

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092723\
 Data File : BP017393.D
 Acq On : 27 Sep 2023 12:36
 Operator : MA/JU
 Sample : PB155590TB 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLEB590

Quant Time: Sep 27 22:29:29 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	36303	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	166904	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.581	164	100601	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.351	188	211762	20.000	ng/u1	0.00
79) Chrysene-d12	21.457	240	177303	20.000	ng/u1	0.00
88) Perylene-d12	23.968	264	188736	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	25255	28.574	ng/uL	0.00
4) Pyridine-d5	3.722	84	369886	149.699	ng/u1	0.00
7) Phenol-d5	7.075	99	447352	150.040	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	282063	156.758	ng/u1	0.00
11) 2-Chlorophenol-d4	7.440	132	365261	154.460	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	357326	154.151	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	174707	144.791	ng/u1	0.00
24) 2-Nitrophenol-d4	9.834	143	188008	143.500	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	344403	139.137	ng/u1	0.00
31) 4-Chloroaniline-d4	10.910	131	507131	141.434	ng/u1	0.00
46) Dimethylphthalate-d6	13.998	166	1118799	155.243	ng/u1	0.00
49) Acenaphthylene-d8	14.281	160	1210063	145.990	ng/u1	0.00
54) 4-Nitrophenol-d4	14.816	143	115487	91.982	ng/u1	0.00
60) Fluorene-d10	15.581	176	929592	149.829	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.710	200	89933	75.350	ng/u1	0.00
73) Anthracene-d10	17.451	188	1432170	152.492	ng/u1	0.00
81) Pyrene-d10	19.698	212	1626111	169.597	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.804	264	1463272	161.028	ng/u1	0.00

Target Compounds Qvalue

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

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