

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031419\
 Data File : BG039933.D
 Acq On : 14 Mar 2019 20:20
 Operator : JU/SJ
 Sample : K1894-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RFK-RMAB-MW12-GW

Quant Time: Mar 15 08:35:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG030419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 04 14:38:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.15	152	40355	20.00	ng	-0.02
21) Naphthalene-d8	10.97	136	164172	20.00	ng	-0.02
39) Acenaphthene-d10	14.77	164	115392	20.00	ng	-0.02
64) Phenanthrene-d10	17.52	188	278830	20.00	ng	-0.02
76) Chrysene-d12	21.80	240	298300	20.00	ng	-0.02
87) Perylene-d12	25.16	264	300384	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.69	112	161861	68.43	ng	-0.02
7) Phenol-d6	7.29	99	150967	42.80	ng	-0.02
23) Nitrobenzene-d5	9.33	82	309688	102.02	ng	-0.02
42) 2,4,6-Tribromophenol	16.26	330	205487	152.78	ng	-0.02
45) 2-Fluorobiphenyl	13.40	172	746731	92.14	ng	-0.02
79) Terphenyl-d14	20.11	244	1241786	91.66	ng	-0.01
Target Compounds						
17) 2-Methylphenol	8.57	107	6808	2.794	ng	# 89
20) 3+4-Methylphenols	8.90	107	7618	2.251	ng	# 82
22) Acetophenone	8.97	105	42214	9.580	ng	# 98
27) 2,4-Dimethylphenol	10.16	122	17749	7.422	ng	# 88
31) Naphthalene	11.02	128	142443	16.387	ng	# 99
32) Benzoic acid	10.24	122	2935m	8.140	ng	

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

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