

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017379.D
 Acq On : 27 Sep 2023 01:32
 Operator : MA/JU
 Sample : PB155793BL
 Misc :
 ALS Vial : 72 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK793

Quant Time: Sep 27 02:25:45 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	159434	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	698397	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.581	164	434879	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.351	188	971445	20.000	ng/u1	0.00
79) Chrysene-d12	21.457	240	883993	20.000	ng/u1	0.00
88) Perylene-d12	23.968	264	957516	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	26648	6.865	ng/uL	0.00
4) Pyridine-d5	3.728	84	384225	35.408	ng/u1	0.00
7) Phenol-d5	7.081	99	462267	35.303	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.263	67	292535	37.019	ng/u1	0.00
11) 2-Chlorophenol-d4	7.440	132	374549	36.065	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	369769	36.322	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	180061	35.663	ng/u1	0.00
24) 2-Nitrophenol-d4	9.834	143	196382	35.821	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	363826	35.126	ng/u1	0.00
31) 4-Chloroaniline-d4	10.910	131	528499	35.224	ng/u1	0.00
46) Dimethylphthalate-d6	13.998	166	1223772	39.282	ng/u1	0.00
49) Acenaphthylene-d8	14.281	160	1291141	36.035	ng/u1	0.00
54) 4-Nitrophenol-d4	14.810	143	146593	27.009	ng/u1	0.00
60) Fluorene-d10	15.581	176	1019292	38.005	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.710	200	110982	20.270	ng/u1	0.00
73) Anthracene-d10	17.451	188	1685658	39.125	ng/u1	0.00
81) Pyrene-d10	19.698	212	1980953	41.439	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.804	264	1848535	40.097	ng/u1	0.00

Target Compounds Qvalue

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

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