

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM061017\  
 Data File : BM010506.D  
 Acq On : 11 Jun 2017 03:34  
 Operator : SJ/MA  
 Sample : I3558-16  
 Misc :  
 ALS Vial : 31 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 F5C89

Manual Integrations  
 APPROVED

mohammad  
 6/12/2017 3:41:00 PM

Quant Time: Jun 12 08:27:25 2017  
 Quant Method : Z:\HPCHEM1\BNA\_M\Methods\SOM-EPA-BM052717.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 12 07:08:51 2017  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc    | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|---------|--------|----------|
| 1) 1,4-Dichlorobenzene-d4      | 7.89  | 152  | 135897   | 20.00   | ng/ul  | 0.00     |
| 18) Naphthalene-d8             | 10.70 | 136  | 486233   | 20.00   | ng/ul  | 0.00     |
| 35) Acenaphthene-d10           | 14.53 | 164  | 340834   | 20.00   | ng/ul  | 0.00     |
| 61) Phenanthrene-d10           | 17.28 | 188  | 926993   | 20.00   | ng/ul  | 0.00     |
| 75) Chrysene-d12               | 21.44 | 240  | 1449176  | 20.00   | ng/ul  | 0.00     |
| 83) Perylene-d12               | 23.78 | 264  | 1473518  | 20.00   | ng/ul  | 0.00     |
| System Monitoring Compounds    |       |      |          |         |        |          |
| 3) 1,4-Dioxane-d8              | 3.36  | 96   | 17295    | 5.22    | ng/uL  | 0.00     |
| 5) Phenol-d5                   | 7.08  | 99   | 292759   | 28.43   | ng/ul  | 0.00     |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.23  | 67   | 161458   | 27.90   | ng/ul  | 0.00     |
| 9) 2-Chlorophenol-d4           | 7.43  | 132  | 248537   | 30.28   | ng/ul  | 0.00     |
| 13) 4-Methylphenol-d8          | 8.62  | 113  | 249877   | 30.07   | ng/ul  | 0.00     |
| 19) Nitrobenzene-d5            | 9.08  | 128  | 113982   | 32.00   | ng/ul  | 0.00     |
| 22) 2-Nitrophenol-d4           | 9.79  | 143  | 146015   | 35.39   | ng/ul  | 0.00     |
| 26) 2,4-Dichlorophenol-d3      | 10.32 | 165  | 303387   | 35.79   | ng/ul  | 0.00     |
| 29) 4-Chloroaniline-d4         | 10.86 | 131  | 304061   | 37.18   | ng/ul  | 0.00     |
| 43) Dimethylphthalate-d6       | 13.94 | 166  | 1105470  | 39.04   | ng/ul  | 0.00     |
| 46) Acenaphthylene-d8          | 14.23 | 160  | 1186672  | 35.92   | ng/ul  | 0.00     |
| 51) 4-Nitrophenol-d4           | 14.77 | 143  | 158949   | 37.52   | ng/ul  | 0.00     |
| 57) Fluorene-d10               | 15.52 | 176  | 966286   | 38.82   | ng/ul  | 0.00     |
| 62) 4,6-Dinitro-2-methylphenol | 15.65 | 200  | 209255   | 31.87   | ng/ul  | 0.00     |
| 70) Anthracene-d10             | 17.38 | 188  | 1702113m | 37.74   | ng/ul  | 0.00     |
| 76) Pyrene-d10                 | 19.67 | 212  | 2268300  | 37.85   | ng/ul  | 0.00     |
| 87) Benzo(a)pyrene-d12         | 23.63 | 264  | 2511239  | 36.60   | ng/ul  | 0.00     |
| Target Compounds               |       |      |          |         |        |          |
| 6) Phenol                      | 7.10  | 94   | 30299    | 2.870   | ng/ul# | 83       |
| 28) Naphthalene                | 10.75 | 128  | 4911542  | 201.391 | ng/ul  | 99       |
| 34) 2-Methylnaphthalene        | 12.35 | 142  | 875516   | 47.349  | ng/ul  | 96       |
| 40) 1,1'-Biphenyl              | 13.36 | 154  | 229108   | 8.286   | ng/ul  | 99       |
| 44) Dimethylphthalate          | 13.99 | 163  | 465901   | 16.729  | ng/ul  | 100      |
| 47) Acenaphthylene             | 14.26 | 152  | 43049    | 1.333   | ng/ul# | 92       |
| 49) Acenaphthene               | 14.59 | 153  | 734656   | 32.448  | ng/ul  | 99       |
| 53) Dibenzofuran               | 14.93 | 168  | 979763   | 29.265  | ng/ul  | 96       |
| 58) Fluorene                   | 15.58 | 166  | 920022   | 33.501  | ng/ul  | 99       |
| 69) Phenanthrene               | 17.32 | 178  | 4746021  | 95.903  | ng/ul  | 99       |
| 71) Anthracene                 | 17.41 | 178  | 1262175  | 24.431  | ng/ul  | 96       |
| 72) Carbazole                  | 17.70 | 167  | 1184976  | 27.098  | ng/ul  | 97       |
| 74) Fluoranthene               | 19.33 | 202  | 3165745  | 52.020  | ng/ul# | 96       |
| 77) Pyrene                     | 19.70 | 202  | 2002973  | 26.822  | ng/ul# | 94       |
| 80) Benzo(a)anthracene         | 21.43 | 228  | 719675   | 8.806   | ng/ul  | 97       |
| 82) Chrysene                   | 21.48 | 228  | 520701   | 7.049   | ng/ul  | 96       |
| 85) Benzo(b)fluoranthene       | 23.06 | 252  | 338344   | 3.946   | ng/ul  | 98       |
| 86) Benzo(k)fluoranthene       | 23.10 | 252  | 154538m  | 1.887   | ng/ul  |          |
| 88) Benzo(a)pyrene             | 23.67 | 252  | 229460   | 2.700   | ng/ul  | 95       |

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

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM061017\  
Data File : BM010506.D  
Acq On : 11 Jun 2017 03:34  
Operator : SJ/MA  
Sample : I3558-16  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampled :  
F5C89

Manual Integrations  
APPROVED

mohammad  
6/12/2017 3:41:00 PM

Quant Time: Jun 12 08:27:25 2017  
Quant Method : Z:\HPCHEM1\BNA\_M\Methods\SOM-EPA-BM052717.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jun 12 07:08:51 2017  
Response via : Initial Calibration

