





Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080322\  
 Data File : BP011251.D  
 Acq On : 03 Aug 2022 18:42  
 Operator : CG/JU  
 Sample : N3900-04  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0C94

Manual IntegrationsAPPROVED

Quant Time: Aug 03 23:01:15 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

Reviewed By :Jagrut Upadhyay 08/04/2022  
 Supervised By :Sohil Jodhani 08/10/2022

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.599  | 152  | 132440   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.393 | 136  | 532430   | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.275 | 164  | 274677   | 20.000 | ng/ul | # 0.00   |
| 64) Phenanthrene-d10          | 17.045 | 188  | 495665   | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.174 | 240  | 630239   | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 23.492 | 264  | 715852   | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.093  | 96   | 7928     | 4.840  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.511  | 84   | 30971    | 7.010  | ng/ul | 0.00     |
| 7) Phenol-d5                  | 6.781  | 99   | 31844    | 2.885  | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.946  | 67   | 212892   | 26.731 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.134  | 132  | 129887   | 14.438 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.322  | 113  | 60876    | 6.703  | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 8.763  | 128  | 118071   | 30.831 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.487  | 143  | 72333    | 20.150 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.022 | 165  | 125712   | 18.364 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.546 | 131  | 206775   | 19.525 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.698 | 166  | 461150   | 24.513 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 13.963 | 160  | 652698   | 27.575 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.504 | 143  | 5568m    | 1.741  | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.275 | 176  | 519479   | 32.452 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.410 | 200  | 45471    | 19.757 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.145 | 188  | 949779   | 43.600 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.404 | 212  | 1049527  | 32.937 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.339 | 264  | 1227256  | 34.404 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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