

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017359.D
 Acq On : 26 Sep 2023 13:37
 Operator : MA/JU
 Sample : PB155787BL
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK787

Quant Time: Sep 27 00:28:01 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	166608	20.000	ng/ul	0.00
20) Naphthalene-d8	10.745	136	747207	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.586	164	485968	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.357	188	1138566	20.000	ng/ul	0.00
79) Chrysene-d12	21.462	240	1056479	20.000	ng/ul	0.00
88) Perylene-d12	23.974	264	1185791	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	25384	6.258	ng/uL	0.00
4) Pyridine-d5	3.728	84	375359	33.101	ng/ul	0.00
7) Phenol-d5	7.081	99	456334	33.349	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	279083	33.796	ng/ul	0.00
11) 2-Chlorophenol-d4	7.446	132	369965	34.089	ng/ul	0.00
15) 4-Methylphenol-d8	8.634	113	370757	34.851	ng/ul	0.00
21) Nitrobenzene-d5	9.122	128	181770	33.650	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	208377	35.526	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	365525	32.985	ng/ul	0.00
31) 4-Chloroaniline-d4	10.910	131	540674	33.682	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	1210379	34.768	ng/ul	0.00
49) Acenaphthylene-d8	14.286	160	1320178	32.972	ng/ul	0.00
54) 4-Nitrophenol-d4	14.816	143	178449	29.422	ng/ul	0.00
60) Fluorene-d10	15.580	176	1042950	34.799	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	151048	23.538	ng/ul	0.00
73) Anthracene-d10	17.457	188	1751213	34.680	ng/ul	0.00
81) Pyrene-d10	19.698	212	2069426	36.222	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.809	264	1979000	34.663	ng/ul	0.00

Target Compounds Qvalue

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

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