

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018036.D
 Acq On : 10 Nov 2023 21:08
 Operator : MA/JU
 Sample : PB156717BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK717

Quant Time: Nov 10 23:10:56 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	7.852	152	40626	20.000	ng/u1	0.00
20) Naphthalene-d8	10.669	136	162699	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.528	164	107035	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.310	188	244425	20.000	ng/u1	0.00
79) Chrysene-d12	21.433	240	238619	20.000	ng/u1	0.00
88) Perylene-d12	23.939	264	279955	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	7987	6.935	ng/uL	0.00
4) Pyridine-d5	3.664	84	99367	33.092	ng/u1	0.00
7) Phenol-d5	7.016	99	126923	32.361	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	80130	34.764	ng/u1	0.00
11) 2-Chlorophenol-d4	7.375	132	108221	34.843	ng/u1	0.00
15) 4-Methylphenol-d8	8.575	113	103321	32.148	ng/u1	0.00
21) Nitrobenzene-d5	9.057	128	51145	34.891	ng/u1	0.00
24) 2-Nitrophenol-d4	9.769	143	58800	35.408	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.298	165	91628	30.958	ng/u1	0.00
31) 4-Chloroaniline-d4	10.851	131	129322	31.861	ng/u1	0.00
46) Dimethylphthalate-d6	13.945	166	367810	37.243	ng/u1	0.00
49) Acenaphthylene-d8	14.228	160	378950	35.230	ng/u1	0.00
54) 4-Nitrophenol-d4	14.792	143	38112	25.826	ng/u1	0.01
60) Fluorene-d10	15.528	176	294510	36.120	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.686	200	54061	30.595	ng/u1	0.00
73) Anthracene-d10	17.410	188	491511	37.181	ng/u1	0.00
81) Pyrene-d10	19.663	212	599918	37.639	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.774	264	619632	38.064	ng/u1	0.00

Target Compounds Qvalue

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

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