

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101023\
 Data File : BG059427.D
 Acq On : 10 Oct 2023 17:18
 Operator : MA/JU
 Sample : 04654-13
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BBHZ9

Quant Time: Oct 10 22:51:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 07 22:45:22 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.220	152	43154	20.000	ng/ul	0.00
20) Naphthalene-d8	11.064	136	200051	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.853	164	128557	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.603	188	288604	20.000	ng/ul	-0.01
79) Chrysene-d12	21.904	240	260810	20.000	ng/ul	-0.01
88) Perylene-d12	25.382	264	306165	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.520	96	5872	4.852	ng/uL	0.00
4) Pyridine-d5	3.960	84	29063	8.342	ng/ul	0.00
7) Phenol-d5	7.350	99	36041	7.564	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.544	67	99234	36.846	ng/ul	0.00
11) 2-Chlorophenol-d4	7.738	132	95841	28.759	ng/ul	-0.01
15) 4-Methylphenol-d8	8.913	113	64011	17.242	ng/ul	-0.01
21) Nitrobenzene-d5	9.419	128	68434	40.185	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.135	143	67129	36.420	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.664	165	121952	36.657	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.205	131	128628	24.340	ng/ul	-0.01
46) Dimethylphthalate-d6	14.254	166	438622	41.646	ng/ul	0.00
49) Acenaphthylene-d8	14.560	160	520069	41.494	ng/ul	0.00
54) 4-Nitrophenol-d4	15.041	143	11105	5.627	ng/ul	-0.02
60) Fluorene-d10	15.846	176	404769	40.262	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.958	200	47258	24.801	ng/ul	0.00
73) Anthracene-d10	17.703	188	626701	44.266	ng/ul	-0.01
81) Pyrene-d10	19.977	212	724134	45.832	ng/ul	-0.01
92) Benzo(a)pyrene-d12	25.141	264	727974	45.185	ng/ul	-0.01

Target Compounds Qvalue

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

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