

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112315\
 Data File : BF083215.D
 Acq On : 23 Nov 2015 20:39
 Operator : UM/IZ
 Sample : G4443-04DL 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20151110MGP206BRV40DL

Quant Time: Nov 24 06:17:21 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	6.64	152	195732	20.00	ng	-0.02
21) Naphthalene-d8	7.94	136	652858	20.00	ng	-0.01
38) Acenaphthene-d10	9.68	164	372943	20.00	ng	-0.02
63) Phenanthrene-d10	11.15	188	700730	20.00	ng	-0.02
75) Chrysene-d12	13.77	240	573017	20.00	ng	-0.03
86) Perylene-d12	15.13	264	430256	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.24	112	113957	8.80	ng	-0.03
7) Phenol-d6	6.31	99	104447	6.24	ng	-0.03
23) Nitrobenzene-d5	7.21	82	285744	20.00	ng	-0.02
41) 2,4,6-Tribromophenol	10.47	330	49311	12.92	ng	-0.02
44) 2-Fluorobiphenyl	9.00	172	477175	17.84	ng	-0.02
78) Terphenyl-d14	12.73	244	352969	14.51	ng	-0.03
Target Compounds						
31) Naphthalene	7.99	128	14899183	422.86	ng	Qvalue # 80
37) 2-Methylnaphthalene	8.65	142	2241272	98.01	ng	98
45) 1,1'-Biphenyl	9.11	154	288311	9.08	ng	99
48) Acenaphthylene	9.54	152	1363920	35.30	ng	99
51) Acenaphthene	9.70	154	368065	15.65	ng	99
54) Dibenzofuran	9.88	168	77446	2.33	ng	# 93
57) Fluorene	10.22	166	447151	16.27	ng	100
70) Phenanthrene	11.18	178	633763	15.70	ng	99
71) Anthracene	11.22	178	127202	3.09	ng	97
72) Carbazole	11.39	167	154085	4.21	ng	97

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

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