

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP062422\  
 Data File : BP010807.D  
 Acq On : 24 Jun 2022 22:55  
 Operator : CG/JU  
 Sample : N3401-15  
 Misc :  
 ALS Vial : 26 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0AG1

Quant Time: Jun 24 23:41:30 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP062322.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 24 22:42:26 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.722  | 152  | 340667   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.516 | 136  | 1391867  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.375 | 164  | 783807   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.139 | 188  | 1567746  | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.251 | 240  | 1556058  | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.621 | 264  | 1615386  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.205  | 96   | 48881    | 5.428  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.610  | 84   | 289241   | 12.066 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.887  | 99   | 234885   | 8.710  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.063  | 67   | 693870   | 39.343 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.251  | 132  | 691485   | 30.906 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.428  | 113  | 437646   | 20.529 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.881  | 128  | 412288   | 46.531 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.604  | 143  | 361016   | 49.248 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.134 | 165  | 666210   | 36.781 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.657 | 131  | 998483   | 33.771 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.798 | 166  | 2328953  | 43.377 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.069 | 160  | 2850942  | 41.604 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.575 | 143  | 62183    | 7.415  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.375 | 176  | 1981919  | 43.015 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.492 | 200  | 176309   | 33.944 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.239 | 188  | 3055661  | 43.808 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.486 | 212  | 3484357  | 41.101 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.468 | 264  | 3557221  | 44.563 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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