

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111317\
 Data File : BG030978.D
 Acq On : 13 Nov 2017 18:17
 Operator : SJ/JU
 Sample : I6301-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JC-02-110917-B

Quant Time: Nov 14 05:47:46 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	42393	20.00	ng	-0.01
21) Naphthalene-d8	10.96	136	172209	20.00	ng	-0.01
38) Acenaphthene-d10	14.76	164	112210	20.00	ng	-0.02
63) Phenanthrene-d10	17.50	188	319083	20.00	ng	-0.01
75) Chrysene-d12	21.77	240	474135	20.00	ng	-0.02
86) Perylene-d12	25.06	264	465244	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.70	112	333951	135.08	ng	0.00
7) Phenol-d6	7.31	99	372052	111.71	ng	0.00
23) Nitrobenzene-d5	9.31	82	286749	98.67	ng	-0.01
41) 2,4,6-Tribromophenol	16.24	330	325124	179.11	ng	-0.01
44) 2-Fluorobiphenyl	13.39	172	821685	105.22	ng	-0.01
78) Terphenyl-d14	20.09	244	1923075	89.56	ng	0.00
Target Compounds						
31) Naphthalene	11.01	128	42769	5.052	ng	97
51) Acenaphthene	14.83	154	16991	2.476	ng	91
54) Dibenzofuran	15.16	168	22098	2.022	ng	91
57) Fluorene	15.81	166	18178	2.048	ng	# 99
70) Phenanthrene	17.54	178	45571	2.743	ng	98

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

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111317\
Data File : BG030978.D
Acq On : 13 Nov 2017 18:17
Operator : SJ/JU
Sample : I6301-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
JC-02-110917-B

Quant Time: Nov 14 05:47:46 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

