

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM092624\  
 Data File : BM047656.D  
 Acq On : 26 Sep 2024 22:47  
 Operator : RC/JU  
 Sample : P4164-02MS  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 P4164-01MS

Quant Time: Sep 27 00:27:50 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-SIM-BM092524.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Wed Sep 25 16:27:08 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.826  | 152  | 3689     | 0.400 | ng/ul  | 0.00     |
| 4) Naphthalene-d8           | 10.616 | 136  | 10667    | 0.400 | ng/ul  | 0.00     |
| 9) Acenaphthene-d10         | 14.454 | 164  | 5304     | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 17.190 | 188  | 10446    | 0.400 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.353 | 240  | 8689     | 0.400 | ng/ul  | 0.00     |
| 23) Perylene-d12            | 23.633 | 264  | 9691     | 0.400 | ng/ul  | # 0.00   |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.269  | 96   | 13596    | 2.869 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 12.205 | 152  | 4563     | 0.320 | ng/ul  | 0.00     |
| 18) Fluoranthene-d10        | 19.210 | 212  | 10911    | 0.410 | ng/ul  | 0.00     |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       |        | Qvalue   |
| 2) 1,4-Dioxane              | 3.303  | 88   | 5263     | 0.928 | ng/ul# | 87       |
| 5) Naphthalene              | 10.666 | 128  | 9915     | 0.341 | ng/ul  | 99       |
| 7) 2-Methylnaphthalene      | 12.282 | 142  | 5776     | 0.331 | ng/ul  | 98       |
| 8) 1-Methylnaphthalene      | 12.497 | 142  | 7012     | 0.383 | ng/ul  | 99       |
| 10) Acenaphthylene          | 14.172 | 152  | 8342     | 0.324 | ng/ul  | 100      |
| 11) Acenaphthene            | 14.519 | 153  | 6274     | 0.355 | ng/ul  | 99       |
| 12) Fluorene                | 15.500 | 166  | 6967     | 0.348 | ng/ul  | 99       |
| 14) Pentachlorophenol       | 16.840 | 266  | 2539     | 0.808 | ng/ul  | 98       |
| 15) Phenanthrene            | 17.233 | 178  | 11803    | 0.381 | ng/ul  | 99       |
| 16) Anthracene              | 17.321 | 178  | 10210    | 0.370 | ng/ul  | 99       |
| 19) Fluoranthene            | 19.242 | 202  | 14847    | 0.406 | ng/ul  | 99       |
| 20) Pyrene                  | 19.600 | 202  | 15825    | 0.407 | ng/ul  | 97       |
| 21) Benzo(a)anthracene      | 21.339 | 228  | 13993    | 0.381 | ng/ul  | 100      |
| 22) Chrysene                | 21.388 | 228  | 15190    | 0.397 | ng/ul  | 100      |
| 24) Benzo(b)fluoranthene    | 22.946 | 252  | 16261    | 0.386 | ng/ul  | 95       |
| 25) Benzo(k)fluoranthene    | 22.990 | 252  | 17051    | 0.402 | ng/ul  | 97       |
| 26) Benzo(a)pyrene          | 23.531 | 252  | 14467    | 0.435 | ng/ul# | 94       |
| 27) Indeno(1,2,3-cd)pyrene  | 25.953 | 276  | 21325    | 0.403 | ng/ul# | 100      |
| 28) Dibenzo(a,h)anthracene  | 25.967 | 278  | 16300    | 0.408 | ng/ul  | 98       |
| 29) Benzo(g,h,i)perylene    | 26.661 | 276  | 17414    | 0.410 | ng/ul  | 98       |
| -----                       |        |      |          |       |        |          |

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

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