

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011949.D
 Acq On : 06 Oct 2022 02:31
 Operator : CG/JU
 Sample : PB148042BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK042

Quant Time: Oct 06 08:48:20 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 05 23:18:55 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	227811	20.000	ng/u1	0.00
20) Naphthalene-d8	10.687	136	973198	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.528	164	597072	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.281	188	1255230	20.000	ng/u1	0.00
79) Chrysene-d12	21.369	240	1199917	20.000	ng/u1	0.00
88) Perylene-d12	23.798	264	1193780	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	40159	7.775	ng/uL	0.00
4) Pyridine-d5	3.723	84	507372	33.534	ng/u1	0.00
7) Phenol-d5	7.040	99	689854	35.077	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	392777	36.030	ng/u1	0.00
11) 2-Chlorophenol-d4	7.411	132	578294	37.249	ng/u1	0.00
15) 4-Methylphenol-d8	8.593	113	551473	33.921	ng/u1	0.00
21) Nitrobenzene-d5	9.046	128	271730	36.657	ng/u1	0.00
24) 2-Nitrophenol-d4	9.769	143	290335	36.609	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.305	165	528054	35.997	ng/u1	0.00
31) 4-Chloroaniline-d4	10.828	131	797682	35.669	ng/u1	0.00
46) Dimethylphthalate-d6	13.940	166	1613118	37.781	ng/u1	0.00
49) Acenaphthylene-d8	14.222	160	1857033	38.152	ng/u1	0.00
54) 4-Nitrophenol-d4	14.728	143	263483	29.800	ng/u1	0.00
60) Fluorene-d10	15.522	176	1459346	39.357	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	200967	26.452	ng/u1	0.00
73) Anthracene-d10	17.381	188	2294851	40.135	ng/u1	0.00
81) Pyrene-d10	19.616	212	2589034	42.604	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.639	264	2472815	41.380	ng/u1	0.00

Target Compounds Qvalue

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

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