

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010838.D
 Acq On : 25 Jun 2022 21:15
 Operator : CG/JU
 Sample : PB145746BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK746

Quant Time: Jun 26 22:54:21 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 24 22:42:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	406096	20.000	ng/u1	0.00
20) Naphthalene-d8	10.510	136	1638771	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.374	164	853048	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.133	188	1574635	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.245	240	1513049	20.000	ng/u1	#-0.01
88) Perylene-d12	23.609	264	1598603	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.210	96	77145	7.186	ng/uL	0.00
4) Pyridine-d5	3.610	84	976020	34.155	ng/u1	0.00
7) Phenol-d5	6.898	99	1151452	35.819	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.057	67	741282	35.259	ng/u1	0.00
11) 2-Chlorophenol-d4	7.251	132	983221	36.865	ng/u1	0.00
15) 4-Methylphenol-d8	8.434	113	908140	35.736	ng/u1	0.00
21) Nitrobenzene-d5	8.881	128	442324	42.399	ng/u1	0.00
24) 2-Nitrophenol-d4	9.598	143	423525	49.071	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.139	165	756542	35.475	ng/u1	0.00
31) 4-Chloroaniline-d4	10.657	131	1228630	35.294	ng/u1	0.00
46) Dimethylphthalate-d6	13.792	166	2186554	37.419	ng/u1	0.00
49) Acenaphthylene-d8	14.063	160	2664288	35.725	ng/u1	0.00
54) 4-Nitrophenol-d4	14.586	143	331378	36.307	ng/u1	0.00
60) Fluorene-d10	15.374	176	1831623	36.526	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.492	200	155325	29.773	ng/u1	0.00
73) Anthracene-d10	17.233	188	2648060	37.798	ng/u1	-0.01
81) Pyrene-d10	19.486	212	2913037	35.339	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.462	264	3117194	39.460	ng/u1	-0.01

Target Compounds	Qvalue

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

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