

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG062215\
 Data File : BG017748.D
 Acq On : 22 Jun 2015 16:52
 Operator : TP/IZ
 Sample : G2653-03
 Misc : LOD-SVOC-SOIL-2PPM
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOD-SOIL2015-01

Quant Time: Jun 23 03:50:41 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG061815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 19 06:05:07 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	78961	20.00	ng	0.00
21) Naphthalene-d8	10.39	136	340690	20.00	ng	0.00
38) Acenaphthene-d10	14.27	164	200860	20.00	ng	0.00
63) Phenanthrene-d10	17.01	188	456029	20.00	ng	-0.02
75) Chrysene-d12	21.22	240	485621	20.00	ng	-0.02
86) Perylene-d12	23.44	264	452611	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	562742	125.07	ng	0.00
7) Phenol-d6	6.79	99	763154	120.67	ng	0.00
23) Nitrobenzene-d5	8.76	82	467301	78.90	ng	-0.02
41) 2,4,6-Tribromophenol	15.76	330	204023	109.54	ng	-0.02
44) 2-Fluorobiphenyl	12.88	172	936132	75.69	ng	-0.02
78) Terphenyl-d14	19.66	244	1504279	70.39	ng	-0.02
Target Compounds						Qvalue
3) Pyridine	3.56	79	6981	1.28	ng	# 61
10) Phenol	6.81	94	11231	1.63	ng	85
35) Caprolactam	11.30	113	2462	0.91	ng	# 76
40) Hexachlorocyclopentadiene	12.41	237	3561	0.98	ng	# 75
53) 2,4-Dinitrophenol	14.37	184	750	7.65	ng	# 67
64) 4,6-Dinitro-2-methylphenol	15.39	198	1668	7.76	ng	86
76) Benzidine	19.28	184	4801	0.30	ng	95

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

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