

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP062722\  
 Data File : BP010889.D  
 Acq On : 27 Jun 2022 22:11  
 Operator : CG/JU  
 Sample : PB145713BL  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :

Quant Time: Jun 27 23:17:37 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP062322.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jun 27 23:12:49 2022  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.710  | 152  | 471276   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.504 | 136  | 1942450  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.363 | 164  | 1014871  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.127 | 188  | 1811790  | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12              | 21.239 | 240  | 1504762  | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12              | 23.604 | 264  | 1630553  | 20.000 | ng/u1 | -0.01    |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.205  | 96   | 74042    | 5.943  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.611  | 84   | 977854   | 29.486 | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 6.881  | 99   | 1109943  | 29.752 | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.052  | 67   | 729354   | 29.894 | ng/u1 | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.240  | 132  | 929204   | 30.021 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.422  | 113  | 877094   | 29.741 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 8.869  | 128  | 415603   | 33.610 | ng/u1 | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.587  | 143  | 381442   | 37.285 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.122 | 165  | 706328   | 27.943 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.645 | 131  | 1190415  | 28.850 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.786 | 166  | 2025588  | 29.137 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.057 | 160  | 2440109  | 27.502 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.575 | 143  | 289973   | 26.705 | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.363 | 176  | 1725631  | 28.925 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.481 | 200  | 153778   | 25.618 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.227 | 188  | 2464003  | 30.567 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.474 | 212  | 2587490  | 31.562 | ng/u1 | -0.01    |
| 92) Benzo(a)pyrene-d12        | 23.451 | 264  | 2569297  | 31.887 | ng/u1 | 0.00     |

| Target Compounds | Qvalue |
|------------------|--------|
| -----            |        |

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

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