

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018118.D
 Acq On : 12 Nov 2023 23:24
 Operator : MA/JU
 Sample : PB156906TB 10X
 Misc :
 ALS Vial : 99 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SLEB906

Quant Time: Nov 13 03:20:14 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 21:55:18 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	81581	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.663	136	315429	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.516	164	180191	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.298	188	395321	20.000	ng/u1	-0.01
79) Chrysene-d12	21.427	240	390159	20.000	ng/u1	0.00
88) Perylene-d12	23.927	264	441441	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	24094	10.418	ng/uL	0.00
4) Pyridine-d5	3.658	84	319972	53.064	ng/u1	0.00
7) Phenol-d5	7.010	99	397019	50.408	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	247560	53.486	ng/u1	0.00
11) 2-Chlorophenol-d4	7.363	132	316900	50.809	ng/u1	-0.01
15) 4-Methylphenol-d8	8.569	113	317747	49.233	ng/u1	0.00
21) Nitrobenzene-d5	9.046	128	153351	53.962	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.763	143	180290	55.999	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.287	165	292065	50.898	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.840	131	414114	52.625	ng/u1	-0.01
46) Dimethylphthalate-d6	13.939	166	902908	54.308	ng/u1	-0.01
49) Acenaphthylene-d8	14.216	160	995585	54.979	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.781	143	128124	51.572	ng/u1	0.00
60) Fluorene-d10	15.522	176	748754	54.548	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	142476	49.854	ng/u1	-0.01
73) Anthracene-d10	17.398	188	1216145	56.881	ng/u1	-0.01
81) Pyrene-d10	19.651	212	1475893	56.632	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.762	264	1528253	59.538	ng/u1	-0.01

Target Compounds Qvalue

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

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