

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG110117\
 Data File : BG030652.D
 Acq On : 2 Nov 2017 2:23
 Operator : SJ/JU
 Sample : I6205-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 102517-TIDO-F-D-S

Quant Time: Nov 02 07:02:41 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	64128	20.00	ng	0.00
21) Naphthalene-d8	10.97	136	304033	20.00	ng	0.00
38) Acenaphthene-d10	14.77	164	211529	20.00	ng	0.00
63) Phenanthrene-d10	17.51	188	480053	20.00	ng	0.00
75) Chrysene-d12	21.79	240	529301	20.00	ng	0.00
86) Perylene-d12	25.09	264	528893	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.72	112	207144	53.96	ng	0.00
7) Phenol-d6	7.31	99	193374	35.89	ng	0.00
23) Nitrobenzene-d5	9.32	82	529279	110.63	ng	0.00
41) 2,4,6-Tribromophenol	16.26	330	288660	105.70	ng	0.00
44) 2-Fluorobiphenyl	13.40	172	1474034	102.20	ng	0.00
78) Terphenyl-d14	20.11	244	2289905	98.42	ng	0.00
Target Compounds						
10) Phenol	7.34	94	15503	2.669	ng	94
49) Dimethylphthalate	14.22	163	57680	3.495	ng	98

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

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