

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN061224\  
 Data File : BN031924.D  
 Acq On : 12 Jun 2024 00:14  
 Operator : MA/JU  
 Sample : P2659-15  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 BHC36

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 06/12/2024  
 Supervised By :mohammad ahmed 06/14/2024

Quant Time: Jun 12 00:45:26 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN052924.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jun 11 22:28:10 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| -----                         |        |      |          |        |        |          |
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.658  | 152  | 349111   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.440 | 136  | 1568764  | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.304 | 164  | 977085   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.051 | 188  | 1777240  | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.286 | 240  | 1231533  | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.215 | 264  | 1433283  | 20.000 | ng/u1  | 0.00     |
|                               |        |      |          |        |        |          |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.181  | 96   | 45166    | 5.352  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.587  | 84   | 344591   | 14.506 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 6.840  | 99   | 1014315  | 33.891 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.005  | 67   | 581499   | 33.217 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.193  | 132  | 824388   | 36.100 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.369  | 113  | 780082   | 31.975 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 8.822  | 128  | 434459   | 36.221 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.534  | 143  | 473989   | 38.254 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.069 | 165  | 817348   | 35.335 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.587 | 131  | 892140   | 25.976 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.722 | 166  | 2413723  | 35.460 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 13.998 | 160  | 2763116  | 35.303 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.522 | 143  | 349952   | 25.677 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.304 | 176  | 1983203  | 34.434 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.434 | 200  | 275796   | 30.873 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.151 | 188  | 2668485  | 35.071 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.451 | 212  | 2363947  | 35.871 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.015 | 264  | 1811860  | 26.474 | ng/u1  | 0.00     |
|                               |        |      |          |        |        |          |
| Target Compounds              |        |      |          |        | Qvalue |          |
| 50) Acenaphthylene            | 14.028 | 152  | 293608   | 3.319  | ng/u1  | 99       |
| 61) Fluorene                  | 15.357 | 166  | 69298    | 1.082  | ng/u1# | 99       |
| 72) Phenanthrene              | 17.092 | 178  | 1718393  | 18.980 | ng/u1  | 100      |
| 74) Anthracene                | 17.186 | 178  | 310703   | 3.387  | ng/u1  | 99       |
| 77) Carbazole                 | 17.463 | 167  | 122797   | 1.570  | ng/u1  | 99       |
| 80) Fluoranthene              | 19.122 | 202  | 3180348  | 40.381 | ng/u1  | 96       |
| 82) Pyrene                    | 19.486 | 202  | 2498340  | 30.917 | ng/u1  | 95       |
| 83) Butylbenzylphthalate      | 20.386 | 149  | 46653    | 1.222  | ng/u1  | 95       |
| 85) Benzo(a)anthracene        | 21.269 | 228  | 1429898  | 17.346 | ng/u1  | 98       |
| 86) Bis(2-ethylhexyl)phtha... | 21.198 | 149  | 230011   | 4.134  | ng/u1  | 100      |
| 87) Chrysene                  | 21.333 | 228  | 1600527  | 20.787 | ng/u1  | 98       |
| 90) Benzo(b)fluoranthene      | 23.292 | 252  | 2028841  | 23.682 | ng/u1  | 99       |
| 91) Benzo(k)fluoranthene      | 23.351 | 252  | 688507m  | 8.190  | ng/u1  |          |
| 93) Benzo(a)pyrene            | 24.074 | 252  | 861659   | 10.739 | ng/u1  | 96       |
| 94) Indeno(1,2,3-cd)pyrene    | 27.504 | 276  | 795705   | 8.606  | ng/u1  | 96       |
| 95) Dibenzo(a,h)anthracene    | 27.545 | 278  | 266511   | 3.531  | ng/u1# | 90       |
| 96) Benzo(g,h,i)perylene      | 28.527 | 276  | 117304m  | 1.596  | ng/u1  |          |
| -----                         |        |      |          |        |        |          |

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

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