

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121219\  
 Data File : BM024098.D  
 Acq On : 13 Dec 2019 19:00  
 Operator : JU  
 Sample : K6167-20  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BFR92

Manual Integrations  
 APPROVED

mohammad  
 12/16/2019 12:26:08 PM

Quant Time: Dec 14 01:12:57 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM112719MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Dec 14 00:26:31 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.47  | 152  | 331642   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.24 | 136  | 1315561  | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.13 | 164  | 769597   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.89 | 188  | 1554533  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.09 | 240  | 1493689  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.23 | 264  | 1828677  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.03  | 96  | 46474   | 6.23  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.67  | 99  | 169164  | 5.82  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.82  | 67  | 486138  | 28.76 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.01  | 132 | 515205  | 22.93 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.19  | 113 | 324417  | 14.15 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.62  | 128 | 293534  | 28.75 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.34  | 143 | 332839  | 28.68 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.87  | 165 | 585945  | 26.66 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.41 | 131 | 25728m  | 1.07  | ng/ul | 0.02 |
| 44) Dimethylphthalate-d6       | 13.55 | 166 | 1855197 | 31.02 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.82 | 160 | 2247973 | 32.08 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.39 | 143 | 58086   | 5.78  | ng/ul | 0.02 |
| 58) Fluorene-d10               | 15.13 | 176 | 1618646 | 32.17 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.28 | 200 | 279826  | 28.67 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 16.98 | 188 | 2445861 | 32.94 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.29 | 212 | 2610154 | 33.82 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.10 | 264 | 3238919 | 33.51 | ng/ul | 0.00 |

## Target Compounds

|                       |       |     |        |       | Ovalue   |
|-----------------------|-------|-----|--------|-------|----------|
| 2) 1,4-Dioxane        | 3.06  | 88  | 45933  | 5.746 | ng/uL 97 |
| 6) Phenol             | 6.70  | 94  | 45393  | 1.518 | ng/ul 94 |
| 45) Dimethylphthalate | 13.60 | 163 | 113852 | 1.945 | ng/ul 97 |

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121219\  
Data File : BM024098.D  
Acq On : 13 Dec 2019 19:00  
Operator : JU  
Sample : K6167-20  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BFR92

Quant Time: Dec 14 01:12:57 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM112719MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sat Dec 14 00:26:31 2019  
Response via : Initial Calibration

Manual Integrations  
APPROVED

mohammad  
12/16/2019 12:26:08 PM

