

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP102121\  
 Data File : BP007616.D  
 Acq On : 23 Oct 2021 06:31  
 Operator : CG/JU  
 Sample : M4282-03  
 Misc :  
 ALS Vial : 76 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BFYA9

Quant Time: Oct 23 07:20:48 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP101421.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 20 10:49:33 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.946  | 152  | 188072   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8            | 10.752 | 136  | 763457   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10          | 14.581 | 164  | 500836   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10          | 17.334 | 188  | 1084004  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12              | 21.422 | 240  | 1152525  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12              | 23.868 | 264  | 1194515  | 20.000 | ng/u1 | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.375  | 96   | 22475    | 4.720  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.793  | 84   | 112566   | 8.636  | ng/u1 | 0.00     |
| 7) Phenol-d5                  | 7.111  | 99   | 122497   | 8.340  | ng/u1 | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.269  | 67   | 304680   | 34.780 | ng/u1 | -0.01    |
| 11) 2-Chlorophenol-d4         | 7.475  | 132  | 354117   | 30.648 | ng/u1 | 0.00     |
| 15) 4-Methylphenol-d8         | 8.652  | 113  | 231688   | 20.479 | ng/u1 | 0.00     |
| 21) Nitrobenzene-d5           | 9.099  | 128  | 222042   | 45.957 | ng/u1 | -0.01    |
| 24) 2-Nitrophenol-d4          | 9.828  | 143  | 239486   | 53.066 | ng/u1 | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.369 | 165  | 435237   | 39.540 | ng/u1 | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.881 | 131  | 495591   | 33.294 | ng/u1 | 0.00     |
| 46) Dimethylphthalate-d6      | 13.993 | 166  | 1434002  | 41.386 | ng/u1 | 0.00     |
| 49) Acenaphthylene-d8         | 14.275 | 160  | 1693112  | 38.881 | ng/u1 | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.775 | 143  | 48870    | 8.790  | ng/u1 | 0.00     |
| 60) Fluorene-d10              | 15.575 | 176  | 1216079  | 41.938 | ng/u1 | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.687 | 200  | 104931   | 24.158 | ng/u1 | 0.00     |
| 73) Anthracene-d10            | 17.434 | 188  | 2031456  | 44.128 | ng/u1 | 0.00     |
| 81) Pyrene-d10                | 19.663 | 212  | 2461108  | 42.715 | ng/u1 | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.710 | 264  | 2514367  | 42.303 | ng/u1 | 0.00     |

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

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

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