

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080322\  
 Data File : BP011254.D  
 Acq On : 03 Aug 2022 20:26  
 Operator : CG/JU  
 Sample : N3900-14  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0CA2

Quant Time: Aug 03 23:01:43 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP080222.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Aug 02 23:52:37 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.599  | 152  | 196096   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.393 | 136  | 859195   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.275 | 164  | 417563   | 20.000 | ng/u1 | # 0.00   |
| 64) Phenanthrene-d10          | 17.045 | 188  | 814170   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.174 | 240  | 892549   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.486 | 264  | 930656   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.093  | 96   | 5541     | 2.285  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.517  | 84   | 23677    | 3.619  | ng/u1 | 0.01     |
| 7) Phenol-d5                  | 6.787  | 99   | 63981    | 3.915  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.946  | 67   | 303827   | 25.765 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.134  | 132  | 232975   | 17.490 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.322  | 113  | 131211   | 9.758  | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.763  | 128  | 176689   | 28.590 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.487  | 143  | 134230   | 23.172 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.022 | 165  | 234484   | 21.227 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.546 | 131  | 214933   | 12.577 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.698 | 166  | 896842   | 31.359 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.963 | 160  | 1125516  | 31.280 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.510 | 143  | 6217     | 1.278  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.275 | 176  | 817031   | 33.575 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.410 | 200  | 71422    | 18.892 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.145 | 188  | 1362167  | 38.068 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.404 | 212  | 1607743  | 35.627 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.339 | 264  | 1712329  | 36.923 | ng/u1 | 0.00     |

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

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

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