

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040522\
 Data File : BG053037.D
 Acq On : 5 Apr 2022 18:33
 Operator : CG/JU
 Sample : N2246-09
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-8-9-2-1

Quant Time: Apr 06 04:15:10 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 05 14:24:42 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.129	152	36189	20.000	ng	#-0.02
21) Naphthalene-d8	10.937	136	136282	20.000	ng	-0.02
39) Acenaphthene-d10	14.744	164	96180	20.000	ng	-0.02
64) Phenanthrene-d10	17.487	188	228266	20.000	ng	-0.02
76) Chrysene-d12	21.770	240	254762	20.000	ng	-0.03
86) Perylene-d12	25.071	264	257007	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.692	112	287710	138.454	ng	0.00
7) Phenol-d6	7.278	99	373057	119.080	ng	-0.01
23) Nitrobenzene-d5	9.287	82	328596	121.016	ng	-0.02
42) 2,4,6-Tribromophenol	16.230	330	212533	164.448	ng	-0.02
45) 2-Fluorobiphenyl	13.375	172	656486	100.667	ng	-0.02
79) Terphenyl-d14	20.096	244	1280891	89.437	ng	-0.02
Target Compounds						
38) 1-Methylnaphthalene	12.805	142	25716	5.105	ng	Qvalue 95
50) Dimethylphthalate	14.186	163	18334	2.565	ng	# 98

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

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