

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120315\
 Data File : BF083476.D
 Acq On : 3 Dec 2015 23:22
 Operator : UM/IZ
 Sample : G4608-05
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CWC-MISC-7-20151130

Quant Time: Dec 04 01:37:34 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 01 01:17:16 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.53	152	169637m	20.00	ng	-0.01
21) Naphthalene-d8	7.81	136	634339	20.00	ng	-0.02
38) Acenaphthene-d10	9.56	164	304504	20.00	ng	-0.02
63) Phenanthrene-d10	11.09	188	467962	20.00	ng	-0.03
75) Chrysene-d12	14.72	240	376001	20.00	ng	-0.05
86) Perylene-d12	16.79	264	300904	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	1036441	91.68	ng	-0.01
7) Phenol-d6	6.20	99	1346752	87.04	ng	-0.02
23) Nitrobenzene-d5	7.10	82	979331	67.87	ng	-0.02
41) 2,4,6-Tribromophenol	10.35	330	227621	74.46	ng	-0.03
44) 2-Fluorobiphenyl	8.90	172	1269324	59.40	ng	-0.02
78) Terphenyl-d14	13.19	244	901515	57.18	ng	-0.03
Target Compounds						Qvalue
48) Acenaphthylene	9.42	152	76704	2.36	ng	97
49) Dimethylphthalate	9.30	163	343229	13.87	ng	99
70) Phenanthrene	11.13	178	933909	33.72	ng	99
71) Anthracene	11.18	178	114410	4.10	ng	97
72) Carbazole	11.38	167	59906	2.39	ng	98
74) Fluoranthene	12.63	202	1069182	37.25	ng	97
77) Pyrene	12.93	202	945935	36.02	ng	98
80) Benzo(a)anthracene	14.69	228	375185	16.29	ng	97
82) Chrysene	14.75	228	412791	18.55	ng	98
85) Indeno(1,2,3-cd)pyrene	18.16	276	142500	8.79	ng	95
87) Benzo(b)fluoranthene	16.27	252	439606m	22.77	ng	
88) Benzo(k)fluoranthene	16.31	252	93838m	5.15	ng	
89) Benzo(a)pyrene	16.71	252	257771	15.54	ng	# 96
90) Dibenzo(a,h)anthracene	18.18	278	39552m	2.74	ng	
91) Benzo(a,h,i)perylene	18.52	276	132226	9.32	ng	# 94

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

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