

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070824\
 Data File : BG061942.D
 Acq On : 8 Jul 2024 14:45
 Operator : MA/JU
 Sample : P3059-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E1GM8

Quant Time: Jul 08 15:31:57 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 03 02:38:46 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.094	152	37055	20.000	ng/u1	0.00
20) Naphthalene-d8	10.908	136	200935	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.716	164	155234	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.459	188	354842	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.719	240	318993	20.000	ng/u1	0.00
88) Perylene-d12	24.951	264	384480	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.482	96	4902	4.971	ng/uL	0.00
4) Pyridine-d5	3.905	84	35516	11.714	ng/u1	0.00
7) Phenol-d5	7.242	99	46803	11.542	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.418	67	87015	33.884	ng/u1	0.00
11) 2-Chlorophenol-d4	7.624	132	98925	35.564	ng/u1	0.00
15) 4-Methylphenol-d8	8.793	113	83443	26.134	ng/u1	0.00
21) Nitrobenzene-d5	9.257	128	61951	40.539	ng/u1	0.00
24) 2-Nitrophenol-d4	9.980	143	71785	41.134	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.521	165	134640	39.999	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.038	131	160005	32.741	ng/u1	0.00
46) Dimethylphthalate-d6	14.116	166	504658	39.543	ng/u1	0.00
49) Acenaphthylene-d8	14.416	160	529693	39.693	ng/u1	0.00
54) 4-Nitrophenol-d4	14.909	143	23396	11.479	ng/u1	0.00
60) Fluorene-d10	15.709	176	444228	41.347	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.820	200	84039	45.096	ng/u1	0.00
73) Anthracene-d10	17.559	188	716531	43.251	ng/u1	0.00
81) Pyrene-d10	19.839	212	805795	42.946	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.739	264	863007	45.259	ng/u1	0.00

Target Compounds	Qvalue

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

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