

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM072622\  
 Data File : BM035924.D  
 Acq On : 26 Jul 2022 13:59  
 Operator : CG/JU  
 Sample : N3801-22  
 Misc :  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 GBHY7

Quant Time: Jul 26 14:55:36 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM072522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jul 25 14:55:39 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.887  | 152  | 391273   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.692 | 136  | 1758661  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.521 | 164  | 1123879  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.262 | 188  | 2388709  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.433 | 240  | 2323200  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.809 | 264  | 2364070  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.275  | 96   | 58534    | 5.629  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.699  | 84   | 230513   | 8.092  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.045  | 99   | 210151   | 6.405  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.216  | 67   | 782624   | 35.779 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.416  | 132  | 698802   | 25.658 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.592  | 113  | 426780   | 16.055 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.051  | 128  | 509572   | 36.566 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.775  | 143  | 472391   | 34.793 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.310 | 165  | 770852   | 31.391 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.833 | 131  | 1267955  | 30.501 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.939 | 166  | 3242852  | 42.359 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.216 | 160  | 3912480  | 40.055 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.715 | 143  | 85218    | 5.421  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.510 | 176  | 2898190  | 41.894 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.627 | 200  | 412537   | 32.129 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.356 | 188  | 4565514  | 42.275 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.650 | 212  | 5314979  | 41.279 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.656 | 264  | 5147125  | 42.885 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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