

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN100224\  
 Data File : BN034336.D  
 Acq On : 02 Oct 2024 14:20  
 Operator : RC/JU  
 Sample : P4016-14  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 DCVW5

Quant Time: Oct 02 14:57:08 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN091124.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Sep 30 14:14:38 2024  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.387  | 152  | 229133   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.140 | 136  | 895037   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.040 | 164  | 517079   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 16.798 | 188  | 999968   | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.027 | 240  | 922372   | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.739 | 264  | 1142260  | 20.000 | ng/ul | 0.00     |

|                               |        |     |         |        |       |      |
|-------------------------------|--------|-----|---------|--------|-------|------|
| System Monitoring Compounds   |        |     |         |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.005  | 96  | 18500   | 3.295  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.393  | 84  | 216401  | 12.966 | ng/ul | 0.00 |
| 7) Phenol-d5                  | 6.587  | 99  | 296469  | 14.579 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 6.740  | 67  | 185884  | 14.733 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4         | 6.922  | 132 | 265079  | 16.560 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8         | 8.099  | 113 | 240304  | 14.361 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5           | 8.528  | 128 | 121196  | 16.989 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.240  | 143 | 122332  | 16.676 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 9.775  | 165 | 225093  | 16.293 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4        | 10.287 | 131 | 325932  | 15.583 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6      | 13.463 | 166 | 694840  | 17.981 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8         | 13.722 | 160 | 783812  | 18.030 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.269 | 143 | 61794   | 7.792  | ng/ul | 0.00 |
| 60) Fluorene-d10              | 15.040 | 176 | 585683  | 17.883 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.175 | 200 | 50995   | 8.553  | ng/ul | 0.00 |
| 73) Anthracene-d10            | 16.892 | 188 | 869019  | 18.485 | ng/ul | 0.00 |
| 81) Pyrene-d10                | 19.204 | 212 | 987022  | 19.206 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12        | 23.557 | 264 | 1102577 | 18.791 | ng/ul | 0.00 |

|                  |        |
|------------------|--------|
| Target Compounds | Qvalue |
| -----            | -----  |

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

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