

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122618\
 Data File : BF111734.D
 Acq On : 27 Dec 2018 00:40
 Operator : JU/SJ
 Sample : J6388-11
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SS010-0612-01

Quant Time: Dec 27 04:00:28 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	152860	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	513478	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	232329	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	430512	20.00	ng	0.01
76) Chrysene-d12	14.07	240	324243	20.00	ng	0.02
87) Perylene-d12	15.56	264	249843	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	864613	96.41	ng	0.02
7) Phenol-d6	6.55	99	1022445	89.59	ng	0.02
23) Nitrobenzene-d5	7.46	82	664673	75.35	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	275356	100.31	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	1157154	72.60	ng	0.00
79) Terphenyl-d14	13.02	244	1123187	65.69	ng	0.02
Target Compounds						
10) Phenol	6.56	94	136511	10.601	ng	Qvalue 97
20) 3+4-Methylphenols	7.31	107	21978	2.187	ng	# 66
32) Benzoic acid	7.90	122	13190	2.699	ng	95
50) Dimethylphthalate	9.65	163	142647	8.396	ng	100
68) Hexachlorobenzene	11.04	284	23834	4.003	ng	92
69) Atrazine	11.13	200	59890	11.929	ng	# 39

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

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