

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101116\
 Data File : BF090521.D
 Acq On : 12 Oct 2016 00:02
 Operator : UM/SJ
 Sample : H5146-09
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S8-0615

Quant Time: Oct 12 02:00:13 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	6.95	152	217362	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	845849	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	494297	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	726386	20.00	ng	0.01
75) Chrysene-d12	14.13	240	471713	20.00	ng	0.00
86) Perylene-d12	15.62	264	387814	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	658064	47.10	ng	0.01
7) Phenol-d6	6.60	99	778538	44.16	ng	0.01
23) Nitrobenzene-d5	7.53	82	1259163	80.74	ng	0.01
41) 2,4,6-Tribromophenol	10.81	330	686001	140.00	ng	0.01
44) 2-Fluorobiphenyl	9.33	172	2345029	79.73	ng	0.01
78) Terphenyl-d14	13.08	244	1776100	92.01	ng	0.00
Target Compounds						
20) 3+4-Methylphenols	7.39	107	71119	4.65	ng	# 82
22) Acetophenone	7.37	105	388122	20.84	ng	# 86
27) 2,4-Dimethylphenol	7.91	122	210130	16.20	ng	90
31) Naphthalene	8.27	128	2136423	50.70	ng	99
32) Benzoic acid	8.15	122	381202	37.23	ng	99
37) 2-Methylnaphthalene	8.97	142	485201	17.15	ng	99

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

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