

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN013020\  
 Data File : BN009485.D  
 Acq On : 30 Jan 2020 21:29  
 Operator : CG/JU  
 Sample : L1300-04  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 C5AN4

Quant Time: Jan 31 01:39:12 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN012220MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jan 22 15:13:20 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.46  | 152  | 177059   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.20 | 136  | 743153   | 20.00 | ng/ul | -0.01    |
| 35) Acenaphthene-d10      | 14.09 | 164  | 466537   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 16.85 | 188  | 972491   | 20.00 | ng/ul | 0.00     |
| 77) Chrysene-d12          | 21.05 | 240  | 898525   | 20.00 | ng/ul | 0.00     |
| 85) Perylene-d12          | 23.17 | 264  | 939904   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |       |
|--------------------------------|-------|-----|--------|-------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.09  | 96  | 7868   | 1.84  | ng/uL | 0.00  |
| 5) Phenol-d5                   | 6.65  | 99  | 152847 | 9.73  | ng/ul | 0.00  |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.81  | 67  | 110156 | 9.88  | ng/ul | 0.00  |
| 9) 2-Chlorophenol-d4           | 7.00  | 132 | 123363 | 10.76 | ng/ul | 0.00  |
| 13) 4-Methylphenol-d8          | 8.16  | 113 | 112883 | 9.03  | ng/ul | 0.00  |
| 19) Nitrobenzene-d5            | 8.60  | 128 | 57675  | 10.67 | ng/ul | 0.00  |
| 22) 2-Nitrophenol-d4           | 9.31  | 143 | 62183  | 10.97 | ng/ul | 0.00  |
| 26) 2,4-Dichlorophenol-d3      | 9.85  | 165 | 116918 | 10.40 | ng/ul | 0.00  |
| 29) 4-Chloroaniline-d4         | 10.36 | 131 | 134533 | 9.91  | ng/ul | 0.00  |
| 43) Dimethylphthalate-d6       | 13.52 | 166 | 375901 | 10.89 | ng/ul | 0.00  |
| 46) Acenaphthylene-d8          | 13.77 | 160 | 438191 | 10.69 | ng/ul | -0.01 |
| 51) 4-Nitrophenol-d4           | 14.33 | 143 | 41667  | 6.48  | ng/ul | 0.00  |
| 57) Fluorene-d10               | 15.09 | 176 | 339658 | 11.21 | ng/ul | 0.00  |
| 62) 4,6-Dinitro-2-methylphenol | 15.23 | 200 | 41885  | 7.11  | ng/ul | 0.00  |
| 70) Anthracene-d10             | 16.94 | 188 | 530583 | 11.79 | ng/ul | 0.00  |
| 78) Pyrene-d10                 | 19.25 | 212 | 654715 | 13.54 | ng/ul | 0.00  |
| 89) Benzo(a)pyrene-d12         | 23.03 | 264 | 643844 | 13.57 | ng/ul | 0.00  |

## Target Compounds

|                          |       |     |        |       | Ovalue |    |
|--------------------------|-------|-----|--------|-------|--------|----|
| 44) Dimethylphthalate    | 13.56 | 163 | 41146  | 1.158 | ng/ul# | 96 |
| 76) Fluoranthene         | 18.91 | 202 | 118970 | 1.891 | ng/ul  | 98 |
| 79) Pyrene               | 19.27 | 202 | 106505 | 1.637 | ng/ul  | 98 |
| 82) Benzo(a)anthracene   | 21.03 | 228 | 66023  | 1.079 | ng/ul  | 97 |
| 84) Chrysene             | 21.09 | 228 | 77172  | 1.295 | ng/ul  | 99 |
| 87) Benzo(b)fluoranthene | 22.54 | 252 | 91192  | 1.534 | ng/ul# | 93 |
| 90) Benzo(a)pyrene       | 23.07 | 252 | 59145  | 1.109 | ng/ul# | 95 |

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

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