

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058975.D
 Acq On : 11 Sep 2023 8:37
 Operator : MA/JU
 Sample : 04195-10
 Misc :
 ALS Vial : 73 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YBY64

Quant Time: Sep 11 09:12:07 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	74344	20.000	ng/ul	# 0.00
20) Naphthalene-d8	11.167	136	365716	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.951	164	271058	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.695	188	702347	20.000	ng/ul	# 0.00
79) Chrysene-d12	22.019	240	671207	20.000	ng/ul	0.00
88) Perylene-d12	25.585	264	772737	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.635	96	9345	5.343	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.430	99	222736	29.710	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.642	67	121364	27.608	ng/ul	0.00
11) 2-Chlorophenol-d4	7.841	132	102162	21.247	ng/ul	0.00
15) 4-Methylphenol-d8	8.999	113	114945	19.846	ng/ul	0.00
21) Nitrobenzene-d5	9.522	128	50923	19.196	ng/ul	0.00
24) 2-Nitrophenol-d4	10.239	143	60793	20.885	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.756	165	112160	18.130	ng/ul	0.00
31) 4-Chloroaniline-d4	11.308	131	112008	13.464	ng/ul	0.00
46) Dimethylphthalate-d6	14.345	166	398419	18.949	ng/ul	0.00
49) Acenaphthylene-d8	14.651	160	433824	18.529	ng/ul	0.00
54) 4-Nitrophenol-d4	15.109	143	65975	18.890	ng/ul	0.00
60) Fluorene-d10	15.932	176	361830	19.391	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.038	200	52383	15.194	ng/ul	0.00
73) Anthracene-d10	17.794	188	604970	19.403	ng/ul	0.00
81) Pyrene-d10	20.062	212	700318	19.696	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.333	264	729755	19.265	ng/ul	0.00

Target Compounds Qvalue

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

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