

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111217\
 Data File : BF100494.D
 Acq On : 12 Nov 2017 17:13
 Operator : SJ/JU
 Sample : I6403-07
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 110817-TIDO-F-U-B

Quant Time: Nov 13 06:51:58 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	143439	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	527618	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	213728	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	323742	20.00	ng	0.00
75) Chrysene-d12	14.06	240	289212	20.00	ng	0.00
86) Perylene-d12	15.54	264	204501	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	270968	30.28	ng	0.00
7) Phenol-d6	6.51	99	204619	18.54	ng	-0.02
23) Nitrobenzene-d5	7.46	82	605998	75.93	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	117925	54.15	ng	-0.01
44) 2-Fluorobiphenyl	9.26	172	1079500	76.91	ng	0.00
78) Terphenyl-d14	13.01	244	730564	58.04	ng	0.00
Target Compounds						
10) Phenol	6.53	94	25462	2.198	ng	96
49) Dimethylphthalate	9.65	163	90319	5.535	ng	98

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

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