

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102317\
 Data File : BF099769.D
 Acq On : 23 Oct 2017 19:58
 Operator : SJ/JU
 Sample : I5950-11
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20170724-C-A-COMP3

Quant Time: Oct 23 23:59:04 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	103875	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	432433	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	190758	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	250923	20.00	ng	0.01
75) Chrysene-d12	13.98	240	176773	20.00	ng	0.00
86) Perylene-d12	15.44	264	191021	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	419242	63.36	ng	0.00
7) Phenol-d6	6.45	99	319802	40.76	ng	0.00
23) Nitrobenzene-d5	7.37	82	609173	90.69	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	197380	113.54	ng	0.00
44) 2-Fluorobiphenyl	9.16	172	1187320	96.03	ng	0.00
78) Terphenyl-d14	12.93	244	723258	89.41	ng	0.01
Target Compounds						
31) Naphthalene	8.11	128	1147040	54.951	ng	100
32) Benzoic acid	7.89	122	12795	2.545	ng	# 81
37) 2-Methylnaphthalene	8.80	142	137991	9.865	ng	98
51) Acenaphthene	9.87	154	40911	3.469	ng	100
54) Dibenzofuran	10.05	168	85043	5.108	ng	99
70) Phenanthrene	11.37	178	76472	5.400	ng	# 88

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

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