

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111317\
 Data File : BG030999.D
 Acq On : 14 Nov 2017 7:50
 Operator : SJ/JU
 Sample : I6436-02
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 153RD-63RD-ST-DRUM

Quant Time: Nov 14 11:26:23 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 10 15:38:25 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	35156	20.00	ng	-0.01
21) Naphthalene-d8	10.95	136	134848	20.00	ng	-0.01
38) Acenaphthene-d10	14.76	164	86958	20.00	ng	-0.01
63) Phenanthrene-d10	17.50	188	252368	20.00	ng	-0.01
75) Chrysene-d12	21.77	240	400321	20.00	ng	-0.01
86) Perylene-d12	25.07	264	400733	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.70	112	302709	147.65	ng	0.00
7) Phenol-d6	7.31	99	338958	122.72	ng	0.00
23) Nitrobenzene-d5	9.31	82	261806	115.05	ng	-0.01
41) 2,4,6-Tribromophenol	16.25	330	302053	214.72	ng	0.00
44) 2-Fluorobiphenyl	13.38	172	762396	125.98	ng	-0.02
78) Terphenyl-d14	20.09	244	1891077	104.31	ng	0.00
Target Compounds						
31) Naphthalene	11.01	128	148613	22.419	ng	99
37) 2-Methylnaphthalene	12.61	142	19931	3.895	ng	98
51) Acenaphthene	14.82	154	12974	2.440	ng	92
54) Dibenzofuran	15.16	168	20621	2.434	ng	98
70) Phenanthrene	17.55	178	33088	2.518	ng	97

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

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