

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP042024\  
 Data File : BP020071.D  
 Acq On : 21 Apr 2024 12:01  
 Operator : MA/JU  
 Sample : P2104-11  
 Misc :  
 ALS Vial : 37 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0AB1

Quant Time: Apr 22 01:44:40 2024  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP041924.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Apr 20 01:32:50 2024  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.640  | 152  | 145330   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.416 | 136  | 608003   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.322 | 164  | 377822   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.175 | 188  | 703288   | 20.000 | ng/u1 | 0.02     |
| 79) Chrysene-d12          | 21.674 | 240  | 617393   | 20.000 | ng/u1 | 0.02     |
| 88) Perylene-d12          | 25.086 | 264  | 674765   | 20.000 | ng/u1 | 0.04     |

|                               |        |     |         |        |       |      |
|-------------------------------|--------|-----|---------|--------|-------|------|
| System Monitoring Compounds   |        |     |         |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.234  | 96  | 13663   | 3.852  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.628  | 84  | 210103  | 20.125 | ng/u1 | 0.00 |
| 7) Phenol-d5                  | 6.828  | 99  | 335207  | 26.465 | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 6.993  | 67  | 218372  | 24.666 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.181  | 132 | 232887  | 26.929 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 8.346  | 113 | 266265  | 26.912 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5           | 8.799  | 128 | 123886  | 28.327 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.510  | 143 | 139030  | 29.409 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.046 | 165 | 256965  | 27.399 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4        | 10.569 | 131 | 334215  | 26.695 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6      | 13.734 | 166 | 753111  | 29.166 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.004 | 160 | 831892  | 27.031 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.581 | 143 | 84884   | 21.595 | ng/u1 | 0.02 |
| 60) Fluorene-d10              | 15.345 | 176 | 594883  | 25.881 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.498 | 200 | 38838   | 9.453  | ng/u1 | 0.01 |
| 73) Anthracene-d10            | 17.280 | 188 | 852985  | 27.072 | ng/u1 | 0.02 |
| 81) Pyrene-d10                | 19.674 | 212 | 1013850 | 30.557 | ng/u1 | 0.01 |
| 92) Benzo(a)pyrene-d12        | 24.839 | 264 | 921849  | 27.967 | ng/u1 | 0.03 |

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

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

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