

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM092816\  
 Data File : BM007604.D  
 Acq On : 28 Sep 2016 15:55  
 Operator : UM/SJ  
 Sample : H4955-07  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 A4VL7

Manual Integrations  
 APPROVED

sohil  
 9/29/2016 6:55:19 PM

Quant Time: Sep 28 18:05:40 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM092316.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 28 13:23:56 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.77  | 152  | 183034   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.56 | 136  | 869029   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.40 | 164  | 708768   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.15 | 188  | 1664046  | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.33 | 240  | 2275207  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.60 | 264  | 2154547  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.30  | 96  | 4701    | 1.26  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.95  | 99  | 86217   | 5.31  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.11  | 67  | 231059  | 26.45 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.31  | 132 | 250915  | 21.27 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.48  | 113 | 182566  | 13.16 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.93  | 128 | 169459  | 27.28 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.65  | 143 | 190558  | 27.00 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.19 | 165 | 345697  | 24.58 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.72 | 131 | 27689m  | 1.88  | ng/ul | 0.02 |
| 43) Dimethylphthalate-d6       | 13.82 | 166 | 1613690 | 31.30 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.10 | 160 | 1790687 | 29.93 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.63 | 143 | 41778   | 4.59  | ng/ul | 0.02 |
| 57) Fluorene-d10               | 15.40 | 176 | 1446276 | 31.99 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.53 | 200 | 290570  | 28.62 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.25 | 188 | 2564004 | 33.53 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.54 | 212 | 3298314 | 36.53 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.46 | 264 | 3281913 | 34.43 | ng/ul | 0.00 |

| Target Compounds | Qvalue |
|------------------|--------|
|------------------|--------|

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

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