

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101618\
 Data File : BE097940.D
 Acq On : 17 Oct 2018 10:17
 Operator : SJ/JU
 Sample : J5393-09DL 2X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

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

Quant Time: Oct 17 16:59:32 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	2951	0.40	ng	-0.01
7) Naphthalene-d8	10.68	136	13018	0.40	ng	-0.01
13) Acenaphthene-d10	14.54	164	9019	0.40	ng	-0.02
19) Phenanthrene-d10	17.29	188	26770	0.40	ng	-0.01
27) Chrysene-d12	21.47	240	27676	0.40	ng	-0.01
34) Perylene-d12	24.03	264	21593	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.46	112	540	0.06	ng	-0.01
5) Phenol-d6	7.05	99	612	0.04	ng	-0.01
8) Nitrobenzene-d5	9.04	82	1814	0.15	ng	0.00
11) 2-Methylnaphthalene-d10	12.28	152	3255	0.15	ng	-0.02
14) 2,4,6-Tribromophenol	16.04	330	612	0.13	ng	0.00
15) 2-Fluorobiphenyl	13.17	172	5566	0.15	ng	0.00
25) Fluoranthene-d10	19.33	212	54240	0.15	ng	0.00
29) Terphenyl-d14	19.92	244	16554	0.26	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.38	88	10959	2.130	ng	Qvalue 93
6) bis(2-Chloroethyl)ether	7.30	93	269	0.023	ng	# 88

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

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