

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052219\  
 Data File : BM020498.D  
 Acq On : 22 May 2019 19:09  
 Operator : JU/SJ  
 Sample : K2814-01  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 A4246

Manual Integrations  
 APPROVED

mohammad

Quant Time: May 23 08:52:55 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051619MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu May 23 07:52:10 2019  
 Response via : Initial Calibration

5/23/2019 2:22:12 PM

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.67  | 152  | 67831    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.47 | 136  | 267331   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.34 | 164  | 165271   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.09 | 188  | 401570   | 20.00 | ng/ul | 0.00     |
| 77) Chrysene-d12          | 21.29 | 240  | 426664   | 20.00 | ng/ul | 0.00     |
| 85) Perylene-d12          | 23.54 | 264  | 489425   | 20.00 | ng/ul | 0.00     |

#### System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.15  | 96  | 5679   | 3.09  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.85  | 99  | 82588  | 15.26 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.02  | 67  | 56551  | 16.44 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.21  | 132 | 63644  | 16.04 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.39  | 113 | 53084  | 13.47 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.85  | 128 | 29847  | 17.16 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.56  | 143 | 34919  | 18.20 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.10 | 165 | 66758  | 16.46 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.63 | 131 | 49116  | 11.72 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.75 | 166 | 216284 | 18.19 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.03 | 160 | 259859 | 17.47 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.57 | 143 | 14609m | 8.46  | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.33 | 176 | 193003 | 18.25 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.48 | 200 | 18722  | 8.18  | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.19 | 188 | 294088 | 17.07 | ng/ul | 0.00 |
| 78) Pyrene-d10                 | 19.49 | 212 | 379092 | 18.45 | ng/ul | 0.00 |
| 89) Benzo(a)pyrene-d12         | 23.40 | 264 | 402601 | 18.10 | ng/ul | 0.00 |

#### Target Compounds

|                            |       |     |         |              | Ovalue |
|----------------------------|-------|-----|---------|--------------|--------|
| 4) Benzaldehyde            | 6.83  | 77  | 6650    | 1.435 ng/ul  | 87     |
| 44) Dimethylphthalate      | 13.80 | 163 | 76454   | 6.500 ng/ul# | 98     |
| 69) Phenanthrene           | 17.13 | 178 | 142209  | 7.432 ng/ul  | 99     |
| 71) Anthracene             | 17.23 | 178 | 24154   | 1.241 ng/ul  | 98     |
| 76) Fluoranthene           | 19.16 | 202 | 256804  | 10.625 ng/ul | 99     |
| 79) Pyrene                 | 19.52 | 202 | 293486  | 11.460 ng/ul | 99     |
| 82) Benzo(a)anthracene     | 21.27 | 228 | 120349m | 4.686 ng/ul  |        |
| 84) Chrysene               | 21.33 | 228 | 105051  | 4.281 ng/ul  | 96     |
| 87) Benzo(b)fluoranthene   | 22.86 | 252 | 148989  | 5.334 ng/ul  | 95     |
| 88) Benzo(k)fluoranthene   | 22.90 | 252 | 45124m  | 1.645 ng/ul  |        |
| 90) Benzo(a)pyrene         | 23.44 | 252 | 121592  | 4.626 ng/ul  | 98     |
| 91) Indeno(1,2,3-cd)pyrene | 25.81 | 276 | 86224   | 2.879 ng/ul  | 97     |
| 93) Benzo(g,h,i)perylene   | 26.52 | 276 | 82227   | 3.316 ng/ul  | 96     |

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

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