

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP072821\  
 Data File : BP006394.D  
 Acq On : 28 Jul 2021 17:31  
 Operator : CG/JU  
 Sample : PB137941BL  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK941

Quant Time: Jul 28 18:01:22 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SFAM-EPA-BP072421.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jul 28 15:50:13 2021  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.781  | 152  | 386965   | 20.000 | ng/ul | -0.01    |
| 20) Naphthalene-d8        | 10.563 | 136  | 1693521  | 20.000 | ng/ul | -0.02    |
| 38) Acenaphthene-d10      | 14.410 | 164  | 1014170  | 20.000 | ng/ul | -0.01    |
| 64) Phenanthrene-d10      | 17.163 | 188  | 2008606  | 20.000 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.257 | 240  | 1896263  | 20.000 | ng/ul | -0.01    |
| 88) Perylene-d12          | 23.604 | 264  | 1771508  | 20.000 | ng/ul | -0.01    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.299  | 96  | 69178   | 6.673  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.705  | 84  | 1011646 | 32.872 | ng/ul | 0.00  |
| 7) Phenol-d5                  | 6.957  | 99  | 1162731 | 31.993 | ng/ul | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.128  | 67  | 751284  | 33.223 | ng/ul | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.316  | 132 | 925053  | 34.260 | ng/ul | -0.01 |
| 15) 4-Methylphenol-d8         | 8.487  | 113 | 943331  | 33.006 | ng/ul | -0.01 |
| 21) Nitrobenzene-d5           | 8.940  | 128 | 457487  | 36.280 | ng/ul | -0.01 |
| 24) 2-Nitrophenol-d4          | 9.657  | 143 | 474616  | 39.419 | ng/ul | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.187 | 165 | 794512  | 33.674 | ng/ul | -0.01 |
| 31) 4-Chloroaniline-d4        | 10.704 | 131 | 1280929 | 32.977 | ng/ul | -0.01 |
| 46) Dimethylphthalate-d6      | 13.834 | 166 | 2635533 | 36.513 | ng/ul | 0.00  |
| 49) Acenaphthylene-d8         | 14.104 | 160 | 3145615 | 35.471 | ng/ul | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.610 | 143 | 447471  | 31.222 | ng/ul | 0.00  |
| 60) Fluorene-d10              | 15.404 | 176 | 2134615 | 35.179 | ng/ul | -0.01 |
| 65) 4,6-Dinitro-2-methylph... | 15.522 | 200 | 323532  | 28.893 | ng/ul | -0.01 |
| 73) Anthracene-d10            | 17.263 | 188 | 3399585 | 37.493 | ng/ul | 0.00  |
| 81) Pyrene-d10                | 19.504 | 212 | 3736057 | 38.305 | ng/ul | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.451 | 264 | 3310028 | 36.769 | ng/ul | -0.01 |

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

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

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