

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102317\
 Data File : BF099767.D
 Acq On : 23 Oct 2017 19:03
 Operator : SJ/JU
 Sample : I5950-09
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20170722-C-A-COMP1

Quant Time: Oct 23 23:47:20 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 23 15:08:33 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	99564	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	412144	20.00	ng	0.00
38) Acenaphthene-d10	9.84	164	194096	20.00	ng	0.00
63) Phenanthrene-d10	11.33	188	308433	20.00	ng	0.00
75) Chrysene-d12	13.97	240	171739	20.00	ng	-0.01
86) Perylene-d12	15.44	264	189746	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	359023	56.61	ng	0.00
7) Phenol-d6	6.44	99	258419	34.37	ng	-0.01
23) Nitrobenzene-d5	7.37	82	554663	86.64	ng	0.00
41) 2,4,6-Tribromophenol	10.64	330	202107	114.26	ng	0.00
44) 2-Fluorobiphenyl	9.16	172	1109362	88.18	ng	0.00
78) Terphenyl-d14	12.92	244	765299	97.38	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.12	128	2427784	122.033	ng	99
37) 2-Methylnaphthalene	8.80	142	259595	19.471	ng	97
45) 1,1'-Biphenyl	9.26	154	61532	3.504	ng	96
51) Acenaphthene	9.87	154	142956	11.912	ng	98
54) Dibenzofuran	10.05	168	191842	11.325	ng	98
57) Fluorene	10.39	166	107909	7.694	ng	100
70) Phenanthrene	11.36	178	163416	9.388	ng	97

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

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