

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122618\
 Data File : BF111735.D
 Acq On : 27 Dec 2018 1:07
 Operator : JU/SJ
 Sample : J6388-12
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SS010-1218-01

Quant Time: Dec 27 04:05:22 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	149963	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	494077	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	215756	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	384672	20.00	ng	0.01
76) Chrysene-d12	14.07	240	286331	20.00	ng	0.02
87) Perylene-d12	15.57	264	217086	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	889560	101.11	ng	0.03
7) Phenol-d6	6.55	99	1066100	95.22	ng	0.02
23) Nitrobenzene-d5	7.46	82	659236	77.67	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	268378	105.27	ng	0.01
45) 2-Fluorobiphenyl	9.26	172	1167637	78.88	ng	0.00
79) Terphenyl-d14	13.02	244	1083152	71.74	ng	0.02
Target Compounds						
10) Phenol	6.56	94	408208	32.312	ng	99
17) 2-Methylphenol	7.16	107	29154	3.736	ng	# 94
20) 3+4-Methylphenols	7.31	107	129925	13.176	ng	# 62
27) 2,4-Dimethylphenol	7.84	122	49464	6.955	ng	93
30) 1,2,4-Trichlorobenzene	8.13	180	23051	2.704	ng	96
32) Benzoic acid	7.92	122	33237	7.068	ng	95
50) Dimethylphthalate	9.65	163	143665	9.106	ng	99
68) Hexachlorobenzene	11.04	284	17811	3.348	ng	94
69) Atrazine	11.13	200	14775	3.294	ng	# 52

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

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