

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017385.D
 Acq On : 27 Sep 2023 04:55
 Operator : MA/JU
 Sample : 04566-10
 Misc :
 ALS Vial : 78 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 B102-08-2

Quant Time: Sep 27 05:34:45 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	146991	20.000	ng/u1	0.00
20) Naphthalene-d8	10.740	136	659223	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.581	164	416514	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.351	188	979432	20.000	ng/u1	0.00
79) Chrysene-d12	21.457	240	957160	20.000	ng/u1	0.00
88) Perylene-d12	23.968	264	1067091	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	30338	8.478	ng/uL	0.00
4) Pyridine-d5	3.728	84	282867	28.274	ng/u1	0.00
7) Phenol-d5	7.075	99	231738	19.196	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	427147	58.629	ng/u1	0.00
11) 2-Chlorophenol-d4	7.440	132	444466	46.420	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	318622	33.948	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	279481	58.643	ng/u1	0.00
24) 2-Nitrophenol-d4	9.834	143	291744	56.378	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	520086	53.197	ng/u1	0.00
31) 4-Chloroaniline-d4	10.910	131	631625	44.599	ng/u1	0.00
46) Dimethylphthalate-d6	13.998	166	1862827	62.432	ng/u1	0.00
49) Acenaphthylene-d8	14.281	160	2049412	59.720	ng/u1	0.00
54) 4-Nitrophenol-d4	14.816	143	73048	14.052	ng/u1	0.00
60) Fluorene-d10	15.581	176	1614239	62.841	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	254308	46.068	ng/u1	0.00
73) Anthracene-d10	17.457	188	2661589	61.273	ng/u1	0.00
81) Pyrene-d10	19.698	212	3400918	65.705	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.810	264	3396318	66.105	ng/u1	0.00

Target Compounds	Qvalue

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

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