

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022120\  
 Data File : BM024994.D  
 Acq On : 21 Feb 2020 19:34  
 Operator : CG/JU  
 Sample : L1582-03MSD  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 DBDJ5MSD

Manual Integrations  
 APPROVED

mohammad  
 2/24/2020 3:34:55 PM

Quant Time: Feb 22 01:38:22 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM020620MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Feb 22 01:15:56 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.34  | 152  | 150861   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.09 | 136  | 592191   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.00 | 164  | 389870   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.75 | 188  | 888231   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 20.97 | 240  | 938341   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.04 | 264  | 1165671  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 2.95  | 96  | 24852   | 7.92  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.55  | 99  | 63086   | 6.29  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.69  | 67  | 191305  | 33.80 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 6.88  | 132 | 211973  | 23.61 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.06  | 113 | 121266  | 14.34 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.48  | 128 | 123852  | 28.92 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.19  | 143 | 138894  | 27.45 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.73  | 165 | 255795  | 25.06 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.26 | 131 | 15434m  | 1.62  | ng/ul | 0.02 |
| 44) Dimethylphthalate-d6       | 13.42 | 166 | 960515  | 32.64 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.68 | 160 | 1103483 | 32.11 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.24 | 143 | 15924m  | 3.33  | ng/ul | 0.01 |
| 58) Fluorene-d10               | 15.00 | 176 | 851326  | 32.67 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.14 | 200 | 143806  | 24.84 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 16.85 | 188 | 1381127 | 34.06 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.16 | 212 | 1710270 | 36.44 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 22.91 | 264 | 2020713 | 34.52 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |         |        | Ovalue |    |
|--------------------------------|-------|-----|---------|--------|--------|----|
| 6) Phenol                      | 6.57  | 94  | 83639   | 7.929  | ng/ul  | 93 |
| 10) 2-Chlorophenol             | 6.91  | 128 | 220186  | 24.053 | ng/ul  | 97 |
| 15) N-Nitroso-di-n-propylamine | 8.15  | 70  | 237229  | 37.664 | ng/ul# | 82 |
| 33) 4-Chloro-3-methylphenol    | 11.40 | 107 | 247617  | 26.474 | ng/ul  | 96 |
| 45) Dimethylphthalate          | 13.47 | 163 | 37753   | 1.230  | ng/ul# | 97 |
| 50) Acenaphthene               | 14.06 | 153 | 759266  | 31.582 | ng/ul  | 99 |
| 53) 4-Nitrophenol              | 14.26 | 109 | 13400m  | 3.431  | ng/ul  |    |
| 55) 2,4-Dinitrotoluene         | 14.38 | 165 | 262339  | 29.610 | ng/ul  | 96 |
| 69) Pentachlorophenol          | 16.42 | 266 | 232358  | 29.522 | ng/ul  | 98 |
| 80) Pyrene                     | 19.19 | 202 | 2181402 | 35.121 | ng/ul  | 96 |

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

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