

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082619\  
 Data File : BM022363.D  
 Acq On : 27 Aug 2019 11:54  
 Operator : HP/JU  
 Sample : K4414-02MS  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleID :  
 ETJQ7MS

Manual Integrations  
 APPROVED

mohammad  
 8/28/2019 11:32:16 PM

Quant Time: Aug 27 18:16:09 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM082219MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Aug 27 17:04:34 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.54  | 152  | 37440    | 20.00 | ng/ul | -0.01    |
| 18) Naphthalene-d8        | 10.31 | 136  | 170721   | 20.00 | ng/ul | -0.02    |
| 35) Acenaphthene-d10      | 14.20 | 164  | 122164   | 20.00 | ng/ul | -0.01    |
| 61) Phenanthrene-d10      | 16.96 | 188  | 276402   | 20.00 | ng/ul | -0.01    |
| 77) Chrysene-d12          | 21.17 | 240  | 232710   | 20.00 | ng/ul | 0.00     |
| 85) Perylene-d12          | 23.37 | 264  | 234954   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |       |
|--------------------------------|-------|-----|--------|-------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.13  | 96  | 3377   | 4.04  | ng/uL | 0.00  |
| 5) Phenol-d5                   | 6.74  | 99  | 17090  | 5.29  | ng/ul | 0.00  |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.90  | 67  | 41288  | 21.97 | ng/ul | -0.01 |
| 9) 2-Chlorophenol-d4           | 7.08  | 132 | 50671  | 19.42 | ng/ul | 0.00  |
| 13) 4-Methylphenol-d8          | 8.27  | 113 | 33410  | 11.93 | ng/ul | 0.00  |
| 19) Nitrobenzene-d5            | 8.71  | 128 | 30888  | 24.27 | ng/ul | -0.01 |
| 22) 2-Nitrophenol-d4           | 9.42  | 143 | 36881  | 24.89 | ng/ul | -0.01 |
| 26) 2,4-Dichlorophenol-d3      | 9.96  | 165 | 68260  | 22.63 | ng/ul | 0.00  |
| 29) 4-Chloroaniline-d4         | 10.52 | 131 | 3218   | 0.88  | ng/ul | 0.03  |
| 43) Dimethylphthalate-d6       | 13.63 | 166 | 266716 | 24.79 | ng/ul | 0.00  |
| 46) Acenaphthylene-d8          | 13.89 | 160 | 283308 | 24.45 | ng/ul | -0.01 |
| 51) 4-Nitrophenol-d4           | 14.49 | 143 | 6091m  | 3.82  | ng/ul | 0.02  |
| 57) Fluorene-d10               | 15.20 | 176 | 230783 | 25.87 | ng/ul | 0.00  |
| 62) 4,6-Dinitro-2-methylphenol | 15.37 | 200 | 38793  | 18.19 | ng/ul | 0.00  |
| 70) Anthracene-d10             | 17.06 | 188 | 405504 | 27.71 | ng/ul | -0.01 |
| 78) Pyrene-d10                 | 19.37 | 212 | 474179 | 39.21 | ng/ul | -0.01 |
| 89) Benzo(a)pyrene-d12         | 23.23 | 264 | 352685 | 27.32 | ng/ul | 0.00  |

## Target Compounds

|                                |       |     |        |        | Ovalue    |
|--------------------------------|-------|-----|--------|--------|-----------|
| 6) Phenol                      | 6.77  | 94  | 20397  | 5.895  | ng/ul 91  |
| 10) 2-Chlorophenol             | 7.11  | 128 | 48349  | 17.853 | ng/ul 97  |
| 15) N-Nitroso-di-n-propylamine | 8.35  | 70  | 51413  | 20.979 | ng/ul 98  |
| 28) Naphthalene                | 10.36 | 128 | 503459 | 53.437 | ng/ul 99  |
| 33) 4-Chloro-3-methylphenol    | 11.64 | 107 | 66442  | 19.694 | ng/ul 98  |
| 44) Dimethylphthalate          | 13.67 | 163 | 14087  | 1.287  | ng/ul# 94 |
| 47) Acenaphthylene             | 13.92 | 152 | 26349  | 2.193  | ng/ul 97  |
| 49) Acenaphthene               | 14.26 | 153 | 202720 | 23.763 | ng/ul 99  |
| 52) 4-Nitrophenol              | 14.51 | 109 | 8775   | 3.313  | ng/ul 88  |
| 54) 2,4-Dinitrotoluene         | 14.62 | 165 | 73872  | 22.734 | ng/ul 86  |
| 68) Pentachlorophenol          | 16.62 | 266 | 55043  | 20.669 | ng/ul 98  |
| 69) Phenanthrene               | 17.00 | 178 | 28734  | 1.738  | ng/ul 98  |
| 79) Pyrene                     | 19.40 | 202 | 567498 | 37.676 | ng/ul# 94 |

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

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