

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112215\
 Data File : BF083177.D
 Acq On : 23 Nov 2015 00:06
 Operator : UM/IZ
 Sample : G4496-01
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 A17-COMP

Quant Time: Nov 23 06:02:25 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF112115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Nov 22 09:22:46 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.65	152	217881	20.00	ng	-0.01
21) Naphthalene-d8	7.94	136	781364	20.00	ng	-0.01
38) Acenaphthene-d10	9.68	164	367178	20.00	ng	-0.02
63) Phenanthrene-d10	11.16	188	652081	20.00	ng	-0.01
75) Chrysene-d12	13.78	240	491609	20.00	ng	-0.02
86) Perylene-d12	15.15	264	443029	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	1991712	138.20	ng	-0.01
7) Phenol-d6	6.32	99	2467265	132.32	ng	-0.02
23) Nitrobenzene-d5	7.22	82	1585866	92.74	ng	-0.01
41) 2,4,6-Tribromophenol	10.48	330	516461	137.49	ng	-0.01
44) 2-Fluorobiphenyl	9.02	172	2237584	84.98	ng	-0.01
78) Terphenyl-d14	12.75	244	1822131	87.28	ng	-0.01
Target Compounds						Qvalue
10) Phenol	6.33	94	56686	2.51	ng	88
49) Dimethylphthalate	9.42	163	654581	23.16	ng	99
74) Fluoranthene	12.37	202	136947	3.56	ng	97
77) Pyrene	12.59	202	114757	3.42	ng	97
80) Benzo(a)anthracene	13.77	228	73111	2.40	ng	98
82) Chrysene	13.81	228	56834	2.02	ng	96
87) Benzo(b)fluoranthene	14.78	252	79890m	2.64	ng	
89) Benzo(a)pyrene	15.10	252	52641	2.11	ng	# 92

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

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