

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM083021\  
 Data File : BM031958.D  
 Acq On : 30 Aug 2021 20:21  
 Operator : CG/JU  
 Sample : PB138565BL  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SBLK565

Quant Time: Aug 31 02:36:43 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     | 7.510  | 152  | 102019   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.269 | 136  | 442423   | 20.000 | ng/u1 | # 0.00   |
| 38) Acenaphthene-d10          | 14.151 | 164  | 331758   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 16.903 | 188  | 730813   | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.109 | 240  | 751850   | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.250 | 264  | 733850   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.134  | 96   | 15878    | 7.181  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.522  | 84   | 227363   | 36.211 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.704  | 99   | 291648   | 34.474 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.869  | 67   | 178330   | 36.433 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.045  | 132  | 244374   | 37.539 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.222  | 113  | 251089   | 34.594 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.663  | 128  | 127474   | 39.669 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.369  | 143  | 153698   | 40.415 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.898  | 165  | 275674   | 36.130 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.422 | 131  | 377370   | 35.062 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.580 | 166  | 996986   | 39.715 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.839 | 160  | 1189189  | 40.000 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.386 | 143  | 157838   | 34.737 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.151 | 176  | 861237   | 39.028 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.298 | 200  | 171765   | 34.205 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.003 | 188  | 1422541  | 41.010 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.309 | 212  | 1642581  | 46.643 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.121 | 264  | 1573548  | 40.799 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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