

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032516\
 Data File : BF085764.D
 Acq On : 25 Mar 2016 16:25
 Operator : UM/SJ
 Sample : H1858-12
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

Quant Time: Mar 25 16:54:24 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 24 14:15:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	OIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	107628	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	387209	20.00	ng	-0.01
38) Acenaphthene-d10	9.86	164	157218	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	243708	20.00	ng	-0.01
75) Chrysene-d12	13.98	240	200101	20.00	ng	0.00
86) Perylene-d12	15.40	264	175771	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.43	112	721205	111.60	ng	0.02
7) Phenol-d6	6.46	99	922125	112.23	ng	0.00
23) Nitrobenzene-d5	7.40	82	715128	104.00	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	203521	136.25	ng	0.00
44) 2-Fluorobiphenyl	9.19	172	1119547	118.97	ng	0.00
78) Terphenyl-d14	12.93	244	724751	74.49	ng	0.00

Target Compounds

						Qvalue
9) Benzaldehyde	6.60	77	18791	3.80	ng	83
15) Benzyl Alcohol	6.98	79	11222	2.05	ng	# 42
19) n-Nitroso-di-n-propylamine	7.40	70	94828	19.27	ng	# 74
25) Isophorone	7.40	82	698566	52.74	ng	# 68
31) Naphthalene	8.13	128	322886	16.55	ng	99
33) 4-Chloroaniline	8.13	127	41098	5.16	ng	# 32
50) 2,6-Dinitrotoluene	9.86	165	21093	8.22	ng	# 19
55) 4-Nitrophenol	10.07	139	4371	2.28	ng	# 10
56) 2,4-Dinitrotoluene	9.86	165	21093	6.46	ng	# 37
66) 4-Bromophenyl-phenylether	10.65	248	13526	5.42	ng	# 1

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

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