

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM121024\  
 Data File : BM048922.D  
 Acq On : 10 Dec 2024 14:21  
 Operator : RC/JU  
 Sample : P5109-04MS  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BHJA2MS

Quant Time: Dec 10 15:22:05 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-SIM-BM120924.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Dec 09 17:52:26 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.586  | 152  | 3630     | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.348 | 136  | 9508     | 0.400 | ng/ul  | 0.00     |
| 9) Acenaphthene-d10         | 14.220 | 164  | 5089     | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 16.964 | 188  | 10414    | 0.400 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.157 | 240  | 8549     | 0.400 | ng/ul  | 0.00     |
| 23) Perylene-d12            | 23.326 | 264  | 8774     | 0.400 | ng/ul  | 0.00     |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.160  | 96   | 1308     | 0.308 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 11.948 | 152  | 4663     | 0.385 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 18.998 | 212  | 11121    | 0.435 | ng/ul  | 0.00     |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 2) 1,4-Dioxane              | 3.193  | 88   | 4280     | 0.872 | ng/ul# | 87       |
| 5) Naphthalene              | 10.398 | 128  | 9080     | 0.373 | ng/ul  | 100      |
| 7) 2-Methylnaphthalene      | 12.025 | 142  | 5327     | 0.358 | ng/ul  | 100      |
| 8) 1-Methylnaphthalene      | 12.240 | 142  | 5520     | 0.364 | ng/ul  | 100      |
| 10) Acenaphthylene          | 13.933 | 152  | 7956     | 0.370 | ng/ul  | 99       |
| 11) Acenaphthene            | 14.280 | 153  | 6250     | 0.384 | ng/ul  | 99       |
| 12) Fluorene                | 15.275 | 166  | 7072     | 0.391 | ng/ul  | 97       |
| 14) Pentachlorophenol       | 16.622 | 266  | 1270     | 0.592 | ng/ul  | 96       |
| 15) Phenanthrene            | 17.006 | 178  | 11707    | 0.410 | ng/ul  | 100      |
| 16) Anthracene              | 17.099 | 178  | 10170    | 0.400 | ng/ul  | 98       |
| 19) Fluoranthene            | 19.025 | 202  | 14430    | 0.407 | ng/ul  | 98       |
| 20) Pyrene                  | 19.388 | 202  | 15542    | 0.407 | ng/ul  | 97       |
| 21) Benzo(a)anthracene      | 21.143 | 228  | 12158    | 0.404 | ng/ul  | 98       |
| 22) Chrysene                | 21.195 | 228  | 15127    | 0.444 | ng/ul  | 99       |
| 24) Benzo(b)fluoranthene    | 22.683 | 252  | 13727    | 0.453 | ng/ul  | 99       |
| 25) Benzo(k)fluoranthene    | 22.727 | 252  | 15616    | 0.445 | ng/ul  | 96       |
| 26) Benzo(a)pyrene          | 23.235 | 252  | 11726    | 0.418 | ng/ul  | 100      |
| 27) Indeno(1,2,3-cd)pyrene  | 25.484 | 276  | 17078    | 0.434 | ng/ul# | 99       |
| 28) Dibenzo(a,h)anthracene  | 25.497 | 278  | 13185    | 0.430 | ng/ul  | 98       |
| 29) Benzo(g,h,i)perylene    | 26.134 | 276  | 14021    | 0.443 | ng/ul  | 98       |
| -----                       |        |      |          |       |        |          |

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM121024\  
Data File : BM048922.D  
Acq On : 10 Dec 2024 14:21  
Operator : RC/JU  
Sample : P5109-04MS  
Misc :  
ALS Vial : 9 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BHJA2MS

Quant Time: Dec 10 15:22:05 2024  
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-SIM-BM120924.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Mon Dec 09 17:52:26 2024  
Response via : Initial Calibration

