

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP072921\  
 Data File : BP006433.D  
 Acq On : 29 Jul 2021 22:15  
 Operator : CG/JU  
 Sample : M3141-13  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BGER1

Quant Time: Jul 30 01:15:23 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  | 324618   | 20.000 | ng/ul | 0.00     |
| 20) Naphthalene-d8            | 10.563 | 136  | 1430205  | 20.000 | ng/ul | 0.00     |
| 38) Acenaphthene-d10          | 14.404 | 164  | 826110   | 20.000 | ng/ul | 0.00     |
| 64) Phenanthrene-d10          | 17.157 | 188  | 1594978  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12              | 21.257 | 240  | 1432273  | 20.000 | ng/ul | 0.00     |
| 88) Perylene-d12              | 23.598 | 264  | 1270143  | 20.000 | ng/ul | 0.00     |
| System Monitoring Compounds   |        |      |          |        |       |          |
| 3) 1,4-Dioxane-d8             | 3.299  | 96   | 34024    | 3.912  | ng/uL | 0.00     |
| 4) Pyridine-d5                | 3.699  | 84   | 498050   | 19.292 | ng/ul | 0.00     |
| 7) Phenol-d5                  | 6.952  | 99   | 572660   | 18.783 | ng/ul | 0.00     |
| 9) Bis-(2-Chloroethyl)eth...  | 7.122  | 67   | 424912   | 22.399 | ng/ul | 0.00     |
| 11) 2-Chlorophenol-d4         | 7.310  | 132  | 441345   | 19.485 | ng/ul | 0.00     |
| 15) 4-Methylphenol-d8         | 8.481  | 113  | 460253   | 19.197 | ng/ul | 0.00     |
| 21) Nitrobenzene-d5           | 8.934  | 128  | 220159   | 20.674 | ng/ul | 0.00     |
| 24) 2-Nitrophenol-d4          | 9.651  | 143  | 213599   | 21.007 | ng/ul | 0.00     |
| 28) 2,4-Dichlorophenol-d3     | 10.181 | 165  | 372115   | 18.675 | ng/ul | 0.00     |
| 31) 4-Chloroaniline-d4        | 10.704 | 131  | 602015   | 18.352 | ng/ul | 0.00     |
| 46) Dimethylphthalate-d6      | 13.828 | 166  | 2282489  | 38.820 | ng/ul | 0.00     |
| 49) Acenaphthylene-d8         | 14.098 | 160  | 1468945  | 20.335 | ng/ul | 0.00     |
| 54) 4-Nitrophenol-d4          | 14.604 | 143  | 197925   | 16.954 | ng/ul | 0.00     |
| 60) Fluorene-d10              | 15.398 | 176  | 1000658  | 20.245 | ng/ul | 0.00     |
| 65) 4,6-Dinitro-2-methylph... | 15.522 | 200  | 194304   | 21.852 | ng/ul | 0.00     |
| 73) Anthracene-d10            | 17.257 | 188  | 1530732  | 21.260 | ng/ul | 0.00     |
| 81) Pyrene-d10                | 19.498 | 212  | 1716901  | 23.306 | ng/ul | 0.00     |
| 92) Benzo(a)pyrene-d12        | 23.451 | 264  | 1396891  | 21.642 | ng/ul | 0.00     |

Target Compounds Qvalue

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

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