

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058969.D
 Acq On : 11 Sep 2023 4:29
 Operator : MA/JU
 Sample : 04195-02
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YBY24

Quant Time: Sep 11 05:17:25 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	71480	20.000	ng/u1	0.00
20) Naphthalene-d8	11.167	136	347237	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.951	164	247041	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.701	188	690552	20.000	ng/u1	# 0.00
79) Chrysene-d12	22.019	240	680781	20.000	ng/u1	0.00
88) Perylene-d12	25.586	264	793463	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.635	96	2819	1.676	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	7.430	99	53036	7.358	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.648	67	33499	7.926	ng/u1	0.00
11) 2-Chlorophenol-d4	7.842	132	36237	7.838	ng/u1	0.00
15) 4-Methylphenol-d8	8.999	113	31189	5.601	ng/u1	0.00
21) Nitrobenzene-d5	9.522	128	17126	6.800	ng/u1	0.00
24) 2-Nitrophenol-d4	10.239	143	20308	7.348	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.762	165	40238	6.851	ng/u1	0.00
31) 4-Chloroaniline-d4	11.308	131	33284	4.214	ng/u1	0.00
46) Dimethylphthalate-d6	14.340	166	153157	7.992	ng/u1	0.00
49) Acenaphthylene-d8	14.651	160	161267	7.557	ng/u1	0.00
54) 4-Nitrophenol-d4	15.110	143	20005	6.285	ng/u1	0.00
60) Fluorene-d10	15.932	176	140982	8.290	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	16.038	200	19416	5.728	ng/u1	0.00
73) Anthracene-d10	17.795	188	237426	7.745	ng/u1	0.00
81) Pyrene-d10	20.063	212	296587	8.224	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.327	264	302336	7.773	ng/u1	-0.01
Target Compounds						
					Qvalue	
16) Acetophenone	9.170	105	15638	1.694	ng/u1	97

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

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