

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG110117\
 Data File : BG030648.D
 Acq On : 1 Nov 2017 23:54
 Operator : SJ/JU
 Sample : I6204-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 102617-TIDO-E-D-S

Quant Time: Nov 02 06:55:14 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 17:41:24 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	64733	20.00	ng	0.00
21) Naphthalene-d8	10.97	136	301476	20.00	ng	0.00
38) Acenaphthene-d10	14.77	164	207682	20.00	ng	0.00
63) Phenanthrene-d10	17.51	188	471862	20.00	ng	0.00
75) Chrysene-d12	21.78	240	511509	20.00	ng	0.00
86) Perylene-d12	25.09	264	511873	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.72	112	224401	57.91	ng	0.00
7) Phenol-d6	7.31	99	201124	36.98	ng	0.00
23) Nitrobenzene-d5	9.33	82	510336	107.57	ng	0.00
41) 2,4,6-Tribromophenol	16.26	330	272975	101.80	ng	0.00
44) 2-Fluorobiphenyl	13.40	172	1414791	99.91	ng	0.00
78) Terphenyl-d14	20.11	244	2197280	97.72	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	14.21	163	34876	2.152	ng	99

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

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