

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101618\
 Data File : BE097936.D
 Acq On : 17 Oct 2018 7:51
 Operator : SJ/JU
 Sample : J5317-10
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 RE-108D2-20181004

Quant Time: Oct 17 08:56:38 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	7.88	152	1760	0.40	ng	0.00
7) Naphthalene-d8	10.68	136	7316	0.40	ng	-0.01
13) Acenaphthene-d10	14.54	164	4423	0.40	ng	-0.01
19) Phenanthrene-d10	17.29	188	12491	0.40	ng	-0.01
27) Chrysene-d12	21.47	240	17283	0.40	ng	-0.01
34) Perylene-d12	24.03	264	17566	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.47	112	919	0.17	ng	0.00
5) Phenol-d6	7.04	99	857	0.10	ng	-0.01
8) Nitrobenzene-d5	9.03	82	2693	0.39	ng	-0.01
11) 2-Methylnaphthalene-d10	12.28	152	4555	0.37	ng	-0.01
14) 2,4,6-Tribromophenol	16.04	330	819	0.34	ng	0.00
15) 2-Fluorobiphenyl	13.17	172	8766	0.48	ng	0.00
25) Fluoranthene-d10	19.33	212	71702	0.42	ng	0.00
29) Terphenyl-d14	19.92	244	20125	0.51	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.38	88	10951	3.569	ng	95
6) bis(2-Chloroethyl)ether	7.30	93	793	0.114	ng	99
9) Naphthalene	10.73	128	2061	0.028	ng	# 90

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

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