

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM081522\  
 Data File : BM036279.D  
 Acq On : 15 Aug 2022 10:30  
 Operator : CG/JU  
 Sample : PB146648BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SBLK648

Quant Time: Aug 15 11:15:36 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM072522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun Aug 14 23:34:05 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.804  | 152  | 167349   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.604 | 136  | 680159   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.451 | 164  | 373612   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.192 | 188  | 732363   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.374 | 240  | 685146   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.709 | 264  | 741832   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.210  | 96   | 27722    | 6.233  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.640  | 84   | 300181   | 24.638 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.992  | 99   | 417277   | 29.734 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.145  | 67   | 271012   | 28.968 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.339  | 132  | 354975   | 30.473 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.533  | 113  | 320304   | 28.173 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.975  | 128  | 165862   | 30.774 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.698  | 143  | 163806   | 31.195 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.239 | 165  | 283724   | 29.875 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.757 | 131  | 492431   | 30.629 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.868 | 166  | 801105   | 31.478 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.145 | 160  | 1016207  | 31.296 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.680 | 143  | 121870   | 23.320 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.439 | 176  | 728259   | 31.667 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.568 | 200  | 93149    | 23.662 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.292 | 188  | 1122655  | 33.906 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.586 | 212  | 1272399  | 33.509 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.556 | 264  | 1313321  | 34.871 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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