

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM041222\  
 Data File : BM034623.D  
 Acq On : 12 Apr 2022 15:53  
 Operator : CG/JU  
 Sample : SSTDICV0.4  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SICV028

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 04/13/2022  
 Supervised By :Yogesh Patel 04/14/2022

Quant Time: Apr 12 18:43:45 2022  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-SIM-BM041222.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Apr 12 15:46:33 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.893  | 152  | 1549     | 0.400 | ng/ul  | -0.02    |
| 4) Naphthalene-d8           | 10.704 | 136  | 4600     | 0.400 | ng/ul  | # 0.00   |
| 9) Acenaphthene-d10         | 14.533 | 164  | 2397     | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 17.287 | 188  | 3436     | 0.400 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.470 | 240  | 3352     | 0.400 | ng/ul  | # 0.00   |
| 23) Perylene-d12            | 23.855 | 264  | 2651     | 0.400 | ng/ul  | # 0.01   |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.256  | 96   | 1036     | 0.386 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.304 | 152  | 2442     | 0.401 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 19.312 | 212  | 3987     | 0.409 | ng/ul  | 0.00     |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 2) 1,4-Dioxane              | 3.290  | 88   | 1040     | 0.391 | ng/ul# | 67       |
| 5) Naphthalene              | 10.753 | 128  | 5502     | 0.413 | ng/ul# | 92       |
| 7) 2-Methylnaphthalene      | 12.381 | 142  | 3348     | 0.421 | ng/ul  | 99       |
| 8) 1-Methylnaphthalene      | 12.590 | 142  | 3169     | 0.394 | ng/ul  | 98       |
| 10) Acenaphthylene          | 14.260 | 152  | 4881     | 0.434 | ng/ul  | 94       |
| 11) Acenaphthene            | 14.597 | 153  | 3694     | 0.410 | ng/ul  | 95       |
| 12) Fluorene                | 15.587 | 166  | 3712     | 0.401 | ng/ul  | 99       |
| 14) Pentachlorophenol       | 16.958 | 266  | 611      | 0.458 | ng/ul  | 95       |
| 15) Phenanthrene            | 17.329 | 178  | 4634     | 0.417 | ng/ul  | 99       |
| 16) Anthracene              | 17.426 | 178  | 3460     | 0.408 | ng/ul  | 99       |
| 19) Fluoranthene            | 19.340 | 202  | 5387     | 0.389 | ng/ul# | 95       |
| 20) Pyrene                  | 19.702 | 202  | 6720     | 0.411 | ng/ul# | 93       |
| 21) Benzo(a)anthracene      | 21.479 | 228  | 2275m    | 0.325 | ng/ul  |          |
| 22) Chrysene                | 21.508 | 228  | 3418     | 0.386 | ng/ul  | 98       |
| 24) Benzo(b)fluoranthene    | 23.124 | 252  | 3635m    | 0.417 | ng/ul  |          |
| 25) Benzo(k)fluoranthene    | 23.177 | 252  | 3736     | 0.428 | ng/ul# | 77       |
| 26) Benzo(a)pyrene          | 23.753 | 252  | 4684     | 0.468 | ng/ul# | 61       |
| 27) Indeno(1,2,3-cd)pyrene  | 26.349 | 276  | 5722     | 0.431 | ng/ul# | 57       |
| 28) Dibenzo(a,h)anthracene  | 26.396 | 278  | 4094     | 0.426 | ng/ul# | 58       |
| 29) Benzo(g,h,i)perylene    | 27.097 | 276  | 6421     | 0.444 | ng/ul# | 94       |
| -----                       |        |      |          |       |        |          |

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

