

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN111124\  
 Data File : BN035003.D  
 Acq On : 11 Nov 2024 14:00  
 Operator : RC/JU  
 Sample : P4636-07  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 CC0P6

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 11/12/2024  
 Supervised By :mohammad ahmed 11/14/2024

Quant Time: Nov 11 14:56:03 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-SIM-BN110924.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Sat Nov 09 00:09:12 2024  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc  | Units  | Dev(Min) |
|-----------------------------|--------|------|----------|-------|--------|----------|
| -----                       |        |      |          |       |        |          |
| Internal Standards          |        |      |          |       |        |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.562  | 152  | 3196     | 0.400 | ng/ul  | -0.01    |
| 4) Naphthalene-d8           | 10.328 | 136  | 7911     | 0.400 | ng/ul  | -0.01    |
| 9) Acenaphthene-d10         | 14.198 | 164  | 5356     | 0.400 | ng/ul  | 0.00     |
| 13) Phenanthrene-d10        | 16.945 | 188  | 11863m   | 0.400 | ng/ul  | 0.00     |
| 17) Chrysene-d12            | 21.142 | 240  | 10576    | 0.400 | ng/ul  | # 0.00   |
| 23) Perylene-d12            | 23.308 | 264  | 10900m   | 0.400 | ng/ul  | 0.00     |
|                             |        |      |          |       |        |          |
| System Monitoring Compounds |        |      |          |       |        |          |
| 3) 1,4-Dioxane-d8           | 3.149  | 96   | 10265    | 3.375 | ng/ul  | 0.00     |
| 6) 2-Methylnaphthalene-d10  | 11.923 | 152  | 2918     | 0.241 | ng/ul  | -0.01    |
| 18) Fluoranthene-d10        | 18.981 | 212  | 7127     | 0.262 | ng/ul  | 0.00     |
|                             |        |      |          |       |        |          |
| Target Compounds            |        |      |          |       |        |          |
|                             |        |      |          |       | Qvalue |          |
| 5) Naphthalene              | 10.377 | 128  | 935      | 0.047 | ng/ul# | 85       |
| 7) 2-Methylnaphthalene      | 12.000 | 142  | 461      | 0.031 | ng/ul  | 98       |
| 8) 1-Methylnaphthalene      | 12.220 | 142  | 590      | 0.040 | ng/ul  | 97       |
| 10) Acenaphthylene          | 13.916 | 152  | 815      | 0.037 | ng/ul# | 74       |
| 12) Fluorene                | 15.253 | 166  | 1021     | 0.048 | ng/ul# | 96       |
| 14) Pentachlorophenol       | 16.603 | 266  | 58       | 0.021 | ng/ul# | 86       |
| 15) Phenanthrene            | 16.987 | 178  | 27014    | 0.852 | ng/ul  | 99       |
| 16) Anthracene              | 17.076 | 178  | 3206     | 0.108 | ng/ul# | 93       |
| 19) Fluoranthene            | 19.009 | 202  | 30059    | 0.851 | ng/ul# | 97       |
| 20) Pyrene                  | 19.372 | 202  | 30106    | 0.822 | ng/ul# | 95       |
| 21) Benzo(a)anthracene      | 21.128 | 228  | 8127     | 0.212 | ng/ul# | 92       |
| 22) Chrysene                | 21.177 | 228  | 10624    | 0.272 | ng/ul# | 93       |
| 24) Benzo(b)fluoranthene    | 22.665 | 252  | 11320m   | 0.295 | ng/ul  |          |
| 25) Benzo(k)fluoranthene    | 22.697 | 252  | 3389m    | 0.084 | ng/ul  |          |
| 26) Benzo(a)pyrene          | 23.214 | 252  | 6104m    | 0.178 | ng/ul  |          |
| 27) Indeno(1,2,3-cd)pyrene  | 25.456 | 276  | 4596m    | 0.110 | ng/ul  |          |
| 28) Dibenzo(a,h)anthracene  | 25.470 | 278  | 1191m    | 0.037 | ng/ul  |          |
| 29) Benzo(g,h,i)perylene    | 26.110 | 276  | 5404m    | 0.169 | ng/ul  |          |
| -----                       |        |      |          |       |        |          |

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

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