

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071118\  
 Data File : BN002008.D  
 Acq On : 11 Jul 2018 13:40  
 Operator : JU/SJ  
 Sample : J3775-12  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 DAZQ1

Manual Integrations  
 APPROVED

Sohil  
 7/12/2018 3:03:38 PM

Quant Time: Jul 11 18:47:53 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN061918.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jul 10 01:03:06 2018  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.70  | 152  | 454640   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.48 | 136  | 2137717  | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.33 | 164  | 1282172  | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.08 | 188  | 2900619  | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.30 | 240  | 2878180  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.54 | 264  | 2593238  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.26  | 96  | 50532   | 4.83  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.88  | 99  | 1217281 | 29.59 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.05  | 67  | 764406  | 30.11 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.25  | 132 | 960424  | 29.44 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.41  | 113 | 961027  | 29.16 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.85  | 128 | 495434  | 31.21 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.57  | 143 | 568386  | 34.59 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.10 | 165 | 1023582 | 29.18 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.62 | 131 | 1032408 | 28.43 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.75 | 166 | 3187703 | 29.44 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.03 | 160 | 4042926 | 30.24 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.54 | 143 | 536044  | 26.87 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.33 | 176 | 2860474 | 30.14 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.45 | 200 | 457382  | 28.95 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.18 | 188 | 4284914 | 30.58 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.49 | 212 | 4746345 | 32.71 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.40 | 264 | 4272000 | 30.29 | ng/ul | 0.01 |

## Target Compounds

|                                |       |     |         |       | Ovalue |    |
|--------------------------------|-------|-----|---------|-------|--------|----|
| 6) Phenol                      | 6.91  | 94  | 65082   | 1.569 | ng/ul  | 96 |
| 44) Dimethylphthalate          | 13.79 | 163 | 793810  | 7.459 | ng/ul  | 99 |
| 69) Phenanthrene               | 17.12 | 178 | 323589  | 1.980 | ng/ul  | 99 |
| 73) Di-n-butylphthalate        | 18.06 | 149 | 389822  | 2.228 | ng/ul  | 99 |
| 74) Fluoranthene               | 19.15 | 202 | 1081404 | 6.296 | ng/ul# | 90 |
| 77) Pyrene                     | 19.52 | 202 | 943052  | 5.236 | ng/ul# | 87 |
| 78) Butylbenzylphthalate       | 20.43 | 149 | 321413  | 4.182 | ng/ul  | 91 |
| 80) Benzo(a)anthracene         | 21.28 | 228 | 604967  | 3.326 | ng/ul  | 98 |
| 81) Bis(2-ethylhexyl)phthalate | 21.23 | 149 | 184421  | 1.609 | ng/ul  | 99 |
| 82) Chrysene                   | 21.33 | 228 | 625465  | 3.696 | ng/ul  | 96 |
| 85) Benzo(b)fluoranthene       | 22.87 | 252 | 972354  | 5.937 | ng/ul# | 90 |
| 86) Benzo(k)fluoranthene       | 22.91 | 252 | 271419m | 1.709 | ng/ul  |    |
| 88) Benzo(a)pyrene             | 23.45 | 252 | 562050  | 3.672 | ng/ul# | 90 |
| 89) Indeno(1,2,3-cd)pyrene     | 25.77 | 276 | 416145  | 2.555 | ng/ul# | 82 |
| 91) Benzo(g,h,i)perylene       | 26.46 | 276 | 370592  | 2.758 | ng/ul# | 91 |

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

