

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011216\
 Data File : BF084428.D
 Acq On : 13 Jan 2016 00:06
 Operator : SJ/UM
 Sample : G4988-09
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RX2

Quant Time: Jan 13 02:28:40 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 04 12:22:40 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	264833	20.00	ng	-0.06
21) Naphthalene-d8	8.13	136	986165	20.00	ng	-0.06
38) Acenaphthene-d10	9.88	164	473077	20.00	ng	-0.07
63) Phenanthrene-d10	11.37	188	836782	20.00	ng	-0.06
75) Chrysene-d12	14.00	240	534223	20.00	ng	-0.07
86) Perylene-d12	15.43	264	459830	20.00	ng	-0.08
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	895975	54.72	ng	-0.04
7) Phenol-d6	6.48	99	667838	32.48	ng	-0.06
23) Nitrobenzene-d5	7.41	82	1801029	101.64	ng	-0.06
41) 2,4,6-Tribromophenol	10.67	330	510234	106.83	ng	-0.07
44) 2-Fluorobiphenyl	9.21	172	2433136	90.92	ng	-0.06
78) Terphenyl-d14	12.96	244	2043348	85.38	ng	-0.06
Target Compounds						Qvalue
31) Naphthalene	8.17	128	8604542	192.31	ng	# 84
37) 2-Methylnaphthalene	8.84	142	2042696	63.60	ng	98
45) 1,1'-Biphenyl	9.30	154	371070	9.87	ng	96
48) Acenaphthylene	9.74	152	199667	4.58	ng	# 94
51) Acenaphthene	9.92	154	1509802	51.59	ng	98
54) Dibenzofuran	10.09	168	1506876	39.38	ng	93
57) Fluorene	10.43	166	1147485	38.82	ng	99
70) Phenanthrene	11.39	178	1616304	36.93	ng	97
71) Anthracene	11.44	178	106871	2.49	ng	98
72) Carbazole	11.60	167	245340	5.85	ng	98
74) Fluoranthene	12.58	202	149680	3.27	ng	94

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

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