

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110520\  
 Data File : BM027704.D  
 Acq On : 05 Nov 2020 03:46  
 Operator : JU/CG  
 Sample : SSTD1.601  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD1.6015

Quant Time: Nov 05 04:18:53 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM110520.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Thu Nov 05 03:44:30 2020  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|-----------------------------|-------|------|----------|--------|--------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 8.40  | 152  | 455      | 0.40   | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 11.24 | 136  | 1542     | 0.40   | ng/ul  | 0.00     |
| 9) Acenaphthene-d10         | 15.01 | 164  | 972      | 0.40   | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 17.72 | 188  | 2228     | 0.40   | ng/ul  | 0.00     |
| 19) Chrysene-d12            | 21.82 | 240  | 2050     | 0.40   | ng/ul  | 0.00     |
| 23) Perylene-d12            | 24.26 | 264  | 1888     | 0.40   | ng/ul  | 0.00     |
| System Monitoring Compounds |       |      |          |        |        |          |
| 3) 1,4-Dioxane-d8           | 3.54  | 96   | 706      | 0.76   | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.80 | 152  | 3944     | 1.59   | ng/ul  | 0.00     |
| 17) Fluoranthene-d10        | 19.71 | 212  | 8807     | 1.50   | ng/ul  | 0.00     |
| Target Compounds            |       |      |          | Ovalue |        |          |
| 2) 1,4-Dioxane              | 3.58  | 88   | 762      | 1.204  | ng/ul# | 91       |
| 5) Naphthalene              | 11.29 | 128  | 6685     | 1.559  | ng/ul  | 93       |
| 7) 2-Methylnaphthalene      | 12.87 | 142  | 4674     | 1.617  | ng/ul  | 99       |
| 8) 1-Methylnaphthalene      | 13.09 | 142  | 4411     | 1.605  | ng/ul  | 98       |
| 10) Acenaphthylene          | 14.74 | 152  | 6713     | 1.503  | ng/ul  | 96       |
| 11) Acenaphthene            | 15.07 | 153  | 5057     | 1.520  | ng/ul  | 98       |
| 12) Fluorene                | 16.04 | 166  | 6052     | 1.564  | ng/ul  | 97       |
| 14) Pentachlorophenol       | 17.36 | 266  | 1107     | 1.541  | ng/ul  | 96       |
| 15) Phenanthrene            | 17.76 | 178  | 9559     | 1.445  | ng/ul  | 94       |
| 16) Anthracene              | 17.85 | 178  | 9635     | 1.512  | ng/ul  | 97       |
| 18) Fluoranthene            | 19.74 | 202  | 11425    | 1.500  | ng/ul# | 96       |
| 20) Pyrene                  | 20.09 | 202  | 11427    | 1.366  | ng/ul  | 96       |
| 21) Benzo(a)anthracene      | 21.80 | 228  | 10709    | 1.414  | ng/ul  | 98       |
| 22) Chrysene                | 21.86 | 228  | 11037    | 1.404  | ng/ul  | 98       |
| 24) Benzo(b)fluoranthene    | 23.52 | 252  | 10649    | 1.368  | ng/ul  | 93       |
| 25) Benzo(k)fluoranthene    | 23.57 | 252  | 10881    | 1.401  | ng/ul# | 92       |
| 26) Benzo(a)pyrene          | 24.16 | 252  | 9962     | 1.446  | ng/ul# | 91       |
| 27) Indeno(1,2,3-cd)pyrene  | 26.80 | 276  | 12096    | 1.514  | ng/ul# | 98       |
| 28) Dibenzo(a,h)anthracene  | 26.81 | 278  | 10134    | 1.523  | ng/ul# | 90       |
| 29) Benzo(g,h,i)perylene    | 27.58 | 276  | 10212    | 1.492  | ng/ul# | 89       |

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

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