

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101618\
 Data File : BE097945.D
 Acq On : 17 Oct 2018 13:22
 Operator : SJ/JU
 Sample : J5317-07DL 2X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 RE-122D2-IMP-20181004DL

Quant Time: Oct 17 17:20:16 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE101618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 16 15:36:44 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	1923	0.40	ng	0.00
7) Naphthalene-d8	10.69	136	8337	0.40	ng	0.00
13) Acenaphthene-d10	14.54	164	5375	0.40	ng	-0.02
19) Phenanthrene-d10	17.29	188	15062	0.40	ng	-0.01
27) Chrysene-d12	21.47	240	16084	0.40	ng	-0.01
34) Perylene-d12	24.03	264	16785	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.47	112	565	0.09	ng	0.00
5) Phenol-d6	7.06	99	510	0.06	ng	0.00
8) Nitrobenzene-d5	9.04	82	1419	0.18	ng	0.00
11) 2-Methylnaphthalene-d10	12.28	152	2543	0.18	ng	-0.02
14) 2,4,6-Tribromophenol	16.04	330	342	0.12	ng	0.00
15) 2-Fluorobiphenyl	13.17	172	5088	0.23	ng	0.00
25) Fluoranthene-d10	19.33	212	35672	0.17	ng	0.00
29) Terphenyl-d14	19.92	244	9076	0.25	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.38	88	8377	2.498	ng	Qvalue 96
6) bis(2-Chloroethyl)ether	7.30	93	153	0.020	ng	# 84

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

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