

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011868.D
 Acq On : 03 Oct 2022 12:50
 Operator : CG/JU
 Sample : PB147991BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK991

Quant Time: Oct 03 23:30:25 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 01 00:53:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	270284	20.000	ng/u1	0.00
20) Naphthalene-d8	10.705	136	1108432	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.540	164	659424	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.292	188	1354504	20.000	ng/u1	0.00
79) Chrysene-d12	21.375	240	1346676	20.000	ng/u1	-0.01
88) Perylene-d12	23.810	264	1428807	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	35080	5.724	ng/uL	0.00
4) Pyridine-d5	3.728	84	451718	25.164	ng/u1	0.00
7) Phenol-d5	7.058	99	635172	27.221	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	362960	28.063	ng/u1	0.00
11) 2-Chlorophenol-d4	7.422	132	533010	28.937	ng/u1	0.00
15) 4-Methylphenol-d8	8.605	113	500738	25.960	ng/u1	0.00
21) Nitrobenzene-d5	9.058	128	255065	30.211	ng/u1	0.00
24) 2-Nitrophenol-d4	9.781	143	270553	29.952	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	466671	27.931	ng/u1	0.00
31) 4-Chloroaniline-d4	10.840	131	713648	28.018	ng/u1	0.00
46) Dimethylphthalate-d6	13.951	166	1417253	30.055	ng/u1	0.00
49) Acenaphthylene-d8	14.228	160	1660349	30.886	ng/u1	0.00
54) 4-Nitrophenol-d4	14.740	143	230333	23.588	ng/u1	0.00
60) Fluorene-d10	15.534	176	1259071	30.745	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	187637	22.887	ng/u1	0.00
73) Anthracene-d10	17.392	188	1985248	32.176	ng/u1	0.00
81) Pyrene-d10	19.628	212	2325714	34.101	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.657	264	2408698	33.677	ng/u1	0.00

Target Compounds Qvalue

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

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