

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN082724\  
 Data File : BN033632.D  
 Acq On : 27 Aug 2024 12:24  
 Operator : MA/JU  
 Sample : SST0.421  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 SST0.4204

Quant Time: Aug 27 23:37:39 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-SIM-BN082724.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Aug 27 23:34:19 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.540  | 152  | 5619     | 0.400 | ng/u1  | 0.00     |
| 4) Naphthalene-d8           | 10.298 | 136  | 15327    | 0.400 | ng/u1  | 0.00     |
| 9) Acenaphthene-d10         | 14.178 | 164  | 7957     | 0.400 | ng/u1  | 0.00     |
| 13) Phenanthrene-d10        | 16.930 | 188  | 16648    | 0.400 | ng/u1  | 0.00     |
| 17) Chrysene-d12            | 21.137 | 240  | 14141    | 0.400 | ng/u1  | 0.00     |
| 23) Perylene-d12            | 23.300 | 264  | 15253    | 0.400 | ng/u1  | 0.00     |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.143  | 96   | 2338     | 0.405 | ng/u1  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 11.898 | 152  | 8672     | 0.394 | ng/u1  | 0.00     |
| 18) Fluoranthene-d10        | 18.965 | 212  | 16318    | 0.387 | ng/u1  | 0.00     |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 2) 1,4-Dioxane              | 3.177  | 88   | 2617     | 0.408 | ng/u1  | 100      |
| 5) Naphthalene              | 10.347 | 128  | 16107    | 0.393 | ng/u1  | 100      |
| 7) 2-Methylnaphthalene      | 11.970 | 142  | 10234    | 0.394 | ng/u1  | 100      |
| 8) 1-Methylnaphthalene      | 12.195 | 142  | 10572    | 0.396 | ng/u1  | 100      |
| 10) Acenaphthylene          | 13.891 | 152  | 14408    | 0.399 | ng/u1  | 100      |
| 11) Acenaphthene            | 14.238 | 153  | 10508    | 0.403 | ng/u1  | 100      |
| 12) Fluorene                | 15.232 | 166  | 11816    | 0.405 | ng/u1  | 100      |
| 14) Pentachlorophenol       | 16.584 | 266  | 1950     | 0.372 | ng/u1  | 100      |
| 15) Phenanthrene            | 16.968 | 178  | 18490    | 0.377 | ng/u1  | 100      |
| 16) Anthracene              | 17.061 | 178  | 16792    | 0.380 | ng/u1  | 100      |
| 19) Fluoranthene            | 18.997 | 202  | 21040    | 0.377 | ng/u1  | 100      |
| 20) Pyrene                  | 19.360 | 202  | 22332    | 0.377 | ng/u1  | 100      |
| 21) Benzo(a)anthracene      | 21.119 | 228  | 20869    | 0.375 | ng/u1  | 100      |
| 22) Chrysene                | 21.172 | 228  | 20937    | 0.372 | ng/u1  | 100      |
| 24) Benzo(b)fluoranthene    | 22.659 | 252  | 22133    | 0.374 | ng/u1  | 100      |
| 25) Benzo(k)fluoranthene    | 22.700 | 252  | 21858    | 0.368 | ng/u1  | 100      |
| 26) Benzo(a)pyrene          | 23.206 | 252  | 17661    | 0.372 | ng/u1  | 100      |
| 27) Indeno(1,2,3-cd)pyrene  | 25.450 | 276  | 26363    | 0.369 | ng/u1  | 100      |
| 28) Dibenzo(a,h)anthracene  | 25.464 | 278  | 20443    | 0.369 | ng/u1  | 100      |
| 29) Benzo(g,h,i)perylene    | 26.098 | 276  | 20645    | 0.370 | ng/u1  | 100      |
| -----                       |        |      |          |       |        |          |

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

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