

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100117\  
 Data File : BM011814.D  
 Acq On : 01 Oct 2017 18:21  
 Operator : SJ/JU  
 Sample : I5477-11  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 ET229

Manual Integrations  
 APPROVED

Sohil  
 10/2/2017 5:51:40 PM

Quant Time: Oct 02 04:31:45 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  | 105218   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.69 | 136  | 493189   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.52 | 164  | 294641   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.26 | 188  | 714286   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.43 | 240  | 874998   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.76 | 264  | 999501   | 20.00 | ng/ul | 0.00     |

| System Monitoring Compounds    | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|--------------------------------|-------|------|----------|-------|-------|----------|
| 3) 1,4-Dioxane-d8              | 3.29  | 96   | 12272m   | 5.20  | ng/uL | 0.00     |
| 5) Phenol-d5                   | 7.04  | 99   | 68683    | 6.64  | ng/ul | 0.00     |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.21  | 67   | 225451   | 30.04 | ng/ul | 0.00     |
| 9) 2-Chlorophenol-d4           | 7.41  | 132  | 194029   | 25.20 | ng/ul | 0.00     |
| 13) 4-Methylphenol-d8          | 8.58  | 113  | 136114   | 16.79 | ng/ul | 0.00     |
| 19) Nitrobenzene-d5            | 9.05  | 128  | 116406   | 30.26 | ng/ul | 0.00     |
| 22) 2-Nitrophenol-d4           | 9.77  | 143  | 131390   | 32.02 | ng/ul | 0.00     |
| 26) 2,4-Dichlorophenol-d3      | 10.30 | 165  | 217638   | 27.79 | ng/ul | 0.00     |
| 29) 4-Chloroaniline-d4         | 10.83 | 131  | 32173    | 3.72  | ng/ul | 0.01     |
| 43) Dimethylphthalate-d6       | 13.93 | 166  | 844795   | 33.43 | ng/ul | 0.00     |
| 46) Acenaphthylene-d8          | 14.22 | 160  | 969773   | 31.73 | ng/ul | 0.00     |
| 51) 4-Nitrophenol-d4           | 14.72 | 143  | 20249m   | 4.69  | ng/ul | 0.00     |
| 57) Fluorene-d10               | 15.52 | 176  | 714809   | 32.85 | ng/ul | 0.00     |
| 62) 4,6-Dinitro-2-methylphenol | 15.63 | 200  | 125493   | 25.94 | ng/ul | 0.00     |
| 70) Anthracene-d10             | 17.36 | 188  | 1179543  | 32.45 | ng/ul | 0.00     |
| 76) Pyrene-d10                 | 19.65 | 212  | 1499689  | 35.91 | ng/ul | 0.00     |
| 87) Benzo(a)pyrene-d12         | 23.62 | 264  | 1551269  | 32.42 | ng/ul | 0.00     |

| Target Compounds | Ovalue |
|------------------|--------|
|------------------|--------|

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

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