

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG110117\
 Data File : BG030649.D
 Acq On : 2 Nov 2017 00:31
 Operator : SJ/JU
 Sample : I6204-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 102617-TIDO-E-U-B

Quant Time: Nov 02 06:57:02 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	68919	20.00	ng	0.00
21) Naphthalene-d8	10.97	136	325521	20.00	ng	0.00
38) Acenaphthene-d10	14.77	164	225815	20.00	ng	0.00
63) Phenanthrene-d10	17.51	188	508529	20.00	ng	0.00
75) Chrysene-d12	21.78	240	545764	20.00	ng	0.00
86) Perylene-d12	25.09	264	547726	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.71	112	242631	58.81	ng	0.00
7) Phenol-d6	7.31	99	241695	41.74	ng	0.00
23) Nitrobenzene-d5	9.32	82	540647	105.54	ng	0.00
41) 2,4,6-Tribromophenol	16.25	330	301679	103.47	ng	0.00
44) 2-Fluorobiphenyl	13.40	172	1520337	98.74	ng	0.00
78) Terphenyl-d14	20.10	244	2307651	96.19	ng	0.00
Target Compounds						
10) Phenol	7.34	94	16891	2.706	ng	92
49) Dimethylphthalate	14.22	163	46874	2.660	ng	97

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

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