

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017587.D
 Acq On : 06 Oct 2023 04:27
 Operator : MA/JU
 Sample : PB155912BL
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK912

Quant Time: Oct 06 06:00:13 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 06 05:55:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	101274	20.000	ng/u1	0.00
20) Naphthalene-d8	10.769	136	417941	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.628	164	263559	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.416	188	613006	20.000	ng/u1	0.00
79) Chrysene-d12	21.551	240	617554	20.000	ng/u1	0.00
88) Perylene-d12	24.127	264	679043	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	20366	8.040	ng/uL	0.00
4) Pyridine-d5	3.722	84	300530	45.325	ng/u1	0.00
7) Phenol-d5	7.087	99	374717	43.976	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	231553	44.534	ng/u1	0.00
11) 2-Chlorophenol-d4	7.452	132	301337	44.914	ng/u1	0.00
15) 4-Methylphenol-d8	8.652	113	298757	43.797	ng/u1	0.00
21) Nitrobenzene-d5	9.140	128	143066	44.495	ng/u1	0.00
24) 2-Nitrophenol-d4	9.857	143	161819	44.621	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	286391	42.598	ng/u1	0.00
31) 4-Chloroaniline-d4	10.940	131	418112	43.656	ng/u1	0.00
46) Dimethylphthalate-d6	14.040	166	974250	47.389	ng/u1	0.00
49) Acenaphthylene-d8	14.322	160	1023060	45.129	ng/u1	0.00
54) 4-Nitrophenol-d4	14.869	143	149830	41.618	ng/u1	0.00
60) Fluorene-d10	15.628	176	822430	47.164	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.769	200	124123	32.479	ng/u1	0.00
73) Anthracene-d10	17.516	188	1385024	48.811	ng/u1	0.00
81) Pyrene-d10	19.769	212	1694072	49.368	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.957	264	1688228	49.264	ng/u1	0.00

Target Compounds Qvalue

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

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