

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM092721\  
 Data File : BM032200.D  
 Acq On : 27 Sep 2021 11:13  
 Operator : CG/JU  
 Sample : SST0.431  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SST0.4031

Quant Time: Sep 27 11:46:39 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-SIM-BM092721.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Sep 27 11:46:23 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.664  | 152  | 7518     | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.456 | 136  | 24060    | 0.400 | ng/ul  | 0.00     |
| 9) Acenaphthene-d10         | 14.308 | 164  | 12520    | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 17.033 | 188  | 25962    | 0.400 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.196 | 240  | 21361    | 0.400 | ng/ul  | 0.00     |
| 23) Perylene-d12            | 23.358 | 264  | 16396    | 0.400 | ng/ul  | 0.00     |
|                             |        |      |          |       |        |          |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 2.993  | 96   | 3188     | 0.386 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.048 | 152  | 12632    | 0.339 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 19.053 | 212  | 23734    | 0.379 | ng/ul  | 0.00     |
|                             |        |      |          |       |        |          |
| Target Compounds            |        |      |          |       | Qvalue |          |
| 2) 1,4-Dioxane              | 3.027  | 88   | 3344     | 0.415 | ng/ul  | 100      |
| 5) Naphthalene              | 10.505 | 128  | 27754    | 0.435 | ng/ul  | 100      |
| 7) 2-Methylnaphthalene      | 12.120 | 142  | 17916    | 0.402 | ng/ul  | 100      |
| 8) 1-Methylnaphthalene      | 12.341 | 142  | 17628    | 0.402 | ng/ul  | 100      |
| 10) Acenaphthylene          | 14.024 | 152  | 19423    | 0.384 | ng/ul  | 100      |
| 11) Acenaphthene            | 14.369 | 153  | 17828    | 0.431 | ng/ul  | 100      |
| 12) Fluorene                | 15.351 | 166  | 20435    | 0.424 | ng/ul  | 100      |
| 14) Pentachlorophenol       | 16.700 | 266  | 2348     | 0.342 | ng/ul  | 100      |
| 15) Phenanthrene            | 17.075 | 178  | 32700    | 0.414 | ng/ul  | 100      |
| 16) Anthracene              | 17.165 | 178  | 26735    | 0.386 | ng/ul  | 100      |
| 19) Fluoranthene            | 19.084 | 202  | 36095    | 0.391 | ng/ul  | 100      |
| 20) Pyrene                  | 19.442 | 202  | 36061    | 0.379 | ng/ul  | 100      |
| 21) Benzo(a)anthracene      | 21.179 | 228  | 29004    | 0.372 | ng/ul  | 100      |
| 22) Chrysene                | 21.230 | 228  | 34053    | 0.380 | ng/ul  | 100      |
| 24) Benzo(b)fluoranthene    | 22.714 | 252  | 32468    | 0.431 | ng/ul  | 100      |
| 25) Benzo(k)fluoranthene    | 22.755 | 252  | 31746    | 0.408 | ng/ul  | 100      |
| 26) Benzo(a)pyrene          | 23.264 | 252  | 24372    | 0.370 | ng/ul  | 100      |
| 27) Indeno(1,2,3-cd)pyrene  | 25.505 | 276  | 29520    | 0.329 | ng/ul# | 100      |
| 28) Dibenzo(a,h)anthracene  | 25.510 | 278  | 23970    | 0.332 | ng/ul  | 100      |
| 29) Benzo(g,h,i)perylene    | 26.164 | 276  | 28161    | 0.353 | ng/ul  | 100      |
| -----                       |        |      |          |       |        |          |

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

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