

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101618\
 Data File : BE097922.D
 Acq On : 16 Oct 2018 22:33
 Operator : SJ/JU
 Sample : J5393-09
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 RE-137-745FT-2018109

Quant Time: Oct 17 08:09:51 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	1773	0.40	ng	0.00
7) Naphthalene-d8	10.69	136	7776	0.40	ng	0.00
13) Acenaphthene-d10	14.56	164	5582	0.40	ng	0.00
19) Phenanthrene-d10	17.29	188	18128	0.40	ng	-0.01
27) Chrysene-d12	21.49	240	27570	0.40	ng	0.00
34) Perylene-d12	24.04	264	15658	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.47	112	955	0.17	ng	0.00
5) Phenol-d6	7.06	99	929	0.11	ng	0.00
8) Nitrobenzene-d5	9.04	82	2892	0.39	ng	0.00
11) 2-Methylnaphthalene-d10	12.28	152	5372	0.42	ng	-0.02
14) 2,4,6-Tribromophenol	16.04	330	1187	0.40	ng	0.00
15) 2-Fluorobiphenyl	13.17	172	9165	0.40	ng	0.00
25) Fluoranthene-d10	19.33	212	115740	0.47	ng	0.00
29) Terphenyl-d14	19.92	244	36936	0.58	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.38	88	16981	5.493	ng	95
6) bis(2-Chloroethyl)ether	7.30	93	404	0.058	ng	94

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

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