

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110824\
 Data File : BG063271.D
 Acq On : 9 Nov 2024 19:17
 Operator : RC/JU
 Sample : P4744-04
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GCP58

Quant Time: Nov 10 00:39:01 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	7.836	152	127584	20.000	ng/u1	0.00
20) Naphthalene-d8	10.621	136	624263	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.469	164	532213	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.219	188	1211951	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.473	240	1236664	20.000	ng/u1	# 0.00
88) Perylene-d12	24.510	264	1294552	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.300	96	12762	4.523	ng/uL	0.00
4) Pyridine-d5	3.705	84	74918	8.033	ng/u1	0.00
7) Phenol-d5	7.007	99	106344	9.237	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.166	67	213450	31.609	ng/u1	0.00
11) 2-Chlorophenol-d4	7.366	132	228531	30.583	ng/u1	-0.01
15) 4-Methylphenol-d8	8.541	113	207093	21.223	ng/u1	0.00
21) Nitrobenzene-d5	8.987	128	140603	32.037	ng/u1	0.00
24) 2-Nitrophenol-d4	9.710	143	182191	32.120	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.245	165	341798	30.705	ng/u1	0.00
31) 4-Chloroaniline-d4	10.756	131	385380	26.580	ng/u1	0.00
46) Dimethylphthalate-d6	13.876	166	1160622	27.840	ng/u1	0.00
49) Acenaphthylene-d8	14.164	160	1275224	31.252	ng/u1	0.00
54) 4-Nitrophenol-d4	14.681	143	37446	6.560	ng/u1	0.00
60) Fluorene-d10	15.468	176	1065860	30.294	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.591	200	205347	26.960	ng/u1	0.00
73) Anthracene-d10	17.319	188	1744873	33.499	ng/u1	0.00
81) Pyrene-d10	19.622	212	2164189	34.905	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.305	264	2197782	35.163	ng/u1	0.00

Target Compounds	Qvalue

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

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