

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101618\
 Data File : BE097929.D
 Acq On : 17 Oct 2018 2:52
 Operator : SJ/JU
 Sample : J5317-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 RE-122D2-20181004

Quant Time: Oct 17 08:35:56 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	1846	0.40	ng	0.00
7) Naphthalene-d8	10.69	136	8065	0.40	ng	0.00
13) Acenaphthene-d10	14.54	164	5395	0.40	ng	-0.01
19) Phenanthrene-d10	17.29	188	17419	0.40	ng	-0.01
27) Chrysene-d12	21.48	240	22907	0.40	ng	-0.01
34) Perylene-d12	24.03	264	18178	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.47	112	1225	0.21	ng	0.00
5) Phenol-d6	7.06	99	1041	0.12	ng	0.00
8) Nitrobenzene-d5	9.03	82	3027	0.40	ng	-0.01
11) 2-Methylnaphthalene-d10	12.28	152	5348	0.40	ng	-0.01
14) 2,4,6-Tribromophenol	16.04	330	1160	0.40	ng	0.00
15) 2-Fluorobiphenyl	13.17	172	10241	0.46	ng	0.00
25) Fluoranthene-d10	19.33	212	111151	0.47	ng	0.00
29) Terphenyl-d14	19.92	244	30859	0.59	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.38	88	16925	5.259	ng	# 90
6) bis(2-Chloroethyl)ether	7.30	93	440	0.060	ng	# 86

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

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