

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111217\
 Data File : BF100493.D
 Acq On : 12 Nov 2017 16:46
 Operator : SJ/JU
 Sample : I6403-06
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 110817-TIDO-F-D-S

Quant Time: Nov 13 06:50:18 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	138702	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	515850	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	218890	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	342296	20.00	ng	0.00
75) Chrysene-d12	14.06	240	289138	20.00	ng	0.00
86) Perylene-d12	15.54	264	201679	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	339361	39.22	ng	0.00
7) Phenol-d6	6.52	99	250777	23.50	ng	-0.01
23) Nitrobenzene-d5	7.46	82	767712	98.39	ng	0.00
41) 2,4,6-Tribromophenol	10.73	330	165223	74.07	ng	0.00
44) 2-Fluorobiphenyl	9.26	172	1334488	92.83	ng	0.00
78) Terphenyl-d14	13.01	244	918404	72.98	ng	0.00
Target Compounds						
31) Naphthalene	8.20	128	55626	2.239	ng	97
49) Dimethylphthalate	9.65	163	77484	4.636	ng	99
51) Acenaphthene	9.97	154	29337	2.119	ng	99
70) Phenanthrene	11.45	178	52907	2.936	ng	97

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

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