

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121018\
 Data File : BG038454.D
 Acq On : 11 Dec 2018 00:53
 Operator : JU/SJ
 Sample : PB115480BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115480BL

Quant Time: Dec 11 03:53:01 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG120418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 04 17:15:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	756	20.00	ng	0.00
21) Naphthalene-d8	0.00	136	0	0.00	ng	-10.89
39) Acenaphthene-d10	0.00	164	0	0.00	ng	-14.71
64) Phenanthrene-d10	0.00	188	0	0.00	ng	-17.46
76) Chrysene-d12	0.00	240	0	0.00	ng	-21.74
87) Perylene-d12	0.00	264	0	0.00	ng	-25.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	182155	4263.88	ng	0.00
7) Phenol-d6	7.22	99	249283	4037.27	ng	0.00
23) Nitrobenzene-d5	9.25	82	157605	0.00	ng	0.00
42) 2,4,6-Tribromophenol	16.19	330	105777	0.00	ng	-0.01
45) 2-Fluorobiphenyl	13.32	172	389852	0.00	ng	0.00
79) Terphenyl-d14	20.06	244	728476	0.00	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.53	88	62	3.167	ng	# 1
4) n-Nitrosodimethylamine	3.85	42	141	5.495	ng	# 1
9) Benzaldehyde	7.21	77	361	9.233	ng	# 4
10) Phenol	7.60	94	276	4.231	ng	# 1
11) bis(2-Chloroethyl)ether	7.60	93	123	2.302	ng	# 1
15) Benzyl Alcohol	8.39	79	2482	48.894	ng	# 57
19) n-Nitroso-di-n-propylamine	9.25	70	21225	494.237	ng	# 76

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

