

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050118\
 Data File : BG034094.D
 Acq On : 2 May 2018 00:57
 Operator : SJ/JU
 Sample : J2133-03
 Misc : 8PPM MDL
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MDL-SOIL-03-QT2-2018

Quant Time: May 02 07:10:45 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 01 19:08:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	57486	20.00	ng	0.00
21) Naphthalene-d8	10.89	136	246855	20.00	ng	0.00
38) Acenaphthene-d10	14.72	164	190484	20.00	ng	0.00
63) Phenanthrene-d10	17.47	188	519483	20.00	ng	0.00
75) Chrysene-d12	21.76	240	584008	20.00	ng	-0.01
86) Perylene-d12	25.05	264	581210	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	480506	135.04	ng	0.00
7) Phenol-d6	7.22	99	721260	130.21	ng	0.00
23) Nitrobenzene-d5	9.24	82	532163	82.71	ng	0.00
41) 2,4,6-Tribromophenol	16.21	330	357524	135.55	ng	0.00
44) 2-Fluorobiphenyl	13.34	172	1121015	85.75	ng	0.00
78) Terphenyl-d14	20.09	244	2202133	82.60	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	12.89	237	31803	7.282	ng	87
53) 2,4-Dinitrophenol	14.81	184	9231	7.438	ng	# 49
69) Pentachlorophenol	17.11	266	31118	6.819	ng	91

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG050118\
Data File : BG034094.D
Acq On : 2 May 2018 00:57
Operator : SJ/JU
Sample : J2133-03
Misc : 8PPM MDL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MDL-SOIL-03-QT2-2018

Quant Time: May 02 07:10:45 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue May 01 19:08:03 2018
Response via : Initial Calibration

