

Data Path : Z:\HPCHEM1\BNA M\DATA\BM072116\  
 Data File : BM006600.D  
 Acq On : 21 Jul 2016 22:50  
 Operator : UM/SJ  
 Sample : H4077-19  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 D9YT3

## Manual Integrations

umangi  
 7/22/2016 6:24:33 PM

Quant Time: Jul 22 03:45:00 2016  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM072016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jul 22 01:56:00 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.61  | 152  | 142697   | 20.00 | ng/ul | 0.00     |
| 7) Naphthalene-d8         | 10.39 | 136  | 658384   | 20.00 | ng/ul | 0.00     |
| 15) Acenaphthene-d10      | 14.26 | 164  | 389785   | 20.00 | ng/ul | 0.00     |
| 23) Phenanthrene-d10      | 17.02 | 188  | 888596   | 20.00 | ng/ul | 0.00     |
| 29) Chrysene-d12          | 21.21 | 240  | 1025945  | 20.00 | ng/ul | 0.00     |
| 34) Perylene-d12          | 23.41 | 264  | 1071700  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 2) 1,4-Dioxane-d8              | 3.13  | 96  | 4325   | 1.23  | ng/uL | 0.00 |
| 3) Phenol-d5                   | 6.79  | 99  | 161163 | 13.28 | ng/ul | 0.00 |
| 4) Bis-(2-Chloroethyl)ether-d  | 6.95  | 67  | 110681 | 14.72 | ng/ul | 0.00 |
| 5) 2-Chlorophenol-d4           | 7.15  | 132 | 133926 | 13.81 | ng/ul | 0.00 |
| 6) 4-Methylphenol-d8           | 8.32  | 113 | 115428 | 12.26 | ng/ul | 0.00 |
| 8) Nitrobenzene-d5             | 8.77  | 128 | 65772  | 13.20 | ng/ul | 0.00 |
| 9) 2-Nitrophenol-d4            | 9.48  | 143 | 70568  | 13.32 | ng/ul | 0.00 |
| 10) 2,4-Dichlorophenol-d3      | 10.02 | 165 | 128988 | 12.88 | ng/ul | 0.00 |
| 12) 4-Chloroaniline-d4         | 10.53 | 131 | 165261 | 14.98 | ng/ul | 0.00 |
| 16) Dimethylphthalate-d6       | 13.67 | 166 | 430220 | 14.08 | ng/ul | 0.00 |
| 17) Acenaphthylene-d8          | 13.95 | 160 | 530502 | 13.60 | ng/ul | 0.00 |
| 20) 4-Nitrophenol-d4           | 14.49 | 143 | 43551m | 8.20  | ng/ul | 0.00 |
| 21) Fluorene-d10               | 15.26 | 176 | 384726 | 14.08 | ng/ul | 0.00 |
| 24) 4,6-Dinitro-2-methylphenol | 15.40 | 200 | 38287  | 8.30  | ng/ul | 0.00 |
| 26) Anthracene-d10             | 17.11 | 188 | 587368 | 13.92 | ng/ul | 0.00 |
| 30) Pyrene-d10                 | 19.42 | 212 | 693153 | 14.83 | ng/ul | 0.00 |
| 37) Benzo(a)pyrene-d12         | 23.27 | 264 | 719043 | 14.59 | ng/ul | 0.00 |

## Target Compounds

|                            |       |     |         |       |        | Ovalue |
|----------------------------|-------|-----|---------|-------|--------|--------|
| 25) Phenanthrene           | 17.06 | 178 | 525010  | 10.51 | ng/ul  | 99     |
| 27) Anthracene             | 17.14 | 178 | 66714   | 1.30  | ng/ul  | 99     |
| 28) Fluoranthene           | 19.08 | 202 | 896229  | 15.27 | ng/ul  | 97     |
| 31) Pyrene                 | 19.44 | 202 | 728737  | 11.73 | ng/ul# | 97     |
| 32) Benzo(a)anthracene     | 21.20 | 228 | 332113  | 5.34  | ng/ul  | 97     |
| 33) Chrysene               | 21.24 | 228 | 382649  | 6.67  | ng/ul  | 98     |
| 35) Benzo(b)fluoranthene   | 22.75 | 252 | 560863  | 8.73  | ng/ul  | 100    |
| 36) Benzo(k)fluoranthene   | 22.79 | 252 | 134863m | 2.17  | ng/ul  |        |
| 38) Benzo(a)pyrene         | 23.31 | 252 | 328034  | 5.25  | ng/ul  | 98     |
| 39) Indeno(1,2,3-cd)pyrene | 25.60 | 276 | 234010  | 3.23  | ng/ul  | 99     |
| 40) Dibenzo(a,h)anthracene | 25.61 | 278 | 63156m  | 1.05  | ng/ul  |        |
| 41) Benzo(g,h,i)perylene   | 26.27 | 276 | 183830  | 2.96  | ng/ul  | 96     |

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

