

Data Path : V:\HPCHEM1\BNA F\DATA\BF121515\
 Data File : BF083839.D
 Acq On : 16 Dec 2015 3:14
 Operator : UM/IZ
 Sample : G4739-04DL 5X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 K206-0-2DL

Quant Time: Dec 16 07:27:04 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	99303	20.00	ng	-0.02
21) Naphthalene-d8	8.19	136	410461	20.00	ng	-0.02
38) Acenaphthene-d10	9.94	164	177645	20.00	ng	-0.02
63) Phenanthrene-d10	11.42	188	294235	20.00	ng	-0.02
75) Chrysene-d12	14.05	240	209332	20.00	ng	-0.02
86) Perylene-d12	15.49	264	170576	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	152094	24.71	ng	-0.02
7) Phenol-d6	6.52	99	177705	22.80	ng	-0.02
23) Nitrobenzene-d5	7.46	82	116965	15.94	ng	-0.02
41) 2,4,6-Tribromophenol	10.73	330	38021	21.00	ng	-0.02
44) 2-Fluorobiphenyl	9.26	172	198743	13.72	ng	-0.02
78) Terphenyl-d14	13.00	244	134023	13.77	ng	-0.02
Target Compounds						Qvalue
48) Acenaphthylene	9.80	152	152229	7.79	ng	99
49) Dimethylphthalate	9.65	163	52240	3.77	ng	99
70) Phenanthrene	11.45	178	306471	19.61	ng	99
71) Anthracene	11.49	178	79261	4.88	ng	97
74) Fluoranthene	12.64	202	493004	30.52	ng	98
77) Pyrene	12.86	202	721872	43.63	ng	98
80) Benzo(a)anthracene	14.04	228	279996m	23.52	ng	
82) Chrysene	14.08	228	234558	20.20	ng	99
85) Indeno(1,2,3-cd)pyrene	16.92	276	77081	6.76	ng	# 100
87) Benzo(b)fluoranthene	15.08	252	196537m	18.06	ng	
88) Benzo(k)fluoranthene	15.09	252	83560m	8.47	ng	
89) Benzo(a)pyrene	15.44	252	111276	11.46	ng	98
90) Dibenzo(a,h)anthracene	16.93	278	21583m	2.24	ng	
91) Benzo(a,h,i)perylene	17.36	276	80410	8.10	ng	99

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

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