

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013125\
 Data File : BP023822.D
 Acq On : 31 Jan 2025 12:38
 Operator : RC/JU
 Sample : Q1197-14
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E2992

Quant Time: Jan 31 13:34:45 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 31 11:40:55 2025
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 7.787 | 152 | 441787 | 20.000 | ng/u1 | 0.00 |
| 20) Naphthalene-d8 | 10.557 | 136 | 1851957 | 20.000 | ng/u1 | 0.00 |
| 38) Acenaphthene-d10 | 14.398 | 164 | 1211392 | 20.000 | ng/u1 | 0.00 |
| 64) Phenanthrene-d10 | 17.192 | 188 | 2471098 | 20.000 | ng/u1 | 0.00 |
| 79) Chrysene-d12 | 21.639 | 240 | 2449487 | 20.000 | ng/u1 | 0.00 |
| 88) Perylene-d12 | 25.016 | 264 | 2629777 | 20.000 | ng/u1 | 0.00 |
| System Monitoring Compounds | | | | | | |
| 3) 1,4-Dioxane-d8 | 3.299 | 96 | 61990 | 5.830 | ng/uL | 0.00 |
| 4) Pyridine-d5 | 3.711 | 84 | 370872 | 12.088 | ng/u1 | 0.00 |
| 7) Phenol-d5 | 6.975 | 99 | 462235 | 11.823 | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth... | 7.116 | 67 | 709973 | 34.307 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4 | 7.328 | 132 | 1001370 | 32.616 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8 | 8.493 | 113 | 794883 | 24.888 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5 | 8.928 | 128 | 529991 | 35.770 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4 | 9.646 | 143 | 580133 | 36.286 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3 | 10.193 | 165 | 1009940 | 34.434 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4 | 10.693 | 131 | 1396083 | 31.829 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6 | 13.804 | 166 | 3531746 | 39.086 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8 | 14.093 | 160 | 3769180 | 36.985 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4 | 14.628 | 143 | 206954 | 11.533 | ng/u1 | 0.00 |
| 60) Fluorene-d10 | 15.404 | 176 | 3111185 | 40.887 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.522 | 200 | 523292 | 34.326 | ng/u1 | 0.00 |
| 73) Anthracene-d10 | 17.292 | 188 | 5198989 | 42.850 | ng/u1 | -0.01 |
| 81) Pyrene-d10 | 19.663 | 212 | 6004647 | 45.750 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12 | 24.792 | 264 | 6145349 | 44.788 | ng/u1 | 0.00 |

Target Compounds Qvalue

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

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