

Data Path : Z:\HPCHEM1\BNA M\DATA\BM032216\  
 Data File : BM004726.D  
 Acq On : 22 Mar 2016 13:05  
 Operator : UM/SJ  
 Sample : SSTD1.661  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD1.661

Manual Integrations  
 APPROVED

Sohil  
 3/23/2016 12:13:05 PM

Quant Time: Mar 22 13:58:09 2016  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM032216.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Mar 22 13:24:54 2016  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|-----------------------------|-------|------|----------|--------|--------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 7.84  | 152  | 1516     | 0.40   | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.62 | 136  | 6054     | 0.40   | ng/ul  | 0.00     |
| 8) Acenaphthene-d10         | 14.47 | 164  | 2774     | 0.40   | ng/ul  | 0.00     |
| 12) Phenanthrene-d10        | 17.21 | 188  | 5364     | 0.40   | ng/ul  | 0.00     |
| 18) Chrysene-d12            | 21.40 | 240  | 4948     | 0.40   | ng/ul  | 0.00     |
| 22) Perylene-d12            | 23.69 | 264  | 3417     | 0.40   | ng/ul  | 0.00     |
| System Monitoring Compounds |       |      |          |        |        |          |
| 2) 1,4-Dioxane-d8           | 3.32  | 96   | 5765     | 7.58   | ng/uL  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.22 | 152  | 11072    | 1.40   | ng/ul  | 0.00     |
| 16) Fluoranthene-d10        | 19.24 | 212  | 18114    | 1.34   | ng/ul  | 0.00     |
| Target Compounds            |       |      |          | Ovalue |        |          |
| 3) 1,4-Dioxane              | 3.35  | 88   | 6700     | 6.66   | ng/ul# | 17       |
| 5) Naphthalene              | 10.67 | 128  | 22246    | 1.53   | ng/ul  | 98       |
| 7) 2-Methylnaphthalene      | 12.28 | 142  | 14751    | 1.42   | ng/ul  | 99       |
| 9) Acenaphthylene           | 14.18 | 152  | 17121    | 1.72   | ng/ul  | 96       |
| 10) Acenaphthene            | 14.52 | 153  | 14448    | 1.58   | ng/ul  | 85       |
| 11) Fluorene                | 15.51 | 166  | 16207    | 1.47   | ng/ul  | 85       |
| 13) Pentachlorophenol       | 16.87 | 266  | 1252     | 2.83   | ng/ul  | 96       |
| 14) Phenanthrene            | 17.24 | 178  | 25201    | 1.72   | ng/ul# | 93       |
| 15) Anthracene              | 17.33 | 178  | 22725m   | 1.84   | ng/ul  |          |
| 17) Fluoranthene            | 19.27 | 202  | 26458    | 1.44   | ng/ul  | 97       |
| 19) Pyrene                  | 19.63 | 202  | 26692    | 1.73   | ng/ul  | 99       |
| 20) Benzo(a)anthracene      | 21.38 | 228  | 19403    | 1.54   | ng/ul  | 95       |
| 21) Chrysene                | 21.43 | 228  | 24543    | 1.60   | ng/ul  | 97       |
| 23) Benzo(b)fluoranthene    | 22.99 | 252  | 20909    | 1.84   | ng/ul  | 99       |
| 24) Benzo(k)fluoranthene    | 23.03 | 252  | 21252m   | 1.98   | ng/ul  |          |
| 25) Benzo(a)pyrene          | 23.59 | 252  | 17670    | 1.78   | ng/ul  | 97       |
| 26) Indeno(1,2,3-cd)pyrene  | 26.01 | 276  | 19649    | 1.80   | ng/ul# | 82       |
| 27) Dibenzo(a,h)anthracene  | 26.03 | 278  | 15717    | 1.91   | ng/ul# | 94       |
| 28) Benzo(g,h,i)perylene    | 26.71 | 276  | 17587    | 1.72   | ng/ul# | 93       |

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

