

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100122\
 Data File : BP011853.D
 Acq On : 01 Oct 2022 10:10
 Operator : CG/JU
 Sample : PB147968BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK968

Quant Time: Oct 03 00:57:58 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	215867	20.000	ng/u1	0.00
20) Naphthalene-d8	10.705	136	990802	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.540	164	676107	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.298	188	1489005	20.000	ng/u1	0.00
79) Chrysene-d12	21.380	240	1526937	20.000	ng/u1	0.00
88) Perylene-d12	23.821	264	1694883	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	30542	6.240	ng/uL	0.00
4) Pyridine-d5	3.728	84	371426	25.907	ng/u1	0.00
7) Phenol-d5	7.052	99	558947	29.993	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	338416	32.761	ng/u1	0.00
11) 2-Chlorophenol-d4	7.422	132	472762	32.136	ng/u1	0.00
15) 4-Methylphenol-d8	8.599	113	487096	31.619	ng/u1	0.00
21) Nitrobenzene-d5	9.058	128	248111	32.876	ng/u1	0.00
24) 2-Nitrophenol-d4	9.781	143	273579	33.883	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	466147	31.212	ng/u1	0.00
31) 4-Chloroaniline-d4	10.840	131	731217	32.116	ng/u1	0.00
46) Dimethylphthalate-d6	13.951	166	1903094	39.362	ng/u1	0.00
49) Acenaphthylene-d8	14.234	160	2012584	36.514	ng/u1	0.00
54) 4-Nitrophenol-d4	14.734	143	330661	33.026	ng/u1	0.00
60) Fluorene-d10	15.534	176	1578090	37.584	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	293519	32.568	ng/u1	0.00
73) Anthracene-d10	17.398	188	2633854	38.832	ng/u1	0.00
81) Pyrene-d10	19.628	212	3090808	39.969	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.663	264	3420987	40.321	ng/u1	0.00

Target Compounds Qvalue

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

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