

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080322\
 Data File : BG054497.D
 Acq On : 3 Aug 2022 13:14
 Operator : CG/JU
 Sample : N3935-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-133

Quant Time: Aug 04 00:43:26 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	7.977	152	14891	20.000	ng	-0.01
21) Naphthalene-d8	10.785	136	58429	20.000	ng	#-0.01
39) Acenaphthene-d10	14.616	164	50184	20.000	ng	0.00
64) Phenanthrene-d10	17.360	188	133490	20.000	ng	#-0.01
76) Chrysene-d12	21.631	240	151317	20.000	ng	# 0.00
86) Perylene-d12	24.828	264	160843	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.562	112	93080	130.509	ng	0.00
7) Phenol-d6	7.178	99	128249	131.231	ng	0.00
23) Nitrobenzene-d5	9.146	82	83543	77.864	ng	0.00
42) 2,4,6-Tribromophenol	16.114	330	134271	127.431	ng	0.00
45) 2-Fluorobiphenyl	13.235	172	273587	76.906	ng	-0.01
79) Terphenyl-d14	19.975	244	649620	85.167	ng	0.00
Target Compounds						
78) Pyrene	19.781	202	19380	2.232	ng	Qvalue 99

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

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