

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092923\
 Data File : BP017443.D
 Acq On : 29 Sep 2023 13:22
 Operator : MA/JU
 Sample : 04497-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBBP1

Quant Time: Sep 29 22:51:18 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	7.957	152	182535	20.000	ng/ul	0.00
20) Naphthalene-d8	10.793	136	743649	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.645	164	464234	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.433	188	1061038	20.000	ng/ul	-0.01
79) Chrysene-d12	21.569	240	981774	20.000	ng/ul	0.00
88) Perylene-d12	24.145	264	1107284	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	18387	4.137	ng/uL	0.00
4) Pyridine-d5	3.740	84	129446	10.419	ng/ul	0.00
7) Phenol-d5	7.110	99	92198	6.150	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	254681	28.150	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.475	132	277388	23.329	ng/ul	0.00
15) 4-Methylphenol-d8	8.681	113	167298	14.354	ng/ul	0.00
21) Nitrobenzene-d5	9.163	128	177560	33.027	ng/ul	0.00
24) 2-Nitrophenol-d4	9.881	143	188119	32.226	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.416	165	330651	29.981	ng/ul	0.00
31) 4-Chloroaniline-d4	10.963	131	363236	22.736	ng/ul	0.00
46) Dimethylphthalate-d6	14.063	166	1081083	32.508	ng/ul	0.00
49) Acenaphthylene-d8	14.345	160	1253441	32.771	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.892	143	26465	4.568	ng/ul	0.00
60) Fluorene-d10	15.651	176	956772	33.418	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	118485	19.813	ng/ul	-0.01
73) Anthracene-d10	17.539	188	1540984	32.747	ng/ul	0.00
81) Pyrene-d10	19.786	212	1836113	34.584	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.974	264	1684140	31.590	ng/ul	-0.01

Target Compounds Qvalue

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

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