

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121318\
 Data File : BF111369.D
 Acq On : 13 Dec 2018 14:29
 Operator : JU/SJ
 Sample : J6334-10DL 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GW-121018DL

Quant Time: Dec 13 16:30:55 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 13:16:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	177923	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	733656	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	393466	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	751040	20.00	ng	0.00
76) Chrysene-d12	13.94	240	576769	20.00	ng	-0.01
87) Perylene-d12	15.39	264	454121	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	343311	32.63	ng	0.00
7) Phenol-d6	6.42	99	299029	21.08	ng	-0.01
23) Nitrobenzene-d5	7.35	82	608405	46.74	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	252133	72.69	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1180344	48.16	ng	0.00
79) Terphenyl-d14	12.90	244	1377863	52.58	ng	0.00
Target Compounds						
31) Naphthalene	8.10	128	1536041	44.938	ng	98
37) 2-Methylnaphthalene	8.79	142	503908	22.805	ng	100
38) 1-Methylnaphthalene	8.89	142	302843	14.201	ng	97

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

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