

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081022\
 Data File : BG054585.D
 Acq On : 10 Aug 2022 12:41
 Operator : CG/JU
 Sample : N3962-11
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-197

Quant Time: Aug 14 11:20:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 13 00:32:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.971	152	18934	20.000	ng	0.00
21) Naphthalene-d8	10.779	136	72121	20.000	ng	# 0.00
39) Acenaphthene-d10	14.616	164	53470	20.000	ng	0.00
64) Phenanthrene-d10	17.360	188	133813	20.000	ng	# 0.00
76) Chrysene-d12	21.625	240	158835	20.000	ng	# 0.00
86) Perylene-d12	24.833	264	166541	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.562	112	90174	99.436	ng	0.00
7) Phenol-d6	7.178	99	116791	93.988	ng	0.00
23) Nitrobenzene-d5	9.140	82	79234	59.828	ng	0.00
42) 2,4,6-Tribromophenol	16.108	330	95254	84.846	ng	0.00
45) 2-Fluorobiphenyl	13.229	172	235143	62.038	ng	0.00
79) Terphenyl-d14	19.969	244	504765	63.044	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.094	88	944	2.617	ng	# 1
19) n-Nitroso-di-n-propyla...	9.140	70	11482	13.837	ng	# 76
51) 2,6-Dinitrotoluene	14.610	165	7151	7.806	ng	# 16
57) 2,4-Dinitrotoluene	14.610	165	7151	5.445	ng	# 21
67) 4-Bromophenyl-phenylether	16.108	248	5793	3.082	ng	# 1
77) Benzidine	19.969	184	6502	2.260	ng	# 88

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

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