

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN052424\  
 Data File : BN031663.D  
 Acq On : 24 May 2024 16:19  
 Operator : MA/JU  
 Sample : P2548-16  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 BH611

Manual Integrations  
 APPROVED

Reviewed By :Yogesh Patel 05/25/2024  
 Supervised By :mohammad ahmed 05/27/2024

Quant Time: May 25 04:42:04 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN052124.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 21 23:42:49 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|--------|--------|----------|
| Internal Standards            |        |      |          |        |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.693  | 152  | 412687   | 20.000 | ng/u1  | 0.00     |
| 20) Naphthalene-d8            | 10.475 | 136  | 1591548  | 20.000 | ng/u1  | 0.00     |
| 38) Acenaphthene-d10          | 14.328 | 164  | 880971   | 20.000 | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.069 | 188  | 1740244  | 20.000 | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.304 | 240  | 1659542  | 20.000 | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.239 | 264  | 1983205  | 20.000 | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |        |        |          |
| 3) 1,4-Dioxane-d8             | 3.205  | 96   | 32436    | 3.115  | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.605  | 84   | 446701   | 15.374 | ng/u1  | 0.00     |
| 7) Phenol-d5                  | 6.858  | 99   | 593068   | 17.458 | ng/u1  | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.028  | 67   | 351343   | 17.244 | ng/u1  | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.222  | 132  | 474158   | 18.113 | ng/u1  | 0.00     |
| 15) 4-Methylphenol-d8         | 8.387  | 113  | 449057   | 17.075 | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 8.840  | 128  | 223759   | 18.577 | ng/u1  | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.557  | 143  | 211779   | 18.412 | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.087 | 165  | 408766   | 18.176 | ng/u1  | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.610 | 131  | 614399   | 18.952 | ng/u1  | 0.00     |
| 46) Dimethylphthalate-d6      | 13.740 | 166  | 1188654  | 21.410 | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 14.016 | 160  | 1454853  | 20.908 | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.522 | 143  | 157772   | 15.019 | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.322 | 176  | 1043674  | 21.692 | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.439 | 200  | 78311    | 10.408 | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.169 | 188  | 1574215  | 21.369 | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.469 | 212  | 1949377  | 18.727 | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 24.039 | 264  | 2092330  | 22.413 | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |        |        |          |
|                               |        |      |          |        | Qvalue |          |
| 72) Phenanthrene              | 17.110 | 178  | 235467   | 2.666  | ng/u1  | 100      |
| 80) Fluoranthene              | 19.127 | 202  | 617679   | 4.801  | ng/u1  | 100      |
| 82) Pyrene                    | 19.492 | 202  | 604481   | 4.581  | ng/u1  | 97       |
| 85) Benzo(a)anthracene        | 21.286 | 228  | 349563   | 3.132  | ng/u1  | 94       |
| 87) Chrysene                  | 21.345 | 228  | 360675   | 3.499  | ng/u1  | 94       |
| 90) Benzo(b)fluoranthene      | 23.310 | 252  | 478616   | 4.136  | ng/u1  | 97       |
| 91) Benzo(k)fluoranthene      | 23.363 | 252  | 207435m  | 1.758  | ng/u1  |          |
| 93) Benzo(a)pyrene            | 24.098 | 252  | 364018   | 3.274  | ng/u1  | 98       |
| 94) Indeno(1,2,3-cd)pyrene    | 27.533 | 276  | 248393   | 1.862  | ng/u1  | 95       |
| 96) Benzo(g,h,i)perylene      | 28.556 | 276  | 238169   | 2.161  | ng/u1  | 96       |

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

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