

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102317\
 Data File : BF099770.D
 Acq On : 23 Oct 2017 20:26
 Operator : SJ/JU
 Sample : I5950-12
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20170725-PA-A-COMP1

Manual Integrations
 APPROVED

Sohil
 10/24/2017 8:41:03 PM

Quant Time: Oct 24 00:08:22 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	99961	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	424656	20.00	ng	0.00
38) Acenaphthene-d10	9.84	164	205271	20.00	ng	0.00
63) Phenanthrene-d10	11.33	188	345115	20.00	ng	0.00
75) Chrysene-d12	13.97	240	180713	20.00	ng	-0.01
86) Perylene-d12	15.44	264	188226	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	426844	67.04	ng	0.00
7) Phenol-d6	6.44	99	322594	42.73	ng	-0.01
23) Nitrobenzene-d5	7.37	82	623321	94.50	ng	0.00
41) 2,4,6-Tribromophenol	10.64	330	237545	126.98	ng	0.00
44) 2-Fluorobiphenyl	9.16	172	1236765	92.96	ng	0.00
78) Terphenyl-d14	12.92	244	906935	109.68	ng	0.00
Target Compounds						
31) Naphthalene	8.10	128	433525	21.149	ng	99
32) Benzoic acid	7.88	122	11738	2.378	ng	87
37) 2-Methylnaphthalene	8.80	142	99216	7.223	ng	97
51) Acenaphthene	9.87	154	62122	4.895	ng	98
54) Dibenzofuran	10.05	168	97967	5.468	ng	98
57) Fluorene	10.39	166	59608	4.019	ng	98
70) Phenanthrene	11.36	178	97796	5.021	ng	98

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

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