

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058984.D
 Acq On : 11 Sep 2023 15:28
 Operator : MA/JU
 Sample : 04195-08
 Misc :
 ALS Vial : 83 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YC1B7

Quant Time: Sep 12 00:50:58 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 11 00:07:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.324	152	68886	20.000	ng/ul	0.00
20) Naphthalene-d8	11.168	136	338504	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.945	164	270072	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.695	188	747158	20.000	ng/ul	# 0.00
79) Chrysene-d12	22.019	240	735405	20.000	ng/ul	0.00
88) Perylene-d12	25.580	264	848140	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.641	96	5419	3.344	ng/uL	0.00
4) Pyridine-d5	4.070	84	2277	0.444	ng/ul	0.00
7) Phenol-d5	7.431	99	127972	18.422	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.642	67	75387	18.508	ng/ul	0.00
11) 2-Chlorophenol-d4	7.836	132	83755	18.799	ng/ul	0.00
15) 4-Methylphenol-d8	8.999	113	94230	17.558	ng/ul	0.00
21) Nitrobenzene-d5	9.517	128	42559	17.333	ng/ul	0.00
24) 2-Nitrophenol-d4	10.239	143	48779	18.105	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.762	165	99926	17.451	ng/ul	0.00
31) 4-Chloroaniline-d4	11.309	131	61957	8.046	ng/ul	0.00
46) Dimethylphthalate-d6	14.340	166	354156	16.906	ng/ul	0.00
49) Acenaphthylene-d8	14.652	160	403195	17.284	ng/ul	0.00
54) 4-Nitrophenol-d4	15.110	143	56205	16.152	ng/ul	0.00
60) Fluorene-d10	15.932	176	350590	18.858	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.032	200	49820	13.584	ng/ul	0.00
73) Anthracene-d10	17.795	188	602270	18.158	ng/ul	0.00
81) Pyrene-d10	20.063	212	724404	18.595	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.333	264	706950	17.004	ng/ul	0.00

Target Compounds	Qvalue

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

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