

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121018\
 Data File : BF111256.D
 Acq On : 10 Dec 2018 21:41
 Operator : JU/SJ
 Sample : J6237-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SAMPLE-2

Quant Time: Dec 11 05:59:56 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 10 01:15:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	597786	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	2222901	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	846920	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	1211730	20.00	ng	0.00
76) Chrysene-d12	13.96	240	634873	20.00	ng	0.00
87) Perylene-d12	15.40	264	526779	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	3516755	92.90	ng	0.02
7) Phenol-d6	6.44	99	4420952	93.97	ng	0.00
23) Nitrobenzene-d5	7.36	82	2830522	71.42	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	509374	67.90	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	3896184	77.99	ng	0.00
79) Terphenyl-d14	12.91	244	1850372	69.91	ng	0.00
Target Compounds						
10) Phenol	6.45	94	304562	5.536	ng	97
50) Dimethylphthalate	9.55	163	356659	5.707	ng	# 96
52) Acenaphthene	9.87	154	193563	3.834	ng	99
58) Fluorene	10.39	166	135343	2.546	ng	98
71) Phenanthrene	11.35	178	373255	6.107	ng	95
75) Fluoranthene	12.53	202	127482	2.283	ng	98
78) Pyrene	12.76	202	103238	2.276	ng	94

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

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