

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122618\
 Data File : BF111736.D
 Acq On : 27 Dec 2018 1:35
 Operator : JU/SJ
 Sample : J6388-13
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SS010-1824-01

Quant Time: Dec 27 04:11:00 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	139753	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	474770	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	214070	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	391319	20.00	ng	0.01
76) Chrysene-d12	14.07	240	300969	20.00	ng	0.02
87) Perylene-d12	15.56	264	229125	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	687122	83.81	ng	0.02
7) Phenol-d6	6.55	99	843982	80.89	ng	0.02
23) Nitrobenzene-d5	7.46	82	516229	63.29	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	225393	89.11	ng	0.01
45) 2-Fluorobiphenyl	9.26	172	963051	65.57	ng	0.00
79) Terphenyl-d14	13.02	244	937166	59.05	ng	0.02
Target Compounds						
10) Phenol	6.56	94	200976	17.071	ng	97
20) 3+4-Methylphenols	7.31	107	53218	5.791	ng	# 59
27) 2,4-Dimethylphenol	7.84	122	17271	2.527	ng	93
32) Benzoic acid	7.90	122	11383	2.519	ng	87
50) Dimethylphthalate	9.65	163	130880	8.361	ng	# 98
68) Hexachlorobenzene	11.04	284	11315	2.091	ng	100

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

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