

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100517\
 Data File : BF099108.D
 Acq On : 5 Oct 2017 17:30
 Operator : SJ/JU
 Sample : I5526-02DL 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 QNWP8-MW30D-GWDL

Quant Time: Oct 05 18:30:07 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 05 17:20:48 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	127207	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	469866	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	242953	20.00	ng	0.00
63) Phenanthrene-d10	11.39	188	422602	20.00	ng	0.00
75) Chrysene-d12	14.03	240	300994	20.00	ng	0.00
86) Perylene-d12	15.50	264	214947	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	38463	4.81	ng	0.00
7) Phenol-d6	6.48	99	30049	3.05	ng	-0.02
23) Nitrobenzene-d5	7.42	82	71392	9.74	ng	-0.01
41) 2,4,6-Tribromophenol	10.69	330	21497	10.81	ng	-0.01
44) 2-Fluorobiphenyl	9.22	172	152208	10.46	ng	-0.01
78) Terphenyl-d14	12.96	244	124566	9.41	ng	-0.01
Target Compounds						
31) Naphthalene	8.19	128	5982640	271.102	ng	98
37) 2-Methylnaphthalene	8.86	142	304834	21.249	ng	99
45) 1,1'-Biphenyl	9.32	154	113843	5.452	ng	97
51) Acenaphthene	9.93	154	256184	16.851	ng	99
54) Dibenzofuran	10.10	168	260029	12.846	ng	99
57) Fluorene	10.45	166	109408	6.725	ng	98
70) Phenanthrene	11.41	178	127979	5.658	ng	98
72) Carbazole	11.62	167	106702	4.708	ng	99

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

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