

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092923\
 Data File : BP017447.D
 Acq On : 29 Sep 2023 15:37
 Operator : MA/JU
 Sample : 04497-10
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHQ2

Quant Time: Sep 29 22:35:27 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 29 22:31:03 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	209260	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.793	136	840562	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.645	164	527773	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.433	188	1227114	20.000	ng/u1	-0.01
79) Chrysene-d12	21.563	240	1236413	20.000	ng/u1	-0.01
88) Perylene-d12	24.145	264	1427928	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	10935	2.146	ng/uL	0.00
4) Pyridine-d5	3.734	84	88803	6.235	ng/u1	0.00
7) Phenol-d5	7.111	99	111865	6.509	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	311193	30.003	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.475	132	336577	24.692	ng/u1	0.00
15) 4-Methylphenol-d8	8.675	113	207934	15.562	ng/u1	-0.01
21) Nitrobenzene-d5	9.163	128	213140	35.075	ng/u1	0.00
24) 2-Nitrophenol-d4	9.881	143	229738	34.818	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.404	165	382716	30.701	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.963	131	377715	20.917	ng/u1	0.00
46) Dimethylphthalate-d6	14.063	166	1317367	34.844	ng/u1	0.00
49) Acenaphthylene-d8	14.345	160	1519226	34.938	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.887	143	37473	5.689	ng/u1	0.00
60) Fluorene-d10	15.651	176	1180455	36.267	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	187889	27.166	ng/u1	-0.01
73) Anthracene-d10	17.539	188	1931850	35.497	ng/u1	0.00
81) Pyrene-d10	19.786	212	2412401	36.080	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.974	264	2345682	34.119	ng/u1	-0.01

Target Compounds	Qvalue

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

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