

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN043024\  
 Data File : BN031326.D  
 Acq On : 30 Apr 2024 18:36  
 Operator : MA/JU  
 Sample : P2161-20  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 C0RD7

Manual Integrations  
 APPROVED

Reviewed By :Jagrut Upadhyay 05/01/2024  
 Supervised By :mohammad ahmed 05/02/2024

Quant Time: Apr 30 23:43:43 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN042924.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Apr 30 23:30:44 2024  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc    | Units  | Dev(Min) |
|-------------------------------|--------|------|----------|---------|--------|----------|
| -----                         |        |      |          |         |        |          |
| Internal Standards            |        |      |          |         |        |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.634  | 152  | 122468m  | 20.000  | ng/u1  | 0.02     |
| 20) Naphthalene-d8            | 10.422 | 136  | 793322   | 20.000  | ng/u1  | 0.03     |
| 38) Acenaphthene-d10          | 14.263 | 164  | 522234   | 20.000  | ng/u1  | 0.00     |
| 64) Phenanthrene-d10          | 17.016 | 188  | 1119047  | 20.000  | ng/u1  | 0.00     |
| 79) Chrysene-d12              | 21.251 | 240  | 942243   | 20.000  | ng/u1  | 0.00     |
| 88) Perylene-d12              | 24.139 | 264  | 863340   | 20.000  | ng/u1  | 0.00     |
| System Monitoring Compounds   |        |      |          |         |        |          |
| 3) 1,4-Dioxane-d8             | 3.134  | 96   | 20678    | 7.861   | ng/uL  | 0.00     |
| 4) Pyridine-d5                | 3.546  | 84   | 188866   | 25.072  | ng/u1  | 0.01     |
| 7) Phenol-d5                  | 6.799  | 99   | 158769   | 16.111  | ng/u1  | 0.01     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.975  | 67   | 450703   | 78.109  | ng/u1  | 0.02     |
| 11) 2-Chlorophenol-d4         | 7.157  | 132  | 467232   | 59.263  | ng/u1  | 0.01     |
| 15) 4-Methylphenol-d8         | 8.328  | 113  | 285735   | 34.988  | ng/u1  | 0.00     |
| 21) Nitrobenzene-d5           | 8.781  | 128  | 293497   | 51.334  | ng/u1  | 0.01     |
| 24) 2-Nitrophenol-d4          | 9.493  | 143  | 308118   | 50.073  | ng/u1  | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.028 | 165  | 519697   | 45.481  | ng/u1  | 0.01     |
| 31) 4-Chloroaniline-d4        | 10.551 | 131  | 727887   | 43.956  | ng/u1  | 0.01     |
| 46) Dimethylphthalate-d6      | 13.681 | 166  | 1812693  | 52.650  | ng/u1  | 0.00     |
| 49) Acenaphthylene-d8         | 13.957 | 160  | 2064940  | 51.690  | ng/u1  | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.481 | 143  | 60669    | 9.710   | ng/u1  | 0.00     |
| 60) Fluorene-d10              | 15.263 | 176  | 1585424  | 53.925  | ng/u1  | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.386 | 200  | 224093   | 38.941  | ng/u1  | 0.00     |
| 73) Anthracene-d10            | 17.116 | 188  | 2511067  | 54.372  | ng/u1  | 0.00     |
| 81) Pyrene-d10                | 19.422 | 212  | 2855933  | 49.768  | ng/u1  | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.945 | 264  | 2236198  | 55.948  | ng/u1  | 0.00     |
| Target Compounds              |        |      |          |         |        |          |
|                               |        |      |          |         | Qvalue |          |
| 12) 2-Chlorophenol            | 7.193  | 128  | 14326    | 1.748   | ng/u1  | 99       |
| 30) Naphthalene               | 10.463 | 128  | 69150    | 1.774   | ng/u1  | 99       |
| 39) 1,2,4,5-Tetrachloroben... | 12.434 | 216  | 1098415  | 81.493  | ng/u1  | 98       |
| 75) 1,2,3,4-Tetrachloroben... | 13.057 | 216  | 1949550  | 137.540 | ng/uL  | 99       |
| 76) Pentachlorobenzene        | 14.581 | 250  | 33638    | 2.420   | ng/uL  | 94       |
| -----                         |        |      |          |         |        |          |

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

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