

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058979.D
 Acq On : 11 Sep 2023 12:02
 Operator : MA/JU
 Sample : 04195-12
 Misc :
 ALS Vial : 78 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YBY74

Quant Time: Sep 12 00:49:28 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	8.323	152	67058	20.000	ng/ul	0.00
20) Naphthalene-d8	11.167	136	343722	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.951	164	262228	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.700	188	702829	20.000	ng/ul	# 0.00
79) Chrysene-d12	22.019	240	950176	20.000	ng/ul	# 0.00
88) Perylene-d12	25.585	264	1081263	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.640	96	7121	4.514	ng/uL	0.00
4) Pyridine-d5	4.063	84	1593	0.319	ng/ul	0.00
7) Phenol-d5	7.436	99	127913	18.916	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.641	67	79956	20.164	ng/ul	0.00
11) 2-Chlorophenol-d4	7.835	132	87508	20.177	ng/ul	0.00
15) 4-Methylphenol-d8	9.005	113	93179	17.836	ng/ul	0.00
21) Nitrobenzene-d5	9.516	128	43058	17.270	ng/ul	0.00
24) 2-Nitrophenol-d4	10.232	143	52935	19.349	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.761	165	98791	16.991	ng/ul	0.00
31) 4-Chloroaniline-d4	11.308	131	40443	5.173	ng/ul	0.00
46) Dimethylphthalate-d6	14.345	166	368849	18.134	ng/ul	0.00
49) Acenaphthylene-d8	14.651	160	411690	18.176	ng/ul	0.00
54) 4-Nitrophenol-d4	15.109	143	55489	16.423	ng/ul	0.00
60) Fluorene-d10	15.938	176	348804	19.323	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.037	200	45494	13.187	ng/ul	0.00
73) Anthracene-d10	17.794	188	588837	18.873	ng/ul	0.00
81) Pyrene-d10	20.068	212	978983	19.450	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.338	264	886267	16.721	ng/ul	0.00

Target Compounds Qvalue

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

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