

Data Path : Z:\HPCHEM1\BNA M\DATA\BM072116\  
 Data File : BM006614.D  
 Acq On : 22 Jul 2016 07:57  
 Operator : UM/SJ  
 Sample : H4076-22  
 Misc :  
 ALS Vial : 31 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 D9YQ5

## Manual Integrations

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 7/22/2016 6:24:59 PM

Quant Time: Jul 22 11:13:41 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 | 7.62  | 152  | 187184   | 20.00 | ng/ul | 0.00     |
| 7) Naphthalene-d8         | 10.39 | 136  | 809909   | 20.00 | ng/ul | 0.00     |
| 15) Acenaphthene-d10      | 14.26 | 164  | 450243   | 20.00 | ng/ul | 0.00     |
| 23) Phenanthrene-d10      | 17.02 | 188  | 1001949  | 20.00 | ng/ul | 0.00     |
| 29) Chrysene-d12          | 21.22 | 240  | 1003149  | 20.00 | ng/ul | 0.00     |
| 34) Perylene-d12          | 23.44 | 264  | 871940   | 20.00 | ng/ul | 0.03     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 2) 1,4-Dioxane-d8              | 3.13  | 96  | 5325   | 1.15  | ng/uL | 0.00 |
| 3) Phenol-d5                   | 6.79  | 99  | 243945 | 15.32 | ng/ul | 0.00 |
| 4) Bis-(2-Chloroethyl)ether-d  | 6.95  | 67  | 157272 | 15.94 | ng/ul | 0.00 |
| 5) 2-Chlorophenol-d4           | 7.15  | 132 | 196396 | 15.43 | ng/ul | 0.00 |
| 6) 4-Methylphenol-d8           | 8.33  | 113 | 138184 | 11.19 | ng/ul | 0.00 |
| 8) Nitrobenzene-d5             | 8.77  | 128 | 95928  | 15.65 | ng/ul | 0.00 |
| 9) 2-Nitrophenol-d4            | 9.49  | 143 | 104980 | 16.10 | ng/ul | 0.00 |
| 10) 2,4-Dichlorophenol-d3      | 10.02 | 165 | 187299 | 15.20 | ng/ul | 0.00 |
| 12) 4-Chloroaniline-d4         | 10.54 | 131 | 131541 | 9.69  | ng/ul | 0.00 |
| 16) Dimethylphthalate-d6       | 13.68 | 166 | 587681 | 16.65 | ng/ul | 0.00 |
| 17) Acenaphthylene-d8          | 13.96 | 160 | 770586 | 17.10 | ng/ul | 0.00 |
| 20) 4-Nitrophenol-d4           | 14.50 | 143 | 82366  | 13.43 | ng/ul | 0.02 |
| 21) Fluorene-d10               | 15.26 | 176 | 542030 | 17.17 | ng/ul | 0.00 |
| 24) 4,6-Dinitro-2-methylphenol | 15.40 | 200 | 36746  | 7.07  | ng/ul | 0.01 |
| 26) Anthracene-d10             | 17.12 | 188 | 805415 | 16.93 | ng/ul | 0.00 |
| 30) Pyrene-d10                 | 19.42 | 212 | 901288 | 19.72 | ng/ul | 0.00 |
| 37) Benzo(a)pyrene-d12         | 23.29 | 264 | 679580 | 16.94 | ng/ul | 0.02 |

## Target Compounds

|                            |       |     |         |      |        | Ovalue |
|----------------------------|-------|-----|---------|------|--------|--------|
| 13) 2-Methylnaphthalene    | 12.06 | 142 | 54725   | 1.79 | ng/ul  | 99     |
| 14) 1-Methylnaphthalene    | 12.29 | 142 | 33994   | 1.16 | ng/ul  | 89     |
| 25) Phenanthrene           | 17.06 | 178 | 200267  | 3.55 | ng/ul  | 97     |
| 28) Fluoranthene           | 19.08 | 202 | 533462  | 8.06 | ng/ul  | 98     |
| 31) Pyrene                 | 19.45 | 202 | 477142  | 7.86 | ng/ul  | 97     |
| 32) Benzo(a)anthracene     | 21.20 | 228 | 262742  | 4.32 | ng/ul  | 93     |
| 33) Chrysene               | 21.26 | 228 | 298837m | 5.33 | ng/ul  |        |
| 35) Benzo(b)fluoranthene   | 22.77 | 252 | 359669  | 6.88 | ng/ul# | 72     |
| 36) Benzo(k)fluoranthene   | 22.81 | 252 | 96309m  | 1.91 | ng/ul  |        |
| 38) Benzo(a)pyrene         | 23.34 | 252 | 223375  | 4.39 | ng/ul# | 76     |
| 39) Indeno(1,2,3-cd)pyrene | 25.65 | 276 | 207838  | 3.53 | ng/ul# | 90     |
| 40) Dibenzo(a,h)anthracene | 25.65 | 278 | 59633   | 1.22 | ng/ul# | 59     |
| 41) Benzo(g,h,i)perylene   | 26.32 | 276 | 195967  | 3.88 | ng/ul# | 90     |

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

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