

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072418\
 Data File : BG035846.D
 Acq On : 24 Jul 2018 14:06
 Operator : JU/SJ
 Sample : J3762-01
 Misc : LOD 8PPM
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOD-MDL-SOIL-01-QT3-2018

Quant Time: Jul 25 02:11:04 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 24 14:48:00 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	39121	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	147627	20.00	ng	0.00
38) Acenaphthene-d10	14.65	164	101201	20.00	ng	0.00
63) Phenanthrene-d10	17.39	188	280781	20.00	ng	0.00
75) Chrysene-d12	21.66	240	294012	20.00	ng	0.00
86) Perylene-d12	24.85	264	286517	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	304302	129.14	ng	0.00
7) Phenol-d6	7.19	99	436275	124.10	ng	0.00
23) Nitrobenzene-d5	9.19	82	291836	82.58	ng	0.00
41) 2,4,6-Tribromophenol	16.14	330	188400	141.24	ng	0.00
44) 2-Fluorobiphenyl	13.27	172	678207	89.78	ng	0.00
78) Terphenyl-d14	20.00	244	1371042	102.79	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	12.82	237	13350	6.664	ng	95
53) 2,4-Dinitrophenol	14.78	184	1361	7.927	ng	# 45
69) Pentachlorophenol	17.05	266	5968	2.921	ng	95

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

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