

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP030322\  
 Data File : BP009256.D  
 Acq On : 03 Mar 2022 12:29  
 Operator : CG/JU  
 Sample : PB142786BL  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SBLK786

Quant Time: Mar 03 13:30:53 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP030222.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Mar 02 18:15:01 2022  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.828  | 152  | 364186   | 20.000 | ng/u1 | # 0.00   |
| 20) Naphthalene-d8        | 10.622 | 136  | 1449741  | 20.000 | ng/u1 | # 0.00   |
| 38) Acenaphthene-d10      | 14.469 | 164  | 882782   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.228 | 188  | 1725360  | 20.000 | ng/u1 | # 0.00   |
| 79) Chrysene-d12          | 21.333 | 240  | 1619734  | 20.000 | ng/u1 | # 0.00   |
| 88) Perylene-d12          | 23.745 | 264  | 1675188  | 20.000 | ng/u1 | -0.01    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.317  | 96  | 56901   | 6.280  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.723  | 84  | 728805  | 32.236 | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 6.993  | 99  | 928843  | 31.484 | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.164  | 67  | 560034  | 31.631 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.358  | 132 | 739813  | 32.075 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.528  | 113 | 707481  | 31.372 | ng/u1 | 0.00  |
| 21) Nitrobenzene-d5           | 8.981  | 128 | 305322  | 36.113 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.705  | 143 | 297166  | 37.084 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.240 | 165 | 663231  | 31.063 | ng/u1 | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.752 | 131 | 966113  | 31.999 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 13.881 | 166 | 1986682 | 32.112 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 14.157 | 160 | 2511888 | 30.728 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.651 | 143 | 242319  | 24.732 | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.463 | 176 | 1704596 | 31.665 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.575 | 200 | 162708  | 23.234 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.328 | 188 | 2574373 | 31.748 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.569 | 212 | 2894014 | 32.504 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.586 | 264 | 2741523 | 32.357 | ng/u1 | -0.01 |

Target Compounds Qvalue

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

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