

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM040422\  
 Data File : BM034494.D  
 Acq On : 04 Apr 2022 19:23  
 Operator : CG/JU  
 Sample : SSTDICV0.4  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SICV021

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 04/05/2022  
 Supervised By :Yogesh Patel 04/09/2022

Quant Time: Apr 05 11:53:00 2022  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-SIM-BM040422.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Apr 05 11:50:53 2022  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.925  | 152  | 2001     | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.735 | 136  | 5701     | 0.400 | ng/ul  | # 0.00   |
| 9) Acenaphthene-d10         | 14.564 | 164  | 2866     | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 17.316 | 188  | 3978     | 0.400 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.498 | 240  | 3686     | 0.400 | ng/ul  | # 0.00   |
| 23) Perylene-d12            | 23.901 | 264  | 3040     | 0.400 | ng/ul  | # 0.00   |
|                             |        |      |          |       |        |          |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.297  | 96   | 1545     | 0.461 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.346 | 152  | 3410     | 0.456 | ng/ul  | 0.01     |
| 18) Fluoranthene-d10        | 19.338 | 212  | 5082     | 0.470 | ng/ul  | 0.00     |
|                             |        |      |          |       |        |          |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       |        | Qvalue   |
| 2) 1,4-Dioxane              | 3.330  | 88   | 1502     | 0.449 | ng/ul# | 72       |
| 5) Naphthalene              | 10.790 | 128  | 7915     | 0.444 | ng/ul# | 91       |
| 7) 2-Methylnaphthalene      | 12.423 | 142  | 4576     | 0.470 | ng/ul  | 97       |
| 8) 1-Methylnaphthalene      | 12.621 | 142  | 4309     | 0.440 | ng/ul  | 100      |
| 10) Acenaphthylene          | 14.287 | 152  | 6787     | 0.512 | ng/ul# | 91       |
| 11) Acenaphthene            | 14.624 | 153  | 5071     | 0.472 | ng/ul  | 95       |
| 12) Fluorene                | 15.615 | 166  | 5025     | 0.454 | ng/ul  | 99       |
| 14) Pentachlorophenol       | 16.983 | 266  | 815      | 0.483 | ng/ul  | 95       |
| 15) Phenanthrene            | 17.358 | 178  | 6125     | 0.474 | ng/ul  | 99       |
| 16) Anthracene              | 17.464 | 178  | 4477     | 0.456 | ng/ul  | 97       |
| 19) Fluoranthene            | 19.366 | 202  | 7040     | 0.455 | ng/ul  | 95       |
| 20) Pyrene                  | 19.729 | 202  | 8619     | 0.477 | ng/ul# | 94       |
| 21) Benzo(a)anthracene      | 21.486 | 228  | 3293m    | 0.411 | ng/ul  |          |
| 22) Chrysene                | 21.530 | 228  | 4143m    | 0.437 | ng/ul  |          |
| 24) Benzo(b)fluoranthene    | 23.164 | 252  | 4459m    | 0.448 | ng/ul  |          |
| 25) Benzo(k)fluoranthene    | 23.214 | 252  | 3847     | 0.434 | ng/ul# | 80       |
| 26) Benzo(a)pyrene          | 23.801 | 252  | 5795     | 0.499 | ng/ul# | 65       |
| 27) Indeno(1,2,3-cd)pyrene  | 26.424 | 276  | 6505     | 0.466 | ng/ul# | 98       |
| 28) Dibenzo(a,h)anthracene  | 26.445 | 278  | 4945     | 0.486 | ng/ul# | 63       |
| 29) Benzo(g,h,i)perylene    | 27.172 | 276  | 7322     | 0.484 | ng/ul  | 94       |
| -----                       |        |      |          |       |        |          |

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

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