

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP113022\
 Data File : BP012891.D
 Acq On : 30 Nov 2022 22:22
 Operator : CG/JU
 Sample : N5759-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHB79

Quant Time: Nov 30 22:53:08 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 29 22:53:34 2022
 Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|-------------------------------|--------|------|----------|--------|-------|----------|
| Internal Standards | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 8.010 | 152 | 584120 | 20.000 | ng/u1 | 0.00 |
| 20) Naphthalene-d8 | 10.822 | 136 | 2691416 | 20.000 | ng/u1 | 0.00 |
| 38) Acenaphthene-d10 | 14.645 | 164 | 1845506 | 20.000 | ng/u1 | 0.00 |
| 64) Phenanthrene-d10 | 17.398 | 188 | 3826189 | 20.000 | ng/u1 | 0.00 |
| 79) Chrysene-d12 | 21.486 | 240 | 2574110 | 20.000 | ng/u1 | 0.00 |
| 88) Perylene-d12 | 23.992 | 264 | 2119167 | 20.000 | ng/u1 | 0.00 |
| System Monitoring Compounds | | | | | | |
| 3) 1,4-Dioxane-d8 | 3.393 | 96 | 60112 | 4.410 | ng/uL | 0.00 |
| 4) Pyridine-d5 | 3.822 | 84 | 167292 | 4.101 | ng/u1 | 0.00 |
| 7) Phenol-d5 | 7.157 | 99 | 335944 | 6.999 | ng/u1 | 0.00 |
| 9) Bis-(2-Chloroethyl)eth... | 7.334 | 67 | 983995 | 33.269 | ng/u1 | 0.00 |
| 11) 2-Chlorophenol-d4 | 7.534 | 132 | 932447 | 25.495 | ng/u1 | 0.00 |
| 15) 4-Methylphenol-d8 | 8.710 | 113 | 648075 | 16.758 | ng/u1 | 0.00 |
| 21) Nitrobenzene-d5 | 9.175 | 128 | 614493 | 37.303 | ng/u1 | 0.00 |
| 24) 2-Nitrophenol-d4 | 9.899 | 143 | 633764 | 36.233 | ng/u1 | 0.00 |
| 28) 2,4-Dichlorophenol-d3 | 10.440 | 165 | 1161685 | 28.464 | ng/u1 | 0.00 |
| 31) 4-Chloroaniline-d4 | 10.957 | 131 | 600544 | 10.746 | ng/u1 | 0.00 |
| 46) Dimethylphthalate-d6 | 14.051 | 166 | 4524716 | 35.619 | ng/u1 | 0.00 |
| 49) Acenaphthylene-d8 | 14.339 | 160 | 4977855 | 33.585 | ng/u1 | 0.00 |
| 54) 4-Nitrophenol-d4 | 14.822 | 143 | 86262 | 3.809 | ng/u1 | 0.00 |
| 60) Fluorene-d10 | 15.639 | 176 | 4030400 | 36.229 | ng/u1 | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.745 | 200 | 536584 | 29.342 | ng/u1 | 0.00 |
| 73) Anthracene-d10 | 17.498 | 188 | 6392832 | 37.956 | ng/u1 | 0.00 |
| 81) Pyrene-d10 | 19.727 | 212 | 6580529 | 43.996 | ng/u1 | 0.00 |
| 92) Benzo(a)pyrene-d12 | 23.833 | 264 | 4002754 | 38.454 | ng/u1 | 0.00 |

Target Compounds Qvalue

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

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