

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092616\  
 Data File : BM007582.D  
 Acq On : 26 Sep 2016 21:43  
 Operator : UM/SJ  
 Sample : H4866-11  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 E43S3

Manual Integrations  
 APPROVED

umangi  
 9/27/2016 1:09:38 PM

Quant Time: Sep 27 03:08:27 2016  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM092316.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 27 01:59:49 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.78  | 152  | 235090   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.56 | 136  | 1084666  | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.41 | 164  | 862715   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.15 | 188  | 2006286  | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.34 | 240  | 2869591  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.60 | 264  | 2734748  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.30  | 96  | 5936m   | 1.24  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.95  | 99  | 105359  | 5.05  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.12  | 67  | 252419  | 22.50 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.31  | 132 | 273767  | 18.07 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.48  | 113 | 208263  | 11.69 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.93  | 128 | 177650  | 22.91 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.65  | 143 | 205582  | 23.34 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.19 | 165 | 355795  | 20.27 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.71 | 131 | 68265   | 3.70  | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.82 | 166 | 1494513 | 23.82 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.10 | 160 | 1703274 | 23.39 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.63 | 143 | 51046m  | 4.61  | ng/ul | 0.01 |
| 57) Fluorene-d10               | 15.40 | 176 | 1348129 | 24.50 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.53 | 200 | 254309  | 20.77 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.25 | 188 | 2342544 | 25.41 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.54 | 212 | 3013320 | 26.46 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.46 | 264 | 3028129 | 25.02 | ng/ul | 0.00 |

Target Compounds Ovalue

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

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092616\  
Data File : BM007582.D  
Acq On : 26 Sep 2016 21:43  
Operator : UM/SJ  
Sample : H4866-11  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
E43S3

Manual Integrations  
APPROVED

umangi  
9/27/2016 1:09:38 PM

Quant Time: Sep 27 03:08:27 2016  
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM092316.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Sep 27 01:59:49 2016  
Response via : Initial Calibration

