

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122116\
 Data File : BG025373.D
 Acq On : 22 Dec 2016 1:18
 Operator : UM/SJ
 Sample : H6103-01
 Misc : LOD 5 PPM
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOD-SOIL-01-QT1-2017

Manual Integrations
 APPROVED

sohil
 12/22/2016 7:04:29 PM

Quant Time: Dec 22 05:08:33 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 21 18:42:01 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.21	152	48243	20.00	ng	0.00
21) Naphthalene-d8	11.04	136	190782	20.00	ng	0.00
38) Acenaphthene-d10	14.84	164	154402	20.00	ng	0.00
63) Phenanthrene-d10	17.59	188	355015	20.00	ng	0.00
75) Chrysene-d12	21.88	240	547188	20.00	ng	0.00
86) Perylene-d12	25.28	264	569327	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.75	112	381886	143.88	ng	0.00
7) Phenol-d6	7.37	99	523832	140.14	ng	0.00
23) Nitrobenzene-d5	9.40	82	392484	90.88	ng	0.00
41) 2,4,6-Tribromophenol	16.33	330	322467	129.72	ng	0.00
44) 2-Fluorobiphenyl	13.46	172	880605	92.34	ng	0.00
78) Terphenyl-d14	20.17	244	1786430	86.56	ng	0.00
Target Compounds						
32) Benzoic acid	10.39	122	7447m	3.11	ng	81
35) Caprolactam	11.98	113	4055m	3.39	ng	
55) 4-Nitrophenol	15.11	139	5672m	3.21	ng	
64) 4,6-Dinitro-2-methylphenol	16.01	198	6285	2.56	ng	
69) Pentachlorophenol	17.25	266	7986m	3.06	ng	

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

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