

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041322\
 Data File : BG053137.D
 Acq On : 13 Apr 2022 14:08
 Operator : CG/JU
 Sample : N2334-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TWO-INCH-WELL

Quant Time: Apr 14 02:56:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 08 15:31:24 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.117	152	34642	20.000	ng	0.00
21) Naphthalene-d8	10.931	136	126410	20.000	ng	# 0.00
39) Acenaphthene-d10	14.743	164	99712	20.000	ng	0.00
64) Phenanthrene-d10	17.487	188	230861	20.000	ng	0.00
76) Chrysene-d12	21.769	240	233805	20.000	ng	# 0.00
86) Perylene-d12	25.065	264	236026	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.685	112	137827	68.201	ng	0.00
7) Phenol-d6	7.277	99	125005	40.111	ng	0.00
23) Nitrobenzene-d5	9.286	82	288555	92.685	ng	0.00
42) 2,4,6-Tribromophenol	16.229	330	229001	144.355	ng	0.00
45) 2-Fluorobiphenyl	13.369	172	664012	91.825	ng	0.00
79) Terphenyl-d14	20.089	244	1189560	85.929	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	10.984	128	20220	2.998	ng	# 54
37) 2-Methylnaphthalene	12.581	142	427323	85.629	ng	# 95
38) 1-Methylnaphthalene	12.799	142	296535	60.875	ng	# 93

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

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