

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
 Data File : BP012557.D
 Acq On : 08 Nov 2022 18:42
 Operator : CG/JU
 Sample : PB148835BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK835

Quant Time: Nov 08 22:44:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 08 22:40:23 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	147170	20.000	ng/u1	0.00
20) Naphthalene-d8	10.569	136	722695	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.428	164	514325	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.186	188	1160206	20.000	ng/u1	0.00
79) Chrysene-d12	21.286	240	984170	20.000	ng/u1	0.00
88) Perylene-d12	23.663	264	880867	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	36472	9.979	ng/uL	0.00
4) Pyridine-d5	3.658	84	146277	13.932	ng/u1	0.00
7) Phenol-d5	6.963	99	757161	57.361	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	410803	52.132	ng/u1	0.00
11) 2-Chlorophenol-d4	7.305	132	566662	58.785	ng/u1	0.00
15) 4-Methylphenol-d8	8.499	113	611278	58.552	ng/u1	0.00
21) Nitrobenzene-d5	8.940	128	291574	62.215	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	335468	68.965	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	626845	59.463	ng/u1	0.00
31) 4-Chloroaniline-d4	10.734	131	75390	4.803	ng/u1	0.01
46) Dimethylphthalate-d6	13.845	166	2018490	57.330	ng/u1	0.00
49) Acenaphthylene-d8	14.116	160	2199082	54.655	ng/u1	0.00
54) 4-Nitrophenol-d4	14.681	143	314554	47.205	ng/u1	0.00
60) Fluorene-d10	15.422	176	1774020	58.850	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	325305	53.854	ng/u1	0.00
73) Anthracene-d10	17.286	188	2923084	58.473	ng/u1	0.00
81) Pyrene-d10	19.533	212	3275591	66.292	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.510	264	2492800	57.918	ng/u1	0.00

Target Compounds Qvalue

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

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