

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101618\
 Data File : BE097939.D
 Acq On : 17 Oct 2018 9:41
 Operator : SJ/JU
 Sample : J5317-13
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 RE-109D3-20181005

Quant Time: Oct 17 16:57:50 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	2049	0.40	ng	-0.01
7) Naphthalene-d8	10.68	136	8717	0.40	ng	-0.01
13) Acenaphthene-d10	14.54	164	5590	0.40	ng	-0.01
19) Phenanthrene-d10	17.29	188	14910	0.40	ng	-0.01
27) Chrysene-d12	21.47	240	18303	0.40	ng	-0.01
34) Perylene-d12	24.03	264	18950	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.47	112	1261	0.20	ng	0.00
5) Phenol-d6	7.05	99	1087	0.11	ng	-0.01
8) Nitrobenzene-d5	9.03	82	3508	0.42	ng	-0.01
11) 2-Methylnaphthalene-d10	12.28	152	5813	0.40	ng	-0.01
14) 2,4,6-Tribromophenol	16.05	330	1238	0.41	ng	0.00
15) 2-Fluorobiphenyl	13.17	172	10062	0.44	ng	0.00
25) Fluoranthene-d10	19.33	212	84421	0.42	ng	0.00
29) Terphenyl-d14	19.92	244	32865	0.78	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.38	88	13789	3.860	ng	91
6) bis(2-Chloroethyl)ether	7.30	93	262	0.032	ng	# 87

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

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