

Data Path : Z:\HPCHEM1\BNA F\DATA\BF082015\
 Data File : BF081133.D
 Acq On : 20 Aug 2015 19:39
 Operator : UM/IZ
 Sample : G3238-01DL 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-A(2)DL

Quant Time: Aug 21 02:37:54 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 20 11:22:13 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.57	152	64105	20.00	ng	-0.01
21) Naphthalene-d8	7.86	136	259371	20.00	ng	-0.01
38) Acenaphthene-d10	9.61	164	117890	20.00	ng	-0.01
63) Phenanthrene-d10	11.08	188	227195	20.00	ng	-0.01
75) Chrysene-d12	13.69	240	201932	20.00	ng	-0.01
86) Perylene-d12	15.03	264	125746	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.13	112	125618	29.47	ng	-0.01
7) Phenol-d6	6.22	99	109135	19.63	ng	-0.01
23) Nitrobenzene-d5	7.14	82	273010	48.08	ng	-0.02
41) 2,4,6-Tribromophenol	10.39	330	62949	61.60	ng	-0.01
44) 2-Fluorobiphenyl	8.94	172	375155	41.70	ng	-0.02
78) Terphenyl-d14	12.66	244	417296	44.30	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	7.89	128	593481	41.05	ng	98
37) 2-Methylnaphthalene	8.57	142	420629	46.06	ng	100
45) 1,1'-Biphenyl	9.04	154	45030	4.34	ng	97
49) Dimethylphthalate	9.34	163	26234	2.70	ng	# 94
51) Acenaphthene	9.64	154	20955	2.35	ng	97
57) Fluorene	10.15	166	33499	3.64	ng	# 96
70) Phenanthrene	11.10	178	58645	4.36	ng	98
72) Carbazole	11.32	167	215812	17.11	ng	99

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

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