

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021419\
 Data File : BG039518.D
 Acq On : 14 Feb 2019 22:36
 Operator : JU/SJ
 Sample : K1479-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OIL-WATER-SEPERATOR

Quant Time: Feb 15 04:54:59 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 29 15:41:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	45346	20.00	ng	-0.02
21) Naphthalene-d8	10.99	136	209042	20.00	ng	-0.02
39) Acenaphthene-d10	14.79	164	148198	20.00	ng	-0.02
64) Phenanthrene-d10	17.54	188	347477	20.00	ng	-0.02
76) Chrysene-d12	21.83	240	373651	20.00	ng	-0.02
87) Perylene-d12	25.21	264	382373	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.72	112	406631	153.48	ng	0.00
7) Phenol-d6	7.32	99	560554	143.73	ng	0.00
23) Nitrobenzene-d5	9.35	82	384854	115.01	ng	-0.01
42) 2,4,6-Tribromophenol	16.28	330	271466	170.11	ng	0.00
45) 2-Fluorobiphenyl	13.42	172	951660	92.12	ng	-0.02
79) Terphenyl-d14	20.13	244	1447236	81.98	ng	-0.02
Target Compounds						
10) Phenol	7.35	94	11651	2.866	ng	85
15) Benzyl Alcohol	8.41	79	44367	14.491	ng	97
20) 3+4-Methylphenols	8.93	107	383295	105.800	ng	89
32) Benzoic acid	10.35	122	262682	117.621	ng	97

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

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