

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM121417\  
 Data File : BM013421.D  
 Acq On : 15 Dec 2017 03:03  
 Operator : SJ/JU  
 Sample : I6775-18 10X  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 F9A63

Manual Integrations  
 APPROVED

Sohil  
 12/15/2017 6:48:53 PM

Quant Time: Dec 15 07:22:20 2017  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM113017.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Dec 15 05:52:21 2017  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.67  | 152  | 236770   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.45 | 136  | 1046954  | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.32 | 164  | 620834   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.07 | 188  | 1331879  | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.26 | 240  | 1147748  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.48 | 264  | 892222   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |      |       |      |
|--------------------------------|-------|-----|--------|------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.24  | 96  | 1197   | 0.25 | ng/uL | 0.01 |
| 5) Phenol-d5                   | 6.86  | 99  | 35981  | 1.76 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.02  | 67  | 22935  | 1.96 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.21  | 132 | 31105  | 1.94 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.39  | 113 | 31371  | 1.79 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.83  | 128 | 16981  | 2.05 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.55  | 143 | 16659  | 1.87 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.09 | 165 | 29255  | 1.60 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.60 | 131 | 22343  | 1.33 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.73 | 166 | 117658 | 2.13 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.01 | 160 | 135326 | 1.98 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.55 | 143 | 9950   | 1.04 | ng/ul | 0.02 |
| 57) Fluorene-d10               | 15.32 | 176 | 94742  | 1.99 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.45 | 200 | 8790   | 1.05 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.16 | 188 | 152144 | 2.28 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.46 | 212 | 146007 | 2.55 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.34 | 264 | 92836  | 2.04 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |         |        | Qvalue |    |
|----------------------------|-------|-----|---------|--------|--------|----|
| 47) Acenaphthylene         | 14.04 | 152 | 104445  | 1.655  | ng/ul# | 97 |
| 68) Pentachlorophenol      | 16.72 | 266 | 50763   | 5.083  | ng/ul  | 92 |
| 69) Phenanthrene           | 17.11 | 178 | 131692  | 1.740  | ng/ul  | 99 |
| 71) Anthracene             | 17.20 | 178 | 176935  | 2.288  | ng/ul  | 99 |
| 74) Fluoranthene           | 19.13 | 202 | 810886  | 8.745  | ng/ul# | 93 |
| 77) Pyrene                 | 19.49 | 202 | 1011271 | 14.503 | ng/ul# | 93 |
| 80) Benzo(a)anthracene     | 21.24 | 228 | 591044  | 8.116  | ng/ul  | 97 |
| 82) Chrysene               | 21.29 | 228 | 665493  | 9.850  | ng/ul  | 96 |
| 85) Benzo(b)fluoranthene   | 22.81 | 252 | 1538770 | 25.584 | ng/ul# | 98 |
| 86) Benzo(k)fluoranthene   | 22.86 | 252 | 474007m | 8.374  | ng/ul  |    |
| 88) Benzo(a)pyrene         | 23.38 | 252 | 668557  | 12.374 | ng/ul# | 98 |
| 89) Indeno(1,2,3-cd)pyrene | 25.70 | 276 | 488089  | 9.105  | ng/ul# | 95 |
| 91) Benzo(g,h,i)perylene   | 26.38 | 276 | 413438  | 9.777  | ng/ul# | 91 |

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

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