

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111723\
 Data File : BG059729.D
 Acq On : 17 Nov 2023 12:44
 Operator : MA/JU
 Sample : PB157066BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Nov 20 16:02:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 17 01:51:38 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.338	152	29713	20.000	ng	0.00
21) Naphthalene-d8	11.187	136	129626	20.000	ng	# 0.00
39) Acenaphthene-d10	14.971	164	103173	20.000	ng	0.00
64) Phenanthrene-d10	17.721	188	257852	20.000	ng	# 0.00
76) Chrysene-d12	22.051	240	241117	20.000	ng	0.00
86) Perylene-d12	25.635	264	296660	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.835	112	239270	135.538	ng	0.00
7) Phenol-d6	7.462	99	345098	131.916	ng	0.00
23) Nitrobenzene-d5	9.542	82	306857	90.410	ng	0.00
42) 2,4,6-Tribromophenol	16.458	330	183333	156.611	ng	0.00
45) 2-Fluorobiphenyl	13.596	172	653318	85.235	ng	0.00
79) Terphenyl-d14	20.294	244	1430720	87.596	ng	0.00
Target Compounds						
19) n-Nitroso-di-n-propyla...	9.542	70	45069	21.877	ng	# 84
51) 2,6-Dinitrotoluene	14.971	165	13391	8.829	ng	# 3
57) 2,4-Dinitrotoluene	14.971	165	13391	5.741	ng	# 6
67) 4-Bromophenyl-phenylether	16.458	248	11878	4.288	ng	# 1
77) Benzidine	20.294	184	40751	7.453	ng	# 96

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

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