

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122618\
 Data File : BF111750.D
 Acq On : 27 Dec 2018 8:33
 Operator : JU/SJ
 Sample : J6492-01DL 5X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 QNWP8-MW26S-GWDL

Quant Time: Dec 27 12:28:08 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 27 11:05:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	110693	20.00	ng	0.00
21) Naphthalene-d8	8.20	136	373366	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	189539	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	379499	20.00	ng	0.00
76) Chrysene-d12	14.06	240	278568	20.00	ng	0.00
87) Perylene-d12	15.55	264	221358	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	56629	8.72	ng	0.00
7) Phenol-d6	6.54	99	44388	5.37	ng	0.00
23) Nitrobenzene-d5	7.46	82	96949	15.11	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	43102	19.25	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	190578	14.66	ng	0.00
79) Terphenyl-d14	13.00	244	179131	12.19	ng	0.00
Target Compounds						
6) Aniline	6.56	93	34728	3.296	ng	# 55
10) Phenol	6.56	94	47764	5.122	ng	# 70
17) 2-Methylphenol	7.16	107	39269	6.817	ng	# 94
20) 3+4-Methylphenols	7.31	107	74747	10.270	ng	# 65
27) 2,4-Dimethylphenol	7.85	122	239341	44.530	ng	# 93
31) Naphthalene	8.24	128	9915619	552.720	ng	# 92
37) 2-Methylnaphthalene	8.90	142	458277	39.264	ng	# 97
38) 1-Methylnaphthalene	9.00	142	470521	41.975	ng	# 99
46) 1,1'-Biphenyl	9.36	154	130895	8.181	ng	# 94
52) Acenaphthene	9.97	154	187835	16.455	ng	# 98
55) Dibenzofuran	10.15	168	255519	14.816	ng	# 92
56) 4-Nitrophenol	10.07	139	7232	3.215	ng	# 58
58) Fluorene	10.49	166	147218	11.846	ng	# 99
71) Phenanthrene	11.45	178	237453	11.798	ng	# 93
73) Carbazole	11.66	167	208007	10.606	ng	# 95

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

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