

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101618\
 Data File : BE097938.D
 Acq On : 17 Oct 2018 9:02
 Operator : SJ/JU
 Sample : J5317-12
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 TT-DUP04-20181004

Quant Time: Oct 17 16:54:55 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	1970	0.40	ng	0.00
7) Naphthalene-d8	10.68	136	8671	0.40	ng	-0.01
13) Acenaphthene-d10	14.54	164	5985	0.40	ng	-0.02
19) Phenanthrene-d10	17.29	188	18689	0.40	ng	-0.01
27) Chrysene-d12	21.48	240	20594	0.40	ng	-0.01
34) Perylene-d12	24.03	264	17237	0.40	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.47	112	1220	0.20	ng	0.00
5) Phenol-d6	7.05	99	1112	0.12	ng	0.00
8) Nitrobenzene-d5	9.03	82	3628	0.44	ng	-0.01
11) 2-Methylnaphthalene-d10	12.28	152	6160	0.43	ng	-0.02
14) 2,4,6-Tribromophenol	16.05	330	1403	0.44	ng	0.00
15) 2-Fluorobiphenyl	13.17	172	11555	0.47	ng	0.00
25) Fluoranthene-d10	19.33	212	117452	0.46	ng	0.00
29) Terphenyl-d14	19.92	244	44238	0.94	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.38	88	12481	3.634	ng	96
6) bis(2-Chloroethyl)ether	7.30	93	999	0.128	ng	99
9) Naphthalene	10.73	128	2968	0.034	ng	# 93
12) 2-Methylnaphthalene	12.35	142	306	0.021	ng	# 89

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

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