

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080322\
 Data File : BG054496.D
 Acq On : 3 Aug 2022 12:33
 Operator : CG/JU
 Sample : N3935-11
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-141

Quant Time: Aug 04 00:43:20 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 28 04:30:50 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.976	152	16076	20.000	ng	#-0.01
21) Naphthalene-d8	10.791	136	65837	20.000	ng	# 0.00
39) Acenaphthene-d10	14.616	164	53763	20.000	ng	0.00
64) Phenanthrene-d10	17.365	188	144608	20.000	ng	# 0.00
76) Chrysene-d12	21.631	240	162062	20.000	ng	# 0.00
86) Perylene-d12	24.839	264	168515	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.567	112	115010	149.370	ng	0.00
7) Phenol-d6	7.177	99	157853	149.617	ng	0.00
23) Nitrobenzene-d5	9.146	82	101907	84.292	ng	0.00
42) 2,4,6-Tribromophenol	16.114	330	151785	134.463	ng	0.00
45) 2-Fluorobiphenyl	13.241	172	331470	86.975	ng	0.00
79) Terphenyl-d14	19.974	244	734964	89.967	ng	0.00
Target Compounds						
78) Pyrene	19.780	202	22752	2.447	ng	Qvalue 100

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

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