

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110824\
 Data File : BG063235.D
 Acq On : 8 Nov 2024 19:15
 Operator : RC/JU
 Sample : P4627-13
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0CF3

Quant Time: Nov 08 22:08:05 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 06 15:45:52 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.837	152	100503	20.000	ng/u1	0.00
20) Naphthalene-d8	10.628	136	463977	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.476	164	396219	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.226	188	975163	20.000	ng/u1	0.00
79) Chrysene-d12	21.480	240	1005011	20.000	ng/u1	0.00
88) Perylene-d12	24.511	264	1058360	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	9305	4.186	ng/uL	0.00
4) Pyridine-d5	3.712	84	29223	3.978	ng/u1	0.00
7) Phenol-d5	7.008	99	67435	7.436	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.167	67	161504	30.361	ng/u1	0.00
11) 2-Chlorophenol-d4	7.373	132	164732	27.985	ng/u1	0.00
15) 4-Methylphenol-d8	8.542	113	135699	17.654	ng/u1	0.00
21) Nitrobenzene-d5	8.988	128	99637	30.546	ng/u1	0.00
24) 2-Nitrophenol-d4	9.711	143	129643	30.751	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.252	165	246646	29.812	ng/u1	0.00
31) 4-Chloroaniline-d4	10.763	131	205665	19.085	ng/u1	0.00
46) Dimethylphthalate-d6	13.877	166	950559	30.628	ng/u1	0.00
49) Acenaphthylene-d8	14.165	160	984575	32.411	ng/u1	0.00
54) 4-Nitrophenol-d4	14.688	143	29286	6.892	ng/u1	0.00
60) Fluorene-d10	15.469	176	842361	32.159	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.593	200	196593	32.078	ng/u1	0.00
73) Anthracene-d10	17.326	188	1578468	37.662	ng/u1	0.00
81) Pyrene-d10	19.623	212	1959472	38.887	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.312	264	1942943	38.023	ng/u1	0.00

Target Compounds	Qvalue

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

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