

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080322\
 Data File : BG054510.D
 Acq On : 3 Aug 2022 22:21
 Operator : CG/JU
 Sample : N3930-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-95

Quant Time: Aug 04 00:45:04 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	7.982	152	18524	20.000	ng	0.00
21) Naphthalene-d8	10.791	136	74526	20.000	ng	# 0.00
39) Acenaphthene-d10	14.616	164	58010	20.000	ng	0.00
64) Phenanthrene-d10	17.365	188	142135	20.000	ng	# 0.00
76) Chrysene-d12	21.631	240	154763	20.000	ng	# 0.00
86) Perylene-d12	24.833	264	163600	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.567	112	95852	108.037	ng	0.00
7) Phenol-d6	7.183	99	132266	108.798	ng	0.01
23) Nitrobenzene-d5	9.145	82	90471	66.108	ng	0.00
42) 2,4,6-Tribromophenol	16.114	330	122318	100.426	ng	0.00
45) 2-Fluorobiphenyl	13.235	172	280923	68.315	ng	-0.01
79) Terphenyl-d14	19.974	244	544386	69.781	ng	0.00
Target Compounds						
78) Pyrene	19.780	202	23342	2.629	ng	Qvalue 99

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

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