

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP082422\  
 Data File : BP011558.D  
 Acq On : 25 Aug 2022 10:43  
 Operator : CG/JU  
 Sample : PB147280BL  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK280

Quant Time: Aug 26 01:02:19 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP081622.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Aug 25 02:14:54 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.552  | 152  | 64217    | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.340 | 136  | 279807   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.228 | 164  | 164487   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 16.998 | 188  | 324979   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.127 | 240  | 281211   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.416 | 264  | 284514   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.064  | 96   | 12913    | 6.240  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.475  | 84   | 114628   | 21.635 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.734  | 99   | 179942   | 30.492 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 6.899  | 67   | 122560   | 31.967 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.081  | 132  | 148293   | 34.166 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.275  | 113  | 146184   | 30.949 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.716  | 128  | 69867    | 34.684 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.434  | 143  | 73317    | 35.762 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 9.969  | 165  | 120636   | 31.878 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.493 | 131  | 141595   | 22.477 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.651 | 166  | 420255   | 36.123 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 13.916 | 160  | 447909   | 34.640 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.457 | 143  | 57873    | 25.819 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.228 | 176  | 339386   | 34.676 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.369 | 200  | 50399    | 28.700 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.098 | 188  | 532770   | 37.016 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.363 | 212  | 559185   | 39.302 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.269 | 264  | 515319   | 37.234 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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