

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121515\  
 Data File : BM003572.D  
 Acq On : 16 Dec 2015 06:04  
 Operator : UM/IZ  
 Sample : SSTDC00.4EC  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD0.454

Manual Integrations  
 APPROVED

MMdadoda  
 12/16/2015 6:39:15 PM

Quant Time: Dec 16 18:31:41 2015  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-SIM-BM121515.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Dec 16 05:49:57 2015  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|-----------------------------|-------|------|----------|--------|--------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 7.71  | 152  | 3627     | 0.40   | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.50 | 136  | 14078    | 0.40   | ng/ul  | 0.00     |
| 8) Acenaphthene-d10         | 14.37 | 164  | 7921     | 0.40   | ng/ul  | 0.00     |
| 12) Phenanthrene-d10        | 17.11 | 188  | 20174    | 0.40   | ng/ul  | 0.00     |
| 18) Chrysene-d12            | 21.31 | 240  | 31884    | 0.40   | ng/ul  | 0.00     |
| 22) Perylene-d12            | 23.57 | 264  | 23008    | 0.40   | ng/ul  | 0.00     |
| System Monitoring Compounds |       |      |          |        |        |          |
| 2) 1,4-Dioxane-d8           | 3.17  | 96   | 3121     | 2.01   | ng/uL  | -0.02    |
| 6) 2-Methylnaphthalene-d10  | 12.09 | 152  | 7090     | 0.42   | ng/ul  | -0.01    |
| 16) Fluoranthene-d10        | 19.14 | 212  | 19303    | 0.45   | ng/ul  | 0.00     |
| Target Compounds            |       |      |          | Ovalue |        |          |
| 3) 1,4-Dioxane              | 3.20  | 88   | 3867     | 1.96   | ng/ul# | 100      |
| 5) Naphthalene              | 10.55 | 128  | 13037    | 0.41   | ng/ul  | 98       |
| 7) 2-Methylnaphthalene      | 12.16 | 142  | 9154     | 0.43   | ng/ul  | 99       |
| 9) Acenaphthylene           | 14.07 | 152  | 12400    | 0.41   | ng/ul  | 95       |
| 10) Acenaphthene            | 14.42 | 153  | 10248    | 0.42   | ng/ul  | 87       |
| 11) Fluorene                | 15.41 | 166  | 12149    | 0.45   | ng/ul  | 84       |
| 13) Pentachlorophenol       | 16.76 | 266  | 1337     | 0.46   | ng/ul  | 96       |
| 14) Phenanthrene            | 17.14 | 178  | 22010m   | 0.42   | ng/ul  |          |
| 15) Anthracene              | 17.24 | 178  | 18641    | 0.41   | ng/ul  | 98       |
| 17) Fluoranthene            | 19.18 | 202  | 26126    | 0.46   | ng/ul  | 92       |
| 19) Pyrene                  | 19.54 | 202  | 27595    | 0.32   | ng/ul  | 95       |
| 20) Benzo(a)anthracene      | 21.29 | 228  | 27234    | 0.35   | ng/ul  | 93       |
| 21) Chrysene                | 21.35 | 228  | 34912    | 0.39   | ng/ul  | 93       |
| 23) Benzo(b)fluoranthene    | 22.89 | 252  | 33538    | 0.43   | ng/ul  | 96       |
| 24) Benzo(k)fluoranthene    | 22.93 | 252  | 29195    | 0.41   | ng/ul  | 96       |
| 25) Benzo(a)pyrene          | 23.46 | 252  | 27043    | 0.41   | ng/ul  | 97       |
| 26) Indeno(1,2,3-cd)pyrene  | 25.84 | 276  | 31781    | 0.41   | ng/ul# | 91       |
| 27) Dibenzo(a,h)anthracene  | 25.85 | 278  | 25291    | 0.41   | ng/ul  | 99       |
| 28) Benzo(g,h,i)perylene    | 26.53 | 276  | 27757    | 0.42   | ng/ul  | 97       |

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

