

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP103122\  
 Data File : BP012422.D  
 Acq On : 01 Nov 2022 12:07  
 Operator : CG/JU  
 Sample : N5270-06  
 Misc :  
 ALS Vial : 43 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BHA81

Quant Time: Nov 01 22:56:54 2022  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP102522.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Nov 01 01:03:58 2022  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.793  | 152  | 336973   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.593 | 136  | 1523786  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.440 | 164  | 995264   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.204 | 188  | 2182839  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.298 | 240  | 2343807  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 23.686 | 264  | 2313445  | 20.000 | ng/u1 | -0.01    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.234  | 96  | 41502   | 4.959  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.664  | 84  | 249733  | 10.388 | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 6.964  | 99  | 239061  | 7.910  | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.128  | 67  | 657338  | 36.432 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.322  | 132 | 652614  | 29.568 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.510  | 113 | 460299  | 19.256 | ng/u1 | 0.00  |
| 21) Nitrobenzene-d5           | 8.963  | 128 | 444776  | 45.011 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.681  | 143 | 463168  | 45.160 | ng/u1 | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.216 | 165 | 773284  | 34.790 | ng/u1 | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.740 | 131 | 1062677 | 32.110 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 13.857 | 166 | 2749319 | 40.353 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 14.134 | 160 | 3128582 | 40.183 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.669 | 143 | 84061   | 6.519  | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.440 | 176 | 2407128 | 41.266 | ng/u1 | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.575 | 200 | 420873  | 37.033 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.304 | 188 | 3880651 | 41.260 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.545 | 212 | 4645829 | 39.480 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 23.533 | 264 | 4692715 | 41.514 | ng/u1 | -0.01 |

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

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

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