

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021720\  
 Data File : BM024917.D  
 Acq On : 17 Feb 2020 21:42  
 Operator : CG/JU  
 Sample : L1486-04MS  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DBCT5MS

Manual Integrations  
 APPROVED

mohammad  
 2/18/2020 3:17:56 PM

Quant Time: Feb 18 08:19:37 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM020620MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Feb 15 06:02:17 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.39  | 152  | 276017   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.14 | 136  | 989240   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.04 | 164  | 588841   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.79 | 188  | 1228117  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.01 | 240  | 1253938  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.10 | 264  | 1338504  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.03  | 96  | 35661   | 6.21  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.59  | 99  | 120312  | 6.56  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.75  | 67  | 358906  | 34.66 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 6.93  | 132 | 414290  | 25.22 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.10  | 113 | 226446  | 14.64 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.53  | 128 | 246695  | 34.49 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.24  | 143 | 276902  | 32.76 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.77  | 165 | 469821  | 27.55 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.30 | 131 | 24466   | 1.53  | ng/ul | 0.01 |
| 44) Dimethylphthalate-d6       | 13.46 | 166 | 1685301 | 37.92 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.72 | 160 | 1895208 | 36.51 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.27 | 143 | 39746m  | 5.50  | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.04 | 176 | 1431062 | 36.36 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.17 | 200 | 261270  | 32.64 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 16.89 | 188 | 2328070 | 41.53 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.20 | 212 | 2618947 | 41.75 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 22.97 | 264 | 2856290 | 42.49 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |         |        | Ovalue |     |
|--------------------------------|-------|-----|---------|--------|--------|-----|
| 6) Phenol                      | 6.62  | 94  | 132847  | 6.883  | ng/ul  | 93  |
| 10) 2-Chlorophenol             | 6.96  | 128 | 408690  | 24.402 | ng/ul  | 99  |
| 15) N-Nitroso-di-n-propylamine | 8.19  | 70  | 370977  | 32.192 | ng/ul# | 84  |
| 33) 4-Chloro-3-methylphenol    | 11.45 | 107 | 410753  | 26.289 | ng/ul  | 97  |
| 45) Dimethylphthalate          | 13.51 | 163 | 52021   | 1.122  | ng/ul  | 99  |
| 50) Acenaphthene               | 14.10 | 153 | 1198753 | 33.013 | ng/ul  | 98  |
| 53) 4-Nitrophenol              | 14.29 | 109 | 31909   | 5.409  | ng/ul  | 93  |
| 55) 2,4-Dinitrotoluene         | 14.42 | 165 | 481811  | 36.005 | ng/ul  | 97  |
| 69) Pentachlorophenol          | 16.46 | 266 | 394187  | 36.222 | ng/ul  | 100 |
| 80) Pyrene                     | 19.23 | 202 | 3210623 | 38.681 | ng/ul# | 95  |

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

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