

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN042321\  
 Data File : BN014388.D  
 Acq On : 23 Apr 2021 21:38  
 Operator : CG/JU  
 Sample : M2065-17  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 BG570

Quant Time: Apr 24 00:35:25 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN041521.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Apr 23 12:06:34 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.68  | 152  | 149864   | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.46 | 136  | 629781   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.32 | 164  | 377722   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.06 | 188  | 772205   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.26 | 240  | 750419   | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.49 | 264  | 786319   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.21  | 96  | 13931   | 3.54  | ng/uL | 0.00 |
| 4) Pyridine-d5                 | 3.61  | 84  | 153431  | 14.71 | ng/ul | 0.00 |
| 7) Phenol-d5                   | 6.85  | 99  | 312380  | 24.93 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)ether-d  | 7.02  | 67  | 197020  | 26.26 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4          | 7.22  | 132 | 270828  | 27.48 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8          | 8.38  | 113 | 243852  | 25.68 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5            | 8.83  | 128 | 142300  | 31.27 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4           | 9.55  | 143 | 118091  | 26.50 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3      | 10.08 | 165 | 243554  | 25.37 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4         | 10.59 | 131 | 326555  | 22.93 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6       | 13.73 | 166 | 783827  | 28.64 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8          | 14.01 | 160 | 1071709 | 32.56 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4           | 14.52 | 143 | 77671   | 14.63 | ng/ul | 0.00 |
| 60) Fluorene-d10               | 15.32 | 176 | 728459  | 30.11 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylphenol | 15.43 | 200 | 40736   | 9.86  | ng/ul | 0.00 |
| 73) Anthracene-d10             | 17.16 | 188 | 1120637 | 31.09 | ng/ul | 0.00 |
| 81) Pyrene-d10                 | 19.47 | 212 | 1255810 | 34.47 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12         | 23.34 | 264 | 1316007 | 31.55 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |       | Ovalue    |
|--------------------------------|-------|-----|--------|-------|-----------|
| 8) Phenol                      | 6.88  | 94  | 20225  | 1.502 | ng/ul 96  |
| 18) 4-Methylphenol             | 8.44  | 108 | 11876  | 1.123 | ng/ul 97  |
| 47) Dimethylphthalate          | 13.77 | 163 | 124348 | 4.377 | ng/ul 96  |
| 67) N-Nitrosodiphenylamine     | 15.59 | 169 | 24150  | 1.005 | ng/ul# 47 |
| 72) Phenanthrene               | 17.11 | 178 | 69560  | 1.543 | ng/ul# 95 |
| 80) Fluoranthene               | 19.13 | 202 | 112920 | 2.453 | ng/ul 98  |
| 82) Pyrene                     | 19.49 | 202 | 125648 | 2.563 | ng/ul 99  |
| 85) Benzo(a)anthracene         | 21.24 | 228 | 51097  | 1.054 | ng/ul# 91 |
| 86) Bis(2-ethylhexyl)phthalate | 21.19 | 149 | 153199 | 5.841 | ng/ul 97  |
| 87) Chrysene                   | 21.30 | 228 | 60331  | 1.284 | ng/ul# 93 |
| 90) Benzo(b)fluoranthene       | 22.82 | 252 | 62611  | 1.190 | ng/ul# 65 |

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

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