

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101216\
 Data File : BF090535.D
 Acq On : 12 Oct 2016 15:10
 Operator : UM/SJ
 Sample : H5153-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AOU2-MW3-100716

Quant Time: Oct 12 16:46:24 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	223730	20.00	ng	-0.01
21) Naphthalene-d8	8.23	136	956050	20.00	ng	-0.01
38) Acenaphthene-d10	9.99	164	505911	20.00	ng	-0.01
63) Phenanthrene-d10	11.48	188	815768	20.00	ng	0.00
75) Chrysene-d12	14.12	240	676639	20.00	ng	-0.01
86) Perylene-d12	15.60	264	503591	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	480894	33.44	ng	0.02
7) Phenol-d6	6.59	99	449144	24.75	ng	0.00
23) Nitrobenzene-d5	7.51	82	1422569	80.70	ng	0.00
41) 2,4,6-Tribromophenol	10.78	330	394060	78.58	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	2487789	82.64	ng	-0.01
78) Terphenyl-d14	13.07	244	2279409	82.33	ng	-0.01
Target Compounds						
6) Aniline	6.61	93	110853	4.82	ng	Qvalue 96
20) 3+4-Methylphenols	7.37	107	150295	9.54	ng	# 84
22) Acetophenone	7.35	105	60833	2.89	ng	# 90
31) Naphthalene	8.26	128	661392	13.89	ng	97
54) Dibenzofuran	10.20	168	363819	8.80	ng	97
57) Fluorene	10.54	166	928395	27.49	ng	100
70) Phenanthrene	11.50	178	501389	11.15	ng	96

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

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