

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091115\
 Data File : BF081594.D
 Acq On : 11 Sep 2015 21:10
 Operator : UM/IZ
 Sample : G3511-15
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-09(0-0.16)

Quant Time: Sep 14 07:00:33 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	55284	20.00	ng	0.00
21) Naphthalene-d8	11.14	136	214040	20.00	ng	-0.01
38) Acenaphthene-d10	15.28	164	98158	20.00	ng	-0.01
63) Phenanthrene-d10	18.79	188	173571	20.00	ng	0.00
75) Chrysene-d12	23.40	240	100556	20.00	ng	-0.01
86) Perylene-d12	24.84	264	97457	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	469604	131.79	ng	0.03
7) Phenol-d6	7.68	99	643105	136.35	ng	0.00
23) Nitrobenzene-d5	9.55	82	429886	96.90	ng	0.00
41) 2,4,6-Tribromophenol	17.19	330	118598	135.70	ng	-0.01
44) 2-Fluorobiphenyl	13.80	172	735431	96.01	ng	-0.01
78) Terphenyl-d14	22.13	244	467724	108.28	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	14.85	163	143700	18.84	ng	# 96
70) Phenanthrene	18.84	178	85988	8.91	ng	97
71) Anthracene	18.96	178	21664	2.19	ng	96
74) Fluoranthene	21.45	202	159241	15.24	ng	86
77) Pyrene	21.80	202	150045	20.14	ng	97
80) Benzo(a)anthracene	23.38	228	87814	13.71	ng	94
82) Chrysene	23.43	228	76876	13.39	ng	94
83) Bis(2-ethylhexyl)phthalate	23.51	149	10438	2.44	ng	# 88
85) Indeno(1,2,3-cd)pyrene	25.89	276	52659	12.50	ng	# 83
87) Benzo(b)fluoranthene	24.49	252	118528m	17.04	ng	
88) Benzo(k)fluoranthene	24.50	252	44581m	7.72	ng	
89) Benzo(a)pyrene	24.79	252	79840	13.54	ng	# 82
90) Dibenzo(a,h)anthracene	25.89	278	13802m	3.03	ng	
91) Benzo(g,h,i)perylene	26.20	276	46113	10.60	ng	# 71

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

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