

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP103122\  
 Data File : BP012420.D  
 Acq On : 01 Nov 2022 10:58  
 Operator : CG/JU  
 Sample : N5270-13  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BHA69

Quant Time: Nov 01 22:56:34 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  | 368714   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.593 | 136  | 1649035  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.440 | 164  | 1071534  | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.204 | 188  | 2342158  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.298 | 240  | 2535647  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 23.686 | 264  | 2516083  | 20.000 | ng/u1 | -0.01    |

|                               |        |     |         |        |       |      |
|-------------------------------|--------|-----|---------|--------|-------|------|
| System Monitoring Compounds   |        |     |         |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.234  | 96  | 39225   | 4.284  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.664  | 84  | 229302  | 8.717  | ng/u1 | 0.00 |
| 7) Phenol-d5                  | 6.969  | 99  | 233263  | 7.053  | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.128  | 67  | 641218  | 32.480 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.322  | 132 | 631132  | 26.133 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 8.505  | 113 | 460038  | 17.588 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5           | 8.963  | 128 | 438820  | 41.036 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.681  | 143 | 462102  | 41.633 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.216 | 165 | 777165  | 32.309 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4        | 10.740 | 131 | 1078208 | 30.104 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6      | 13.857 | 166 | 3039822 | 41.441 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.134 | 160 | 3288093 | 39.225 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.669 | 143 | 106760  | 7.690  | ng/u1 | 0.00 |
| 60) Fluorene-d10              | 15.440 | 176 | 2641277 | 42.057 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.575 | 200 | 504771  | 41.394 | ng/u1 | 0.00 |
| 73) Anthracene-d10            | 17.304 | 188 | 4474806 | 44.341 | ng/u1 | 0.00 |
| 81) Pyrene-d10                | 19.545 | 212 | 5415864 | 42.542 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12        | 23.539 | 264 | 5624966 | 45.754 | ng/u1 | 0.00 |

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

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

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