

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP121123\  
 Data File : BP018735.D  
 Acq On : 11 Dec 2023 19:59  
 Operator : MA/JU  
 Sample : 04929-05  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 C0AQ3

Quant Time: Dec 11 22:02:04 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Dec 11 22:02:13 2023  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.840  | 152  | 294552   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.634 | 136  | 826064   | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.469 | 164  | 521277   | 20.000 | ng/u1 | 0.00     |
| 64) Phenanthrene-d10      | 17.222 | 188  | 1229174  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.304 | 240  | 1370092  | 20.000 | ng/u1 | 0.00     |
| 88) Perylene-d12          | 23.674 | 264  | 1635821  | 20.000 | ng/u1 | 0.00     |

|                               |        |     |         |        |       |      |
|-------------------------------|--------|-----|---------|--------|-------|------|
| System Monitoring Compounds   |        |     |         |        |       |      |
| 3) 1,4-Dioxane-d8             | 3.264  | 96  | 37345   | 4.320  | ng/uL | 0.00 |
| 4) Pyridine-d5                | 3.681  | 84  | 483559  | 20.886 | ng/u1 | 0.00 |
| 7) Phenol-d5                  | 7.010  | 99  | 633712  | 21.439 | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth...  | 7.175  | 67  | 389520  | 22.234 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4         | 7.369  | 132 | 501083  | 21.725 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8         | 8.552  | 113 | 490167  | 20.490 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5           | 9.004  | 128 | 231063  | 31.428 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4          | 9.722  | 143 | 225616  | 29.598 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3     | 10.257 | 165 | 357027  | 22.906 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4        | 10.775 | 131 | 468571  | 23.003 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6      | 13.887 | 166 | 1141723 | 24.540 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8         | 14.163 | 160 | 1241206 | 24.228 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4          | 14.669 | 143 | 171851  | 22.503 | ng/u1 | 0.00 |
| 60) Fluorene-d10              | 15.463 | 176 | 985991  | 25.248 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.581 | 200 | 145966  | 19.805 | ng/u1 | 0.00 |
| 73) Anthracene-d10            | 17.316 | 188 | 1594036 | 24.668 | ng/u1 | 0.00 |
| 81) Pyrene-d10                | 19.557 | 212 | 2062822 | 19.342 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12        | 23.521 | 264 | 2311324 | 24.233 | ng/u1 | 0.00 |

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

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

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