

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM062217\  
 Data File : BM010664.D  
 Acq On : 22 Jun 2017 17:46  
 Operator : SJ/JU  
 Sample : I3627-01  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 F5A03

Manual Integrations  
 APPROVED

mohammad  
 6/26/2017 8:25:21 AM

Quant Time: Jun 23 05:02:12 2017  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM062017.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 23 01:03:57 2017  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|--------|--------|----------|
| 1) 1,4-Dichlorobenzene-d4      | 7.46  | 152  | 52005    | 20.00  | ng/ul  | 0.00     |
| 18) Naphthalene-d8             | 10.23 | 136  | 230761   | 20.00  | ng/ul  | 0.00     |
| 35) Acenaphthene-d10           | 14.12 | 164  | 162553   | 20.00  | ng/ul  | 0.00     |
| 61) Phenanthrene-d10           | 16.87 | 188  | 415325   | 20.00  | ng/ul  | 0.00     |
| 75) Chrysene-d12               | 21.08 | 240  | 487617   | 20.00  | ng/ul  | 0.00     |
| 83) Perylene-d12               | 23.22 | 264  | 405617   | 20.00  | ng/ul  | 0.00     |
| System Monitoring Compounds    |       |      |          |        |        |          |
| 3) 1,4-Dioxane-d8              | 3.08  | 96   | 4507     | 4.97   | ng/uL  | 0.00     |
| 5) Phenol-d5                   | 6.65  | 99   | 104380   | 21.07  | ng/ul  | 0.00     |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.81  | 67   | 62704    | 22.17  | ng/ul  | 0.00     |
| 9) 2-Chlorophenol-d4           | 7.00  | 132  | 80690    | 23.22  | ng/ul  | 0.00     |
| 13) 4-Methylphenol-d8          | 8.17  | 113  | 92358    | 22.49  | ng/ul  | 0.00     |
| 19) Nitrobenzene-d5            | 8.61  | 128  | 42477    | 25.81  | ng/ul  | 0.00     |
| 22) 2-Nitrophenol-d4           | 9.32  | 143  | 50100    | 26.42  | ng/ul  | 0.00     |
| 26) 2,4-Dichlorophenol-d3      | 9.85  | 165  | 101557   | 25.43  | ng/ul  | 0.00     |
| 29) 4-Chloroaniline-d4         | 10.37 | 131  | 107654   | 25.58  | ng/ul  | 0.00     |
| 43) Dimethylphthalate-d6       | 13.53 | 166  | 400532   | 28.07  | ng/ul  | 0.00     |
| 46) Acenaphthylene-d8          | 13.80 | 160  | 435192   | 26.22  | ng/ul  | 0.00     |
| 51) 4-Nitrophenol-d4           | 14.32 | 143  | 45709    | 18.20  | ng/ul  | 0.00     |
| 57) Fluorene-d10               | 15.12 | 176  | 332852   | 26.98  | ng/ul  | 0.00     |
| 62) 4,6-Dinitro-2-methylphenol | 15.24 | 200  | 54907    | 21.53  | ng/ul  | 0.00     |
| 70) Anthracene-d10             | 16.97 | 188  | 545133m  | 27.18  | ng/ul  | 0.00     |
| 76) Pyrene-d10                 | 19.27 | 212  | 650671   | 28.73  | ng/ul  | 0.00     |
| 87) Benzo(a)pyrene-d12         | 23.09 | 264  | 574241   | 29.29  | ng/ul  | 0.00     |
| Target Compounds               |       |      |          |        |        |          |
| 6) Phenol                      | 6.67  | 94   | 11627    | 2.279  | ng/ul# | 90       |
| 44) Dimethylphthalate          | 13.58 | 163  | 152052   | 10.871 | ng/ul  | 99       |
| 74) Fluoranthene               | 18.94 | 202  | 104581   | 3.675  | ng/ul# | 96       |
| 77) Pyrene                     | 19.30 | 202  | 242207   | 8.600  | ng/ul  | 98       |
| 80) Benzo(a)anthracene         | 21.07 | 228  | 116465   | 4.101  | ng/ul  | 99       |
| 82) Chrysene                   | 21.12 | 228  | 118834m  | 4.694  | ng/ul  |          |
| 85) Benzo(b)fluoranthene       | 22.59 | 252  | 253895   | 9.643  | ng/ul  | 99       |
| 86) Benzo(k)fluoranthene       | 22.62 | 252  | 71585m   | 2.868  | ng/ul  |          |
| 88) Benzo(a)pyrene             | 23.13 | 252  | 114911   | 4.741  | ng/ul# | 98       |
| 89) Indeno(1,2,3-cd)pyrene     | 25.33 | 276  | 69922    | 2.781  | ng/ul  | 99       |
| 91) Benzo(g,h,i)perylene       | 25.97 | 276  | 43405    | 2.169  | ng/ul# | 94       |

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

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