

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG041619\
 Data File : BG040521.D
 Acq On : 17 Apr 2019 3:35
 Operator : JU/SJ
 Sample : PB118938TB
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB118938TB

Quant Time: Apr 17 04:59:13 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG032719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 28 05:47:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	2448	20.00	ng	-0.03
21) Naphthalene-d8	0.00	136	0	0.00	ng	-10.88
39) Acenaphthene-d10	14.21	164	57	20.00	ng	-0.49
64) Phenanthrene-d10	0.00	188	0	0.00	ng	-17.45
76) Chrysene-d12	0.00	240	0	0.00	ng	-21.73
87) Perylene-d12	0.00	264	0	0.00	ng	-25.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	369867	2511.73	ng	-0.02
7) Phenol-d6	7.20	99	520062	2555.46	ng	-0.03
23) Nitrobenzene-d5	9.22	82	309467	0.00	ng	-0.03
42) 2,4,6-Tribromophenol	0.00	330	0	0.00	ng	
45) 2-Fluorobiphenyl	13.30	172	748129	194482.37	ng	-0.03
79) Terphenyl-d14	20.02	244	1308847	0.00	ng	-0.03
Target Compounds						
9) Benzaldehyde	7.20	77	974	6.977	ng	# 2
11) bis(2-Chloroethyl)ether	7.57	93	505	2.854	ng	# 1
15) Benzyl Alcohol	8.36	79	5404	32.972	ng	# 52
19) n-Nitroso-di-n-propylamine	9.22	70	38579	264.399	ng	# 74
46) 1,1'-Biphenyl	13.30	154	1367	303.525	ng	# 1
48) 2-Nitroaniline	13.29	65	1185	1016.919	ng	# 1
49) Acenaphthylene	13.52	152	318	54.291	ng	# 36
50) Dimethylphthalate	13.29	163	63	14.122	ng	# 76
57) 2,4-Dinitrotoluene	14.38	165	60	43.109	ng	# 24
60) Diethylphthalate	14.62	149	57	12.486	ng	# 57
63) Azobenzene	15.47	77	61	14.166	ng	# 48

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

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