

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092117\  
 Data File : BM011678.D  
 Acq On : 22 Sep 2017 05:56  
 Operator : SJ/JU  
 Sample : I5284-05  
 Misc :  
 ALS Vial : 29 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 CB3D0

Manual Integrations  
 APPROVED

Sohil  
 9/22/2017 4:06:40 PM

Quant Time: Sep 22 09:26:13 2017  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM090817MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Sep 22 08:34:40 2017  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.96  | 152  | 578204   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.76 | 136  | 2691813  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.59 | 164  | 1514476  | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.33 | 188  | 3225789  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.50 | 240  | 2603202  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.87 | 264  | 2056734  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.39  | 96  | 14119   | 1.10  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.11  | 99  | 373357  | 6.73  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.28  | 67  | 328534  | 8.78  | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.49  | 132 | 316268  | 7.56  | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.66  | 113 | 293217  | 6.76  | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.12  | 128 | 171160  | 8.38  | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.85  | 143 | 138785  | 6.53  | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.38 | 165 | 276098  | 6.46  | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.90 | 131 | 73112   | 1.55  | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.99 | 166 | 819564  | 6.40  | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.29 | 160 | 1304822 | 8.43  | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.79 | 143 | 27979m  | 1.16  | ng/ul | 0.01 |
| 58) Fluorene-d10               | 15.58 | 176 | 924353  | 8.33  | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.69 | 200 | 30867   | 1.64  | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.43 | 188 | 1399722 | 8.81  | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.72 | 212 | 1544114 | 12.67 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.71 | 264 | 919952  | 9.39  | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |       | Ovalue |    |
|--------------------------------|-------|-----|--------|-------|--------|----|
| 45) Dimethylphthalate          | 14.04 | 163 | 211620 | 1.726 | ng/ul# | 98 |
| 70) Phenanthrene               | 17.37 | 178 | 473788 | 2.635 | ng/ul  | 99 |
| 77) Fluoranthene               | 19.38 | 202 | 875215 | 4.732 | ng/ul# | 91 |
| 80) Pyrene                     | 19.74 | 202 | 760083 | 5.004 | ng/ul# | 91 |
| 83) Benzo(a)anthracene         | 21.48 | 228 | 377098 | 2.631 | ng/ul  | 97 |
| 84) Bis(2-ethylhexyl)phthalate | 21.39 | 149 | 126790 | 1.182 | ng/ul# | 95 |
| 85) Chrysene                   | 21.53 | 228 | 344620 | 2.652 | ng/ul  | 99 |
| 88) Benzo(b)fluoranthene       | 23.14 | 252 | 390083 | 3.153 | ng/ul# | 87 |
| 91) Benzo(a)pyrene             | 23.76 | 252 | 241946 | 2.072 | ng/ul# | 93 |
| 92) Indeno(1,2,3-cd)pyrene     | 26.29 | 276 | 151014 | 1.175 | ng/ul# | 93 |
| 94) Benzo(g,h,i)perylene       | 27.04 | 276 | 112778 | 1.084 | ng/ul# | 88 |

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

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