

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN082322\  
 Data File : BN021330.D  
 Acq On : 23 Aug 2022 15:56  
 Operator : CG/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD0.4297

Quant Time: Aug 24 00:59:12 2022  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-SIM-BN082222.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Aug 22 18:06:52 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.097  | 152  | 2598     | 0.400 | ng/u1  | 0.00     |
| 4) Naphthalene-d8           | 10.927 | 136  | 8073     | 0.400 | ng/u1  | # 0.00   |
| 9) Acenaphthene-d10         | 14.749 | 164  | 4283     | 0.400 | ng/u1  | 0.00     |
| 13) Phenanthrene-d10        | 17.497 | 188  | 9094     | 0.400 | ng/u1  | # 0.00   |
| 17) Chrysene-d12            | 21.688 | 240  | 7659     | 0.400 | ng/u1  | # 0.00   |
| 23) Perylene-d12            | 24.222 | 264  | 6155     | 0.400 | ng/u1  | 0.00     |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.385  | 96   | 1312     | 0.391 | ng/u1  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.511 | 152  | 4911     | 0.402 | ng/u1  | 0.00     |
| 18) Fluoranthene-d10        | 19.524 | 212  | 9509     | 0.403 | ng/u1  | 0.00     |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 2) 1,4-Dioxane              | 3.422  | 88   | 1378     | 0.392 | ng/u1  | 96       |
| 5) Naphthalene              | 10.977 | 128  | 9047     | 0.354 | ng/u1  | 99       |
| 7) 2-Methylnaphthalene      | 12.588 | 142  | 5707     | 0.357 | ng/u1  | 98       |
| 8) 1-Methylnaphthalene      | 12.803 | 142  | 6065     | 0.404 | ng/u1  | 100      |
| 10) Acenaphthylene          | 14.472 | 152  | 7582     | 0.343 | ng/u1  | 99       |
| 11) Acenaphthene            | 14.809 | 153  | 5973     | 0.353 | ng/u1  | 98       |
| 12) Fluorene                | 15.795 | 166  | 6653     | 0.352 | ng/u1  | 99       |
| 14) Pentachlorophenol       | 17.143 | 266  | 390      | 0.388 | ng/u1  | 97       |
| 15) Phenanthrene            | 17.540 | 178  | 11661    | 0.346 | ng/u1  | 99       |
| 16) Anthracene              | 17.633 | 178  | 9161     | 0.339 | ng/u1  | 99       |
| 19) Fluoranthene            | 19.552 | 202  | 12548    | 0.346 | ng/u1  | 98       |
| 20) Pyrene                  | 19.919 | 202  | 12270    | 0.338 | ng/u1  | 98       |
| 21) Benzo(a)anthracene      | 21.670 | 228  | 9464     | 0.357 | ng/u1  | 99       |
| 22) Chrysene                | 21.726 | 228  | 11982    | 0.351 | ng/u1  | 100      |
| 24) Benzo(b)fluoranthene    | 23.442 | 252  | 11060    | 0.418 | ng/u1  | 97       |
| 25) Benzo(k)fluoranthene    | 23.491 | 252  | 10754    | 0.378 | ng/u1  | 97       |
| 26) Benzo(a)pyrene          | 24.108 | 252  | 9346     | 0.375 | ng/u1  | 97       |
| 27) Indeno(1,2,3-cd)pyrene  | 26.894 | 276  | 11217    | 0.384 | ng/u1# | 97       |
| 28) Dibenzo(a,h)anthracene  | 26.924 | 278  | 9280     | 0.386 | ng/u1  | 95       |
| 29) Benzo(g,h,i)perylene    | 27.722 | 276  | 10987    | 0.378 | ng/u1  | 98       |
| -----                       |        |      |          |       |        |          |

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

