

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP072622\  
 Data File : BP011102.D  
 Acq On : 26 Jul 2022 10:28  
 Operator : CG/JU  
 Sample : PB146402BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK402

Quant Time: Jul 26 14:36:31 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP071422.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jul 26 14:31:36 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.663  | 152  | 434995   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.457 | 136  | 2019464  | 20.000 | ng/u1 | # 0.00   |
| 38) Acenaphthene-d10          | 14.328 | 164  | 1239205  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.092 | 188  | 2426920  | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.210 | 240  | 1694684  | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.557 | 264  | 1431007  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.164  | 96   | 74883    | 6.206  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.569  | 84   | 1054015  | 33.796 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.840  | 99   | 1244606  | 34.856 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.005  | 67   | 864619   | 37.656 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.193  | 132  | 1036597  | 35.927 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.381  | 113  | 1057962  | 37.010 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.828  | 128  | 537200   | 36.502 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.546  | 143  | 567790   | 38.990 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.081 | 165  | 913081   | 34.286 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.604 | 131  | 1522517  | 34.678 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.751 | 166  | 3451497  | 40.752 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.016 | 160  | 3883160  | 38.859 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.545 | 143  | 515332   | 30.064 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.328 | 176  | 2719704  | 37.778 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.451 | 200  | 431196   | 35.906 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.192 | 188  | 4047416  | 39.229 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.451 | 212  | 4141900  | 46.536 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.404 | 264  | 2777699  | 40.415 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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