

Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\
 Data File : BG032329.D
 Acq On : 26 Jan 2018 14:09
 Operator : SJ/JU
 Sample : PB106006TB
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB106006TB

Quant Time: Jan 26 14:44:00 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	26930	20.00	ng	0.00
21) Naphthalene-d8	11.60	136	107255	20.00	ng	0.00
38) Acenaphthene-d10	15.35	164	84566	20.00	ng	-0.01
63) Phenanthrene-d10	18.09	188	255199	20.00	ng	0.00
75) Chrysene-d12	22.41	240	330485	20.00	ng	-0.01
86) Perylene-d12	26.10	264	350632	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	6.15	112	164651	111.82	ng	0.00
7) Phenol-d6	7.83	99	204194	110.11	ng	0.00
23) Nitrobenzene-d5	9.92	82	123801	78.09	ng	0.00
41) 2,4,6-Tribromophenol	16.83	330	167662	119.81	ng	0.00
44) 2-Fluorobiphenyl	13.98	172	487526	71.48	ng	0.00
78) Terphenyl-d14	20.62	244	1158864	77.43	ng	-0.01
Target Compounds						
19) n-Nitroso-di-n-propylamine	9.92	70	17107	14.868	ng	# 76
32) Benzoic acid	10.83	122	61	5.302	ng	# 1
50) 2,6-Dinitrotoluene	15.36	165	10961	7.136	ng	# 18
56) 2,4-Dinitrotoluene	15.36	165	10961	5.300	ng	# 16
64) 4,6-Dinitro-2-methylphenol	16.56	198	63	4.919	ng	# 29
66) 4-Bromophenyl-phenylether	16.83	248	14276	3.932	ng	# 1
76) Benzidine	20.62	184	16535	2.001	ng	# 89

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

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