

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN112622\  
 Data File : BN022883.D  
 Acq On : 26 Nov 2022 16:15  
 Operator : CG/JU  
 Sample : SSTDCCC0.4  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SSTD0.4341

Quant Time: Nov 28 03:48:35 2022  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-SIM-BN112622.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Nov 28 03:16:56 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 8.070  | 152  | 7598     | 0.400 | ng/u1  | 0.00     |
| 4) Naphthalene-d8           | 10.892 | 136  | 19539    | 0.400 | ng/u1  | 0.00     |
| 9) Acenaphthene-d10         | 14.710 | 164  | 9543     | 0.400 | ng/u1  | 0.00     |
| 13) Phenanthrene-d10        | 17.454 | 188  | 18529    | 0.400 | ng/u1  | 0.00     |
| 17) Chrysene-d12            | 21.639 | 240  | 12124    | 0.400 | ng/u1  | 0.00     |
| 23) Perylene-d12            | 24.130 | 264  | 8543     | 0.400 | ng/u1  | 0.00     |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.412  | 96   | 3979     | 0.436 | ng/u1  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.476 | 152  | 12013    | 0.403 | ng/u1  | 0.00     |
| 18) Fluoranthene-d10        | 19.481 | 212  | 18555    | 0.414 | ng/u1  | 0.00     |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 2) 1,4-Dioxane              | 3.450  | 88   | 4025     | 0.423 | ng/u1  | 97       |
| 5) Naphthalene              | 10.942 | 128  | 25052    | 0.422 | ng/u1  | 100      |
| 7) 2-Methylnaphthalene      | 12.548 | 142  | 15109    | 0.408 | ng/u1  | 99       |
| 8) 1-Methylnaphthalene      | 12.762 | 142  | 15278    | 0.405 | ng/u1  | 100      |
| 10) Acenaphthylene          | 14.432 | 152  | 19504    | 0.415 | ng/u1  | 100      |
| 11) Acenaphthene            | 14.770 | 153  | 15598    | 0.419 | ng/u1  | 98       |
| 12) Fluorene                | 15.756 | 166  | 17000    | 0.403 | ng/u1  | 99       |
| 14) Pentachlorophenol       | 17.099 | 266  | 1417     | 0.366 | ng/u1# | 100      |
| 15) Phenanthrene            | 17.496 | 178  | 26423    | 0.416 | ng/u1  | 100      |
| 16) Anthracene              | 17.589 | 178  | 22303    | 0.409 | ng/u1  | 100      |
| 19) Fluoranthene            | 19.509 | 202  | 25601    | 0.409 | ng/u1  | 100      |
| 20) Pyrene                  | 19.871 | 202  | 25004    | 0.407 | ng/u1  | 99       |
| 21) Benzo(a)anthracene      | 21.622 | 228  | 18705    | 0.404 | ng/u1  | 99       |
| 22) Chrysene                | 21.674 | 228  | 21460    | 0.422 | ng/u1  | 99       |
| 24) Benzo(b)fluoranthene    | 23.364 | 252  | 18274    | 0.422 | ng/u1  | 98       |
| 25) Benzo(k)fluoranthene    | 23.416 | 252  | 19188    | 0.419 | ng/u1  | 99       |
| 26) Benzo(a)pyrene          | 24.016 | 252  | 13770    | 0.418 | ng/u1  | 99       |
| 27) Indeno(1,2,3-cd)pyrene  | 26.741 | 276  | 18238    | 0.465 | ng/u1# | 100      |
| 28) Dibenzo(a,h)anthracene  | 26.768 | 278  | 14574    | 0.468 | ng/u1  | 98       |
| 29) Benzo(g,h,i)perylene    | 27.540 | 276  | 16500    | 0.483 | ng/u1  | 99       |
| -----                       |        |      |          |       |        |          |

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

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