

Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\
 Data File : BG032328.D
 Acq On : 26 Jan 2018 13:31
 Operator : SJ/JU
 Sample : PB105983TB
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB105983TB

Quant Time: Jan 26 14:08:00 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 24 14:02:41 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.72	152	27967	20.00	ng	0.00
21) Naphthalene-d8	11.60	136	113992	20.00	ng	0.00
38) Acenaphthene-d10	15.36	164	86722	20.00	ng	0.00
63) Phenanthrene-d10	18.09	188	257870	20.00	ng	0.00
75) Chrysene-d12	22.41	240	333948	20.00	ng	-0.01
86) Perylene-d12	26.09	264	358016	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	6.15	112	187718	122.76	ng	0.00
7) Phenol-d6	7.83	99	234092	121.55	ng	0.00
23) Nitrobenzene-d5	9.91	82	142516	84.58	ng	0.00
41) 2,4,6-Tribromophenol	16.83	330	182413	127.11	ng	0.00
44) 2-Fluorobiphenyl	13.98	172	546207	78.10	ng	0.00
78) Terphenyl-d14	20.63	244	1246369	82.41	ng	0.00
Target Compounds						
15) Benzyl Alcohol	9.05	79	2810	2.008	ng	# 59
19) n-Nitroso-di-n-propylamine	9.91	70	19956	16.701	ng	# 75
50) 2,6-Dinitrotoluene	15.36	165	11197	7.108	ng	# 18
56) 2,4-Dinitrotoluene	15.36	165	11197	5.280	ng	# 16
64) 4,6-Dinitro-2-methylphenol	16.83	198	637	5.231	ng	# 7
66) 4-Bromophenyl-phenylether	16.83	248	15295	4.169	ng	# 1
76) Benizidine	20.63	184	17312	2.073	ng	98

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

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