

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110918\
 Data File : BF110646.D
 Acq On : 9 Nov 2018 23:52
 Operator : JU/SJ
 Sample : J5884-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3-S

Quant Time: Nov 11 04:01:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 09 15:13:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	80630	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	324449	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	141312	20.00	ng	0.00
64) Phenanthrene-d10	11.48	188	213332	20.00	ng	0.00
76) Chrysene-d12	14.13	240	145850	20.00	ng	-0.01
87) Perylene-d12	15.66	264	115253	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	693552	139.72	ng	0.01
7) Phenol-d6	6.57	99	898510	141.55	ng	0.00
23) Nitrobenzene-d5	7.51	82	558926	99.21	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	170723	119.90	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	868595	87.79	ng	0.00
79) Terphenyl-d14	13.06	244	564746	72.52	ng	0.00
Target Compounds						
10) Phenol	6.59	94	57254	7.779	ng	# 65
50) Dimethylphthalate	9.69	163	69004	6.544	ng	# 99
52) Acenaphthene	10.02	154	48612	5.624	ng	99
58) Fluorene	10.53	166	33907	3.491	ng	98

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

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