

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080422\
 Data File : BG054528.D
 Acq On : 4 Aug 2022 21:20
 Operator : CG/JU
 Sample : N3931-11
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RVE-122

Quant Time: Aug 05 05:03:29 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	20374	20.000	ng	0.00
21) Naphthalene-d8	10.791	136	78508	20.000	ng	# 0.00
39) Acenaphthene-d10	14.622	164	57390	20.000	ng	0.00
64) Phenanthrene-d10	17.371	188	136899	20.000	ng	# 0.00
76) Chrysene-d12	21.637	240	143116	20.000	ng	# 0.00
86) Perylene-d12	24.851	264	122983	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.573	112	119255	122.210	ng	0.01
7) Phenol-d6	7.189	99	157100	117.492	ng	0.02
23) Nitrobenzene-d5	9.151	82	106318	73.747	ng	0.00
42) 2,4,6-Tribromophenol	16.120	330	131400	109.048	ng	0.00
45) 2-Fluorobiphenyl	13.241	172	319837	78.619	ng	0.00
79) Terphenyl-d14	19.980	244	575225	79.735	ng	0.00

Target Compounds Qvalue

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

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