

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101618\
 Data File : BE097916.D
 Acq On : 16 Oct 2018 18:52
 Operator : SJ/JU
 Sample : J5393-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 TT-DUP05-20181008

Quant Time: Oct 17 15:47:41 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	2001	0.40	ng	0.00
7) Naphthalene-d8	10.69	136	8276	0.40	ng	0.00
13) Acenaphthene-d10	14.56	164	5404	0.40	ng	0.00
19) Phenanthrene-d10	17.29	188	14447	0.40	ng	-0.01
27) Chrysene-d12	21.48	240	17162	0.40	ng	-0.01
34) Perylene-d12	24.04	264	17598	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.47	112	1189	0.19	ng	0.00
5) Phenol-d6	7.06	99	1008	0.11	ng	0.00
8) Nitrobenzene-d5	9.04	82	3133	0.40	ng	0.00
11) 2-Methylnaphthalene-d10	12.28	152	5407	0.39	ng	-0.01
14) 2,4,6-Tribromophenol	16.04	330	929	0.32	ng	0.00
15) 2-Fluorobiphenyl	13.18	172	8862	0.40	ng	0.00
25) Fluoranthene-d10	19.33	212	75730	0.39	ng	0.00
29) Terphenyl-d14	19.92	244	22594	0.57	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.38	88	10770	3.087	ng	93

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

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