

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG022122\
 Data File : BG052436.D
 Acq On : 21 Feb 2022 14:15
 Operator : CG/JU
 Sample : N1600-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AA0

Quant Time: Feb 21 14:57:04 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG020722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 08 01:21:20 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.221	152	33713	20.000	ng	0.00
21) Naphthalene-d8	11.064	136	132554	20.000	ng	# 0.00
39) Acenaphthene-d10	14.870	164	86038	20.000	ng	0.00
64) Phenanthrene-d10	17.620	188	193516	20.000	ng	0.00
76) Chrysene-d12	21.931	240	201872	20.000	ng	0.00
86) Perylene-d12	25.397	264	224927	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.753	112	197316	102.006	ng	0.00
7) Phenol-d6	7.369	99	275452	92.291	ng	0.00
23) Nitrobenzene-d5	9.395	82	190607	62.495	ng	0.00
42) 2,4,6-Tribromophenol	16.357	330	115310	93.287	ng	0.00
45) 2-Fluorobiphenyl	13.490	172	386506	62.422	ng	0.00
79) Terphenyl-d14	20.210	244	624771	54.651	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	7.392	94	9859	3.325	ng	91
37) 2-Methylnaphthalene	12.709	142	17862	3.461	ng	94
38) 1-Methylnaphthalene	12.926	142	12163	2.432	ng	# 93

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

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