730  | 1.04 | ng/ul  | 98     |
| 33) Chrysene             | 21.24 | 228 | 82408  | 1.27 | ng/ul  | 96     |
| 35) Benzo(b)fluoranthene | 22.76 | 252 | 122946 | 1.73 | ng/ul# | 98     |
| 38) Benzo(a)pyrene       | 23.32 | 252 | 70205  | 1.01 | ng/ul  | 96     |

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

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