

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017367.D
 Acq On : 26 Sep 2023 18:10
 Operator : MA/JU
 Sample : 04450-11
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJD3

Quant Time: Sep 27 00:29:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 26 04:52:29 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	125136	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	534452	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.586	164	331090	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.357	188	744846	20.000	ng/u1	0.00
79) Chrysene-d12	21.457	240	689224	20.000	ng/u1	0.00
88) Perylene-d12	23.974	264	761678	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	18493	6.070	ng/uL	0.00
4) Pyridine-d5	3.734	84	137860	16.186	ng/u1	0.00
7) Phenol-d5	7.081	99	98112	9.546	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	305010	49.177	ng/u1	0.00
11) 2-Chlorophenol-d4	7.446	132	299347	36.724	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	185868	23.262	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	193868	50.176	ng/u1	0.00
24) 2-Nitrophenol-d4	9.834	143	196315	46.794	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	336750	42.486	ng/u1	0.00
31) 4-Chloroaniline-d4	10.910	131	409713	35.684	ng/u1	0.00
46) Dimethylphthalate-d6	13.998	166	1213211	51.151	ng/u1	0.00
49) Acenaphthylene-d8	14.281	160	1335475	48.956	ng/u1	0.00
54) 4-Nitrophenol-d4	14.822	143	21189	5.128	ng/u1	0.00
60) Fluorene-d10	15.581	176	1050063	51.425	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	134967	32.150	ng/u1	0.00
73) Anthracene-d10	17.457	188	1735854	52.547	ng/u1	0.00
81) Pyrene-d10	19.698	212	2094107	56.185	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.810	264	1989144	54.241	ng/u1	0.00

Target Compounds Qvalue

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

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