

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP052323\  
 Data File : BP015262.D  
 Acq On : 24 May 2023 03:42  
 Operator : CG/JU  
 Sample : 02861-19  
 Misc :  
 ALS Vial : 33 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 HOFM0

Quant Time: May 24 04:19:57 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP051123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 23 01:25:41 2023  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 8.116  | 152  | 225225   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.945 | 136  | 1016344  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.757 | 164  | 654405   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.510 | 188  | 1432089  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.580 | 240  | 1041414  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 24.133 | 264  | 1055505  | 20.000 | ng/u1 | 0.00     |

|                               |        |     |         |        |       |      |
|-------------------------------|--------|-----|---------|--------|-------|------|
| System Monitoring Compounds   |        |     |         |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.428  | 96  | 25075   | 4.326  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.869  | 84  | 118870  | 7.234  | ng/u1 | 0.00 |
| 7) Phenol-d5                  | 7.263  | 99  | 141164  | 6.751  | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.434  | 67  | 392046  | 32.461 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.640  | 132 | 401997  | 26.145 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 8.828  | 113 | 275711  | 16.546 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5           | 9.293  | 128 | 266188  | 33.330 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 10.022 | 143 | 295385  | 33.240 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.557 | 165 | 499633  | 31.159 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4        | 11.081 | 131 | 468016  | 19.727 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6      | 14.157 | 166 | 1870423 | 37.824 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.451 | 160 | 1961868 | 34.688 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.951 | 143 | 47267   | 4.922  | ng/u1 | 0.00 |
| 60) Fluorene-d10              | 15.745 | 176 | 1527253 | 36.601 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.863 | 200 | 266097  | 29.484 | ng/u1 | 0.00 |
| 73) Anthracene-d10            | 17.610 | 188 | 2431793 | 37.083 | ng/u1 | 0.00 |
| 81) Pyrene-d10                | 19.827 | 212 | 2692338 | 44.575 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12        | 23.968 | 264 | 2111811 | 40.305 | ng/u1 | 0.00 |

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

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

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