

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012458.D
 Acq On : 02 Nov 2022 19:25
 Operator : CG/JU
 Sample : PB148617BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK617

Quant Time: Nov 03 00:10:22 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 03 00:09:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	334811	20.000	ng/u1	0.00
20) Naphthalene-d8	10.587	136	1518669	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.439	164	1040374	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.198	188	2301928	20.000	ng/u1	0.00
79) Chrysene-d12	21.298	240	2146846	20.000	ng/u1	0.00
88) Perylene-d12	23.680	264	1968012	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	45770	5.504	ng/uL	0.00
4) Pyridine-d5	3.658	84	583219	24.416	ng/u1	0.00
7) Phenol-d5	6.963	99	832909	27.736	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	510298	28.465	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	657072	29.963	ng/u1	0.00
15) 4-Methylphenol-d8	8.504	113	683765	28.789	ng/u1	0.00
21) Nitrobenzene-d5	8.957	128	343281	34.857	ng/u1	0.00
24) 2-Nitrophenol-d4	9.675	143	392055	38.355	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	678964	30.650	ng/u1	0.00
31) 4-Chloroaniline-d4	10.734	131	984324	29.842	ng/u1	0.00
46) Dimethylphthalate-d6	13.857	166	2296578	32.247	ng/u1	0.00
49) Acenaphthylene-d8	14.128	160	2522362	30.992	ng/u1	0.00
54) 4-Nitrophenol-d4	14.663	143	379082	28.124	ng/u1	0.00
60) Fluorene-d10	15.434	176	1954380	32.052	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	304726	25.426	ng/u1	0.00
73) Anthracene-d10	17.298	188	3162209	31.882	ng/u1	0.00
81) Pyrene-d10	19.539	212	3635471	33.729	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.527	264	3097873	32.216	ng/u1	0.00

Target Compounds	Qvalue

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

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