

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080921\  
 Data File : BP006606.D  
 Acq On : 09 Aug 2021 10:17  
 Operator : CG/JU  
 Sample : PB138195BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK195

Quant Time: Aug 09 11:21:26 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP072421.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Aug 07 13:59:25 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.752  | 152  | 213930   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.534 | 136  | 953548   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.381 | 164  | 530602   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.134 | 188  | 999167   | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.233 | 240  | 907070   | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.563 | 264  | 890392   | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.275  | 96   | 31747    | 5.539  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.681  | 84   | 455030   | 26.745 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.934  | 99   | 532176   | 26.487 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.099  | 67   | 341764   | 27.338 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.287  | 132  | 426246   | 28.555 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.463  | 113  | 424183   | 26.846 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.910  | 128  | 213463   | 30.065 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.628  | 143  | 217351   | 32.061 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.157 | 165  | 350633   | 26.394 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.681 | 131  | 561918   | 25.692 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.804 | 166  | 1077778  | 28.540 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.075 | 160  | 1368768  | 29.501 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.598 | 143  | 170756   | 22.773 | ng/u1 | 0.01     |
| 60) Fluorene-d10              | 15.375 | 176  | 881630   | 27.771 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.504 | 200  | 106617   | 19.141 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.234 | 188  | 1359243  | 30.136 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.480 | 212  | 1435676  | 30.772 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.416 | 264  | 1435188  | 31.719 | ng/u1 | 0.00     |

Target Compounds Qvalue

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

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP080921\  
Data File : BP006606.D  
Acq On : 09 Aug 2021 10:17  
Operator : CG/JU  
Sample : PB138195BL  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SBLK195

Quant Time: Aug 09 11:21:26 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP072421.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sat Aug 07 13:59:25 2021  
Response via : Initial Calibration

