

Data Path : V:\HPCHEM1\BNA F\DATA\BF112015\
 Data File : BF083084.D
 Acq On : 20 Nov 2015 16:23
 Operator : UM/IZ
 Sample : G4436-01DL 10X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PM-1350DL

Quant Time: Nov 20 17:47:50 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF111915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 20 01:33:39 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	143782	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	535955	20.00	ng	-0.01
38) Acenaphthene-d10	9.74	164	265587	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	442134	20.00	ng	0.00
75) Chrysene-d12	13.85	240	371953	20.00	ng	-0.01
86) Perylene-d12	15.23	264	326756	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	114986	12.42	ng	0.00
7) Phenol-d6	6.34	99	155214	12.73	ng	-0.01
23) Nitrobenzene-d5	7.26	82	83065	7.10	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	17132	6.08	ng	-0.01
44) 2-Fluorobiphenyl	9.07	172	181625	9.21	ng	-0.01
78) Terphenyl-d14	12.81	244	130821	8.30	ng	0.00
Target Compounds						Qvalue
48) Acenaphthylene	9.60	152	60256	2.06	ng	98
70) Phenanthrene	11.24	178	537278	20.14	ng	99
71) Anthracene	11.29	178	123281	4.59	ng	98
72) Carbazole	11.45	167	47548	2.02	ng	99
74) Fluoranthene	12.43	202	899252	32.35	ng	97
77) Pyrene	12.66	202	829658	31.68	ng	98
80) Benzo(a)anthracene	13.85	228	449302	19.01	ng	96
82) Chrysene	13.88	228	384797m	17.75	ng	
85) Indeno(1,2,3-cd)pyrene	16.53	276	138342	7.84	ng	96
87) Benzo(b)fluoranthene	14.85	252	415271m	18.18	ng	
88) Benzo(k)fluoranthene	14.87	252	160793m	9.24	ng	
89) Benzo(a)pyrene	15.17	252	275127m	15.01	ng	
90) Dibenzo(a,h)anthracene	16.53	278	35143m	2.20	ng	
91) Benzo(a,h,i)perylene	16.91	276	121007	7.77	ng	# 93

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

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