

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP072921\  
 Data File : BP006432.D  
 Acq On : 29 Jul 2021 21:41  
 Operator : CG/JU  
 Sample : M3141-04  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BGEP4

Quant Time: Jul 30 01:15:12 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP072421.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jul 29 11:29:11 2021  
 Response via : Initial Calibration

| Compound                      | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| -----                         |        |      |          |        |       |          |
| Internal Standards            |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4     | 7.781  | 152  | 313352   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.563 | 136  | 1320104  | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.404 | 164  | 739120   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.157 | 188  | 1382699  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.257 | 240  | 1163159  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 23.598 | 264  | 999979   | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.299  | 96   | 57065    | 6.798  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.699  | 84   | 834765   | 33.497 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 6.952  | 99   | 1046705  | 35.566 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.122  | 67   | 685322   | 37.426 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.311  | 132  | 815056   | 37.278 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.487  | 113  | 837524   | 36.188 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 8.934  | 128  | 408616   | 41.571 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.652  | 143  | 413296   | 44.036 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.187 | 165  | 677927   | 36.861 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.704 | 131  | 1055034  | 34.844 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.828 | 166  | 2093179  | 39.791 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.098 | 160  | 2622093  | 40.571 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.610 | 143  | 316829   | 30.333 | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.404 | 176  | 1740029  | 39.347 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.522 | 200  | 217314   | 28.192 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.257 | 188  | 2556905  | 40.965 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.498 | 212  | 2689481  | 44.954 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.451 | 264  | 1999927  | 39.356 | ng/ul | 0.00     |

|                  |       |    |       |             |        |
|------------------|-------|----|-------|-------------|--------|
| Target Compounds |       |    |       |             | Qvalue |
| 8) Phenol        | 6.981 | 94 | 61370 | 2.040 ng/ul | 98     |

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

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