

Data Path : V:\HPCHEM1\BNA F\DATA\BF121515\
 Data File : BF083840.D
 Acq On : 16 Dec 2015 3:42
 Operator : UM/IZ
 Sample : G4739-07DL 5X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 K1206-0-2DL

Quant Time: Dec 16 07:33:09 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 14 01:30:12 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	90152	20.00	ng	-0.02
21) Naphthalene-d8	8.19	136	348444	20.00	ng	-0.02
38) Acenaphthene-d10	9.94	164	150459	20.00	ng	-0.02
63) Phenanthrene-d10	11.42	188	249697	20.00	ng	-0.02
75) Chrysene-d12	14.05	240	197645	20.00	ng	-0.02
86) Perylene-d12	15.49	264	151736	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	122555	21.93	ng	-0.02
7) Phenol-d6	6.52	99	146384	20.69	ng	-0.02
23) Nitrobenzene-d5	7.46	82	90543	14.54	ng	-0.02
41) 2,4,6-Tribromophenol	10.73	330	26901	17.55	ng	-0.02
44) 2-Fluorobiphenyl	9.26	172	160543	13.11	ng	-0.02
78) Terphenyl-d14	13.00	244	109234	11.89	ng	-0.02
Target Compounds						Qvalue
48) Acenaphthylene	9.80	152	121735	7.35	ng	99
49) Dimethylphthalate	9.65	163	24407	2.08	ng	99
70) Phenanthrene	11.45	178	287241	21.65	ng	99
71) Anthracene	11.49	178	66312	4.81	ng	98
74) Fluoranthene	12.64	202	422899	30.85	ng	98
77) Pyrene	12.87	202	647050	41.42	ng	99
80) Benzo(a)anthracene	14.04	228	228060m	20.29	ng	
82) Chrysene	14.08	228	190284	17.36	ng	99
85) Indeno(1,2,3-cd)pyrene	16.92	276	62512	5.81	ng	# 100
87) Benzo(b)fluoranthene	15.08	252	156335m	16.15	ng	
88) Benzo(k)fluoranthene	15.09	252	66399m	7.57	ng	
89) Benzo(a)pyrene	15.44	252	75849	8.78	ng	99
91) Benzo(a,h,i)perylene	17.36	276	63360	7.18	ng	98

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

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