

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
 Data File : BP017510.D
 Acq On : 03 Oct 2023 10:59
 Operator : MA/JU
 Sample : PB155865BL
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK865

Quant Time: Oct 03 23:25:56 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 03 22:59:42 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	132351	20.000	ng/u1	0.00
20) Naphthalene-d8	10.775	136	561208	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	359086	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.422	188	826189	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	776735	20.000	ng/u1	0.00
88) Perylene-d12	24.139	264	829659	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	22525	6.991	ng/uL	0.00
4) Pyridine-d5	3.716	84	355886	39.507	ng/u1	0.00
7) Phenol-d5	7.093	99	451774	41.562	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	262919	40.079	ng/u1	0.00
11) 2-Chlorophenol-d4	7.457	132	357935	41.518	ng/u1	0.00
15) 4-Methylphenol-d8	8.657	113	363577	43.022	ng/u1	0.00
21) Nitrobenzene-d5	9.146	128	176504	43.504	ng/u1	0.00
24) 2-Nitrophenol-d4	9.863	143	199813	45.357	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	364802	43.830	ng/u1	0.00
31) 4-Chloroaniline-d4	10.945	131	500023	41.473	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	1140518	44.337	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	1279132	43.235	ng/u1	0.00
54) 4-Nitrophenol-d4	14.875	143	160804	35.881	ng/u1	0.00
60) Fluorene-d10	15.639	176	996132	44.981	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200	98493	21.151	ng/u1	0.00
73) Anthracene-d10	17.522	188	1652330	45.094	ng/u1	0.00
81) Pyrene-d10	19.780	212	1997988	47.567	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.974	264	1833923	45.911	ng/u1	0.00

Target Compounds	Qvalue

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

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