

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040522\
 Data File : BG053038.D
 Acq On : 5 Apr 2022 19:12
 Operator : CG/JU
 Sample : N2246-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

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

Quant Time: Apr 06 04:15:18 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	8.125	152	36050	20.000	ng	-0.02
21) Naphthalene-d8	10.939	136	137798	20.000	ng	-0.02
39) Acenaphthene-d10	14.746	164	95664	20.000	ng	-0.02
64) Phenanthrene-d10	17.489	188	224410	20.000	ng	-0.01
76) Chrysene-d12	21.772	240	249780	20.000	ng	-0.03
86) Perylene-d12	25.073	264	255421	20.000	ng	#-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.693	112	283297	136.856	ng	0.00
7) Phenol-d6	7.280	99	369859	118.514	ng	0.00
23) Nitrobenzene-d5	9.289	82	320830	116.856	ng	-0.01
42) 2,4,6-Tribromophenol	16.232	330	207158	161.154	ng	-0.01
45) 2-Fluorobiphenyl	13.371	172	648239	99.939	ng	-0.02
79) Terphenyl-d14	20.092	244	1230029	87.598	ng	-0.02
Target Compounds						
38) 1-Methylnaphthalene	12.801	142	20865	4.097	ng	# 81
50) Dimethylphthalate	14.188	163	14331	2.016	ng	96

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

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