

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080824\
 Data File : BG062347.D
 Acq On : 9 Aug 2024 4:32
 Operator : MA/JU
 Sample : P3434-05
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E05T6

Quant Time: Aug 09 05:18:10 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG080524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 09 01:00:03 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	65167	20.000	ng/u1	0.00
20) Naphthalene-d8	10.748	136	310515	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.573	164	236089	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.316	188	503993	20.000	ng/u1	0.00
79) Chrysene-d12	21.551	240	494458	20.000	ng/u1	0.00
88) Perylene-d12	24.635	264	576299	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.423	96	2773	1.835	ng/uL	0.00
4) Pyridine-d5	3.828	84	45649	9.601	ng/u1	0.00
7) Phenol-d5	7.112	99	86560	13.916	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.288	67	50644	13.227	ng/u1	0.00
11) 2-Chlorophenol-d4	7.488	132	61761	13.859	ng/u1	0.00
15) 4-Methylphenol-d8	8.651	113	71987	13.895	ng/u1	0.00
21) Nitrobenzene-d5	9.109	128	31590	16.499	ng/u1	0.00
24) 2-Nitrophenol-d4	9.832	143	35126	19.502	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.361	165	78407	14.988	ng/u1	0.00
31) 4-Chloroaniline-d4	10.878	131	99338	13.092	ng/u1	0.00
46) Dimethylphthalate-d6	13.979	166	275336	15.116	ng/u1	0.00
49) Acenaphthylene-d8	14.267	160	313008	15.156	ng/u1	0.00
54) 4-Nitrophenol-d4	14.749	143	36601	14.353	ng/u1	0.00
60) Fluorene-d10	15.565	176	255005	15.597	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.671	200	26286	17.611	ng/u1	0.00
73) Anthracene-d10	17.410	188	375403	15.443	ng/u1	0.00
81) Pyrene-d10	19.689	212	483672	16.457	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.424	264	506070	16.427	ng/u1	0.00

Target Compounds	Qvalue

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

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