

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110116\
 Data File : BG024615.D
 Acq On : 2 Nov 2016 6:22
 Operator : UM/SJ
 Sample : H5489-03
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-69-0.5-1.0

Quant Time: Nov 02 07:20:24 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG102516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 26 11:14:13 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.27	152	119806	20.00	ng	0.00
21) Naphthalene-d8	11.09	136	475831	20.00	ng	0.00
38) Acenaphthene-d10	14.88	164	362476	20.00	ng	0.00
63) Phenanthrene-d10	17.63	188	800399	20.00	ng	0.00
75) Chrysene-d12	21.92	240	1101492	20.00	ng	-0.01
86) Perylene-d12	25.35	264	1148426	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.80	112	862439	117.98	ng	0.00
7) Phenol-d6	7.40	99	1501271	149.72	ng	0.00
23) Nitrobenzene-d5	9.44	82	1107609	105.98	ng	0.00
41) 2,4,6-Tribromophenol	16.36	330	404523	74.70	ng	0.00
44) 2-Fluorobiphenyl	13.51	172	2113249	96.90	ng	0.00
78) Terphenyl-d14	20.21	244	3410029	92.51	ng	0.00
Target Compounds						
10) Phenol	7.43	94	24549	2.38	ng	84
49) Dimethylphthalate	14.32	163	644955	25.06	ng	99

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

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