

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\  
 Data File : BM024504.D  
 Acq On : 17 Jan 2020 20:23  
 Operator : JU  
 Sample : L1139-05  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C0D51

Manual Integrations  
 APPROVED

mohammad  
 1/21/2020 7:59:32 AM

Quant Time: Jan 18 00:26:38 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Jan 18 00:06:51 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.51  | 152  | 257514   | 20.00 | ng/ul | 0.05     |
| 18) Naphthalene-d8        | 10.25 | 136  | 876169   | 20.00 | ng/ul | 0.02     |
| 36) Acenaphthene-d10      | 14.10 | 164  | 634041   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.86 | 188  | 1502410  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.06 | 240  | 1591259  | 20.00 | ng/ul | -0.01    |
| 86) Perylene-d12          | 23.18 | 264  | 1712892  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.06  | 96  | 20623   | 3.91  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.65  | 99  | 92438   | 4.84  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.82  | 67  | 258201  | 24.35 | ng/ul | 0.01 |
| 9) 2-Chlorophenol-d4           | 7.01  | 132 | 346646  | 21.26 | ng/ul | 0.01 |
| 13) 4-Methylphenol-d8          | 8.17  | 113 | 200836m | 11.41 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.60  | 128 | 229891  | 37.34 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.32  | 143 | 271662  | 40.75 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.86  | 165 | 467992  | 31.19 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.37 | 131 | 29830   | 1.79  | ng/ul | 0.01 |
| 44) Dimethylphthalate-d6       | 13.53 | 166 | 1893249 | 38.11 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.79 | 160 | 2166360 | 37.89 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.31 | 143 | 57840   | 6.98  | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.11 | 176 | 1655246 | 37.85 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.23 | 200 | 313655  | 36.57 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 16.95 | 188 | 2771687 | 39.67 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.26 | 212 | 3326319 | 39.11 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.04 | 264 | 3543633 | 39.89 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |         | Ovalue  |           |
|--------------------------------|-------|-----|---------|---------|-----------|
| 20) Nitrobenzene               | 8.65  | 77  | 2104832 | 141.278 | ng/ul 97  |
| 23) 2-Nitrophenol              | 9.35  | 139 | 19766   | 2.711   | ng/ul# 87 |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.27 | 216 | 881974  | 42.281  | ng/ul 99  |
| 45) Dimethylphthalate          | 13.57 | 163 | 118494  | 2.346   | ng/ul 100 |
| 73) 1,2,3,4-Tetrachlorobenzene | 12.90 | 216 | 3336468 | 156.174 | ng/uL 98  |
| 74) Pentachlorobenzene         | 14.43 | 250 | 83545   | 3.694   | ng/uL 98  |

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

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