

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP041922\  
 Data File : BP009968.D  
 Acq On : 20 Apr 2022 08:53  
 Operator : CG/JU  
 Sample : PB144232BL  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK232

Quant Time: Apr 21 01:01:49 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP040722.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Apr 20 01:32:14 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.952  | 152  | 120549   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.757 | 136  | 533531   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.581 | 164  | 387922   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.333 | 188  | 885153   | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.416 | 240  | 987011   | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.874 | 264  | 977606   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.375  | 96   | 16534    | 6.121  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.787  | 84   | 214461   | 28.797 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.105  | 99   | 279811   | 26.439 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.275  | 67   | 190924   | 30.302 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.481  | 132  | 227329   | 28.476 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.652  | 113  | 243249   | 27.533 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.110  | 128  | 118007   | 30.671 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.834  | 143  | 129046   | 30.593 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.375 | 165  | 224214   | 25.402 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.887 | 131  | 342464   | 27.241 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.987 | 166  | 934776   | 30.677 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.275 | 160  | 978607   | 27.108 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.769 | 143  | 147796   | 26.744 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.575 | 176  | 772361   | 30.130 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.687 | 200  | 137952   | 26.308 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.433 | 188  | 1292028  | 30.486 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.663 | 212  | 1637301  | 30.373 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.715 | 264  | 1559347  | 30.629 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

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