

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM080921\  
 Data File : BM031623.D  
 Acq On : 10 Aug 2021 03:52  
 Operator : CG/JU  
 Sample : PB138111BS  
 Misc :  
 ALS Vial : 31 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SLCS111

Manual Integrations  
 APPROVED

mohammad  
 8/10/2021 4:36:46 PM

Quant Time: Aug 10 06:05:20 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-SIM-BM072821.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Aug 09 10:16:43 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.580  | 152  | 1278     | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.342 | 136  | 4657     | 0.400 | ng/ul  | 0.00     |
| 9) Acenaphthene-d10         | 14.211 | 164  | 2469     | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 16.963 | 188  | 4674     | 0.400 | ng/ul  | # 0.00   |
| 17) Chrysene-d12            | 21.157 | 240  | 3724     | 0.400 | ng/ul  | # 0.00   |
| 23) Perylene-d12            | 23.315 | 264  | 3785     | 0.400 | ng/ul  | # 0.00   |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.186  | 96   | 1019     | 0.718 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 11.934 | 152  | 3014     | 0.384 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 18.996 | 212  | 5329     | 0.429 | ng/ul  | 0.00     |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 2) 1,4-Dioxane              | 3.217  | 88   | 2765     | 1.928 | ng/ul# | 75       |
| 5) Naphthalene              | 10.387 | 128  | 5292     | 0.381 | ng/ul  | 100      |
| 7) 2-Methylnaphthalene      | 12.010 | 142  | 3484     | 0.368 | ng/ul  | 98       |
| 8) 1-Methylnaphthalene      | 12.231 | 142  | 3143     | 0.336 | ng/ul  | 100      |
| 10) Acenaphthylene          | 13.931 | 152  | 4400     | 0.418 | ng/ul  | 100      |
| 11) Acenaphthene            | 14.276 | 153  | 3437     | 0.388 | ng/ul  | 99       |
| 12) Fluorene                | 15.269 | 166  | 3679     | 0.361 | ng/ul  | 98       |
| 14) Pentachlorophenol       | 16.623 | 266  | 1177     | 0.786 | ng/ul  | 97       |
| 15) Phenanthrene            | 17.005 | 178  | 5652     | 0.381 | ng/ul  | 99       |
| 16) Anthracene              | 17.099 | 178  | 5357     | 0.386 | ng/ul  | 99       |
| 19) Fluoranthene            | 19.027 | 202  | 6586     | 0.414 | ng/ul# | 95       |
| 20) Pyrene                  | 19.389 | 202  | 6838     | 0.424 | ng/ul# | 93       |
| 21) Benzo(a)anthracene      | 21.140 | 228  | 6329     | 0.421 | ng/ul  | 97       |
| 22) Chrysene                | 21.193 | 228  | 6411     | 0.412 | ng/ul  | 99       |
| 24) Benzo(b)fluoranthene    | 22.672 | 252  | 6491m    | 0.381 | ng/ul  |          |
| 25) Benzo(k)fluoranthene    | 22.714 | 252  | 6482     | 0.368 | ng/ul# | 87       |
| 26) Benzo(a)pyrene          | 23.220 | 252  | 5796     | 0.388 | ng/ul# | 82       |
| 27) Indeno(1,2,3-cd)pyrene  | 25.449 | 276  | 7287     | 0.377 | ng/ul# | 95       |
| 28) Dibenzo(a,h)anthracene  | 25.461 | 278  | 5831     | 0.375 | ng/ul# | 90       |
| 29) Benzo(g,h,i)perylene    | 26.098 | 276  | 6182     | 0.375 | ng/ul# | 89       |
| -----                       |        |      |          |       |        |          |

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

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