

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP020325\  
 Data File : BP023878.D  
 Acq On : 03 Feb 2025 13:53  
 Operator : RC/JU  
 Sample : Q1247-10  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A6337

Quant Time: Feb 03 14:46:37 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP012925.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Feb 03 11:19:36 2025  
 Response via : Initial Calibration

| Compound                  | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|---------------------------|--------|------|----------|--------|-------|----------|
| -----                     |        |      |          |        |       |          |
| Internal Standards        |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4 | 7.775  | 152  | 335073   | 20.000 | ng/u1 | 0.00     |
| 20) Naphthalene-d8        | 10.552 | 136  | 1435747  | 20.000 | ng/u1 | 0.00     |
| 38) Acenaphthene-d10      | 14.398 | 164  | 901785   | 20.000 | ng/u1 | 0.01     |
| 64) Phenanthrene-d10      | 17.186 | 188  | 1845689  | 20.000 | ng/u1 | 0.00     |
| 79) Chrysene-d12          | 21.627 | 240  | 1819362  | 20.000 | ng/u1 | -0.01    |
| 88) Perylene-d12          | 24.992 | 264  | 1918377  | 20.000 | ng/u1 | -0.02    |

|                               |        |     |         |        |       |       |
|-------------------------------|--------|-----|---------|--------|-------|-------|
| System Monitoring Compounds   |        |     |         |        |       |       |
| 3) 1,4-Dioxane-d8             | 3.299  | 96  | 23911   | 2.965  | ng/uL | 0.00  |
| 4) Pyridine-d5                | 3.711  | 84  | 186761  | 8.026  | ng/u1 | 0.00  |
| 7) Phenol-d5                  | 6.958  | 99  | 278616  | 9.396  | ng/u1 | 0.00  |
| 9) Bis-(2-Chloroethyl)eth...  | 7.111  | 67  | 351050  | 22.366 | ng/u1 | 0.00  |
| 11) 2-Chlorophenol-d4         | 7.316  | 132 | 528526  | 22.697 | ng/u1 | 0.00  |
| 15) 4-Methylphenol-d8         | 8.481  | 113 | 456389  | 18.841 | ng/u1 | 0.00  |
| 21) Nitrobenzene-d5           | 8.922  | 128 | 235607  | 20.511 | ng/u1 | 0.00  |
| 24) 2-Nitrophenol-d4          | 9.640  | 143 | 234993  | 18.959 | ng/u1 | 0.01  |
| 28) 2,4-Dichlorophenol-d3     | 10.181 | 165 | 492723  | 21.670 | ng/u1 | 0.01  |
| 31) 4-Chloroaniline-d4        | 10.681 | 131 | 903152  | 26.560 | ng/u1 | 0.00  |
| 46) Dimethylphthalate-d6      | 13.798 | 166 | 1558825 | 23.175 | ng/u1 | 0.00  |
| 49) Acenaphthylene-d8         | 14.087 | 160 | 1702932 | 22.447 | ng/u1 | 0.00  |
| 54) 4-Nitrophenol-d4          | 14.610 | 143 | 110261  | 8.254  | ng/u1 | 0.00  |
| 60) Fluorene-d10              | 15.404 | 176 | 1380045 | 24.363 | ng/u1 | 0.01  |
| 65) 4,6-Dinitro-2-methylph... | 15.516 | 200 | 159666  | 14.022 | ng/u1 | 0.00  |
| 73) Anthracene-d10            | 17.286 | 188 | 2220721 | 24.505 | ng/u1 | 0.00  |
| 81) Pyrene-d10                | 19.657 | 212 | 2579585 | 26.461 | ng/u1 | 0.00  |
| 92) Benzo(a)pyrene-d12        | 24.757 | 264 | 2271236 | 22.691 | ng/u1 | -0.02 |

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

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

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