

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM060917\  
 Data File : BM010472.D  
 Acq On : 10 Jun 2017 02:15  
 Operator : SJ/MA  
 Sample : I3521-13  
 Misc :  
 ALS Vial : 27 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 F5C00

Manual Integrations  
 APPROVED

mohammad  
 6/12/2017 11:18:31 AM

Quant Time: Jun 10 04:13:20 2017  
 Quant Method : Z:\HPCHEM1\BNA\_M\Methods\SOM-EPA-BM052717.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Jun 10 02:32:05 2017  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.90  | 152  | 284755   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.70 | 136  | 1034386  | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.54 | 164  | 735214   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.29 | 188  | 1801897m | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.45 | 240  | 2114657  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.79 | 264  | 1856704  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.38  | 96  | 27200   | 3.92  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.09  | 99  | 475332  | 22.03 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.25  | 67  | 260945  | 21.52 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.44  | 132 | 411068  | 23.90 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.62  | 113 | 424545  | 24.39 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.09  | 128 | 201131  | 26.54 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.80  | 143 | 242904  | 27.68 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.33 | 165 | 527031  | 29.23 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.87 | 131 | 489839  | 28.16 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.95 | 166 | 1742944 | 28.53 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.24 | 160 | 1928498 | 27.06 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.77 | 143 | 181353  | 19.84 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.53 | 176 | 1530661 | 28.51 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.66 | 200 | 273093  | 21.40 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.39 | 188 | 2456442 | 28.02 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.67 | 212 | 2942515 | 33.65 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.64 | 264 | 2337399 | 27.04 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |         |        | Qvalue |     |
|----------------------------|-------|-----|---------|--------|--------|-----|
| 6) Phenol                  | 7.12  | 94  | 39252   | 1.774  | ng/ul# | 88  |
| 44) Dimethylphthalate      | 14.00 | 163 | 1513738 | 25.198 | ng/ul  | 100 |
| 74) Fluoranthene           | 19.33 | 202 | 792061m | 6.696  | ng/ul  |     |
| 77) Pyrene                 | 19.70 | 202 | 958493  | 8.796  | ng/ul# | 94  |
| 80) Benzo(a)anthracene     | 21.43 | 228 | 582325  | 4.883  | ng/ul  | 98  |
| 82) Chrysene               | 21.49 | 228 | 730536  | 6.777  | ng/ul  | 98  |
| 85) Benzo(b)fluoranthene   | 23.07 | 252 | 1245481 | 11.528 | ng/ul  | 100 |
| 86) Benzo(k)fluoranthene   | 23.11 | 252 | 435597m | 4.220  | ng/ul  |     |
| 88) Benzo(a)pyrene         | 23.69 | 252 | 534724  | 4.993  | ng/ul  | 99  |
| 89) Indeno(1,2,3-cd)pyrene | 26.17 | 276 | 437203  | 3.231  | ng/ul  | 97  |
| 90) Dibenzo(a,h)anthracene | 26.17 | 278 | 124838  | 1.071  | ng/ul# | 85  |
| 91) Benzo(g,h,i)perylene   | 26.91 | 276 | 351632  | 3.155  | ng/ul# | 97  |

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

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