

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012918\  
 Data File : BM013993.D  
 Acq On : 30 Jan 2018 12:34  
 Operator : SJ/JU  
 Sample : J1244-10ME 5X  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 F6B51ME

Manual Integrations  
 APPROVED

Sohil  
 1/30/2018 3:35:11 PM

Quant Time: Jan 30 13:46:39 2018  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jan 30 07:49:34 2018  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.57  | 152  | 77447    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.34 | 136  | 314844   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.22 | 164  | 205594   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 16.97 | 188  | 510031   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.18 | 240  | 642225   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.36 | 264  | 636022   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |      |       |      |
|--------------------------------|-------|-----|---------|------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.13  | 96  | 1291    | 0.66 | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.76  | 99  | 28095m  | 4.73 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.92  | 67  | 15415   | 4.28 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.12  | 132 | 20010   | 4.13 | ng/ul | 0.01 |
| 13) 4-Methylphenol-d8          | 8.30  | 113 | 20669   | 4.52 | ng/ul | 0.01 |
| 19) Nitrobenzene-d5            | 8.75  | 128 | 13101m  | 5.36 | ng/ul | 0.02 |
| 22) 2-Nitrophenol-d4           | 9.45  | 143 | 11999   | 4.65 | ng/ul | 0.01 |
| 26) 2,4-Dichlorophenol-d3      | 10.00 | 165 | 21227   | 4.26 | ng/ul | 0.02 |
| 29) 4-Chloroaniline-d4         | 10.52 | 131 | 25385   | 5.87 | ng/ul | 0.02 |
| 43) Dimethylphthalate-d6       | 13.65 | 166 | 97958   | 5.78 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 13.92 | 160 | 117200  | 5.42 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.52 | 143 | 3570    | 1.30 | ng/ul | 0.06 |
| 57) Fluorene-d10               | 15.22 | 176 | 87458   | 5.87 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.39 | 200 | 6383    | 2.08 | ng/ul | 0.02 |
| 70) Anthracene-d10             | 17.07 | 188 | 143650m | 5.78 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.38 | 212 | 171011  | 5.72 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.23 | 264 | 170701  | 5.84 | ng/ul | 0.00 |

## Target Compounds

|                          |       |     |         |        | Ovalue |     |
|--------------------------|-------|-----|---------|--------|--------|-----|
| 28) Naphthalene          | 10.39 | 128 | 247847  | 15.823 | ng/ul  | 96  |
| 34) 2-Methylnaphthalene  | 12.02 | 142 | 158559  | 13.577 | ng/ul  | 93  |
| 40) 1,1'-Biphenyl        | 13.05 | 154 | 44265   | 2.587  | ng/ul  | 90  |
| 49) Acenaphthene         | 14.29 | 153 | 193514  | 14.133 | ng/ul  | 94  |
| 53) Dibenzofuran         | 14.63 | 168 | 183849  | 8.925  | ng/ul  | 98  |
| 58) Fluorene             | 15.28 | 166 | 179780  | 10.655 | ng/ul  | 97  |
| 69) Phenanthrene         | 17.02 | 178 | 909258  | 32.072 | ng/ul  | 100 |
| 71) Anthracene           | 17.12 | 178 | 110426m | 3.776  | ng/ul  |     |
| 74) Fluoranthene         | 19.04 | 202 | 536628  | 15.857 | ng/ul# | 97  |
| 77) Pyrene               | 19.41 | 202 | 358991  | 9.195  | ng/ul# | 96  |
| 80) Benzo(a)anthracene   | 21.17 | 228 | 79841   | 2.057  | ng/ul  | 99  |
| 82) Chrysene             | 21.22 | 228 | 63459m  | 1.783  | ng/ul  |     |
| 85) Benzo(b)fluoranthene | 22.73 | 252 | 42684   | 1.093  | ng/ul# | 95  |

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

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