

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF111218\
 Data File : BF110698.D
 Acq On : 13 Nov 2018 3:05
 Operator : JU/SJ
 Sample : J5904-09
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DFTPP

Quant Time: Nov 13 07:17:40 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 09 15:13:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	88848	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	359363	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	150619	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	238529	20.00	ng	0.00
76) Chrysene-d12	14.13	240	166827	20.00	ng	-0.01
87) Perylene-d12	15.66	264	121255	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	231626	42.35	ng	0.00
7) Phenol-d6	6.56	99	175339	25.07	ng	-0.01
23) Nitrobenzene-d5	7.51	82	615750	98.68	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	103202	68.00	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	1080680	102.47	ng	0.00
79) Terphenyl-d14	13.06	244	702282	78.84	ng	0.00
Target Compounds						
19) n-Nitroso-di-n-propylamine	7.51	70	80463	18.021	ng	# 82
25) Isophorone	7.51	82	615746	60.205	ng	# 67
51) 2,6-Dinitrotoluene	9.99	165	19259	7.790	ng	# 21
57) 2,4-Dinitrotoluene	9.99	165	19259	5.991	ng	# 17
67) 4-Bromophenyl-phenylether	10.78	248	6138	2.328	ng	# 1

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

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