

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017378.D
 Acq On : 27 Sep 2023 00:58
 Operator : MA/JU
 Sample : PB155788BL
 Misc :
 ALS Vial : 71 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK788

Quant Time: Sep 27 02:24:02 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 26 04:52:29 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	142121	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	629425	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.581	164	403236	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.351	188	899724	20.000	ng/u1	0.00
79) Chrysene-d12	21.457	240	801806	20.000	ng/u1	0.00
88) Perylene-d12	23.969	264	877871	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	27269	7.881	ng/uL	0.00
4) Pyridine-d5	3.728	84	380951	39.383	ng/u1	0.00
7) Phenol-d5	7.075	99	456191	39.083	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.264	67	284746	40.423	ng/u1	0.00
11) 2-Chlorophenol-d4	7.440	132	373780	40.375	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	365478	40.274	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	180499	39.667	ng/u1	0.00
24) 2-Nitrophenol-d4	9.834	143	196856	39.842	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.358	165	359322	38.493	ng/u1	0.00
31) 4-Chloroaniline-d4	10.910	131	532314	39.366	ng/u1	0.00
46) Dimethylphthalate-d6	13.999	166	1172856	40.602	ng/u1	0.00
49) Acenaphthylene-d8	14.281	160	1273391	38.328	ng/u1	0.00
54) 4-Nitrophenol-d4	14.816	143	144649	28.743	ng/u1	0.00
60) Fluorene-d10	15.581	176	1003920	40.369	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	103163	20.344	ng/u1	0.00
73) Anthracene-d10	17.451	188	1621664	40.640	ng/u1	0.00
81) Pyrene-d10	19.698	212	1902515	43.878	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.804	264	1769444	41.864	ng/u1	0.00

Target Compounds Qvalue

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

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