

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101216\
 Data File : BF090536.D
 Acq On : 12 Oct 2016 15:37
 Operator : UM/SJ
 Sample : H5153-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AOU2-MW3T-100716

Quant Time: Oct 12 17:46:53 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 11 16:14:43 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	213606	20.00	ng	-0.01
21) Naphthalene-d8	8.23	136	907590	20.00	ng	-0.01
38) Acenaphthene-d10	9.99	164	472324	20.00	ng	-0.01
63) Phenanthrene-d10	11.48	188	788779	20.00	ng	0.00
75) Chrysene-d12	14.12	240	651379	20.00	ng	-0.01
86) Perylene-d12	15.60	264	445781	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	485367	35.35	ng	0.02
7) Phenol-d6	6.59	99	465458	26.86	ng	0.00
23) Nitrobenzene-d5	7.51	82	1263188	75.49	ng	0.00
41) 2,4,6-Tribromophenol	10.78	330	370920	79.22	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	2360064	83.97	ng	-0.01
78) Terphenyl-d14	13.07	244	2201672	82.60	ng	-0.01
Target Compounds						Qvalue
6) Aniline	6.61	93	104829	4.77	ng	90
22) Acetophenone	7.35	105	57529	2.88	ng	# 93
31) Naphthalene	8.26	128	909787	20.12	ng	97
37) 2-Methylnaphthalene	8.95	142	69645	2.29	ng	97
54) Dibenzofuran	10.20	168	409031	10.59	ng	96
57) Fluorene	10.54	166	1033050	32.77	ng	99
70) Phenanthrene	11.50	178	606382	13.95	ng	96

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

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