

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100618\  
 Data File : BG037207.D  
 Acq On : 6 Oct 2018 2:51  
 Operator : JU/SJ  
 Sample : SSTDCCC040  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTDCCC040EC

Quant Time: Oct 06 05:33:25 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Sep 24 10:41:23 2018  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.03  | 152  | 28537    | 20.00 | ng    | -0.01    |
| 21) Naphthalene-d8        | 10.84 | 136  | 117458   | 20.00 | ng    | -0.02    |
| 38) Acenaphthene-d10      | 14.66 | 164  | 79677    | 20.00 | ng    | -0.02    |
| 63) Phenanthrene-d10      | 17.40 | 188  | 233384   | 20.00 | ng    | -0.02    |
| 75) Chrysene-d12          | 21.67 | 240  | 256322   | 20.00 | ng    | -0.02    |
| 86) Perylene-d12          | 24.86 | 264  | 252829   | 20.00 | ng    | -0.05    |

|                             |       |     |        |        |    |       |
|-----------------------------|-------|-----|--------|--------|----|-------|
| System Monitoring Compounds |       |     |        |        |    |       |
| 5) 2-Fluorophenol           | 5.60  | 112 | 196554 | 132.09 | ng | -0.02 |
| 7) Phenol-d6                | 7.19  | 99  | 268581 | 116.66 | ng | -0.02 |
| 23) Nitrobenzene-d5         | 9.19  | 82  | 171718 | 78.29  | ng | -0.02 |
| 41) 2,4,6-Tribromophenol    | 16.14 | 330 | 110666 | 116.09 | ng | -0.02 |
| 44) 2-Fluorobiphenyl        | 13.28 | 172 | 393290 | 73.52  | ng | -0.02 |
| 78) Terphenyl-d14           | 20.01 | 244 | 854281 | 87.64  | ng | -0.02 |

|                                |       |     |       |        |    |      |
|--------------------------------|-------|-----|-------|--------|----|------|
| Target Compounds               |       |     |       | Qvalue |    |      |
| 2) 1,4-Dioxane                 | 3.49  | 88  | 57    | 0.084  | ng | # 1  |
| 4) n-Nitrosodimethylamine      | 3.84  | 42  | 98    | 0.112  | ng | # 1  |
| 6) Aniline                     | 7.56  | 93  | 237   | 0.079  | ng | # 1  |
| 9) Benzaldehyde                | 7.18  | 77  | 130   | 0.085  | ng | # 4  |
| 10) Phenol                     | 7.57  | 94  | 444   | 0.187  | ng | # 1  |
| 11) bis(2-Chloroethyl)ether    | 7.56  | 93  | 237   | 0.108  | ng | # 1  |
| 15) Benzyl Alcohol             | 8.35  | 79  | 2913  | 1.559  | ng | # 61 |
| 24) Nitrobenzene               | 9.20  | 77  | 284   | 0.121  | ng | # 43 |
| 25) Isophorone                 | 9.66  | 82  | 515   | 0.128  | ng | # 67 |
| 45) 1,1'-Biphenyl              | 13.52 | 154 | 80    | 0.013  | ng | # 40 |
| 47) 2-Nitroaniline             | 13.27 | 65  | 803   | 0.484  | ng | # 1  |
| 51) Acenaphthene               | 14.66 | 154 | 158   | 0.035  | ng | # 4  |
| 57) Fluorene                   | 16.14 | 166 | 4128  | 0.718  | ng | # 80 |
| 62) Azobenzene                 | 16.15 | 77  | 610   | 0.098  | ng | # 28 |
| 64) 4,6-Dinitro-2-methylphenol | 16.14 | 198 | 166   | 0.136  | ng | # 1  |
| 65) n-Nitrosodiphenylamine     | 16.15 | 169 | 4468  | 0.693  | ng | # 40 |
| 73) Di-n-butylphthalate        | 18.38 | 149 | 70    | 0.006  | ng | # 76 |
| 76) Benzidine                  | 20.01 | 184 | 14246 | 1.839  | ng | # 93 |
| 77) Pyrene                     | 20.01 | 202 | 2924  | 0.190  | ng | # 60 |
| 79) Butylbenzylphthalate       | 20.70 | 149 | 116   | 0.019  | ng | # 20 |
| 80) Benzo(a)anthracene         | 21.66 | 228 | 1335  | 0.089  | ng | # 78 |
| 82) Chrysene                   | 21.72 | 228 | 419   | 0.029  | ng | # 48 |
| 83) Bis(2-ethylhexyl)phthalate | 21.55 | 149 | 289   | 0.034  | ng | # 52 |
| 84) Di-n-octyl phthalate       | 22.75 | 149 | 407   | 0.029  | ng | # 1  |
| 87) Benzo(b)fluoranthene       | 23.87 | 252 | 243   | 0.018  | ng | # 4  |
| 88) Benzo(k)fluoranthene       | 23.92 | 252 | 68    | 0.005  | ng | # 3  |
| 89) Benzo(a)pyrene             | 24.70 | 252 | 55    | 0.004  | ng | # 59 |

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

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