

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122222\  
 Data File : BP013254.D  
 Acq On : 22 Dec 2022 20:43  
 Operator : CG/JU  
 Sample : N6015-08  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :

Quant Time: Dec 22 21:49:14 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Dec 21 23:38:37 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.946  | 152  | 684502   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.757 | 136  | 2905052  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.587 | 164  | 1774921  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.345 | 188  | 3366570  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.427 | 240  | 2293402  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.910 | 264  | 2721335  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.334  | 96   | 68169    | 4.258  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.764  | 84   | 933097   | 20.518 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.105  | 99   | 1272448  | 22.822 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.275  | 67   | 830313   | 24.687 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.475  | 132  | 1059118  | 24.131 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.657  | 113  | 940634   | 20.590 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.116  | 128  | 527977   | 24.736 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.840  | 143  | 599041   | 24.929 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.381 | 165  | 1105311  | 23.682 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.898 | 131  | 1205029  | 18.288 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.992 | 166  | 3165197  | 23.203 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.281 | 160  | 3642750  | 24.302 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.787 | 143  | 326293   | 13.298 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.581 | 176  | 2702973  | 23.422 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.698 | 200  | 332647   | 15.354 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.445 | 188  | 3642418  | 23.982 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.674 | 212  | 3463664  | 26.311 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.745 | 264  | 3300047  | 24.329 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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