

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM083021\  
 Data File : BM031995.D  
 Acq On : 31 Aug 2021 21:30  
 Operator : CG/JU  
 Sample : M3517-04  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C0AB7

Quant Time: Sep 01 01:04:52 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM082821.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Aug 30 09:21:30 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.504  | 152  | 87782    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.263 | 136  | 380214   | 20.000 | ng/u1 | # 0.00   |
| 38) Acenaphthene-d10          | 14.145 | 164  | 282197   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 16.904 | 188  | 598102   | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.109 | 240  | 567727   | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.244 | 264  | 495271   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.134  | 96   | 9593     | 5.042  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.528  | 84   | 54969    | 10.174 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.699  | 99   | 73969    | 10.162 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.863  | 67   | 53380    | 12.674 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.046  | 132  | 66481    | 11.868 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.216  | 113  | 67875    | 10.868 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.657  | 128  | 38174    | 13.823 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.363  | 143  | 37568    | 11.495 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.898  | 165  | 77594    | 11.833 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.416 | 131  | 110047   | 11.897 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.575 | 166  | 297198   | 13.918 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.833 | 160  | 346771   | 13.713 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.398 | 143  | 27649    | 7.154  | ng/u1 | 0.01     |
| 60) Fluorene-d10              | 15.145 | 176  | 261827   | 13.949 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.292 | 200  | 19215    | 4.676  | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 16.998 | 188  | 420587   | 14.815 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.304 | 212  | 473501   | 17.806 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.109 | 264  | 373572   | 14.352 | ng/u1 | 0.00     |

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

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

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