

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092723\
 Data File : BP017398.D
 Acq On : 27 Sep 2023 15:26
 Operator : MA/JU
 Sample : PB155590TB 10X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLEB590

Quant Time: Sep 27 22:30:48 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	157674	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	698258	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.581	164	438829	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.357	188	961892	20.000	ng/u1	0.00
79) Chrysene-d12	21.457	240	845319	20.000	ng/u1	0.00
88) Perylene-d12	23.974	264	908723	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	25314	6.594	ng/uL	0.00
4) Pyridine-d5	3.722	84	375944	35.031	ng/u1	0.00
7) Phenol-d5	7.075	99	450574	34.794	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	291125	37.252	ng/u1	0.00
11) 2-Chlorophenol-d4	7.440	132	365300	35.567	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	361105	35.867	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	175509	34.768	ng/u1	0.00
24) 2-Nitrophenol-d4	9.834	143	188445	34.380	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	349062	33.708	ng/u1	0.00
31) 4-Chloroaniline-d4	10.910	131	515106	34.339	ng/u1	0.00
46) Dimethylphthalate-d6	13.998	166	1127716	35.873	ng/u1	0.00
49) Acenaphthylene-d8	14.281	160	1227845	33.960	ng/u1	0.00
54) 4-Nitrophenol-d4	14.816	143	125979	23.002	ng/u1	0.00
60) Fluorene-d10	15.581	176	970657	35.866	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	109489	20.196	ng/u1	0.00
73) Anthracene-d10	17.457	188	1539180	36.080	ng/u1	0.00
81) Pyrene-d10	19.698	212	1779059	38.918	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.804	264	1631520	37.290	ng/u1	0.00

Target Compounds Qvalue

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

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