

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111716\
 Data File : BF091216.D
 Acq On : 17 Nov 2016 17:03
 Operator : UM/SJ
 Sample : H5686-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-1

Quant Time: Nov 18 03:38:54 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 08 12:59:29 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.61	152	179731	20.00	ng	-0.05
21) Naphthalene-d8	7.90	136	759756	20.00	ng	-0.05
38) Acenaphthene-d10	9.66	164	421203	20.00	ng	-0.03
63) Phenanthrene-d10	11.14	188	676289	20.00	ng	-0.03
75) Chrysene-d12	13.77	240	475141	20.00	ng	-0.05
86) Perylene-d12	15.14	264	343689	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	731841	67.25	ng	-0.05
7) Phenol-d6	6.27	99	635292	43.01	ng	-0.05
23) Nitrobenzene-d5	7.20	82	1194678	85.78	ng	-0.05
41) 2,4,6-Tribromophenol	10.45	330	487315	127.97	ng	-0.03
44) 2-Fluorobiphenyl	8.99	172	1975904	80.76	ng	-0.05
78) Terphenyl-d14	12.73	244	1809776	95.05	ng	-0.03
Target Compounds						Qvalue
37) 2-Methylnaphthalene	8.62	142	632320	27.07	ng	99
51) Acenaphthene	9.69	154	154778	6.77	ng	99
54) Dibenzofuran	9.87	168	167422	5.44	ng	# 88
57) Fluorene	10.20	166	263501	10.48	ng	# 96
70) Phenanthrene	11.16	178	307466	8.55	ng	99

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

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