

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111117\
 Data File : BF100477.D
 Acq On : 12 Nov 2017 8:48
 Operator : SJ/JU
 Sample : I6402-04
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 110717-TIDO-F-U-S

Quant Time: Nov 13 06:12:13 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 10 15:44:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	134541	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	493302	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	200978	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	305972	20.00	ng	0.00
75) Chrysene-d12	14.06	240	270569	20.00	ng	0.00
86) Perylene-d12	15.54	264	192840	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	374343	44.60	ng	0.00
7) Phenol-d6	6.52	99	288018	27.83	ng	-0.01
23) Nitrobenzene-d5	7.47	82	807433	108.21	ng	0.00
41) 2,4,6-Tribromophenol	10.73	330	163635	79.90	ng	0.00
44) 2-Fluorobiphenyl	9.26	172	1367164	103.58	ng	0.00
78) Terphenyl-d14	13.01	244	938642	79.71	ng	0.00
Target Compounds						
10) Phenol	6.53	94	32578	2.998	ng	98
49) Dimethylphthalate	9.65	163	79119	5.156	ng	99

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

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