

Data Path : Z:\HPCHEM1\BNA M\DATA\BM032516\  
 Data File : BM004771.D  
 Acq On : 25 Mar 2016 10:35  
 Operator : UM/SJ  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD0.446

Manual Integrations  
 APPROVED

Sohil  
 3/28/2016 7:33:09 PM

Quant Time: Mar 25 11:09:09 2016  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM032216.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Mar 25 03:18:48 2016  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|-----------------------------|-------|------|----------|--------|--------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 7.84  | 152  | 1393     | 5.00   | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.62 | 136  | 6212     | 5.00   | ng/ul  | 0.00     |
| 8) Acenaphthene-d10         | 14.47 | 164  | 3943     | 5.00   | ng/ul  | 0.00     |
| 12) Phenanthrene-d10        | 17.21 | 188  | 8634     | 5.00   | ng/ul  | 0.00     |
| 18) Chrysene-d12            | 21.40 | 240  | 8041     | 5.00   | ng/ul  | -0.01    |
| 22) Perylene-d12            | 23.69 | 264  | 6345     | 5.00   | ng/ul  | -0.01    |
| System Monitoring Compounds |       |      |          |        |        |          |
| 2) 1,4-Dioxane-d8           | 3.32  | 96   | 1375     | 2.04   | ng/uL  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.20 | 152  | 3647     | 0.52   | ng/ul  | -0.01    |
| 16) Fluoranthene-d10        | 19.24 | 212  | 8476     | 0.49   | ng/ul  | 0.00     |
| Target Compounds            |       |      |          | Ovalue |        |          |
| 3) 1,4-Dioxane              | 3.35  | 88   | 1645     | 2.00   | ng/ul# | 22       |
| 5) Naphthalene              | 10.67 | 128  | 6383     | 0.44   | ng/ul  | 96       |
| 7) 2-Methylnaphthalene      | 12.28 | 142  | 4566     | 0.49   | ng/ul  | 100      |
| 9) Acenaphthylene           | 14.18 | 152  | 7562     | 0.52   | ng/ul  | 96       |
| 10) Acenaphthene            | 14.52 | 153  | 5257     | 0.40   | ng/ul# | 84       |
| 11) Fluorene                | 15.51 | 166  | 6264     | 0.46   | ng/ul  | 87       |
| 13) Pentachlorophenol       | 16.87 | 266  | 760      | 0.66   | ng/ul  | 95       |
| 14) Phenanthrene            | 17.24 | 178  | 10102    | 0.42   | ng/ul# | 93       |
| 15) Anthracene              | 17.33 | 178  | 9634m    | 0.50   | ng/ul  |          |
| 17) Fluoranthene            | 19.27 | 202  | 11438    | 0.47   | ng/ul  | 94       |
| 19) Pyrene                  | 19.63 | 202  | 11997    | 0.43   | ng/ul  | 98       |
| 20) Benzo(a)anthracene      | 21.38 | 228  | 10568    | 0.55   | ng/ul# | 91       |
| 21) Chrysene                | 21.43 | 228  | 9971m    | 0.38   | ng/ul  |          |
| 23) Benzo(b)fluoranthene    | 23.00 | 252  | 9583     | 0.42   | ng/ul  | 89       |
| 24) Benzo(k)fluoranthene    | 23.04 | 252  | 8327     | 0.36   | ng/ul# | 94       |
| 25) Benzo(a)pyrene          | 23.59 | 252  | 8249     | 0.43   | ng/ul# | 91       |
| 26) Indeno(1,2,3-cd)pyrene  | 26.02 | 276  | 7271     | 0.35   | ng/ul# | 84       |
| 27) Dibenzo(a,h)anthracene  | 26.04 | 278  | 5721     | 0.35   | ng/ul# | 85       |
| 28) Benzo(g,h,i)perylene    | 26.73 | 276  | 5954     | 0.30   | ng/ul# | 87       |

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

