

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058978.D
 Acq On : 11 Sep 2023 11:21
 Operator : MA/JU
 Sample : 04195-11
 Misc :
 ALS Vial : 77 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YBY69

Quant Time: Sep 12 00:49:20 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	8.325	152	88540	20.000	ng/ul	0.00
20) Naphthalene-d8	11.169	136	415824	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.953	164	307500	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.697	188	792424	20.000	ng/ul	# 0.00
79) Chrysene-d12	22.021	240	769699	20.000	ng/ul	0.00
88) Perylene-d12	25.587	264	877445	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.643	96	8380	4.023	ng/uL	0.00
4) Pyridine-d5	4.071	84	177	0.027	ng/ul	0.00
7) Phenol-d5	7.432	99	145207	16.263	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.644	67	88101	16.828	ng/ul	0.00
11) 2-Chlorophenol-d4	7.838	132	96350	16.826	ng/ul	0.00
15) 4-Methylphenol-d8	9.007	113	86251	12.504	ng/ul	0.00
21) Nitrobenzene-d5	9.524	128	51705	17.142	ng/ul	0.00
24) 2-Nitrophenol-d4	10.241	143	62386	18.850	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.764	165	113690	16.163	ng/ul	0.00
31) 4-Chloroaniline-d4	11.316	131	21374	2.260	ng/ul	0.00
46) Dimethylphthalate-d6	14.342	166	408370	17.121	ng/ul	0.00
49) Acenaphthylene-d8	14.653	160	438794	16.520	ng/ul	0.00
54) 4-Nitrophenol-d4	15.111	143	62014	15.652	ng/ul	0.00
60) Fluorene-d10	15.934	176	377872	17.851	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.040	200	56249	14.461	ng/ul	0.00
73) Anthracene-d10	17.797	188	617664	17.558	ng/ul	0.00
81) Pyrene-d10	20.064	212	747550	18.335	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.335	264	674537	15.682	ng/ul	0.00

Target Compounds Qvalue

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

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