

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP111423\  
 Data File : BP018159.D  
 Acq On : 14 Nov 2023 17:17  
 Operator : MA/JU  
 Sample : 05337-05  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 GCDJ0

Quant Time: Nov 15 00:39:21 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP111023.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Nov 13 15:50:18 2023  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units   | Dev(Min) |
|---------------------------|--------|------|----------|--------|---------|----------|
| Internal Standards        |        |      |          |        |         |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.787  | 152  | 98033    | 20.000 | ng/ul   | 0.00     |
| 20) Naphthalene-d8        | 10.604 | 136  | 398125   | 20.000 | ng/ul # | 0.00     |
| 38) Acenaphthene-d10      | 14.469 | 164  | 214549   | 20.000 | ng/ul   | 0.00     |
| 64) Phenanthrene-d10      | 17.251 | 188  | 488985   | 20.000 | ng/ul   | 0.00     |
| 79) Chrysene-d12          | 21.380 | 240  | 489965   | 20.000 | ng/ul   | -0.01    |
| 88) Perylene-d12          | 23.851 | 264  | 545637   | 20.000 | ng/ul   | -0.02    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.164  | 96  | 9052    | 3.257  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.622  | 84  | 20247   | 2.794  | ng/ul | 0.02  |
| 7) Phenol-d5                  | 6.963  | 99  | 41820   | 4.419  | ng/ul | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.140  | 67  | 156161  | 28.077 | ng/ul | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.311  | 132 | 140266  | 18.715 | ng/ul | 0.00  |
| 15) 4-Methylphenol-d8         | 8.516  | 113 | 86086   | 11.100 | ng/ul | 0.00  |
| 21) Nitrobenzene-d5           | 8.993  | 128 | 89018   | 24.818 | ng/ul | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.704  | 143 | 95634   | 23.534 | ng/ul | 0.00  |
| 28) 2,4-Dichlorophenol-d3     | 10.234 | 165 | 149261  | 20.609 | ng/ul | 0.00  |
| 31) 4-Chloroaniline-d4        | 10.793 | 131 | 141851  | 14.282 | ng/ul | 0.00  |
| 46) Dimethylphthalate-d6      | 13.893 | 166 | 579370  | 29.267 | ng/ul | 0.00  |
| 49) Acenaphthylene-d8         | 14.163 | 160 | 570570  | 26.463 | ng/ul | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.775 | 143 | 6612    | 2.235  | ng/ul | 0.04  |
| 60) Fluorene-d10              | 15.469 | 176 | 469577  | 28.731 | ng/ul | 0.00  |
| 65) 4,6-Dinitro-2-methylph... | 15.628 | 200 | 73597   | 20.820 | ng/ul | -0.04 |
| 73) Anthracene-d10            | 17.351 | 188 | 808697  | 30.579 | ng/ul | 0.00  |
| 81) Pyrene-d10                | 19.604 | 212 | 1040288 | 31.786 | ng/ul | -0.01 |
| 92) Benzo(a)pyrene-d12        | 23.686 | 264 | 1036884 | 32.681 | ng/ul | -0.02 |

Target Compounds Qvalue

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

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