

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042115\
 Data File : BG016603.D
 Acq On : 22 Apr 2015 4:31
 Operator : TP/IZ
 Sample : G1937-02
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TRC-MW2

Quant Time: Apr 22 05:28:05 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 22 02:10:07 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	57276	20.00	ng	0.00
21) Naphthalene-d8	10.42	136	239107	20.00	ng	0.00
38) Acenaphthene-d10	14.28	164	133465	20.00	ng	0.00
63) Phenanthrene-d10	17.02	188	295651	20.00	ng	0.00
75) Chrysene-d12	21.21	240	328409	20.00	ng	0.00
86) Perylene-d12	23.42	264	302289	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	169710	47.49	ng	0.00
7) Phenol-d6	6.82	99	151260	30.30	ng	0.00
23) Nitrobenzene-d5	8.79	82	299675	77.17	ng	0.00
41) 2,4,6-Tribromophenol	15.77	330	98733	108.92	ng	0.00
44) 2-Fluorobiphenyl	12.90	172	661857	82.13	ng	0.00
78) Terphenyl-d14	19.66	244	1010533	74.13	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	10.47	128	76524	6.13	ng	99
49) Dimethylphthalate	13.74	163	45435	4.60	ng	100
51) Acenaphthene	14.34	154	79437	9.53	ng	99
54) Dibenzofuran	14.68	168	24467	2.08	ng	100
57) Fluorene	15.33	166	35434	3.43	ng	99
70) Phenanthrene	17.06	178	35463	2.11	ng	99
72) Carbazole	17.43	167	200567	11.75	ng	97

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

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