

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091224\
 Data File : BG062767.D
 Acq On : 12 Sep 2024 18:23
 Operator : JU/RC
 Sample : PB163313BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK313

Quant Time: Sep 12 23:03:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 10 01:28:17 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	45694	20.000	ng/u1	0.00
20) Naphthalene-d8	10.696	136	206570	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.533	164	168962	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.276	188	449146	20.000	ng/u1	0.00
79) Chrysene-d12	21.519	240	487744	20.000	ng/u1	0.00
88) Perylene-d12	24.586	264	552073	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.357	96	6235	6.168	ng/uL	0.00
4) Pyridine-d5	3.763	84	98796	30.704	ng/u1	0.00
7) Phenol-d5	7.071	99	100442	25.068	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.229	67	57821	24.069	ng/u1	0.00
11) 2-Chlorophenol-d4	7.435	132	76685	26.275	ng/u1	0.00
15) 4-Methylphenol-d8	8.604	113	87358	26.163	ng/u1	0.00
21) Nitrobenzene-d5	9.057	128	40665	27.544	ng/u1	0.00
24) 2-Nitrophenol-d4	9.773	143	49489	30.650	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.320	165	96331	27.395	ng/u1	0.00
31) 4-Chloroaniline-d4	10.825	131	111953	23.162	ng/u1	0.00
46) Dimethylphthalate-d6	13.933	166	329180	25.300	ng/u1	0.00
49) Acenaphthylene-d8	14.221	160	371354	25.698	ng/u1	0.00
54) 4-Nitrophenol-d4	14.726	143	48297	21.851	ng/u1	0.00
60) Fluorene-d10	15.526	176	315628	26.920	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.637	200	50154	19.413	ng/u1	0.00
73) Anthracene-d10	17.376	188	580459	27.384	ng/u1	0.00
81) Pyrene-d10	19.668	212	761903	27.759	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.374	264	771475	26.465	ng/u1	0.00

Target Compounds Qvalue

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

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