

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012491.D
 Acq On : 03 Nov 2022 16:03
 Operator : CG/JU
 Sample : PB148620BL
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 03 23:04:37 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	7.781	152	241431	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	1136350	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.434	164	786727	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.198	188	1775284	20.000	ng/u1	0.00
79) Chrysene-d12	21.292	240	1605877	20.000	ng/u1	0.00
88) Perylene-d12	23.680	264	1393320	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	37454	6.246	ng/uL	0.00
4) Pyridine-d5	3.658	84	278991	16.197	ng/u1	0.00
7) Phenol-d5	6.957	99	716932	33.108	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	424493	32.838	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	549046	34.720	ng/u1	0.00
15) 4-Methylphenol-d8	8.504	113	591631	34.544	ng/u1	0.00
21) Nitrobenzene-d5	8.951	128	286352	38.859	ng/u1	0.00
24) 2-Nitrophenol-d4	9.675	143	326182	42.646	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	576271	34.766	ng/u1	0.00
31) 4-Chloroaniline-d4	10.734	131	482627	19.555	ng/u1	0.00
46) Dimethylphthalate-d6	13.851	166	1912052	35.503	ng/u1	0.00
49) Acenaphthylene-d8	14.128	160	2096067	34.057	ng/u1	0.00
54) 4-Nitrophenol-d4	14.663	143	302715	29.699	ng/u1	0.00
60) Fluorene-d10	15.434	176	1660864	36.020	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	234421	25.362	ng/u1	0.00
73) Anthracene-d10	17.298	188	2763363	36.126	ng/u1	0.00
81) Pyrene-d10	19.539	212	3135663	38.892	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.527	264	2471327	36.301	ng/u1	0.00

Target Compounds	Qvalue

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

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