

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111117\
 Data File : BG030945.D
 Acq On : 11 Nov 2017 15:53
 Operator : SJ/JU
 Sample : I6389-11
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 3N-16

Quant Time: Nov 13 04:12:58 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	37043	20.00	ng	-0.01
21) Naphthalene-d8	10.96	136	160995	20.00	ng	0.00
38) Acenaphthene-d10	14.76	164	111202	20.00	ng	-0.01
63) Phenanthrene-d10	17.50	188	301274	20.00	ng	-0.01
75) Chrysene-d12	21.77	240	413051	20.00	ng	-0.01
86) Perylene-d12	25.07	264	443364	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.70	112	224595	103.97	ng	0.00
7) Phenol-d6	7.31	99	283686	97.48	ng	0.00
23) Nitrobenzene-d5	9.31	82	162467	59.80	ng	-0.01
41) 2,4,6-Tribromophenol	16.25	330	155355	86.36	ng	0.00
44) 2-Fluorobiphenyl	13.39	172	455085	58.80	ng	-0.01
78) Terphenyl-d14	20.09	244	933631	49.91	ng	0.00
Target Compounds						
10) Phenol	7.33	94	12171	4.027	ng	91
49) Dimethylphthalate	14.21	163	40189	4.180	ng	98

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

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