

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120419\
 Data File : BF118101.D
 Acq On : 4 Dec 2019 22:11
 Operator : JU
 Sample : K6098-06
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OWL-C7-GW

Quant Time: Dec 05 03:03:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 03 15:16:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	176654	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	629611	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	337921	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	646187	20.00	ng	0.00
76) Chrysene-d12	14.07	240	415295	20.00	ng	0.00
87) Perylene-d12	15.57	264	439557	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1070675	107.28	ng	0.01
7) Phenol-d6	6.50	99	920371	76.46	ng	0.00
23) Nitrobenzene-d5	7.44	82	981733	92.69	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	616803	172.41	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	2014367	106.94	ng	0.00
79) Terphenyl-d14	13.02	244	2530631	111.37	ng	0.00
Target Compounds						
10) Phenol	6.52	94	27237	2.177	ng	95
50) Dimethylphthalate	9.64	163	354633	14.966	ng	99
52) Acenaphthene	9.96	154	49744	2.582	ng	100

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

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