

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\
 Data File : BF110742.D
 Acq On : 14 Nov 2018 1:51
 Operator : JU/SJ
 Sample : J5925-17 20X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-5-GW

Manual Integrations
 APPROVED

Sohil
 11/14/2018 1:52:32 PM

Quant Time: Nov 14 07:08:58 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 12:43:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	71496	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	261187	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	101096	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	168768	20.00	ng	0.00
76) Chrysene-d12	14.13	240	156631	20.00	ng	0.00
87) Perylene-d12	15.66	264	101239	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	8884	2.02	ng	0.01
7) Phenol-d6	6.58	99	7422	1.32	ng	0.00
23) Nitrobenzene-d5	7.51	82	13367	2.95	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	3752	3.68	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	29613	4.18	ng	0.00
79) Terphenyl-d14	13.06	244	24257	2.90	ng	-0.01
Target Compounds						
37) 2-Methylnaphthalene	8.95	142	477380	59.391	ng	99
38) 1-Methylnaphthalene	9.05	142	344511m	44.567	ng	
52) Acenaphthene	10.02	154	21565	3.487	ng	# 55
58) Fluorene	10.54	166	47779	6.876	ng	# 82
71) Phenanthrene	11.51	178	125833	13.917	ng	96

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

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