

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042622\
 Data File : BG053334.D
 Acq On : 27 Apr 2022 00:03
 Operator : CG/JU
 Sample : N2548-11
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-11

Quant Time: Apr 27 03:39:09 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Apr 24 01:13:30 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.105	152	20057	20.000	ng	0.00
21) Naphthalene-d8	10.919	136	77144	20.000	ng	# 0.00
39) Acenaphthene-d10	14.726	164	60896	20.000	ng	0.00
64) Phenanthrene-d10	17.475	188	180202	20.000	ng	0.00
76) Chrysene-d12	21.757	240	218533	20.000	ng	0.00
86) Perylene-d12	25.047	264	219826	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.667	112	145539	124.386	ng	0.00
7) Phenol-d6	7.265	99	192546	106.712	ng	0.00
23) Nitrobenzene-d5	9.263	82	162133	85.336	ng	-0.01
42) 2,4,6-Tribromophenol	16.218	330	142423	147.005	ng	0.00
45) 2-Fluorobiphenyl	13.351	172	362893	82.172	ng	0.00
79) Terphenyl-d14	20.077	244	1072797	82.910	ng	0.00

Target Compounds	Qvalue

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

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