

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031519\
 Data File : BG039953.D
 Acq On : 15 Mar 2019 22:48
 Operator : JU/SJ
 Sample : K1930-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-3

Quant Time: Mar 16 04:10:54 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	2066	20.00	ng	-0.02
21) Naphthalene-d8	0.00	136	0	0.00	ng	-10.99
39) Acenaphthene-d10	0.00	164	0	0.00	ng	-14.79
64) Phenanthrene-d10	0.00	188	0	0.00	ng	-17.53
76) Chrysene-d12	0.00	240	0	0.00	ng	-21.83
87) Perylene-d12	0.00	264	0	0.00	ng	-25.20
System Monitoring Compounds						
5) 2-Fluorophenol	5.69	112	321358	2653.63	ng	-0.02
7) Phenol-d6	7.29	99	472039	2614.17	ng	-0.02
23) Nitrobenzene-d5	9.33	82	289717	0.00	ng	-0.02
42) 2,4,6-Tribromophenol	16.26	330	172224	0.00	ng	-0.02
45) 2-Fluorobiphenyl	13.40	172	610454	0.00	ng	-0.02
79) Terphenyl-d14	20.11	244	879987	0.00	ng	-0.01
Target Compounds						
6) Aniline	7.31	93	629	2.659	ng	# 1
9) Benzaldehyde	7.29	77	1184	10.165	ng	# 4
10) Phenol	7.32	94	28028	143.282	ng	99
11) bis(2-Chloroethyl)ether	7.67	93	391	2.564	ng	# 1
15) Benzyl Alcohol	8.47	79	4332	30.244	ng	# 54
19) n-Nitroso-di-n-propylamine	9.33	70	38249	294.292	ng	# 68

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

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