

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102323\
 Data File : BP017747.D
 Acq On : 23 Oct 2023 10:16
 Operator : MA/JU
 Sample : PB156466BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK466

Quant Time: Oct 24 04:56:36 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 19 01:32:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	114304	20.000	ng/u1	0.00
20) Naphthalene-d8	10.716	136	456917	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.586	164	284404	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.374	188	573506	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.521	240	464779	20.000	ng/u1	0.00
88) Perylene-d12	24.074	264	471057	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.257	96	22869	7.443	ng/uL	0.00
4) Pyridine-d5	3.687	84	287495	35.738	ng/u1	0.00
7) Phenol-d5	7.045	99	340185	35.765	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	232161	39.522	ng/u1	0.00
11) 2-Chlorophenol-d4	7.404	132	282775	36.450	ng/u1	0.00
15) 4-Methylphenol-d8	8.604	113	260745	34.908	ng/u1	0.00
21) Nitrobenzene-d5	9.092	128	124330	33.580	ng/u1	0.00
24) 2-Nitrophenol-d4	9.810	143	138340	33.817	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.339	165	243226	30.310	ng/u1	0.00
31) 4-Chloroaniline-d4	10.892	131	322694	30.728	ng/u1	0.00
46) Dimethylphthalate-d6	14.004	166	799832	33.501	ng/u1	0.00
49) Acenaphthylene-d8	14.280	160	861748	32.366	ng/u1	0.00
54) 4-Nitrophenol-d4	14.833	143	68641	20.177	ng/u1	0.00
60) Fluorene-d10	15.586	176	688159	33.344	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.733	200	93991	24.881	ng/u1	0.00
73) Anthracene-d10	17.474	188	1045770	35.683	ng/u1	0.00
81) Pyrene-d10	19.733	212	1186314	40.712	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.903	264	939872	35.823	ng/u1	0.00

Target Compounds Qvalue

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

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