

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP112723\  
 Data File : BP018405.D  
 Acq On : 27 Nov 2023 17:47  
 Operator : MA/JU  
 Sample : 05341-14  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0QF3

Quant Time: Nov 27 22:24:46 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP112223.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Nov 27 21:50:01 2023  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.863  | 152  | 448184   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.663 | 136  | 1704364  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.492 | 164  | 912787   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.245 | 188  | 1619301  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.374 | 240  | 1537918  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 23.880 | 264  | 1814184  | 20.000 | ng/u1 | 0.01     |

|                               |        |     |         |        |       |      |
|-------------------------------|--------|-----|---------|--------|-------|------|
| System Monitoring Compounds   |        |     |         |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.269  | 96  | 42050   | 3.514  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.687  | 84  | 246208  | 7.487  | ng/u1 | 0.00 |
| 7) Phenol-d5                  | 7.022  | 99  | 190740  | 4.657  | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.193  | 67  | 591878  | 24.165 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.387  | 132 | 628042  | 18.262 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 8.569  | 113 | 360355  | 10.902 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5           | 9.022  | 128 | 386584  | 27.925 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.745  | 143 | 344400  | 26.938 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.281 | 165 | 658276  | 21.342 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4        | 10.798 | 131 | 853412  | 20.514 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6      | 13.910 | 166 | 2079552 | 25.604 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.186 | 160 | 2379641 | 25.713 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.686 | 143 | 28541   | 2.545  | ng/u1 | 0.00 |
| 60) Fluorene-d10              | 15.486 | 176 | 1722283 | 25.352 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.598 | 200 | 134923  | 19.573 | ng/u1 | 0.00 |
| 73) Anthracene-d10            | 17.345 | 188 | 2464408 | 28.474 | ng/u1 | 0.00 |
| 81) Pyrene-d10                | 19.592 | 212 | 2781821 | 22.727 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12        | 23.715 | 264 | 3019759 | 28.711 | ng/u1 | 0.02 |

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

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

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