

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091223\
 Data File : BG059016.D
 Acq On : 12 Sep 2023 18:39
 Operator : MA/JU
 Sample : 04175-09
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EZYK3

Quant Time: Sep 13 00:16:39 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 12 23:54:19 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.309	152	102865	20.000	ng/ul	0.00
20) Naphthalene-d8	11.153	136	497805	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.937	164	374296	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.686	188	895656	20.000	ng/ul	# 0.00
79) Chrysene-d12	22.005	240	765773	20.000	ng/ul	0.00
88) Perylene-d12	25.571	264	887621	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.632	96	6538	2.702	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.422	99	153767	14.824	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.627	67	98507	16.195	ng/ul	0.00
11) 2-Chlorophenol-d4	7.827	132	90369	13.584	ng/ul	0.00
15) 4-Methylphenol-d8	8.991	113	107203	13.377	ng/ul	0.00
21) Nitrobenzene-d5	9.508	128	53918	14.932	ng/ul	0.00
24) 2-Nitrophenol-d4	10.224	143	21957	5.542	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.747	165	62840	7.463	ng/ul	0.00
31) 4-Chloroaniline-d4	11.294	131	170930	15.095	ng/ul	0.00
46) Dimethylphthalate-d6	14.331	166	509218	17.539	ng/ul	0.00
49) Acenaphthylene-d8	14.637	160	585365	18.105	ng/ul	0.00
54) 4-Nitrophenol-d4	15.113	143	648	0.134	ng/ul	0.00
60) Fluorene-d10	15.924	176	508294	19.727	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.029	200	608	0.138	ng/ul	0.00
73) Anthracene-d10	17.786	188	774917	19.490	ng/ul	0.00
81) Pyrene-d10	20.054	212	849301	20.937	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.319	264	846942	19.465	ng/ul	0.00
Target Compounds						
					Qvalue	
16) Acetophenone	9.155	105	29597	2.228	ng/ul	98

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

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