

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101719\
 Data File : BG043265.D
 Acq On : 17 Oct 2019 17:05
 Operator : JU
 Sample : K5337-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 HHT-WATER

Quant Time: Oct 17 17:55:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 09 01:11:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	49929	20.00	ng	-0.02
21) Naphthalene-d8	10.85	136	199005	20.00	ng	-0.02
39) Acenaphthene-d10	14.67	164	151585	20.00	ng	-0.02
64) Phenanthrene-d10	17.41	188	336405	20.00	ng	0.00
76) Chrysene-d12	21.69	240	364206	20.00	ng	0.00
87) Perylene-d12	24.91	264	426756	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	216151	77.26	ng	-0.02
7) Phenol-d6	7.21	99	203872	50.60	ng	-0.02
23) Nitrobenzene-d5	9.20	82	275657	67.33	ng	-0.02
42) 2,4,6-Tribromophenol	16.16	330	282392	108.34	ng	0.00
45) 2-Fluorobiphenyl	13.29	172	719849	68.84	ng	-0.02
79) Terphenyl-d14	20.02	244	1399505	81.88	ng	0.00
Target Compounds						
10) Phenol	7.24	94	17197	4.652	ng	93
50) Dimethylphthalate	14.11	163	183570	16.163	ng	# 98
74) Di-n-butylphthalate	18.37	149	920376	50.737	ng	100

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

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