

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091115\
 Data File : BF081590.D
 Acq On : 11 Sep 2015 18:42
 Operator : UM/IZ
 Sample : G3511-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-01(0-0.16)

Quant Time: Sep 14 05:50:53 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 11 23:14:51 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.25	152	53782	20.00	ng	0.00
21) Naphthalene-d8	11.14	136	217342	20.00	ng	-0.01
38) Acenaphthene-d10	15.28	164	100197	20.00	ng	-0.01
63) Phenanthrene-d10	18.79	188	187441	20.00	ng	0.00
75) Chrysene-d12	23.40	240	113503	20.00	ng	-0.01
86) Perylene-d12	24.84	264	104227	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	365893	105.55	ng	0.02
7) Phenol-d6	7.67	99	511840	111.55	ng	-0.01
23) Nitrobenzene-d5	9.54	82	337571	74.94	ng	-0.01
41) 2,4,6-Tribromophenol	17.19	330	94201	105.59	ng	-0.01
44) 2-Fluorobiphenyl	13.80	172	554153	70.87	ng	-0.01
78) Terphenyl-d14	22.13	244	408022	83.68	ng	0.00
Target Compounds						Qvalue
48) Acenaphthylene	14.93	152	40615	3.44	ng	95
49) Dimethylphthalate	14.85	163	140598	18.06	ng	# 96
70) Phenanthrene	18.84	178	151055	14.50	ng	97
71) Anthracene	18.96	178	31868	2.98	ng	99
74) Fluoranthene	21.45	202	322510	28.59	ng	88
77) Pyrene	21.80	202	274431	32.64	ng	97
80) Benzo(a)anthracene	23.38	228	134207	18.57	ng	# 88
82) Chrysene	23.43	228	118430	18.28	ng	92
83) Bis(2-ethylhexyl)phthalate	23.51	149	39852	8.26	ng	# 92
85) Indeno(1,2,3-cd)pyrene	25.89	276	78584m	16.53	ng	
87) Benzo(b)fluoranthene	24.49	252	207201m	27.85	ng	
88) Benzo(k)fluoranthene	24.52	252	15955m	2.58	ng	
89) Benzo(a)pyrene	24.79	252	115352	18.29	ng	# 81
90) Dibenzo(a,h)anthracene	25.90	278	20484m	4.20	ng	
91) Benzo(g,h,i)perylene	26.20	276	73136	15.73	ng	# 73

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

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