

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011990.D
 Acq On : 07 Oct 2022 03:11
 Operator : CG/JU
 Sample : PB148079BL
 Misc :
 ALS Vial : 79 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK079

Quant Time: Oct 07 04:04:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 07 03:22:44 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	212794	20.000	ng/u1	0.00
20) Naphthalene-d8	10.681	136	908404	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.522	164	560737	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.275	188	1197081	20.000	ng/u1	0.00
79) Chrysene-d12	21.363	240	1098128	20.000	ng/u1	0.00
88) Perylene-d12	23.792	264	1100286	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	40211	8.334	ng/uL	0.00
4) Pyridine-d5	3.717	84	508958	36.012	ng/u1	0.00
7) Phenol-d5	7.040	99	701153	38.167	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	406052	39.877	ng/u1	0.00
11) 2-Chlorophenol-d4	7.405	132	574671	39.628	ng/u1	0.00
15) 4-Methylphenol-d8	8.587	113	545348	35.911	ng/u1	0.00
21) Nitrobenzene-d5	9.040	128	280199	40.496	ng/u1	0.00
24) 2-Nitrophenol-d4	9.763	143	289901	39.161	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.305	165	510835	37.307	ng/u1	0.00
31) 4-Chloroaniline-d4	10.822	131	792314	37.956	ng/u1	0.00
46) Dimethylphthalate-d6	13.934	166	1588523	39.615	ng/u1	0.00
49) Acenaphthylene-d8	14.210	160	1841559	40.286	ng/u1	0.00
54) 4-Nitrophenol-d4	14.722	143	266897	32.142	ng/u1	0.00
60) Fluorene-d10	15.516	176	1439531	41.338	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	193583	26.717	ng/u1	0.00
73) Anthracene-d10	17.375	188	2290777	42.010	ng/u1	0.00
81) Pyrene-d10	19.610	212	2559841	46.029	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.633	264	2409877	43.753	ng/u1	0.00

Target Compounds Qvalue

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

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