

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP062722\  
 Data File : BP010888.D  
 Acq On : 27 Jun 2022 21:37  
 Operator : CG/JU  
 Sample : PB145835BL  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK835

Quant Time: Jun 27 23:17:28 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 | 7.711  | 152  | 555546   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.505 | 136  | 2296109  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.363 | 164  | 1228303  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.128 | 188  | 2171175  | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12          | 21.239 | 240  | 1621237  | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12          | 23.604 | 264  | 1686774  | 20.000 | ng/u1 | -0.01    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.205  | 96  | 90561   | 6.166  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.611  | 84  | 1186106 | 30.341 | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 6.881  | 99  | 1392337 | 31.660 | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.052  | 67  | 905007  | 31.467 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.240  | 132 | 1165153 | 31.934 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.422  | 113 | 1110452 | 31.942 | ng/u1 | 0.00  |
| 21) Nitrobenzene-d5           | 8.869  | 128 | 531776  | 36.381 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.587  | 143 | 508895  | 42.082 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.128 | 165 | 921072  | 30.826 | ng/u1 | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.646 | 131 | 1518285 | 31.129 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 13.787 | 166 | 2702805 | 32.123 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 14.057 | 160 | 3272386 | 30.473 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.575 | 143 | 384039  | 29.222 | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.363 | 176 | 2245423 | 31.098 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.481 | 200 | 210540  | 29.269 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.228 | 188 | 3170199 | 32.818 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.475 | 212 | 3138998 | 35.539 | ng/u1 | -0.01 |
| 92) Benzo(a)pyrene-d12        | 23.451 | 264 | 2841177 | 34.086 | ng/u1 | 0.00  |

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

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

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