

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
 Data File : BG062784.D
 Acq On : 13 Sep 2024 12:26
 Operator : JU/RC
 Sample : P3952-12
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0B51

Quant Time: Sep 13 23:00:33 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	7.899	152	44278	20.000	ng/u1	0.00
20) Naphthalene-d8	10.689	136	190533	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.526	164	152206	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.270	188	401617	20.000	ng/u1	0.00
79) Chrysene-d12	21.518	240	454136	20.000	ng/u1	0.00
88) Perylene-d12	24.579	264	551630	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	5299	5.410	ng/uL	0.00
4) Pyridine-d5	3.768	84	33110	10.619	ng/u1	0.00
7) Phenol-d5	7.070	99	32444	8.356	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.229	67	63916	27.457	ng/u1	0.00
11) 2-Chlorophenol-d4	7.434	132	73800	26.095	ng/u1	0.00
15) 4-Methylphenol-d8	8.604	113	65796	20.335	ng/u1	0.00
21) Nitrobenzene-d5	9.050	128	41274	30.309	ng/u1	0.00
24) 2-Nitrophenol-d4	9.773	143	48270	32.411	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.313	165	103436	31.892	ng/u1	0.00
31) 4-Chloroaniline-d4	10.824	131	112849	25.313	ng/u1	0.00
46) Dimethylphthalate-d6	13.933	166	395757	33.765	ng/u1	0.00
49) Acenaphthylene-d8	14.220	160	414696	31.857	ng/u1	0.00
54) 4-Nitrophenol-d4	14.726	143	15072	7.570	ng/u1	0.00
60) Fluorene-d10	15.519	176	348925	33.036	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.636	200	69291	29.994	ng/u1	0.00
73) Anthracene-d10	17.370	188	686144	36.201	ng/u1	0.00
81) Pyrene-d10	19.661	212	912565	35.709	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.367	264	1062163	36.466	ng/u1	-0.01

Target Compounds	Qvalue

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

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