

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017344.D
 Acq On : 26 Sep 2023 04:55
 Operator : MA/JU
 Sample : PB155729BL
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK729

Quant Time: Sep 26 05:26:06 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	220835	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	963236	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.587	164	598386	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.357	188	1319266	20.000	ng/u1	0.00
79) Chrysene-d12	21.457	240	1202593	20.000	ng/u1	0.00
88) Perylene-d12	23.974	264	1324320	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	32251	5.999	ng/uL	0.00
4) Pyridine-d5	3.728	84	475546	31.639	ng/u1	0.00
7) Phenol-d5	7.081	99	578269	31.883	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	361823	33.056	ng/u1	0.00
11) 2-Chlorophenol-d4	7.446	132	468226	32.549	ng/u1	0.00
15) 4-Methylphenol-d8	8.640	113	468045	33.193	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	225308	32.355	ng/u1	0.00
24) 2-Nitrophenol-d4	9.840	143	255474	33.787	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	454452	31.813	ng/u1	0.00
31) 4-Chloroaniline-d4	10.910	131	664718	32.122	ng/u1	0.00
46) Dimethylphthalate-d6	14.004	166	1453913	33.917	ng/u1	0.00
49) Acenaphthylene-d8	14.287	160	1594989	32.352	ng/u1	0.00
54) 4-Nitrophenol-d4	14.816	143	198080	26.523	ng/u1	0.00
60) Fluorene-d10	15.581	176	1241544	33.642	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	151472	20.371	ng/u1	0.00
73) Anthracene-d10	17.457	188	1974025	33.738	ng/u1	0.00
81) Pyrene-d10	19.698	212	2309454	35.512	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.810	264	2163897	33.937	ng/u1	0.00

Target Compounds Qvalue

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

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