

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100117\  
 Data File : BM011812.D  
 Acq On : 01 Oct 2017 17:08  
 Operator : SJ/JU  
 Sample : I5477-01  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 ET211

Manual Integrations  
 APPROVED

Sohil  
 10/2/2017 5:51:36 PM

Quant Time: Oct 02 04:10:34 2017  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM092717.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Oct 02 02:22:21 2017  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.88  | 152  | 103589   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.69 | 136  | 466286   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.52 | 164  | 280853   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.26 | 188  | 675692   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.43 | 240  | 869075   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.77 | 264  | 977477   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.29  | 96  | 12432m  | 5.35  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.04  | 99  | 67868   | 6.66  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.21  | 67  | 229035  | 31.00 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.41  | 132 | 199000  | 26.25 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.58  | 113 | 132248  | 16.57 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.05  | 128 | 120052  | 33.01 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.77  | 143 | 130673  | 33.69 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.30 | 165 | 226837  | 30.64 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.84 | 131 | 21809m  | 2.66  | ng/ul | 0.02 |
| 43) Dimethylphthalate-d6       | 13.93 | 166 | 897560  | 37.26 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.22 | 160 | 1001205 | 34.37 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.72 | 143 | 21399m  | 5.20  | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.52 | 176 | 747450  | 36.03 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.63 | 200 | 134097  | 29.30 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.36 | 188 | 1252949 | 36.44 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.65 | 212 | 1577961 | 38.04 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.62 | 264 | 1684641 | 36.00 | ng/ul | 0.00 |

Target Compounds Ovalue

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

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