

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111117\
 Data File : BF100467.D
 Acq On : 12 Nov 2017 4:23
 Operator : SJ/JU
 Sample : I6408-02
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-A19R(17.5-19.5)

Quant Time: Nov 13 05:43:51 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 10 15:44:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	133291	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	492091	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	201101	20.00	ng	0.00
63) Phenanthrene-d10	11.43	188	314206	20.00	ng	0.00
75) Chrysene-d12	14.06	240	283528	20.00	ng	0.00
86) Perylene-d12	15.54	264	199543	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	971653	116.85	ng	0.01
7) Phenol-d6	6.53	99	1199286	116.96	ng	0.00
23) Nitrobenzene-d5	7.46	82	716477	96.26	ng	0.00
41) 2,4,6-Tribromophenol	10.73	330	221978	108.32	ng	0.00
44) 2-Fluorobiphenyl	9.26	172	1053924	79.80	ng	0.00
78) Terphenyl-d14	13.01	244	804931	65.23	ng	0.00
Target Compounds						
10) Phenol	6.54	94	59551	5.532	ng	94
37) 2-Methylnaphthalene	8.90	142	35471	2.254	ng	94
49) Dimethylphthalate	9.65	163	85020	5.537	ng	# 94

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

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