

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122315\
 Data File : BM003727.D
 Acq On : 23 Dec 2015 02:43
 Operator : UM/SJ
 Sample : G4912-08
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 STOCKPILE-2(0-2)

Quant Time: Dec 23 04:58:05 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM120915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 09 18:52:47 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	41024	20.00	ng	-0.03
21) Naphthalene-d8	10.48	136	160131	20.00	ng	-0.04
38) Acenaphthene-d10	14.34	164	80451	20.00	ng	-0.03
63) Phenanthrene-d10	17.09	188	166329	20.00	ng	-0.03
75) Chrysene-d12	21.29	240	171921	20.00	ng	-0.03
86) Perylene-d12	23.53	264	176889	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.28	112	322972	139.97	ng	-0.02
7) Phenol-d6	6.87	99	378962	118.27	ng	-0.02
23) Nitrobenzene-d5	8.85	82	268524	103.97	ng	-0.03
41) 2,4,6-Tribromophenol	15.84	330	114992	143.12	ng	-0.03
44) 2-Fluorobiphenyl	12.96	172	587743	99.53	ng	-0.03
78) Terphenyl-d14	19.73	244	679060	93.15	ng	-0.03
Target Compounds						Qvalue
31) Naphthalene	10.53	128	255572	30.50	ng	100
37) 2-Methylnaphthalene	12.15	142	26279	4.57	ng	97
51) Acenaphthene	14.41	154	15348	3.03	ng	99
54) Dibenzofuran	14.74	168	24795	3.39	ng	94
57) Fluorene	15.40	166	20022	3.30	ng	98
70) Phenanthrene	17.13	178	58882	6.31	ng	99
72) Carbazole	17.50	167	19858	2.46	ng	99

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

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