

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG111517\
 Data File : BG031028.D
 Acq On : 16 Nov 2017 1:17
 Operator : SJ/JU
 Sample : I6481-03
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SCMUA-SLUDGE-CAKE

Quant Time: Nov 16 08:15:23 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	36706	20.00	ng	-0.01
21) Naphthalene-d8	10.96	136	136923	20.00	ng	-0.01
38) Acenaphthene-d10	14.76	164	98381	20.00	ng	-0.02
63) Phenanthrene-d10	17.50	188	281731	20.00	ng	-0.02
75) Chrysene-d12	21.77	240	399198	20.00	ng	-0.02
86) Perylene-d12	25.06	264	395035	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.70	112	103856	48.52	ng	0.00
7) Phenol-d6	7.31	99	121038	41.97	ng	0.00
23) Nitrobenzene-d5	9.31	82	90686	39.25	ng	-0.02
41) 2,4,6-Tribromophenol	16.24	330	107438	67.51	ng	-0.01
44) 2-Fluorobiphenyl	13.38	172	316247	46.19	ng	-0.02
78) Terphenyl-d14	20.09	244	746268	41.28	ng	-0.01
Target Compounds						
10) Phenol	7.33	94	14637	4.887	ng	90
20) 3+4-Methylphenols	8.91	107	38733	13.061	ng	95

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

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