

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM071321\  
 Data File : BM030924.D  
 Acq On : 13 Jul 2021 11:01  
 Operator : CG/JU  
 Sample : SST00.202  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SST00.2002

Quant Time: Jul 13 12:13:40 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM071321.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jul 13 12:12:03 2021  
 Response via : Initial Calibration

| Internal Standards          | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|-----------------------------|-------|------|----------|--------|--------|----------|
| 1) 1,4-Dichlorobenzene-d4   | 7.67  | 152  | 931      | 0.40   | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.44 | 136  | 2997     | 0.40   | ng/ul  | 0.00     |
| 9) Acenaphthene-d10         | 14.30 | 164  | 1908     | 0.40   | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 17.05 | 188  | 4082     | 0.40   | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.24 | 240  | 3341     | 0.40   | ng/ul  | 0.00     |
| 23) Perylene-d12            | 23.43 | 264  | 3337     | 0.40   | ng/ul  | 0.00     |
| System Monitoring Compounds |       |      |          |        |        |          |
| 3) 1,4-Dioxane-d8           | 3.24  | 96   | 189      | 0.19   | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.03 | 152  | 1054     | 0.22   | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 19.08 | 212  | 2348     | 0.25   | ng/ul  | 0.00     |
| Target Compounds            |       |      |          | Ovalue |        |          |
| 2) 1,4-Dioxane              | 3.28  | 88   | 208      | 0.198  | ng/ul  | 91       |
| 5) Naphthalene              | 10.49 | 128  | 1785     | 0.209  | ng/ul# | 95       |
| 7) 2-Methylnaphthalene      | 12.11 | 142  | 1256     | 0.218  | ng/ul  | 99       |
| 8) 1-Methylnaphthalene      | 12.33 | 142  | 1232     | 0.220  | ng/ul  | 98       |
| 10) Acenaphthylene          | 14.02 | 152  | 1908     | 0.224  | ng/ul  | 97       |
| 11) Acenaphthene            | 14.36 | 153  | 1372     | 0.199  | ng/ul  | 96       |
| 12) Fluorene                | 15.35 | 166  | 1633     | 0.214  | ng/ul  | 96       |
| 14) Pentachlorophenol       | 16.70 | 266  | 258      | 0.194  | ng/ul  | 94       |
| 15) Phenanthrene            | 17.09 | 178  | 2561     | 0.187  | ng/ul  | 96       |
| 16) Anthracene              | 17.18 | 178  | 2484     | 0.207  | ng/ul  | 95       |
| 19) Fluoranthene            | 19.11 | 202  | 3064     | 0.208  | ng/ul# | 96       |
| 20) Pyrene                  | 19.47 | 202  | 3081     | 0.208  | ng/ul  | 97       |
| 21) Benzo(a)anthracene      | 21.22 | 228  | 2930     | 0.220  | ng/ul  | 97       |
| 22) Chrysene                | 21.27 | 228  | 3027     | 0.203  | ng/ul  | 99       |
| 24) Benzo(b)fluoranthene    | 22.78 | 252  | 3225     | 0.218  | ng/ul  | 83       |
| 25) Benzo(k)fluoranthene    | 22.82 | 252  | 3230     | 0.212  | ng/ul# | 81       |
| 26) Benzo(a)pyrene          | 23.34 | 252  | 2814     | 0.224  | ng/ul# | 80       |
| 27) Indeno(1,2,3-cd)pyrene  | 25.62 | 276  | 3789     | 0.231  | ng/ul# | 99       |
| 28) Dibenzo(a,h)anthracene  | 25.64 | 278  | 3074     | 0.234  | ng/ul# | 87       |
| 29) Benzo(g,h,i)perylene    | 26.28 | 276  | 3195     | 0.225  | ng/ul# | 92       |

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

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